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Jean Claude Dutailly

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MATHEMATICS IN PHYSICS

Jean Claude Dutailly

29 june 2015

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Introduction

A century after the introduction of Relativity and Quantum Mechanics these theories are still the topic of hot controversies. If this is not a serious concern for most physicists, there is no doubt that this is the symptom of a weakness in the tools that they use daily, and is at the core of the lack of progress in Theoretical Physics. The attempts to solve the conundrum have been directed in two directions. Most of them are based on assuming new properties to the physical world, either additional dimensions, special behavior of particles of fields, alternate universe,... Others on the introduction of new formal systems, involving highly abstract concepts, to base the computational practices on firmer ground. In this book I propose another way : I keep the objects of conventional physics (particles, fields, space, time,...) and their properties, as almost all physicists understand them (so there is 4 dimensions, no special particle or field, no GTU,...), but I focus on their representation by mathematical objects. It is clear that we cannot do the Physics of the XXI^o century with the Mathematics of the XIX^o century - when Einstein introduced Relativity, Differential Geometry was in infancy - but conversely the race to the introduction of new mathematical tools is void if it does not go with a deep understanding of their adequation with the physical concepts that they are supposed to represent. It is then possible to find from where comes Quantum Mechanics, to show that its main axioms are actually theorems which can be proven, to introduce in a simple and elegant way General Relativity, spinors, and the like wave representation of particles, and to work with them and the usual principles of Physics. Moreover it provides a must needed understanding of the gravitational field, showing that it has rotational and transversal components which play different roles. We propose also an explanation of bosons as discontinuities of the fields.

So the first purpose of this book is to show how mathematical tools can be used to represent rigorously and efficiently some of the key concepts of modern Physics, in Relativity and Quantum Mechanics, and through this representation, show how these concepts, and physical laws which seem often paradoxical, can be clearly understood. So the book addresses both Mathematicians who want to use their knowledge to get an overview of some of the main issues of modern physics, and Physicists who want to improve their grasp of mathematical tools in their work. Its purpose is to help graduated and undergraduate students to understand the place and the role of mathematics in Physics. Why do we have

Hilbert spaces in Quantum Mechanics ? What is the meaning of probabilist representations ? How physical concepts such as motion, moment, matter or force fields, are related to mathematical objects ? How physical principles, such as the Principle of Relativity or the Principle of Least Action, can be translated into Mathematics and be efficiently used in Physics ? How can we understand the entanglement of particles ? What is a spinor ?

In this book the reader will see how to deal with manifolds, fiber bundles, connections, Clifford algebras, group representations, generalized functions or Lagrange equations. There are many books which deal with these topics, usually for physicists, with the purpose to make understandable in a nut shell what are, after all, some of the most abstract parts of Mathematics. We will not choose this path, not by some pedantic pretense, but because for a scientist the most general approach, which requires few but key concepts, is easier than a pragmatic one based upon the acceptance of many computational rules. So we will, from the beginning, introduce the mathematical tools, usually in their most general definition, into the representation of physical phenomena and show how their properties fit with what we can understand of these phenomena, and how they help to solve some classical problems. This will be illustrated by the building, step by step, of a formal model which incorporates all the bricks to show how they work. We will use many mathematical definitions or theorems. The most important will be recalled, and for the proofs and a more comprehensive understanding I refer to a previous book ("Mathematics for Theoretical Physics") which is freely available.

But Mathematics can offer only tools : to use them efficiently it is necessary to understand what they mean, and why and how they can be adapted to represent physical phenomena. To do so one needs a new, candid, look at these phenomena, just as Einstein did in his celebrated 1905 article : space and time are not necessarily how we are used to see them, and more than often one needs to pause before jumping to Mathematics. Actually the indiscriminate use of formal systems can be hazardous, and we will see in this books some examples of common mathematical representations, seen as granted because they are used over and over, which are not grounded in any legitimate conceptual frame, or even do not reflect the actual practices of the workers in natural sciences. Many physicists are indeed disturbed by the way Mathematics invade Theoretical Physics, with the continuous flow of more and more sophisticated mathematical theories, difficult to understand but for a few specialists, and which are estranged from experimental work. But most of them also comply with the mandatory, at least in the academic world, use of formal practices which seem to be against any logical justification, or are even non sensible, such as minimal substitution rules or divergent integral. Computation in Physics, as it is commonly done, has quite often the appearance of magic, justified only by the usual "experiments have proven that it works", or more blatantly "the great masters have said that it is right". Perhaps, but as a consequence it leaves the students with the feeling that Physics is more an exercise in rote knowledge than in rational understanding, and it opens the way to interpretations : if the formal system

is strange, it is perhaps that nature itself is strange. A quick Google search for “quantum mechanics interpretations” provides more than 5 millions links, and there are more than 50 elaborate theories, the multiuniverse having the largest support in the scientific community. So one cannot say that modern Physics answer clearly our questions about nature. And it is not true that experiments have proven the rightfulness of the common practices. The discrepancy between what the theories predict and what is observed is patched with the introduction of new concepts, whose physical realization is more and more difficult to check : collapse of the wave function, Higgs boson, dark matter, brown energy,...

This is not the purpose of this book to add another interpretation to the existing long list. There will be few assumptions about the physical world, clear stated, and they are well in line with what Physicists know and most Scientists agree upon. There will be no extra-dimensions, string theory, branes, supersymmetry,...Not that such theories should be discarded, or will be refuted, but only because they are not necessary to get a solid picture of the basic concepts in Physics. And indeed we do not answer to all questions in this book, some issues are still open, but I hope that their meaning will be clearer, leading the way to a better and stronger understanding of the real world.

The first chapter is devoted to a bit of philosophy. From many discussions with scientists I felt that it is appropriate. Because the book is centered on the relation between Mathematics and Physics, it is necessary to have a good understanding of what is meant by physical laws, theories, validation by experiments, models, representations,...Philosophy has a large scope, so it deals also with knowledge : epistemology helps us to sort out the different meanings of what we call knowledge, the status of Science and Mathematics, how the Sciences improve and theories are replaced by new ones. This chapter will not introduce any new Philosophy, just provide a summary of what scientists should know from the works of professional philosophers.

The second chapter is dedicated to Quantum Mechanics (QM). This is mandatory, because QM has dominated theoretical Physics for almost a century, with many disturbing and confusing issues. It is at the beginning of the book because, as we will see, actually QM is not a physical theory per se, it does not require any assumption about how Nature works. QM is a theory which deals with the way one represents the world : its axioms, which appear as physical laws, are actually mathematical theorems, which are the consequences of the use by Physicists of mathematical models to make their computations and collect their data from experiments. This is not surprising that measure has such a prominent place in QM : it is all about the measures, that is the image of the world that physicists build, and not about the world itself. And this is the first, and newest, example of how the use of Mathematics can be misleading.

The third chapter is dedicated to the Geometry of the Universe. By this we do not mean how the whole universe is, which is the topic of Cosmology. Cosmology is a branch of Physics of its own, which raises issues of an epistemological nature, and is, from my point of view, speculative, even if it is grounded in Astrophysics. We will only evoke some points of Cosmology in passing in this

book. By Geometry of the Universe I mean here the way we represent locations of points, components of vectors and tensors, and the consequences which follow for the rules in a change of representation. This will be done in the relativist framework, and more precisely in the framework of General Relativity. It is less known, seen usually as a difficult topic, but, as we will see, some of the basic concepts of Relativity are easier to understand when we quit the usual, and misleading, representations, and are not very complicated when one uses the right mathematical tools, notably fiber bundles.

The fourth chapter addresses Kinematics, which, by the concept of moment, is the gate between forces and geometry. Relativity requires a brand new vision of these concepts, which has been engaged, but neither fully or consistently. Rotation in particular has a different meaning in the 4 dimensional space than in the usual euclidean space, and a revision of rotational moment requires the introduction of a new framework. Spinors are not new in Physics, we will see what they mean, in Physics and in Mathematics, with Clifford algebras. This leads naturally to the introduction of the spin, which has a clear and simple interpretation, and to the representation of particles by fields of spinors, which incorporates in a single quantity the motion, translational and rotational, and the kinematics characteristics of material objects.

The fifth chapter addresses Force Fields. After a short reminder of the Standard Model we will see how charges of particles and force fields can be represented, with the concept of connections on fiber bundles. We will not deal with all the intricacies of the Standard Model, but focus on the principles and main mechanisms. The integration of Gravity, not in a Great Unification Theory, but with tools similar to the other forces and in parallel with them, opens a fresh vision on important issues in General Relativity. In particular it appears that the common and exclusive use of the Levi-Civita connection and scalar curvature introduces useless complications but, more importantly, misses important features of the gravitational field.

The sixth chapter is dedicated to lagrangians. They are the work horses of Theoretical Physics, and we will review the problems, physical and mathematical, that they involve, and how to deal with them. We will see why a lagrangian cannot incorporate explicitly some variables, and build a simple lagrangian with 6 variables, which can be used in most of the problems.

The seventh chapter is dedicated to continuous models. Continuous processes are not the rule in the physical world, but are the simplest to represent and understand. We will see how the material introduced in the previous chapters can be used, how the methods of Variational Calculus can be used in solving two models, for a field of particles and for a single particle. They help to understand the concepts of currents and Energy-Momentum tensors.

The eighth chapter is dedicated to discontinuous processes. They are common in the real world but their study is difficult. We will see how one can deal with discontinuities in Mathematics, and introduce a new interpretation of the bosons, the particle-like objects which carry the force in the Standard Model. And we end with a short survey of the principal tools used in discontinuous processes, notably path integrals and gas of interacting particles.

Chapter 1

WHAT IS SCIENCE ?

Science has acquired a unique status in our societies. It is seen by the laymen as the premier gate to the truth in this world, both feared and respected. Who could not be amazed by its technical prowess ? How many engineers, technicians, daily put their faith in its laws ? For many scientists their work has a distinctive quality, which puts them in another class than novelists, theologians, or artists. Even when dealing with some topics as government, traditions, religion,... they mark their territory by claiming the existence of Social Sciences, such as Economics, Sociology or Political Sciences, endowed with methods and procedures which stand them apart, and lest us say, above the others who engage in narratives on the same topics. But what are the bases for such pretense ? After all, many scientific assertions are controversial, when they impact our daily lives (from the climate warming to almost any drug), but not least in the scientific community itself. The latter is natural and even sound - controversy is consubstantial to science - however it has attained a more bitter tone in the last years, fueled by the fierce competition between its servants, but also by the frustrations of many scientists, mostly in Physics, at a scientifically correct corpus with too many loopholes. A common answer to the discontents is to refer them to the all powerful experimental proofs, but these are more and more difficult to reach and to interpret : how many people could sensibly discuss the discovery of the Higgs boson ?

To put some light on these issues, the natural way is to look towards Philosophy, and more precisely Epistemology, which is its branch that deals with knowledge. After all, for thousands of years philosophers have been the architects of knowledge. It started with the Greeks, mainly Aristotle who provided the foundations, was frozen with the scholastic interpretation, was revitalized by Descartes who brought in experimental knowledge, was challenged by the British empiricists Hume, Locke, Berkeley, achieved its full rigor with Kant, and the American pragmatists (Peirce, James, Putnam) added the concept of revision of knowledge. Poincaré made precise the role of formalism in scientific theory, and Popper introduced, with the concept of falsifiability, a key element in the relation between experiment and formal theories. But since the middle

of the XX^o century epistemology seems to have drifted away from science, and philosophers tend to think that actually, philosophy and science have little to share. This feeling is shared by many scientists (Stephen Weinberg in “Dreams of a Final Theory”). This is a pity as modern sciences need more than ever a demanding investigation of their foundations.

Without pretending to create a new epistemology and using all the basic work done by philosophers, I will try to draw a schematic view of epistemology, using a format and words which may be more familiar to the scientific reader. The purpose is here to set the ground, starting from questions such as What is knowledge ? How does it appear, is formatted, transformed, challenged ? What are the relations between experimentation and intuition ? We will see what are the specificities of scientific knowledge, how scientific theories are built and improved, what is the role of measures and facts, what is the meaning of the mathematical formalism in our theories. These are the topics of this first chapter.

1.1 WHAT IS KNOWLEDGE ?

First, a broad description of what is, and what is not knowledge.

Knowledge is different from perception : the most basic element of knowledge is the belief (a state of mind) of an individual with regard to a subject. It can be initiated, or not, by a sensitive perception or by the measure of a physical phenomenon.

Knowledge is not necessarily justified : it can be a certain perception, or a plausible perception (“I think that I have seen...”), or a pure stated belief (“God exists”), or a hypothesis.

Knowledge is shared beliefs : if individual states of minds can be an interesting topic, knowledge is concerned with beliefs which can be shared with other human beings. So knowledge is expressed in conventional formats, which are generally accepted by a community of people interested by a topic. This is not a matter of the tongue which is used, it supposes the existence of common conventions, which enables the transmission of knowledge without loss of meaning.

Knowledge is a construct : this is more than an accumulation of beliefs, knowledge can be learnt and taught and for this purpose it uses basic concepts and rules, organized more or less tightly in fields addressing similar topics.

1.1.1 Circumstantial assertions

The most basic element of knowledge can be defined as a **circumstantial individual assertion**, which can be formatted as comprised of :

- the author of the assertion
- the specific case (the circumstances) about which the assertion is made. Even if it is often implicit, it is assumed that the circumstances, people, background,.. are known, this is a crucial part of the assertion.
- the content of the assertion itself : it can be simply a logical assertion (it has the value true or false) or be expressed in a value using a code or a number.

The assertion can be **justified** or not. The author may himself think that his assertion is only plausible, it is a hypothesis. An assertion can be justified by being shared by several persons. A stronger form of justification is a **factual justification**, when everybody who wants to check can share by himself the assertion : the assertion is justified by evidence. In Sciences factual justifications are grounded in measures, done according to precise and agreed upon procedures.

Examples of circumstantial individual assertions :

“Alice says that yesterday Bob had a blue hat”, “I think that this morning the temperature was in the low 15 °C”, “I believe that the cure of Alice is the result of a miracle”...

Knowledge, and specially scientific knowledge, is more than individual circumstantial assertions : it is a method to build narratives from assertions. It

proceeds by enlargement, by going from individuals to a community, from circumstantial to universal, and by linking together assertions.

1.1.2 Rational narrative

By combining together several assertions one can build a narrative, and any kind of theory is based upon such construct. To be convincing, or only useful, a narrative must meet several criteria, which makes it rational. Rationality is different from justification : *it addresses the syntax of the narrative*, the rules that the combination of different assertions must follow in the construct, and does not consider a priori the validity of the assertions. The generally accepted rules come from logic. Aristotle has exposed the basis of logic but, since then, it has become a field of research on its own (for more see Maths.Part 1).

Formal logic deals with logical assertions, that is assertions which can take the value true (T) or false (F) exclusively. Any assertion can be put in this format.

Propositional logic builds propositions by linking assertions with four operators \wedge (and), \vee (or), \neg (not), \Rightarrow (implies). For each value T or F of the assertions the propositions resulting from the application of the operators take a precise value, T or F. For instance the proposition : $P = (A \Rightarrow B)$ is F if $A = T$ and $B = F$, and $P = T$ otherwise. Then one can combine propositions in the same way, and explore all their possible values by table-truth, which are just tables listing the propositions in columns, and all their possible values in rows.

Demonstration in formal logic uses propositions, built as above, and deduces true propositions from a collection of propositions deemed true (the axioms). To do this it lists axioms, then row after row, new true propositions using a rule of inference : if A is T, and $(A \Rightarrow B)$ is T, then B is T. The last, true, proposition is then proven.

These two kinds of propositional logic can be formalized in the Boolean calculus, and automated.

Propositions deal with circumstantial assertions. To enlarge the scope of formal logic, predicates are propositions which enables the use of variables, belonging to some fixed collection. Assertions and propositions are then linked with the use of two additional operators : \forall (whatever the value of the variable in the collection), \exists (there is a value of the variable in the collection). Notice that, in first order predicates, these operators act only on variables, which are previously listed, and not on predicates themselves. One can build table-truth in the same way as above, for all combinations of the variable. Demonstrations can be done in a similar way, with rules of inference which are a bit more complicated.

The Gödel's completeness theorem says that any true predicate can be proven, and conversely that only true predicates can be proven. The Gödel's compactness theorem says in addition that if a formula can be proven from a set of predicates, it can also be proven by a finite set of predicates : there is

always a demonstration using a finite number of steps and predicates. These two theorems show that, so formalized, formal logic is fully consistent, and can be accepted as a sound and solid basis to build rational narratives.

This is only a sketch of logic, which has been developed in a sophisticated system, important in computer theory. Several alternate formal logics have been proposed, but they lead to more complicated, and less efficient, systems, and so are not commonly used. Other systems called also logic, have been proposed in special fields, such as Quantum Mechanics (see Josef Jauch and Charles Francis for more) and information theory. Actually they are Formal Systems, similar to the Theories of Sets or Arithmetic in Mathematics : they do not introduce any new Calculus of Predicates, but use Mathematical Logic acting on a set of axioms and propositions.

Using the basic rules of formal logic, one can build a rational narrative, in any field. Notice that in the predicates the collections to which variables must belong are not sets, such as defined in Mathematics, and no special property is assumed about them. A variable can be a citizen, belonging to a country and indeed many laws could be formulated using formal logic.

Formal logic is not concerned about the justification or the veracity of the assertions. It tells only what can be logically deduced from a set of assertions, and of course can be used to refute propositions which cannot be right, given their premises. For instance the narrative :

$\forall X$ human being, $((X \text{ is ill}) \wedge (X \text{ prays}) \wedge (\text{God wills})) \Rightarrow (X \text{ is cured})$

is rational. It is F only if there is a X such that the first part is T and X is not cured. And one can deduce that God's will is F in this case. Without the proposition (God wills) it would be irrational.

Rational narrative are the ingredient of mystery books : at the end the detective comes with a set of assertions to unveil the criminal. A rational narrative can provide a plausible explanation, and a rational, justified, narrative, is the basis for a judgement in a court of law.

Scientific knowledge of course requires rational narratives, but it is more than that. A plausible explanation is rooted in the specific circumstances in which it has occurred : there is no reason why, under the same circumstances, the same facts would happen. To go further one needs a feature which is called necessitation by philosophers, and this requires to go from the circumstantial to the universal. And scientific knowledge is justified, which means that the evidences which support the explanation can be provided in a controlled way.

1.2 SCIENTIFIC KNOWLEDGE

1.2.1 Scientific laws

Let us take some examples of scientific laws :

A material body which is not submitted to any force keeps its motion

For any ideal gas contained in a vessel there is a relation $PV = nRT$ between its pressure, volume, and temperature

For any conductive material submitted to an electric field there is a relation $U = RI$ between the potential U and the intensity of the current

Any dominant allele is transmitted to the descenders

Scientific laws are assertions, which have two key characteristics :

i) They are universal : they are valid whenever the circumstances are met. A plausible explanation if true in specific circumstances, a scientific law is true whenever some circumstances are met. Thus in formal logic they should be preceded by the operator \forall . This is a strong feature, because if it is false in only one circumstance then it is false : it is **falsifiable**. This falsifiability, which has been introduced by Popper, is a key criterion of scientificity.

ii) They are justifiable : what they express is linked to physical phenomena which can be reproduced, and the truth of the law can then be checked by anybody. In a justified plausible explanation, the evidences are specific and exist only in one realization. For a scientific law the evidences can be supplied at will, by following procedures. This feature, introduced by Kant, distinguishes scientific narratives from metaphysical narratives.

One subtle point of falsifiability, by checking a prediction, is that it requires the possibility, at least theoretically, to test and check any value of each initial assertion before the prediction. Take the explanation that we have seen above :

$\forall X$ human being, $((X \text{ is ill}) \wedge (X \text{ prays}) \wedge (\text{God wills})) \Rightarrow (X \text{ is cured})$

For any occurrence, three of the assertions can be checked, and so one could assume that the value of the fourth (God's will) is defined by the final outcome in each occurrence, and we would have a scientific law. However falsifiability requires that one could test for different values of the God's will before measuring the outcome, so we do not have a scientific law. The requirement is obvious in this example but we have less obvious cases in Physics. Take the two slits experiment and the narrative :

$(\text{particles are targeted to a screen with two slits}) \wedge (\text{particles behave as waves}) \Rightarrow (\text{we see a pattern of interferences})$

Without the capability to predict which of the two, contradictory, behaviors, is chosen, we cannot have a scientific law.

These criteria are valid for any science. The capability to describe the circumstances, to reproduce or at least to observe similar occurrences, to check and whenever possible to measure the facts, are essential in any science. However falsifiability is usually a difficult criterion to meet in Social Sciences, even if one

strives to control the environment, but this is close to impossible in Archeology or History, where the circumstances in which events happened are difficult or impossible to reproduce, and are usually not well known. The extinction of the dinosaurs by the consequences of the fall of an asteroid is a plausible explanation, it seems difficult to make a law of it.

1.2.2 Probability in scientific laws

The universality of scientific laws opens the way to probabilistic formalization : because one can reproduce, in similar or identical manner, the circumstances, one can compute the probability of a given outcome. But this is worth some clarification because it is closely linked to a big issue : are all physical processes determinist ?

In Social Sciences, which involve the behavior of individuals, the assumption of free will negates the possibility of determinist laws : the behavior of a man or woman cannot be determined by his or her biological, social or economic characteristics. This has been a lasting issue for philosophers such as Spinoza, with a following in Marxist ideas. However, from our point of view here, no scientific law has been proven, just more or less strong correlations between variables, which can be used in empirical studies (such as market studies).

In the other fields, the discrepancy between the outcomes can be imputed to the fact that the circumstances are similar, but not identical :

- the measures are imperfect
- the properties of the objects (such as their shape) are not exactly what is assumed
- some phenomena are neglected, because it is assumed that their effect is small, but it is non null and unknown

This is a common case in Engineering, where phenomenological laws are usually sufficient for their practical use (for instance for assessing the strength of materials). In Biology the Mendel's heredity laws provide another example. As an extreme example, consider the distribution of the height of people in a given population. One could assume that, for a given individual, the biological processes at work are such that, for a given genetic structure, diet, way of life,... the height is determined (or quite so). The distribution that one observes is the result of the distribution of the factors which are neglected, and it can be made more precise, for instance just by the distinction between male and female.

And similarly, at a macroscopic scale, probabilist laws are commonly used to represent physical processes which involve a great number of interacting microsystems (such as in Thermodynamics) whose behavior cannot be individually measured, or discontinuous processes such as the breakdown of a material, an earth-quake,...which are assumed to be the result of slow continuous processes.

In all these cases a probabilist law does not imply that the process which is represented is not determinist, just that all the factors involved have not been accounted for. I don't think that any geologist believes that earth-quakes are a pure random process.

However one knows of physical elementary processes which, in our state of knowledge, seem to be not determinist : the tunnel effect in semi-conductors, the disintegration of a nucleus or a particle, or conversely the spontaneous creation of a particle, ..

Quantum Mechanics (QM) makes an extensive use of probability laws, and some of its interpretations postulate that at some scale physical laws are fundamentally not determinist. Up to now QM is still the only theory which can represent efficiently elementary non determinist phenomena. However, as we will see in the next chapter, the probabilist feature of the main axioms of QM does not come from some random behavior of natural objects, but from the discrepancy between the measures which can be done and their representation in our theories.

1.2.3 Models

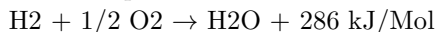
To implement a scientific laws, either to check it or to use it for practical purpose (to predict an outcome), scientists and engineers use models. A **model** can be seen as the general representation of the law. It comprises :

- a system : the area in which the system is located and the time frame during which it is observed, the list of the objects and of their properties which are considered
- the circumstances if they are specific (temperature, interference with the exterior of the system,...)
- the variables representing the properties, associated each to a mathematical object with more specific mathematical properties if necessary (a scalar can be positive, a function can be continuous,...)
- the procedures used to collect and analyze the data, notably if statistical methods are used.

Building and using models are a crucial part of the scientific work. Economists are familiar with the denomination models, either theoretical or as a forecasting tool. If they are not known by the name, any engineer or theoretical physicist use them, either to compute solutions of a problem from well established laws, or to explore the consequences of more general hypotheses. A model is a representation, usually simplified, of part of the reality, built from concepts, assumptions and accepted laws. The simplification helps to focus on the purpose, trading accuracy for efficiency. Models provide both a framework in which to make the computations, using some formalism in an ideal representation, and a practical procedure to organize the collection and analysis of the data. They are the embodiment of scientific laws, implemented in more specific circumstances, but still with a large degree of generality which enables to transpose the results from one realization to another. Actually most, if not all, scientific laws can be expressed in the framework of a model.

Models use a formalism, that is a way to represent the properties in terms of variables, which can take different values according to the specific realizations of the model, and which are used to make computations to predict a result. The

main purpose of the formalism is efficiency, because it enables to use rules and theorems well established in a more specific field. If the variables are logic, then formal logic provides an adequate formalism. Usually in Physics the formalism is mathematical, but other formalisms exist. The most illuminating example is the atomic representation used in Chemistry. A set of symbols such as :



tells us almost everything which is useful to understand and work with most of chemical experiments.

However the role of Mathematics in the formalism used in Physics leads us to have a look about the status of Mathematics itself in Science.

1.2.4 The status of Mathematics

It is usually acknowledged that Euclide founded Mathematics, with his Geometry, based on the definition of simple objects (points, lines,...) which are idealization of physical objects, a small number of axioms, and logic as the computational motor. For millennia it has been seen as the embodiment of rationality, and Mathematics has been developed in a patchwork of different fields : Algebra, Analysis, Differential Geometry... extending the scope of objects, endowed with more sophisticated properties. In the XIX^o century mathematicians felt the need to unify this patchwork and to found a clean Mathematics, grounded in as few axioms as possible. This was also the consequence of discoveries, such as non euclidean geometries by Lobatchevski, and of paradoxes in the newly borne Cantor's set theory. And this was also the beginning of many controversies, which are not totally closed at this day.

However this endeavour (promoted by Hilbert) lead to the creation of Mathematical Logic. This is actually a vibrant field of Mathematics of its own, which aims at scrutinizing Mathematics with respect to its consistency. It became clear that, in order to progress, it was necessary to distinguish in the patchwork some mathematical theories, and the focus has been put on Arithmetic and Set Theory, as they are the starting point for all the other fields of Mathematics. Without attempting to give even an overview of Mathematical Logic, three main features emerge from its results :

- the need to define objects specific to each field (natural numbers, sets) through their properties which are then enshrined in the axioms of the theories;
- the fact that these objects are of an abstract nature, in the meaning that they cannot be seen simply as the idealized realization of some physical objects, as points, lines,... were in Euclidean Geometry;
- and this fact is compounded by the need to assume properties which cannot be the realization of physical objects : the key example is the axiom of infinity in the Set Theory which postulates the existence of a set with an infinite number of elements.

So Mathematics is essentially different from formal logic (even if it uses it to work on these objects) : it relies on the prior definitions of objects and precise

axioms, and deal only with these objects and those which can be constructed from them. Formal logic is only syntax, Mathematics assumes a semantic part.

On these bases several sets of axioms have been proposed both for Arithmetic (Peano) and the Set Theory (Zermello, Frankael). They provide efficient systems, which have been generally accepted at the time and still nowadays, with some variants. However two results came as a big surprise:

- Gödel proved in 1931, with complements given by Gentzen in 1936 and Ackerman in 1940, that in any formal system powerful enough to represent Arithmetic, there are propositions which are true but cannot be proven.

- Church proved in 1936 that there cannot exist a fixed procedure to prove any problem in Arithmetic in a finite time (this is not a decidable theory).

The incompleteness Gödel's theorem is commonly misunderstood. Its meaning is that, to represent Arithmetic with all its usual properties that we know, we need a minimum set of axioms, but one could then add an infinite number of other axioms, which would not be inconsistent with the theory : they are true, because they are axioms, and they cannot be proven, because they are independent from the other axioms.

The Church's theorem is directly linked with computers (formalized as Turing's machines) : it cannot exist a program which would solve automatically all problems in Arithmetic.

Many similar or more sophisticated results have been proven in different fields of Mathematics. For our purpose here, several conclusions can be drawn :

- i) Mathematics can be seen as a science : it deals with objects and properties, using formal logic, to deduce laws which are scientific by the fact that they are always true for any realization of the objects. It has the great privilege to invent its own objects, however this comes with a price : the definition is not unique, other properties could be added or specified without harming Mathematics.

- ii) The choice of the right axioms is not dictated by necessity, but by efficiency. Mathematics, as we know it, has not been created from scratch by an axiomatic construct, it is the product of centuries of work, sometimes not rigorous, and the axioms which emerge today are the ones which have been proven efficient for our needs. But perhaps, one day, we will find necessary to enlarge the set of axioms, as it has been done with the axiom of infinity.

- iii) Because the objects are not simple idealization of physical realizations, and because there is no automated procedure to prove theorems, and so to extend Mathematics, it appears that it is a true product of the human mind. All mathematicians (as Poincaré noticed) have known these short periods of illumination, when intuition prevails over deduction, to find the right path to the truth. It seems that an artificial intelligence could not have arrived to the creativity that Mathematics requires.

1.3 THEORIES

Scientific laws are an improvement over circumstantial explanations, because they have the character of necessity and they are related to physical observable phenomena. Often philosophers view laws of nature as something which has to be discovered, as a new planet, hidden from our knowledge or perception. But science is more than a collection of laws, it has higher goals, it aims at providing a plausible explanation for as many cases as possible. Early on appeared the want to unify these laws, either to induce a cross fertilization process, or by the more holistic concern to understand what is the real world that they describe : *Science should provide more than efficient tools, it should explain what it is.*

Scientific laws rely on the definition of objects (material body, force, ideal gas,...) which have properties (motion, volume, pressure,...) related to observable physical phenomena and also represented by mathematical objects (scalar, functions,...). These concepts have emerged in each field, and have been organized in Theories : Mechanics, Fluid Mechanics, Thermodynamics, Electromagnetism, Theory of Fields,... and a similar process has been at work in Chemistry or Biology. And of course the want to unify further these fields has appeared. However the endeavour has not gone as well as in Mathematics. Many scientists are quite pleased with their tools and do not feel the need to go beyond what they use and know. A pervasive mood exists in Physics that the focus shall be put on experiments : if it works then it is true, whatever the way the computations are done. In an empiricist vision the concepts are nothing more than what is measured : a scientific law is essentially the repeated occurrences of observed facts, and one can accept a patchwork of laws. QM has greatly strengthened this approach, at first by casting a deep doubt about concepts which were thought to be strong (such as location, speed, matter,...) and the generalization of probabilist laws, and then by promoting the use of new concepts (fields, wave function, superposition of states...) which, from the beginning, were deemed to have no physical meaning, at least that we could understand. However the need for a more unified and consistent vision exist, even if it is met by unsatisfying construct, and one goes from a patchwork of scientific laws to theories.

1.3.1 What is a scientific theory ?

A scientific theory aims at giving a unified vision of a field, a framework in which scientific laws can be expressed, and a formalism which enables to deduce new laws that can be checked. So it comprises :

- a set of concepts, objects related to physical realizations, to which are attached properties which can be measured. These properties can be seen as defining the objects.
- a set of fundamental laws, or first principles : expressed in general terms, they are based on the observation of the physical world, and grounded in experiments, but they can or cannot be checked directly.
- a formalism, which provides the framework of models, and the computational tools to deduce new laws, forecast the results of experiments and check

the laws.

Examples :

The atomist theory in Chemistry. Compounds are made of a combination of 118 elements with distinct chemical properties, chemical reactions occur without loss of elements and an exchange of energy, ruled by thermodynamics.

The Newton's Mechanics. Material bodies are composed of material points, in a solid they stay at a constant distance from each other. The motion of material bodies is represented in the Galilean Geometry, it depends on their inertia and on forces which are exerted by contact or at a distance, according to fundamental laws.

Special Relativity. The universe is a four dimensional affine space endowed with a fixed Lorentz metric. Material bodies move along world lines at a constant velocity and their kinematics is characterized by their energy-momentum vector. The speed of light is constant for any inertial observer.

The properties are crucial because, for each situation, they can identify generic objects with similar properties, and associate to these objects a set of well defined values, which can be measured in each occurrence : "all insects have three pairs of legs", "material bodies travel along a world line in the 4 dimensional universe", "for any gas there is a temperature T". But by themselves they do not have a predictive power. In some cases the value of the variable comes from the definition itself (the number of legs of an insect), but usually it does not provide the value of the variable (the temperature of a gas).

As said before, the formalism used is not necessarily mathematical, but it acquires a special importance. This is a matter of much controversies but it is clear that major steps in the theories would have been impossible without prior progresses in the formalism which is used : Chemistry with the atomist representation, Mechanics with differential and integral calculus, General Relativity with differential geometry, and even Economics with Statistics. The use of more powerful mathematical tools, and similarly of computational techniques, increases our capacity to check predictions, but also to build the theories. Inspired by Thermodynamics and QM, it has been proposed to give to Information Theory an unifying role in Physics. A step further, considering that many structures used in different fields have similar features, the Category Theory, a branch of Mathematics developed around 1945 (Eilenberg, Mac Lane) has been used as a formalism in Physics, notably in Quantum Computing (Heyting algebras).

Fundamental laws can be not justified experimentally, their validity stems essentially from the consequences which can be deduced from them. From this point of view this is the theory as a whole which is falsifiable : if any law that can be deduced in the framework of the theory is proven false, then this is the entire theory which is at risk. And actually this has been a recurring event : Maxwell's laws and Galilean Geometry had to be revised after the Michelson and Morley experiments, the Atomist theory has had to integrate radio-activity,...The process has not gone smoothly, and usually patches are proposed to sustain the existing theory. And indeed a good part of the job of

scientists is to improve the theories, meaning to propose new theories which are then checked. What are the criteria in this endeavour ?

1.3.2 The criteria to improve scientific theories

Simplicity

The first criterion is simplicity. This is an extension of the Occam's razor rule : whenever we face several possible explanations, the fewer assumptions are made, the better. With our description of scientific theories it is easy to see what are the parameters to look for improvements. There must be as few kinds of objects as possible, themselves differentiated by a small number of properties or variables. There are 118 elements with distinct chemical properties, their nuclei are comprised of 12 fermions, there are millions of eukaryotes, but their main distinctive characteristics come from their DNA, organized in a small number of chromosomes, which are a combination of 4 bases. The electric and magnetic fields have been unified by the Maxwell's laws, and the unifications of all force fields including gravitation is the Graal of physicists. Similarly there should be as few fundamental hypotheses as possible. The Galilean system was not more accurate or legitimate (the assertions that Earth circles the Sun or that Sun moves around the Earth are both valid) than the Ptolemaic system, but it was simpler and provided a general theory to compute the trajectories of bodies around a star and paved the way to the Newton's gravitation law.

Enlarge the scope of phenomena addressed by the theory

The second criterion is the scope of the field which is addressed by the theory. Science is imperialist : it strives to find a rational explanation to everything. Lead by the Occam's razor rule it looks for more fundamental objects and theories, from which all the others could be deduced. This is a fact, and a legitimate endeavour. It has been developed in the different forms of positivism. In its earlier version (A.Comte) science had to deal only with and proceed from empirical evidence, scientific knowledge could be built by a logic formalization, which leads to a hierarchy of sciences giving preeminence to mathematics. In its more modern version positivism embraces the idea of the unity of science, that there is, underlying the various scientific disciplines, basically one science about one real world. Actually this is more complicated.

Starting with mathematics, as we have seen it could be seen as a science. True, mathematicians can invent their own objects. Quite often a narrative in Mathematics starts as "Let be a set such that...", but the first step required is to prove that such a set exists (as an example the definition of the tensorial product of vector space from an universal property). And if this is not possible one has to add another axiom (such as infinite sets), and support the consequences.

In natural sciences it is a sound requirement that there is a strong, unified background, explaining and reflecting the unity of the physical world. But in the different fields theories usually do not proceed from the most elementary laws.

The atomic representation used in Chemistry precedes quantum field theories of particles. Biology acknowledges the role of chemical reactions, but its basic concepts are not embedded in chemistry. We do not have in Physics a theory which would be general and powerful enough to account for everything. And anyway in most practical cases specific theories suffice. They use a larger set of assumptions, which are simplified cases of general laws (Galilean Geometry replacing Relativist Geometry, Newton's laws substituted to General Relativity) or phenomenological laws based on experimental data. In doing this the main motivation of scientists is efficiency : they do not claim the independence of their fields, but acknowledge the necessity of simpler theories for their work. However one cannot ignore that this move from one level to the other may cover a part of mystery. We still do not understand what is life. We do not have any determinist model of irreversible elementary process (such as the disintegration of a particle).

Economics is by far the social science which has achieved the higher level of formalization, in theoretical studies, empirical predictive tools, and in the definition of a set of concepts which give a rigorous basis for the collection and organization of data. Through the accounting apparatus, at the company level, the state level as well as many specialized fields (welfare, health care, R&D,...) one can have a reliable and quantified explanation of facts, and be able to assess the potential consequences of decisions. Because of the stakes involved these concepts are controversial, but this is not an exclusivity of Economics ¹. Actually what hampers Economics, and more generally the Social Sciences, is the difficulty of experimentation. Most of the work of scientists in these fields relies on data about specific occurrences, past or related to a few number of cases. The huge number of factors involved, most of which cannot be controlled, weakens any prediction, and the frailty of phenomenological laws in return limits the power of the falsifiability check. But this does not prevent us to try.

So we are still far away from a theory of everything. But the imperialism of science is legitimate, and we should go with the Hilbert's famous saying : "Wir müssen wissen, wir werden wissen". It is backed by the pressing want of people to have explanations, even when they are not always willing to accept them. As a consequence it increases the pressure on scientists and more generally on those who claim to have knowledge. As G.B.Shaw said "All professions are a conspiracy against the laity". So it is a sound democratic principle that scientists should be kept accountable to the people who fund their work.

Conservative pragmatism

The third criterion in the choice of theories is that any new theory should account for the ones that it claims to replace. What one can call a conservative pragmatism. Sciences can progress by jumps, but most often they are revisions of present theories, which become embedded in new ones and are seen as special

¹Actually some philosophers (who qualify themselves as feminists, such as Antony) deny that science is objective, and is very much an instrument of oppression (in Turri about Quine).

case occurring in more common circumstances. This process, well studied by G.Bachelard, is most obvious in Relativity : Special Relativity encompasses Galilean Geometry, valid when the speeds are weak, and General Relativity encompasses Special Relativity, valid when gravitation does not vary too much. Old theories have been established on an extended basis of experimental data, and backed by strong evidences which cannot be dismissed easily. New evidences appear in singular and exceptional occurrences and this leads to a quest for more difficult, and expansive, experimentations, which require more complex explanations. This is unavoidable but has drawbacks and the path is not without risks. The complexity of the proofs is often contrary to the first criterion - simplicity - all the more so when the new theory involves new objects with assumed, non checked, properties. The obvious examples are dark matter, or the Higgs boson. Of course it has happened in the past, with the nucleus, the neutrino, ... but it is difficult to feel comfortable in piling up enigma : the purpose of science is to provide answers, not to explain a mystery by a riddle. And when the new enigma requires more powerful tools the race may turn into a justification in itself.

1.4 FOUR QUESTIONS ABOUT SCIENCE

1.4.1 Is there a scientific method ?

It is commonly believed that one distinctive feature of the scientific work is that it proceeds according to a specific method. There is no doubt that the prerequisite of any scientific result is that it is justified for the scientific community. So the specificity of a scientific method would be guaranteed by higher ethical and professional standards. This claim is commonly associated to the “peer review” process : any result is deemed scientific if it has been approved for publication by at least two boffins of the field. Knowing the economics of this process, this criterion seems less reliable than what is usually required for an evidence in a court of justice, as recent troubles with published results show. The comparison is not fortuitous. For people who have dedicated years of their life to develop or to teach ideas, it is neither easy nor natural to challenge their beliefs, and all the more so when these beliefs are supported by the highest authorities in the field. Science has become a very competitive area, with great fame and financial stakes. Assume that fierce competition has increased the pressure to innovate is a bit optimistic. The real pressure comes from outside the scientific community, when quick economic return can be expected from a new discovery. This is no surprise that Computer Sciences or Biology have made gigantic progresses, meanwhile Particle Physics is still praising a Standard Model 40 years old. In any business, if the introduction of a new product was submitted to the anonymous judgment of your competitors, there would be no innovation. Only the interest of the customers should matter, but in Science this is a very distant concern, as well as the more direct interest of students who strive to understand theories that are reputed impossible to understand.

More generally this leads to question the existence of a science in fields such as History, Archeology,... Clearly there are criteria for the justification of assertions in these fields, which are more or less agreed upon by their communities, but it seems difficult that these assertions would ever be granted the status of scientific laws, at best they are plausible explanations.

So, and in agreement with most philosophers, I consider that scientific knowledge cannot be characterized by its method.

1.4.2 Is there a Scientific Truth ?

A justified assertion can be accepted as truth in a Court of justice. But not that many people would endorse a scientific truth, and probably few scientists as well. Scientific theories are backed by a huge amount of checked evidences, and justified by their power to provide plausible explanations for a large scope of occurrences. So in many ways they are closer to the truth than most conceivable human assertions, but the purpose of science is not the quest for the truth, because science is a work in progress and doubt is a necessary condition for this progress. A striking example of this complex relation between science and truth is Marxism : Karl Marx made very valuable observations about the relations

between technology, economic and political organizations, and claimed to have founded a new science, which enables people to make history. The fact that his followers accepted his claims to be the truth had dramatic consequences ².

1.4.3 Science and Reality

Science requires the existence of a real world, which does not depend on our minds, without which it would be impossible to conceive universal assertions. Moreover it assumes that this reality is unified, in a way that enables us to know its different faces, if any. Perhaps this is most obvious in social sciences : communities have very different organizations, beliefs and customs, but we strive to study them through common concepts because we see them as special occurrences of Human civilizations, with common needs and constraints. However this does not mean that we know what is reality : what we can achieve is the most accurate and plausible representation of reality, but it will stay temporary, subject to revision, and adjusted to the capability of our minds.

Because this representation is made through a formalization, the language which is used acquires a special importance. Some scientists resent this fact, perceived as an undue race towards abstraction, meanwhile they believe that empirical research should stay at the core of scientific progress. Actually the issue stems less from the use of more sophisticated mathematics than from the reluctance to adjust the concepts upon which the theories are based to take full advantage of the new tools. It is disconcerting to see physical concepts such as fields, particles, mass, energy, momentum,.. mixed freely with highly technical topological or algebraic tools. The discrepancy between the precision of the mathematical concepts and the crudeness of the physical concepts is source of confusion, and defiance. But the revision of the concepts will not come from the accumulation of empirical data, whatever the sophistication of the computational methods, it will come from fresh ideas.

From where do come these fresh ideas ? As we have seen above, and clearly in the case of mathematics, they are not the result of inference : a theory, with its collection of concepts and related formalism, has for purpose to provide models to explain specific occurrences. A continuous enlargement of the scope of experimental research provides more reliable laws, or conversely the proof of the failure of the theory, but it does not creates a new theory. New theories require a revision of the concepts, which may imply, but not necessarily, new hypotheses which are then checked. Innovation is not a linear, predictable process, it keeps some mystery, which, probably, is related to the genuine difference between computers and human intelligence.

1.4.4 Dogmatism and Hubris in Science

As the criteria for the validation of Scientific knowledge began to emerge, the implementation of the same criteria leads to two opposite dogmatisms, and their

²This aspect of marxism as the pretense of a science has been explained in my article published in 1982 in "les temps modernes".

unavoidable hubris. And what is strange is that, in some areas of the present days Physics, these opposite succeeded to be packaged together, for the worst.

The first dogmatism is the identification of the real world with the concepts. This is what Euclide and generations of mathematicians did for millennia : a point, a line, exist really, as well as parallels lines : after all they are nothing more than the idealization of tangible objects whose properties can be studied as suited. The overwhelming place taken by the mathematical formalism and the power it gives to compute complicated predictions lead to believe in the adequation between models and the real world. If it can be computed, then it exists. And if something cannot be computed, it is not worth to be considered. The first challenges to this dogmatism appear with Relativity, then the Physics in the atomic world. Scientists had been used to consider natural a 3 dimensional euclidean universe, with an external time. The jump to a 4 dimensional representation, and worst a curved Universe, seemed intractable. If the Universe integrates time, do the past and future events exist all together ? Still today, even for some professionals physicists, it seems difficult to address these questions. They do not realize that, after all, the idea of an infinite, flat Universe, existing for ever, is also a controversial representation. Similarly Mechanics and its admirable mathematical apparatus, seemed to breakdown when confronted to experiments in the atomic world : particles cannot pass the test of the two slits experiments, electrons could not keep a stable orbit around the nucleus, even Chemistry was challenged with the non conservation of matter and elements. Of course Engineers had for centuries a more pragmatic approach to the problem, the clean idea of continuous, non dissipative, motion had been replaced by phenomenological laws which could deal with deformable solids, fluid, and gas. But this was only Engineering...

The second dogmatism appeared, and triumphed, in reaction to the disarray caused by this discrepancy between a comprehensive and consistent vision and the experiments. Since the facts are the ultimate jury in checking a Scientific Theory, let us put the measures at the starting point in the elaboration of the theories. And because experimentation is overall a matter of statistical evaluation, it is natural to give to probability the place that it should have had from the beginning. There is nothing wrong in acknowledging the actual practices of scientific experiments. After all a Scientific Law is no more than the repetition of occurrences. The formalism of Statistical Mechanics was available, and soon, with the support of some mathematical justification, Quantum Physics had been born, and stated in axioms, rules and computational methods. The central issue, pushed by the supporters of the first dogmatism, was then to find a physical justification to the new formalism. As of today there has not been a unique answer. For some physicists Quantum Mechanics belong to a realm inaccessible to human understanding, a modern Metaphysics that it is vain to discuss, even if it can be marginally justified by mathematical considerations in simple cases. For others the want to find an interpretation is stronger, and the past century has been heralded with hundred of interpretations. They succeed actually in merging the two dogmatisms : if QM is stated in bizarre, non intuitive rules, it is because Reality itself is bizarre : it is discreet, non determinist.

We retrieve the identification of the formalism, as convoluted as it is, with the real world, but at the price of an obvious lack of agreement in the Scientific community, and at best a muddled picture. One of the strangest example of this new dogmatism is given in Cosmology : because we can model the Universe, it is possible to compute the whole Universe, and adding some QM, even consider the wave function of the Universe, which could then assess the probability of occurrences of the parallel universes...

Dogmatism and hubris go together. The criterion of factual justification is replaced by the forced identification of the real world with the formalism : if the computation works, it is because this is how the physical reality is. Humility is not the most significant feature of the Human mind, happily so. We need concepts, broad, easy to understand, illuminating and consistent representations which can be implemented and developed, which can be understood, learnt and taught. They can only be the product of intuition, of the imagination of the Human brain, they will never come from a batch of data. These ideas must be kept in check by the facts, not suppressed by the facts. But in the same time we must keep in mind that these are our concepts, our ideas, and that reality is still there, waiting to be probed, not enlisted to our cause. This leads to the reintroduction of the Observer in Physics, an object to which the rest of the book will give a significant place, and to which it could be dedicated.

1.5 FUNDAMENTAL PRINCIPLES IN PHYSICS

Whatever the theory in Physics there are some fundamental principles which are generally accepted.

1.5.1 Principle of Relativity

Scientific laws in Physics require measures of physical phenomena. Each object identified in a model has properties which are associated to mathematical objects, and the measure of these properties implies that it is possible to associate figures, real scalars, to the properties. There are many ways to do this, and because Scientific laws are universal, it shall be possible to do the measures in a consistent way, in precise protocols, and because it shall be possible to check the law in different occurrences, the protocol must tell how to adapt the measures to different circumstances.

The Principle of Relativity is used with different meanings in the literature. Here I will state it as “Scientific laws do not depend on the observer”. Which is the logical consequence of the definition of Scientific laws : they should be checked for any occurrence, as long as the proper protocols are followed, whoever do the experiment (the observers), whenever and wherever they are located. It has strong and important consequences in the mathematical formalization of the theories.

In any model the quantities which are measured are represented as mathematical objects, which have their own properties, and these properties are a defining part of the model, notably because they impose the format to collect the data. For instance in the Newton’s law $\vec{F} = m\vec{\gamma}$ the quantities \vec{F} , $\vec{\gamma}$ are vectors, and we must know how their components change when one uses one frame or another. Similarly the laws should not depend on the units in which the quantities are expressed. As a general rule, if a law is expressed as a relation $Y = L(X)$ between variables X, Y and there are relations $X' = R(X)$, $Y' = S(Y)$ where R, S are fixed maps, given by the protocols under which two observers proceed and thus known, then the law L' shall be such that : $Y' = L'(X') \Leftrightarrow L' = S \circ L \circ R^{-1}$. This is of special interest when R, S vary according to some parameters, because the last relation must be met whatever the value of the parameter. This is the starting point for the gauge theories in Physics.

The Principle of Relativity assumes that there are observers. In its common meaning an observer is the scientist which makes the measures. But in a Theory it requires that one defines the properties of an observer : this is a concept as the others, and it is not always obvious to define precisely and in a consistent way what are these properties. One key property of observers is that they have free will, and this implies notably that they can change freely the conditions of an experiment (as the universality of scientific laws requires) : they can choose different units, spatial location of their devices, repeat the same experiment over and over,... Free will implies also that the observers are not subjected to the laws which rule the system they observe, however they are also subjected

to physical laws but it is assumed that these laws do not interfere with the experiment they review. This raises some issues in Relativity, and a big issue in Cosmology, which is a theory of the whole Universe.

1.5.2 Principle of Conservation of Energy and Momentum

The principle is usually stated as “In any physical process the total quantity of energy and momentum of a system is conserved”.

But its interpretation raises many questions.

The first is about the definition of energy and momentum. They come from the intuitive notion that every physical object carries with it a capacity either to resist to a change, or to cause a change in other objects. So energy and momentum are attached to each object of the system : it is one of their properties. For localized objects such as material bodies, these quantities are localized as well. For objects which are spread over a vast area (fluids, force fields), energy and momentum are defined as density, related to some measure of volume of the area. Then the principle reads as the sum of energy and momentum for all the objects of the system is conserved.

The second is the distinction between energy and momentum. The former is expressed as a scalar, the latter as a vectorial like quantity. And both are intuitively linked.

For a material body the momentum is related to the motion. Motion is a purely geometric concept, corresponding to the location and disposition of a material body, and their change with time. The motion of a material body cannot be changed without an external action, and this resistance, characterizing its inertia, is related to the mass of the body and its motion, and represented by kinematic quantities : the translational momentum and the rotational momentum, which have the general form $[\text{mass}] \times [\text{motion}]$. So momenta for material body are expressed by kinematic quantities, which express their inertia, and they require the introduction, besides the representation of the motion, of other variables similar to mass. Motion is a derivative of location, and it can be expressed at any given time : momenta are continuous variables, as long as the motion is continuous.

The instantaneous effort which is required to change the momenta is a force : $F = \frac{d}{dt} [\text{Momentum}]$, which equals the force of inertia of the body. The total effort done in a physical process during which the momenta have been changed is the energy transferred to the body, expressed as $\Delta([\text{Momentum}] \times [\text{Motion}])$ essentially because its most visible form is the kinetic energy of material bodies, which is also the source of heat at a microscopic level. So energy can be deduced from the momenta, but only the total variation of energy in a process has a clear meaning : in Thermodynamics energy is a state variable, defined up to a constant, and which can be positive or negative (depending on which object is considered).

If the picture is fairly clear for localized body, what about objects which are spread, such as fluids or fields ? For a fluid it is assumed that the momentum, and then the energy, are transferred by contact. But it is acknowledged that

there are forces which act without any contact, and then we have to assume that, if the transfers occurs, force fields also carry momentum, evidenced by their action on material bodies, and exchange energy.

This leads to the definition of potential energy : the energy which can be given or received by a physical object in the transformation of a system. So it has no meaning out of the context of a given system.

The third issue is about the measure of energy and momentum. The mass, translational as well as rotational, is assumed to be an intrinsic characteristic of material bodies. The measure of motion depends on the observer, so the measure of momentum varies with the observer. The variation of energy depends on the initial and final states.

The concepts of energy, momentum, and of evolution of a system, require a clear definition of the time, which can depend on the observer. These issues were not solved until the advent of Relativity, which also gives a relation between energy and momentum.

One feature to notice about this Principle is that it does not assume that the evolution is continuous : there are clearly two states of the system, differentiated by a time elapsed between the measures, but the process can be continuous or discontinuous. Then this is the difference between the values of energy and momentum at the beginning and at the end of the process which matters.

1.5.3 Principle of Least Action

As there are quantities which are globally conserved in a physical process, based upon experience, it is assumed that any system has privileged states, called states of equilibrium, from which it does not move without a change in its environment, for instance an external action. This concept still holds when one considers the evolution in time : equilibrium does not imply that the state of the system is frozen, it can move along a path from which it does not differ easily. This is the formalization of the idea that an isolated system is in the state of least energy, or will follow an evolution such that it will use the least energy. States of equilibrium can be achieved by a continuous or a discontinuous process : the Principle of Least Action does not tell how a state of equilibrium is reached, only what are its characteristics.

From Mechanics, this principle is usually represented in Physics by the fact that a scalar functional, the action, is stationary for the values corresponding to the state of equilibrium : $\ell \left(L \left(z^i, z_{\alpha}^i, z_{\alpha\beta}^i, \dots \right) \right)$ where $Z = z^i, z_{\alpha}^i, z_{\alpha\beta}^i, \dots$ are the variables and their partial derivatives and L a scalar function (the scalar lagrangian).

It comes from Analytic Mechanics where $L = \sum_i \frac{1}{2} m_i v_i^2 - U$ is the total energy of the system (Kinetic and potential) and the lagrangian has the general meaning of a density of energy / momentum, as described above.

The stationarity means that for any (small) changes δZ of the value of the variables around the equilibrium Z_0 the value of the functional ℓ is unchanged. So this is not necessarily a maximum or a minimum, even local. And a state of

equilibrium is not necessarily unique.

Whenever the variables are maps or functions defined over the area Ω of a manifold endowed with a volume measure ϖ the functional is assumed to be an integral :

$$\int_{\Omega} L \left(z^i(m), z_{\alpha}^i(m), z_{\alpha\beta}^i(m), \dots \right) \varpi(m)$$

More simply when the variables are functions of the time only, varying from t_1 to t_2 , the action reads:

$$\int_{t_1}^{t_2} L \left(z^i(t), z_{\alpha}^i(t), z_{\alpha\beta}^i(t), \dots \right) dt$$

This formulation is extensively used, and most of the laws in Physics can be expressed this way. The Principle does not tell anything about the lagrangian, in which lies the physical content. There are constraints on its expression, as we will see, due to the Principle of Relativity, but the choice of the right lagrangian is mostly an art, which of course must be checked by the consequences that can be deduced.

The Principle seems to introduce a paradox in that the values taken by the variables at any moment depend on the values on the whole evolution of the system, that is on the values which will be taken in the future. But this paradox stems from the model itself : at the very beginning the physicist assumes that the variables which are measured or computed belong to some class of objects which are defined all over Ω . So the variables are the maps and not the values that they take for each value of their arguments. We will come back on this point in the presentation of QM.

1.5.4 Second Principle of Thermodynamics and Entropy

The universality of scientific laws implies that experiments are reproducible, time after time, which requires either that the circumstances stay the same, or can be reproduced identically. This can be achieved only to some degree, controlled by checking all the parameters which could influence the results. It is assumed that the parameters which are not directly involved in the law which is tested are not significant, or keep a steady value, in time as well as in the domains which are exterior to the area which is studied. So universality implies some continuity of the phenomena.

Many discontinuous phenomena at a macroscopic scale can be explained as the result of continuous processes at a smaller scale : an earthquake is the result of the slow motion of tectonic plates. Others involve the transition between phases, which are themselves states of equilibrium, and can be explained, as we will see, by the interaction of microsystems. However discontinuous processes exist, but they are never totally discontinuous : what happens is the transition between different continuous processes. Brownian motion is modelled by patching continuous paths. And indeed totally discontinuous functions are a mathematical curiosity, not easy to build. So the maps involved in physical models can be safely assumed to be continuous, except at isolated points.

If a transition occurs between two states of equilibrium the Principle of Least action can be implemented for each of them. But this leaves several issues.

The first is about the concept of equilibrium itself. As we have seen above, stationary does not mean frozen, the Principle of Least Action encompasses systems whose state varies with the time, so it addresses also processes. But in Physics there are common restrictions imposed on these processes. In Thermodynamics equilibrium is identified with reversible processes, seen as slow processes : at any moment the system is close to equilibrium. In Theoretical Physics these are processes whose evolution is ruled by equations which are invariant by time reversal : if $X(t)$ is solution, then the replacement of t by $-t$ is still a solution. A reversible process is determinist (there is only one path to go from a state to another) but the converse is not true. The Second Principle of Thermodynamics is a way to study processes which do not meet these restrictions.

In Thermodynamics the Second Principle is based upon the equation :

$$dU = TdS - pdV + \sum_c \mu_c dN_c$$

where the internal energy U , the entropy S , the volume V and the number of moles of chemical species N_c are variables which characterize the state of the system. The key point is that they do not depend on the path which has been followed to arrive at a given state. In the evolution between states:

$$dU = \delta Q + \delta W$$

where $\delta Q, \delta W$ are the quantity of heat and work exchanged by the system with its surroundings during any evolution. The variable temperature T is a true thermodynamic variable : it has a meaning only at a macroscopic scale. The symbol d represents a differential, meaning that the corresponding state variables are differentiable, and thus continuous, and δ a variation, which can be discontinuous.

For a system in any process :

$$dS \geq \frac{\delta Q}{T}$$

so that for isolated systems $dS \geq 0$: their entropy can only increase and this defines an arrow of time. We have an equality only in reversible processes.

The Thermodynamics formulation can be generalized to the evolution of systems comprised of many interacting microsystems. The model, proposed first by Boltzmann and Gibbs, has been used with many variants, notably by E.T. Jayes in his Principle of Maximum Entropy in relation with Information Theory. Its most common formalization is the following. A system is comprised of N (a large number) identical microsystems. Their states are represented by a random variable $X = (X_a)_{a=1}^m$ valued in a domain Ω with an unknown probability law $\Pr(X_1 = x_1, \dots, X_N = x_N) = \rho(x_1, \dots, x_N)$. There are m macroscopic variables $(Y_k)_{k=1}^m$ which can be measured for the whole system, whose value depend on the states of the microsystems : $Y_k = f_k(x_1, \dots, x_N)$. Knowing the values $(\hat{Y}_k)_{k=1}^m$ observed, the problem is to estimate ρ .

The Principle of Maximum Entropy states that the law ρ is such that the integral :

$$S = \int_{\Omega} -\rho(x_1, \dots, x_N) \ln \rho(x_1, \dots, x_N) dx_1 \dots dx_N$$

over the domain Ω of the x_a is maximum, under the constraints :

$$\hat{Y}_k = f_k(x_1, \dots, x_N)$$

$$\int_{\Omega} \rho dx = 1$$

The solution of this problem leads to the introduction of m new variables $(\theta_k)_{k=1}^m$ (the Lagrange parameters) dual of the observables Y_k which are truly thermodynamic, they have no meaning for the microsystems. Temperature is the dual of energy.

So formulated we have a classic problem of Statistics, and we can give a more precise definition of a reversible process. If the process is such that :

- the state of a microsystem does not depend on the state of other microsystems, only on the state of the global system
- the collection $(Y_k)_{k=1}^m$ is a complete statistic (one cannot expect to have more information on the system by adding another macroscopic variable)

then it is not difficult, using the Pitman-Koopman-Darmois theorem, to show that the solution given by the Principle of Maximum Entropy is indeed a good maximum of likelihood estimator, and validates it, at least for computational purposes.

This scheme has been extended in the framework of QM, the quantity $-Tr[\rho \ln \rho]$ called the information entropy, becoming a functional and ρ an operator on the space of states.

So, for reversible processes, there is a model which is well grounded, and entropy has a clear meaning. It leads to the introduction of additional variables (notably temperature) which can be measured but have a meaning only for the whole system. But it is usually acknowledged that there is no satisfying general model for non reversible processes, or processes which involve disequilibrium (see G.Röpke for more).

However the concept of entropy has a real significance to answer the issue of the choice of a state of equilibrium, if several are possible in the implementation of the Principle of Least Action : it tells why one of them is privileged, and the driving force is the entropy. Moreover if disequilibrium processes are of importance at a macroscopic scale, due to their prevalence in practical problems, this is less so from a theoretical point of view. There is no obvious reason to focus on processes which are modelled by equations invariant by time reversal, and actually they are not in Quantum Theory of Fields. The issue of determinism, closely related to discontinuous processes, seem more important.

1.5.5 Principle of Locality

It can be stated as : “the outcome of any physical process occurring at a location depends only of the values of the involved physical quantities at this location”. So it prohibits actions at a distance. This is obvious in the lagrangian formulation of the Principle of Least Action : the integral is computed from data whose values are taken at each point m (but one can conceive of other functionals ℓ).

Any physical theory assumes the existence of material objects, whose main characteristic is that they are localized : they are at a definite place at a given time. To account for phenomena such as electromagnetism or gravity, the principle requires the existence of physical objects, the force fields, which have a

value at any point. Thus this principle is consubstantial to the distinction matter / fields. It does not prohibit by itself the existence of objects which are issued from fields and behave like matter (the bosons). And similarly it does not forbid the representation of material bodies in a formalism which is defined at any point : in Mechanics the trajectory of a material point is a map $x(t)$ defined over a period of time. But these features appear in the representation of the objects, and do not imply physical action at a distance. The validity of this principle has been challenged by the entanglement of states of bosons, but it seems difficult to accept that it is false, as most of the Physics use it.

Because any measurement involves a physical process, the principle of locality implies that the measures shall be done locally, that is by observers at each location. This does not preclude the observers to exchange their information, but requires a procedure to collect and compare these measures. This procedure is part of the system, and the laws that they represent. As it has been said before, the observer, if he is not by himself submitted to the phenomena that he measures (he has free will), has distinctive characteristics which must be accounted for in the formulation of a law.

The rest of this book will be in some way a practical illustration of this first chapter. We will successfully expose the Geometry of General Relativity, the Kinematics of material bodies, the Force fields, the Interactions Fields / Particles, the Bosons. Starting from facts, common or scientific known facts, we will make assumptions, then, using the right mathematical formalism and Fundamental principles, we will predict scientific laws, as theorems. And this is the experimental verification of these laws that will provide the validity of the theory. So this is very different, almost the opposite, of what is usually done in Physics Books, such as Feynman's, where the starting point is almost always an experiment. The next chapter, dedicated to Quantum Theory, is purely mathematical but, as we will see, it starts also by the construction of its own objects : physical models.

Chapter 2

QUANTUM MECHANICS

Quantum Physics encompasses several theories, with three distinct areas:

i) Quantum Mechanics (QM) proper, which, since the seminal von Neumann's book, is expressed as a collection of axioms, such as summarized by Weinberg :

- Physical states of a system are represented by vectors ψ in a Hilbert space H , defined up to a complex number (a ray in a projective Hilbert space)
- Observables are represented by Hermitian operators
- The only values that can be observed for an operator are one of its eigen values λ_k corresponding to the eigen vector ψ_k
- The probability to observe λ_k if the system is in the state ψ is proportional to $|\langle\psi, \psi_k\rangle|^2$
- If two systems with Hilbert space H_1, H_2 interact, the states of the total system are represented in $H_1 \otimes H_2$

and, depending on the authors, the Schrödinger's equation.

ii) Wave Mechanics, which states that particles can behave like fields which propagate, and conversely force fields can behave like pointwise particles. Moreover particles are endowed with a spin. In itself it constitutes a new theory, with the introduction of new concepts, for which QM is the natural formalism. Actually this is essentially a theory of electromagnetism, and is formalized in Quantum Electrodynamics (QED).

iii) The Quantum Theory of Fields (QTF) is a theory which encompasses theoretically all the phenomena at the atomic or subatomic scale, but has been set up mainly to deal with the other forces (weak and strong interactions) and the organization of elementary particles. It uses additional concepts (such as gauge fields) and formalism and computation rules (Feynman diagrams, path integrals).

I will address in this chapter *QM only*. It would seem appropriate to begin the Physics part of this book by QM, as it has been dominant and pervasive since 70 years. But actually it is the converse : the place of this chapter comes

from the fact that QM is not a physical theory. This is obvious with a look at the axioms : they do not define any physical object, or physical property (if we except the Schrödinger's equation which is or not part of the corpus). They are deemed valid for any system and, actually, they would not be out of place in a book on Economics. These axioms, which are used commonly, are not Physical Laws, and indeed they are not falsifiable (how could we check that an observable is a Hermitian operator ?). Some, whose wording is general, could be seen as Fundamental Laws, similar to the Principle of Least Action, but others have an almost supernatural precision (the eigen vectors). Nevertheless they are granted with a total infallibility, supported by an unshakable faith, lauded by the media as well as the Highest Academic Authorities, reputed to make incredibly precise predictions. Their power is limited only by a scale which is not even mentioned and which is impossible to compute.

This strange status, quite unique in Science, is at the origin of the search for interpretations, and for the same reason, makes so difficult any sensible discussion on the topic. Actually these axioms have emerged slowly from the practices of great physicists, kept without any change in the last decenniums, and endorsed by the majority, mostly because, from their first Physics 101 to the software that they use, it is part of their environment. I will not enter into a debate about the interpretations of these axioms, but it is necessary to evoke the attempts which have been made to address directly their foundations.

In seminal books and articles, von Neumann and Birkhoff have proposed a new direction to understand and justify these axioms. Their purpose was, from general considerations, to set up a Formal System, actually similar to what is done in Mathematics for Arithmetic or Sets Theory, in which the assertions done in Physics can be expressed and used in the predictions of experiments, and so granting to Physics a status which would be less speculative and more respecting of the facts as they can actually be established. This work has been pursued, notably by Jauch, Haag, Varadarajan and Francis in the recent years. An extension which accounts for Relativity has been proposed by Wightman and has been developed as an Axiomatic Quantum Field Theory (Haag, Araki, Halvorson, Borchers, Doplicher, Roberts, Schroer, Fredenhagen, Buchholz, Summers, Longo,...). It assumes the existence of the formalism of Hilbert space itself, so the validity of most of the axioms, and emphasizes the role to the algebra of operators. Since all the information which can be extracted from a system goes through operators, it can be conceived to define the system itself as the set of these operators. This is a more comfortable venue, as it is essentially mathematical, which has been studied by several authors (Bratelli and others). Recently this approach has been completed by attempts to link QM with Information Theory, either in the framework of Quantum Computing, or through the use of the Categories Theory.

These works share some philosophical convictions, supported with a strength depending on the authors, but which are nonetheless present :

i) A deep mistrust with regard to realism, the idea that there is a real world, which can be understood and described through physical concepts such as particles, location,...At best they are useless, at worst they are misleading.

ii) A great faith in the mathematical formalism, which should ultimately replace the concepts.

iii) The preeminence of experimentation over theories : experimental facts are seen as the unique source of innovation, physical laws are essentially the repeated occurrences of events whose correlation must be studied by statistical methods, the imperative necessity to consider the conditions in which the experiments can or cannot be made.

As any formal system, the axiomatic QM defines its own objects, which are basically the assertions that a physicist can make from the results of experiments (“the yes-no experiments” of Jauch), and set up a system of rules of inference according to which other assertions can be made, with a special attention given to the possibility to make simultaneous measures, and the fact that any measure is the product of a statistical estimation. With the addition of some axioms, which obviously cannot reflect any experimental work (it is necessary to introduce infinity), the formal system is then identified, by a kind of structural isomorphism, with the usual Hilbert space and its operators of Mathematics. And from there the axioms of QM are deemed to be safely grounded.

One can be satisfied or not by this approach. But some remarks can be done.

In many ways this attempt is similar to the one by which mathematicians tried to give an ultimate, consistent and logical basis to Mathematics. Their attempt has not failed, but have shown the limits of what can be achieved : the necessity to detach the objects of the formal system from any idealization of physical objects, the non unicity of the axioms, and the fact that they are justified by experience and efficiency and not by a logical necessity. The same limits are obvious in axiomatic QM. If to acknowledge the role of experience and efficiency in the foundations of the system should not be disturbing, the pretense to enshrine them in axioms, not refutable and not subject to verification, places a great risk to the possibility of any evolution. And indeed the axioms have not changed for more than 50 years, without stopping the controversies about their meaning. The unavoidable replacement of physical concepts, identification of physical objects and their properties, by formal and abstract objects, which is consistent with the philosophical premises, is specially damaging in Physics. Because there is always a doubt about the meaning of the objects (for instance it is quite impossible to find the definition of a “state”) the implementation of the system sums up practically to a set of “generally accepted computations”, it makes its learning and teaching perilous (the Feynmann’s affirmation that it cannot be understood), and eventually to the recurring apparitions of “unidentified physical objects” whose existence is supposed to fill the gap. In many ways the formal system has replaced the Physical Theories, that is a set of objects, properties and behaviors, which can be intuitively identified and understood. The Newton’s laws of motion are successful, not only because they can be checked, but also because it is easy to understand them. This is not the case for the decoherence of the wave function...

Nevertheless, this attempt is right in looking for the origin of these axioms in the critique (in the Kantian meaning) of the method specific to Physics. But it

is aimed at the wrong target : the concepts are not the source of the problems, they are and will stay necessary because they make the link between formalism and real world, and are the field in which new ideas can germinate. And the solution is not in a sanctification of the experiments, which are too diverse to be submitted to any analytical method. Actually these attempts have missed a step, which always exists between the concepts and the collection of data : the mathematical formalization itself, in models. Models, because they use a precise formalism, can be easily analyzed and it is possible to show that, indeed, they have specific properties of their own, which do not come from the reality they represent, but from their mathematical properties and the way they are used. The objects of an axiomatic QM, if one wishes to establish one, are then clearly identified, without disturbing the elaboration or the implementation of theories. The axioms can then be proven, they can also be safely used, as we will show in this book.

QM is about the representation of physical phenomena, and not a representation of these phenomena (as can be Wave Mechanics, QED or QTF). It expresses properties of the data which can be extracted from measures of physical phenomena but not properties of physical objects. To sum up : QM is not about how the physical world works, it is about how it looks.

2.1 HILBERT SPACE

2.1.1 Representation of a system

We have seen the central role played by models in the practical implementation of a theory to specific situations. It will be our starting point.

Let us start with common Analytic Mechanics. A system, meaning a delimited area of space comprising material bodies, is represented by scalar generalized coordinates $q = (q_1, \dots, q_N)$ its evolution by the derivatives $q' = (q'_1, \dots, q'_N)$. By extension q can be the coordinates of a point Q of some manifold M to account for additional constraints, and then the state of the system at a given time is fully represented by a point of the vector bundle $TM : W = (Q, V_Q)$. By mathematical transformations the derivatives q' can be exchanged with conjugate momenta, and the state of the system is then represented in the phase space, with a symplectic structure. But we will not use this addition and stay at the very first step, that is the representation of the system by (q, q') .

Trouble arises when one considers the other fundamental objects of Physics : force fields. By definition their value is defined all over the space x time. So in the previous representation one should account, at a given time, for the value of the fields at each point, and introduce unaccountably infinitely many coordinates. This issue has been at the core of many attempts to improve Analytic Mechanics.

But let us consider two facts :

- Analytic Mechanics, as it is usually used, is aimed at representing the evolution of the system over a whole period of time $[0, T]$, as it is clear in the Lagrangian formalism : the variables are accounted, together, for the duration of the experiment;

- the state of the system is represented by a map $W : [0, T] \rightarrow (Q, V_Q)$: the knowledge of this map sums up all that can be said on the system, the map itself represents the state of the system.

Almost all the problems in Physics involve a model which comprises the following :

- i) a set of physical objects (material bodies or particles, force fields) in a delimited area Ω of space x time (it can be in the classical or the relativist framework) called the system;

- ii) the state of the system is represented by a fixed finite number N of variables $X = (X_k)_{k=1}^N$ which can be maps defined on Ω , with their derivatives;

so that the state of the system is defined by a finite number of maps, which usually belong themselves to infinite dimensional vector spaces.

And it is legitimate to substitute the maps to the coordinates in Ω . We still have infinite dimensional vector spaces, but by proceeding first to an aggregation by maps, the vector space is more manageable, and we have some mathematical tools to deal with it. But we need to remind the definition of a manifold, a structure that we will use abundantly in the following (more in Maths.15.1.1).

2.1.2 Manifold

Let M be a set, E a topological vector space, an atlas, denoted $A = (O_i, \varphi_i, E)_{i \in I}$ is a collection of :

subsets $(O_i)_{i \in I}$ of M such that $\cup_{i \in I} O_i = M$ (this is a cover of M)
 maps $(\varphi_i)_{i \in I}$ called **charts**, such that :

i) $\varphi_i : O_i \rightarrow U_i :: \xi = \varphi_i(m)$ is bijective and ξ are the coordinates of M in the chart

ii) U_i is an open subset of E

iii) $\forall i, j \in I : O_i \cap O_j \neq \emptyset :$

$\varphi_i(O_i \cap O_j), \varphi_j(O_i \cap O_j)$ are open subsets of E , and there is a bijective, continuous map, called a transition map :

$$\varphi_{ij} : \varphi_i(O_i \cap O_j) \rightarrow \varphi_j(O_i \cap O_j)$$

Notice that no mathematical structure of any kind is required on M . A topological structure can be imported on M , by telling that all the charts are continuous, and conversely if there is a topological structure on M the charts must be compatible with it. But the set M has no algebraic structure : a combination such as $am + bm'$ has no meaning.

Two atlas $A = (O_i, \varphi_i, E)_{i \in I}, A' = (O'_j, \varphi'_j, E)_{j \in J}$ of M are said to be compatible if their union is still an atlas. Which implies that :

$\forall i \in I, j \in J : O_i \cap O'_j \neq \emptyset : \exists \varphi_{ij} : \varphi_i(O_i \cap O'_j) \rightarrow \varphi'_j(O_i \cap O'_j)$ which is a homeomorphism

The relation A, A' are compatible atlas of M , is a relation of equivalence. A class of equivalence is a **structure of manifold** on the set M .

The key points are :

- there can be different structures of manifold on the same set. On \mathbb{R}^4 there are unaccountably many non equivalent structures of smooth manifolds (this is special to \mathbb{R}^4 : on $\mathbb{R}^n, n \neq 4$ all the smooth structures are equivalent !).

- all the interesting properties on M come from E : the dimension of M is the dimension of E (possibly infinite); if E is a Fréchet space we have a Fréchet manifold, if E is a Banach space we have a Banach manifold and then we can have differentials, if E is a Hilbert space we have a Hilbert manifold, but these additional properties require that the transition maps φ_{ij} meet additional properties.

- for many sets several charts are required (a sphere requires at least two charts) but an atlas can have only one chart, then the manifold structure is understood as the same point M will be defined by a set of compatible charts.

The usual, euclidean, 3 dimensional space of Physics is an affine space. It has a structure of manifold, which can use an atlas with orthonormal frames, or with curved coordinates (spherical or cylindrical). Passing from one system of coordinates to another is a change of charts, and represented by transition maps φ_{ij} .

2.1.3 Fundamental theorem

In this chapter we will consider models which meet the following conditions:

Condition 1 i) The system is represented by a fixed finite number N of variables $(X_k)_{k=1}^N$

ii) Each variable belongs to an open subset O_k of a separable Fréchet real vector space V_k

iii) At least one of the vector spaces $(V_k)_{k=1}^N$ is infinite dimensional

iv) For any other model of the system using N variables $(X'_k)_{k=1}^N$ belonging to open subset O'_k of V_k , and for $X_k, X'_k \in O_k \cap O'_k$ there is a continuous map : $X'_k = F_k(X_k)$

Remarks :

i) The variables must be vectorial. This condition is similar to the superposition principle which is assumed in QM. This is one of the most important condition. By this we mean that the associated physical phenomena can be represented as vectors (or tensors, or scalars). The criterion, to check if this is the case, is : if the physical phenomenon can be represented by X and X' , does the phenomenon corresponding to any linear combination $\alpha X + \beta X'$ has a physical meaning ?

Are usually vectorial variables : the speed of a material point, the electric or magnetic field, a force, a moment,...and the derivatives, which are, by definition, vectors.

Are not usually vectorial variables : qualitative variables (which take discrete values), a point in the euclidean space or on a circle, or any surface. The point can be represented by coordinates, but these coordinates are not the physical object, which is the material point. For instance in Analytic Mechanics the coordinates $q = (q_1, \dots, q_N)$ are not a geometric quantity : usually a linear combination $\alpha q + \beta q'$ has no physical meaning. The issue arises because physicists are used to think in terms of coordinates (in euclidean or relativist Lorentz frame) which leads to forget that the coordinates are just a representation of an object which, even in its mathematical form (a point in an affine space) is not vectorial.

So this condition, which has a simple mathematical expression, has a deep physical meaning : it requires to understand clearly why the properties of the physical phenomena can be represented by a vectorial variable, and reaches the most basic assumptions of the theory. The status, vectorial or not, of a quantity is not something which can be decided at will by the Physicist : it is part of the Theory which he uses to build his model. We will see numerous examples, both ways, of this in the next chapters.

However we will see that the addition of a variable which is not a vector can be very useful (Theorem 24).

ii) The variables are assumed to be independent, in the meaning that there is no given relation such that $\sum_k X_k = 1$. Of course usually the model is used with the purpose to compute or check relations between the variables, but these relations do not matter here. Actually to check the validity of a model one considers all the variables, those which are given and those which can be computed, they are all subject to measures and this is the comparison, after the experiment, between computed values and measured values which provides the validation. So

in this initial stage of specification of the model there is no distinction between the variables, which are on the same footing.

Similarly there is no distinction between variables internal and external to the system : if the evolution of a variable is determined by the observer or by phenomena out of the system (it is external) its value must be measured to be accounted for in the model, so it is on the same footing as any other variable. And it is assumed that the value of all variables can be measured (we will come back on this point in the next section).

The derivative $\frac{dX_k}{dt}$ (or partial derivative at any order) of a variable X_k is considered as an independent variable, as it is usually done in Analytic Mechanics and in the mathematical formalism of r-jets.

iii) The variables can be restricted to take only some range (for instance it must be positive). The vector spaces are infinite dimensional whenever the variables are functions. The usual case is when they represent the evolution of the system with the time t : then X_k is the function itself : $X_k : \mathbb{R} \rightarrow O_k :: X_k(t)$. What we consider here are variables which cover the whole evolution of the system over the time, and not only just a snapshot $X_k(t)$ at a given time. But the condition encompasses other cases, notably fields F which are defined over a domain Ω . *The variables are the maps $F_k : \Omega \rightarrow O_k$ and not their values $F_k(\xi)$ at a given point $\xi \in \Omega$.*

iv) A Fréchet space is a Hausdorff, complete, topological space endowed with a countable family of semi-norms (Maths.971). It is locally convex and metric. This, quite complicated, mathematical definition is required because we will prove a theorem, and as usual in Mathematics we need to be precise in stating the conditions of its validity.

Are Fréchet spaces :

- any Banach vector space : the spaces of bounded functions, the spaces $L^p(E, \mu, \mathbb{C})$ of integrable functions on a measured space (E, μ) (Maths.2270), the spaces $L^p(M, \mu, E)$ of integrable sections of a vector bundle (valued in a Banach E) (Maths.2276)
- the spaces of continuously differentiable sections on a vector bundle (Maths.2310), the spaces of differentiable functions on a manifold (Maths.2314).

A topological vector space is separable if it has a dense countable subset (Maths.590) which, for a Fréchet space, is equivalent to be second countable (Maths.698). A totally bounded ($\forall r > 0$ there is a finite number of balls which cover V), or a connected locally compact Fréchet space, is separable (Maths.702, 703). The spaces $L^p(\mathbb{R}^n, dx, \mathbb{C})$ of integrable functions for $1 \leq p < \infty$, the spaces of continuous functions on a compact domain, are separable (Lieb).

Thus this somewhat complicated specification encompasses most of the usual cases.

In the following of this book we will see examples of these spaces : they are mostly maps : $X : \Omega \rightarrow E$ from a relatively compact subset Ω of a manifold M to a finite dimensional vector space, endowed with a norm. Then the space of maps such that $\int_{\Omega} \|X(m)\| \varpi(m) < \infty$ where ϖ is a measure on M (a volume measure) is an infinite dimensional, separable, Fréchet space.

v) The condition iv addresses the case when the variables are defined over

connected domains. But it implicitly tells that any other set of variables which represent the same phenomena are deemed compatible with the model. This point is addressed more precisely in another section, with the change of variables.

The set of all potential states of the system is then given by the set $S = \left\{ (X_k)_{k=1}^N, X_k \in O_k \right\}$. If there is some relation between the variables, stated by a physical law or theory, its consequence is to restrict the domain in which the state of the system will be found, but as said before we stay at the step before any experiment, so O_k represents the set of all possible values of X_k .

Theorem 2 *For any system represented by a model meeting the conditions 1, there is a separable, infinite dimensional, Hilbert space H , defined up to isomorphism, such that \mathcal{S} can be embedded as an open subset $\Omega \subset H$ which contains 0 and a convex subset.*

Proof. i) Each value of the set \mathcal{S} of variables defines a state of the system, denoted X , belonging to the product $O = \prod_{k=1}^N O_k \subset V = \prod_{k=1}^N V_k$. The couple (O, X) , together with the property iv) defines the structure of a Fréchet manifold M on the set \mathcal{S} , modelled on the Fréchet space $V = \prod_{k=1}^N V_k$. The coordinates are the values $(x_k)_{k=1}^N$ of the functions X_k . This manifold is infinite dimensional. Any Fréchet space is metric, so V is a metric space, and M is metrizable.

ii) As M is a metrizable manifold, modelled on an infinite dimensional separable Fréchet space, the Henderson's theorem (Henderson - corollary 5, Maths.1386) states that it can be embedded as a open subset Ω of an infinite dimensional separable Hilbert space H , defined up to isomorphism. Moreover this structure is smooth, the set $H - \Omega$ is homeomorphic to H , the border $\partial\Omega$ is homeomorphic to Ω and its closure $\overline{\Omega}$.

iii) Translations by a field vector are isometries. Let us denote $\langle \cdot \rangle_H$ the scalar product on H (this is a bilinear symmetric positive definite form). The map : $\Omega \rightarrow \mathbb{R} :: \langle \psi, \psi \rangle_H$ is bounded from below and continuous, so it has a minimum (possibly not unique) ψ_0 in Ω . By translation of H with ψ_0 we can define an isomorphic structure, and then assume that 0 belongs to Ω . There is a largest convex subset of H contained in Ω , defined as the intersection of all the convex subset contained in Ω . Its interior is an open convex subset C . It is not empty : because 0 belongs to Ω which is open in H , there is an open ball $B_0 = (0, r)$ contained in Ω . ■

So the state of the system can be represented by a single vector ψ in a Hilbert space.

From a practical point of view, often V itself can be taken as the product of Hilbert spaces, notably of square summable functions such as $L^2(\mathbb{R}, dt)$ which are separable Hilbert spaces and then the proposition is obvious.

If the variables belong to an open O' such that $O \subset O'$ we would have the same Hilbert space, and an open Ω' such that $\Omega \subset \Omega'$. V is open so we have a largest open $\Omega_V \subset H$ which contains all the Ω .

Notice that this is a real vector space.

The interest of Hilbert spaces lies with Hilbertian basis, and we now see how to relate such basis of H with a basis of the vector space V . It will enable us to show a linear chart of the manifold M .

2.1.4 Basis

Theorem 3 For any basis $(e_i)_{i \in I}$ of V contained in O , there are unique families $(\varepsilon_i)_{i \in I}, (\phi_i)_{i \in I}$ of independent vectors of H , a linear isometry $\Upsilon : V \rightarrow H$ such that :

$$\begin{aligned} \forall X \in O : \Upsilon(X) &= \sum_{i \in I} \langle \phi_i, \Upsilon(X) \rangle_H \varepsilon_i \in \Omega \\ \forall i \in I : \varepsilon_i &= \Upsilon(e_i) \\ \forall i, j \in I : \langle \phi_i, \varepsilon_j \rangle_H &= \delta_{ij} \\ &\text{and } \Upsilon \text{ is a compatible chart of } M. \end{aligned}$$

Proof. i) Let $(e_i)_{i \in I}$ be a basis of V such that $e_i \in O$ and $V_0 = \text{Span}(e_i)_{i \in I}$. Thus $O \subset V_0$.

Any vector of V_0 reads : $X = \sum_{i \in I} x_i e_i$ where only a finite number of x_i are non null. Or equivalently the following map is bijective :

$$\pi_V : V_0 \rightarrow \mathbb{R}_0^I :: \pi_V \left(\sum_{i \in I} x_i e_i \right) = x = (x_i)_{i \in I}$$

where the set $\mathbb{R}_0^I \subset \mathbb{R}^I$ is the subset of maps $I \rightarrow \mathbb{R}$ such that only a finite number of components x_i are non null.

(O, X) is an atlas of the manifold M and M is embedded in H , let us denote $\Xi : O \rightarrow \Omega$ a homeomorphism accounting for this embedding.

The inner product on H defines a positive kernel :

$$K : H \times H \rightarrow \mathbb{R} :: K(\psi_1, \psi_2) = \langle \psi_1, \psi_2 \rangle_H$$

Then $K_V : O \times O \rightarrow \mathbb{R} :: K_V(X, Y) = K(\Xi(X), \Xi(Y))$ defines a positive kernel on O (Math.1196).

K_V defines a definite positive symmetric bilinear form on V_0 , denoted $\langle \rangle_V$, by :

$$\left\langle \sum_{i \in I} x_i e_i, \sum_{i \in I} y_i e_i \right\rangle_V = \sum_{i, j \in I} x_i y_j K_{ij} \text{ with } K_{ij} = K_V(e_i, e_j)$$

which is well defined because only a finite number of monomials $x_i y_j$ are non null. It defines a norm on V_0 .

ii) Let : $\varepsilon_i = \Xi(e_i) \in \Omega$ and $H_0 = \text{Span}(\varepsilon_i)_{i \in I}$ the set of finite linear combinations of vectors $(\varepsilon_i)_{i \in I}$. It is a vector subspace (Math.901) of H . The family $(\varepsilon_i)_{i \in I}$ is linearly independent, because, for any finite subset J of I , the determinant

$$\det [\langle \varepsilon_i, \varepsilon_j \rangle_H]_{i, j \in J} = \det [K_V(e_i, e_j)]_{i, j \in J} \neq 0.$$

Thus $(\varepsilon_i)_{i \in I}$ is a non Hilbertian basis of H_0 .

H_0 can be defined similarly by the bijective map :

$$\pi_H : H_0 \rightarrow \mathbb{R}_0^I :: \pi_H \left(\sum_{i \in I} y_i \varepsilon_i \right) = y = (y_i)_{i \in I}$$

iii) By the Gram-Schmidt procedure (which works for infinite sets of vectors) it is always possible to built an orthonormal basis $(\tilde{\varepsilon}_i)_{i \in I}$ of H_0 starting with the

vectors $(\varepsilon_i)_{i \in I}$ indexed on the same set I (as H is separable I can be assimilated to \mathbb{N}).

$\ell^2(I) \subset \mathbb{R}^I$ is the set of families $y = (y_i)_{i \in I} \subset \mathbb{R}^I$ such that :

$$\sup \left(\sum_{i \in J} (y_i)^2 \right) < \infty \text{ for any countable subset } J \text{ of } I.$$

$$\mathbb{R}_0^I \subset \ell^2(I)$$

The map : $\chi : \ell^2(I) \rightarrow H_1 :: \chi(y) = \sum_{i \in I} y_i \tilde{\varepsilon}_i$ is an isomorphism to the closure $H_1 = \overline{Span(\tilde{\varepsilon}_i)_{i \in I}} = \overline{H_0}$ of H_0 in H (Math.1121). H_1 is a closed vector subspace of H , so it is a Hilbert space. The linear span of $(\tilde{\varepsilon}_i)_{i \in I}$ is dense in H_1 , so it is a Hilbertian basis of H_1 (Math.1122).

Let $\pi : H \rightarrow H_1$ be the orthogonal projection on H_1 : $\|\psi - \pi(\psi)\|_H = \min_{u \in H_1} \|\psi - u\|_H$ then :

$$\psi = \pi(\psi) + o(\psi) \text{ with } o(\psi) \in H_1^\perp \text{ which implies : } \|\psi\|^2 = \|\pi(\psi)\|^2 + \|o(\psi)\|^2$$

There is a open convex subset, containing 0, which is contained in Ω so there is $r > 0$ such that :

$$\|\psi\| < r \Rightarrow \psi \in \Omega \text{ and as } \|\psi\|^2 = \|\pi(\psi)\|^2 + \|o(\psi)\|^2 < r^2$$

$$\text{then } \|\psi\| < r \Rightarrow \pi(\psi), o(\psi) \in \Omega$$

$$o(\psi) \in H_1^\perp, H_0 \subset H_1 \Rightarrow o(\psi) \in H_0^\perp$$

$$\Rightarrow \forall i \in I : \langle \varepsilon_i, o(\psi) \rangle_H = 0 = K_V(\Xi^{-1}(\varepsilon_i), \Xi^{-1}(o(\psi))) = K_V(e_i, \Xi^{-1}(o(\psi)))$$

$$\Rightarrow \Xi^{-1}(o(\psi)) = 0 \Rightarrow o(\psi) = 0$$

$$H_1^\perp = 0 \text{ thus } H_1 \text{ is dense in } H \text{ (Math.1115), and as it is closed : } H_1 = H$$

$(\tilde{\varepsilon}_i)_{i \in I}$ is a Hilbertian basis of H and

$$\forall \psi \in H : \psi = \sum_{i \in I} \langle \tilde{\varepsilon}_i, \psi \rangle_H \tilde{\varepsilon}_i \text{ with } \sum_{i \in I} |\langle \tilde{\varepsilon}_i, \psi \rangle_H|^2 < \infty$$

$$\Leftrightarrow (\langle \tilde{\varepsilon}_i, \psi \rangle_H)_{i \in I} \in \ell^2(I)$$

H_0 is the interior of H , it is the union of all open subsets contained in H , so $\Omega \subset H_0$

$H_0 = Span((\tilde{\varepsilon}_i)_{i \in I})$ thus the map :

$$\tilde{\pi}_H : H_0 \rightarrow \mathbb{R}_0^I :: \tilde{\pi}_H(\sum_{i \in I} \tilde{y}_i \tilde{\varepsilon}_i) = \tilde{y} = (\tilde{y}_i)_{i \in I}$$

is bijective and : $\tilde{\pi}_H(H_0) = \tilde{R}_0 \subset \mathbb{R}_0^I \subset \ell^2(I)$

$$\text{Moreover : } \forall \psi \in H_0 : \tilde{\pi}_H(\psi) = (\langle \tilde{\varepsilon}_i, \psi \rangle_H)_{i \in I} \in \mathbb{R}_0^I$$

Thus :

$$\forall X \in O : \Xi(X) = \sum_{i \in I} \langle \tilde{\varepsilon}_i, \Xi(X) \rangle_H \tilde{\varepsilon}_i \in \Omega$$

$$\text{and } \tilde{\pi}_H(\Xi(X)) = (\langle \tilde{\varepsilon}_i, \Xi(X) \rangle_H)_{i \in I} \in \tilde{R}_0$$

$$\forall i \in I, e_i \in O \Rightarrow \Xi(e_i) = \varepsilon_i = \sum_{j \in I} \langle \tilde{\varepsilon}_j, \varepsilon_i \rangle_H \tilde{\varepsilon}_j$$

$$\text{and } \tilde{\pi}_H(\varepsilon_i) = (\langle \tilde{\varepsilon}_j, \varepsilon_i \rangle_H)_{j \in I} \in \tilde{R}_0$$

$$\text{iv) Let be : } \tilde{e}_i = \Xi^{-1}(\tilde{\varepsilon}_i) \in V_0 \text{ and } \mathcal{L}_V \in GL(V_0; V_0) :: \mathcal{L}_V(e_i) = \tilde{e}_i$$

We have the following diagram :

$$\begin{array}{ccccc} & \Xi & & \mathcal{L}_H^{-1} & \\ e_i & \rightarrow & \varepsilon_i & \rightarrow & \tilde{\varepsilon}_i \\ & \searrow & & & \downarrow \\ & \mathcal{L}_V & \searrow & & \downarrow \\ & & & \searrow & \downarrow \\ & & & & \tilde{e}_i \end{array} \quad \Xi^{-1}$$

$$\langle \tilde{e}_i, \tilde{e}_j \rangle_V = \langle \Xi(\tilde{e}_i), \Xi(\tilde{e}_j) \rangle_H = \langle \tilde{\varepsilon}_i, \tilde{\varepsilon}_j \rangle_H = \delta_{ij}$$

So $(\tilde{e}_i)_{i \in I}$ is an orthonormal basis of V_0 for the scalar product K_V

$$\forall X \in V_0 : X = \sum_{i \in I} \tilde{x}_i \tilde{e}_i = \sum_{i \in I} \langle \tilde{e}_i, X \rangle_V \tilde{e}_i \text{ and } (\langle \tilde{e}_i, X \rangle_V)_{i \in I} \in \mathbb{R}_0^I$$

The coordinates of $X \in O$ in the basis $(\tilde{e}_i)_{i \in I}$ are $(\langle \tilde{e}_i, X \rangle_V)_{i \in I} \in \mathbb{R}_0^I$

The coordinates of $\Xi(X) \in H_0$ in the basis $(\tilde{\varepsilon}_i)_{i \in I}$ are $(\langle \tilde{\varepsilon}_i, \Xi(X) \rangle_H)_{i \in I} \in \mathbb{R}_0^I$

$$\langle \tilde{\varepsilon}_i, \Xi(X) \rangle_H = \langle \Xi(\tilde{e}_i), \Xi(X) \rangle_H = \langle \tilde{e}_i, X \rangle_V$$

Define the maps :

$$\tilde{\pi}_V : V_0 \rightarrow \mathbb{R}_0^I :: \tilde{\pi}_V \left(\sum_{i \in I} \tilde{x}_i \tilde{e}_i \right) = \tilde{x} = (\tilde{x}_i)_{i \in I}$$

$$\Upsilon : V_0 \rightarrow H_0 :: \Upsilon = \tilde{\pi}_H^{-1} \circ \tilde{\pi}_V^{-1}$$

which associates to each vector of V the vector of H with the same components in the orthonormal bases, then :

$$\forall X \in O : \Upsilon(X) = \Xi(X)$$

and Υ is a bijective, linear map, which preserves the scalar product, so it is continuous and is an isometry.

v) There is a bijective linear map : $\mathcal{L}_H \in GL(H_0; H_0)$ such that : $\forall i \in I : \varepsilon_i = \mathcal{L}_H(\tilde{\varepsilon}_i)$.

$(\tilde{\varepsilon}_i)_{i \in I}$ is a basis of H_0 thus $\varepsilon_i = \sum_{j \in I} [\mathcal{L}_H]_i^j \tilde{\varepsilon}_j$ where only a finite number of coefficients $[\mathcal{L}_H]_i^j$ is non null.

$$\text{Let us define : } \varpi_i : H_0 \rightarrow \mathbb{R} :: \varpi_i \left(\sum_{j \in I} \psi_j \varepsilon_j \right) = \psi_i$$

This map is continuous at $\psi = 0$ on H_0 :

$$\text{take } \psi \in H_0, \|\psi\| \rightarrow 0$$

$$\text{then } \psi = \sum_{i \in I} \langle \tilde{\varepsilon}_i, \psi \rangle_H \tilde{\varepsilon}_i \text{ and } \tilde{\psi}_j = \langle \tilde{\varepsilon}_j, \psi \rangle_H \rightarrow 0$$

$$\text{so if } \|\psi\| < r \text{ then } \|\psi\|^2 = \sum_{j \in I} |\tilde{\psi}_j|^2 < r^2 \text{ and } \forall j \in I : |\tilde{\psi}_j| < r$$

$$\psi_i = \sum_{j \in J} [\mathcal{L}_H]_i^j \tilde{\psi}_j \Rightarrow |\psi_i| < \varepsilon \sum_{j \in I} \max |[\mathcal{L}_H]_i^j| \text{ and } \left(|[\mathcal{L}_H]_i^j| \right)_{j \in I} \text{ is}$$

bounded $\Rightarrow |\psi_i| \rightarrow 0$

Thus ϖ_i is continuous and belongs to the topological dual H'_0 of H_0 . It can be extended as a continuous map $\overline{\varpi}_i \in H'$ according to the Hahn-Banach theorem (Maths.958). Because H is a Hilbert space, there is a vector $\phi_i \in H$ such that : $\forall \psi \in H : \overline{\varpi}_i(\psi) = \langle \phi_i, \psi \rangle_H$ so that :

$$\begin{aligned} \forall X \in O : \Upsilon(X) = \Xi(X) &= \sum_{i \in I} \psi_i \varepsilon_i \\ &= \sum_{i \in I} \langle \phi_i, \psi \rangle_H \varepsilon_i = \sum_{i \in I} \langle \phi_i, \Xi(X) \rangle_H \varepsilon_i \end{aligned}$$

$\forall i \in I :$

$$\Xi(e_i) = \varepsilon_i = \Upsilon(e_i) = \sum_{j \in I} \langle \phi_j, \varepsilon_i \rangle_H \varepsilon_j \Rightarrow \langle \phi_j, \varepsilon_i \rangle_H = \delta_{ij}$$

$$\Xi(\tilde{e}_i) = \sum_{j \in I} \langle \phi_j, \Xi(\tilde{e}_i) \rangle_H \varepsilon_j = \tilde{\varepsilon}_i = \sum_{j \in I} \langle \phi_j, \tilde{\varepsilon}_i \rangle_H \varepsilon_j$$

vi) The map $\Upsilon : O \rightarrow \Omega$ is a linear chart of M , using two orthonormal bases : it is continuous, bijective so it is an homeomorphism, and is obviously compatible with the chart Ξ . ■

Remarks

i) Because $(\tilde{\varepsilon}_i)_{i \in I}$ is a Hilbertian basis of the separable infinite dimensional Hilbert space H , I is a countable set which can be identified to \mathbb{N} . The assumption

about $(e_i)_{i \in I}$ is that it is a Hamel basis, which is the most general because any vector space has one. From the proposition above we see that this basis must be of cardinality \aleph_0 . Hamel bases of infinite dimensional normed vector spaces must be uncountable, however our assumption about V is that it is a Fréchet space, which is a metrizable but not a normed space. If V is a Banach vector space then, according to the Mazur theorem, it implies that there it has an infinite dimensional vector subspace W which has a Shauder basis: $\forall X \in W : X = \sum_{i \in I} x_i e_i$ where the sum is understood in the topological limit. Then the same reasoning as above shows that the closure of W is itself a Hilbert space. Moreover it has been proven that any separable Banach space is homeomorphic to a Hilbert space, and most of the applications will concern spaces of integrable functions (or sections of vector bundle endowed with a norm) which are separable Fréchet spaces.

One interesting fact is that we assume that the variables belong to an open subset O of V . The main concern is to allow for variables which can take values only in some bounded domain. But this assumption addresses also the case of a Banach vector space which is “hollowed out” : O can be itself a vector subspace (in an infinite dimensional vector space a vector subspace can be open), for instance generated by a countable subbasis of a Hamel basis, and we assume explicitly that the basis $(e_i)_{i \in I}$ belongs to O .

ii) For $O = V$ we have a largest open Ω_V and a linear map $\Upsilon : V \rightarrow \Omega_V$ with domain V .

iii) To each (Hamel) basis on V is associated a linear chart Υ of the manifold, such that a point of M has the same coordinates both in V and H . So Υ depends on the choice of the basis, and similarly the positive kernel K_V depends on the basis.

iv) In the proof we have introduced a map : $K_V : O \times O \rightarrow \mathbb{R} :: K_V(X, Y)$ which is not bilinear, but is definite positive in a precise way. It plays an important role in several following demonstrations. From a physical point of view it can be seen as related to the probability of transition between two states X, Y often used in QM.

2.1.5 Complex structure

The variables X and vector space V are real and H is a real Hilbert space. The condition that the vector space V is real is required only in Proposition 2 to prove the existence of a Hilbert space, because the Henderson’s theorem holds only for real structures. However, as it is easily checked, if H exists, all the following theorems hold even if H is a complex Hilbert space. This is specially useful when the space V over which the maps X are defined is itself a complex Hilbert space, as this is often the case.

Moreover it can be useful to endow H with the structure of a complex Hilbert space : the set does not change but one distinguishes real and imaginary components, and the scalar product is given by a Hermitian form. Notice that this is a convenience, not a necessity.

Theorem 4 Any real separable infinite dimensional Hilbert space can be endowed with the structure of a complex separable Hilbert space

Proof. H has a infinite countable Hilbertian basis $(\varepsilon_\alpha)_{\alpha \in \mathbb{N}}$ because it is separable.

A complex structure is defined by a linear map $J \in \mathcal{L}(H; H)$ such that $J^2 = -Id$. Then the operation $i \times \psi$ is defined by $i\psi = J(\psi)$.

Define :

$$J(\varepsilon_{2\alpha}) = \varepsilon_{2\alpha+1}; J(\varepsilon_{2\alpha+1}) = -\varepsilon_{2\alpha}$$

$$\forall \psi \in H : i\psi = J(\psi)$$

$$\text{So : } i(\varepsilon_{2\alpha}) = \varepsilon_{2\alpha+1}; i(\varepsilon_{2\alpha+1}) = -\varepsilon_{2\alpha}$$

The bases $\varepsilon_{2\alpha}$ or $\varepsilon_{2\alpha+1}$ are complex bases of H :

$$\psi = \sum_{\alpha} \psi^{2\alpha} \varepsilon_{2\alpha} + \psi^{2\alpha+1} \varepsilon_{2\alpha+1} = \sum_{\alpha} (\psi^{2\alpha} - i\psi^{2\alpha+1}) \varepsilon_{2\alpha}$$

$$= \sum_{\alpha} (-i\psi^{2\alpha} + \psi^{2\alpha+1}) \varepsilon_{2\alpha+1}$$

$$\|\psi\|^2 = \sum_{\alpha} |\psi^{2\alpha} - i\psi^{2\alpha+1}|^2$$

$$= \sum_{\alpha} |\psi^{2\alpha}|^2 + |\psi^{2\alpha+1}|^2 + i(-\overline{\psi^{2\alpha}} \psi^{2\alpha+1} + \psi^{2\alpha} \overline{\psi^{2\alpha+1}})$$

$$= \sum_{\alpha} |\psi^{2\alpha}|^2 + |\psi^{2\alpha+1}|^2 + i(-\psi^{2\alpha} \psi^{2\alpha+1} + \psi^{2\alpha} \psi^{2\alpha+1})$$

Thus $\varepsilon_{2\alpha}$ is a Hilbertian complex basis

H has a structure of complex vector space that we denote $H_{\mathbb{C}}$

The map $T : H \rightarrow H_{\mathbb{C}} : T(\psi) = \sum_{\alpha} (\psi^{2\alpha} - i\psi^{2\alpha+1}) \varepsilon_{2\alpha}$ is linear and continuous

The map $\overline{T} : H \rightarrow H_{\mathbb{C}} : \overline{T}(\psi) = \sum_{\alpha} (\psi^{2\alpha} + i\psi^{2\alpha+1}) \varepsilon_{2\alpha}$ is antilinear and continuous

$$\text{Define : } \gamma(\psi, \psi') = \langle \overline{T}(\psi), T(\psi') \rangle_H$$

γ is sesquilinear

$$\gamma(\psi, \psi') = \langle \sum_{\alpha} (\psi^{2\alpha} + i\psi^{2\alpha+1}) \varepsilon_{2\alpha}, \sum_{\alpha} (\psi'^{2\alpha} - i\psi'^{2\alpha+1}) \varepsilon_{2\alpha} \rangle_H$$

$$= \sum_{\alpha} (\psi^{2\alpha} + i\psi^{2\alpha+1}) (\psi'^{2\alpha} - i\psi'^{2\alpha+1})$$

$$= \sum_{\alpha} \psi^{2\alpha} \psi'^{2\alpha} + \psi^{2\alpha+1} \psi'^{2\alpha+1} + i(\psi^{2\alpha+1} \psi'^{2\alpha} - \psi^{2\alpha} \psi'^{2\alpha+1})$$

$$\gamma(\psi, \psi) = 0 \Rightarrow \langle \psi, \psi \rangle_H = 0 \Rightarrow \psi = 0$$

Thus γ is definite positive ■

2.1.6 Decomposition of the Hilbert space

V is the product $V = V_1 \times V_2 \dots \times V_N$ of vector spaces, thus the proposition implies that the Hilbert space H is also the direct product of Hilbert spaces $H_1 \times H_2 \dots \times H_N$ or equivalently $H = \bigoplus_{k=1}^N H_k$ where H_k are Hilbert vector subspaces of H . More precisely :

Theorem 5 If the model is comprised of N continuous variables $(X_k)_{k=1}^N$, each belonging to a separable Fréchet vector space V_k , then the real Hilbert space H of states of the system is the Hilbert sum of N Hilbert space $H = \bigoplus_{k=1}^N H_k$ and any vector ψ representing a state of the system is uniquely the sum of N vectors ψ_k , each image of the value of one variable X_k in the state ψ

Proof. By definition $V = \prod_{k=1}^N V_k$. The set $V_k^0 = \{0, \dots, V_k, \dots, 0\} \subset V$ is a vector subspace of V . A basis of V_k^0 is a subfamily $(e_i)_{i \in I_k}$ of a basis $(e_i)_{i \in I}$ of V . V_k^0 has for image by the continuous linear map Υ a closed vector subspace H_k of H . Any vector X of V reads : $X \in \prod_{k=1}^N V_k : X = \sum_{k=1}^N \sum_{i \in I_k} x^i e_i$ and it has for image by $\Upsilon : \psi = \Upsilon(X) = \sum_{k=1}^N \sum_{i \in I_k} x^i \varepsilon_i = \sum_{k=1}^N \psi_k$ with $\psi_k \in H_k$. This decomposition of $\Upsilon(X)$ is unique.

Conversely, the family $(e_i)_{i \in I_k}$ has for image by Υ the set $(\varepsilon_i)_{i \in I_k}$ which are linearly independent vectors of H_k . It is always possible to build an orthonormal basis $(\tilde{\varepsilon}_i)_{i \in I_k}$ from these vectors as done previously. H_k is a closed subspace of H , so it is a Hilbert space. The map : $\hat{\pi}_k : \ell^2(I_k) \rightarrow H_k :: \hat{\pi}_k(x) = \sum_{i \in I_k} x^i \tilde{\varepsilon}_i$ is an isomorphism of Hilbert spaces and : $\forall \psi \in H_k : \psi = \sum_{i \in I_k} \langle \tilde{\varepsilon}_i, \psi \rangle_H \tilde{\varepsilon}_i$.

$$\forall \psi_k \in H_k, \psi_l \in H_l, k \neq l : \langle \psi_k, \psi_l \rangle_H = \langle \Upsilon^{-1}(\psi_k), \Upsilon^{-1}(\psi_l) \rangle_E = 0$$

Any vector $\psi \in H$ reads : $\psi = \sum_{k=1}^N \pi_k(\psi)$ with the orthogonal projection $\pi_k : H \rightarrow H_k :: \pi_k(\psi) = \sum_{i \in I_k} \langle \tilde{\varepsilon}_i, \psi \rangle_H \tilde{\varepsilon}_i$ so H is the Hilbert sum of the H_k ■

As a consequence the definite positive kernel of (V, Υ) decomposes as :

$$\begin{aligned} & K((X_1, \dots, X_N), (X'_1, \dots, X'_N)) \\ &= \sum_{k=1}^N K_k(X_k, X'_k) \\ &= \sum_{k=1}^N \langle \Upsilon(X_k), \Upsilon(X'_k) \rangle_{H_k} \end{aligned}$$

This decomposition comes handy when we have to translate relations between variables into relations between vector states, notably if they are linear. But it requires that we keep the real Hilbert space structure.

2.1.7 Discrete variables

It is common in a model to have discrete variables $(D_k)_{k=1}^K$, taking values in a finite discrete set. They correspond to different cases:

i) the discrete variables identify different elementary systems (such as different populations of particles) which coexist simultaneously in the same global system, follow different rules of behavior, but interact together. We will see later how to deal with these cases (tensorial product).

ii) the discrete variables identify different populations, whose interactions are not relevant. Actually one could consider as many different systems but, by putting them together, one increases the size of the samples of data and improve the statistical estimations. They are not of great interest here, in a study of formal models.

iii) the discrete variables represent different kinds of behaviors, which cannot be strictly identified with specific populations. Usually a discrete variable is then used as a proxy for a quantitative parameter which tells how close the system is from a specific situation.

We will focus on this third case. The system is represented as before by quantitative variables X , whose possible values belong to some set M , which

has the structure of an infinite dimensional manifold. The general idea in the third case is that the possible states of the system can be regrouped in two distinct subsets. That we formalize in the following assumption : the set O of possible states of the system has two connected components O_1, O_2

Theorem 6 *If the condition of the theorem 2 are met, and the set O of possible states of the system has two connected components O_1, O_2 then there is a continuous function $f : H \rightarrow [0, 1]$ such that $f(\Upsilon(X)) = 1$ in O_1 and $f(\Upsilon(X)) = 0$ in O_2*

Proof. The connected components O_1, O_2 of a topological space are closed, so O_1, O_2 are disjoint and both open and closed in V (Maths.624). Using a linear continuous map Υ then Ω has itself two connected components, $\Omega_1 = \Upsilon^{-1}(O_1), \Omega_2 = \Upsilon^{-1}(O_2)$ both open and closed, and disjoint. H is metric, so it is normal (Maths.705). Ω_1, Ω_2 are disjoint and closed in H . Then, by the Urysohn's Theorem (Maths.596) there is a continuous function f on H valued in $[0, 1]$ such that $f(\psi) = 1$ in H_1 and $f(\psi) = 0$ in H_2 . ■

The set of continuous, bounded functions is a Banach vector space, so it is always possible, in these conditions, to replace a discrete variable by a quantitative variable with the same features.

2.2 OBSERVABLES

The key point in the conditions 1 above is that the variables are maps, which take an infinite number of values (usually non countable). So the variables would require the same number of data to be totally known, which is impossible. The physicist estimates the variable by statistical methods. But any practical method involves a first step : the scope of all maps is reduced from V to a smaller subset W , so that any map of W can be characterized by a finite number of parameters. The procedure sums up to replace X by another variable $\Phi(X)$ that we will call an observable, which is then estimated from a finite batch of data. The mechanism of estimating the variables $X \subset V$ is then the following :

- the observer collects data, as a set $Y = \{x_p\}_{p=1}^N$ of values assumed to be taken by the variable X , in the mathematical format fitted to X (scalars, vectors,...for different values of the arguments)

- he proceeds to the estimation \hat{X} of the map $\Phi(X)$ by statistical adjustment to the data $\{x_p\}_{p=1}^N$. Because there are a finite number of parameters (the coordinates of $\Phi(X)$ in W) this is possible

- the estimation is : $\hat{X} = \varphi(Y) \in W$: this is a map which is a simplified version of X .

The procedure of the replacement of X by $\Phi(X)$, called the choice of a **specification**, is done by the physicist, and an observable is not unique. However we make three general assumptions about Φ :

Definition 7 *i) an **observable** is a linear map : $\Phi \in L(V; V)$*

ii) the range of an observable is a finite dimensional vector W subspace of V : $W \subset V, \dim \Phi(W) < \infty$

iii) $\forall X \in O, \Phi(X)$ is an admissible value, that is $\Phi(O) \subset O$.

Using the linear chart Υ given by any basis, to Φ one can associate a map :

$$\hat{\Phi} : H \rightarrow H :: \hat{\Phi} = \Upsilon \circ \Phi \circ \Upsilon^{-1} \tag{2.1}$$

and $\hat{\Phi}$ is an operator on H . And conversely.

The image of W by Υ is a finite dimensional vector subspace $H_\Phi = \Upsilon(W)$ of H , so it is closed and a Hilbert space : $\hat{\Phi} \in \mathcal{L}(H; H_\Phi)$

$$\begin{array}{ccccccc}
 & & \Phi & & & & \\
 & V & \rightarrow & \rightarrow & \rightarrow & W & \\
 & \downarrow & & & & \downarrow & \\
 \Upsilon & \downarrow & & & & \downarrow & \Upsilon \\
 & \downarrow & & \hat{\Phi} & & \downarrow & \\
 & H & \rightarrow & \rightarrow & \rightarrow & H_\Phi &
 \end{array}$$

2.2.1 Primary observables

The simplest specification for an observable is, given a basis $(e_i)_{i \in I}$, to define Φ as the projection on the subspace spanned by a finite number of vectors of the basis. For instance if X is a function $X(t)$ belonging to some space such as : $X(t) = \sum_{n \in \mathbb{N}} a_n e_n(t)$ where $e_n(t)$ are fixed functions, then a primary observable would be $Y_J(X(t)) = \sum_{n=0}^N a_n e_n(t)$ meaning that the components $(a_n)_{n > N}$ are discarded and the data are used to compute $(a_n)_{n=0}^N$. To stay at the most general level, we define :

Definition 8 A *primary observable* $\Phi = Y_J$ is the projection of $X = \{X_k, k = 1 \dots N\}$ on the vector subspace V_J spanned by the vectors $(e_i)_{i \in J} \equiv (e_i^k)_{i \in J_k}$ where

$$J = \prod_{k=1}^N J_k \subset I = \prod_{k=1}^N I_k \text{ is a finite subset of } I \text{ and } (\varepsilon_i)_{i \in I} = \prod_{k=1}^N (e_i^k)_{i \in I_k} \text{ is a basis of } V.$$

So the procedure can involve simultaneously several variables. It requires the choice of a finite set of independent vectors of V .

Theorem 9 To any primary observable Y_J is associated uniquely a self-adjoint, compact, trace-class operator \hat{Y}_J on $H : Y_J = \Upsilon^{-1} \circ \hat{Y}_J \circ \Upsilon$ such that the measure $Y_J(X)$ of the primary observable Y_J , if the system is in the state $X \in O$, is

$$Y_J(X) = \sum_{i \in I} \langle \phi_i, \hat{Y}_J(\Upsilon(X)) \rangle_H e_i$$

Proof. i) We use the notations and definitions of the previous section. The family of variables $X = (X_k)_{k=1}^N$ define the charts : $\Xi : O \rightarrow \Omega$ and the basis $(e_i)_{i \in I}$ defines the bijection $\Upsilon : V \rightarrow H$

$$\forall X = \sum_{i \in I} x_i e_i \in O :$$

$$\Upsilon(X) = \sum_{i \in I} x_i \Upsilon(e_i) = \sum_{i \in I} x_i \varepsilon_i = \sum_{i \in I} \langle \phi_i, \Upsilon(X) \rangle_H \varepsilon_i$$

$$\Leftrightarrow x_i = \langle \phi_i, \Upsilon(X) \rangle_H$$

$$\forall i, j \in I : \langle \phi_i, \varepsilon_j \rangle_H = \delta_{ij}$$

ii) The primary observable Y_J is the map :

$$Y_J : V \rightarrow V_J :: Y_J(X) = \sum_{j \in J} x_j e_j$$

This is a projection : $Y_J^2 = Y_J$

$Y_J(X) \in O$ so it is associated to a vector of H :

$$\Upsilon(Y_J(X)) = \Upsilon\left(\sum_{j \in J} x_j e_j\right) = \sum_{j \in J} \langle \phi_j, \Upsilon(Y_J(X)) \rangle_H \varepsilon_j$$

$$= \sum_{j \in J} \langle \phi_j, \Upsilon(X) \rangle_H \varepsilon_j$$

iii) $\forall X \in O : \Upsilon(Y_J(X)) \in H_J$ where H_J is the vector subspace of H spanned by $(\varepsilon_j)_{j \in J}$. It is finite dimensional, thus it is closed in H . There is a unique map (Math.1111) :

$$\hat{Y}_J \in \mathcal{L}(H; H) :: \hat{Y}_J^2 = \hat{Y}_J, \hat{Y}_J = \hat{Y}_J^*$$

\widehat{Y}_J is the orthogonal projection from H onto H_J . It is linear, self-adjoint, and compact because its range is a finite dimensional vector subspace. As a projection : $\|\widehat{Y}_J\| = 1$.

\widehat{Y}_J is a Hilbert-Schmidt operator (Maths.1143) : take the Hilbertian basis $\tilde{\varepsilon}_i$ in H :

$$\sum_{i \in I} \|\widehat{Y}_J(\tilde{\varepsilon}_i)\|^2 = \sum_{ij \in J} |\langle \phi_j, \tilde{\varepsilon}_i \rangle|^2 \|\varepsilon_j\|^2 = \sum_{j \in J} \|\phi_j\|^2 \|\varepsilon_j\|^2 < \infty$$

\widehat{Y}_J is a trace class operator (Maths.1147) with trace $\dim H_J$

$$\sum_{i \in I} \langle \widehat{Y}_J(\tilde{\varepsilon}_i), \tilde{\varepsilon}_i \rangle = \sum_{ij \in J} \langle \phi_j, \tilde{\varepsilon}_i \rangle \langle \varepsilon_j, \tilde{\varepsilon}_i \rangle = \sum_{j \in J} \langle \phi_j, \varepsilon_j \rangle = \sum_{j \in J} \delta_{jj} = \dim H_J$$

$$\text{iv) } \forall \psi \in H_J : \widehat{Y}_J(\psi) = \psi$$

$$\forall X \in O : \Upsilon(Y_J(X)) \in H_J$$

$$\forall X \in O : \Upsilon(Y_J(X)) = \widehat{Y}_J(\Upsilon(X)) \Leftrightarrow Y_J(X) = \Upsilon^{-1} \circ \widehat{Y}_J(\Upsilon(X)) \Leftrightarrow Y_J = \Upsilon^{-1} \circ \widehat{Y}_J \circ \Upsilon$$

$$\text{v) The value of the observable reads : } Y_J(X) = \sum_{i \in I} \langle \phi_i, \widehat{Y}_J(\Upsilon(X)) \rangle e_i$$

■

2.2.2 von Neumann algebras

Taking the axioms of QM as a starting point, one can define a system itself by the set of its observables. In this framework the focus has been given to von Neumann algebras, which are special sets of operators on a Hilbert space. These mathematical models are convenient to explore further the behavior of systems or some sensitive issues such as the continuity of the operators. But this approach has a fundamental drawback : it leads further from an understanding of the physical foundations of the theory itself. To tell that a system should be represented by a von Neumann algebra does not explain more why a state should be represented in a Hilbert space at the beginning. But as we will see now, von Neumann algebras appear naturally in our framework.

There is a bijective correspondence between the projections, meaning the maps $P \in \mathcal{L}(H; H) : P^2 = P, P = P^*$ and the closed vector subspaces of H (Maths.1111). Then P is the orthogonal projection on the vector subspace. So the operators \widehat{Y}_J for any finite subset J of I are the orthogonal projections on the finite dimensional, and thus closed, vector subspace H_J spanned by $(\varepsilon_j)_{j \in J}$.

We will enlarge the family of primary observables in several steps, in keeping the same basis $(e_i)_{i \in I}$ of V .

1. For any given basis $(e_i)_{i \in I}$ of V , we extend the definition of these operators \widehat{Y}_J to any finite or infinite, subset of I by taking \widehat{Y}_J as the orthogonal projection on the closure $\overline{H_J}$ in H of the vector subspace H_J spanned by $(\varepsilon_j)_{j \in J}$. $\overline{H_J} = \overline{\text{Span}(\varepsilon_j)_{j \in J}}$.

Theorem 10 *The operators $\{\widehat{Y}_J\}_{J \subset I}$ are self-adjoint and commute*

Proof. Because they are projections the operators \widehat{Y}_J are such that : $\widehat{Y}_J^2 = \widehat{Y}_J, \widehat{Y}_J^* = \widehat{Y}_J$

\widehat{Y}_J has for eigen values :

1 for $\psi \in \overline{H_J}$

0 for $\psi \in (\overline{H_J})^\perp$

For any subset J of I , by the Gram-Schmidt procedure one can built an orthonormal basis $(\tilde{\varepsilon}_i)_{i \in J}$ of H_J starting with the vectors $(\varepsilon_i)_{i \in J}$ and an orthonormal basis $(\tilde{\varepsilon}_i)_{i \in J^c}$ of H_{J^c} starting with the vectors $(\varepsilon_i)_{i \in J^c}$

Any vector $\psi \in H$ can be written :

$$\psi = \sum_{j \in I} x_j \tilde{\varepsilon}_j = \sum_{j \in J} x_j \tilde{\varepsilon}_j + \sum_{j \in J^c} x_j \tilde{\varepsilon}_j \text{ with } (x_j)_{j \in I} \in \ell^2(I)$$

$\overline{H_J}$ is defined as $\sum_{j \in J} x_j \tilde{\varepsilon}_j$ with $(x_j)_{j \in J} \in \ell^2(J)$ and similarly $\overline{H_{J^c}}$ is defined as $\sum_{j \in J^c} x_j \tilde{\varepsilon}_j$ with $(x_j)_{j \in J^c} \in \ell^2(J^c)$

$$\text{So } \widehat{Y}_J \text{ can be defined as : } \widehat{Y}_J \left(\sum_{j \in I} x_j \tilde{\varepsilon}_j \right) = \sum_{j \in J} x_j \tilde{\varepsilon}_j$$

For any subsets $J_1, J_2 \subset I$:

$$\widehat{Y}_{J_1} \circ \widehat{Y}_{J_2} = \widehat{Y}_{J_1 \cap J_2} = \widehat{Y}_{J_2} \circ \widehat{Y}_{J_1}$$

$$\widehat{Y}_{J_1 \cup J_2} = \widehat{Y}_{J_1} + \widehat{Y}_{J_2} - \widehat{Y}_{J_1 \cap J_2} = \widehat{Y}_{J_1} + \widehat{Y}_{J_2} - \widehat{Y}_{J_1} \circ \widehat{Y}_{J_2}$$

So the operators commute. ■

2. Let us define $W = \text{Span} \left\{ \widehat{Y}_i \right\}_{i \in I}$ the vector subspace of $\mathcal{L}(H; H)$ comprised of finite linear combinations of \widehat{Y}_i (as defined in 1 above). The elements $\left\{ \widehat{Y}_i \right\}_{i \in I}$ are linearly independent and constitute a basis of W .

The operators $\widehat{Y}_j, \widehat{Y}_k$ are mutually orthogonal for $j \neq k$:

$$\widehat{Y}_j \circ \widehat{Y}_k(\psi) = \langle \phi_k, \psi \rangle \langle \phi_j, \varepsilon_k \rangle \varepsilon_j = \langle \phi_k, \psi \rangle \delta_{jk} = \delta_{jk} \widehat{Y}_j(\psi)$$

Let us define the scalar product on W :

$$\left\langle \sum_{i \in I} a_i \widehat{Y}_i, \sum_{i \in I} b_i \widehat{Y}_i \right\rangle_W = \sum_{i \in I} a_i b_i$$

$$\left\| \sum_{i \in I} a_i \widehat{Y}_i \right\|_W^2 = \sum_{i \in I} a_i^2 \left\| \widehat{Y}_i \right\|_W^2 = \sum_{i \in I} a_i^2$$

W is isomorphic to \mathbb{R}_0^I and its closure in $\mathcal{L}(H; H)$: $\overline{W} = \overline{\text{Span} \left\{ \widehat{Y}_i \right\}_{i \in I}}$ is isomorphic to $\ell^2(I)$, and has the structure of a Hilbert space with :

$$\overline{W} = \left\{ \sum_{i \in I} a_i \widehat{Y}_i, (a_i)_{i \in I} \in \ell^2(I) \right\}$$

3. Let us define A as the algebra generated by any finite linear combination or products of elements \widehat{Y}_J, J finite or infinite, and \overline{A} as the closure of A in $\mathcal{L}(H; H)$: $\overline{A} = \overline{\text{Span} \left\{ \widehat{Y}_J \right\}_{J \subset I}}$ with respect to the strong topology, that is in norm.

Theorem 11 \overline{A} is a commutative von Neumann algebra of $\mathcal{L}(H, H)$

Proof. It is obvious that A is a *-subalgebra of $\mathcal{L}(H, H)$ with unit element $Id = \widehat{Y}_I$.

Because its generators are projections, \overline{A} is a von Neumann algebra (Maths.1190).

The elements of $A = Span \left\{ \widehat{Y}_J \right\}_{J \subset I}$ that is of finite linear combination of \widehat{Y}_J commute

$$Y, Z \in \overline{A} \Rightarrow \exists (Y_n)_{n \in \mathbb{N}}, (Z_n)_{n \in \mathbb{N}} \in A^{\mathbb{N}} : Y_n \rightarrow_{n \rightarrow \infty} Y, Z_n \rightarrow_{n \rightarrow \infty} Z$$

The composition is a continuous operation.

$$Y_n \circ Z_n = Z_n \circ Y_n \Rightarrow \lim (Y_n \circ Z_n) = \lim (Z_n \circ Y_n) = \lim Y_n \circ \lim Z_n = \lim Z_n \circ \lim Y_n = Z \circ Y = Y \circ Z$$

So \overline{A} is commutative.

\overline{A} is identical to the bicommutant of its projections, that is to \overline{A}'' (Maths.1189)

■

This result is of interest because commutative von Neumann algebras are classified : they are isomorphic to the space of functions $f \in L^\infty(E, \mu)$ acting by pointwise multiplication $\varphi \rightarrow f\varphi$ on functions $\varphi \in L^2(E, \mu)$ for some set E and measure μ (not necessarily absolutely continuous). They are the topic of many studies, notably in ergodic theory. The algebra \overline{A} depends on the choice of a basis $(e_i)_{i \in I}$ and, as can be seen in the formulation through $(\tilde{\varepsilon}_i)_{i \in I}$, is defined up to a unitary transformation.

In the formalization of QM, it is usual to define a system by a von Neumann algebra of operators on a Hilbert space. We see here how such an algebra appears naturally. However the algebra \overline{A} is commutative, and this property is the consequence of the choice of a unique basis $(e_i)_{i \in I}$. It would not hold for primary observables defined through different bases : they do not even constitute an algebra. Any von Neumann algebra is the closure of the linear span of its projections (Maths.1190), and any projection can be defined through a basis, thus one can say that the “observables” (with their usual definition) of a system are the collection of all primary observables (as defined here) for all bases of V . This is a crucial issue in the axiomatic interpretation of QM, but the concept of observables introduced here allows us to deal with this issue and understand how probabilities enter the picture. But, before that, we need to see what can be said about more general observables, which are not just primary.

2.2.3 Secondary observables

Beyond primary observables, general observables Φ can be studied using spectral theory.

1. A spectral measure defined on a measurable space E with σ -algebra σ_E and acting on the Hilbert space H is a map : $P : \sigma_E \rightarrow \mathcal{L}(H; H)$ such that (Maths.1240) :

- i) $P(\varpi)$ is a projection
- ii) $P(E) = Id$

iii) $\forall \psi \in H$ the map: $\varpi \rightarrow \langle P(\varpi)\psi, \psi \rangle_H = \|P(\varpi)\psi\|^2$ is a finite positive measure on (E, σ_E) .

One can show (Maths.1242) that there is a bijective correspondence between the spectral measures on H and the maps: $\chi: \sigma_E \rightarrow H$ such that:

- i) $\chi(\varpi)$ is a closed vector subspace of H
- ii) $\chi(E) = H$
- iii) $\forall \varpi, \varpi' \in \sigma_E, \varpi \cap \varpi' = \emptyset: \chi(\varpi) \cap \chi(\varpi') = \{0\}$

then $P(\varpi)$ is the orthogonal projection on $\chi(\varpi)$, denoted: $\widehat{\pi}_{\chi(\varpi)}$

Thus, for any fixed $\psi \neq 0 \in H$ the function $\widehat{\chi}_\psi: \sigma_E \rightarrow \mathbb{R} :: \widehat{\chi}_\psi(\varpi) = \frac{\langle \widehat{\pi}_{\chi(\varpi)}\psi, \psi \rangle}{\|\psi\|^2} = \frac{\|\widehat{\pi}_{\chi(\varpi)}\psi\|^2}{\|\psi\|^2}$ is a probability law on (E, σ_E) .

2. An application of standard theorems on spectral measures (Maths.1243, 1245) tells that, for any bounded measurable function $f: E \rightarrow \mathbb{R}$, the spectral integral: $\int_E f(\xi) \widehat{\pi}_{\chi(\xi)}$ defines a continuous operator $\widehat{\Phi}_f$ on H . $\widehat{\Phi}_f$ is such that:

$$\forall \psi, \psi' \in H: \langle \widehat{\Phi}_f(\psi), \psi' \rangle = \int_E f(\xi) \langle \widehat{\pi}_{\chi(\xi)}(\psi), \psi' \rangle$$

And conversely (Math.1252), for any continuous normal operator $\widehat{\Phi}$ on H , that is such that:

$$\widehat{\Phi} \in \mathcal{L}(H; H): \widehat{\Phi} \circ \widehat{\Phi}^* = \widehat{\Phi}^* \circ \widehat{\Phi} \text{ with the adjoint } \widehat{\Phi}^*$$

there is a unique spectral measure P on $(\mathbb{R}, \sigma_{\mathbb{R}})$ such that $\widehat{\Phi} = \int_{Sp(\widehat{\Phi})} sP(s)$

where $Sp(\widehat{\Phi}) \subset \mathbb{R}$ is the spectrum of $\widehat{\Phi}$.

So there is a map $\chi: \sigma_{\mathbb{R}} \rightarrow H$ where $\sigma_{\mathbb{R}}$ is the Borel algebra of \mathbb{R} such that:

$\chi(\varpi)$ is a closed vector subspace of H

$$\chi(\mathbb{R}) = Id$$

$$\forall \varpi, \varpi' \in \sigma_{\mathbb{R}}, \varpi \cap \varpi' = \emptyset \Rightarrow \chi(\varpi) \cap \chi(\varpi') = \{0\}$$

$$\text{and } \widehat{\Phi} = \int_{Sp(\widehat{\Phi})} s\widehat{\pi}_{\chi(s)}$$

The spectrum $Sp(\widehat{\Phi})$ is a non empty compact subset of \mathbb{R} . If $\widehat{\Phi}$ is normal then $\lambda \in Sp(\widehat{\Phi}) \Leftrightarrow \bar{\lambda} \in Sp(\widehat{\Phi}^*)$.

For any fixed $\psi \neq 0 \in H$ the function $\widehat{\mu}_\psi: \sigma_{\mathbb{R}} \rightarrow \mathbb{R} :: \widehat{\mu}_\psi(\varpi) = \frac{\langle \widehat{\pi}_{\chi(\varpi)}\psi, \psi \rangle}{\|\psi\|^2} = \frac{\|\widehat{\pi}_{\chi(\varpi)}\psi\|^2}{\|\psi\|^2}$ is a probability law on $(\mathbb{R}, \sigma_{\mathbb{R}})$.

3. We will define:

Definition 12 A *secondary observable* is a linear map $\Phi \in L(V; V)$ valued in a finite dimensional vector subspace of V , such that $\widehat{\Phi} = \Upsilon \circ \Phi \circ \Upsilon^{-1}$ is a normal operator: $\widehat{\Phi} \circ \widehat{\Phi}^* = \widehat{\Phi}^* \circ \widehat{\Phi}$ with the adjoint $\widehat{\Phi}^*$

Theorem 13 Any secondary observable Φ is a compact, continuous map, its associated map $\widehat{\Phi} = \Upsilon \circ \Phi \circ \Upsilon^{-1}$ is a compact, self-adjoint, Hilbert-Schmidt and trace class operator.

$\hat{\Phi} = \sum_{p=1}^n \lambda_p Y_{J_p}$ where $(Y_{J_p})_{p=1}^n$ are primary observables associated to a basis $(e_i)_{i \in I}$ of V and $(J_p)_{p=1}^n$ are disjoint finite subsets of I

Proof. i) $\hat{\Phi}(H)$ is a finite dimensional vector subspace of H . So :

$\hat{\Phi}$ has 0 for eigen value, with an infinite dimensional eigen space H_c .

$\hat{\Phi}, \hat{\Phi}$ are compact and thus continuous (Maths.912).

ii) As $\hat{\Phi}$ is continuous and normal, there is a unique spectral measure P on $(\mathbb{R}, \sigma_{\mathbb{R}})$ such that $\hat{\Phi} = \int_{Sp(\hat{\Phi})} sP(s)$ where $Sp(\hat{\Phi}) \subset \mathbb{R}$ is the spectrum of $\hat{\Phi}$. As $\hat{\Phi}$ is compact, by the Riesz theorem (Maths.1142) its spectrum is either finite or is a countable sequence converging to 0 (which may or not be an eigen value) and, except possibly for 0, is identical to the set $(\lambda_p)_{p \in \mathbb{N}}$ of its eigen values (Maths.1020). For each distinct eigen value the eigen spaces H_p are orthogonal and H is the direct sum $H = \bigoplus_{p \in \mathbb{N}} H_p$. For each non null eigen value λ_p the eigen space H_p is finite dimensional.

Let λ_0 be the eigen value 0 of $\hat{\Phi}$. So : $\hat{\Phi} = \sum_{p \in \mathbb{N}} \lambda_p \hat{\pi}_{H_p}$ and any vector of H reads : $\psi = \sum_{p \in \mathbb{N}} \psi_p$ with $\psi_p = \hat{\pi}_{H_p}(\psi)$

Because $\hat{\Phi}(H)$ is finite dimensional, the spectrum is finite and the non null eigen values are $(\lambda_p)_{p=1}^n$, the eigen space corresponding to 0 is $H_c = (\bigoplus_{p=1}^n H_p)^\perp$

$\forall \psi \in H : \psi = \psi_c + \sum_{p=1}^n \psi_p$ with $\psi_p = \hat{\pi}_{H_p}(\psi), \psi_c = \hat{\pi}_{H_c}(\psi)$

$$\hat{\Phi} = \sum_{p=1}^n \lambda_p \hat{\pi}_{H_p}$$

Its adjoint reads : $\hat{\Phi}^* = \sum_{p \in \mathbb{N}} \bar{\lambda}_p \hat{\pi}_{H_p} = \sum_{p \in \mathbb{N}} \lambda_p \hat{\pi}_{H_p}$ because H is a real Hilbert space

$\hat{\Phi}$ is then self-adjoint, Hilbert-Schmidt and trace class, as the sum of the trace class operators $\hat{\pi}_{H_p}$.

iii) The observable reads :

$\Phi = \sum_{p=1}^n \lambda_p \pi_p$ where $\pi_p = \Upsilon^{-1} \circ \hat{\pi}_{H_p} \circ \Upsilon$ is the projection on a finite dimensional vector subspace of V :

$$\pi_p \circ \pi_q = \Upsilon^{-1} \circ \hat{\pi}_{H_p} \circ \Upsilon \circ \Upsilon^{-1} \circ \hat{\pi}_{H_q} \circ \Upsilon = \Upsilon^{-1} \circ \hat{\pi}_{H_p} \circ \hat{\pi}_{H_q} \circ \Upsilon = \delta_{pq} \Upsilon^{-1} \circ \hat{\pi}_{H_p} \circ \Upsilon = \delta_{pq} \pi_p$$

$\Phi \circ \pi_p = \lambda_p \pi_p$ so $\pi_p(V) = V_p$ is the eigen space of Φ for the eigen value λ_p and the subspaces $(V_p)_{p=1}^n$ are linearly independent.

By choosing any basis $(e_i)_{i \in J_p}$ of V_p , and $(e_i)_{i \in J^c}$ with $J^c = \mathcal{C}_I(\bigoplus_{p=1}^n J_p)$ for the basis of $V_c = Span((e_i)_{i \in J^c})$

$$X = Y_{J^c}(X) + \sum_{p=1}^n Y_{J_p}(X)$$

the observable Φ reads : $\Phi = \sum_{p=1}^n \lambda_p Y_{J_p}$ ■

We have :

$$\begin{aligned} Y_{J_p}(X) &= \sum_{i \in J_p} \left\langle \phi_i, \hat{Y}_{J_p}(\Upsilon(X)) \right\rangle_H e_i \\ \Phi(X) &= \sum_{p=1}^n \lambda_p \sum_{i \in J_p} \left\langle \phi_i, \hat{Y}_{J_p}(\Upsilon(X)) \right\rangle_H e_i \\ &= \sum_{i \in I} \left\langle \phi_i, \sum_{p=1}^n \lambda_p \hat{Y}_{J_p}(\Upsilon(X)) \right\rangle_H e_i \\ &= \sum_{i \in I} \left\langle \phi_i, \hat{\Phi}(\Upsilon(X)) \right\rangle_H e_i \end{aligned}$$

$\Phi, \widehat{\Phi}$ have invariant vector spaces, which correspond to the direct sum of the eigen spaces.

The probability law $\widehat{\mu}_\psi : \sigma_{\mathbb{R}} \rightarrow \mathbb{R}$ reads :

$$\widehat{\mu}_\psi(\varpi) = \Pr(\lambda_p \in \varpi) = \frac{\|\widehat{\pi}_{H_p}(\psi)\|^2}{\|\psi\|^2}$$

To sum up :

Theorem 14 *For any primary or secondary observable Φ , there is a basis $(e_i)_{i \in I}$ of V , a compact, self-adjoint, Hilbert-Schmidt and trace class operator $\widehat{\Phi}$ on the associated Hilbert space H such that :*

$$\widehat{\Phi} = \Upsilon \circ \Phi \circ \Upsilon^{-1}$$

if the system is in the state $X = \sum_{i \in I} \langle \phi_i, \Upsilon(X) \rangle_H e_i$ the value of the observable is : $\Phi(X) = \sum_{i \in I} \langle \phi_i, \widehat{\Phi}(\Upsilon(X)) \rangle_H e_i$

$\widehat{\Phi}$ has a finite set of eigen values, whose eigen spaces (except possibly for 0) are finite dimensional and orthogonal. The vectors corresponding to the eigen value 0 are never observed, so it is convenient to represent the Hilbert space H through a basis of eigen vectors, each of them corresponding to a definite state, which usually can be identified. This is a method commonly used in Quantum Mechanics, however the vector has also a component in the eigen space corresponding to the null eigen value, which is not observed but exists. Conversely any observable (on V) can be defined through an operator on H with the required properties (compact, normal, it is then self-adjoint). We will come back on this point in the following, when a group is involved.

2.2.4 Efficiency of an observable

A crucial factor for the quality and the cost of the estimation procedure is the number of parameters to be estimated, which is closely related to the dimension of the vector space $\Phi(V)$, which is finite. The error made by the choice of $\Phi(X)$ when the system is in the state X is : $o_\Phi(X) = X - \Phi(X)$. If two observables Φ, Φ' are such that $\Phi(V), \Phi'(V)$ have the same dimension, one can say that Φ is more efficient than Φ' if : $\forall X : \|o_\Phi(X)\|_V \leq \|o_{\Phi'}(X)\|_V$

To assess the efficiency of a secondary observable Φ it is legitimate to compare Φ to the primary observable Y_J with a set J which has the same cardinality as the dimension of $\oplus_{p=1}^n H_p$.

The error with the choice of Φ is :

$$o_\Phi(X) = X - \Phi(X) = Y_c(\psi) + \sum_{p=1}^n (1 - \lambda_p) Y_p(\psi)$$

$$\|o_\Phi(X)\|_V^2 = \|Y_c(\psi)\|_V^2 + \sum_{p=1}^n (1 - \lambda_p)^2 \|Y_p(\psi)\|^2$$

$$\widehat{o}_\Phi(\Upsilon(X)) = \Upsilon(X) - \widehat{\Phi}(\Upsilon(X)) = \widehat{\pi}_{H_c}(\psi) + \sum_{p=1}^n (1 - \lambda_p) \widehat{\pi}_{H_p}(\psi)$$

$$\|\widehat{o}_\Phi(\Upsilon(X))\|^2 = \|\widehat{\pi}_{H_c}(\psi)\|^2 + \sum_{p=1}^n (1 - \lambda_p)^2 \|\widehat{\pi}_{H_p}(\psi)\|^2 = \|o_\Phi(X)\|_V^2$$

$$\text{And for } Y_J : \|\widehat{o}_{Y_J}(\Upsilon(X))\|^2 = \|\widehat{\pi}_{H_c}(\psi)\|^2 \text{ because } \lambda_p = 1$$

So :

Theorem 15 *For any secondary observable there is always a primary observable which is at least as efficient.*

This result justifies the restriction, in the usual formalism, of observables to operators belonging to a von Neumann algebra.

2.2.5 Statistical estimation and primary observables

At first the definition of a primary observable seems naive, and the previous results will seem obvious to the specialists of Axiomatic QM. After all the definition of a primary observable requires only the choice of a finite number of independent vectors of V . We have already seen that a primary observable is always better than a, more sophisticated, secondary observable. But we have also to compare a primary observable to what is practically done in an experiment, where we have to estimate a map from a batch of data.

Consider a model with variables X , maps, belonging to a Hilbert space H (to keep it simple), from a set M to a normed vector space E , endowed with a scalar product $\langle \rangle_E$. The physicist has a batch of data, that is a finite set $\{x_p \in E, p = 1..N\}$ of N measures of X done at different points $\Omega = \{m_p \in M, p = 1..N\}$: of $M : x_p = X(m_p)$. The estimated map \hat{X} should be a solution of the collection of equations : $x_p = X(m_p)$ where x_p, m_p are known.

The **evaluation maps**, that we will encounter several times, is the collection of maps $\mathcal{E}(m)$ on H :

$$\mathcal{E}(m) : H \rightarrow E :: \mathcal{E}(m)Y = Y(m)$$

Because H and E are vector spaces $\mathcal{E}(m)$ is a linear map : $\mathcal{E}(m) \in L(H; E)$, depending on both H and E . It can be continuous or not.

The set of solutions of the equations, that is of maps Y of H such that $\forall m_p \in \Omega : Y(m_p) = x_p$ is :

$$A = \bigcap_{m_p \in \Omega} \mathcal{E}(m_p)^{-1}(x_p)$$

$$Y \in A \Leftrightarrow \forall m \in \Omega : Y(m) = X(m)$$

It is not empty because it contains at least X . Its closed convex hull is the set B in H (Maths.361) :

$$\forall Z \in B : \exists \alpha \in [0.1], Y, Y' \in A : Z = \alpha Y + (1 - \alpha) Y'$$

$$\Rightarrow \forall m \in \Omega : Z(m) = x_p$$

B is the smallest closed set of H such that all its elements Z are solutions of the equations : $\forall p = 1..N : Z(m_p) = x_p$.

If we specify an observable, we restrict X to a finite dimensional subspace $H_J \subset H$. With the evaluation map \mathcal{E}_J on H_J we can consider the same procedure, but then usually $A_J = \emptyset$. The simplification of the map to be estimated as for consequence that there is no solution to the equations. So the physicist uses a statistical method, that is a map which associates to each batch of data $X(\Omega)$ a map $\varphi(X(\Omega)) = \hat{X} \in H_J$. Usually \hat{X} is such that it minimizes the sum

of the distance between points in E : $\sum_{m \in \Omega} \left\| \widehat{X}(m) - x_p \right\|_E$ (other additional conditions can be required).

The primary observable Φ gives another solution : $\Phi(X)$ is the orthogonal projection of X on the Hilbert space H_J , it is such that it minimizes the distance between maps :

$$\forall Z \in H_J : \|X - Z\|_H \geq \|X - \Phi(X)\|_H.$$

$\Phi(X)$ always exist, and does not depend on the choice of an estimation procedure φ . $\Phi(X)$ minimizes the distance between maps in H , meanwhile $\varphi(X(\Omega))$ minimizes distance between points in E . Usually $\varphi(X(\Omega))$ is different from $\Phi(X)$ and $\Phi(X)$ is a better estimate than \widehat{X} : *a primary observable is actually the best statistical estimator* for a given size of the sample. But it requires the explicit knowledge of the scalar product and H_J . This can be practically done in some significant cases (see for an example J.C.Dutailly *Estimation of the probability of transitions between phases*).

Knowing the estimate \widehat{X} provided by a statistical method φ , we can implement the previous procedure to the set $\widehat{X}(\Omega)$ and compute the set of solutions : $\widehat{A} = \bigcap_{m_p \in \Omega} \mathcal{E}_J(m_p)^{-1}(\widehat{X}(m))$. It is not empty. Its closed convex hull \widehat{B} in H_J can be considered as the domain of confidence of \widehat{X} : they are maps which take the same values as \widehat{X} in Ω and as a consequence give the same value to $\sum_{m \in \Omega} \left\| \widehat{X}(m) - x_p \right\|_E$.

Because \widehat{B} is closed and convex there is a unique orthogonal projection Y of X on \widehat{B} (Maths.1107) and :

$$\forall Z \in \widehat{B} : \|X - Z\|_H \geq \|X - Y\|_H \Rightarrow \left\| X - \widehat{X} \right\|_H \geq \|X - Y\|_H$$

so Y is a better estimate than $\varphi(X(\Omega))$, and can be computed if we know the scalar product on H .

We see clearly the crucial role played by the choice of a specification. But it leads to a more surprising result, of deep physical meaning.

2.2.6 Quantization of singularities

A classic problem in Physics is to prove the existence of a singular phenomenon, appearing only for some values of the parameters m . To study this problem we use a model similar to the previous one, with the same notations. But here the variable X is comprised of two maps, X_1, X_2 with unknown, disconnected, domains M_1, M_2 : $M = M_1 + M_2$. The first problem is to estimate X_1, X_2 .

With a statistical process $\varphi(X(\Omega))$ it is always possible to find estimations $\widehat{X}_1, \widehat{X}_2$ of X_1, X_2 . The key point is to distinguish in the set Ω the points which belong to M_1 and M_2 . There are $\frac{1}{2}(2^N - 2) = 2^{N-1} - 1$ distinct partitions of Ω in two subsets $\Omega_1 + \Omega_2$, on each subset the statistical method φ gives the estimates :

$$\widehat{Y}_1 = \varphi(X(\Omega_1)), \widehat{Y}_2 = \varphi(X(\Omega_2))$$

Denote : $\rho(\Omega_1, \Omega_2)$
 $= \sum_{m_p \in \Omega_1} \|X(m_p) - \varphi(X(\Omega_1))(m_p)\| + \sum_{m_p \in \Omega_2} \|X(m_p) - \varphi(X(\Omega_2))(m_p)\|$

A partition (Ω_1, Ω_2) is said to be a better fit than (Ω'_1, Ω'_2) if :

$$\rho(\Omega_1, \Omega_2) \leq \rho(\Omega'_1, \Omega'_2)$$

Then $\widehat{X}_1 = \varphi(X(\Omega_1))$, $\widehat{X}_2 = \varphi(X(\Omega_2))$ is the solution for the best partition.

So there is a procedure, which provides always the best solution given the data and φ , but it does not give M_1, M_2 precisely, their estimation depends on the structure of M.

However it is a bit frustrating, if we want to test a law, because the procedure provides always a solution, even if actually there is no such partition of X. And this can happen. If we define the sets as above with the evaluation map : $\mathcal{E}_J(m) : H_J \rightarrow E :: \mathcal{E}(m)Y = Y(m)$

$$A_k = \cap_{m_p \in \Omega_k} \mathcal{E}(m_p)^{-1}(\widehat{X}_k(m_p)) \subset H_J \text{ for } k = 1, 2. \text{ It is not empty}$$

because it contains at least \widehat{X}_k .

B_k the closed convex hull of A_k in H_J

$$\text{Then : } \forall Y \in B_k, m \in \Omega_k : Y(m) = \widehat{X}_k(m)$$

If $B_1 \cap B_2 \neq \emptyset$ there is at least one map, which can be defined uniquely on M, belongs to H_J and is equivalent to $\widehat{X}_1, \widehat{X}_2$.

This issue is of importance because many experiments aim at proving the existence of a special behavior. We need, in addition, a test of the hypothesis (denoted H_0) : there is a partition (and then the best solution would be $\widehat{X}_1, \widehat{X}_2$) against the hypothesis (denoted H_1) there is no partition : there is a unique map $\widehat{X} \in H_J$ for the domain Ω . The simplest test is to compare $\sum_{m_p \in \Omega} \|X(m_p) - \varphi(\Omega)(m_p)\|$ to $\rho(\Omega_1, \Omega_2)$. If $\varphi(\Omega)$ gives results as good as $\widehat{X}_1, \widehat{X}_2$ we can reject the hypothesis. Notice that it accounts for the properties assumed for the maps in H_J . For instance if H_J is comprised uniquely of continuous maps, then $\varphi(X(\Omega))$ is continuous, and clearly distinct from the maps $\widehat{X}_1, \widehat{X}_2$ continuous only on M_1, M_2 .

It is quite obvious that the efficiency of this test decreases with N : the smaller N, the greater the chance to accept H_0 . Is there a way to control the validity of an experiment ? The Theory of Tests, a branch of Statistics, studies this kind of problems.

The problem is, given a sample of points $\Omega = (m_p)_{p=1}^N$ and the corresponding values $x = (x_p)_{p=1}^N$, decide if they obey to a simple (X , Hypothesis H_1) or a double (X_1, X_2 , Hypothesis H_0) distribution law.

The choice of the points $(m_p)_{p=1}^N$ in a sample is assumed to be random : all the points m of M have the same probability to be in Ω , but the size of M_1, M_2 can be different, so it could give a different chance for a point of M_1 or M_2 to be in the sample. Let us say that :

$$\Pr(m \in M_1 | H_0) = 1 - \lambda, \Pr(m \in M_2 | H_0) = \lambda, \Pr(m \in M | H_1) = 1$$

(all the probabilities are for a sample of a given size N)

Then the probability for any vector of E to have a given value x depends only on the map X : this is the number of points m of M for which $X(m) = x$. For instance if there are two points m with $X(m) = x$ then x has two times the probability to appear, and if X is more concentrated in an area of E , this area has more probability to appear. Let us denote this value $\rho(x) \in [0, 1]$.

Rigorously (Maths.869), with a measure dx on E , μ on M , $\rho(x)dx$ is the pull-back of the measure μ on M . For any ϖ belonging to the Borel algebra σE of E :

$$\int_{\varpi} \rho(x) dx = \int_{\mathcal{E}(m)^{-1}(\varpi)} \mu(m) \Leftrightarrow \rho(x) dx = X^* \mu$$

If H_1 is true, the probability $\Pr(x|H_1) = \rho(x)$ depends only on the value x , that is of the map X .

If H_0 is true the probability depends on the maps and if $m \in M_1$ or $m \in M_2$ ($M = M_1 + M_2$)

$$\Pr(x|H_0 \wedge m \in M_1) = \rho_1(x)$$

$$\Pr(x|H_0 \wedge m \in M_2) = \rho_2(x)$$

$$\Rightarrow \Pr(x|H_0) = (1 - \lambda) \rho_1(x) + \lambda \rho_2(x)$$

Moreover we have with some measure dx on E :

$$\int_E \rho(x) dx = \int_E \rho_1(x) dx = \int_E \rho_2(x) dx = 1$$

The likelihood function is the probability of a given batch of data. It depends on the hypothesis :

$$L(x|H_0) = \Pr(x_1, x_2, \dots, x_N|H_0) = \prod_{p=1}^N ((1 - \lambda) \rho_1(x_p) + \lambda \rho_2(x_p))$$

$$L(x|H_1) = \Pr(x_1, x_2, \dots, x_N|H_1) = \prod_{p=1}^N \rho(x_p)$$

The Theory of Tests gives us some rules (see Kendall t.II). A critical region is an area $w \subset E^N$ such that H_0 is rejected if $x \in w$.

One considers two risks :

- the risk of type I is to wrongly reject H_0 . It has the probability : $\alpha = \Pr(x \in w|H_0)$

- the risk of type II is to wrongly accept H_0 . It has the probability : $1 - \beta = \Pr(x \in E^N - w|H_0)$ called the power of the test thus :

$$\beta = \Pr(x \in w|H_1)$$

A simple rule, proved by Neyman and Pearson, says that the best critical region w is defined by :

$$w = \left\{ x : \frac{L(x|H_0)}{L(x|H_1)} \leq k \right\}$$

the scalar k being defined by : $\alpha = \Pr(x \in w|H_0)$. So we are left with a single parameter α , which can be seen as the rigor of the test.

The critical area $w \subset E^N$ is then :

$$w = \left\{ x \in E^N : \prod_{p=1}^N \frac{((1-\lambda)\rho_1(x_p) + \lambda\rho_2(x_p))}{\rho(x_p)} \leq k \right\}$$

with :

$$\alpha = \int_w \prod_{p=1}^N ((1 - \lambda) \rho_1(\xi_p) + \lambda \rho_2(\xi_p)) (d\xi)^N$$

It provides a reliable method to build a test, but requires to know, or to estimate, $\rho, \rho_1, \rho_2, \lambda$.

In most of the cases encountered, actually one looks for an anomaly.

H_1 is unchanged, there is only one map X , defined over M . Then : $\Pr(x|H_1) = \rho(x)$

H_0 becomes :

$$M = M_1 + M_2$$

$$\Pr(m \in M_1|H_0) = 1 - \lambda, \Pr(m \in M_2|H_0) = \lambda$$

On M_1 the variable is X :

$$\Pr(x_p|H_0 \wedge m_p \in M_1) = \rho(x) \Rightarrow \Pr(x_p|H_0) = (1 - \lambda) \rho(x)$$

On M_2 the variable becomes X_2

$$\Pr(x_p|H_0 \wedge m_p \in M_2) = \rho_2(x) \Rightarrow \Pr(x_p|H_0) = \lambda \rho_2(x)$$

And w is :

$$w = \left\{ x \in E^N : \prod_{p=1}^N \frac{((1-\lambda)\rho(x_p) + \lambda\rho_2(x_p))}{\rho(x_p)} \leq k \right\}$$

$$w = \left\{ x \in E^N : \prod_{p=1}^N \left(1 - \lambda + \lambda \frac{\rho_2(x_p)}{\rho(x_p)} \right) \leq k \right\}$$

$$\alpha = \int_w \prod_{p=1}^N ((1 - \lambda) \rho(x_p) + \lambda \rho_2(x_p)) (dx)^N$$

$$\beta = \Pr(x \in w|H_1) = \int_w \left(\prod_{p=1}^N \rho(x_p) \right) (dx)^N$$

If there is one observed value such that $\rho(x_p) = 0$ then H_0 should be accepted. But, because ρ, ρ_2 are not well known, and the imprecision of the experiments, H_0 would be proven if $\frac{L(x|H_0)}{L(x|H_1)} > k$ for a great number of experiments. So we can say that H_0 is scientifically proven if :

$$\forall (x_1, x_2, \dots, x_N) : \prod_{p=1}^N \left((1 - \lambda) + \lambda \frac{\rho_2(x_p)}{\rho(x_p)} \right) > k$$

By taking $x_1 = x_2 = \dots = x_N = x$:

$$\forall x : (1 - \lambda) + \lambda \frac{\rho_2(x)}{\rho(x)} > k^{1/N}$$

$$\frac{\rho_2(x)}{\rho(x)} > (k^{1/N} + \lambda - 1) / \lambda$$

$$\text{When } N \rightarrow \infty : k^{1/N} \rightarrow 1 \Rightarrow \frac{\rho_2(x)}{\rho(x)} > 1$$

So a necessary condition to have a chance to say that a singularity has been reliably proven is that : $\forall x : \frac{\rho_2(x)}{\rho(x)} > 1$.

The function $\frac{\rho_2(x)}{\rho(x)}$ can be called the Signal to Noise Ratio, by similarity with the Signal Theory. Notice that we have used very few assumptions about the variables. And we can state :

Theorem 16 *In a system represented by variables X which are maps defined on a set M and valued in a vector space E , a necessary condition for a singularity*

to be detected is that the Signal to Noise Ratio is greater than 1 for all values of the variables in E .

This result can be seen in another way : if a signal is acknowledged, then necessarily it is such that $\frac{\rho_2(x)}{\rho(x)} > 1$. Any other signal would be interpreted as related to the imprecision of the measure. So there is a threshold under which phenomena are not acknowledged, and their value is necessarily above this threshold. The singular phenomena are quantized. We will see an application of this result about the Planck's law.

2.3 PROBABILITY

One of the main purposes of the model is to know the state X , represented by some vector $\psi \in H$. The model is fully determinist, in that the values of the variables X are not assumed to depend on a specific event : there is no probability law involved in its definition. However the value of X which will be acknowledged at the end of the experiment, when all the data have been collected and analyzed, differs from its actual value. The discrepancy stems from the usual imprecision of any measure, but also more fundamentally from the fact that we estimate a vector in an infinite dimensional vector space from a batch of data, which is necessarily finite. We will focus on this later aspect, that is on the discrepancy between an observable $\Phi(X)$ and X .

In any practical physical experiment the estimation of X requires the choice of an observable. We have seen that the most efficient solution is to choose a primary observable which, furthermore, provides the best statistical estimator. However usually neither the map Φ nor the basis $(e_i)_{i \in I}$ are explicit, even if they do exist. So we can look at the discrepancy $X - \Phi(X)$ from a different point of view : for a given, fixed, value of the state X , what is the uncertainty which stems from the choice of Φ among a large class of observables ? This sums up to assess the risk linked to the choice of a specification for the estimation of X .

2.3.1 Primary observables

Let us start with primary observables : the observable Φ is some projection on a finite dimensional vector subspace of V .

The bases of the vector space V_0 (such that $O \subset V_0$) have the same cardinality, so we can consider that the set I does not depend on a choice of a basis (actually one can take $I = \mathbb{N}$). The set 2^I is the largest σ -algebra on I . The set $(I, 2^I)$ is measurable (Maths.802).

For any fixed $\psi \neq 0 \in H$ the function

$$\hat{\mu}_\psi : 2^I \rightarrow \mathbb{R} :: \hat{\mu}_\psi(J) = \frac{\langle \hat{Y}_J \psi, \psi \rangle}{\|\psi\|^2} = \frac{\|\hat{Y}_J \psi\|^2}{\|\psi\|^2}$$

is a probability law on $(I, 2^I)$: it is positive, countably additive and $\hat{\mu}_\psi(I) = 1$ (Maths.11.4.1).

If we see the choice of a finite subset $J \in 2^I$ as an event in a probabilist point of view, for a given $\psi \neq 0 \in H$ the quantity $\hat{Y}_J(\psi)$ is a random variable, with a distribution law $\hat{\mu}_\psi$

The operator \hat{Y}_J has two eigen values : 1 with eigen space $\hat{Y}_J(H)$ and 0 with eigen space $\hat{Y}_{J^c}(H)$. Whatever the primary observable, the value of $\Phi(X)$ will be $Y_J(X)$ for some J , that is an eigen vector of the operator $\Phi = Y_J$, and the probability to observe $\Phi(X)$, if the system is in the state X , is :

$$\Pr(\Phi(X) = Y_J(X)) = \Pr(J|\psi) = \hat{\mu}_\psi(J) = \frac{\|\hat{Y}_J \psi\|^2}{\|\psi\|^2} = \frac{\|\hat{\Phi}(Y(X))\|_H^2}{\|Y(X)\|_H^2}$$

This result still holds if another basis had been chosen : $\Phi(X)$ will be $Y_J(X)$ for some J , expressed in the new basis, but with a set J of same cardinality. And some specification must always be chosen. So we have :

Theorem 17 For any primary observable Φ , the value $\Phi(X)$ which is measured is an eigen vector of the operator Φ , and the probability to measure a value $\Phi(X)$ if the system is in the state X is :

$$\Pr(\Phi(X)|X) = \frac{\|\widehat{\Phi}(\Upsilon(X))\|_H^2}{\|\Upsilon(X)\|_H^2}$$

2.3.2 Secondary observables

For a secondary observable, as defined previously :

$$\Phi = \sum_{p=1}^n \lambda_p Y_{J_p}$$

$$\widehat{\Phi} = \sum_{p=1}^n \lambda_p \widehat{\pi}_{H_p}$$

The vectors decompose as :

$$X = Y_{J^c}(X) + \sum_{p=1}^n X_p$$

$$\text{with } X_p = Y_{J_p}(X) = \sum_{i \in J_p} \langle \phi_i, \widehat{Y}_{J_p}(\Upsilon(X)) \rangle_H e_i \in V_p$$

$$\Upsilon(X) = \psi = \psi_c + \sum_{p=1}^n \psi_p \text{ with } \psi_p = \widehat{\pi}_{H_p}(\psi), \psi_c = \widehat{\pi}_{H^c}(\psi)$$

where ψ_p is an eigen vector of $\widehat{\Phi}$, X_p is an eigen vector of Φ both for the eigen value λ_p

and

$$\Phi(X) = \sum_{p=1}^n \lambda_p X_p$$

$$\widehat{\Phi}(\psi) = \sum_{p=1}^n \lambda_p \psi_p$$

If, as above, we see the choice of a finite subset $J \in 2^I$ as an event in a probabilist point of view then the probability that $\Phi(X) = \lambda_p X_p$ if the system is in the state X , is given by $\Pr(J_p|X) = \frac{\|\widehat{Y}_p \psi\|^2}{\|\psi\|^2} = \frac{\|\psi_p\|^2}{\|\psi\|^2}$

And we have :

Theorem 18 For any secondary observable Φ , the value $\Phi(X)$ which is observed if the system is in the state X is a linear combination of eigen vectors X_p of Φ for the eigen value λ_p : $\Phi(X) = \sum_{p=1}^n \lambda_p X_p$

The probability that $\Phi(X) = \lambda_p X_p$ is:

$$\Pr(\Phi(X) = \lambda_p X_p | X) = \frac{\|\Upsilon(X_p)\|_H^2}{\|\Upsilon(X)\|_H^2}$$

Which can also be stated as : $\Phi(X)$ can take the values $\lambda_p X_p$, each with the probability $\frac{\|\psi_p\|^2}{\|\psi\|^2}$, then $\Phi(X)$ reads as an expected value. This is the usual way it is expressed in QM.

The interest of these results comes from the fact that we do not need to explicit any basis, or even the set I . And we do not involve any specific property of the estimator of X , other than Φ is an observable. The operator $\widehat{\Phi}$ sums up the probability law.

Of course this result can be seen in another way : as only $\Phi(X)$ can be accessed, one can say that the system takes only the states $\Phi(\lambda_p X_p)$, with a probability $\frac{\|\psi_p\|^2}{\|\psi\|^2}$. This gives a probabilistic behavior to the system (X becoming

a random variable) which is not present in its definition, but is closer to the usual interpretation of QM.

This result can be illustrated by a simple example. Let us take a single continuous variable x , which takes its values in \mathbb{R} . It is clear that any physical measure will at best give a rational number $Y(x) \in \mathbb{Q}$ up to some scale. There are only countably many rational numbers for unaccountably many real scalars. So the probability to get $Y(x) \in \mathbb{Q}$ should be zero. The simple fact of the measure gives the paradox that rational numbers have an incommensurable weight, implying that each of them has some small, but non null, probability to appear. In this case I can be assimilated to \mathbb{Q} , the subsets J are any finite collection of rational numbers.

2.3.3 Wave function

The wave function is a central object in QM, but it has no general definition and is deemed non physical (except in the Bohm's interpretation). Usually this is a complex valued function, defined over the space of configuration of the system : the set of all possible values of the variables representing the system. If it is square integrable, then it belongs to a Hilbert space, and can be assimilated to the vector representing the state. Because its arguments comprise the coordinates of objects such as particles, it has a value at each point, and the square of the module of the function is proportional to the probability that the measure of the variable takes the values of the arguments at this point. Its meaning is relatively clear for systems comprised of particles, but less so for systems which include force fields, because the space of configuration is not defined. But we will see now how it can be precisely defined in our framework.

Theorem 19 *In a system modelled by N variables, collectively denoted X , which are maps $X : M \rightarrow F$ from a common measured set M to a finite dimensional normed vector space F and belonging to an open subset of an infinite dimensional, separable, real Fréchet vector space V , such that the evaluation map $\mathcal{E}(m) : V \rightarrow F :: \mathcal{E}(m)(X) = X(m)$ which assigns at any X its value in a fixed point m in M is measurable : then for any state X of the system there is a function $W : M \times F \rightarrow \mathbb{R}$ such that $W(m, y) = \Pr(\Phi(X)(m) = y|X)$ is the probability that the measure of the value of any primary observable $\Phi(X)$ at m is y .*

Proof. The conditions 1 apply, there is a Hilbert space H and an isometry $\Upsilon : V \rightarrow H$.

To the primary observable $\Phi : V \rightarrow V_J$ is associated the self-adjoint operator $\widehat{\Phi} = \Upsilon \circ \Phi \circ \Upsilon^{-1}$

We can apply the theorem 17: the probability to measure a value $\Phi(X) = Y$ if the system is in the state X is :

$$\Pr(\Phi(X) = Y|X) = \frac{\|\widehat{\Phi}(\Upsilon(X))\|_H^2}{\|\Upsilon(X)\|_H^2} = \pi(Y)$$

Because only the maps belonging to V_J are observed it provides a probability law π on the set $V_J : \pi : V_\sigma \rightarrow [0, 1]$ where V_σ is the Borel algebra of V_J .

The evaluation map : $\mathcal{E}_J(m) : V_J \rightarrow F :: \mathcal{E}_J(m)(Y) = Y(m)$ assigns at any $Y \in V_J$ its value in the fixed point m in M .

If $y \in F$ is a given vector of F , the set of maps in V_J which gives the value y in m is : $\varpi(m, y) = \mathcal{E}_J(m)^{-1}(y) \subset V_J$.

The probability that the observable takes the value y at m $\Phi(X)(m) = y$ is

$$\begin{aligned} \pi(\varpi(m, y)) &= \pi\left(\mathcal{E}_J(m)^{-1}(x)\right) \\ &= \frac{1}{\|\Upsilon(X)\|_H^2} \int_{Y \in \varpi(m, y)} \left\| \widehat{\Phi}(\Upsilon(Y)) \right\|_H^2 \pi(Y) = W(m, y) \quad \blacksquare \end{aligned}$$

If M is endowed with a positive measure μ and X is a scalar function, the space V of square integrable maps $\int_\Omega |X(m)|^2 \mu(m) < \infty$ is a separable Hilbert space H , then the conditions 1 are met and H can be identified with the space of the states.

$$W(m, y) = \frac{1}{\|X\|_H^2} \int_{Y \in \varpi(m, y)} |Y|_H^2 = \left(\int_\Omega |X|^2 \mu \right)^{-1} \mu(Y^{-1}(m, y))$$

No structure, other than the existence of the measure μ , is required on M . But of course if the variables X include derivatives M must be at least a differentiable manifold.

W can be identified with the square of the wave function of QM.

2.4 CHANGE OF VARIABLES

In the conditions 1 we have noticed that, in the model, the variables could be defined over different connected domains. Actually one can go further and consider the change of variables, which leads to a theorem similar to the well known Wigner's theorem. The problem appears in Physics in two different ways, which reflect the interpretations of Scientific laws. As we will see in the following chapters these considerations are fundamental in Physics, notably they are the basis for the theories of gauge.

2.4.1 Two ways to define the same state of a system

The first way : from a theoretical model

In the first way the scientist has built a theoretical model, using known concepts and their usual representation by mathematical objects. A change of variables appears notably when :

i) The variables are the components of a geometric quantity (a vector, a tensor,...) expressed in some basis. According to the general Principle of Relativity, the state of the system shall not depend on the observers (those measuring the coordinates). For instance it should not matter if the state of a system is measured in different units. The data change, but according to rules which depend on the mathematical representation which is used, and not on the system itself. In a change of basis coordinates change but they represent the same vectorial quantity. We will see another example with interacting, indistinguishable systems.

ii) The variables are maps, depending on arguments which are themselves coordinates of some event : $X_k = X_k(\xi_1, \dots, \xi_{p_k})$. Similarly these coordinates ξ can change according to some rules, while the variable X_k represents the same event. A simple example that we will develop later on is a simple function of the time $X_k(t)$ such that the time t can be expressed in different units, or with different origin : $X_k(t)$ and $X'_k(t) = X_k(t + \theta)$ represent the same state.

By definition in both cases there is a continuous bijective map $U : V \rightarrow V'$ such that X and $X' = U(X)$ represent the same state of the system. This is the way mathematicians see a change of variables, and is usually called the passive way by physicists.

Any primary or secondary observable Φ is a linear map $\Phi \in L(V; W)$ into a finite dimensional vector subspace W . For the new variable the observable is $\Phi' \in L(V; W')$. Both $W, W' \subset V$ but W' is not necessarily identical to W . However the assumption that $X' = U(X)$ and X represents the same state of the system implies that for any measure of the state we have a similar relation : $\Phi' \circ U(X) = U \circ \Phi(X) \Leftrightarrow \Phi' \circ U = U \circ \Phi$. This is actually the true meaning of "represent the same state". This means that actually one makes the measures according to a fixed procedure, given by Φ , on variables which vary with U . Because U is a bijection on V : $\Phi' = U \circ \Phi \circ U^{-1}$.

The second way : from experimental measures

In the second way the scientist makes measures with a device that can be adjusted according to different values of a parameter, say θ : the simplest example is using different units, but often it is the orientation of the device which can be changed. And the measures $Y(\theta)$ which are taken are related to the choice of parameter for the device. If the results of experiments show that $Y(\theta) = Q(\theta)Y(\theta_0)$ with a bijective map $Q(\theta)$ and θ_0 some fixed value of the parameter one can assume that this experimental relation is a feature of the system itself.

Physicists distinguish a passive transformation, when only the device changes, and an active transformation, when actually the experiment involves a physical change on the system. In a passive transformation we come back to the first way and it is legitimate to assume that we have actually the same state, represented by different data, reflecting some mathematical change in their expression, even if the observable, which is valued in a finite dimensional space, does not account for all the possible values of the variables. In an active transformation (for instance in the Stern-Gerlach experiment one changes the orientation of a magnetic field to which the particles are submitted) one can say that there is some map U acting on the space V of the states of the system, such that the measure is done by a unique procedure $\tilde{\Phi}$ on a state X which is changed by a map $U(\theta)$. So that the measures are $Y(\theta) = \tilde{\Phi} \circ U(\theta) X$ and the relation $Y(\theta) = Q(\theta)Y(\theta_0)$ reads : $\tilde{\Phi} \circ U(\theta)(X) = U(\theta) \circ \tilde{\Phi}(X)$. So this is very similar to the first case, where θ represents the choice of a frame.

In both cases there is the general idea that the state of the system is represented by some fixed quantity, which can be measured in different procedures, so that there is a relation, given by the way one goes from one procedure to the others, between the measures of the state. In the first way the conclusion comes from the mathematical definition in a theoretical model : this is a simple mathematical deduction using the Principle of Relativity. In the second way there is an assumption : that one can extend the experimental facts, necessarily limited to a finite number of data, to the whole set of possible values of the variable.

The Theorem 2 is based on the existence of a Fréchet manifold structure on the set of possible values of the maps X . The same manifold structure can be defined by different, compatible, atlas. So the choice of other variables can lead to the same manifold, and the fixed quantity that we identify with a state is just a point on the manifold, and the change of variables is a change of charts between compatible atlas. The variables must be related by transition maps, that is continuous bijections, but additional conditions are required, depending on the manifold structure considered. For instance for differentiable manifolds the transition maps must be differentiable. We will request that the transition maps preserve the positive kernel, which plays a crucial role in Fréchet manifolds.

2.4.2 Fundamental theorem for a change of variables

We will summarize these features in the following :

Condition 20

i) The same system is represented by the variables $X = (X_1, \dots, X_N)$ and $X' = (X'_1, \dots, X'_{N'})$ which belong to open subsets O, O' of the infinite dimensional, separable, Fréchet vector space V .

ii) There is a continuous map $U : V \rightarrow V$, bijective on (O, O') , such that X and $X' = U(X)$ represent the same state of the system

iii) U preserves the positive kernel on V^1

iv) For any observable Φ of X , and Φ' of $X' : \Phi' \circ U = U \circ \Phi$

The map U shall be considered as part of the model, as it is directly related to the definition of the variables, and is assumed to be known. There is no hypothesis that it is linear.

Theorem 21 Whenever a change of variables on a system meets the conditions 20 above,

i) there is a unitary, linear, bijective map $\widehat{U} \in \mathcal{L}(H; H)$ such that : $\forall X \in O : \widehat{U}(\Upsilon(X)) = \Upsilon(U(X))$ where H is the Hilbert space and Υ is the linear map : $\Upsilon : V \rightarrow H$ associated to X, X'

ii) U is necessarily a bijective linear map.

For any observables Φ, Φ' :

iii) $W' = \Phi'(V)$ is a finite dimensional vector subspace of V , isomorphic to $W = \Phi(V) : W' = U(W)$

iv) the associated operators $\widehat{\Phi} = \Upsilon \circ \Phi \circ \Upsilon^{-1}, \widehat{\Phi}' = \Upsilon \circ \Phi' \circ \Upsilon^{-1}$ are such that : $\widehat{\Phi}' = \widehat{U} \circ \widehat{\Phi} \circ \widehat{U}^{-1}$ and $H'_{\Phi'} = \widehat{\Phi}'(H)$ is a vector subspace of H isomorphic to $H_{\Phi} = \widehat{\Phi}(H)$

Proof. i) Let $V_0 = O \cup O'$. This is an open set and we can apply the theorem 2. There is a homeomorphism $\Xi : V_0 \rightarrow H_0$ where H_0 is an open subset of a Hilbert space H . For a basis $(e_i)_{i \in I}$ of $\text{Span}(V_0)$ there is an isometry Υ such that :

$$\Upsilon : V_0 \rightarrow H_0 :: \Upsilon(Y) = \sum_{i \in I} \langle \phi_i, \Upsilon(Y) \rangle_H \varepsilon_i$$

such that :

$$\forall i \in I : \varepsilon_i = \Upsilon(e_i);$$

$$\forall i, j \in I : \langle \phi_i, \varepsilon_j \rangle_H = \delta_{ij};$$

ii) Υ defines a positive kernel on $V_0 : K_V(Y_1, Y_2) = \langle \Upsilon Y_1, \Upsilon Y_2 \rangle_H$

The sets (V_0, Υ, H) and $(V_0, \Upsilon U, H)$ are two realizations triple of K_V . Then there is an isometry φ such that :

$$\Upsilon U = \varphi \circ \Upsilon \text{ (Maths.1200).}$$

¹The positive kernel plays a role similar to the probability of transition between states of the Wigner's Theorem.

$$\begin{aligned}\langle UX_1, UX_2 \rangle_V &= \langle \Upsilon UX_1, \Upsilon UX_2 \rangle_H = \langle \varphi \circ \Upsilon X_1, \varphi \circ \Upsilon X_2 \rangle_H \\ &= \langle \Upsilon X_1, \Upsilon X_2 \rangle_H = \langle X_1, X_2 \rangle\end{aligned}$$

So U preserves the scalar product on V

Let be : $\widehat{U} = \Upsilon \circ U \circ \Upsilon^{-1}$

$$\begin{aligned}\langle \widehat{U}\psi_1, \widehat{U}\psi_2 \rangle_H &= \langle \Upsilon \circ U \circ (\Upsilon^{-1}\psi_1), \Upsilon \circ U \circ (\Upsilon^{-1}\psi_2) \rangle_H \\ &= \langle U \circ (\Upsilon^{-1}\psi_1), U \circ (\Upsilon^{-1}\psi_2) \rangle_V = \langle (\Upsilon^{-1}\psi_1), (\Upsilon^{-1}\psi_2) \rangle_V \\ &= \langle \psi_1, \psi_2 \rangle_H\end{aligned}$$

So \widehat{U} preserves the scalar product on H

iii) As seen in Theorem 2 starting from the basis $(\varepsilon_i)_{i \in I}$ of H one can define a Hermitian basis $(\tilde{\varepsilon}_i)_{i \in I}$ of H , an orthonormal basis $(\tilde{e}_i)_{i \in I}$ of V for the scalar product $K_V = \langle \cdot \rangle_V$ with $\tilde{e}_i = \Upsilon^{-1}(\tilde{\varepsilon}_i)$

U is defined for any vector of V , so for $(\tilde{e}_i)_{i \in I}$ of V .

$$\text{Define : } \widehat{U}(\tilde{\varepsilon}_i) = \widehat{U}(\Upsilon(\tilde{e}_i)) = \Upsilon(U(\tilde{e}_i)) = \tilde{\varepsilon}'_i$$

The set of vectors $(\tilde{\varepsilon}'_i)_{i \in I}$ is an orthonormal basis of H :

$$\langle \tilde{\varepsilon}'_i, \tilde{\varepsilon}'_j \rangle_H = \langle \widehat{U}(\Upsilon(\tilde{e}_i)), \widehat{U}(\Upsilon(\tilde{e}_j)) \rangle_H = \langle \tilde{e}_i, \tilde{e}_j \rangle_V = \delta_{ij}$$

The map : $\chi : \ell^2(I) \rightarrow H :: \chi(y) = \sum_{i \in I} y_i \tilde{\varepsilon}'_i$ is an isomorphism (same as in Theorem 2) and $(\tilde{\varepsilon}'_i)_{i \in I}$ is a Hilbertian basis of H . So we can write :

$$\forall \psi \in H : \psi = \sum_{i \in I} \psi^i \tilde{\varepsilon}'_i, \widehat{U}(\psi) = \sum_{i \in I} \psi^i \tilde{\varepsilon}'_i$$

$$\text{and : } \psi^i = \langle \tilde{\varepsilon}_i, \psi \rangle_H = \langle \widehat{U}(\tilde{\varepsilon}_i), \widehat{U}(\psi) \rangle_H = \langle \tilde{\varepsilon}'_i, \sum_{j \in I} \psi^j \tilde{\varepsilon}'_j \rangle_H = \psi^i$$

$$\text{Thus the map } \widehat{U} \text{ reads : } \widehat{U} : H \rightarrow H :: \widehat{U}(\sum_{i \in I} \psi^i \tilde{\varepsilon}_i) = \sum_{i \in I} \psi^i \tilde{\varepsilon}'_i$$

It is linear, continuous and unitary : $\langle \widehat{U}(\psi_1), \widehat{U}(\psi_2) \rangle = \langle \psi_1, \psi_2 \rangle$ and \widehat{U} is invertible

$U = \Upsilon^{-1} \circ \widehat{U} \circ \Upsilon$ is linear and bijective

iv) For any primary or secondary observable Φ there is a self-adjoint, Hilbert-Schmidt and trace class operator $\widehat{\Phi}$ on the associated Hilbert space H such that : $\widehat{\Phi} = \Upsilon \circ \Phi \circ \Upsilon^{-1}$. For the new variable the observable is $\Phi' \in L(V; W')$ and $W' \subset V$ is not necessarily identical to W . It is associated to the operator : $\widehat{\Phi}' = \Upsilon \circ \Phi' \circ \Upsilon^{-1}$. W and W' are finite dimensional vector subspaces of V .

$$\begin{array}{ccccccc} & & \Phi & & U & & \Phi' \\ W & \leftarrow & \leftarrow & \leftarrow & V & \rightarrow & \rightarrow & \rightarrow & V & \rightarrow & \rightarrow & \rightarrow & W' \\ \downarrow & & & & \downarrow & & & & \downarrow & & & & \downarrow \\ \downarrow & \Upsilon & & & \downarrow & \Upsilon & \Upsilon & \downarrow & & \Upsilon & & & \downarrow \\ \downarrow & & \widehat{\Phi} & & \downarrow & \widehat{U} & \downarrow & & \widehat{\Phi}' & & & & \downarrow \\ H_\Phi & \leftarrow & \leftarrow & \leftarrow & H & \rightarrow & \rightarrow & \rightarrow & H & \rightarrow & \rightarrow & \rightarrow & H_{\Phi'}\end{array}$$

Because U is a bijection on $V : \Phi' \circ U = U \circ \Phi \Rightarrow \Phi' = U \circ \Phi \circ U^{-1}$ and V is globally invariant by U

$$\Phi'(V) = W' = U \circ \Phi \circ U^{-1}(V) = U \circ \Phi(V) = U(W)$$

thus W' is a vector subspace of V isomorphic to W

$$\widehat{\Phi}' = \Upsilon \circ \Phi' \circ \Upsilon^{-1} = \Upsilon \circ U \circ \Phi \circ U^{-1} \circ \Upsilon^{-1} = \widehat{U} \circ \Upsilon \circ \Phi \circ \Upsilon^{-1} \circ \widehat{U}^{-1} = \widehat{U} \circ \widehat{\Phi} \circ \widehat{U}^{-1}$$

Let us denote : $\widehat{\Phi}(H) = H_{\Phi}$, $\widehat{\Phi}'(H) = H_{\Phi'}$
 $\widehat{U}(H) = H$ because it is a unitary map
 $\widehat{\Phi}'(H) = \widehat{U} \circ \widehat{\Phi} \circ \widehat{U}^{-1}(H) = \widehat{U} \circ \widehat{\Phi}(H) = \widehat{U}(H_{\Phi}) = H_{\Phi'}$
 thus $H_{\Phi'}$ is a vector subspace of H isomorphic to H_{Φ} ■

As a consequence the map U is necessarily linear, even if this was not assumed in the conditions 20 : variables which are not linearly related (in the conditions 20) cannot represent the same state.

As \widehat{U} is unitary, it cannot be self adjoint or trace class (except if $U = Id$). So it differs from an observable.

2.4.3 Change of units

A special case of this theorem is the choice of units to measure the variables. A change of units is a map : $X'_k = \alpha_k X_k$ with fixed scalars $(\alpha_k)_{k=1}^N$. As we must have :

$$\langle U(X_1), U(X_2) \rangle_V = \langle X_1, X_2 \rangle_V = \sum_{k=1}^N \alpha_k^2 \langle X_1, X_2 \rangle_V = \langle X_1, X_2 \rangle_V \Rightarrow \sum_{k=1}^N \alpha_k^2 = 1$$

which implies for any single variable X_k : $\alpha_k = 1$. So the variables in the model should be dimensionless quantities. This is in agreement with the elementary rule that any formal theory should not depend on the units which are used.

More generally whenever one has a law which relates quantities which are not expressed in the same units, there should be some fundamental constant involved, to absorb the discrepancy between the units. For instance some Physical laws involve an exponential, such as the wave equation for a plane wave :

$$\psi = \exp i \left(\langle \vec{k}, \vec{r} \rangle - \omega t \right)$$

They require that the argument in the exponential is dimensionless, and because \vec{r} is a length and t a time we should have a fundamental constant with the dimension of a speed (in this case c).

But also it implies that there should be some “universal system of units” (based on a single quantity) in which all quantities of the theory can be measured. In Physics this is the Planck’s system which relate the units of different quantities through the values of the fundamental constants c , G (gravity), R (Boltzmann constant), \hbar , and the charge of the electron (see Wikipedia for more).

Usually the variables are defined with respect to some frame, then the rules for a change of frame have a special importance and are a defining feature of the model. When the rules involve a group, the previous theorem can help to precise the nature of the abstract Hilbert space H and from there the choice of the maps X .

2.4.4 Group representation

The theory of group representation is a key tool in Physics. We will remind some basic results here, see Maths.23 for a comprehensive study of this topic.

The left action of a group G on a set E is a map $\lambda : G \times E \rightarrow E :: \lambda(g, x)$ such that $\lambda(gg', x) = \lambda(g, \lambda(g', x))$, $\lambda(1, x) = x$. And similarly for a right action $\rho(x, g)$.

The representation of a group G is a couple (E, f) of a vector space E and a continuous map $f : G \rightarrow GL(E; E)$ (the set of linear invertible maps from E to E) such that :

$$\forall g, g' \in G : f(g \cdot g') = f(g) \circ f(g'); f(1) = Id \Rightarrow f(g^{-1}) = f(g)^{-1}$$

A representation is **faithful** if f is bijective.

A vector subspace F is **invariant** if $\forall u \in F, g \in G : f(g)u \in F$

A representation is **irreducible** if there is no other invariant subspace than $E, 0$.

A representation is not unique : from a given representation one can build many others. The sum of the representations $(E_1, f_1), (E_2, f_2)$ is $(E_1 \oplus E_2, f_1 + f_2)$.

A representation is **unitary** if there is a scalar product on E and $f(g)$ is unitary : $\forall u, v \in F, g \in G : \langle f(g)u, f(g)v \rangle = \langle u, v \rangle$

If two groups G, G' are isomorphic by ϕ , then a representation (E, f) of G provides a representation of G' :

$$\phi : G' \rightarrow G :: \forall g, g' \in G' : \phi(g \cdot g') = \phi(g) \cdot \phi(g'); \phi(1_{G'}) = 1_G \Rightarrow \phi(g^{-1}) = \phi(g)^{-1}$$

$$f : G \rightarrow GL(E; E)$$

$$\text{Define } f' : G' \rightarrow GL(E; E) :: f'(g') = f(\phi(g'))$$

$$f'(g'_1 \cdot g'_2) = f(\phi(g'_1 \cdot g'_2)) = f(\phi(g'_2)) \circ f(\phi(g'_1)) = f'(g'_1) \circ f'(g'_2)$$

A **Lie group** is a group endowed with the structure of a manifold. On the tangent space T_1G at its unity (that we will denote 1) there is an algebraic structure of **Lie algebra**, that we will also denote generally T_1G , endowed with a bracket $[]$ which is a bilinear antisymmetric map on T_1G .

If G is a Lie group with Lie algebra T_1G and (E, f) a representation of G , then $(E, f'(1))$ is a representation of the Lie algebra T_1G :

$$f'(1) \in \mathcal{L}(T_1G; \mathcal{L}(E; E))$$

$$\forall X, Y \in T_1G : f'(1)([X, Y]) = f'(1)(X) \circ f'(1)(Y) - f'(1)(Y) \circ f'(1)(X)$$

The converse, from the Lie algebra to the group, holds if G is simply connected, otherwise a representation of the Lie algebra provides usually multiple valued representations of the group (we will see important examples later).

Any Lie group G has the **adjoint representation** (T_1G, Ad) over its Lie algebra.

Any irreducible representation of a commutative (abelian) group is unidimensional.

Any unitary representation of a compact or finite group is reducible in the sum of orthogonal, finite dimensional, irreducible unitary representations.

Any representation of a group on a finite dimensional vector space becomes a representation on a set of matrices by choosing a basis. The representations of the common groups of matrices are tabulated. In the standard representation

(K^n, ι) of a group G of $n \times n$ matrices on a field K the map ι is the usual action of matrices on column vectors in the space K^n . If G is a Lie group then the standard representation of its Lie algebra is the representation (K^n, ι) by matrices, deduced by derivation.

Two representations $(E, f), (F, \rho)$ of the same group G are **equivalent** if there is an isomorphism $\phi : E \rightarrow F$ such that :

$$\forall g \in G : f(g) = \phi^{-1} \circ \rho(g) \circ \phi$$

Then from a basis $(e_i)_{i \in I}$ of E one deduces a basis $|e_i\rangle$ of F by : $|e_i\rangle = \phi(e_i)$. Because ϕ is an isomorphism $|e_i\rangle$ is a basis of F . Moreover the matrix of the action of G is in this basis the same as for (E, f) :

$$\begin{aligned} \rho(g)|e_i\rangle &= \sum_{j \in J} [\rho(g)]_j^i |e_j\rangle = \rho(g) \phi(e_i) = \phi \circ f(g)(e_i) \\ &= \phi \left(\sum_{j \in I} [f(g)]_i^j e_j \right) = \sum_{p \in I} [f(g)]_i^j \phi(e_j) = \sum_{p \in I} [f(g)]_i^j |e_j\rangle \\ [\rho(g)] &= [f(g)] \end{aligned}$$

If K is a subgroup of G , and (E, f) a representation of G , then (E, f) is a subrepresentation of K .

The vector subspaces F of E which are invariant by K provide representations (F, f) of K .

2.4.5 Application to groups of transformations

Change of variable parametrized by a group

This is the usual case in Physics. The second point of view that we have noticed above is clear when U is defined by a group. The system is represented by fixed variables, and the measures are taken according to procedures which change with g and we have :

$$\Phi(g)(X) = U(g) \circ \Phi(1)(X)$$

$\Phi \in L(V; W)$ and $U(g)$ is a bijection so X and $\Phi(1)(X)$ are in bijective correspondence and X must belong to $W \subset V$: we reduce the definition of the states at what can be observed. And to assume that this is true for any observable leads to redefine X as in the first way, but this requires an additional assumption.

Theorem 22 *If the conditions 20 are met, and (V, U) is a representation of the group G , then:*

i) (H, \hat{U}) is a unitary representation of the group G with $\hat{U}(g) = \Upsilon \circ U(g) \circ \Upsilon^{-1}$

ii) For any observable $\Phi \in L(V; W)$ the vector space $W \subset V$ is invariant by U and (W, U) is a representation of G , and for the associated operator $\hat{\Phi} = \hat{U}(g) \circ \Phi \circ \hat{U}(g)^{-1} \in L(H; H_\Phi)$, (H_Φ, \hat{U}) is a finite dimensional unitary representation of the group G .

If G is a Lie group, and U continuous, then :

iii) U is smooth, \hat{U} is differentiable and $(\hat{U}'(1), H)$ is an anti-symmetric representation of the Lie algebra T_1G of G

iv) For any observable $\Phi \in L(V; W)$ $(H_\Phi, \widehat{U}'(1))$ is an anti-symmetric representation of the Lie algebra T_1G of G

If (F, f) is a unitary representation of G , equivalent to (H_Φ, \widehat{U}) , and Φ a primary or secondary observable, then :

v) The results of measures of Φ for two values $1, g$ and the same state of the system are related by :

$\Phi \circ U(1)(X) = \sum_{j \in J} X^j(1) e_j, \Phi \circ U(g)(X) = \sum_{j \in J} X^j(g) e_j$ for some basis $(e_i)_{i \in I}$ of V

$X^j(g) = \sum_{k \in J} [f(g)]_k^j X^k(1)$ where $[f(g)]$ is the matrix of $f(g)$ in orthonormal bases of F

vi) If moreover G is a Lie group and U, f continuous, then the action $U'(1)(\kappa_a)$ of $U'(1)$ for vectors κ_a of T_1G are expressed by the same matrices $[K_a]$ of the action $f'(1)(\kappa_a)$:

$$f'(1)(\kappa_a)(f_j) = \sum_{k \in J} [K_a]_j^k f_k \rightarrow U'(1)(\kappa_a)(e_j) = \sum_{k \in J} [K_a]_j^k e_k$$

$$\text{and similarly for the observable } \Phi : \Phi \circ U'(1)(\kappa_a)(e_j) = \sum_{k \in J} [K_a]_j^k e_k$$

Proof. i) The map $U : G \rightarrow GL(V; V)$ is such that $U(g \cdot g') = U(g) \circ U(g')$; $U(1) = Id$ where G is a group and 1 is the unit in G .

Then $U(g)$ is necessarily invertible, because $U(g^{-1}) = U(g)^{-1}$

$\widehat{U} : G \rightarrow \mathcal{L}(H; H) :: \widehat{U} = \Upsilon \circ U \circ \Upsilon^{-1}$ is such that :

$$\widehat{U}(g \cdot g') = \Upsilon \circ U(g \cdot g') \circ \Upsilon^{-1} = \Upsilon \circ U(g) \circ U(g') \circ \Upsilon^{-1} = \Upsilon \circ U(g) \circ \Upsilon^{-1} \circ \Upsilon \circ U(g') \circ \Upsilon^{-1} = \widehat{U}(g) \circ \widehat{U}(g')$$

$$\widehat{U}(1) = \Upsilon \circ U(1) \circ \Upsilon^{-1} = Id$$

So (H, \widehat{U}) is a unitary representation of the group G ($\widehat{U}(g)$ is bijective, thus invertible).

ii) For any observable $\Phi \circ U(g) = U(g) \circ \Phi, \widehat{\Phi} = \widehat{U}(g) \circ \widehat{\Phi} \circ \widehat{U}(g)^{-1}$

Let us take $Y \in W = \Phi(V) : \exists X \in V : Y = \Phi(X)$

$$U(g)Y = U(g)(\Phi(X)) = \Phi(U(g)X) \in \Phi(V)$$

And similarly

$$\widehat{Y} \in \widehat{\Phi}(H) : \exists \psi \in H : \widehat{Y} = \widehat{\Phi}(\psi)$$

$$\widehat{U}(g)\widehat{Y} = \widehat{U}(g)(\widehat{\Phi}(\psi)) = \widehat{\Phi}(\widehat{U}(g)\psi) \in \widehat{\Phi}(H)$$

thus $W, H_\Phi = \widehat{\Phi}(H)$ are invariant by U, \widehat{U}

The scalar product on H holds on the finite dimensional subspace $\widehat{\Phi}(H)$, which is a Hilbert space.

iii) If G is a Lie group and the map $U : G \rightarrow \mathcal{L}(V; V)$ continuous, then it is smooth (Maths.1789), \widehat{U} is differentiable and $(\widehat{U}'(1), H)$ is an anti-symmetric representation of the Lie algebra T_1G of G :

$$\forall \kappa \in T_1G : (\widehat{U}'(1)\kappa)^* = -(\widehat{U}'(1)\kappa)$$

$\widehat{U}(\exp \kappa) = \exp \widehat{U}'(1)\kappa$ where the first exponential is taken on T_1G and the second on $\mathcal{L}(H; H)$ (Maths.1886).

iv) $\widehat{\Phi}$ is a primary or secondary observable, and so is $\widehat{\Phi} \circ U(g)$, then $\widehat{\Phi} \circ \widehat{U}(g) = \widehat{U}(g) \circ \widehat{\Phi}$ is a self-adjoint, compact operator, and by the Riesz theorem (Math.1142) its spectrum is either finite or is a countable sequence converging to 0 (which may or not be an eigen value) and, except possibly for 0, is identical to the set $(\lambda_p(g))_{p \in \mathbb{N}}$ of its eigen values (Maths.1020). For each distinct eigen value the eigen spaces $H_p(g)$ are orthogonal and H is the direct sum $H = \bigoplus_{p \in \mathbb{N}} H_p(g)$. For each non null eigen value $\lambda_p(g)$ the eigen space $H_p(g)$ is finite dimensional. For a primary observable the eigen values are either 1 or 0.

Because $H_{\widehat{\Phi}}$ is finite dimensional, for each value of g there is an orthonormal basis $(\tilde{\varepsilon}_i(g))_{i \in J}$ of $H_{\widehat{\Phi}}$ comprised of a finite number of vectors which are eigen vectors of $\widehat{\Phi} \circ \widehat{U}(g) : \widehat{\Phi} \circ \widehat{U}(g) (\tilde{\varepsilon}_j(g)) = \lambda_j(g) \tilde{\varepsilon}_j(g)$

Any vector of $H_{\widehat{\Phi}}$ reads :

$$\psi = \sum_{j \in J} \psi^j(g) \tilde{\varepsilon}_j(g) \text{ and}$$

$\widehat{\Phi} \circ \widehat{U}(g) = \sum_{p \in \mathbb{N}} \lambda_p(g) \widehat{\pi}_{H_p(g)}$ with the orthogonal projection $\widehat{\pi}_{H_p(g)}$ on $H_p(g)$.

And, because any measure belongs to $H_{\widehat{\Phi}}$ it is a linear combination of eigen vectors

$$\begin{aligned} \widehat{\Phi} \circ U(g)(X) &= \Upsilon^{-1} \circ \widehat{\Phi} \circ \widehat{U}(g) \circ \Upsilon(X) = \Upsilon^{-1} \left(\sum_{j \in J} \lambda_j(g) \psi^j(g) \tilde{\varepsilon}_j(g) \right) \\ &= \sum_{j \in J} \lambda_j(g) \psi^j \Upsilon^{-1}(\tilde{\varepsilon}_j(g)) = \sum_{j \in J} \lambda_j(g) \psi^j e_j(g) \end{aligned}$$

for some basis $(e_i)_{i \in I}$ of V : $e_j(g) = \Upsilon^{-1}(\tilde{\varepsilon}_j(g))$ and $\widehat{\Phi} \circ U(g)(e_j(g)) = \lambda_j e_j(g)$

That we can write :

$$\begin{aligned} \widehat{\Phi} \circ U(g)(X) &= \sum_{j \in J} \lambda_j \psi^j(g) e_j(g) = \sum_{j \in J} X^j(g) e_j(g) = U(g) \circ \widehat{\Phi}(X) \\ \widehat{\Phi}(X) &= U(g^{-1}) \left(\sum_{j \in J} X^j(g) e_j(g) \right) \end{aligned}$$

v) If the representations $(H_{\widehat{\Phi}}, \widehat{U})$, (F, f) are equivalent (which happens if they have the same finite dimension) there is an isomorphism $\phi : H_{\widehat{\Phi}} \rightarrow F$ which can be defined by taking an orthonormal basis $(\tilde{\varepsilon}_i(g_0))_{i \in J}$, $(f_j(g_0))_{j \in J}$ in each vector space, for some fixed $g_0 \in G$ that we can take $g_0 = 1$: $\phi \left(\sum_{i \in J} \psi^i \tilde{\varepsilon}_i(1) \right) = \sum_{i \in J} \psi^i f_i(1) \Leftrightarrow \phi(\tilde{\varepsilon}_j(1)) = f_j(1)$

To a change of g corresponds a change of orthonormal basis, both in $H_{\widehat{\Phi}}$ and F , given by the known unitary map $f(g) : f_j(g) = f(g)(f_j(1)) = \sum_{k \in J} [f(g)]_j^k f_k(1)$ and thus we have the same matrix for $\widehat{U}(g)$:

$$\tilde{\varepsilon}_j(g) = \widehat{U}(g)(\tilde{\varepsilon}_j(1)) = \phi^{-1} \circ f(g) \circ \phi(\tilde{\varepsilon}_j(1)) = \phi^{-1} \circ f(g)(f_j(1)) = \sum_{k \in J} [f(g)]_j^k \tilde{\varepsilon}_k(1)$$

$$\begin{array}{ccccccc}
& & U(g) & & \Phi & & \\
V & \rightarrow & \rightarrow & V & \rightarrow & \rightarrow & W \\
\downarrow & & & \downarrow & & & \downarrow \\
\downarrow & \Upsilon & & \Upsilon & & \Upsilon & \downarrow \\
\downarrow & & \widehat{U}(g) & \downarrow & \widehat{\Phi} & \downarrow & \widehat{U}(g) \\
H & \rightarrow & \rightarrow & H & \rightarrow & \rightarrow & H_{\Phi} \rightarrow \rightarrow H_{\Phi} \\
& & & & & \downarrow & \\
& & & & & \phi & \downarrow \\
& & & & & \downarrow & \\
& & & & & F & \rightarrow \rightarrow F \\
& & & & & & f(g) \\
& & & & & & \downarrow \\
& & & & & & \phi \\
& & & & & & \downarrow \\
& & & & & & F
\end{array}$$

$$\tilde{\varepsilon}_j(g) = \widehat{U}(g)(\tilde{\varepsilon}_j(1)) = \sum_{k \in J} [f(g)]_j^k \tilde{\varepsilon}_k(1)$$

$$e_j(g) = \Upsilon^{-1}(\tilde{\varepsilon}_j(g)) = \Upsilon^{-1}\left(\sum_{k \in J} [f(g)]_j^k \tilde{\varepsilon}_k(1)\right)$$

$$= \sum_{k \in J} [f(g)]_j^k \Upsilon^{-1}(\tilde{\varepsilon}_k(1)) = \sum_{k \in J} [f(g)]_j^k e_k(1)$$

$$e_j(g) = \Upsilon^{-1} \circ \widehat{U}(g) \circ \Upsilon(e_j(1)) = U(g)(e_j(1))$$

Thus the matrix of $U(g)$ to go from 1 to g is $[f(g)]$

$$\Phi(X) = U(g^{-1})\left(\sum_{j \in J} X^j(g) e_j(g)\right)$$

$$\Phi \circ U(g)(X) = \sum_{j \in J} X^j(g) e_j(g) = \sum_{j \in J} X^j(g) \sum_{k \in J} [f(g^{-1})]_j^k e_k(1)$$

$$\Phi \circ U(g_0)(X) = \sum_{k \in J} X^k(1) e_k(1) \Rightarrow \sum_{j \in J} X^j(g) [f(g^{-1})]_j^k = X^k(1)$$

$$X^j(g) = \sum_{k \in J} [f(g)]_j^k X^k(1)$$

The measures $\Phi \circ U(g)(X)$ transform with the known matrix $f(g)$.

vi) $(H_{\Phi}, \widehat{U}'(1)), (F, f'(1))$ are equivalent, anti-symmetric (or anti-hermitian

for complex vector spaces) representations of the Lie algebra T_1G . If $(\kappa_a)_{a=1}^m$ is a basis of T_1G then $f'(1)$, which is a linear map, is defined by the values of $f'(1)(\kappa_a) \in L(F; F)$.

$$\begin{array}{cccc}
& & \widehat{U}'(1)(\kappa) & \\
H_{\Phi} & \rightarrow & \rightarrow & H_{\Phi} \\
\downarrow & & & \downarrow \\
\phi & \downarrow & & \phi \downarrow \\
\downarrow & & f'(1)(\kappa) & \downarrow \\
F & \rightarrow & \rightarrow & F
\end{array}$$

$$\widehat{U}'(1)(\kappa)(\psi) = \phi^{-1} \circ f'(1)(\kappa) \circ \phi(\psi)$$

If we know the values of the action of $f'(1)(\kappa_a)$ on any orthonormal basis $(f_j)_{j \in J}$ of F :

$$f'(1)(\kappa_a)(f_j) = \sum_{k \in J} [K_a]_j^k f_k$$

we have the value of $\widehat{U}'(1)(\kappa_a)$ for the corresponding orthonormal basis $(\widehat{\varepsilon}_j)_{j \in J}$ of H_{Φ}

$$\widehat{U}'(1)(\kappa_a)(\widehat{\varepsilon}_j) = \widehat{U}'(1)(\kappa_a) \phi^{-1}(f_j) = \phi^{-1} \circ f'(1)(\kappa_a)(f_j)$$

$$= \phi^{-1}\left(\sum_{k \in J} [K_a]_j^k f_k\right) = \sum_{k \in J} [K_a]_j^k \widehat{\varepsilon}_k$$

So $\widehat{U}'(1)$ is represented in an orthonormal basis of H_{Φ} by the same matrices $[K_a]$

And similarly :

$$\widehat{U}(g) = \Upsilon \circ U(g) \circ \Upsilon^{-1} \Rightarrow \widehat{U}'(1)(\kappa) = \Upsilon \circ U'(1)(\kappa) \circ \Upsilon^{-1}$$

$$U'(1)(\kappa_a)(e_j) = \Upsilon \circ U'(1)(\kappa_a) \circ \Upsilon^{-1}(e_j) = \Upsilon \circ U'(1)(\kappa_a)(\widehat{e}_j) = \Upsilon \left(\sum_{k \in J} [K_a]_j^k \widehat{e}_k \right) =$$

$$\sum_{k \in J} [K_a]_j^k e_k$$

vii) Because $\Phi \circ U(g) = U(g) \circ \Phi \Rightarrow \Phi \circ U'(1)(\kappa_a) = U'(1)(\kappa_a) \circ \Phi :$

$$\Phi \circ U'(1)(\kappa_a)(e_j) = \sum_{k \in J} [K_a]_j^k \Phi(e_k) \quad \blacksquare$$

This result is specially important in Physics. Any unitary representation of a compact or finite group is reducible in the sum of orthogonal, finite dimensional, irreducible unitary representations. As a consequence the space V of the variables X has the same structure. If, as it can be assumed, the state of the system stays in the same irreducible representation, it can belong only to some specific finite dimensional spaces, defined through the representation or an equivalent representation of G . X depends only on a finite number of parameters, This is the starting point of quantization.

Notice that the nature of the space E does not matter, only the matrices $[f(g)], [K]$.

Usually in Physics the changes are not parametrized by the group, but by a vector of the Lie algebra (for instance rotations are not parametrized by a matrix but by a vector representing the rotation), which gives a special interest to the two last results.

The usual geometric representations, based on frames defined through a point and a set of vectors, such as in Galilean Geometry and Special Relativity, have been generalized by the formalism of fiber bundles, which encompasses also General Relativity, and is the foundation of gauge theories. Gauge theories use abundantly group transformations, so they are a domain of choice to implement the previous results. We will see them in the following chapters.

Fourier transform

If G is an abelian group we have more. Irreducible representations of abelian groups are unidimensional, and any unitary representation of an abelian group is the sum of projections on unidimensional vector subspaces which, for infinite dimensional representations, takes the form of spectral integrals. More precisely, there is a bijective correspondence between the unitary representation of an abelian group G and the spectral measures on the Pontryagin dual \widehat{G} , which is the space of continuous maps $\vartheta : G \rightarrow T$ where T is the set of complex numbers of module 1 (Maths.1932). This can be made less abstract if G is a topological, locally compact group. Then it has a Haar measure μ and the representation (H, \widehat{U}) is equivalent to $(L^2(G, \mu, \mathbb{C}), \mathcal{F})$ that is to the Fourier transform \mathcal{F} on complex valued, square integrable, functions on G (Maths.2421).

If $\varphi \in L^2(G, \mu, \mathbb{C}) \cap L^1(G, \mu, \mathbb{C}) :$

$$\mathcal{F}(\varphi)(\vartheta) = \int_G \varphi(g) \overline{\vartheta(g)} \mu(g)$$

$$\mathcal{F}^*(h)(g) = \int_{\widehat{G}} h(\vartheta) \vartheta(g) \nu(\vartheta) \text{ for a unique Haar measure } \nu \text{ on } \widehat{G} \text{ and } \mathcal{F}^* = \mathcal{F}^{-1}$$

If G is a compact group then we have Fourier series on a space of periodic functions, and if G is a non compact, finite dimensional Lie group, G is isomorphic to some vector space E and we have the usual Fourier transform on functions on E .

These cases are important from a practical point of view as it is possible to replace the abstract Hilbert space H by more familiar spaces of functions, and usually one can assume that the space V is itself some Hilbert space. The previous tools (observables,...) are then directly available.

The most usual application is about periodic phenomena : whenever a system is inclosed in some box, it can be usually assumed that they are periodic (and null out of the box). Then the representation is naturally through Fourier series and we have convenient Hilbert bases.

One parameter groups

An important case, related to the previous one, is when the variables X depend on a scalar real argument, and the model is such that $X(t), X'(t') = X(t + \theta)$, with any fixed θ , represent the same state. The associated operator is parametrized by a scalar and we have a map :

$$\begin{aligned}\widehat{U} : \mathbb{R}_+ &\rightarrow \mathcal{GL}(H, H) \text{ such that :} \\ \widehat{U}(t + t') &= \widehat{U}(t) \circ \widehat{U}(t') \\ \widehat{U}(0) &= Id\end{aligned}$$

Then we have a one parameter semi-group. If moreover the map \widehat{U} is strongly continuous (that is $\lim_{\theta \rightarrow 0} \|\widehat{U}(\theta) - Id\| = 0$), it can be extended to \mathbb{R} . (\widehat{U}, H) is a unitary representation of the abelian group $(\mathbb{R}, +)$. We have a one parameter group, and because \widehat{U} is a continuous Lie group morphism it is differentiable with respect to θ (Maths.1784).

Any strongly continuous one parameter group of operators on a Banach vector space admits an infinitesimal generator $S \in \mathcal{L}(H; H)$ such that : $\widehat{U}(t) = \sum_{n=0}^{\infty} \frac{t^n}{n!} S^n = \exp tS$ (Maths.1033). By derivation with respect to t we get : $\frac{d}{ds} \widehat{U}(s) |_{t=s} = (\exp tS) \circ S \Rightarrow S = \frac{d}{ds} \widehat{U}(s) |_{t=0}$

Because $\widehat{U}(t)$ is unitary S is anti-hermitian :

$$\begin{aligned}\langle \widehat{U}(t) \psi, \widehat{U}(t) \psi' \rangle_H &= \langle \psi, \psi' \rangle_H \\ \Rightarrow \left\langle \frac{d}{dt} \widehat{U}(t) \psi, \widehat{U}(t) \psi' \right\rangle_H + \left\langle \widehat{U}(t) \psi, \frac{d}{dt} \widehat{U}(t) \psi' \right\rangle_H &= 0 \Rightarrow S = -S^*\end{aligned}$$

S is normal and has a spectral resolution P :

$$S = \int_{S_p(S)} sP(s)$$

S is anti-hermitian so its eigen-values are pure imaginary : $\lambda = -\bar{\lambda}$. $\widehat{U}(t)$ is not compact and S is not compact, usually its spectrum is continuous, so it is not associated to any observable.

We will see in the next section a striking application of this case.

2.4.6 Extension to manifolds

Several extensions of the theorem 2 can be considered. One problem that we will meet in the next chapters is the following. In a model variables X are maps defined on a manifold M , valued in a fixed vector space, and belong to a space V of maps with the required properties. But a variable Y is defined through X : $Y(m) = f(X(m))$ and belongs to a manifold $N(X)$ depending on X . So the conditions 1 do not apply.

To address this kind of problem we need to adapt our point of view. We have seen the full mathematical definition of a manifold in the first section. A manifold M is a class of equivalence : the same point m of M can be defined by several charts, maps $\varphi : E \rightarrow M$ from a vector space E to M , with different coordinates : $m = \varphi_a(\xi_a) = \varphi_b(\xi_b)$ so that it defines classes of equivalence between sets of coordinates : $\xi_a \sim \xi_b \Leftrightarrow \varphi_a(\xi_a) = \varphi_b(\xi_b)$. These classes of equivalence are made clear by the transitions maps $\chi_{ba} : E \rightarrow E$, which are bijective : $\xi_a \sim \xi_b \Leftrightarrow \xi_b = \chi_{ba}(\xi_a)$. And these transitions maps are the key characteristic of the manifold. To a point m of M corresponds a class of equivalence of coordinates.

So let us consider a system represented by a model which meets the following :

Condition 23 *The model is comprised of :*

i) A finite number of variables, collectively denoted X , which are maps valued in a vector space E and meeting the conditions 1 : they belong to an open subset O of a separable, infinite dimensional Fréchet space V .

ii) A variable Y , valued in a set F , defined by a map : $f : O \rightarrow F :: Y = f(X)$

iii) A collection of linear continuous bijective maps $\mathfrak{U} = (U_a \in GL(V; V))_{a \in A}$, comprising the identity, closed under composition : $\forall a, b \in A : U_a \circ U_b \in \mathfrak{U}$

iv) On V and F the equivalence relation :

$$R : X \sim X' \Leftrightarrow \exists a \in A : X' = U_a(X) : f(X) = f(X')$$

Denote the set $N = \{Y = f(X), X \in O\}$. The quotient set : N/R is comprised of classes of equivalence of points Y which can be defined by related coordinates. This is a manifold, which can be discrete and comprising only a finite number of points. One can also see the classes of equivalence of N/R as representing states of the system, defined equivalently by the variable $X, X' = U_a(X)$.

Notice that f is unique, no condition is required on E other than to be a vector space, and nothing on F . Usually the maps U_a are defined by : $U_a(X) = \chi_a \circ X$ where the maps $\chi_a \in GL(E; E)$ are bijective on E (not F or V) but only the continuity of U_a can be defined.

We have the following result :

Theorem 24 *For a system represented by a model meeting the conditions 23 :*

i) V can be embedded as an open of a Hilbert space H with a linear isometry $\Upsilon : V \rightarrow H$, to each U_a is associated the unitary operator $\widehat{U}_a = \Upsilon \circ U_a \circ \Upsilon^{-1}$ on

H , each class of equivalence $[V]_y$ of R on V is associated to a class of equivalence $[H]_y$ in H of :

$\widehat{R} : \psi \sim \psi' \Leftrightarrow \exists a \in A : \psi' = \widehat{U}_a(\psi)$. $[V]_y$ is a partition of V and $[H]_y$ of H .

ii) If (V, U) is a representation of a Lie group G , then (H, \widehat{U}) is a unitary representation of G and each $[H]_y$ is invariant by the action of G .

Proof. i) R defines a partition of V , we can label each class of equivalence by the value of Y , and pick one element X_y in each class :

$$\begin{aligned} [V]_y &= \{X \in O : f(X) \sim f(X_y) = y\} \equiv \{X \in O : \exists a \in A : X = U_a(X_y)\} \\ &\equiv \{X \in O : X = U_a(X_y), a \in A\} \end{aligned}$$

The variables X meet the conditions 1, O can be embedded as an open of a Hilbert space H and there is linear isomorphism : $\Upsilon : V \rightarrow H$

In $[V]_y$ the variables $X, X' = U_a(X)$ define the same state and we can implement the theorem 21. $\widehat{U}_a = \Upsilon \circ U_a \circ \Upsilon^{-1}$ is an unitary operator on H

$$\forall X \in [V]_y : \widehat{U}_a \circ \Upsilon(X_y) = \Upsilon \circ U_a(X_y) = \Upsilon(X)$$

The set $[H]_y = \Upsilon([V]_y) = \{\psi \in H : \psi = \widehat{U}_a(\Upsilon(X_y)), a \in A\}$ is the class of equivalence of :

$$\widehat{R} : \psi \sim \psi' \Leftrightarrow \exists a \in A : \psi' = \widehat{U}_a(\psi)$$

R defines a partition of $V : V = \cup_y [V]_y$ and \widehat{R} defines a partition of $H : H = \cup_y [H]_y$

ii) If (V, U) is a representation of a Lie group G then $[V]_y$ is the orbit of X_y , (H, \widehat{U}) is a unitary representation of G

Each $[H]_y$ is invariant by G . The vector subspace $[F]_y$ spanned by $[H]_y$ is invariant by G , so $([F]_y, \widehat{U})$ is a representation of G . ■

As a consequence of the last result : if U is a compact group, then the representation (H, \widehat{U}) is the sum of irreducible, orthogonal, finite dimensional representations. For each value of Y the subset $[H]_y$ is invariant by the action of G , so it must belong to one of the irreducible representations, as well as $[F]_y$. The maps X , for a given value of Y , belong to a finite dimensional vector space, and depend on a finite number of parameters. This is the usual meaning of the quantization of X .

2.5 THE EVOLUTION OF THE SYSTEM

In many models involving maps, the variables X_k are functions of the time t , which represents the evolution of the system. So this is a privileged argument of the functions. So far we have not made any additional assumption about the model : the open Ω of the Hilbert space contains all the possible values but, due to the laws to which it is subject, only some solutions will emerge, depending on the initial conditions. They are fixed by the value $X(0)$ of the variables at some origin 0 of time. They are specific to each realization of the system, but we should expect that the model and the laws provide a general solution, that is a map : $X(0) \rightarrow X$ which determines X for each specific occurrence of $X(0)$. It will happen if the laws are determinist. One says that the problem is well posed if for any initial conditions there is a unique solution X , and that X depends continuously on $X(0)$. We will give a more precise meaning of determinism by enlarging the conditions 1 as follows :

Condition 25 : *The model representing the system meets the conditions 1. Moreover :*

i) V is an infinite dimensional separable Fréchet space V of maps : $X = (X_k)_{k=1}^N :: R \rightarrow E$ where R is an open subset of \mathbb{R} and E a normed vector space

ii) $\forall t \in R$ the evaluation map : $\mathcal{E}(t) : V \rightarrow E : \mathcal{E}(t) X = X(t)$ is continuous

The laws for the evolution of the system are such that the variables $(X_k)_{k=1}^N$, which define the possible states considered for the system (that we call the admissible states) meet the conditions :

iii) The initial state of the system, defined at $t = 0 \in R$, belongs to an open subset A of E

iv) For any solutions X, X' belonging to O if the set $\varpi = \{t, X(t) = X'(t)\}$ has a non null Lebesgue measure then $X = X'$.

The last condition iv) means that the system is semi determinist : to the same initial conditions can correspond several different solutions, but if two solutions are equal on some interval then they are equal almost everywhere.

The condition ii) is rather technical and should be usually met. Practically it involves some relation between the semi-norms on V and the norm on E (this is why we need a norm on E) : when two variables X, X' are close in V , then their values $X(t), X'(t)$ must be close for almost all t . More precisely, because $\mathcal{E}(t)$ is linear, the continuity can be checked at $X = 0$ and reads:

$\forall t \in R, \forall X \in O : \forall \varepsilon > 0, \exists \eta : d(X, 0)_V < \eta \Rightarrow \|X(t)\|_E < \varepsilon$ where d is the metric on V

In all usual cases (such as L^p spaces or spaces of differentiable functions) $d(X, 0)_V \rightarrow 0 \Rightarrow \forall t \in R : \|X(t)\|_E \rightarrow 0$ and the condition ii) is met, but this is not a general result.

Notice that :

- the variables X can depend on any other arguments besides t as previously
- E can be infinite dimensional but must be normed
- no continuity condition is imposed on X .

2.5.1 Fundamental theorems for the evolution of a system

If the model meets the conditions 25 then it meets the conditions 1 : there is a separable, infinite dimensional, Hilbert space H , defined up to isomorphism, such that the states (admissible or not) \mathcal{S} belonging to O can be embedded as an open subset $\Omega \subset H$ which contains 0 and a convex subset. Moreover to any basis of V is associated a bijective linear map $\Upsilon : V \rightarrow H$.

Theorem 26 *If the conditions 25 are met, then there are :*

- i) a Hilbert space F , an open subset $\tilde{A} \subset F$
- ii) a map $\Theta : R \rightarrow \mathcal{L}(F; F)$ such that $\Theta(t)$ is unitary and, for the admissible states $X \in O \subset V$:
 $X(0) \in \tilde{A} \subset F$
 $\forall t : X(t) = \Theta(t)(X(0)) \in F$
- iii) for each value of t an isometry $\hat{\mathcal{E}}(t) \in \mathcal{L}(H; F)$ such that for the admissible states $X \in O \subset V$:
 $\forall X \in O : \hat{\mathcal{E}}(t)\Upsilon(X) = X(t)$
 where H is the Hilbert space and Υ is the linear chart associated to X and any basis of V

Proof. i) Define the equivalence relation on V :

$$\mathcal{R} : X \sim X' \Leftrightarrow X(t) = X'(t) \text{ for almost every } t \in R$$

and take the quotient space V/\mathcal{R} , then the set of admissible states is a set \tilde{O} such that :

$$\tilde{O} \in O \subset V$$

$$\forall X \in \tilde{O} : X(0) \in A$$

$$\forall X, X' \in \tilde{O}, \forall t \in R : X(t) = X'(t) \Rightarrow X = X'$$

ii) Define :

$$\forall t \in R : \tilde{F}(t) = \left\{ X(t), X \in \tilde{O} \right\} \text{ thus } \tilde{F}(0) = A$$

A is a subset of E . There are families of independent vectors belonging to A , and a largest family $(f_j)_{j \in J}$ of independent vectors. It generates a vector space $F(0)$ which is a vector subspace of E , containing A .

$$\forall u \in F(0) : \exists (x_j)_{j \in J} \in \mathbb{R}_0^J : u = \sum_{j \in J} x_j f_j$$

The map :

$$\tilde{\Theta}(t) : \tilde{F}(0) \rightarrow \tilde{F}(t) :: \tilde{\Theta}(t)u = \mathcal{E}(t) \circ \mathcal{E}(0)^{-1}u$$

is bijective and continuous

The set $F(t) = \tilde{\Theta}(t)F(0) \subset E$ is well defined by linearity :

$$\tilde{\Theta}(t) \left(\sum_{j \in J} x_j f_j \right) = \sum_{j \in J} x_j \tilde{\Theta}(t)(f_j)$$

The map $\tilde{\Theta}(t) : F(0) \rightarrow F(t)$ is linear, bijective, continuous on an open subset A , thus continuous, and the spaces $F(t)$ are isomorphic, vector subspaces of E , containing $\tilde{F}(t)$.

Define $(\varphi_j)_{j \in J}$ the largest family of independent vectors of

$$\left\{ \tilde{\Theta}(t)(f_j), t \in R \right\}. \text{ This is a family of independent vectors of } E, \text{ which}$$

generates a subspace \tilde{F} of E , containing each of the $F(t)$ and thus each of the

$\tilde{F}(t)$. Moreover each of the φ_j is the image of a unique vector f_j for some $t_j \in R$.

The map $\tilde{\Theta}(t)$ is then a continuous linear map $\tilde{\Theta}(t) \in \mathcal{L}(\tilde{F}; \tilde{F})$

iii) The conditions of proposition 1 are met for O and V, so there are a Hilbert space H and a linear map $\Upsilon : O \rightarrow \Omega$

Each of the φ_j is the image of a unique vector f_j for some $t \in R$, and thus there is a uniquely defined family $(X_j)_{j \in J}$ of \tilde{O} such that $X_j(t_j) = \varphi_j$.

Define on \tilde{F} the bilinear symmetric definite positive form with coefficients :

$$\begin{aligned} \langle \varphi_j, \varphi_k \rangle_{\tilde{F}} &= K_V \left(\mathcal{E}(t_j)^{-1} \varphi_j, \mathcal{E}(t_k)^{-1} \varphi_k \right) \\ &= \left\langle \Upsilon \mathcal{E}(t_j)^{-1} \varphi_j, \Upsilon \mathcal{E}(t_k)^{-1} \varphi_k \right\rangle_H = \langle X_j, X_k \rangle_H \end{aligned}$$

By the Gram-Schmidt procedure we can build an orthonormal basis $(\tilde{\varphi}_j)_{j \in J}$ of \tilde{F} : $\tilde{F} = \text{Span}(\tilde{\varphi}_j)_{j \in J}$ and the Hilbert vector space: $F = \left\{ \sum_{j \in J} \tilde{x}_j \tilde{\varphi}_j, (\tilde{x}_j)_{j \in J} \in \ell^2(J) \right\}$ which is a vector space containing \tilde{F} (but is not necessarily contained in E).

iv) The map $\tilde{\Theta}(t) \in \mathcal{L}(\tilde{F}; \tilde{F})$ is a linear homomorphism, \tilde{F} is dense in F, thus $\tilde{\Theta}(t)$ can be extended to a continuous operator $\Theta(t) \in \mathcal{L}(F; F)$ (Math.1003).

$\tilde{\Theta}(t)$ is unitary on \tilde{F} : $\langle u, v \rangle_{\tilde{F}} = K_V \left(\mathcal{E}(0)^{-1} u, \mathcal{E}(0)^{-1} v \right)$ so $\Theta(t)$ is unitary on F.

iv) Define the map :

$$\hat{\mathcal{E}}(t) : \Omega \rightarrow F :: \hat{\mathcal{E}}(t) \Upsilon(X) = X(t)$$

where $\Omega \subset H$ is the open associated to V and O.

For $X \in \tilde{O}$:

$$\hat{\mathcal{E}}(t) \Upsilon(X) = X(t) = \tilde{\Theta}(t) X = \mathcal{E}(t) \circ \mathcal{E}(0)^{-1} X$$

$$\hat{\mathcal{E}}(t) = \mathcal{E}(t) \circ \mathcal{E}(0)^{-1} \circ \Upsilon^{-1}$$

$\hat{\mathcal{E}}(t)$ is linear, continuous, bijective on Ω , it is an isometry :

$$\left\langle \hat{\mathcal{E}}(t) \psi, \hat{\mathcal{E}}(t) \psi' \right\rangle_F = \langle X(t), X'(t) \rangle_F = \langle \Upsilon X, \Upsilon X' \rangle_H = \langle \psi, \psi' \rangle_H$$

v) $A = \tilde{F}(0)$ is an open subset of $F(0)$, which is itself an open vector subspace of F. Thus A can be embedded as an open subset \tilde{A} of F. ■

When X depends on other arguments ξ , the result reads :

$$\forall t, \forall \xi : X(t, \xi) = \Theta(t) (X(0, \xi)) \in F$$

Indeed the basic feature which is used is :

$$\forall X, X' \in \tilde{O}, \forall t \in R : X(t) = X'(t) \Rightarrow X = X'$$

$$\text{which means : } \forall t, \forall \xi : X(t, \xi) = X'(t, \xi) \Leftrightarrow X = X'$$

As a consequence the model is determinist, up to the equivalence between maps almost everywhere equal. But the operator $\Theta(t)$ depends on t and not necessarily continuously, so the problem is not necessarily well posed. Notice that each solution $X(t)$ belong to E, but the Hilbert space F can be larger than E. Moreover the result holds if the conditions apply to some variables only.

But we have a stronger result.

Theorem 27 *If the model representing the system meets the conditions 1 and moreover :*

i) V is an infinite dimensional separable Fréchet space V of maps : $X = (X_k)_{k=1}^N :: R \rightarrow E$ where E is a normed vector space

ii) $\forall t \in \mathbb{R}$ the evaluation map : $\mathcal{E}(t) : V \rightarrow E : \mathcal{E}(t) X = X(t)$ is continuous

iii) the variables $X'_k(t) = X_k(t + \theta)$ and $X_k(t)$ represent the same state of the system, for any $t' = t + \theta$ with a fixed $\theta \in \mathbb{R}$

then :

i) there is a continuous map $S \in \mathcal{L}(V; V)$ such that :

$$\mathcal{E}(t) = \mathcal{E}(0) \circ \exp tS$$

$$\forall t \in \mathbb{R} : X(t) = (\exp tS \circ X)(0) = \left(\sum_{n=0}^{\infty} \frac{t^n}{n!} S^n X \right)(0)$$

and the operator $\hat{S} = \Upsilon \circ S \circ \Upsilon^{-1}$ associated to S is anti-hermitian

ii) there are a Hilbert space F , an open $\tilde{A} \subset F$, a continuous anti-hermitian map $\tilde{S} \in \mathcal{L}(F; F)$ such that :

$$\forall X \in O \subset V : X(0) \in \tilde{A} \subset F$$

$$\forall t : X(t) = \left(\exp t\tilde{S} \right) (X(0)) \in F$$

iii) The maps X are smooth and :

$$\frac{d}{ds} X(s) |_{s=t} = \tilde{S} X(t)$$

Proof. i) We have a change of variables U depending on a parameter $\theta \in \mathbb{R}$ which reads with the evaluation map : $\mathcal{E} : \mathbb{R} \times V \rightarrow F :: \mathcal{E}(t) X = X(t) :$

$$\forall t, \theta \in \mathbb{R} : \mathcal{E}(t)(U(\theta) X) = \mathcal{E}(t + \theta)(X) \Leftrightarrow \mathcal{E}(t)U(\theta) = \mathcal{E}(t + \theta) = \mathcal{E}(\theta)U(t):$$

U defines a one parameter group of linear operators:

$$U(\theta + \theta') X(t) = X(t + \theta + \theta') = U(\theta) \circ U(\theta') X(t)$$

$$U(0) X(t) = X(t)$$

It is obviously continuous at $\theta = 0$ so it is continuous.

ii) The conditions 1 are met, so there are a Hilbert space H , a linear chart Υ , and $\hat{U} : \mathbb{R} \rightarrow \mathcal{L}(H; H)$ such that $\hat{U}(\theta)$ is linear, bijective, unitary :

$$\forall X \in O : \hat{U}(\theta)(\Upsilon(X)) = \Upsilon(U(\theta)(X))$$

$$\hat{U}(\theta + \theta') = \Upsilon \circ U(\theta + \theta') \circ \Upsilon^{-1} = \Upsilon \circ U(\theta) \circ U(\theta') \circ \Upsilon^{-1} = \Upsilon \circ U(\theta) \circ \Upsilon^{-1} \circ \Upsilon \circ U(\theta') \circ \Upsilon^{-1} = \hat{U}(\theta) \circ \hat{U}(\theta')$$

$$\hat{U}(0) = \Upsilon \circ U(0) \circ \Upsilon^{-1} = Id$$

The map : $\hat{U} : \mathbb{R} \rightarrow \mathcal{L}(H; H)$ is uniformly continuous with respect to θ , it defines a one parameter group of unitary operators. So there is an anti-hermitian operator \hat{S} with spectral resolution P such that :

$$\hat{U}(\theta) = \sum_{n=0}^{\infty} \frac{\theta^n}{n!} \hat{S}^n = \exp \theta \hat{S}$$

$$\frac{d}{ds} \hat{U}(s) |_{\theta=s} = \left(\exp \theta \hat{S} \right) \circ \hat{S}$$

$$\hat{S} = \int_{Sp(S)} s P(s)$$

$$\left\| \hat{U}(\theta) \right\| = 1 \leq \exp \left\| \theta \hat{S} \right\|$$

iii) $S = \Upsilon^{-1} \circ \hat{S} \circ \Upsilon$ is a continuous map on the largest vector subspace V_0 of V which contains O , which is a normed vector space with the norm induced by the positive kernel.

$\|S\| \leq \|\Upsilon^{-1}\| \|\widehat{S}\| \|\Upsilon\| = \|\widehat{S}\|$ because Υ is an isometry.

So the series $\sum_{n=0}^{\infty} \frac{\theta^n}{n!} S^n$ converges in V_0 and :

$$U(\theta) = \Upsilon^{-1} \circ \widehat{U}(\theta) \circ \Upsilon = \sum_{n=0}^{\infty} \frac{\theta^n}{n!} S^n = \exp \theta S$$

$$\forall \theta, t \in \mathbb{R} : U(\theta) X(t) = X(t + \theta) = (\exp \theta S) X(t)$$

$$\mathcal{E}(t) \exp \theta S = \mathcal{E}(t + \theta)$$

Exchange θ, t and take $\theta = 0$:

$$\mathcal{E}(\theta) \exp tS = \mathcal{E}(t + \theta)$$

$$\mathcal{E}(0) \exp tS = \mathcal{E}(t) \in \mathcal{L}(V; E)$$

which reads :

$$\forall t \in \mathbb{R} : U(t) X(0) = X(t) = (\exp tS) X(0)$$

(U, V_0) is a continuous representation of $(\mathbb{R}, +)$, U is smooth and X is smooth

:

$$\frac{d}{ds} U(s) X(0) |_{s=t} = \frac{d}{ds} X(s) |_{s=t} = SX(t)$$

$$\Leftrightarrow \frac{d}{ds} \mathcal{E}(s) |_{s=t} = S\mathcal{E}(t)$$

The same result holds whatever the size of O in V , so S is defined over V .

iv) The set : $F(t) = \{X(t), X \in V\}$ is a vector subspace of E .

Each map is fully defined by its value at one point :

$$\forall t \in \mathbb{R} : X(t) = (\exp tS \circ X)(0)$$

$$X(t) = X'(t) \Rightarrow \forall \theta : X(t + \theta) = X'(t + \theta) \Leftrightarrow X = X'$$

So the conditions 4 are met.

$$\Theta(t) : F(0) \rightarrow F(t) :: \Theta(t) u = \mathcal{E}(t) \circ \mathcal{E}(0)^{-1} u = \mathcal{E}(0) \circ \exp tS \circ \mathcal{E}(0)^{-1} u$$

The map $\Theta(\theta) : F \rightarrow F$ defines a one parameter group, so it has an infinitesimal generator $\widetilde{S} \in \mathcal{L}(F; F) : \Theta(\theta) = \exp \theta \widetilde{S}$ and because $\Theta(\theta)$ is unitary \widetilde{S} is anti-hermitian.

$$\frac{d}{ds} \Theta(s) X(0) |_{s=t} = \frac{d}{ds} X(s) |_{s=t} = \widetilde{S} X(t) \quad \blacksquare$$

As a consequence such a model is necessarily determinist, and the system is represented by smooth maps whose evolution is given by a unique operator. It is clear that the conditions 25 are then met, so this case is actually a special case of the previous one. Notice that, even if X was not assumed to be continuous, smoothness is a necessary result. This result can seem surprising, but actually the basic assumption about a translation in time means that the laws of evolution are smooth, and as a consequence the variables depend smoothly on the time. And conversely this implies that, whenever there is some discontinuity in the evolution of the system, the conditions above cannot hold : time has a specific meaning, related to a change in the environment.

Comments

The conditions above depend deeply on how the time is understood in the model. We have roughly two cases :

A) t is a parameter used only to identify a temporal location. In Galilean Geometry the time is independent from the spatial coordinates for any observer and one can consider a change of coordinates such as : $t' = t + \theta$ with any constant θ . The variables X, X' such that $X'(t') = X(t + \theta)$ represent the

same system. Similarly in Relativist Geometry the universe can be modelled as a manifold, and a change of coordinates with affine parameters, $\xi' = \xi + \theta$ with a fixed 4 vector θ , is a change of charts. The components of any quantity defined on the tensorial tangent bundle change according to the jacobian $\left[\frac{\partial \xi'}{\partial \xi} \right]$ which is the identity, so the corresponding variables represent the same system. Then we are usually in the conditions of the Theorem 27, and this is the basis of the Schrödinger equation.

B) t is a parameter used to measure the duration of a phenomenon, usually the time elapsed since some specific event, and it is clear that the origin of time matters and the variables X, X' such that $X'(t') = X(t + \theta)$ do not represent the same system. This is the case in more specific models, such as in Engineering. The proposition 27 does not hold, but the proposition 26 holds if the model is determinist.

The conditions 25 require at least that all the variables which are deemed significant are accounted for. As it has been discussed in the previous chapter, usually probabilist laws appear because some of them are missing. The Theorem 26 precises this issue : by denoting the missing variables Y , one needs to enlarge the vector space E , and similarly F . The map $\Theta(t)$ still exists, but it encompasses the couples $(X(t), Y(t))$. The dispersion of the observed values of $X(t)$ are then imputed to the distribution of the unknown values $Y(t)$.

2.5.2 Observables

When a system is studied through its evolution, the observables can be considered from two different points of view :

- in the movie way : the estimation of the parameters is done at the end of the period considered, from a batch of data corresponding to several times (which are not necessarily the same for all variables). So this is the map X which is estimated through an observable $X \rightarrow \Phi(X)$.

- in the picture way : the estimation is done at different times (the same for all the variables which are measured). So there are the values $X(t)$ which are estimated. Then the estimation of $X(t)$ is given by $\varphi(X(t)) = \varphi(\mathcal{E}(t) X)$, with φ a linear map from E to a finite dimensional vector space, which usually does not depend on t (the specification stays the same).

In the best scenario the two methods should give the same result, which reads :

$$\varphi(\mathcal{E}(t) X) = \mathcal{E}(t) (\Phi X) \Leftrightarrow \varphi = \mathcal{E}(t) \circ \Phi \circ \mathcal{E}(t)^{-1}$$

But usually, when it is possible, the first way gives a better statistical estimation.

2.5.3 Phases Transitions

There is a large class of problems which involve transitions in the evolution of a system. They do not involve the maps X , which belong to the same family as above, but the values $X(t)$ which are taken over a period of time in some vector

space E . There are distinct subsets of E , that we will call **phases** (to avoid any confusion with states which involves the map X), between which the state of the system goes during its evolution, such as the transition solid / gas or between magnetic states. The questions which arise are then : what are the conditions, about the initial conditions or the maps X , for the occurrence of such an event ? Can we forecast the time at which such event takes place ?

Staying in the general model meeting the conditions 25, the first issue is the definition of the phases. The general idea is that they are significantly different states, and it can be formalized by : the set $\{X(t), t \in R, X \in O\}$ is disconnected, it comprises two disjoint subsets E_1, E_2 closed in E .

If the maps $X : R \rightarrow F$ are continuous and R is an interval of \mathbb{R} (as we will assume) then the image $X(R)$ is connected, the maps X cannot be continuous, and we cannot be in the conditions of proposition 27 (a fact which is interesting in itself), but we can be in the case of proposition 26. This is a difficult but also very common issue : in the real life such discontinuous evolutions are the rule. However, as we have seen, in the physical world discontinuities happen only at isolated points : the existence of a singularity is what makes interesting a change of phase. If the transition points are isolated, there is an open subset of R which contains each of them, a finite number of them in each compact subset of R , and at most a countable number of transition points. A given map X is then continuous (with respect to t) except in a set of points $(\theta_\alpha)_{\alpha \in A}, A \subset \mathbb{N}$. If $X(0) \in E_1$ then the odd transition points $\theta_{2\alpha+1}$ mark a transition $E_1 \rightarrow E_2$ and the opposite for the even points $\theta_{2\alpha}$.

If the conditions 25 are met then Θ is continuous except in $(\theta_\alpha)_{\alpha \in A}$, the transition points do not depend on the initial state $X(0)$, but the phase on each segment does. Then it is legitimate to assume that there is some probability law which rules the occurrence of a transition. We will consider two cases.

The simplest assumption is that the probability of the occurrence of a transition at any time t is constant. Then it depends only on the cumulated lengths of the periods $T_1 = \sum_{\alpha=0} [\theta_{2\alpha}, \theta_{2\alpha+1}], T_2 = \sum_{\alpha=0} [\theta_{2\alpha+1}, \theta_{2\alpha+2}]$ respectively.

Let us assume that $X(0) \in E_1$ then the changes $E_1 \rightarrow E_2$ occur for $t = \theta_{2\alpha+1}$, the probability of transitions read :

$$\begin{aligned} \Pr(X(t+\varepsilon) \in E_2 | X(t) \in E_1) &= \Pr(\exists \alpha \in \mathbb{N} : t + \varepsilon \in [\theta_{2\alpha+1}, \theta_{2\alpha+2}]) \\ &= T_2 / (T_1 + T_2) \\ \Pr(X(t+\varepsilon) \in E_1 | X(t) \in E_2) &= \Pr(\exists \alpha \in \mathbb{N} : t + \varepsilon \in [\theta_{2\alpha}, \theta_{2\alpha+1}]) \\ &= T_1 / (T_1 + T_2) \\ \Pr(X(t) \in E_1) &= T_1 / [R]; \Pr(X(t) \in E_2) = T_2 / [R] \end{aligned}$$

The probability of a transition at t is : $T_2 / (T_1 + T_2) \times T_1 / (T_1 + T_2) + T_1 / (T_1 + T_2) \times T_2 / (T_1 + T_2) = 2T_1T_2 / (T_1 + T_2)^2$. It does not depend of the initial phase, and depends only on Θ . This probability law can be checked from a batch of data about the values of T_1, T_2 for each observed transition.

However usually the probability of a transition depends on the values of the variables. The phases are themselves characterized by the value of $X(t)$,

so a sensible assumption is that the probability of a transition increases with the proximity of the other phase . Using the Hilbert space structure of F it is possible to address practically this case.

If E_1, E_2 are *closed convex subsets* of F , which is a Hilbert space, there is a unique map $\pi_1 : F \rightarrow E_1$. The vector $\pi_1(x)$ is the unique $y \in E_1$ such that $\|x - y\|_F$ is minimum. The map π_1 is continuous and $\pi_1^2 = \pi_1$. And similarly for E_2 .

The quantity $r = \|X(t) - \pi_1(X(t))\|_F + \|X(t) - \pi_2(X(t))\|_F$ = the distance to the other subset than where $X(t)$ lies, so one can assume that the probability of a transition at t is $f(r)$ where $f : \mathbb{R} \rightarrow [0, 1]$ is a probability density. The probability of a transition depends only on the state at t , but one cannot assume that the transitions points θ_α do not depend on X .

The result holds if E_1, E_2 are closed *vector subspaces* of F such that $E_1 \cap E_2 = \{0\}$. Then

$$X(t) = \pi_1(X(t)) + \pi_2(X(t))$$

$$\text{and } \|X(t)\|^2 = \|\pi_1(X(t))\|^2 + \|\pi_2(X(t))\|^2$$

$\frac{\|\pi_1(X(t))\|^2}{\|X(t)\|^2}$ can be interpreted as the probability that the system at t is in the phase E_1 .

One important application is forecasting a transition for a given map X . From the measure of $X(t)$ one can compute for each t the quantity $r(t) = \|X(t) - \pi_1(X(t))\|_F + \|X(t) - \pi_2(X(t))\|_F$ and, if we know f , we have the probability of a transition at t . The practical problem is then to estimate f from the measure of r over a past period $[0, T]$. A very simple, non parametric, estimator can be built when X are maps depending only of t (see J.C.Dutailly *Estimation of the probability of transitions between phases*). It can be used to forecast the occurrence of events such as earth quakes.

2.6 INTERACTING SYSTEMS

2.6.1 Representation of interacting systems

In the propositions above no assumption has been done about the interaction with exterior variables. If the values of some variables are given (for instance to study the impact of external factors with the system) then they shall be fully integrated into the set of variables, at the same footing as the others.

A special case occurs when one considers two systems S_1, S_2 , which are similarly represented, meaning that that we have the same kind of variables, defined as identical mathematical objects and related significance. To account for the interactions between the two systems the models are of the form :

$$\begin{array}{ccc}
 \lrcorner & S_1 & \ulcorner \\
 X_1 & & Z_1 \\
 V_1 & \times & W_1 \\
 & \downarrow \Upsilon_1 & \\
 & \psi_1 & \\
 & H_1 & \\
 & & \lrcorner & S_2 & \ulcorner \\
 & & X_2 & & Z_2 \\
 & & V_2 & \times & W_2 \\
 & & & \downarrow \Upsilon_2 & \\
 & & & \psi_2 & \\
 & & & H_2 & \\
 & & \lrcorner & S_{1+2} & \ulcorner \\
 & & X_1 & & X_2 \\
 & & V_1 & \times & V_2 \\
 & & & & \\
 & & \psi_1 & & \psi_2 \\
 & & H_1 & \times & H_2
 \end{array}$$

X_1, X_2 are the variables (as above X denotes collectively a set of variables) characteristic of the systems S_1, S_2 , and Z_1, Z_2 are variables representing the interactions. Usually these variables are difficult to measure and to handle. One can consider the system S_{1+2} with the direct product $X_1 \times X_2$, but doing so we obviously miss the interactions Z_1, Z_2 .

We see now how it is possible to build a simpler model which keeps the features of S_1, S_2 and accounts for their interactions.

We consider the models without interactions (so with only X_1, X_2) and we assume that they meet the conditions 1. For each model $S_k, k = 1, 2$ there are a linear map : $\Upsilon_k : V_k \rightarrow H_k :: \Upsilon_k (X_k) = \psi_k = \sum_{i \in I_k} \langle \phi_{ki}, \psi_k \rangle e_{ki}$
a positive kernel : $K_k : V_k \times V_k \rightarrow \mathbb{R}$

Let us denote S the new model. Its variables will be collectively denoted Y, valued in a Fréchet vector space V'. There will be another Hilbert space H', and a linear map $\Upsilon' : V' \rightarrow H'$ similarly defined. As we have the choice of the model, we will impose some properties to Y and V' in order to underline both that they come from S_1, S_2 and that they are interacting.

Condition 28 *i) The variable Y can be deduced from the value of X_1, X_2 : there must be a bilinear map : $\Phi : V_1 \times V_2 \rightarrow V'$*

ii) Φ must be such that whenever the systems S_1, S_2 are in the states ψ_1, ψ_2 then S is in the state ψ' and

$$\Upsilon'^{-1}(\psi') = \Phi(\Upsilon_1^{-1}(\psi_1), \Upsilon_2^{-1}(\psi_2))$$

iii) The positive kernel is a defining feature of the models, so we want a positive kernel K' of (V', Υ') such that :

$$\forall X_1, X'_1 \in V_1, \forall X_2, X'_2 \in V_2 :$$

$$K'(\Phi(X_1, X_2), \Phi(X'_1, X'_2)) = K_1(X_1, X'_1) \times K_2(X_2, X'_2)$$

We will prove the following :

Theorem 29 *Whenever two systems S_1, S_2 interact, there is a model S encompassing the two systems and meeting the conditions 28 above. It is obtained by taking the tensor product of the variables specific to S_1, S_2 . Then the Hilbert space of S is the tensorial product of the Hilbert spaces associated to each system.*

Proof. First let us see the consequences of the conditions if they are met.

The map : $\varphi : H_1 \times H_2 \rightarrow H' :: \varphi(\psi_1, \psi_2) = \Phi(\Upsilon_1^{-1}(\psi_1), \Upsilon_2^{-1}(\psi_2))$ is bilinear. So, by the universal property of the tensorial product, there is a unique map $\widehat{\varphi} : H_1 \otimes H_2 \rightarrow H'$ such that : $\varphi = \widehat{\varphi} \circ \iota$ where $\iota : H_1 \times H_2 \rightarrow H_1 \otimes H_2$ is the tensorial product (Maths.369).

The condition iii) reads :

$$\begin{aligned} & \langle \Upsilon_1(X_1), \Upsilon_1(X'_1) \rangle_{H_1} \times \langle \Upsilon_2(X_2), \Upsilon_2(X'_2) \rangle_{H_2} \\ &= \langle (\Upsilon' \circ \Phi(\Upsilon_1(X_1), \Upsilon_2(X_2)), \Upsilon' \circ \Phi(\Upsilon_1(X'_1), \Upsilon_2(X'_2))) \rangle_{H'} \\ & \langle \psi_1, \psi'_1 \rangle_{H_1} \times \langle \psi_2, \psi'_2 \rangle_{H_2} = \langle \varphi(\psi_1, \psi_2), \varphi(\psi'_1, \psi'_2) \rangle_{H'} \\ &= \langle \widehat{\varphi}(\psi_1 \otimes \psi_2), \widehat{\varphi}(\psi'_1 \otimes \psi'_2) \rangle_{H'} \end{aligned}$$

The scalar products on H_1, H_2 extend in a scalar product on $H_1 \otimes H_2$, endowing the latter with the structure of a Hilbert space with :

$$\langle (\psi_1 \otimes \psi_2), (\psi'_1 \otimes \psi'_2) \rangle_{H_1 \otimes H_2} = \langle \psi_1, \psi'_1 \rangle_{H_1} \langle \psi_2, \psi'_2 \rangle_{H_2}$$

and then the reproducing kernel is the product of the reproducing kernels (Maths.1208).

So we must have : $\langle \widehat{\varphi}(\psi_1 \otimes \psi_2), \widehat{\varphi}(\psi'_1 \otimes \psi'_2) \rangle_{H'} = \langle \psi_1 \otimes \psi_2, \psi'_1 \otimes \psi'_2 \rangle_{H_1 \otimes H_2}$ and $\widehat{\varphi}$ must be an isometry : $H_1 \otimes H_2 \rightarrow H'$

So by taking $H' = H_1 \otimes H_2$ and $V' = V_1 \otimes V_2$ we meet the conditions. ■

The conditions above are a bit abstract, but are logical and legitimate in the view of the Hilbert spaces. They lead to a natural solution, which is not unique and makes sense only if the systems are defined by similar variables. The measure of the tensor S can be addressed as before, the observables being linear maps defined in the tensorial products $V_1 \otimes V_2, H_1 \otimes H_2$ and valued in finite dimensional vector subspaces of these tensor products.

Comments

A key point in this representation is the difference between the simple direct product : $V_1 \times V_2$ and the tensorial product $V_1 \otimes V_2$, an issue about which there is much confusion.

The knowledge of the states (X_1, X_2) of both systems requires two vectors of I components each, that is $2 \times I$ scalars, and the knowledge of the state S requires a vector of I^2 components. So the measure of S requires more data, and brings more information, because it encompasses all the interactions. Moreover *a tensor is not necessarily the tensorial product of vectors* (if it is so it is said to be **decomposable**), it is the sum of such tensors. There is no canonical map $: V_1 \otimes V_2 \rightarrow V_1 \times V_2$. So *there is no simple and unique way to associate two vectors (X_1, X_2) to one tensor S* . This seems paradoxical, as one could imagine that both systems can always be studied, and their states measured, even if they are interacting. But the simple fact that we consider interactions means that the measure of the state of one of the system shall account for the conditions in which the measure is done, so it shall precise the value of the state of the other system and of the interactions Z_1, Z_2 .

If a model is arbitrary, its use must be consistent : if the scientist assumes that there are interactions, they must be present somewhere in the model, as variables for the computations as well as data to be collected. They can be dealt with in two ways. Either we opt for the two systems model, and we have to introduce the variables Z_1, Z_2 representing the interactions, then we have two separate models as in the first section. The study of their interactions can be a topic of the models, but this is done in another picture and requires additional hypotheses about the laws of the interactions. Or, if we intend to account for both systems and their interactions in a single model, we need a representation which supports more information that can bring $V_1 \times V_2$. The tensorial product is one way to enrich the model, this is the most economical and, as far as one follows the guidelines i),ii),iii) above, the only one. The complication in introducing general tensors is the price that we have to pay to account for the interactions. This representation does not, in any way, imply anything about *how* the systems interact, or even if they interact at all (in this case S is always decomposable). As usual the choice is up to the scientist, based upon how he envisions the problem at hand. But he has to live with his choice.

This issue is at the root of the paradoxes of entanglement. With many variants it is an experiment which involves two objects, which interact at the beginning, then are kept separated and non interacting, and eventually one measures the state of one of the two objects, from which the state of the other can be deduced with some probability. If we have two objects which interact at some point, with a significant result because it defines a new state, and we compare their states, then we must either incorporate the interactions, or consider that they constitute a single system and use the tensorial product. The fact that the objects cease to interact at some point does not matter : they are considered together if we compare their states. The interactions must be accounted for, one way or another and, when an evolution is considered, this is the map which represents the whole of the evolution which is significant, not its value at some time.

A common interpretation of this representation is to single out decomposable tensors $\Psi = \psi_1 \otimes \psi_2$, called “pure states”, so that actual states would be a superposition of pure states (a concept popularized by the famous Schrödinger’s

cat). It is clear that in an interacting system the pure states are an abstraction, which actually would represent two non interacting systems, so their superposition is an artificial construction. It can be convenient in simple cases, where the states of each system can be clearly identified, or in complicated models to represent quantities which are defined over the whole system as we will see later. But it does not imply any mysterious feature, notably any probabilist behavior, for the real systems. A state of the two interacting systems is represented by a single tensor, and a tensor is not necessarily decomposable, but it is a sum of decomposable tensors.

2.6.2 Homogeneous systems

The previous result can be extended to N (a number that we will assumed to be fixed) similar systems (that we will call **microsystems**), represented by the same model, interacting together. For each microsystem, identified by a label s , the Hilbert space H and the linear map Υ are the same, the state S of the total system can be represented as a vector belonging to the tensorial product $\mathbf{V}_N = \otimes_{s=1}^N V$, associated to a tensor Ψ belonging to the tensorial product $\mathbf{H}_N = \otimes_{s=1}^N H$. The linear maps $\Upsilon \in \mathcal{L}(V; H)$ can be uniquely extended as maps $\Upsilon_N \in \mathcal{L}(\mathbf{V}_N; \mathbf{H}_N)$ such that (Maths.423) :

$$\Upsilon_N(X_1 \otimes \dots \otimes X_N) = \Upsilon(X_1) \otimes \dots \otimes \Upsilon(X_N)$$

The state of the system is then totally defined by the value of tensors S, Ψ , with I^N components.

We have general properties on these tensorial products (Maths.1208).

If $(\tilde{\varepsilon}_i)_{i \in I}$ is a Hilbertian basis of H then $E_{i_1 \dots i_N} = \tilde{\varepsilon}_{i_1} \otimes \dots \otimes \tilde{\varepsilon}_{i_N}$ is a Hilbertian basis of $\otimes_{s=1}^N H$. The scalar product is defined by linear extension of

$$\langle \Psi, \Psi' \rangle_{\mathbf{H}_N} = \langle \psi_1, \psi'_1 \rangle_H \times \dots \times \langle \psi_N, \psi'_N \rangle_H$$

for decomposable tensors : $\Psi = \psi_1 \otimes \dots \otimes \psi_N, \Psi' = \psi'_1 \otimes \dots \otimes \psi'_N$.

The subspaces $\otimes_{s=1}^p H \otimes \tilde{\varepsilon}_i \otimes_{s=p+2}^N H$ are orthogonal and $\otimes_{s=1}^N H \simeq \ell^2(I^N)$

Any operator on H can be extended on $\otimes_{s=1}^N H$ with similar properties : a self adjoint, unitary or compact operator extends uniquely as a self adjoint, unitary or compact operator (Maths.1211).

In the general case the label matters : the state $S = X_1 \otimes \dots \otimes X_N$ is deemed different from $S = X_{\sigma(1)} \otimes \dots \otimes X_{\sigma(N)}$ where $(X_{\sigma(p)})_{p=1}^N$ is a permutation of $(X_s)_{s=1}^N$. If the microsystems have all the same behavior they are, for the observer, indistinguishable. Usually the behavior is related to a parameter analogous to a size, so in such cases the microsystems are assumed to have the same size. We will say that these interacting systems are homogeneous :

Definition 30 *A homogeneous system is a system comprised of a fixed number N of microsystems, represented in the same model, such that any permutation of the N microsystems gives the same state of the total system.*

We have the following result :

Proposition 31 *The states Ψ of homogeneous systems belong to an open subset of a subspace \mathbf{h} of the Hilbert space $\otimes_{s=1}^N H$, defined by :*

i) a class of conjugacy $\mathfrak{S}(\lambda)$ of the group of permutations $\mathfrak{S}(N)$, defined itself by a decomposition of N in p parts :

$$\lambda = \{0 \leq n_p \leq \dots \leq n_1 \leq N, n_1 + \dots + n_p = N\}.$$

ii) p distinct vectors $(\tilde{\varepsilon}_j)_{j=1}^p$ of a Hermitian basis of H which together define a subspace H_J

iii) The space \mathbf{h} of tensors representing the states of the system is then :

either the symmetric tensors belonging to : $\odot_{n_1} H_J \otimes \odot_{n_2} H_J \dots \otimes \odot_{n_p} H_J$

or the antisymmetric tensors belonging to : $\wedge_{n_1} H_J \otimes \wedge_{n_2} H_J \dots \otimes \wedge_{n_p} H_J$

Proof. i) In the representation of the general system the microsystems are identified by some label $s = 1 \dots N$. An exchange of labels $U(\sigma)$ is a change of variables, represented by an action of the group of permutations $\mathfrak{S}(N)$: U is defined uniquely by linear extension of $U(\sigma)(X_1 \otimes \dots \otimes X_N) = X_{\sigma(1)} \otimes \dots \otimes X_{\sigma(N)}$ on decomposable tensors.

We can implement the Theorem 22 proven previously. The tensors ψ representing the states of the system belong to a Hilbert space $\mathbf{H}_N \subset \otimes_{s=1}^N H$ such that (\mathbf{H}_N, \hat{U}) is a unitary representation of $\mathfrak{S}(N)$. Which implies that \mathbf{H}_N is invariant by \hat{U} . The action of \hat{U} on $\otimes_{s=1}^N H$ is defined uniquely by linear extension of

$$\hat{U}(\sigma)(\psi_1 \otimes \dots \otimes \psi_N) = \psi_{\sigma(1)} \otimes \dots \otimes \psi_{\sigma(N)} \text{ on decomposable tensors.}$$

$$\Psi \in \otimes_{s=1}^N H \text{ reads in a Hilbert basis } (\tilde{\varepsilon}_i)_{i \in I} \text{ of } H :$$

$$\Psi = \sum_{i_1 \dots i_N \in I} \Psi^{i_1 \dots i_N} \tilde{\varepsilon}_{i_1} \otimes \dots \otimes \tilde{\varepsilon}_{i_N} \text{ and :}$$

$$\hat{U}(\sigma)\Psi = \sum_{i_1 \dots i_N \in I} \Psi^{i_1 \dots i_N} \hat{U}(\sigma)(\tilde{\varepsilon}_{i_1} \otimes \dots \otimes \tilde{\varepsilon}_{i_N}) = \sum_{i_1 \dots i_N \in I} \Psi^{i_1 \dots i_N} \tilde{\varepsilon}_{\sigma(i_1)} \otimes \dots \otimes \tilde{\varepsilon}_{\sigma(i_N)}$$

$$= \sum_{i_1 \dots i_N \in I} \Psi^{\sigma(i_1) \dots \sigma(i_N)} \tilde{\varepsilon}_{i_1} \otimes \dots \otimes \tilde{\varepsilon}_{i_N}$$

$$\langle \hat{U}(\sigma)\Psi, \hat{U}(\sigma)\Psi' \rangle = \langle \Psi, \Psi' \rangle$$

$$\Leftrightarrow \sum_{i_1 \dots i_N \in I} \Psi^{\sigma(i_1) \dots \sigma(i_N)} \Psi'^{\sigma(i_1) \dots \sigma(i_N)} = \sum_{i_1 \dots i_N \in I} \Psi^{i_1 \dots i_N} \Psi'^{i_1 \dots i_N}$$

The only vector subspaces of $\otimes_{s=1}^N H$ which are invariant by \hat{U} and on which \hat{U} is unitary are spaces of symmetric or antisymmetric tensors :

$$\text{symmetric : } \Psi^{\sigma(i_1) \dots \sigma(i_N)} = \Psi^{i_1 \dots i_N}$$

$$\text{antisymmetric : } \Psi^{\sigma(i_1) \dots \sigma(i_N)} = \epsilon(\sigma) \Psi^{i_1 \dots i_N}$$

ii) $\mathfrak{S}(N)$ is a finite, compact group. Its unitary representations are the sum of orthogonal, finite dimensional, unitary, irreducible representations (Maths.1948).

Let $\mathbf{h} \subset \otimes_{s=1}^N H$ be an irreducible, finite dimensional, representation of \hat{U} . Then $\forall \sigma \in \mathfrak{S}(N) : \hat{U}(\sigma)\mathbf{h} \subset \mathbf{h}$

iii) Let J a finite subset of I with $\text{card}(J) \geq N$, H_J the associated Hilbert space, $\hat{Y}_J : H \rightarrow H_J$ the projection, and $\hat{Y}_{J_N} = \otimes_N \hat{Y}_J$ be the extension of \hat{Y}_J to $\otimes_{s=1}^N H$:

$$\hat{Y}_{J_N} \left(\sum_{i_1 \dots i_N \in I} \Psi^{i_1 \dots i_N} \tilde{\varepsilon}_{i_1} \otimes \dots \otimes \tilde{\varepsilon}_{i_N} \right) = \sum_{i_1 \dots i_N \in J} \Psi^{i_1 \dots i_N} \tilde{\varepsilon}_{i_1} \otimes \dots \otimes \tilde{\varepsilon}_{i_N}$$

Then :

$$\forall \sigma \in \mathfrak{S}(N) : \hat{U}(\sigma)\hat{Y}_{J_N} \left(\sum_{i_1 \dots i_N \in I} \Psi^{i_1 \dots i_N} \tilde{\varepsilon}_{i_1} \otimes \dots \otimes \tilde{\varepsilon}_{i_N} \right)$$

$$= \sum_{i_1 \dots i_N \in J} \Psi^{\sigma(i_1) \dots \sigma(i_N)} \tilde{\varepsilon}_{i_1} \otimes \dots \otimes \tilde{\varepsilon}_{i_N} = \hat{Y}_{J_N} \hat{U}(\sigma) \Psi$$

So if \mathbf{h} is invariant by \hat{U} then $\hat{Y}_{J_N} \mathbf{h}$ is invariant by \hat{U} . If (\mathbf{h}, \hat{U}) is an irreducible representation then the only invariant subspace are 0 and \mathbf{h} itself, so necessarily $\mathbf{h} \subset \hat{Y}_{J_N} (\otimes_{s=1}^N H)$ for $\text{card}(J) = N$. Which implies : $\mathbf{h} \subset \otimes_N H_J$ with $H_J = \hat{Y}_J H$ and $\text{card}(J) = N$.

iv) There is a partition of $\mathfrak{S}(N)$ in conjugacy classes $\mathfrak{S}(\lambda)$ which are subgroups defined by a decomposition of N in p parts :

$\lambda = \{0 \leq n_p \leq \dots \leq n_1 \leq N, n_1 + \dots + n_p = N\}$. Notice that there is an order on the sets $\{\lambda\}$. Each element of a conjugacy class is then defined by a repartition of the integers $\{1, 2, \dots, N\}$ in p subsets of n_k items (this is a Young Tableau) (Maths. 5.2.2). A class of conjugacy is an abelian subgroup of $\mathfrak{S}(N)$: its irreducible representations are unidimensional.

The irreducible representations of $\mathfrak{S}(N)$ are then defined by a class of conjugacy, and the choice of a vector.

\mathbf{h} is a Hilbert space, thus it has a Hilbertian basis, composed of decomposable tensors which are of the kind $\tilde{\varepsilon}_{j_1} \otimes \dots \otimes \tilde{\varepsilon}_{j_N}$ where $\tilde{\varepsilon}_{j_k}$ are chosen among the vectors of a Hermitian basis $(\tilde{\varepsilon}_j)_{j \in J}$ of H_J

If $\tilde{\varepsilon}_{j_1} \otimes \dots \otimes \tilde{\varepsilon}_{j_N} \in H, \forall \sigma \in \mathfrak{S}(N) : \hat{U}(\sigma) \tilde{\varepsilon}_{j_1} \otimes \dots \otimes \tilde{\varepsilon}_{j_N} = \tilde{\varepsilon}_{j_{\sigma(1)}} \otimes \dots \otimes \tilde{\varepsilon}_{j_{\sigma(N)}} \in \mathbf{h}$ and because the representation is irreducible the basis of \mathbf{h} is necessarily composed from a set of $p \leq N$ vectors $\tilde{\varepsilon}_j$ by action of $\hat{U}(\sigma)$

Conversely : for any Hermitian basis $(\tilde{\varepsilon}_i)_{i \in I}$ of H , any subset J of cardinality N of I , any conjugacy class λ , any family of vectors $(\tilde{\varepsilon}_{j_k})_{k=1}^p$ chosen in $(\tilde{\varepsilon}_i)_{i \in J}$, the action of \hat{U} on the tensor :

$$\Psi_\lambda = \otimes_{n_1} \tilde{\varepsilon}_{j_1} \otimes_{n_2} \tilde{\varepsilon}_{j_2} \dots \otimes_{n_p} \tilde{\varepsilon}_{j_p}, j_1 \leq j_2 \dots \leq j_p$$

$$\text{gives the same tensor if } \sigma \in \mathfrak{S}(\lambda) : \hat{U}(\sigma) \Psi_\lambda = \Psi_\lambda$$

gives a different tensor if $\sigma \in \mathfrak{S}(\lambda^c)$ the conjugacy class complementary to $\mathfrak{S}(\lambda) : \mathfrak{S}(\lambda^c) = \mathfrak{C}_{\mathfrak{S}(N)}^{\mathfrak{S}(\lambda)}$

so it provides an irreducible representation by :

$$\forall \Psi \in \mathbf{h} : \Psi = \sum_{\sigma \in \mathfrak{S}(\lambda^c)} \Psi^\sigma \hat{U}(\sigma) (\otimes_{n_1} \tilde{\varepsilon}_{j_1} \otimes_{n_2} \tilde{\varepsilon}_{j_2} \dots \otimes_{n_p} \tilde{\varepsilon}_{j_p})$$

where the components Ψ^σ are labeled by the vectors of a basis of \mathbf{h} . The dimension of \mathbf{h} is given by the cardinality of $\mathfrak{S}(\lambda^c)$ that is : $\frac{N!}{n_1! \dots n_p!}$. All the vector spaces \mathbf{h} of the same conjugacy class (but different vectors $\tilde{\varepsilon}_i$) have the same dimension, thus they are isomorphic.

v) A basis of \mathbf{h} is comprised of tensorial products of N vectors of a Hilbert basis of H . So we can give the components of the tensors of \mathbf{h} with respect to $\otimes_{s=1}^N H$. We have two non equivalent representation :

By symmetric tensors : \mathbf{h} is then isomorphic to $\odot_{n_1} H_J \otimes \odot_{n_2} H_J \dots \otimes \odot_{n_p} H_J$ where the symmetric tensorial product \odot and the space of n order symmetric tensor on H_J is $\odot_n H_J$

By antisymmetric tensors : \mathbf{h} is then isomorphic to $\wedge_{n_1} H_J \otimes \wedge_{n_2} H_J \dots \otimes \wedge_{n_p} H_J$ and the space of n order antisymmetric tensor on H_J is $\wedge_n H_J$

The result extends to V_N by : $S = \Upsilon_N^{-1}(\Psi)$ ■

Remarks

i) For each choice of a class of conjugacy, and each choice of the vectors $(\tilde{\varepsilon}_j)_{j=1}^p$ which defines H_J , we have a different irreducible representation with vector space \mathbf{h} . Different classes of conjugacy gives non equivalent representations. But different choices of the Hermitian basis $(\tilde{\varepsilon}_j)_{j \in I}$ and the subset J of I , for a given class of conjugacy, give equivalent representations, and they can be arbitrary. So, for a given system, the set of states is characterized by a subset J of N elements in any basis of H , and by a class of conjugacy.

A change of the state of the system can occur either inside the same vector space \mathbf{h} , or between irreducible representations: $\mathbf{h} \rightarrow \mathbf{h}'$. As we will see in the next chapters usually the irreducible representation is fixed by other variables (such that energy) and a change of irreducible representation implies a discontinuous process. The states of the total system are quantized by the interactions.

ii) $\otimes_{n_1} \tilde{\varepsilon}_{j_1} \otimes_{n_2} \tilde{\varepsilon}_{j_2} \dots \otimes_{n_p} \tilde{\varepsilon}_{j_p}$ can be seen as representing a configuration where n_k microsystems are in the same state $\tilde{\varepsilon}_{j_k}$. The class of conjugacy, characterized by the integers n_p , correspond to the distribution of the microsystems between fixed states.

iii) If O is a convex subset then S belongs to a convex subset, and the basis can be chosen such that $\forall \Psi \in \mathbf{h}$ is a linear combination $(y_k)_{k=1}^q$ of the generating tensors with $y_k \in [0, 1]$, $\sum_{k=1}^q y_k = 1$. S can then be identified to the expected value of a random variable which would take one of the value $\otimes_{n_1} X_1 \otimes_{n_2} X_2 \dots \otimes_{n_p} X_p$, which corresponds to n_k microsystems having the state X_k . As exposed above the identification with a probabilist model is formal : there is no random behavior assumed for the physical system.

iv) In the probabilist picture one can assume that each microsystem behaves independently, and has a probability π_j to be in the state represented by $\tilde{\varepsilon}_j$ and $\sum_{j=1}^N \pi_j = 1$. Then the probability that we have $(n_k)_{k=1}^p$ microstates in the states $(\tilde{\varepsilon}_k)_{k=1}^p$ is $\frac{N!}{n_1! \dots n_p!} (\pi_{j_1})^{n_1} \dots (\pi_{j_p})^{n_p}$.

v) The set of symmetric tensor $\odot_n H_J$ is a closed vector subspace of $\otimes_n H_J$, this is a Hilbert space, $\dim \odot_n H_J = C_{p+n-1}^{p-1}$ with Hilbertian basis $\frac{1}{\sqrt{n!}} \odot_{j \in J} \tilde{\varepsilon}_j = \frac{1}{\sqrt{n!}} S_n (\otimes_{j \in J} \tilde{\varepsilon}_j)$ where the symmetrizer is :

$$S_n \left(\sum_{(i_1 \dots i_n)} \psi^{i_1 \dots i_n} \tilde{\varepsilon}_{i_1} \otimes \dots \otimes \tilde{\varepsilon}_{i_n} \right) = \sum_{(i_1 \dots i_n)} \psi^{i_1 \dots i_n} \sum_{\sigma \in \mathfrak{S}(n)} \tilde{\varepsilon}_{\sigma(1)} \otimes \dots \otimes \tilde{\varepsilon}_{\sigma(n)}$$

A tensor is symmetric iff : $\Psi \in \odot_n H_J \Leftrightarrow S_n(\Psi) = n! \Psi$ (Maths. 7.2.1, 13.5.2).

The set of antisymmetric tensor $\wedge_n H_J$ is a closed vector subspace of $\otimes_n H_J$, this is a Hilbert space, $\dim \wedge_n H_J = C_p^n$ with Hilbertian basis $\frac{1}{\sqrt{n!}} \wedge_{j \in J} \tilde{\varepsilon}_j = \frac{1}{\sqrt{n!}} A_n (\otimes_{j \in J} \tilde{\varepsilon}_j)$ with the antisymmetrizer :

$$A_n \left(\sum_{(i_1 \dots i_n)} \psi^{i_1 \dots i_n} \tilde{\varepsilon}_{i_1} \otimes \dots \otimes \tilde{\varepsilon}_{i_n} \right) = \sum_{(i_1 \dots i_n)} \psi^{i_1 \dots i_n} \sum_{\sigma \in \mathfrak{S}(n)} \epsilon(\sigma) \tilde{\varepsilon}_{\sigma(1)} \otimes \dots \otimes \tilde{\varepsilon}_{\sigma(n)}$$

A tensor is antisymmetric iff : $\Psi \in \wedge_n H_J \Leftrightarrow A_n(\Psi) = n! \Psi$ (Maths. 7.2.2, 13.5.2)

v) for $\theta \in \mathfrak{S}(N) : \widehat{U}(\theta) \Psi$ is usually different from Ψ

2.6.3 Global observables of homogeneous systems

The previous definitions of observables can be extended to homogeneous systems. An observable is defined on the total system, this is a map : $\Phi : \mathbf{V}_N \rightarrow W$ where W is a finite dimensional vector subspace of \mathbf{V}_N , but not necessarily a tensorial vector product of spaces. To Φ is associated the self-adjoint operator $\widehat{\Phi} = \Upsilon \circ \Phi \circ \Upsilon^{-1}$ and $H_\Phi = \widehat{\Phi} \left(\otimes_{s=1}^N H \right) \subset \otimes_{s=1}^N H$.

Theorem 32 *Any observable of a homogeneous system is of the form :*

$\Phi : \mathbf{V}_N \rightarrow W$ where W is generated by vectors Φ_λ associated to each class of conjugacy of $\mathfrak{S}(N)$

The value of $\Phi(X_1 \otimes \dots \otimes X_N) = \varphi(X_1, \dots, X_N) \Phi_\lambda$ where φ is a scalar linear symmetric map, if the system is in a state corresponding to λ

Proof. The space W must be invariant by U and H_Φ invariant by \widehat{U} . If the system is in a state belonging to \mathbf{h} for a class of conjugacy λ , then $H_\Phi = \widehat{\Phi}\mathbf{h}$ and $(\widehat{\Phi}\mathbf{h}, \widehat{U})$ is an irreducible representation of the abelian subgroup $\mathfrak{S}(\lambda)$ corresponding to λ . It is necessarily unidimensional and $\Phi(X_1 \otimes \dots \otimes X_N)$ is proportional to a unique vector. The observable being a linear map, the function φ is a linear map of the components of the tensor. ■

There is no way to estimate the state of each microsystem. From a practical point of view, this is a vector $\gamma = \widehat{\Phi} \left(\otimes_{n_1} \tilde{\varepsilon}_{j_1} \otimes_{n_2} \tilde{\varepsilon}_{j_2} \dots \otimes_{n_p} \tilde{\varepsilon}_{j_p} \right)$ which is measured, and from it $\lambda, (\tilde{\varepsilon}_{j_k})_{k=1}^p$ are estimated.

In the probabilist picture the expected value of γ is :

$$\langle \gamma \rangle = z(\pi_1, \dots, \pi_N)$$

with

$$z(\pi_1, \dots, \pi_N) = \sum_{\lambda} \frac{N!}{n_1! \dots n_p!} \sum_{1 \leq j_1 \leq \dots \leq j_p \leq N} (\pi_{j_1})^{n_1} \dots (\pi_{j_p})^{n_p} \widehat{\Phi} \left(\otimes_{n_1} \varepsilon_{j_1} \dots \otimes_{n_p} \varepsilon_{j_p} \right)$$

We have a classic statistical problem : estimate the π_i from a statistic given by the measure of γ . If the statistic $\widehat{\Phi}$ is sufficient, meaning that π_i depends only on γ , as F is finite dimensional whatever the number of microsystems, the Pitman-Koopman-Darmois theorem tells us that the probability law is exponential, then an estimation by the maximum likelihood gives the principle of Maximum Entropy with entropy :

$$E = - \sum_{j=1}^N \pi_j \ln \pi_j$$

In the usual interpretation of the probabilist picture, it is assumed that the state of each microsystem can be measured independently. Then the entropy $E = - \sum_{j=1}^N \pi_j \ln \pi_j$ can be seen as a measure of the heterogeneity of the system. And, contrary to a usual idea, the interactions between the micro-systems do not lead to the homogenization of their states, but to their quantization : the states are organized according to the classes of conjugacy.

2.6.4 Evolution of homogeneous systems

The evolution of homogeneous systems raises many interesting issues. The assumptions are a combination of the previous conditions.

Theorem 33 For a model representing the evolution of a homogeneous system comprised of a fixed number N of microsystems $s = 1 \dots N$ which are represented by the same model, with variables $(X_s)_{s=1}^N$ such that, for each microsystem :

i) the variables X_s are maps : $X_s :: R \rightarrow E$ where R is an open subset of \mathbb{R} and E a normed vector space, belonging to an open subset O of an infinite dimensional Fréchet space V

ii) $\forall t \in R$ the evaluation map : $\mathcal{E}(t) : O \rightarrow E : \mathcal{E}(t) X_s = X_s(t)$ is continuous

iii) $\forall t \in R : X_s(t) = X'_s(t) \Rightarrow X_s = X'_s$

There is a map : $S : R \rightarrow \otimes_N F$ such that $S(t)$ represents the state of the system at t . $S(t)$ takes its value in a vector space $f(t)$ such that $(\mathbf{f}(t), \widehat{U}_F)$, where \widehat{U}_F is the permutation on $\otimes_N F$, is an irreducible representation of $\mathfrak{S}(N)$

The crucial point is that the homogeneity is understood as the microsystems follow the same laws, but at a given time they do not have necessarily the same state.

Proof. i) Implement the Theorem 2 for each microsystem : there is a common Hilbert space H associated to V and a continuous linear map $\Upsilon : V \rightarrow H :: \psi_s = \Upsilon(X_s)$

ii) Implement the Theorem 31 on the homogeneous system, that is for the whole of its evolution. The state of the system is associated to a tensor $\Psi \in \mathbf{h}$ where \mathbf{h} is defined by a Hilbertian basis $(\tilde{\varepsilon}_i)_{i \in I}$ of H , a finite subset J of I , a conjugacy class λ and a family of p vectors $(\tilde{\varepsilon}_{jk})_{k=1}^p$ belonging to $(\tilde{\varepsilon}_i)_{i \in J}$. The vector space \mathbf{h} stays the same whatever t .

iii) Implement the Theorem 26 on the evolution of each microsystem : there is a common Hilbert space F , a map : $\widehat{\mathcal{E}} : R \rightarrow \mathcal{L}(H; F)$ such that : $\forall X_s \in O : \widehat{\mathcal{E}}(t) \Upsilon(X_s) = X_s(t)$ and $\forall t \in R$, $\widehat{\mathcal{E}}(t)$ is an isometry

Define $\forall i \in I : \varphi_i : R \rightarrow F :: \varphi_i(t) = \widehat{\mathcal{E}}(t) \tilde{\varepsilon}_i$

iv) $\widehat{\mathcal{E}}(t)$ can be uniquely extended in a continuous linear map :

$\widehat{\mathcal{E}}_N(t) : \otimes_N H \rightarrow \otimes_N F$ such that : $\widehat{\mathcal{E}}_N(t) (\otimes_N \psi_s) = \otimes_N X_s(t)$

$\widehat{\mathcal{E}}_N(t) (\otimes_{s=1}^N \tilde{\varepsilon}_{i_s}) = \otimes_{s=1}^N \varphi_{i_s}(t)$

$\widehat{\mathcal{E}}_N(t)$ is an isometry, so $\forall t \in R : \{\otimes_{s=1}^N \varphi_{i_s}(t), i_s \in I\}$ is a Hilbertian basis of $\otimes_N F$

v) Define as the state of the system at $t : S(t) = \widehat{\mathcal{E}}_N(t) (\Psi) \in \otimes_N F$

Define : $\forall \sigma \in \mathfrak{S}(N) : \widehat{U}_F(\sigma) \in \mathcal{L}(\otimes_N F; \otimes_N F)$ by linear extension of :

$\widehat{U}_F(\sigma) (\otimes_{s=1}^N \varphi_{i_s}(t)) = \otimes_{s=1}^N \varphi_{\sigma(i_s)}(t) = \widehat{\mathcal{E}}_N(t) \widehat{U}(\sigma) (\otimes_{s=1}^N \tilde{\varepsilon}_{i_s})$

$\forall \Psi \in \mathbf{h} : \Psi = \sum_{\sigma \in \mathfrak{S}(\lambda^c)} \Psi^\sigma \widehat{U}(\sigma) (\otimes_{n_1} \tilde{\varepsilon}_{j_1} \otimes_{n_2} \tilde{\varepsilon}_{j_2} \dots \otimes_{n_p} \tilde{\varepsilon}_{j_p})$

$S(t) = \sum_{\sigma \in \mathfrak{S}(\lambda^c)} \Psi^\sigma \widehat{\mathcal{E}}_N(t) \circ \widehat{U}(\sigma) (\otimes_{n_1} \tilde{\varepsilon}_{j_1} \otimes_{n_2} \tilde{\varepsilon}_{j_2} \dots \otimes_{n_p} \tilde{\varepsilon}_{j_p})$

$S(t) = \sum_{\sigma \in \mathfrak{S}(\lambda^c)} \Psi^\sigma \widehat{U}_F(\sigma) \otimes_{n_1} \varphi_{j_1}(t) \otimes_{n_2} \varphi_{j_2}(t) \dots \otimes_{n_p} \varphi_{j_p}(t)$

$\forall \theta \in \mathfrak{S}(\lambda) : \widehat{U}_F(\theta) (\otimes_{n_1} \varphi_{j_1}(t) \otimes_{n_2} \varphi_{j_2}(t) \dots \otimes_{n_p} \varphi_{j_p}(t))$

$= \otimes_{n_1} \varphi_{j_1}(t) \otimes_{n_2} \varphi_{j_2}(t) \dots \otimes_{n_p} \varphi_{j_p}(t)$

$\forall \theta \in \mathfrak{S}(\lambda^c) : \widehat{U}_F(\theta) (\otimes_{n_1} \varphi_{j_1}(t) \otimes_{n_2} \varphi_{j_2}(t) \dots \otimes_{n_p} \varphi_{j_p}(t))$

$\neq (\otimes_{n_1} \varphi_{j_1}(t) \otimes_{n_2} \varphi_{j_2}(t) \dots \otimes_{n_p} \varphi_{j_p}(t))$
and the tensors are linearly independent

So $\left\{ \widehat{U}_F(\sigma) (\otimes_{n_1} \varphi_{j_1}(t) \otimes_{n_2} \varphi_{j_2}(t) \dots \otimes_{n_p} \varphi_{j_p}(t)), \sigma \in \mathfrak{S}(\lambda^c) \right\}$ is an orthonormal basis of

$$\mathbf{f}(t) = \text{Span} \left\{ \widehat{U}_F(\sigma) (\otimes_{n_1} \varphi_{j_1}(t) \otimes_{n_2} \varphi_{j_2}(t) \dots \otimes_{n_p} \varphi_{j_p}(t)), \sigma \in \mathfrak{S}(\lambda^c) \right\}$$

$$\mathbf{f}(t) = \widehat{\mathcal{E}}_N(t)(\mathbf{h})$$

Let $\widetilde{\mathbf{f}}(t) \subset \mathbf{f}(t)$ be any subspace globally invariant by $\left\{ \widehat{U}_F(\theta), \theta \in \mathfrak{S}(N) \right\}$:

$$\widehat{U}_F(\theta) \widetilde{\mathbf{f}}(t) \in \widetilde{\mathbf{f}}(t)$$

$\widehat{\mathcal{E}}_N(t)$ is an isometry, thus a bijective map

$$\widetilde{\mathbf{h}} = \widehat{\mathcal{E}}_N(t)^{-1} \widetilde{\mathbf{f}}(t) \Leftrightarrow \widetilde{\mathbf{f}}(t) = \widehat{\mathcal{E}}_N(t) \widetilde{\mathbf{h}}$$

$$\widehat{U}_F(\theta) \widehat{\mathcal{E}}_N(t) \widetilde{\mathbf{h}} \in \widehat{\mathcal{E}}_N(t) \widetilde{\mathbf{h}}$$

$$\forall \Psi \in \mathbf{h} : \widehat{U}_F(\theta) \widehat{\mathcal{E}}_N(t) \Psi = \widehat{\mathcal{E}}_N(t) \widehat{U}_F(\theta) \Psi$$

$$\Rightarrow \widehat{\mathcal{E}}_N(t) \widehat{U}_F(\theta) \widetilde{\mathbf{h}} \in \widehat{\mathcal{E}}_N(t) \widetilde{\mathbf{h}}$$

$$\Rightarrow \widehat{U}_F(\theta) \widetilde{\mathbf{h}} \in \widetilde{\mathbf{h}}$$

So $(\mathbf{f}(t), \widehat{U}_F)$ is an irreducible representation of $\mathfrak{S}(N)$ ■

For each t the space $\mathbf{f}(t)$ is defined by a Hilbertian basis $(f_i)_{i \in I}$ of F , a finite subset J of I , a conjugacy class $\lambda(t)$ and a family of p vectors $(f_{j_k}(t))_{k=1}^p$ belonging to $(f_i)_{i \in J}$. The set J is arbitrary but defined by \mathbf{h} , so it does not depend on t . For a given class of conjugacy different families of vectors $(f_{j_k}(t))_{k=1}^p$ generate equivalent representations and isomorphic spaces, by symmetrization or antisymmetrization. So for a given system one can pick up a fixed ordered family $(f_j)_{j=1}^N$ of vectors in $(f_i)_{i \in I}$ such that for each class of conjugacy $\lambda = \{0 \leq n_p \leq \dots \leq n_1 \leq N, n_1 + \dots + n_p = N\}$ there is a unique vector space \mathbf{f}_λ defined by $\otimes_{n_1} f_1 \otimes_{n_2} f_2 \dots \otimes_{n_p} f_p$. Then if $S(t) \in \mathbf{f}_\lambda$:

$$S(t) = \sum_{\sigma \in \mathfrak{S}(\lambda^c)} S^\sigma(t) \widehat{U}_F(\sigma) (\otimes_{n_1} f_1 \otimes_{n_2} f_2 \dots \otimes_{n_p} f_p)$$

and at all time $S(t) \in \otimes_N F_J$.

The vector spaces \mathbf{f}_λ are orthogonal. With the orthogonal projection π_λ on \mathbf{f}_λ :

$$\forall t \in R : S(t) = \sum_\lambda \pi_\lambda S(t)$$

$$\|S(t)\|^2 = \sum_\lambda \|\pi_\lambda S(t)\|^2$$

The distance between $S(t)$ and a given \mathbf{f}_λ is well defined and :

$$\|S(t) - \pi_\lambda S(t)\|^2 = \|S(t)\|^2 - \|\pi_\lambda S(t)\|^2$$

Whenever S , and thus Θ , is continuous, the space \mathbf{f}_λ stays the same. As we have seen previously one can assume that, in all practical cases, Θ is continuous but for a countable set $\{t_k, k = 1, 2, \dots\}$ of isolated points. Then the different spaces \mathbf{f}_λ can be seen as phases, each of them associated with a class of conjugacy λ . And there are as many possible phases as classes of conjugacy. So, in a probabilist picture, one can assume that the probability for the system to be in a phase $\lambda : \Pr(S(t) \in \mathbf{f}_\lambda)$ is a function of $\frac{\|\pi_\lambda S(t)\|^2}{\|S(t)\|^2}$. It can be estimated as seen previously from data on a past period, with the knowledge of both λ and $\frac{\|\pi_\lambda S(t)\|^2}{\|S(t)\|^2}$.

2.7 CORRESPONDENCE WITH QUANTUM MECHANICS

It is useful to compare the results proven in the present paper to the axioms of QM as they are usually expressed.

2.7.1 Hilbert space

QM : 1. *The states of a physical system can be represented by rays in a complex Hilbert space H . Rays meaning that two vectors which differ by the product by a complex number of module 1 shall be considered as representing the same state.*

In Theorem 2 we have proven that in a model meeting precise conditions the states of the system can be represented as vectors in an infinite dimensional, separable, real Hilbert space. We have seen that it is always possible to endow the Hilbert space with a complex structure, but this is not a necessity. Moreover the Hilbert space is defined up to an isometry, so notably up to the product by a fixed complex scalar of module 1. We will see in the following how and why rays appear (this is specific to the representation of particles with electromagnetic fields).

In Quantum Physics a great attention is given to the Principle of Superposition. This Principle is equivalent to the condition that the variables of the system (and then its state) belong to a vector space. There is a distinction between pure states, which correspond to actual measures, and mixed states which are linear combination of pure states, usually not actually observed. There has been a great effort to give a physical meaning to these mixed states. Here the concept of pure states appears only in the tensors representing interacting systems, with the usual, but clear, explanation. In Quantum Mechanics some states of a system cannot be achieved (through a preparation for instance) as a combination of other states, and then super-selection rules are required to sort out these specific states. Here there is a simple explanation : because the set H_0 is not the whole of H it can happen that a linear combination of states is not inside H_0 . The remedy is to enlarge the model to account for other physical phenomena, if it appears that these states have a physical meaning.

Actually the main difference comes from the precise conditions of the Theorem 2. The variables must be maps, but also belong to a vector space. Thus for instance it does not apply to the model of a solid body represented by its trajectory $x(t)$ and its speed $v(t)$: the variable $x(t)$ is a map : $x : \mathbb{R} \rightarrow M$ valued in a manifold (an affine space in Galilean geometry). So it is necessary to adapt the model, using the fiber bundle formalism, and this leads to a deep redefinition of the concept of motion (including rotation) and to the spinors. And as it has been abundantly said, the state is defined by maps over the evolution of the system, and not pointwise.

2.7.2 Observables

QM : 2. *To any physical measure Φ , called an observable, which can be done on the system, is associated a continuous, linear, self-adjoint operator $\hat{\Phi}$ on H .*

We have proven that this operator is also compact and trace-class. The main result is that we have here a clear understanding of the concept of observable, rooted in the practical way the data are analyzed and assigned to the value of the variables, with the emphasize given to the procedure of specification, an essential step in any statistical analysis and which is usually overlooked. From primary observables it is possible to define von Neumann algebras of operators, which are necessarily commutative when a fixed basis has been chosen. As the choice of a privileged basis can always be done, one can say that there is always a commutative von Neumann algebra associated to a system. But, as it can be seen, these von Neumann algebras do not play any role in the proofs of the theorems. Their introduction can be useful, but they are not a keystone in our framework.

This is the opposite in the axiomatic interpretations of QM which define the system itself from the existence of a von Neumann algebra. However such interpretation is, eventually, based on the same assumption as any other interpretation of QM : the postulate that for any system there is a quantity, the state, which has a physical meaning and can be represented in a Hilbert space. Nothing preclude the choice of a privileged Hilbertian basis for this Hilbert space (as it is always done in any practical computation in QM), with respect to which the operators can be defined, and then the algebra is commutative. Which nullifies the emphasize given to the commutation of operators. Or at least it should be given another interpretation than the simultaneity of measures.

In QM a great emphasize is given to the commutation of observables, linked to the physical possibility to measure simultaneously two variables. This concept does not play any role here, for the strong reason that we consider maps with a domain over the whole extension, spatial and temporal, of the system, there is no assumption about how the measures are done, so the simultaneity of measures is not considered. In our picture the variables and their properties are the model, they are listed explicitly and it is assumed that there is some way to estimate their value, without any consideration of the time at which the measures are done. So the question of simultaneous measures does not arise, and the product of observables itself has no clear meaning and no use. If a variable is added, we have another model, the variable gets the same status as the others, and it is assumed that it can be measured.

2.7.3 Measure

QM : 3. *The result of any physical measure is one of the eigen-values λ of the associated operator $\hat{\Phi}$. After the measure the system is in the state represented by the corresponding eigen vector ψ_λ*

This is one of the most puzzling axiom. We have here a clear interpretation of this result, with primary observables, and there is always a primary observable which is at least as efficient than a secondary observable.

In our picture there is no assumption about how the measures are done, and particularly if they have or not an impact on the state of the system. If it is assumed that this is the case, a specific variable should be added to the model. Its value can be measured directly or estimated from the value of the other variables, but this does not make a difference : it is a variable as the others. We will see an example in the following chapters.

There is no assumption about the times at which the measures are taken, when the model represents a process the measures can be taken at the beginning, during the process, or at the end. The variables which are estimated are maps, and the estimation of maps requires more than one value of the arguments. The estimation is done by a statistical method which uses all the available data. From this point of view our picture is closer to what is done in the laboratories, than to the idealized vision of simultaneous measures, which should be taken all together at each time, and would be impossible because of the perturbation caused by the measure.

Actually the importance granted to the simultaneity of measures, magnified by Dirac, is somewhat strange. It is also problematic in the Relativist picture. It is clear that some measures cannot be done, at the atomic scale, without disturbing the state of the system that is studied, but this does not preclude to use the corresponding variables in a model, or give them a special status. Before the invention of radar the artillerymen used efficient models even if they were not able to measure the speed of their shells. And in a collider it is assumed that the speed and the location of particles are known when they collide.

2.7.4 Probability

QM : 4. *The probability that the measure is λ is equal to $|\langle \psi_\lambda, \psi \rangle|^2$ (with normalized eigen vectors). If a system is in a state represented by a normalized vector ψ , and an experiment is done to test whether it is in one of the states $(\psi_n)_{n=1}^N$ which constitutes an orthonormal set of vectors, then the probability of finding the system in the state ψ_n is $|\langle \psi_n, \psi \rangle|^2$.*

The first part is addressed by the theorem 17. The second part has no direct equivalent in our picture but can be interpreted as follows : a measure of the primary observable has shown that $\psi \in H_J$, then the probability that it belongs to $H_{J'}$ for any subset $J' \subset J$ is $\left\| \widehat{Y}_{J'}(\psi) \right\|^2$. It is a computation of conditional probabilities :

Proof. The probability that $\psi \in H_K$ for any subset $K \subset I$ is $\left\| \widehat{Y}_K(\psi) \right\|^2$. The probability that $\psi \in H_{J'}$ knowing that $\psi \in H_J$ is :

$$\Pr(\psi \in H_{J'} | \psi \in H_J) = \frac{\Pr(\psi \in H_{J'} \wedge \psi \in H_J)}{\Pr(\psi \in H_J)} = \frac{\Pr(\psi \in H_{J'})}{\Pr(\psi \in H_J)} = \frac{\|\widehat{Y}_{J'}(\psi)\|^2}{\|\widehat{Y}_J(\psi)\|^2} = \|\widehat{Y}_{J'}(\psi)\|^2 \text{ because } \widehat{Y}_{J'}(\psi) = \psi \text{ and } \|\psi\| = 1 \blacksquare$$

Moreover we have seen how the concept of wave functions can be introduced, and its meaning, for models where the variables are maps defined on the same set. Of course the possibility to define such a function does not imply that it is related to a physical phenomenon.

2.7.5 Interacting systems

QM : 5. *When two systems interact, the vectors representing the states belong to the tensorial product of the Hilbert states.*

This is the topic of the theorem 28. We have seen how it can be extended to N systems, and the consequences that entails for homogeneous systems. If the number of microsystems is not fixed, the formalism of Fock spaces can be used but would require a mathematical apparatus that is beyond the scope of this book.

There is a fierce debate about the issue of locality in physics, mainly related to the entanglement of states for interacting particles. It should be clear that the formal system that we have built is global : more so, it is its main asset. While most of the physical theories are local, with the tools which have been presented we can deal with variables which are global, and get some strong results without many assumptions regarding the local laws.

2.7.6 Wigner's theorem

QM : 6. *If the same state is represented by two rays R, R' , then there is an operator \widehat{U} , unitary or antiunitary, on the Hilbert space H such that if the state ψ is in the ray R then $\widehat{U}\psi$ is in the ray R' .*

This the topic of the theorem 21. The issue unitary / antiunitary exists in the usual presentation of QM because of the rays. In our picture the operator is necessarily unitary, which is actually usually the case.

2.7.7 Schrödinger equation

QM : 7. *The vector representing the state of a system which evolves with time follows the equation : $i\hbar \frac{\partial \psi}{\partial t} = \widehat{H}\psi$ where \widehat{H} is the Hamiltonian of the system.*

This is actually the topic of the theorem 27 and the result holds for the variables X in specific conditions, including in the General Relativity context. The imaginary i does not appear because the Hilbert space is real. As for Planck's constant of course it cannot appear in a formal model. However as said before all quantities must be dimensionless, as it is obvious in the equivalent expression $\psi(t) = \exp \frac{t}{i\hbar} \widehat{H}\psi(0)$. Thus it is necessary either to involve some constant, or

that all quantities (including the time t) are expressed in a universal system of units. This is commonly done by using the Planck's system of units. Which is more important is that the theorems (and notably the second) precise fairly strong conditions for their validity. In many cases the Schrödinger's equation, because of its linearity, seems "to good to be true". We can see why.

2.7.8 The scale issue

The results presented here hold whenever the model meets the conditions 1. So it is valid whatever the scale. But it is clear that the conditions are not met in many models used in classic physics, notably in Analytic Mechanics (the variables q are not vectorial quantities). Moreover actually in the other cases it can often be assumed that the variables belong themselves to Hilbert spaces. The results about observables and eigen values are then obvious, and those about the evolution of the system, for interacting systems or for gauge theories keep all their interest.

The "Quantic World", with its strange properties does not come from specific physical laws, which would appear below some scale, but from the physical properties of the atomic world themselves. And of course these cannot be addressed in the simple study of formal models.

So the results presented here, which are purely mathematical, give a consistent and satisfying explanation of the basic axioms of Quantum Mechanics, without the need for any exotic assumptions. They validate, and in many ways make simpler and safer, the use of techniques used for many years. Moreover, as it is easy to check, most of these results do not involve any physics at all : they hold for any scientific theory which is expressed in a mathematical formalism. From my point of view they bring a definitive answer to the issue of the interpretation of QM : the interpretations were sought in the physical world, but actually there is no such interpretation to be found. There is no physical interpretation because QM is not a physical theory.

The results presented go beyond the usual axioms of QM : on the conditions to detect an anomaly, on the quantization of a variable $Y = f(X)$, on the phases transitions. And other results can probably be found. So the method should give a fresh view of the foundations of QM in Physics.

Chapter 3

GEOMETRY

Almost all, if not all, measures rely on measures of lengths and times. These concepts are expressed in theories about the geometry of the universe, meaning of the container in which live the objects of physics. The issue here is not a model of the Universe, seen in its totality, which is the topic of Cosmology, but a model which tells us how to measure lengths and times, and how to compare measures done by different observers. Such a model is a prerequisite to any physical theory. Geometry, as a branch of Mathematics, is the product of this quest of a theory of the universe, and naturally a physical geometry is formalized with the tools of Mathematical Geometry. There are several Geometries used in Physics : Galilean Geometry, Special Relativity (SR) and General Relativity (GR).

In this first section we will see how such a geometry can be built, from simple observations. We will go directly to the General Relativity model. This is the one which is the most general and will be used in the rest of the book. It is said to be difficult, but actually these difficulties can be overcome with the right formalism. Moreover it forces us to leave usual representations, which are often deceptive.

3.1 MANIFOLD STRUCTURE

3.1.1 The Universe has the structure of a manifold

The first question is how do we measure a location ?

For the spatial location one can use a chart, meaning any procedure which relates a point to a system of coordinates. We need three scalar coordinates. There are many maps which are in use : on Earth the geographic coordinates with longitude, latitude and elevation, in Astronomy the celestial coordinates system, and in an experiment in the laboratory the position with respect to a trajectory. Of course a system of Cartesian coordinates, which measure distances with respect to a point (which itself must be located) provides a system of coordinates, even if it is rarely used.

For the temporal location one uses the coincidence with any agreed upon event. For millennia men used the position of celestial bodies for this purpose. Say "See you at Stonehenge at the spring's equinox" and you will be understood. Of course one can use a clock, but the purpose of a clock is to measure elapsed time, so one needs a clock and a starting point, which are agreed upon, to locate an event in time.

When necessary, one can use several charts to cover an area. The key point is that the charts are compatible : there are mathematical functions, transitions maps, which relate the coordinates of the same point in different charts.

A collection of compatible charts, each defined in an open subset of a vector space and valued in a given set M , is an atlas. A collection of compatible atlas over a set M defines the structure of a manifold. The charts define over M a topology, deduced from the vector space. The manifold is differentiable (resp. smooth) if the transition maps are differentiable (resp. smooth) (Maths.15.1.1).

This leads to :

Proposition 34 *The Universe can be represented as a four dimensional real manifold M*

In Galilean Geometry the manifold is the product of \mathbb{R} with a 3 dimensional affine space, and in SR this is a 4 dimensional affine space (affine spaces have a manifold structure).

We will limit ourselves to an area Ω of the universe, which can be large, where there is no singularity such as black hole, so that one can assume that one chart suffices. We will represent such a chart as a map :

$$\varphi_M : \mathbb{R}^4 \rightarrow \Omega :: \varphi_M (\xi^0, \xi^1, \xi^2, \xi^3) = m$$

which is assumed to be bijective and smooth, where $\xi = (\xi^0, \xi^1, \xi^2, \xi^3)$ are the coordinates of m in the chart φ_M .

We will assume that Ω is a relatively compact open in M , so that the manifold structure on M is the same as on Ω , and Ω is bounded.

A change of chart is represented by a bijective smooth map (the transition map) :

$$\chi : \mathbb{R}^4 \rightarrow \mathbb{R}^4 :: \eta^\alpha = \chi^\alpha (\xi^0, \xi^1, \xi^2, \xi^3)$$

such that the new map $\tilde{\varphi}_M$ and the initial map φ_M locate the same point :

$$\tilde{\varphi}_M(\chi^\alpha(\xi^0, \xi^1, \xi^2, \xi^3), \alpha = 0, \dots, 3) = \varphi_M(\xi^0, \xi^1, \xi^2, \xi^3)$$

Notice that there is no algebraic structure on $M : am + bm'$ has no meaning. This is illuminating in GR, but still holds in SR or Galilean Geometry. There is a clear distinction between coordinates, which are scalars depending on the choice of a chart, and the point they locate on the manifold (affine space or not).

3.1.2 The tangent vector space

At each point of a smooth manifold M one can define a set which has the structure of a vector space, with the same dimension as M . The best way to see it is to differentiate the map φ_M with respect to the coordinates (this is close to the mathematical construct). To any vector $u \in \mathbb{R}^4$ is associated the vector $u_m = \sum_{\alpha=0}^3 u^\alpha \partial_\alpha \varphi_M(\xi^0, \xi^1, \xi^2, \xi^3)$ which is denoted $u_m = \sum_{\alpha=0}^3 u^\alpha \partial \xi_\alpha$.

The basis $(\partial \xi_\alpha)_{\alpha=0}^3$ associated to a chart, called a **holonomic basis**, depends on the chart, but the vector space at m denoted $T_m M$ does not depend on the chart. With this vector space structure one can define a dual space $T_m M^*$ and holonomic dual bases denoted $d\xi^\alpha$ with : $d\xi^\alpha(\partial \xi_\beta) = \delta_\beta^\alpha$, and any other tensorial structure (see Math.16).

As one can see in the definition of the holonomic basis, the tangent space is generated by small displacements along one coordinate, around a point m . So, physically, locally the manifold is close to an affine space with a chosen origin m , and locally GR and SR look the same. This is similar to what we see on Earth : locally it looks flat.

So there are essential distinctions between :

- a point on the manifold, which is a geometric object (it does not depend of coordinates, even if it can be represented by coordinates in a chart) but has no vectorial structure attached (the linear combination of points has no meaning);
- a vector in the tangent bundle, which is also a geometric quantity (it exists independently of its measure by components in a basis) but has an algebraic structure : the linear combination of vectors is well defined.

Some physical properties of objects can be represented by vectors, other cannot, and the distinction comes from the fundamental assumptions of the theory. It is enshrined in the theory itself.

The vector spaces $T_m M$ depend on m , and there is no canonical (meaning independent of the choice of a specific tool) procedure to compare vectors belonging to the tangent spaces at two different points. These vectors u_m can be considered as a couple of a location m and a vector u , which can be defined in a holonomic basis or not, and all together they constitute the tangent bundle TM .

However because the manifolds are actually affine spaces, in SR and Galilean Geometry the tangent spaces at different points share the same structure (which is the underlying tangent vector space), and only in these cases they can be assimilated to \mathbb{R}^4 . This is the origin of much confusion on the subject, and the

motivation to start in the GR context where the concepts are clearly differentiated.

3.1.3 Vector fields

A vector field on M is a map : $V : M \rightarrow TM :: V(m) = \sum_{\alpha=0}^3 v^\alpha(m) \partial\xi_\alpha$ which associates to any point m a vector of the tangent space $T_m M$. The vector does not depend on the choice of a basis or a chart, so its components change in a change of chart as (Math.16.1.2) :

$$v^\alpha(m) \rightarrow \tilde{v}^\alpha(m) = \sum_{\beta=0}^3 [J(m)]_\beta^\alpha v^\beta(m)$$

where $[J(m)] = \left[\frac{\partial \eta^\alpha}{\partial \xi^\beta}(m) \right]$ is the 4x4 matrix called the jacobian

Similarly a one form on M is a map $\varpi : M \rightarrow TM^* :: \varpi(m) = \sum_{\alpha=0}^3 \varpi_\alpha(m) d\xi^\alpha$ and the components change as :

$$\varpi_\alpha(m) \rightarrow \tilde{\varpi}_\alpha(m) = \sum_{\beta=0}^3 [K(m)]_\alpha^\beta \varpi_\beta(m) \text{ and } [K(m)] = [J(m)]^{-1}$$

The sets of vector fields, denoted $\mathfrak{X}(TM)$, and of one forms, denoted $\mathfrak{X}(TM^*)$ or $\Lambda_1(M; \mathbb{R})$ are infinite dimensional vector spaces (with pointwise operations).

A **curve** on a manifold is a one dimensional submanifold : this is a geometric structure, and there is a vector space associated to each point of the curve, which is a one dimensional vector subspace of $T_m M$.

A **path** on a manifold is a map : $p : \mathbb{R} \rightarrow M :: m = p(\tau)$ where p is a differentiable map such that $p'(\tau) \neq 0$. Its image is a curve L_p , and p defines a bijection between \mathbb{R} (or any interval of \mathbb{R}) and the curve (this is a chart of the curve), the curve is a 1 dimensional submanifold embedded in M . The same curve can be defined by different paths. The tangent is the map : $p'(t) : \mathbb{R} \rightarrow T_{p(t)} M :: \frac{dp}{d\tau} \in T_{p(\tau)} L_p$. In a change of parameter in the path : $\tilde{\tau} = f(\tau)$ (which is a change of chart) for the same point : $m = \tilde{p}(\tilde{\tau}) = p(f(\tau))$ the new tangent is proportional to the previous one : $\frac{dm}{d\tau} = \frac{d\tilde{p}}{d\tilde{\tau}} \frac{d\tilde{\tau}}{d\tau} \Leftrightarrow \frac{dm}{d\tilde{\tau}} = \frac{1}{f'} \frac{dm}{d\tau}$

For any smooth vector field there is a collection of smooth paths (the **integrals** of the field) such that the tangent at any point of the curve is the vector field. There is a unique **integral line** which goes through a given point. The **flow** of a vector field V is the map (Math.14.3.5):

$\Phi_V : \mathbb{R} \times M \rightarrow M :: \Phi_V(\tau, a)$ such that $\Phi_V(\cdot, a) : \mathbb{R} \rightarrow M :: m = \Phi_V(\tau, a)$ is the integral path going through a :

$$\forall \theta \in \mathbb{R} : \frac{\partial}{\partial \tau} \Phi_V(\tau, a) |_{\tau=\theta} = V(\Phi_V(\theta, a)) \quad (3.1)$$

and $\Phi_V(\cdot, a)$ is a local diffeomorphism

$$\forall \tau, \tau' \in \mathbb{R} : \Phi_V(\tau + \tau', a) = \Phi_V(\tau, \Phi_V(\tau', a)) \quad (3.2)$$

$$\Phi_V(0, a) = a \quad (3.3)$$

$$\forall \tau \in \mathbb{R} : \Phi_V(-\tau, \Phi_V(\tau, a)) = a \quad (3.4)$$

Notice that, for a given vector field, the parameter τ is defined up to a constant, so *it is uniquely defined* with the condition $\Phi_V(0, a) = a$.

In general the flow is defined only for an interval of the parameter, but this restriction does not exist if Ω is relatively compact.

A map $f : C \rightarrow E$ from a curve to a Banach vector space E can be extended to a map $F : \Omega \rightarrow E$ (Maths.1467). So any smooth path can be considered as the integral of some vector field (not uniquely defined), and it is convenient to express a path as the flow of a vector field.

3.1.4 Fundamental symmetry breakdown

The idea that the Universe could be 4 dimensional is not new. R.Penrose remarked in his book “The road to reality” that Galileo considered this possibility. The true revolution of Relativity has been to acknowledge that, if the physical universe is 4 dimensional, it becomes necessary to dissociate the abstract representation of the world, the picture given by a mathematical model, from the actual representation of the world as it can be seen through measures. And this dissociation goes through the introduction of a new object in Physics : the observer. Indeed, if the physical Universe is 4 dimensional, the location of a point is absolute : there is a unique material body, in space and time, which can occupy a location. Then, does that mean that past and future exist together ? Can we say that this apple, which is falling, is somewhere in the Universe, still on the tree ? To avoid the conundrum and all the paradoxes that it entails, the solution is to acknowledge that, if there is a unique reality, actually the reality which is scientifically accessible, because it enables experiments and measures, is specific : it depends on the observer. This does not mean that it would be wrong to represent the reality in its entirety, as it can be done with charts, frames or other abstract mathematical objects. They are necessary to give a consistent picture, and more bluntly, to give a picture that is accessible to our mind. But we cannot identify this abstract representation, common to everybody, with the world as it is, because the one in which I can move, act and measure, is my world. This is one of the reasons that motivate the introduction of Geometry in this book through GR : it is common to introduce subtle concepts such as location and velocity through a frame, which is evoked in passing, as if it was obvious, standing somewhere at the disposition of the public. There is nothing like this. I can build my frame, my charts, and from there conceive that it can be extended, and compared to what other Physicists have done. But comparison requires first dissociation, and this is more easily done in a context to which we are less used to, by years of schematic representations.

The four coordinates are not equivalent : the measure of the time ξ^0 cannot be done with the same procedures as the other coordinates, and one cannot move along in time. This is the fundamental symmetry breakdown.

One assumes that *a given observer* can tell if two events A, B occur in his present time (they are simultaneous), and that the relation “two events are simultaneous” is a relation of equivalence between events. Then the observer

can label each class of equivalence of events by the time of his clock. Which can be expressed by telling that for each observer, there is a function : $f_o : M \rightarrow \mathbb{R} :: f_o(m) = t$ which assigns a time t , *with respect to the clock of the observer*, at any point of the universe (or at least Ω). The points : $\Omega(t) = \{m = f_o(t), m \in \Omega\}$ correspond to the **present** of the observer. No assumption is made about the clock, and different clocks can be used, with the condition that, as for any chart, it is possible to convert the time given by a clock to the time given by another clock (both used by the same observer).

In Galilean Geometry instantaneous communication is possible, so it is possible to define a universal time, to which any observer can refer to locate his position, and the present does not depend on the observer. The manifold M can be assimilated to the product $\mathbb{R} \times \mathbb{R}^3$. The usual representation of material bodies moving in the same affine space is a bit misleading, actually one should say that this affine space $\mathbb{R}^3(t)$ changes continuously, in the same way, for everybody. Told this way we see that Galilean Geometry relies on a huge assumption about the physical universe.

In Relativist Geometry instantaneous communication is impossible, so it is impossible to synchronize all the clocks. However a given observer can synchronize the clocks which correspond to his present, this is the meaning of the function f_o . Different procedures have been proposed for this purpose, the simplest uses electromagnetic signals which are bounced by the target. But this process is specific to each observer, and there is a priori no way to compare the time of clocks synchronized to two different observers, and located at two different spatial locations. We will see how this can be done.

Whenever there is, on a manifold, a map such that f_o , with $f'_o(m) \neq 0$, it defines on M a foliation : there is a collection of hypersurfaces (3 dimensional submanifolds) $\Omega_3(t)$, and the vectors u of the tangent spaces on $\Omega_3(t)$ are such that $f'_o(m)u = 0$, meanwhile the vectors which are transversal to $\Omega_3(t)$ (corresponding to paths which cross the hypersurface only once) are such that $f'_o(m)u > 0$ for any path with t increasing. So there are two faces on $\Omega_3(t)$: one for the incoming paths, and the other one for the outgoing paths. The hypersurfaces $\Omega_3(t)$ are diffeomorphic : they can be deduced from each other by a differentiable bijection, which is the flow of a vector field. Conversely if there is such a foliation one can define a unique function f_o with these properties (Maths.1502). The successions of present “spaces” for any observer is such a foliation, so our representation is consistent. And we state :

Proposition 35 *For any observer there is a function*

$$f_o : M \rightarrow \mathbb{R} :: f_o(m) = t \text{ with } f'_o(m) \neq 0 \quad (3.5)$$

which defines in any area Ω of the Universe a foliation by hypersurfaces

$$\Omega_3(t) = \{m = f_o(t), m \in \Omega\} \quad (3.6)$$

which represents the location of the events occurring at a given time t on his clock.

An observer can then define a chart of M , by taking the time on his clock, and the coordinates of a point x in the 3 dimensional hypersurfaces $\Omega_3(t)$. However we need a way to build consistently these spatial coordinates.

3.1.5 Trajectories of material bodies

The Universe is a container where physical objects live, and the manifold provides a way to measure a location. This is a 4 dimensional manifold which includes the time, but that does not mean that everything is frozen on the manifold : *the universe does not change, but its content changes*. As bodies move in the universe, their representation are paths on the manifold. And the fundamental symmetry breakdown gives a special meaning to the coordinate with respect to which the changes are measured. *Time is not only a parameter to locate an event, it is also a variable which defines the rates of change in the present of an observer*.

Material bodies and particles

The common definition of a material body in Physics is that of a set of material points which are related. A **material point** is assumed to have a location corresponding to a point of the manifold. According to the relations between material points of the same body we have rigid solids (the distance between two points is constant), deformable solids (the deformation tensor is locally given by the matrix of the transformation of a frame), fluids (the speed of material points are given by a vector field). These relations are formulated by phenomenological laws, they are essential in practical applications, but not in a theoretical study. So we will consider material bodies which have no internal structures, or whose internal structure can be neglected, that we will call **particles**. A particle then can be an electron, a nucleus, a molecule, or even a star system, according to the scale of the study. As in Mechanics a particle is a material point, and its location can be assimilated to a point from a geometrical point of view. We will see later how one can extend the concept of solid bodies to the relativist context.

World line and proper time

As required in any scientific theory a particle must be defined by its properties, and the first is that it occupies a precise location at any time. The successive locations of the material body define a curve and the particle travels on this curve according to a specific path called its world line. Any path can be defined by the flow of a vector such that the derivative with respect to the parameter is the tangent to the curve. The parameter called the proper time is then defined uniquely, up to the choice of an origin. The derivative with respect to the proper time is called the velocity. By definition *this is a vector*, defined at each point of the curve, and belonging to the tangent space to M . So the velocity has a definition which is independent of any basis.

Remark : For brevity I will call velocity the 4-vector, also usually called 4-velocity, and spatial speed the common 3 vector.

Observers are assumed to have similarly a world line and a proper time (they have other properties, notably they define a frame).

To sum up :

Definition 36 *Any particle or observer travels in the universe on a curve according to a specific path , $p : \mathbb{R} \rightarrow M :: m = p(\tau)$ called the **world line**, parametrized by the **proper time** τ , defined uniquely up to an origin. The derivative of the world line with respect to the proper time is a vector, the **velocity**, u . So that :*

$$u(\theta) = \left. \frac{dp}{d\tau} \right|_{\tau=\theta} \in T_{p(\theta)}M \quad (3.7)$$

$$p(\tau) = \Phi_u(\tau, a) \text{ with } a = \Phi_u(0, a) = p(0) \quad (3.8)$$

Observers are assumed to have clocks, that they use to measure their temporal location with respect to some starting point. The basic assumption is the following :

Proposition 37 *For any observer his proper time is the time on his clock.*

So the proper time of a particle can be seen formally as the time on the clock of an observer who would be attached to the particle.

We will strive to denote t the time of an observer (specific to an observer) and τ any other proper time. So for a given observer :

$$t = \tau$$

$$p_o : \mathbb{R} \rightarrow M :: m = p_o(t)$$

$$u(\theta) = \left. \frac{dp}{dt} \right|_{t=\theta} \in T_{p(\theta)}M$$

$$p_o(t) = \Phi_u(\tau, a) \text{ with } a = \Phi_u(0, a) = p(0)$$

The observer uses the time on his clock to locate temporally any event : this is the purpose of the function f_o and of the foliation $\Omega_3(t)$. The curve on which any particle travels meets only once each hypersurface $\Omega_3(t)$: it is seen only once. This happens at a time t :

$$f_o(p(\tau)) = t = f_o(\Phi_u(\tau, a))$$

So there is some relation between t and the proper time τ . It is specific, both to the observer and to the particle. It is bijective and both increases simultaneously, so that : $\frac{d\tau}{dt} > 0$.

The travel of the particle on the curve can be represented by the time of an observer. We will call then this path a **trajectory**.

A clock measures the elapsed time. It seems legitimate to assume that, in the procedure, one chooses clocks which run at the same rate. But, to do this, one needs some way to compare this rate. The most natural is to use a scalar measure of the velocity $\frac{d}{d\tau}p_o(\tau)$, and to assume that it is the same : material bodies would travel along their world lines at the same speed. But, as velocities are 4 dimensional vectors, one needs a special scalar product.

3.1.6 Metric on the manifold

Lorentz metric

A scalar product is defined by a bilinear symmetric form g acting on vectors of the tangent space, at each point of the manifold, thus by a tensor field called a **metric**. In a holonomic basis g reads :

$$g(m) = \sum_{\alpha\beta=0}^3 g_{\alpha\beta}(m) d\xi^\alpha \otimes d\xi^\beta \text{ with } g_{\alpha\beta} = g_{\beta\alpha} \quad (3.9)$$

The matrix of g is symmetric and invertible, if we assume that the scalar product is not degenerate. It is diagonalizable, and its eigen values are real. One wants to account for the symmetry breakdown, so these eigen values cannot have all the same sign (a direction is privileged). One knows that the hypersurface $\Omega_3(t)$ are Riemannian : there is a definite positive scalar product (acting on the 3 dimensional vector space tangent to $\Omega_3(t)$), and that transversal vectors correspond to the velocities of material bodies. So there are only two solutions for the signs of the eigen values of $[g(m)]$: either $(-,+,+,+)$ or $(+,-,-,-)$ which provides both a **Lorentz metric**. The scalar product, in an orthonormal basis $(\varepsilon_i)_{i=0}^3$ at m reads :

$$\text{signature } (3,1) : \langle u, v \rangle = u^1v^1 + u^2v^2 + u^3v^3 - u^0v^0 \quad (3.10)$$

$$\text{signature } (1,3) : \langle u, v \rangle = -u^1v^1 - u^2v^2 - u^3v^3 + u^0v^0 \quad (3.11)$$

Such a scalar product defines by restriction on each hypersurface $\Omega_3(t)$ a positive or a negative definite metric, which applies to spatial vectors (tangent to $\Omega_3(t)$) and provides, up to sign, the usual euclidean metric. So that both signatures are acceptable.

Which leads to :

Proposition 38 *The manifold M representing the Universe is endowed with a non degenerate metric, called the **Lorentz metric**, with signature either $(3,1)$ or $(1,3)$ defined at each point.*

This reasoning is a legitimate assumption, which is consistent with all the other concepts and assumptions, this is not the proof of the existence of such a metric. Such a proof comes from the formula in a change of frames between observers, which can be checked experimentally.

Notice that on a finite dimensional, connected, Hausdorff manifold, there is always a definite positive metric (Maths.1385). There is no relation between this metric and a Lorentz metric. Not all manifolds can have a Lorentz metric, the conditions are technical (see Giachetta p.224 for more) but one can safely assume that they are met in a limited region Ω .

Gauge group

The existence of a metric implies that, at any point, there are orthonormal bases $(\varepsilon_i)_{i=0}^3$ with the property :

Definition 39 $\langle \varepsilon_i, \varepsilon_j \rangle = \eta_{ij}$ for the signature $(3,1)$ and $\langle \varepsilon_i, \varepsilon_j \rangle = -\eta_{ij}$ for the signature $(1,3)$

with the matrix :

$$\text{Notation 40 } [\eta] = \begin{bmatrix} -1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \end{bmatrix}$$

so that the vector ε_0 is the time vector.

An orthonormal basis, at each point, is a **gauge**. The choice of an orthonormal basis depends on the observer : he has freedom of gauge. One goes from one gauge to another by a linear map χ which preserves the scalar product. They constitute a group, called the **gauge group**. In any basis these maps are represented by a matrix $[\chi]$ such that :

$$[\chi]^t [\eta] [\chi] = [\eta] \quad (3.12)$$

The group denoted equivalently $O(3,1)$ or $O(1,3)$, does not depend on the signature (replace $[\eta]$ by $-[\eta]$).

$O(3,1)$ is a 6 dimensional Lie group with Lie algebra $\mathfrak{o}(3,1)$ whose matrices $[h]$ are such that : $[h]^t [\eta] + [\eta] [h] = 0$. (Maths.24.5.3). The Lie algebra is a vector space and we will use the basis :

$$\begin{aligned} [\kappa_1] &= \begin{bmatrix} 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & -1 \\ 0 & 0 & 1 & 0 \end{bmatrix}; [\kappa_2] = \begin{bmatrix} 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 1 \\ 0 & 0 & 0 & 0 \\ 0 & -1 & 0 & 0 \end{bmatrix}; [\kappa_3] = \begin{bmatrix} 0 & 0 & 0 & 0 \\ 0 & 0 & -1 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{bmatrix} \\ [\kappa_4] &= \begin{bmatrix} 0 & 1 & 0 & 0 \\ 1 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{bmatrix}; [\kappa_5] = \begin{bmatrix} 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 \\ 1 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{bmatrix}; [\kappa_6] = \begin{bmatrix} 0 & 0 & 0 & 1 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 1 & 0 & 0 & 0 \end{bmatrix} \end{aligned}$$

so that any matrix of $\mathfrak{o}(3,1)$ can be written :

$[\kappa] = [J(r)] + [K(w)]$ with

$$[J(r)] = \begin{bmatrix} 0 & 0 & 0 & 0 \\ 0 & 0 & -r_3 & r_2 \\ 0 & r_3 & 0 & -r_1 \\ 0 & -r_2 & r_1 & 0 \end{bmatrix}; [K(w)] = \begin{bmatrix} 0 & w_1 & w_2 & w_3 \\ w_1 & 0 & 0 & 0 \\ w_2 & 0 & 0 & 0 \\ w_3 & 0 & 0 & 0 \end{bmatrix}$$

The exponential of these matrices read (Maths.493) :

$$\exp [K(w)] = I_4 + \frac{\sinh \sqrt{w^t w}}{\sqrt{w^t w}} K(w) + \frac{\cosh \sqrt{w^t w} - 1}{w^t w} K(w) K(w)$$

$$\exp [K (w)] = \begin{bmatrix} \cosh \sqrt{w^t w} & w^t \frac{\sinh \sqrt{w^t w}}{\sqrt{w^t w}} \\ w \frac{\sinh \sqrt{w^t w}}{\sqrt{w^t w}} & I_3 + \frac{\cosh \sqrt{w^t w} - 1}{w^t w} w w^t \end{bmatrix}$$

$$\exp [J (r)] = I_4 + \frac{\sin \sqrt{r^t r}}{\sqrt{r^t r}} J(r) + \frac{1 - \cos \sqrt{r^t r}}{r^t r} J(r) J(r) = \begin{bmatrix} 1 & 0 \\ 0 & R \end{bmatrix}$$

where R a 3×3 matrix of $O(3)$.

The group $O(3)$ has two connected components : the subgroup $SO(3)$ with determinant = 1, and the subset $O_1(3)$ with determinant -1.

$O(3, 1)$ has four connected components which can be distinguished according to the sign of the determinant and their projection under the compact subgroup $SO(3) \times \{I\}$.

Any matrix of $SO(3, 1)$ can be written as the product : $[\chi] = \exp [K (w)] \exp [J (r)]$ (or equivalently $\exp [J (r')] \exp [K (w')]$). So we have the 4 cases :

- $SO_0(3, 1)$: with determinant 1: $[\chi] = \exp K(w) \times \begin{bmatrix} 1 & 0 \\ 0 & R \end{bmatrix}$
- $SO_1(3, 1)$: with determinant 1: $[\chi] = \exp K(w) \times \begin{bmatrix} -1 & 0 \\ 0 & -R \end{bmatrix}$
- $SO_2(3, 1)$ with determinant = -1: $[\chi] = \exp K(w) \times \begin{bmatrix} -1 & 0 \\ 0 & R \end{bmatrix}$
- $SO_3(3, 1)$ with determinant = -1: $[\chi] = \exp K(w) \times \begin{bmatrix} 1 & 0 \\ 0 & -R \end{bmatrix}$

where R a 3×3 matrix of $SO(3)$, so that $-R \in O_1(3)$

$SO_k(3, 1)$, $k = 1, 2, 3$ are generated by the product of any element of $SO_0(3, 1)$ by either :

$$\text{the time reversal matrix : } T = \begin{bmatrix} -1 & 0 \\ 0 & I_3 \end{bmatrix}$$

$$\text{or the space reversal matrix : } S = \begin{bmatrix} 1 & 0 \\ 0 & -I_3 \end{bmatrix}$$

Orientation

Is the universe orientable ? In a universe where all observers have the same time, the simple existence of stereoisomers which do not have the same chemical properties suffices to answer positively. In a space-time universe one needs a process with an outcome which discriminates an orientation. All chemical reactions starting with a balanced mix of stereoisomers produce an equally balanced mix (stereoisomers have the same level of energy). However there are experiments involving the weak interactions which show the required property. So we can state that the 4 dimensional universe is orientable, and then we can distinguish orientation preserving gauge transformations. The right group to consider is $SO(3, 1)$.

The relativist universe is no longer isotropic : all directions are not equivalent. At any point m one can discriminate the vectors $v \in T_m M$ according to the value of the scalar product $\langle v, v \rangle$.

Definition 41 *Time like* vectors are vectors v such that $\langle v, v \rangle < 0$ with the signature $(3,1)$ and $\langle v, v \rangle > 0$ with the signature $(1,3)$

Space like vectors are vectors v such that $\langle v, v \rangle > 0$ with the signature $(3,1)$ and $\langle v, v \rangle < 0$ with the signature $(1,3)$

Moreover the subset of time like vectors has two disconnected components (this is no longer true in universes with more than one “time component” (Maths.307)). This has two important consequences.

A change of gauge, physically, implies some transport of the frame (one does not jump from one point to another) : we have a map $\chi : I \rightarrow SO(3,1)$ such that at each point of the path $p_o : I \rightarrow M$ defined on an interval I of \mathbb{R} , $\chi(t)$ is an isometry. The path which is followed matters. In particular it is connected. The frame $(\varepsilon_i)_{i=0}^3$ is transported by $\tilde{\varepsilon}_i(\tau) = \chi(t) \varepsilon_i(0)$. So $\{[\chi(\tau)], t \in I\}$, image of the connected interval I by a continuous map is a connected subset of $SO(3,1)$, and because $\chi(0) = Id$ it must be the component of the identity. So the right group to consider is the **connected component of the identity** $SO_0(3,1)$.

Because the subset of time like vectors has two disconnected components one can discriminate these components and, in accordance with the assumptions about the velocity of material bodies, it is logical to consider that their velocity is **future oriented**. And one can distinguish gauge transformations which preserve this time orientation.

Definition 42 *We will assume that the future orientation is given in a gauge by the vector ε_0 . So a vector u is time like and future oriented if :*

$$\begin{aligned} \langle u, u \rangle < 0, \langle u, \varepsilon_0 \rangle < 0 \text{ with the signature } (3,1) \\ \langle u, u \rangle > 0, \langle u, \varepsilon_0 \rangle > 0 \text{ with the signature } (1,3) \end{aligned}$$

A matrix $[\chi]$ of $SO_0(3,1)$ preserves the time orientation iff $[\chi]_0^0 > 0$ and this will always happen if $[\chi] = \exp[K(w)] \exp[J(r)]$ that is if $[\chi] \in SO_0(3,1)$.

A gauge transformation which preserves both the time orientation, and the global orientation must preserve also the spatial orientation.

3.1.7 Velocities have a constant Lorentz norm

The velocity $\frac{dp_o}{dt}$ is a vector which is defined independently of any basis, for any observer it is transversal to $\Omega_3(t)$. It is legitimate to say that it is future oriented, and so it must be time-like. One of the basic assumptions of Relativity is that it has a constant length, as measured by the metric, identical for all observers. So it is possible to use the norm of the velocity to define a standard rate at which the clocks run.

Because the proper time of any material body can be defined as the time on the clock of an observer attached to the body this proposition is extended to any particle.

The time is not measured with the same unit as the lengths, used for the spatial components of the velocity. The ratio ξ^i/t has the dimension of a spatial speed. So we make the general assumption that for any observer or particle the velocity is such that $\left\langle \frac{dp}{d\tau}, \frac{dp}{d\tau} \right\rangle = -c^2$ where τ is the proper time. Notice that c is a constant, with no specific value.

And we sum up :

Proposition 43 *The velocity $\frac{dp}{d\tau}$ of any particle or observer is a time like, future oriented vector with Lorentz norm*

$$\left\langle \frac{dp}{d\tau}, \frac{dp}{d\tau} \right\rangle = -c^2 \quad (3.13)$$

(with signature (3,1) or c^2 with signature (1,3)) where c is a fundamental constant.

3.1.8 Standard chart of an observer

As a consequence :

Theorem 44 *For any observer there is a vector field $\varepsilon_0 \in \mathfrak{X}(TM)$ which is future oriented, with length $\langle \varepsilon_0(m), \varepsilon_0(m) \rangle = -1$, normal to $\Omega_3(t)$ and such that : $\varepsilon_0(p_0(t)) = \frac{1}{c} \frac{dp_0}{dt}$ where $\frac{dp_0}{dt}$ is the velocity of the observer at each point of his world line.*

Proof. For an observer the function $f : \Omega \rightarrow \mathbb{R}$ has for derivative a one form $f'(m) \neq 0$ such that $\forall v \in T_m \Omega_3(t) : f'(m)v = 0$. Using the metric, it is possible to associate to $f'(m)$ a vector : $\varepsilon_0(m) = \text{grad}f : \langle \varepsilon_0(m), v \rangle = f'(m)v$. Thus $\varepsilon_0(m)$ is normal to $\Omega_3(t)$. Along the world line of the observer $\varepsilon_0(m)$ is in the direction of the velocity of the observer. And it is always possible to choose $\varepsilon_0(m)$ such that it is future oriented and with length $\langle \varepsilon_0(m), \varepsilon_0(m) \rangle = -1$ ■

As a consequence :

Theorem 45 *$\Omega_3(t)$ are space like hypersurfaces, with unitary normal $\varepsilon_0 \in \mathfrak{X}(TM)$*

Using the vector field ε_0 , and any any chart φ_Ω of $\Omega(0)$ there is a standard chart associated to an observer.

Definition 46 *The **standard chart** on M of any observer is defined as :*

$$\begin{aligned} \varphi_o : \mathbb{R}^4 &\rightarrow \Omega :: \varphi_o(\xi^0, \xi^1, \xi^2, \xi^3) = \Phi_{\varepsilon_0}(ct, x) \\ \xi^0 &= ct, \varphi_\Omega(\xi^1, \xi^2, \xi^3) = x \text{ in any chart of } \Omega(0) \end{aligned}$$

c is required in $\Phi_{\varepsilon_0}(ct, x)$ so that :

$$\xi^0 = ct \quad (3.14)$$

which makes all the coordinates homogeneous in units [Length].

The holonomic basis associated to this chart is such that :

$$\partial\xi_0 = \frac{\partial\varphi_o}{\partial\xi^0} = \frac{1}{c} \frac{\partial}{\partial t} \Phi_{\varepsilon_0}(ct, x) = \varepsilon_0$$

$$\varepsilon_0(m) = \partial\xi_0 \quad (3.15)$$

For any point $m = \varphi_o(\xi^0, \xi^1, \xi^2, \xi^3) \Phi_{\varepsilon_0}(ct, x)$ the point x is the point where the integral curve of ε_0 passing by m crosses $\Omega_3(0)$.

So the main characteristic of an observer can be summed in the vector field ε_0 (which is equivalently deduced from the function f_o).

3.1.9 Trajectory and speed of a particle

A particle follows a world line $q(\tau)$, parametrized by its proper time. Any observer sees only one instance of the particle, located at the point where the world line crosses the hypersurface $\Omega_3(t)$ so we have a relation between τ and t . This relation identifies the respective location of the observer and the particle on their own world lines. With the standard chart of the observer it is possible to measure the velocity of the particle at any location, and of course at the location where it belongs to $\Omega_3(t)$.

The trajectory (parametrized by t) of any particle in the standard chart of an observer is :

$$q(t) = \Phi_{\varepsilon_0}(ct, x(t)) = \varphi_o(ct, \xi^1(t), \xi^2(t), \xi^3(t))$$

By differentiation with respect to t :

$$\begin{aligned} \frac{dq}{dt} &= c\varepsilon_0(q(t)) + \frac{\partial}{\partial x} \Phi_{\varepsilon_0}(ct, x(t)) \frac{\partial x}{\partial t} \\ \frac{\partial}{\partial x} \Phi_{\varepsilon_0}(ct, x(t)) \frac{\partial x}{\partial t} &= \sum_{\alpha=1}^3 \frac{d\xi_\alpha}{dt} \partial\xi_\alpha \in T_m\Omega_3(t) \text{ so is orthogonal to } \varepsilon_0(q(t)) \end{aligned}$$

Definition 47 The *spatial speed* of a particle on its trajectory with respect to an observer is the vector of $T_{q(t)}\Omega_3(t)$:

$$\vec{v} = \frac{\partial}{\partial x} \Phi_{\varepsilon_0}(ct, x(t)) \frac{\partial x}{\partial t} = \sum_{\alpha=1}^3 \frac{d\xi_\alpha}{dt} \partial\xi_\alpha$$

Thus for any particle in the standard chart of an observer :

$$V(t) = \frac{dq}{dt} = c\varepsilon_0(q(t)) + \vec{v} \quad (3.16)$$

For the observer in the standard chart we had :

$$\frac{dp_0}{dt} = c\varepsilon_0(p_0(t)) \Leftrightarrow \vec{v} = 0$$

Notice that the velocity, and the spatial speed, are measured *in the chart of the observer at the point $q(t)$ where is the particle*. Because we have defined a standard chart it is possible to measure the speed of a particle located at a point $q(t)$ which is different from the location of the observer. And we can express the relation between τ and t .

Theorem 48 *The proper time τ of any particle and the corresponding time of any observer t are related by :*

$$\frac{d\tau}{dt} = \sqrt{1 - \frac{\|\vec{v}\|^2}{c^2}} \quad (3.17)$$

where \vec{v} is the spatial speed of the particle, with respect to the observer and measured in his standard chart.

The velocity of the particle is :

$$\frac{dp}{d\tau} = \frac{1}{\sqrt{1 - \frac{\|\vec{v}\|^2}{c^2}}} (\vec{v} + c\varepsilon_0(m)) \quad (3.18)$$

Proof. i) Let be a particle A with world line :

$$p : \mathbb{R} \rightarrow M :: m = p(\tau) = \Phi_u(\tau, a) \text{ with } a = \Phi_u(0, a) = p(0)$$

In the standard chart $\Phi_{\varepsilon_0}(ct, x)$ of the observer O its trajectory is :

$$q : \mathbb{R} \rightarrow M :: m = q(t) = \Phi_{\varepsilon_0}(ct, x(t))$$

So there is a relation between t, τ :

$$m = p(\tau) = \Phi_u(\tau, a) = q(t) = \Phi_{\varepsilon_0}(ct, x(t))$$

By differentiation with respect to t :

$$\frac{d}{dt} q(t) = c\varepsilon_0(p_A(t)) + \vec{v}$$

$$\frac{dq}{dt} = \vec{v} + c\varepsilon_0(m)$$

$$\frac{dq}{dt} = \frac{dp}{d\tau} \frac{d\tau}{dt}$$

$$\left\langle \frac{dp}{d\tau}, \frac{dp}{d\tau} \right\rangle = -c^2$$

$$\left\langle \frac{dq}{dt}, \frac{dq}{dt} \right\rangle = -c^2 \left(\frac{d\tau}{dt} \right)^2$$

$$\left\langle \frac{dq}{dt}, \frac{dq}{dt} \right\rangle = \langle \vec{v}, \vec{v} \rangle_3 - c^2 \text{ because } \varepsilon_0(m) \perp \Omega_3(t)$$

$$\|\vec{v}\|^2 - c^2 = -c^2 \left(\frac{d\tau}{dt} \right)^2$$

$$\text{and because } \frac{d\tau}{dt} > 0 : \frac{d\tau}{dt} = \sqrt{1 - \frac{\|\vec{v}\|^2}{c^2}}$$

ii) The velocity of the particle is :

$$\frac{dp}{d\tau} = \frac{dq}{dt} \frac{dt}{d\tau} = \frac{1}{\sqrt{1 - \frac{\|\vec{v}\|^2}{c^2}}} (\vec{v} + c\varepsilon_0(m)) \quad \blacksquare$$

As a consequence :

$$\|\vec{v}\| < c \quad (3.19)$$

$V(t) = \frac{dp}{dt}$ is the measure of the motion of the particle with respect to the observer : it can be seen as the relative velocity of the particle with respect to the observer. It involves \vec{v} which has the same meaning as usual, but we see that in Relativity one goes from the 4 velocity $u = \frac{dp}{d\tau}$ (which has an absolute meaning)

to the relative velocity $V(t) = \frac{dp}{dt} = \frac{dp}{d\tau} \frac{d\tau}{dt} = u \sqrt{1 - \frac{\|\vec{v}\|^2}{c^2}}$ by a scalar. If we have two particles A,B, with their path $q_A(\tau_A), q_B(\tau_B)$ can we define their relative

motion, for instance of B relative to A ? The simplest way to do it in relativity is to consider A as an observer, then $V_{B/A}(\tau_A) = \frac{dq_B}{d\tau_A} = u_B \sqrt{1 - \frac{\|\vec{v}_{B/A}\|^2}{c^2}}$ which is defined *in the chart* associated to the observer A.

3.2 FIBER BUNDLES

As said before, the location of a particle is absolute : this is the point in the physical Universe that it occupies at some time. But the measure of this location is relative to the observer, starting with the time at which the particle is at a given place. Similarly the velocity of a particle or an observer is absolute : in its definition there is no reference to a chart or a frame. This is an essential point in Relativity. State that the velocity of a particle is absolute confers to the variable a specific status : it is a geometric vector. In the remarks following the Theorem 2 in the QM Chapter, we noticed that the status - vector or not - of a variable is not arbitrary : it is part of the assumptions of the theory. Velocity is an intrinsic property of material bodies and particles, the measure of this velocity depends on the observer : it is relative.

This remark extends to all measures. A physical measure in itself has no meaning if one does not know how it has been done. The label “done by the observer O” is necessary. So we cannot contend ourselves with maps $X : M \rightarrow E$. We need a way to attach a tag, identifying the way the measure has been done, to the value of the variable. The mathematical tool to achieve that is the fiber bundle formalism. This is more than a sophisticated mathematical theory, it embodies the relation between measure (the value) and conditions of the measure (the gauge).

3.2.1 Fiber bundles theory

(see Math.Part VI)

General fiber bundle

A fiber bundle, denoted $P(M, F, \pi_P)$, is a manifold P, which is locally the product of two manifolds, the base M and the standard fiber F, with a projection : $\pi_P : P \rightarrow M$ which is a surjective submersion. The subset of P : $\pi_P^{-1}(m)$ is the fiber over m . It is usually defined over a collection of open subsets of M, patched together, but we will assume that on the area Ω there is only one component (the fiber bundles are assumed to be trivial). A **trivialization** is a map :

$$\varphi_P : M \times F \rightarrow P :: p = \varphi_P(m, v)$$

and any element of P is projected on M : $\forall v \in F : \pi_P(\varphi_P(m, v)) = m$. So it is similar to a chart, but the arguments are points of the manifolds.

A **section p** on P is defined by a map : $v : M \rightarrow F$ and $\mathbf{p} = \varphi_P(m, v(m))$. The set of sections is denoted $\mathfrak{X}(P)$.

A fiber bundle can be defined by different trivializations. In a **change of trivialization** the *same* element p is defined by a different map φ_P : this is very similar to the charts for manifold.

$$p = \varphi_P(m, v) = \tilde{\varphi}_P(m, \tilde{v})$$

and there is a necessary relation between v and \tilde{v} (m stays always the same) depending on the kind of fiber bundle.

Principal bundle

If $F = G$ is a Lie group then P is a **principal bundle** : its elements are a value $g(m)$ of G localized at a point m . There is a right action of G on P : $p = \varphi_P(m, g) \rightarrow p \cdot g' = \varphi_P(m, g \cdot g')$

p will usually define the basis used to measure vectors, so p is commonly called a gauge. There is a special gauge which can be defined at any point (it will usually be the gauge of the observer) : the **standard gauge**, the element of the fiber bundle such that : $\mathbf{p}(m) = \varphi_P(m, 1)$.

A change of trivialization is induced by a map : $\chi : M \rightarrow G$ that is by a section $\chi \in \mathfrak{X}(P)$ and :

$p = \varphi_P(m, g) = \tilde{\varphi}_P(m, \chi(m) \cdot g) = \tilde{\varphi}_P(m, \tilde{g}) \Leftrightarrow \tilde{g} = \chi(m) \cdot g$ ($\chi(m)$ acts on the left)

$\chi(m)$ can be identical over M (the change is said to be global) or depends on m (the change is local).

The expression of the elements of a section change as :

$$\sigma \in \mathfrak{X}(P) :: \sigma = \varphi_P(m, \sigma(m)) = \tilde{\varphi}_P(m, \tilde{\sigma}(m)) \Leftrightarrow \tilde{\sigma}(m) = \chi(m) \cdot \sigma(m)$$

$$\sigma(m) = \varphi_P(m, \sigma(m)) = \tilde{\varphi}_P(m, \chi(m) \cdot \sigma(m)) \quad (3.20)$$

A change of trivialization induces a **change of standard gauge** :

$$\mathbf{p}(m) = \varphi_P(m, 1) = \tilde{\varphi}_P(m, \chi(m))$$

$$\rightarrow \tilde{\mathbf{p}}(m) = \tilde{\varphi}_P(m, 1) = \tilde{\varphi}_P\left(m, \chi(m) \cdot \chi(m)^{-1}\right) = \mathbf{p}(m) \cdot \chi(m)^{-1}$$

$$\mathbf{p}(m) = \varphi_P(m, 1) \rightarrow \tilde{\mathbf{p}}(m) = \mathbf{p}(m) \cdot \chi(m)^{-1} : \quad (3.21)$$

$$\sigma(m) = \varphi_P(m, \sigma(m)) = \tilde{\varphi}_P(m, \chi(m) \cdot \sigma(m)) \quad (3.22)$$

Vector bundle

If $F=V$ is a vector space then P is a **vector bundle** and it has at each point the structure of a vector space :

$$w_m = \varphi_P(m, w), w'_m = \varphi_P(m, w'), \alpha, \beta \in \mathbb{R} :$$

$$\alpha w_m + \beta w'_m = \varphi_P(m, \alpha w + \beta w')$$

A holonomic basis is defined by a basis $(\varepsilon_i)_{i \in I}$ of V and : $\varepsilon_i(m) = \varphi_P(m, \varepsilon_i)$.

Usually vector bundles are defined as associated vector bundles. The principal bundle defines locally a standard with respect to which the measure is done. The result belong to a fixed set, but its value is labeled by the standard which is used and related to a point of a manifold.

Associated fiber bundle

Whenever there is a manifold F , a left action λ of G on F , one can built an **associated fiber bundle** denoted $P[F, \lambda]$ comprised of couples :

$$(p, v) \in P \times F \text{ with the equivalence relation : } (p, v) \sim (p \cdot g, \lambda(g^{-1}, v))$$

It is convenient to define these couples by using the standard gauge on P :

$$(\mathbf{p}(m), v) = (\varphi_P(m, 1), v) \sim (\varphi_P(m, g), \lambda(g^{-1}, v)) \quad (3.23)$$

A standard gauge is nothing more than the use of an arbitrary standard, represented by 1, with respect to which the measure is done. This is not a section : the standard gauge is the embodiment of the free will of the observer, who can choose the way he proceeds to the measure, and is not fixed by any physical law. A change of standard gauge $\mathbf{p}(m) = \varphi_P(m, 1) \rightarrow \tilde{\mathbf{p}}(m) = \mathbf{p}(m) \cdot \chi(m)^{-1}$ in the principal bundle impacts all associated fiber bundles (this is similar to the change of units) :

$$v_p = (\mathbf{p}(m), v) = (\mathbf{p}(m) \cdot \chi(m)^{-1}, \lambda(\chi(m), v)) \quad (3.24)$$

Similarly for the components of a section :

$$\mathbf{v} \in \mathfrak{X}(P[V, \lambda]) :: \mathbf{v}(m) = (\mathbf{p}(m), v(m)) = (\mathbf{p}(m) \cdot \chi(m)^{-1}, \lambda(\chi(m), v))$$

If F is a vector space V and $[V, \rho]$ a representation of the group G then we have an **associated vector bundle** $P[V, \rho]$ which has locally the structure of a vector space. It is convenient to define a **holonomic basis** $(\varepsilon_i(m))_{i=1}^n$ from a basis $(\varepsilon_i)_{i=1}^n$ of V by : $\varepsilon_i(m) = (\mathbf{p}(m), \varepsilon_i)$ then any vector of $P[V, \rho]$ reads :

$$v_m = (\mathbf{p}(m), v) = \left(\mathbf{p}(m), \sum_{i=1}^n v^i \varepsilon_i \right) = \sum_{i=1}^n v^i \varepsilon_i(m) \quad (3.25)$$

A change of standard gauge $\mathbf{p}(m) = \varphi_P(m, 1) \rightarrow \tilde{\mathbf{p}}(m) = \mathbf{p}(m) \cdot \chi(m)^{-1}$ in the principal bundle impacts all associated vector bundles.

The holonomic basis of a vector bundle changes as :

$$\begin{aligned} \varepsilon_i(m) &= (\mathbf{p}(m), \varepsilon_i) \rightarrow \\ \tilde{\varepsilon}_i(m) &= (\tilde{\mathbf{p}}(m), \varepsilon_i) = (\mathbf{p}(m) \cdot \chi(m)^{-1}, \varepsilon_i) \\ &\sim \left((\mathbf{p}(m) \cdot \chi(m)^{-1}) \cdot \chi(m), \rho(\chi(m)^{-1}) \varepsilon_i \right) \\ &= (\mathbf{p}(m), \rho(\chi(m)^{-1}) \varepsilon_i) = \rho(\chi(m)^{-1}) \varepsilon_i(m) \end{aligned}$$

$$\mathbf{p}(m) = \varphi_P(m, 1) \rightarrow \tilde{\mathbf{p}}(m) = \mathbf{p}(m) \cdot \chi(m)^{-1} : \quad (3.26)$$

$$\varepsilon_i(m) = (\mathbf{p}(m), \varepsilon_i) \rightarrow \tilde{\varepsilon}_i(m) = \rho(\chi(m))^{-1} \varepsilon_i(m) \quad (3.27)$$

so that the components of a vector in the holonomic basis change as :

$$\begin{aligned} v_m &= \sum_{i=1}^n v^i \varepsilon_i(m) = \sum_{i=1}^n \tilde{v}^i \tilde{\varepsilon}_i(m) = \sum_{i=1}^n \tilde{v}^i \rho(\chi(m))^{-1} \varepsilon_i(m) \\ \Rightarrow \tilde{v}^i &= \sum_j [\rho(\chi(m))]_j^i v^j \end{aligned}$$

$$\mathbf{p}(m) = \varphi_P(m, 1) \rightarrow \tilde{\mathbf{p}}(m) = \mathbf{p}(m) \cdot \chi(m)^{-1} : \quad (3.28)$$

$$v^i \rightarrow \tilde{v}^i = \sum_j [\rho(\chi(m))]_j^i v^j \quad (3.29)$$

The set of sections on $P[V, \rho]$, denoted $\mathfrak{X}(P[V, \rho])$, is an infinite dimensional vector space. In a change of standard gauge the components of a section change as :

$$\begin{aligned} v \in \mathfrak{X}(P[V, \rho]) &:: v(m) = \sum_{i=1}^n v^i(m) \varepsilon_i(m) = \sum_{i=1}^n \tilde{v}^i(m) \tilde{\varepsilon}_i(m) \\ \Leftrightarrow \tilde{v}^i(m) &= \sum_j \left[\rho \left(\chi(m)^{-1} \right) \right]_j^i v^j(m) \end{aligned}$$

so that $(\mathfrak{X}(P[V, \rho]), \rho)$ is an infinite dimensional representations of the group G .

I have given with great precision the rules in a change of gauge, as they will be used quite often (and are a source of constant mistakes !). They are necessary to ensure that a quantity is intrinsic : if it is geometric, its measure must change according to the rules. And conversely if it changes according to the rules, then one can say that it is intrinsic (this is similar to tensors). A quantity which is a vector of a fiber bundle is geometric with regard the conditions 1 of the 2nd chapter. Because this is a source of confusion, I will try to stick to these precise terms :

- a section = a point of a fiber bundle whose value is defined for each $m \in M$, this is a geometric object

- a gauge = a point of the principal bundle of P , this is a geometric object, which does not depend on a trivialization

- a standard gauge = a specific element of P , whose definition depends of the trivialization. This is not a section.

- a change of trivialization does not change the points of P , the gauge or the sections, but change the standard gauge and the way the points of P are defined with respect to the standard gauge

- it is equivalent to define a change of trivialization by the change of maps $\varphi_P \rightarrow \tilde{\varphi}_P$ or by the change of standard gauge : $\mathbf{p}(m) = \varphi_P(m, 1) \rightarrow \tilde{\mathbf{p}}(m) = \tilde{\varphi}_P(m, 1)$

Notice that the elements of a section stay the same, but their definition changes, meanwhile the holonomic bases are defined by different elements. This is very similar to what we have in any vector space in a change of basis : the vectors of the basis change, the other vectors stay the same, but their components change.

Scalar product and norm

Whenever there is a scalar product (bilinear symmetric or Hermitian two form) $\langle \rangle$ on a vector space V , so that (V, ρ) is a *unitary* representation of the group G : $\langle \rho(g)v, \rho(g)v' \rangle = \langle v, v' \rangle$, the scalar product can be extended on the associated vector bundle $P[V, \rho]$:

$$\langle (\mathbf{p}(m), v), (\mathbf{p}(m), v') \rangle_{P[V, \rho]} = \langle v, v' \rangle_W \quad (3.30)$$

If this scalar product is definite positive, with any measure μ on the manifold M (usually the Lebesgue measure associated to a volume form as in the relativist context), one can define the spaces of integrable sections $L^q(M, \mu, P[V, \rho])$ of

$P[V, \rho]$ (by taking the integral of the norm pointwise). For $q = 2$ they are Hilbert spaces, and unitary representation of the group G . Notice that the signature of the scalar product is that of the product defined on $P[V, \rho]$, *the metric on M is not involved*.

If there is a **norm** on V , that is a map :

$$\|\cdot\| : V \rightarrow \mathbb{R}_+$$

such that :

$$\|X\| \geq 0$$

$$\|X\| = 0 \Leftrightarrow X = 0$$

$$\|kX\| = |k| \|X\|$$

$$\|X + X'\| \leq \|X\| + \|X'\|$$

which does not depend on ρ :

$$\forall g \in G : \|\rho(g) X\| = \|X\|$$

then one can define a norm pointwise on $P[V, \rho]$:

$$\|(p(m), v)\| = \|v\|$$

$$(p(m), v) \sim (p(m) \rho(g^{-1}), \rho(g)v)$$

$$\|(p(m) \rho(g^{-1}), \rho(g)v)\| = \|\rho(g)v\| = \|v\|$$

and the space of integrable maps :

$$L^1(\mathfrak{X}(P[V, \rho])) = \{X \in \mathfrak{X}(P[V, \rho]), \int_{\Omega} \|X\| \mu < \infty\}$$

is a separable Fréchet space if Ω is a compact subset.

We have several fiber bundles in the Geometry of the Universe that we have defined. The simplest is the usual tangent bundle TM over M , which is a vector bundle associated to the choice of an invertible map at each point (the gauge group is $SL(\mathbb{R}, 4)$). But we have another one through the standard chart of an observer ;

Definition 49 *For any observer there is a **fiber bundle structure** $M_o(\mathbb{R}, \Omega(0), \pi_o)$ on M with base \mathbb{R} and :*

$$\text{projection} : \pi_o(m) = f_o(m)$$

$$\text{trivialization} : \Phi_{\varepsilon_0} : \mathbb{R} \times \Omega(0) \rightarrow \Omega :: \Phi_{\varepsilon_0}(ct, x) = m$$

3.2.2 Standard gauges associated to an observer

Following the Principle of Locality any physical map, used to measure the *components of a vector* at a point m in M , must be done at m , that is in a local frame. One property of the observers is that they have freedom of gauge : along their travel on their world line $p_o(t)$ they can choose a gauge, by choosing 4 orthogonal vectors.

For the time vector the observer has actually no choice : this is necessarily the vector ε_0 in the direction of his velocity $\frac{dp_o}{dt}$. And this vector has been extended as the vector field ε_0 orthogonal to $\Omega_3(t)$. The 3 other vectors of an orthonormal basis, corresponding to the space, belong to the space tangent at $\Omega_3(t)$, they are orthogonal to ε_0 and we assume that the observer can define these 3 vectors at any point of his present $\Omega_3(t)$. This can be achieved by a

system of communication (not instantaneous) between observers who would be located at each point.

This is equivalent to assume that, for each observer, there is a principal bundle $P_o(M, SO_0(3, 1), \pi_p)$, a gauge $\mathbf{p}(m) = \varphi_P(m, 1)$ and an associated vector bundle $P_o[\mathbb{R}^4, \iota]$ where (\mathbb{R}^4, ι) is the standard representation of $SO_0(3, 1)$. It defines at each point an holonomic orthonormal basis : $\varepsilon_i(m) = (\mathbf{p}(m), \varepsilon_i)$. To sum up :

Proposition 50 For each observer there is :

a principal **fiber bundle structure** $\mathbf{P}_o(M, SO_0(3, 1), \pi_p)$ on M with fiber the connected component of unity $SO_0(3, 1)$, which defines at each point a **standard gauge** : $\mathbf{p}(m) = \varphi_P(m, 1)$

an **associated vector bundle structure** $P_o[\mathbb{R}^4, \iota]$ where (\mathbb{R}^4, ι) is the standard representation of $SO_0(3, 1)$, which defines at any point $m \in \Omega$ the **standard basis** $\varepsilon_i(m) = (\mathbf{p}(m), \varepsilon_i), i=0..3$ where $\varepsilon_0(m)$ is orthogonal to the hypersurfaces $\Omega_3(t)$ to which m belongs.

Notice that these structures depend on the observer. Starting with the principal bundle P_o , a change of gauge can be defined at any point by a section $\chi \in \mathfrak{X}(P_o)$ as seen above, with an impact on any associated bundle.

A standard basis is such that its time vector is $\varepsilon_0(m)$, so at the location of the observer it is in the direction of his velocity. Standard bases are not unique : their time vector is the same, but their space vectors can be rotated in $\Omega_3(t)$. Because they constitute an euclidean orthonormal basis the rotation is given by a matrix of $SO(3)$.

3.2.3 Formulas for a change of observer

Theorem 51 For any two observers O, A the components of the vectors of the standard orthonormal basis of A , expressed in the standard basis of O are expressed by the matrix of $SO_0(3, 1)$, where \vec{v} is the instantaneous spatial speed of A with respect to O and R a matrix of $SO(3)$:

$$[\chi] = \begin{bmatrix} \frac{1}{\sqrt{1 - \frac{\|v\|^2}{c^2}}} & \frac{\frac{v^t}{c}}{\sqrt{1 - \frac{\|v\|^2}{c^2}}} \\ \frac{\frac{v}{c}}{\sqrt{1 - \frac{\|v\|^2}{c^2}}} & \left(\frac{1}{\sqrt{1 - \frac{\|v\|^2}{c^2}}} - 1 \right) \frac{vv^t}{\|v\|^2} \end{bmatrix} \begin{bmatrix} 1 & 0 \\ 0 & R \end{bmatrix} \quad (3.31)$$

Proof. Let be :

O be an observer (this will be main observer) with associated vector field ε_0 , proper time t and world line $p_o(t)$

A be another observer with associated vector field ε'_0 , proper time τ

Both observers use their standard chart φ_o, φ_A and their standard orthonormal basis, whose time vector is in the direction of their velocity. The location of A on his world line is the point m such that A belongs to the hypersurface $\Omega_3(t)$

The velocity of A at m :

$\frac{dp_A}{d\tau} = c\varepsilon'_0(m)$ by definition of the standard basis of A

$\frac{dp_A}{d\tau} = \frac{1}{\sqrt{1 - \frac{\|\vec{v}\|^2}{c^2}}} (\vec{v} + c\varepsilon_0(m))$ as measured in the standard basis of O

The matrix $[\chi]$ to go from the orthonormal basis $(\varepsilon_i(m))_{i=0}^3$ to $(\varepsilon'_i(m))_{i=0}^3$ belongs to $SO_0(3,1)$. It reads :

$$[\chi(t)] = \begin{bmatrix} \cosh \sqrt{w^t w} & w^t \frac{\sinh \sqrt{w^t w}}{\sqrt{w^t w}} \\ w \frac{\sinh \sqrt{w^t w}}{\sqrt{w^t w}} & I_3 + \frac{\cosh \sqrt{w^t w} - 1}{w^t w} w w^t \end{bmatrix} \begin{bmatrix} 1 & 0 \\ 0 & R \end{bmatrix}$$

for some $w \in \mathbb{R}^3, R \in SO(3)$

The elements of the first column of $[\chi(t)]$ are the components of $\varepsilon'_0(m)$, that is of $\frac{1}{c} \frac{dp_A}{d\tau}$ expressed in the basis of O :

$$\cosh \sqrt{w^t w} = \frac{1}{\sqrt{1 - \frac{\|\vec{v}\|^2}{c^2}}}$$

$$w \frac{\sinh \sqrt{w^t w}}{\sqrt{w^t w}} = \frac{\vec{v}}{c} \frac{1}{\sqrt{1 - \frac{\|\vec{v}\|^2}{c^2}}}$$

$$w = k \vec{v} \Rightarrow w^t w = k^2 \|\vec{v}\|^2$$

which leads to the classic formula with

$$w = \frac{v}{\|\vec{v}\|} \arg \tanh \left\| \frac{v}{c} \right\| = \frac{1}{2} \frac{v}{\|\vec{v}\|} \ln \left(\frac{c + \|\vec{v}\|}{c - \|\vec{v}\|} \right) \sim \frac{1}{2} \frac{v}{\|\vec{v}\|} \ln \left(1 + 2 \frac{\|\vec{v}\|}{c} \right) \sim \frac{v}{c} \blacksquare$$

Some key points to understand these formulas :

- They hold for any observers O, A, who use their standard orthonormal basis (the time vector is oriented in the direction of their velocity). There is no condition such as inertial frames.

- The points of M where O and A are located can be different, but they belong both to the same hypersurface $\Omega_3(t)$. The spatial speed \vec{v} is a vector belonging to the space tangent at $\Omega_3(t)$ at the location m of A (and not at the location of O at t).

- The formulas are related to the standard orthonormal bases $(\varepsilon_i(m))_{i=0}^3$ of O and $(\varepsilon'_i(m))_{i=0}^3$ located at the point m of $\Omega_3(t)$ where A is located.

- These formulas apply to the components of vectors in the standard orthonormal bases.

These formulas have been verified with a great accuracy, and the experiments show that c is the speed of light. This is an example of a theory which is checked by the consequences that can be drawn from its basic assumptions.

We will see below how these formulas apply in Special Relativity.

If we take $\frac{v}{c} \rightarrow 0$ we get $[\chi] = \begin{bmatrix} 1 & 0 \\ 0 & R \end{bmatrix}$, that is a rotation of the usual space.

The Galilean Geometry is an approximation of SR when the speeds are small with respect to c . Then the velocities are $\frac{d\mu_A}{d\tau} = (\vec{v} + c\varepsilon_0)$ with a common vector ε_0 .

3.2.4 The Tetrad

The principal fiber bundle P_G

So far we have defined a chart φ_o and a fiber bundle structure P_o for an observer : the construct is based on the motion of the observer, and his capability to extend his frame over the hypersurfaces $\Omega_3(t)$. With the formulas above we see how one can go from one observer to another, and thus relate the different fiber bundles P_o . The computations in a change of frame can be done with measures done by the observers, and have been checked experimentally. So it is legitimate to assume that there is a more general structure of principal bundle, denoted $\mathbf{P}_G(M, SO_0(3,1), \pi_G)$, over M . In this representation the bases used by any observer is just a choice of specific trivialization, or equivalently of standard gauge, and one goes from one trivialization to another with the matrix $[\chi]$.

If it is easy to define such a mathematical structure, it is necessary to understand its physical meaning. There is no way to relate an element $p \in P_G$ at a point m of M to a physical phenomenon : a trivialization is, like a chart, arbitrary, but the relations in a change of trivialization can be measured by their consequences on the bases of associated bundle (as can be checked the relations between coordinates in a change of chart). In particular a gauge, denoted $\varphi_G(m, 1)$ is arbitrary, but it will be convenient to keep the notation in this book, and to precise its meaning : we state that the standard gauge $\mathbf{p}(m) = \varphi_G(m, 1)$ is the gauge such that the associated basis $\varepsilon_i(m) = (\mathbf{p}(m), \varepsilon_i)$ is the standard basis chosen by an observer. So the standard gauge is actually defined by the standard basis. Such a basis, for any observer, is orthonormal and its time vector is time like, future oriented, and defined by the velocity of the observer. Any observer can choose another orthonormal basis, which is deduced by an action of $SO_0(3,1)$, and thus by an element of P_G . So a change of trivialization, or standard gauge, is equivalent to a change of standard basis. If the change of basis involves a change of the time vector $\varepsilon_0(m)$ we assume that this is a change of observer.

Proposition 52 *There is a unique structure of principal bundle*

$\mathbf{P}_G(M, SO_0(3,1), \pi_G)$ with base M , standard fiber $SO_0(3,1)$. A change of observer is given by a change of trivialization on P_G .

The standard gauge $\mathbf{p}(m) = \varphi_G(m, 1)$ is, for any observer, associated to his standard basis $\varepsilon_i(m) = (\mathbf{p}(m), \varepsilon_i)$.

As a consequence the change of observer is a change of gauge, given by a section χ (global or not) of \mathbf{P}_G , the vectors of the standard basis transform according to the matrix $[\chi]$. The transformation holds at any point. Moreover the operation is associative : the combination of relative motions is represented by the product of the matrices. This is convenient, and a big change with what is done in SR with Cartesian frames as we will see.

Tetrad

The vectors of a standard basis (the **tetrad**) can be expressed in the holonomic basis of any chart (of an observer or not).

$$\varepsilon_i(m) = \sum_{\alpha=0}^3 P_i^\alpha(m) \partial \xi_\alpha \Leftrightarrow \partial \xi_\alpha = \sum_{i=0}^3 P_\alpha^i(m) \varepsilon_i(m) \quad (3.32)$$

where $[P]$ is a real invertible matrix (which has no other specific property, it does not belong to $SO(3,1)$) and we denote

Notation 53 $[P'] = [P]^{-1} = [P_\alpha^i]$.

The dual of $(\partial \xi_\alpha)_{\alpha=0}^3$ is $(d\xi^\alpha)_{\alpha=0}^3$ with the defining relation :
 $d\xi^\alpha (\partial \xi_\beta) = \delta_\beta^\alpha$.

The dual $(\varepsilon^i(m))_{i=0}^3$ is :

$$\varepsilon^i(m) = \sum_{\alpha=0}^3 P_\alpha^i d\xi^\alpha \Leftrightarrow d\xi^\alpha = \sum_{i=0}^3 P_i^\alpha \varepsilon^i(m) \quad (3.33)$$

$$\varepsilon^i(m) (\varepsilon_j(m)) = \sum_{\alpha,\beta=0}^3 P_\alpha^i P_j^\beta d\xi^\alpha (\partial \xi_\beta) = \sum_{\alpha=0}^3 P_\alpha^i P_j^\alpha = \delta_j^i$$

Notice that, if it is usual to associate to a vector $u \in T_m M$ a covector : $u^* = \sum_{\alpha,\beta} g_{\alpha\beta} u^\beta d\xi^\alpha$, then $u^*(v) = \langle u, v \rangle$ so that $(\varepsilon_i)^* \neq \varepsilon^i : \varepsilon_i^*(v) = \eta_{ii} v^i \neq \varepsilon^i(v) = v^i$.

In the fiber bundle representation the vectors of the tetrad are variables which are vectors $\varepsilon_i \in \mathfrak{X}(TM)$ or covectors $\varepsilon^i \in \mathfrak{X}(TM^*)$. The quantities $(P_i^\alpha(m))_{i=1}^3$ (called vierbein) and $(P_\alpha^i(m))_{i=1}^3$ are the components of the vectors $\varepsilon_i(m)$ or the covectors $\varepsilon^i(m)$ in any chart. They can be measured, if one has a chart. They depend on the observer, change with the location m and in a change of chart as the components of a vector or a covector. The quantities $\varepsilon_i(m)$ are geometric quantities. They are one of the variables in any model in GR : as such they replace the metric g . However it is obvious that $[P]$ is defined, in any chart, up to a matrix of $SO(3,1)$, so there is some freedom in the choice of the gauge, and we will see the consequences in the specification of a lagrangian.

In a change of gauge on the principal bundle $P_G : \mathbf{p}(m) = \varphi_P(m,1) \rightarrow \tilde{\mathbf{p}}(m) = \mathbf{p}(m) \cdot \chi(m)^{-1}$ the holonomic basis becomes with $[\chi(m)] \in SO_0(3,1)$
 $\varepsilon_i(m) = (\mathbf{p}(m), \varepsilon_i) \rightarrow \tilde{\varepsilon}_i(m) = [\chi(m)]^{-1} \varepsilon_i(m)$
 $\sum_{\alpha=0}^3 \tilde{P}_i^\alpha(m) \partial \xi_\alpha = \left[\chi(m)^{-1} \right]_i^j \sum_{\alpha=0}^3 P_j^\alpha(m) \partial \xi_\alpha$

$$\mathbf{p}(m) = \varphi_P(m, 1) \rightarrow \tilde{\mathbf{p}}(m) = \mathbf{p}(m) \cdot \chi(m)^{-1} : [P] \rightarrow [\tilde{P}] = [\chi(m)]^{-1} [P] \quad (3.34)$$

With respect to the standard chart of the observer :

$$\begin{aligned} \varepsilon_0(p_o(t)) = \partial \xi_0 &\Rightarrow P'_0{}^i = \delta_0^i \\ \alpha = 1, 2, 3 : \frac{\partial}{\partial \xi^\alpha} \varphi_o(\xi^0, \xi^1, \xi^2, \xi^3) = \partial \xi_\alpha &= \frac{\partial}{\partial x} \Phi_{\varepsilon_0}(ct, x) \frac{\partial x}{\partial \xi^\alpha} \in T_m \Omega_3(t) \Rightarrow \\ P'_\alpha{}^0 &= 0 \\ \text{so } [P'] &= \begin{bmatrix} 1 & 0 & 0 & 0 \\ 0 & P'_{11} & P'_{12} & P'_{13} \\ 0 & P'_{21} & P'_{22} & P'_{23} \\ 0 & P'_{31} & P'_{32} & P'_{33} \end{bmatrix}; [P] = \begin{bmatrix} 1 & 0 & 0 & 0 \\ 0 & P_{11} & P_{12} & P_{13} \\ 0 & P_{21} & P_{22} & P_{23} \\ 0 & P_{31} & P_{32} & P_{33} \end{bmatrix} \end{aligned}$$

Metric

The scalar product can be computed from the components of the tetrad. By definition :

$$g_{\alpha\beta}(m) = \langle \partial \xi_\alpha, \partial \xi_\beta \rangle = \sum_{ij=0}^3 \eta_{ij} [P']_\alpha^i [P']_\beta^j$$

The induced metric on the cotangent bundle (Maths.1608) is denoted with upper indexes :

$$g^* = \sum_{\alpha\beta} g^{\alpha\beta} \partial \xi_\alpha \otimes \partial \xi_\beta$$

and its matrix is $[g]^{-1}$:

$$g^{\alpha\beta}(m) = \sum_{ij=0}^3 \eta^{ij} [P]_i^\alpha [P]_j^\beta$$

$$[g]^{-1} = [P] [\eta] [P]^t \Leftrightarrow [g] = [P']^t [\eta] [P'] \quad (3.35)$$

It does not depend on the gauge on P_G :

$$[\tilde{g}] = [\tilde{P}']^t [\eta] [\tilde{P}'] = [P']^t [\chi(m)^{-1}]^t [\eta] [\chi(m)^{-1}] [P'] = [P']^t [\eta] [P']$$

So in the standard chart of the observer : $g^{00} = -1$.

$$[g] = [P']^t [\eta] [P'] = \begin{bmatrix} -1 & 0 \\ 0 & [g]_3 \end{bmatrix}$$

and $[g]_3$ is definite positive.

The metric defines a **volume form** on M (Maths.1609). Its expression in any chart is, by definition :

$$\begin{aligned} \varpi_4(m) &= \varepsilon_0 \wedge \varepsilon_1 \wedge \varepsilon_2 \wedge \varepsilon_3 = \sqrt{|\det [g]|} d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3 \\ [g] &= [P']^t [\eta] [P'] \Rightarrow \det [g] = -(\det [P'])^2 \Rightarrow \sqrt{|\det [g]|} = \det [P'] \\ &\text{assuming that the standard basis of } P_G[\mathbb{R}^4, e] \text{ is direct.} \end{aligned}$$

$$\varpi_4 = \det [P'] d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3 \quad (3.36)$$

Induced metric

The metric on M induces a metric on any submanifold but it is not necessarily non degenerated (Maths.19.3.1).

On hypersurfaces the metric is non degenerated if the unitary normal n is such that $\langle n, n \rangle \neq 0$ (Maths.1642). The induced volume form is (Maths.1644) :

$$\mu_3 = i_n \varpi_4 = \det [P'] d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3 (n)$$

For $\Omega_3(t)$ the unitary normal n is ε_0 , the induced metric is Riemannian and the volume form ϖ_3 is :

$$\begin{aligned} \varpi_3 &= i_{\varepsilon_0} \varpi_4 = \det [P'] d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3 (\varepsilon_0) \\ &= \det [P'] d\xi^0 (\varepsilon_0) \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3 \\ &= \det [P'] d\xi^1 \wedge d\xi^2 \wedge d\xi^3 \end{aligned}$$

$$\varpi_3 = \det [P'] d\xi^1 \wedge d\xi^2 \wedge d\xi^3 \quad (3.37)$$

and conversely :

$$\varpi_4 = \varepsilon_0 \wedge \varpi_3 = \det [P'] d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3$$

ϖ_3 is defined with respect to the coordinates ξ^1, ξ^2, ξ^3 but the measure depends on $\xi^0 = t$.

For a curve C , represented by any path : $p : \mathbb{R} \rightarrow C :: m = p(\theta)$ the condition is $\left\langle \frac{dp}{d\theta}, \frac{dp}{d\theta} \right\rangle \neq 0$. The volume form on any curve defined by a path : $q : \mathbb{R} \rightarrow M$ with tangent $V = \frac{dq}{d\theta}$ is $\sqrt{|\langle V, V \rangle|} d\theta$. So on the trajectory $q(t)$ of a particle it is

$$\varpi_1(t) = \sqrt{-\langle V, V \rangle} dt \quad (3.38)$$

For a particle there is the privileged parametrization by the proper time, and as $\left\langle \frac{dp}{d\tau}, \frac{dp}{d\tau} \right\rangle = -c^2$ the length between two points A,B is :

$$\ell_p = \int_{\tau_A}^{\tau_B} \sqrt{-\left\langle \frac{dp}{d\tau}, \frac{dp}{d\tau} \right\rangle} d\tau = \int_{\tau_A}^{\tau_B} c d\tau = c(\tau_B - \tau_A)$$

This is an illustration of the idea that all world lines correspond to a travel at the same speed.

3.2.5 From particles to material bodies

In Mechanics a material body is comprised of “material points” that is elements of matter whose location is a single geometric point, and change with time in a consistent way : their trajectories do not cross, and the body keeps some cohesion, which is represented by a deformation tensor for deformable solid bodies. In Relativity the material points, particles in our terminology, follow independent world lines, which do not cross, and thus can be represented by a field of vectors u , future oriented with length $\langle u, u \rangle = -c^2$, such that, at some time 0, the particles are all together in a compact subset $\varpi(0)$ of a 3 dimensional space like submanifold. Then the location of any particle of the material body is

given by $\Phi_u(\tau, a)$ where τ is its proper time and $a \in \varpi(0)$ its location at $\tau = 0$. The area swept by the body is $\widehat{\omega} = \{\Phi_u(\tau, a) : \tau \in \mathbb{R}, a \in \omega(0)\}$ and we have the function : $f : \widehat{\omega} \rightarrow \mathbb{R} : f(m) = \tau : \exists a \in \omega(0) : \Phi_u(\tau, a) = m$. The function f defines a foliation in diffeomorphic compact 3 dimensional hypersurfaces $\omega(\tau)$ which can be seen as the state of the material body at τ (Maths.1503). So $\Phi_u(\tau, a)$ can be seen as a chart of $\widehat{\omega}$, and the material body has a unique proper time τ . We can then give a definition of a material body which is independent of any observer.

Definition 54 A *material body* is defined by a field of vectors u , future oriented with length $\langle u, u \rangle = -c^2$, and a compact subset $\varpi(0)$ of a 3 dimensional space like submanifold. The body is located at its proper time τ on the set $\omega(\tau)$ diffeomorphic to $\varpi(0)$.

The vector field $u \in \mathfrak{X}(TM)$ does not depend on a chart, but for any observer O the trajectories of the material points of the body will follow a vector field $V \in \mathfrak{X}(TM)$: the curves do not change, but they are traveled with the parameter t and not τ .

A material point can be labeled by its position $y \in \omega(0)$ then its location p is

$\Phi_u(\tau, a)$ at the proper time τ along the world line going through y

$\Phi_{\varepsilon_0}(t, x(t)) = \varphi_O(t, x(t))$ at the time t of the observer in the chart of

the observer given by his vector field ε_0

and $\tau, t, y, x(t)$ are related by :

$$u = \frac{dp}{d\tau}$$

$$V(t) = \frac{dp}{dt} = u \frac{\sqrt{-\langle V, V \rangle}}{c} = c\varepsilon_0 + \vec{v}$$

$$\vec{v} = \frac{dx}{dt}$$

$$x(0) = a$$

At $t = 0$ for the observer the set of points of the material body is $\widehat{\omega}(0) = \widehat{\omega} \cap \Omega(0)$

At $t > 0$ for the observer the set of points of the material body is $\widehat{\omega}(t) = \widehat{\omega} \cap \Omega(t)$

$$\widehat{\omega}(t) = \{\varphi_O(t, x(t)), x(0) \in \widehat{\omega}(0)\}$$

$$= \{\Phi_V(t, x(0)), x(0) \in \widehat{\omega}(0)\}$$

$$= \{\Phi_u(\tau(t), a), a \in \widehat{\omega}(0)\}$$

$$\text{and } \omega(\tau(t)) = \{\Phi_u(\tau(t), y), y \in \omega(0)\}$$

So the characterization of a material body is observer dependant : they do not see the same body.

However we will assume :

Proposition 55 For any material body there are observers with proper time t such that, at $t = 0$ they observe the entire material body : $\omega(0) \subset \Omega(0)$

Then at any given time t : $\widehat{\omega}(t) = \omega(\tau(t))$. This is a legitimate assumption, which will be mainly used to compute the characteristics of material bodies.

In Mechanics a solid is a material body such that the distance between any two of its points is constant. $\omega(\tau)$ is a Riemannian manifold, with the metric g_τ induced by the metric g on M . g_τ defines the length of a curve on $\omega(\tau)$ and the distance between two of its points is then the minimum of the length of all the lines which join the points. Because $\omega(\tau)$ is compact such a minimum exists, however the metric g_τ itself depends on the point and τ , so this concept of solid cannot be extended in Relativity.

In Galilean Geometry the local deformation of a material body is studied through the change of a frame attached to each point, with respect to a fixed common frame. In the previous framework the local deformation can be measured through the derivative of the flow Φ_u . Choose a chart of $\varpi(0) : a = \phi(\eta_1, \eta_2, \eta_3)$ then the derivatives $u = \frac{\partial \Phi_u}{\partial \tau}$, $u_i = \frac{\partial \Phi_u}{\partial \eta_i}$, $i = 1, 2, 3$ provides a basis at each point of $\omega(\tau)$ and the components of these vectors with respect to any chart of M provide a deformation tensor. We will see in the next chapter how one can improve these points.

3.2.6 Special Relativity

All the results of this chapter hold in Special Relativity. This theory, which is still the geometric framework of QTF and Quantum Physics, adds two assumptions : the Universe M can be represented as an affine space, and the metric does not depend on the location (these assumptions are independent). As consequences :

- the underlying vector space \vec{M} (the Minkovski space) is common to all observers : the vectors of all tangent spaces to M belong to \vec{M}
- one can define orthonormal bases which can be freely transported and compared from a location to another
 - because the scalar product of vectors does not depend on the location, at each point one can define time-like and space-like vectors, and a future orientation (this condition relates the mathematical and the physical representations, and \vec{M} is not simply \mathbb{R}^4)
- there are fixed charts $(O, (\varepsilon_i)_{i=0}^3)$, called frames, comprised of an origin (a location O in M : a point) and an orthonormal basis $(\varepsilon_i)_{i=0}^3$. There is necessarily one vector such that $\langle \varepsilon_i, \varepsilon_i \rangle = -1$. It is possible to define, non unique, orthonormal bases such that ε_0 is timelike and future oriented.
- the coordinates of a point m , in any frame $(O, (\varepsilon_i)_{i=0}^3)$, are the components of the vector OM . The transition maps which give the coordinates of m in another frame $(A, (\tilde{\varepsilon}_i)_{i=0}^3)$ are then given by the formulas :

$$\begin{aligned}
 OM &= \sum_{i=0}^3 x_i \varepsilon_i \\
 AM &= \sum_{i=0}^3 \tilde{x}_i \tilde{\varepsilon}_i \\
 OM &= OA + AM = \sum_{i=0}^3 L_i \varepsilon_i + \sum_{i=0}^3 \tilde{x}_i \tilde{\varepsilon}_i \\
 \tilde{\varepsilon}_i &= \sum_{j=0}^3 [\chi]_i^j \varepsilon_j, [\chi] \in SO(3,1)
 \end{aligned}$$

However one needs to go from this abstract representation to a physical definition of frames.

Observers can label points which are in their present with their proper time. The role of the function $f(m) = t$ is crucial, because it defines the 3 dimensional hypersurfaces $\Omega(t)$. They are not necessarily hyperplanes, but they must be space like and do not cross each other : a point m cannot belong to 2 different hypersurfaces. These hypersurfaces define the vector field $\varepsilon_0(m)$ to which belongs the velocity of the observer (up to c). In SR one can compare vectors at different points, and usually the vectors $\varepsilon_0(m)$ are different from one location to another. They are identical only if $\Omega(t)$ are hyperplanes normal to a vector ε_0 , which implies that the world line of the observer is a straight line, and because the proper time is the parameter of the flow, if the motion of the observer is a translation at a constant spatial speed. These observers are called **inertial**. Notice that this definition is purely geometric and does not involve gravitation or inertia : the motion of an observer is absolute, and inertial observers are such that their velocity is a constant vector.

Observers can define a standard chart $\varphi_o(\xi^0, \xi^1, \xi^2, \xi^3)$ with $\xi^0 = ct, \partial\xi_0 = \varepsilon_0(m)$ with the flow of $\varepsilon_0(m)$ and a chart φ_Ω of $\Omega(0)$ which provides the coordinates ξ^1, ξ^2, ξ^3 of the point x where the integral curves of ε_0 passing through m crosses $\Omega(0)$. The general results hold and such a chart can always be defined. However this chart is usually not defined by a frame $(O, (\varepsilon_i)_{i=0}^3)$: the vectors of the basis must be constant, and notably ε_0 so this is possible only if the observer is inertial : *a frame can be associated to an observer only if this is an inertial observer.*

For inertial observers the integral curves are straight lines parallel to ε_0 . Any spatial basis $(\varepsilon_i)_{i=1}^3$ of $\Omega(0)$ can be transported on $\Omega(t)$. The standard chart is then similar to a frame in the 4 dimensional affine space $(O(0), (\varepsilon_i)_{i=0}^3)$ with origin $O(0)$, the 3 spatial vectors $(\varepsilon_i)_{i=1}^3$ and the time vector ε_0 . The coordinates of a point $m \in \Omega_3(t)$ are :

$$\overrightarrow{O(0)m} = ct\varepsilon_0 + \sum_{i=1}^3 \xi^i \varepsilon_i \text{ where } \overrightarrow{O(t)m} = \sum_{i=1}^3 \xi^i \varepsilon_i$$

and the velocity of a particle with trajectory $q(t)$, as measured by O, is :

$$V = c\varepsilon_0 + \vec{v} \text{ with } \vec{v} = \frac{d}{dt} \overrightarrow{O(t)q(t)} = \sum_{i=1}^3 \frac{d\xi^i}{dt} \varepsilon_i \text{ because } \frac{d\varepsilon_i}{dt} = 0$$

If there is another inertial observer with standard chart defined by a frame $(A(0), (\tilde{\varepsilon}_i)_{i=0}^3)$ the coordinates of $A(t)$, as seen by $O(t)$, are :

$$\overrightarrow{O(0)A(t)} = ct\varepsilon_0 + \sum_{i=1}^3 \xi^i(t) \varepsilon_i \text{ where } \overrightarrow{O(t)A(t)} = \sum_{i=1}^3 \xi^i(t) \varepsilon_i$$

$$\text{The spatial speed of A is : } \vec{v} = \frac{d}{dt} \overrightarrow{O(t)A(t)} = \sum_{i=1}^3 \frac{d\xi^i}{dt} \varepsilon_i$$

We can then implement the general results for the change of basis : $\varepsilon_i \rightarrow \tilde{\varepsilon}_i$.

As for the change of coordinates we have :

$$\begin{aligned} \overrightarrow{O(0)m} &= ct\varepsilon_0 + \sum_{i=1}^3 \xi^i \varepsilon_i \\ \overrightarrow{A(0)m} &= c\tau\tilde{\varepsilon}_0 + \sum_{i=1}^3 \tilde{\xi}^i \tilde{\varepsilon}_i \end{aligned}$$

$$\begin{aligned} [\tilde{\xi}^i] &= [\chi]^{-1} [\xi^i] \\ \overrightarrow{O(0)m} &= \sum_{i=0}^3 \xi^i \varepsilon_i = \overrightarrow{O(0)A(0)} + \overrightarrow{A(0)m} = L + \sum_{i=0}^3 \tilde{\xi}^i \tilde{\varepsilon}_i \end{aligned}$$

with a constant vector $\overrightarrow{O(0)A(0)} = L$.

So the transformation of the coordinates is given by the product of a fixed translation and a fixed rotation in the Minkovski space. The set of such transformations is a group, called the Poincaré's group.

This result holds only for two inertial observers, and we need a physical mean to tell what are these observers. The usual answer is that they do not feel a change in the inertial forces to which they are submitted. This is similar to the Galilean observers of Classic Mechanics. For non inertial observers the general formulas hold, but the charts cannot be defined through frames as in an affine space.

The concept of material body presented above holds. But if $\varpi(0)$ belongs to a hyperplane then the $\varpi(\tau)$ will be hyperplanes only if they are all orthonormal to a common vector, that is if the vector field which defines the material body is a constant vector : the body must be in a uniform translation (and not rotate on itself).

The formulas of the Lorentz transformations have a tremendous importance in all aspects of Relativist Physics, they are of a constant use, as well as the Poincaré's group which is the starting point in the identification of particles. However any demonstration based on frames, as it is usually done, holds only for inertial observers. A physical theory which is valid only in the study of bodies in uniform translation would be of little interest. As we have proven in this chapter Relativist Geometry can be explained, in a rigorous and quite simple way, without the need of inertial observers. And these are required only for the use of frames. It would be a pity to loose the deep import of Relativity in order to keep a familiar, but not essential, mathematical tool. As a consequence the role assigned to the Poincaré's group must be revisited.

3.3 SOME ISSUES ABOUT RELATIVITY

It is useful to review here some issues which arise frequently about Relativity.

3.3.1 Preferred frames

Relativity is often expressed as “all inertial frames are equivalent for the Physical Laws”. We have seen above that actually inertial frames are required only to define coordinates in affine space : this is a non issue in GR, and in SR it is possible to achieve the usual results with the use of standard charts which are not given by orthogonal frames. But, beyond this point, this statement is misleading.

The Theory of Relativity is more specific than the Principle of Relativity, it involves inertia and gravitation (that we will see in the next chapters), but this is at first a Theory about the Geometry of the Universe, and it shows that the geometric measures (of lengths and time) are specific to each observer. The Universe which is Scientifically accessible - meaning by the way of measures, data and figures - depends on the observer. We can represent the Universe with 4 dimensions, conceive of a 4 dimensional manifold which extends over the past and the future, but we must cope with the fact that we are stuck into our present, and it is different for each of us. The reintegration of the observer in Physics is one of the most important feature of Relativity, and the true meaning of the celebrated formulas for a change of frames. An observer is an object in Physics, and as such some properties are attached to it, among them the free will : the possibility to choose the way he proceeds to an experiment, without being himself included in the experiment. But as a consequence the measures are related to his choice.

Mathematics give powerful tools to represent manifolds, in any dimensions. And it seems easy to formulate any model using any chart as it is commonly done. This is wrong from a physical point of view. There is no banal chart or frame : it is always linked to an observer, there is a preferred chart, and so a preferred frame for an observer. It is not related to inertia : it is a matter of geometry, and a consequence of the fundamental symmetry breakdown. The observer has no choice in the selection of the time vector of his orthonormal basis, if he wants to change the vector, he has to change his velocity, and this is why the formulas in a change of frames are between two different observers moving with respect to each other. And not any change is possible : an observer cannot travel in the past, or faster than light. These features are clear when one sticks to a chart of an observer, as we will do in this book. Not only they facilitate the computations, they are a reminder of the physical meaning of the chart. This precision is specially important in the fiber bundle formalism, which is, from this point of view, a wise precaution as compared to the usual formalism using undifferentiated charts.

3.3.2 Time travel

The distinction between future and past oriented vectors come from the existence of the Lorentz metric. As it is defined everywhere, it exists everywhere, and along any path. It is not difficult to see that the border between the two kinds of vectors is for null vectors $\langle u, u \rangle = 0$. So a particle which would have a path such that its velocity is past oriented should, at some point, have a null velocity, and, with regard to another observer located at the same point, travel at the speed of light. Afterwards his velocity would be space like ($\langle u, u \rangle > 0$) before being back time like but past oriented. So the first issue lies in the possibility of travel at the speed of light. Clearly this would be a discontinuity on the path and "Scotty engages the drive" from Star Trek has some truth.

But the main issue with time travel lies in the fact that, if ever we would be able to come back to the location where we have been in the past (meaning a point of the universe located in our past), we would not find our old self. The idea that we exist in the past assumes that we exist at any time along our world line, as a frozen copy of ourselves. This possibility is sometimes invoked, but it raises another one : what makes us feel that each instant of time is different ? If we do not travel physically along our world line, what does move ? And of course this assumption raises many other issues in Physics...

3.3.3 Twins paradox

The paradox is well known : one of the twins embarks in a rocket and travels for some time, then comes back and finds that he is younger than his twin who has stayed on Earth. This paradox is true (and has been checked with particles) and comes from two relativist features : the Universe is 4 dimensional, and the definition of the proper time of an observer.

Because the Universe is 4 dimensional, to go from a point A to a point B there are several curves. Each curve can be travelled according to different paths. We have assumed that observers move along a curve according to a specific path, their world line, and then :

$$\ell_{AB} = c(\tau_B - \tau_A)$$

Because the curves are different, the elapsed proper time is usually different.

The proper time is the time measured by a clock attached to the observer, it is his biological time. Assuming that all observers travel along their world lines with a velocity such that at $\left\langle \frac{dp_\alpha}{d\tau}, \frac{dp_\alpha}{d\tau} \right\rangle = -c^2$ is equivalent to say that, with respect to their clock, they age at the same rate. So if they travel along different curves there is no reason that the total duration of their travel would be the same.

Whom of the two twins would have aged the most ? It is not easy to do the computation in GR, but simpler in the SR context.

We can define a fixed frame $(O, (\varepsilon_i)_{i=0}^3)$ with origin O at the time $t = 0$, A is spatially immobile with respect to this frame, moves along the time axis and his coordinates are then : $OA : p_A(\tau_A) = c\tau_A\varepsilon_0$

The twin B moves in the direction of the first axis. His coordinates are then : $OB : p_B(\tau_B) = c\tau_B\varepsilon_0 + x_B(\tau_B)\varepsilon_1$

The spatial speed of B with respect to A is : $\frac{dOB}{d\tau_A} = V(\tau_B)\varepsilon_1$

The velocity of B is : $u_B = \frac{dOB}{d\tau_B} = \frac{1}{\sqrt{1-\frac{V^2}{c^2}}}(V\varepsilon_1 + c\varepsilon_0)$

To be realistic we must assume that B travels at a constant acceleration, but needs to brake before reaching first his turning point, then A. In the first phase we have for instance :

$$V = \gamma c\tau_B \text{ with } \gamma = \frac{1}{\sqrt{1-\frac{V^2}{c^2}}}$$

$$p_B(\tau_B) = \int_0^{\tau_B} \frac{c}{\sqrt{1-(\gamma t)^2}} (\gamma t\varepsilon_1 + \varepsilon_0) dt = \frac{c}{\gamma} \left[\sqrt{1-y^2}\varepsilon_1 + \varepsilon_0 \arcsin y \right]_0^{\gamma\tau_B}$$

A full computation gives : $\frac{\tau_A}{\tau_B} = \frac{\arcsin v_M}{v_M}$ where v_M is the maximum speed in the travel, which gives for $v_M = c$: $\frac{\tau_A}{\tau_B} = 1.57$ that is less than what is commonly assumed.

The Sagnac effect, used in accelerometers, is based on the same idea : two laser beams are sent in a loop in opposite direction : their 4 dimensional paths are not the same, and the difference in the 4 dimensional lengths can be measured by interferometry.

3.3.4 The expansion of the Universe

A manifold by itself can have some topological properties. It can be compact. It can have holes, defined through homotopy (Maths.10.4.1) : there is a hole if there are curves in M which cannot be continuously deformed to be reduced to a point. A hole does not imply some catastrophic feature : a doughnut has a hole. Thus it does not imply that the charts become singular. But there are only few purely topological features which can be defined on a manifold, and they are one of the topic of Differential Geometry. In particular a manifold has no shape to speak of.

The metric on M is an addition to the structure of the Universe. It is a mathematical feature from which more features can be defined on M, such that curvature (we will see it in another chapter). It is also a physical feature and in GR the metric, and so the curvature of M at a point, depends on the distribution of matter. It is customary (see Wald) to define singularities in the Universe by singularities of geodesics, but geodesics are curves whose definition depends on the metric. A singularity for the metric, as Black holes or Bing Bang, is not necessarily a singular point for the manifold itself.

GR has open the possibility to build cosmological models, representing the totality of the Universe. It is clear that Cosmology requires some revision of the usual concepts of Physics (what is an independent observer ?) and even of the premises of our epistemology (do we have free will ?). A requirement which is rarely fulfilled in the common Cosmological Theories. From some general reasoning and Astronomical observations, it is generally assumed that the Universe has the structure of a fiber bundle with base \mathbb{R} (a warped Universe) which can be seen as the generalization of M_o , that we have defined above for an

observer. Thus there is some universal time (the projection from M to \mathbb{R}) and a foliation of M in hypersurfaces similar to $\Omega_3(t)$, which represent the present for the observers who are located on them (see Wald and Peebles for more on this topic). This is what we have defined as a material body : the part of the universe on which stands all matter would be a single body moving together since the Big Bang (the image of an inflating balloon). So there would not be any physical content before or after this $\Omega_3(t)$ (inside the balloon), but nothing can support this interpretation, or the converse, and probably it will never be.

The Riemannian metric $\varpi_3(t)$ on each $\Omega_3(t)$ is induced by the metric on M , and therefore depends on the universal time t . In the most popular models it comes that the distance between two points on $\Omega_3(t)$, measured by the Riemannian metric, increases with t , and this is the foundation of the narrative about an expanding universe, which is supported by astronomical observations. But, assuming that these models are correct, this needs to be well understood. The change of the metric on $\Omega_3(t)$ makes that the volume form $\varpi_3(t)$ increases, but the hypersurfaces $\Omega_3(t)$ belong to the same manifold M , which does not change with time. The physical universe would be a deformable body, whose volume increases inside the unchanged container. Moreover it is generally assumed that material points, belonging to the same material body but traveling on their own world lines, stick together : they are not affected by this dilation, only the vacuum which separates material bodies.

Chapter 4

KINEMATICS

Kinematics is the next step after Geometry. We go further in the physical world, and try to understand what are the relations between the motion of material bodies and the forces which are applied to them. All material bodies manifest some resistance to a change of their motion, either in direction or in speed. This feature is the inertia, and is measured by different quantities which incorporate, in one way or another, the mass of the material body. The mass m is a characteristic of the material body : it does not change with the motion, the forces or the observer. And from motion and mass are defined key quantities : the momenta.

The Newton's law : $\vec{F} = m\vec{\gamma}$ is expressed, more appropriately by : $\vec{F} = \frac{d\vec{p}}{dt}$ where \vec{p} is the momentum. The inertial forces are, by construct, the opposite of the forces which are necessary to change the momenta of a material body. So, Kinematics is, in many ways, the Physics of inertia. The issue of the origin of these inertial forces, which appear everywhere and with a great strength, will be seen in the next chapter. Let us see now how one goes from Geometry, that is motion, to Kinematics, that is inertia.

The study of rotations and rotational momenta in the 4 dimensional Universe will lead to a new representation of the momenta, based on Spinors, which are vectors in a 4 dimensional abstract space, to the introduction of antiparticles and of spin.

4.1 TRANSLATIONAL AND ROTATIONAL MOMENTA

4.1.1 Translational Momentum in the relativist context

In Newtonian Physics the bridge between Geometry and Kinetics is hold by momenta. And, because motion of material bodies involves both a translational motion and a rotational motion, there are linear momenta and rotational momenta.

In Galilean Geometry the linear momentum of a particle is simply : $\vec{p} = \mu \vec{v}$ with a constant scalar μ which is the inertial mass. Its natural extension in the relativist context is the quadri-vector : $P = \mu u$ where u is the velocity. This generalization has two consequences :

- the quadri-vector P is intrinsic : its definition as well as the value of the scalar product $\langle P, P \rangle = -\mu^2 c^2$ do not depend on an observer
- but its measure depends on the observer. In his standard basis it reads :

$$P = \mu \frac{1}{\sqrt{1 - \frac{\|\vec{v}\|^2}{c^2}}} (\vec{v} + c\varepsilon_0(m))$$

In the relativist context location and motion are absolute. If the Universe has a physical meaning, then each of its points is singular, and this is clearly represented by a manifold. The proper time, and the derivative of the location with respect to the proper time, are defined without any reference to a frame, so the vector velocity u is absolute, and this property has been used to compute the rules in a change of frames. If motion is absolute, *its measure is relative*, depends on the observer and its measure changes according to geometric rules, because they are geometric quantities. The spatial speed appears when an observer has been chosen. The definition of the momentum by $P = \mu u$ is consistent with the idea that the kinematic features of a particle are intrinsic, and can be represented by a quantity which does not depend on an observer (even if its measure depends on it).

This is a big change from the Newtonian definitions : the momentum $\vec{p} = \mu \vec{v}$ as well as the kinetic energy $\frac{1}{2}\mu \|\vec{v}\|^2$ are relative and depend on the observer.

If we keep the concept of Energy as measured by [Momentum] \times [Speed] then the energy of the particle $\langle P, u \rangle = -\mu c^2$ is constant and, for an observer, is split between a part related to the spatial speed $\mu \frac{\|\vec{v}\|^2}{1 - \frac{\|\vec{v}\|^2}{c^2}}$, corresponding to a kinetic energy, and a part which is stored in the particle $-\mu \frac{c^2}{1 - \frac{\|\vec{v}\|^2}{c^2}}$. But, if one wants to keep the principle of conservation of energy, one has to accept that mass itself can be transformed into energy, according to the famous relation $E = \mu c^2$.

However there are several interpretations of these concepts. Physicists like to keep a concept of momentum linked to the spatial velocity and, with a fixed mass, define the linear momentum as : $\vec{p}_r = \mu \frac{\vec{v}}{\sqrt{1 - \frac{\|\vec{v}\|^2}{c^2}}}$, that is the spatial

part of P . Then they define the Energy E of the particle by : $E^2 = c^2 \|\vec{p}_r\|^2 + \mu^2 c^4$ that is one part corresponding to a kinetic energy, and another one to an energy at rest. This sums up to define the energy by rewriting Pc with the two components :

$$Pc = (c\vec{p}_r, E) \Rightarrow \langle Pc, Pc \rangle = -\mu^2 c^4 = c^2 \|\vec{p}_r\|^2 - E^2$$

The introduction of c in $c\vec{p}_r$ is necessary in order to have the same unit Energy in both parts. In this formulation Pc is a 4 vector, and its components change according to the Lorentz formula, so E depends on the observer. The advantages of this expression is that for small speed it gives :

$$\begin{aligned} E &\simeq \sqrt{c^2 \mu^2 \|\vec{v}\|^2 \left(1 + \frac{1}{2} \frac{\|\vec{v}\|^2}{c^2}\right) + \mu^2 c^4} \\ &= \mu c^2 \sqrt{\left(1 + \frac{\|\vec{v}\|^2}{c^2}\right) \left(1 + \frac{1}{2} \frac{\|\vec{v}\|^2}{c^2}\right)} \simeq \mu c^2 \sqrt{1 + \frac{\|\vec{v}\|^2}{c^2}} \simeq \frac{1}{2} \mu \|\vec{v}\|^2 + \mu c^2 \end{aligned}$$

and it can be adapted to massless particles such as photons.

So if one can say, with Feynman (in *Lectures in Physics*), that the conservation of energy is a law without exception, the trouble is that in the Relativist context the definition of energy itself depends on the observer and is fairly subtle...We will see that, if it is possible to give a satisfying definition of the energy of a particle, this is more complicated for fields, and further more for bosons.

The only physical quantity which has a clear meaning, is independent of an observer, and is characteristic of a particle, is Pc , which has the dimension of energy, and that we will call **relativist momentum**. The mass at rest, independent from an observer, is $\mu = \frac{1}{c} \sqrt{-\langle Pc, Pc \rangle}$. The usual energy is computed by taking the component of Pc along the direction of $\varepsilon_0(m)$. So it depends on the observer. These quantities are individualized and localized : they are linked to the particle and its position.

Moreover material bodies have a rotation, and a rotational momentum \vec{J} which depends on the mass and the shape of the body. Momenta are characteristics features of material bodies, and the bases for the definition of forces. According to the Newton's law: $\vec{F} = \frac{d\vec{p}}{dt}$ and $\vec{\gamma} = \frac{d\vec{J}}{dt}$ for a torque $\vec{\gamma}$. Rotation and rotational momentum are topics which are more complicated than it seems, in the relativist context in particular. Moreover at the atomic scale particles show properties which look like rotation, and specific features, which have lead to the concepts of spins and spinors. These are the main topics of this chapter.

4.1.2 The issues of the concept of rotation

Rotation in Galilean Geometry

The concept of rotation is well defined in Mathematics : this is the operation which transforms the *orthonormal basis* of a vector space into another, and in Galilean Geometry is represented by a matrix of the group $SO(3)$. This is a

compact, 3 dimensional Lie group, of matrices such that $R^t R = I$. Then an instantaneous rotation, that is the derivative of a rotation with respect to time, is represented by an element of its Lie algebra $so(3)$ which is the vector space of 3×3 real antisymmetric matrices. If we take the following matrices as basis of $so(3)$:

$$\kappa_1 = \begin{bmatrix} 0 & 0 & 0 \\ 0 & 0 & -1 \\ 0 & 1 & 0 \end{bmatrix}; \kappa_2 = \begin{bmatrix} 0 & 0 & 1 \\ 0 & 0 & 0 \\ -1 & 0 & 0 \end{bmatrix}; \kappa_3 = \begin{bmatrix} 0 & -1 & 0 \\ 1 & 0 & 0 \\ 0 & 0 & 0 \end{bmatrix}$$

then any matrix of $so(3)$ reads :

$\sum_{i=1}^3 r^i [\kappa_i] = [j(r)]$ with the operator

$$j : \mathbb{R}^3 \rightarrow L(\mathbb{R}, 3) :: [j(r)] = \begin{bmatrix} 0 & -r_3 & r_2 \\ r_3 & 0 & -r_1 \\ -r_2 & r_1 & 0 \end{bmatrix} \quad (4.1)$$

The operator j is very convenient to represent quantities which are rotated¹ and has many nice algebraic properties (see formulas in the annex) and we will use it often in this book.

For any vector $u : \sum_{ij=1}^3 [j(r)]_j^i u^j \varepsilon_i = \vec{r} \times \vec{u}$ with the cross product \times .

Because $SO(3)$ is a compact Lie group the exponential is surjective :

$\forall g \in SO(3), \exists \kappa = \sum_{i=1}^3 r^i [\kappa_i] \in so(3) : g = \exp \kappa$

It is easy to show that :

$$[g] = \exp [j(r)] = I_3 + [j(r)] \frac{\sin \sqrt{r^t r}}{\sqrt{r^t r}} + [j(r)] [j(r)] \frac{1 - \cos \sqrt{r^t r}}{r^t r}$$

The representation of an instantaneous rotation by a vector $\vec{r} \in \mathbb{R}^3$ is further supported by two facts.

The axis of rotation, which is by definition the unique eigen vector of $[g]$ with eigen value 1 and norm 1 in the standard representation of $SO(3)$, has for

components $\begin{bmatrix} r^1 \\ r^2 \\ r^3 \end{bmatrix} / \sqrt{r^t r}$

Similarly one can define the angle θ of the rotation resulting from a given matrix, and $\theta = \sqrt{r^t r}$, which is also the instantaneous rotational speed with the period $T = 2\pi / \sqrt{r^t r}$

Proof. For any vector u of norm 1 : $\langle u, [g]u \rangle = \cos \theta$ where θ is an angle which depends on u and $[g] = \exp [j(r)]$. With the formula above, and using $[j(r)] [j(r)] = [r] [r]^t - \langle r, r \rangle I$ and $\langle u, [j(r)]u \rangle = 0$ we get :

$$\langle u, [g]u \rangle = 1 + \left(\langle u, r \rangle^2 - \langle r, r \rangle \right) \frac{1 - \cos \sqrt{r^t r}}{r^t r}$$

which is minimum for $\langle u, r \rangle = 0$ that is for the vectors orthogonal to the axis, and :

$$\cos \theta = \cos \sqrt{r^t r}$$

For a rotation at constant speed $\varpi : \theta(t) = \varpi t = t\sqrt{r^t r}$ so $T = 2\pi / \sqrt{r^t r}$ ■

¹It is similar to the Levi-Civita tensor ϵ but, in my opinion, much easier to use.

So we have a very satisfying representation of geometric rotations : a rotation $R(t)$ can be defined by a single vector, which can be easily related to the motion as it can be perceived. However this mathematical representation is not faithful. The same rotation can be defined equally by the opposite axis, and the opposite angle. This is related to the mathematical fact that $SO(3)$ is not the only group which has $so(3)$ as Lie algebra. The more general group is the Spin group $Spin(3)$ which has also for elements the scalars $+1$ and -1 , so that $R(t)$, corresponding to $(r, \sqrt{r^t r})$ and $-R(t)$, corresponding to $(-r, -\sqrt{r^t r})$ can represent the same physical rotation. So, actually, the group which should be used to represent rotations in Galilean Geometry is $Spin(3)$, which makes the distinction between the two rotations, and not $SO(3)$. This is not a problem in Mathematics, but in Physics the distinction matters : in the real world one goes from one point to another along a path, by a continuous transformation which preserves the orientation of a vector, thus the orientation of \vec{r} matters ². A single vector of \mathbb{R}^3 cannot by itself properly identify a physical rotation, one needs an additional parameter which is ± 1 (to tell which one of the two orientations of \vec{r} is chosen, with respect to a direction, the spatial speed on the path).

But to represent rotation of material bodies by geometric rotations, as above, raises several issues.

We could expect that the total rotational momentum of a body is the sum of the rotational momentum of its components, as it happens with the translational momentum. But material points have no attached frame in Mechanics. So actually the rotational momentum is defined only at the level of the body, through a geometric rotation, and it has a meaning only for material bodies which keep some shape, represented by fixed relations between the positions of the material points. So this is doable for solids, but even for them there is a problem. If the solid has a cylindrical symmetry, by definition it is impossible to measure a geometric rotation around the axis, however the physical rotation can be measured by a rotational inertia. And some force fields, such as magnetism, can exercise a pointwise action, represented by a torque, so that the implementation of the Newton's law to the rotational momentum becomes muddled for particles without structure, like the atoms.

So, even if in Mechanics it is convenient to assimilate physical rotation with geometric rotation, they are not the same and the kinematics of rotating bodies is more complicated. Notice also that the representation of rotation by a single vector in Galilean Geometry relies on a fortuitous fact : the lie algebra $so(3)$ is a 3 dimensional vector space, as the space \mathbb{R}^3 , so that it is easy to set up an isomorphism j of vector spaces.

To represent the motion (translation + rotation) of a solid body we have another discrepancy between the physical and the mathematical definitions. In Classic Mechanics the motions of solids are defined by two vectors (\vec{v} , $\vec{\omega}$) and 6

²In his book "The road to reality" Penrose gives a nice, simple trick with a belt and book to show this fact.

parameters : the spatial speed \vec{v} of the center of mass G and the instantaneous rotation \vec{r} , assimilated to 2 vectors of \mathbb{R}^3 . In Mathematics the transformation between two frames is represented by the group of displacements, which is the semi-product of $SO(3)$ with the group of translations. Using the isomorphism j between $SO(3)$ and \mathbb{R}^3 a displacement can be represented by two vectors (\vec{L}, \vec{r}) but \vec{L} is the translation of the origins of the frames : $\vec{L} = \vec{OG}$ and not the speed $\vec{v} = \frac{d\vec{OG}}{dt}$.

The group of displacements in Relativist Geometry

In Relativist Geometry the focus is usually put on the Poincaré's group, the semi product of the group $SO(3, 1)$ of rotations and of the 4 dimensional translations. This is the simple generalization of the group of displacements of Galilean Geometry. However this raises several issues.

The Poincaré's group represents the transformations between coordinates in Cartesian frames. So its use is valid only in SR, and for inertial observers. There has been attempts to extend the concept to the group of isometries (that is maps $f : M \rightarrow M$ such that the derivative $f'(m)$ preserves the scalar product) but this is difficult. And there is a more fundamental objection : the physical comparison between bases located at different points should involve a transport of the vectors (one does not jump from a point to another), but in SR or GR the universe is no longer isotropic, so the path which is followed for the transport matters (which is obvious in the Lorentz formulas, which involve \vec{v}).

The concept of solid is not generalizable in Relativity, thus one cannot expect to represent the motion of a body by the transformation to go from the frame of an observer to a frame which would be attached to the solid.

And actually, even in Galilean Geometry the point of view of the Physicists (who represent a motion by (\vec{v}, \vec{r})) and of the mathematicians (who define the displacement by (\vec{L}, \vec{r})) are not the same : even if they both use two vectors and the same number of parameters, they have not the same meaning. In Relativity the discrepancy is worse : an element of the Poincaré's group is defined by 10 parameters. Geometric rotations in the 4 dimensional space have a very different meaning than rotations in the 3 dimensional space. This is obvious with a look at the matrices of $SO(3, 1)$ which read : $[\chi] = \exp [K (w)] \exp [J (r)]$. The second term has the meaning of a rotation in the space, but the first term (usually called the boost) involves the translational motion, as shown in the formula : $w = \frac{v}{\|v\|} \arg \tanh \left\| \frac{v}{c} \right\|$.

For a physicist the motion of a material body is related to the instantaneous change of its location and disposition (as it is done in Galilean Geometry with (\vec{v}, \vec{r})) and not to the transformation between fixed frames (the vector \vec{L} has no physical interest in the matter). So the Lorentz matrix (defined by two 3 dimensional vectors (w, r) related to (\vec{r}, \vec{v})) only is significant from this point of view, and w can be clearly (even if it is in a complicated way) related to \vec{v} . Moreover we have seen that the formulas hold for any observer.

Our purpose here is to find a way to represent kinematic characteristics of material bodies, by vectorial quantities. This is not to find the formulas in a change of coordinates (that we have), so the Poincaré's group is of no use. But the Lorentz group is essential because it gives the rules for the transformation of the components of vectors.

Momenta in the fiber bundle representation

As we have seen fiber bundles provide an efficient representation of the geometry of the Universe, notably in the GR picture. So it is legitimate to look at what it can provide on this issue. The aim is to represent the kinematic characteristics of material bodies, incorporating both their geometric motion (translation and rotation) and inertial features, in a single quantity, which can be implemented for particles, that is without the need to involve a fixed structure of the body. This quantity should be intrinsic, meaning that its value, measured in a frame attached to the particle, should stay constant along its world line, in accordance with the assumption that the motion is absolute and the mass is constant. Meanwhile its value, as measured by an observer, would change according to the observer, as a geometric quantity (with the same gauge group). The gauge group should be the spin group $Spin(3,1)$ (or $Spin(1,3)$, they are isomorphic, we will denote both Spin), as we noticed before.

Let us denote the principal bundle $P_G(M, Spin, \pi_G)$, the standard fiber a manifold E , and γ the action of the Spin group. Then the quantity would be an element \mathbf{S} of $P_G[E, \gamma]$.

\mathbf{S} , for a given particle, should stay constant on its world line. The world line is some map $p : \mathbb{R} \rightarrow M :: m = p(\tau)$ and the trajectory for an observer $q : \mathbb{R} \rightarrow M :: m = q(t) = p(\tau(t))$.

If S represents an intrinsic kinematic characteristic of the particle, it should stay constant along its worldline $p(\tau)$: there is a privileged frame $g(\tau) \in Spin$ such that $\mathbf{S} = (\varphi_G(p(\tau), g(\tau)), S)$ with $S = Ct$

The value measured by the observer in his gauge $\varphi_G(q(t), \mathbf{1})$ is then :

$$(\varphi_G(q(t), \mathbf{1}), S(t))$$

$$\sim \varphi_G(q(t), g(\tau(t))), \gamma(g^{-1}, S(t)) = (\varphi_G(p(\tau), g(\tau)), S)$$

which is equivalent to say that $S(t) = \gamma(g(\tau(t)), S)$ with S constant.

The relation $S(t) = \gamma(g(\tau(t)), S)$ is just the consequence of our very general assumptions. The measure of $S(t)$ varies locally according to the observer (the measure of motion is relative) but its intrinsic value does not change (motion is absolute). But to get a full profit of this representation we have to adopt an entirely new point of view. We cannot any longer view the particle as living in M and spinning in its tangent space. Actually the particle lives in E , which happens to be associated to P_G . Its trajectory is a curve in E , which projects on a curve in M . E can be seen as the physical world (at least a part of it), that we can represent through networks of frames in M . So $S(t)$ cannot be seen properly as a motion, it is only a characteristic of the particle (such as mass and charge). Experience shows that it can be measured through geometric frames but this does not imply the existence of a real spinning motion of the particle. In some

way this is what physicists do, intuitively, in Galilean geometry : the rotation, the rotational moment, are not represented as elements of the group or the Lie algebra, but as vectors (it happens that it is the same vector space as M , but this is fortuitous).

The issue that we face is then to precise E and γ . We will make the following, reasonable, assumptions :

i) E is some vector space, so that we have an associated vector bundle $P_G[E, \gamma]$

ii) it implies that (E, γ) is a representation of Spin.

iii) this representation should be finite dimensional (we consider here the value of $S(t)$ at some point).

iv) if (E, γ) is a representation of Spin, then $(E, \gamma'(1))$ is a representation of its Lie algebra, both are subsets of the Clifford algebra $Cl(3, 1)$ so that, if γ is a linear map, then $\gamma'(1) = \gamma$ and this is not a big leap forward to assume that (E, γ) is a representation of the Clifford algebra itself.

Then the quantity S is a vector of E , called a **spinor**. Spinors have been first introduced by Dirac in an equation, obtained by some magic, but truly pure intuition, which is still the basis of most Quantum Electrodynamics and Quantum Theory of Fields.

4.2 CLIFFORD ALGEBRAS AND SPINORS

Spinors, as well as the spin, cannot be properly understood without a look at their mathematical berth, which is Clifford algebra. This is a fascinating algebraic structure on vector spaces which is seen in details in Maths.Part 9. The results which will be used in this book are summarized in this section, the proofs are given in the Annex.

4.2.1 Clifford algebra and Spin groups

Clifford Algebras

A Clifford algebra $Cl(F, \langle \rangle)$ is an algebraic structure, which can be defined on any vector space $(F, \langle \rangle)$ on a field K (\mathbb{R} or \mathbb{C}) endowed with a bilinear *symmetric* form $\langle \rangle$. The set $Cl(F, \langle \rangle)$ is defined from K and F and a product, denoted \cdot , with the property that for any two *vectors* u, v :

$$\forall u, v \in F : u \cdot v + v \cdot u = 2 \langle u, v \rangle \quad (4.2)$$

A Clifford algebra is then a set which is larger than F : it includes *all vectors* of F , plus scalars, and any linear combinations of products of vectors of F . A Clifford algebra on a n dimensional vector space is a 2^n dimensional vector space on K , and an algebra with \cdot . Clifford algebras built on vector spaces on the same field, with same dimension and bilinear form with same signature are isomorphic. On a 4 dimensional real vector space $(F, \langle \rangle)$ endowed with a Lorentz metric there are two structures of Clifford Algebra, denoted $Cl(3, 1)$ and $Cl(1, 3)$, depending on the signature of the metric, and they are not isomorphic. In the following we will state the results for $Cl(3, 1)$, and for $Cl(1, 3)$ only when they are different.

The easiest way to work with a Clifford algebra is to use an orthonormal basis of F . On any 4 dimensional real vector space $(F, \langle \rangle)$ with a bilinear symmetric form of signature (3,1) or (1,3) we will denote:

Notation 56 $(\varepsilon_i)_{i=0}^3$ is an orthonormal basis with scalar product: $\langle \varepsilon_i, \varepsilon_i \rangle = \eta_{ii}$

So we have the relation:

$$\varepsilon_i \cdot \varepsilon_j + \varepsilon_j \cdot \varepsilon_i = 2\eta_{ij} \quad (4.3)$$

Then a basis of the Clifford algebra is a set comprised of 1 and all ordered products of $\varepsilon_i, i = 0..3$.

In any orthonormal basis there is a fourth vector which is such that $\varepsilon_i \cdot \varepsilon_i = -1$ (for the signature (3,1)) or $+1$ (for the signature (1,3)). We will label this vector ε_0 .

Spin group

The group $Pin(3,1)$ is the subset of the Clifford algebra $Cl(3,1)$:

$$Pin(3,1) = \{u_1 \cdot u_2 \dots \cdot u_k, \langle u_p, u_p \rangle = \pm 1, u_p \in F\}. Pin(3,1) \text{ is a Lie group,}$$

Spin(3,1) is its subgroup where we have an even number of vectors :

$$Spin(3,1) = \{u_1 \cdot u_2 \dots \cdot u_{2k}, \langle u_p, u_p \rangle = \pm 1, u_p \in F\}$$

and similarly for $Pin(1,3)$ and $Spin(1,3)$.

Notice that the *scalars* ± 1 belong to the groups. The identity element is the scalar 1.

$Pin(3,1)$ and $Pin(1,3)$ are not isomorphic. $Spin(3,1)$ and $Spin(1,3)$ are isomorphic.

Adjoint map

For any $s \in Pin(3,1)$, the map, called the **adjoint map** :

$$\mathbf{Ad}_s : Cl(3,1) \rightarrow Cl(3,1) :: \mathbf{Ad}_s w = s \cdot w \cdot s^{-1} \quad (4.4)$$

is such that

$$\forall w \in F : \mathbf{Ad}_s w \in F \quad (4.5)$$

and it preserves the scalar product on F :

$$\forall u, v \in F, s \in Pin(3,1) : \langle \mathbf{Ad}_s u, \mathbf{Ad}_s v \rangle_E = \langle u, v \rangle_E \quad (4.6)$$

Moreover :

$$\forall s, s' \in Pin(3,1) : \mathbf{Ad}_s \circ \mathbf{Ad}_{s'} = \mathbf{Ad}_{s \cdot s'} \quad (4.7)$$

Because the action \mathbf{Ad}_s of $Spin(3,1)$ on F gives another vector of F and preserves the scalar product, it can be represented by a 4×4 orthogonal matrix. Using any orthonormal basis $(\varepsilon_i)_{i=0}^3$ of F , then \mathbf{Ad}_s is represented by a matrix $\Pi(\mathbf{Ad}_s) = [h(s)] \in SO(3,1)$. To two elements $\pm s \in Spin(3,1)$ correspond a single matrix $[h(s)]$. $Spin(3,1)$ is the double cover (as manifold) of $SO(3,1)$. $Spin(3,1)$ has two connected components (which contains either +1 or -1) and its connected component, that we will denote for brevity also $Spin(3,1)$, is simply connected and is the universal cover group of $SO_0(3,1)$. So with the Spin group one can define two physical rotations, corresponding to opposite signs.

Lie algebra of the Spin group

Theorem 57 *The elements of the Lie algebra $T_1 Spin(3,1)$ belong to the Clifford algebra and can be written as the linear combination of elements $\varepsilon_i \cdot \varepsilon_j$:*

As any algebra $Cl(F, \langle \rangle)$ is a Lie algebra with the bracket :

$$\forall w, w' \in Cl(F, \langle \rangle) : [w, w'] = w \cdot w' - w' \cdot w$$

and the Lie algebra $T_1 Spin(3,1)$ of $Spin(3,1)$ is a subset of $Cl(3,1)$ (Maths.532).

The derivative $\Pi'(1) : T_1 Spin(3,1) \rightarrow so(3,1)$ is an isomorphism of Lie algebras. The inverse map : $\Pi'(1)^{-1} : so(3,1) \rightarrow T_1 Spin(3,1)$ is an isomorphism

of Lie algebras which reads (Maths.534) with any orthonormal basis $(\varepsilon_i)_{i=0}^3$ of F :

$$\Pi'(1)^{-1} : so(3,1) \rightarrow T_1Spin(3,1) :: \Pi'(1)^{-1}([\kappa]) = \frac{1}{4} \sum_{i,j=0}^3 ([\kappa] [\eta])_j^i \varepsilon_i \cdot \varepsilon_j$$

and any element of $T_1Spin(3,1)$ is such expressed in the basis of $Cl(F, \langle \rangle)$: it is the linear combinations of the ordered products of all the four vectors of a basis.

With any orthonormal basis and the following choices of basis $(\vec{\kappa}_a)_{a=1}^6$ of $T_1Spin(3,1)$ then $\Pi'(1)^{-1}$ takes a simple form :

$$\Pi'(1)^{-1}([\kappa_1]) = \vec{\kappa}_1 = \frac{1}{2} \varepsilon_3 \cdot \varepsilon_2,$$

$$\Pi'(1)^{-1}([\kappa_2]) = \vec{\kappa}_2 = \frac{1}{2} \varepsilon_1 \cdot \varepsilon_3,$$

$$\Pi'(1)^{-1}([\kappa_3]) = \vec{\kappa}_3 = \frac{1}{2} \varepsilon_2 \cdot \varepsilon_1,$$

$$\Pi'(1)^{-1}([\kappa_4]) = \vec{\kappa}_4 = \frac{1}{2} \varepsilon_0 \cdot \varepsilon_1,$$

$$\Pi'(1)^{-1}([\kappa_5]) = \vec{\kappa}_5 = \frac{1}{2} \varepsilon_0 \cdot \varepsilon_2,$$

$$\Pi'(1)^{-1}([\kappa_6]) = \vec{\kappa}_6 = \frac{1}{2} \varepsilon_0 \cdot \varepsilon_3$$

where $([\kappa_a])_{a=1}^6$ is the basis of $so(3,1)$ already noticed such that : $[\kappa] = K(w) + J(r) = \sum_{a=1}^3 r^a [\kappa_a] + w^a [\kappa_{a+3}]$
 $a = 1, 2, 3 : \vec{\kappa}_a = -\frac{1}{2} \varepsilon(a, i, j) \varepsilon_i \cdot \varepsilon_j, a = 4, 5, 6 : \vec{\kappa}_a = \frac{1}{2} \varepsilon_0 \cdot \varepsilon_{a-3}$ is a basis of $T_1Spin(3,1)$

We will use extensively the convenient (the order of the indices matters) :

Notation 58 for both $Cl(3,1), Cl(1,3)$:

$$v(r, w) = \frac{1}{2} (w^1 \varepsilon_0 \cdot \varepsilon_1 + w^2 \varepsilon_0 \cdot \varepsilon_2 + w^3 \varepsilon_0 \cdot \varepsilon_3 + r^3 \varepsilon_2 \cdot \varepsilon_1 + r^2 \varepsilon_1 \cdot \varepsilon_3 + r^1 \varepsilon_3 \cdot \varepsilon_2) \quad (4.8)$$

With this notation, whatever the orthonormal basis $(\varepsilon_i)_{i=0}^3$, any element X of the Lie algebras $T_1Spin(3,1)$ or $T_1Spin(1,3)$ reads :

$$X = v(r, w) = \sum_{a=1}^3 r^a \vec{\kappa}_a + w^a \vec{\kappa}_{a+3} \quad (4.9)$$

with $(r, w) \in \mathbb{R}^3 \times \mathbb{R}^3$ then $X = v(r, w)$ is the image of :

$$\Pi'(1)(v(r, w)) = K(w) + J(r) \in so(3,1) \text{ if } X \in T_1Spin(3,1)$$

$$\Pi'(1)(v(r, w)) = -(K(w) + J(r)) \in so(1,3) \text{ if } X \in T_1Spin(1,3)$$

$(Cl(3,1), \mathbf{Ad})$ and $(T_1Spin(3,1), \mathbf{Ad})$ are representations of $Spin(3,1)$. One property that we will use often is the following. A change of orthonormal basis : $\varepsilon_i \rightarrow \tilde{\varepsilon}_i$ can be expressed by an element $s \in Spin(3,1)$: $\varepsilon_i \rightarrow \tilde{\varepsilon}_i = \mathbf{Ad}_s \varepsilon_i$. Then the vectors $v \in F$ stay the same, but their components in the new basis change according to a matrix $[h(s)]$ of $SO(3,1)$:

$$v = \sum_{i=1}^4 v^i \varepsilon_i = \sum_{i=1}^4 \tilde{v}^i \tilde{\varepsilon}_i = \sum_{i=1}^4 \tilde{v}^i \mathbf{Ad}_s \varepsilon_i = \sum_{i=1}^4 \tilde{v}^i [h(s)]_i^j \varepsilon_j$$

The change impacts also the elements of the Lie algebra $T_1Spin(3,1)$:

$$X = \sum_a X_a \vec{\kappa}_a = \sum_a \tilde{X}_a \widetilde{\vec{\kappa}}_a$$

$$\widetilde{\vec{\kappa}}_a = \tilde{\varepsilon}_i \cdot \tilde{\varepsilon}_j = s \cdot \varepsilon_i \cdot s^{-1} \cdot s \cdot \varepsilon_j \cdot s^{-1} = \mathbf{Ad}_s(\varepsilon_i \cdot \varepsilon_j) = \mathbf{Ad}_s(\vec{\kappa}_a)$$

$\tilde{X}_a = \sum_{b=1}^6 [\mathbf{Ad}_s]_a^b X_b$ where $[\mathbf{Ad}_s]$ is a 6×6 matrix (not the same as the 4×4 matrix $h(s)$ of $SO(3, 1)$).

With signature (3,1), with the operator j introduced previously and :

Notation 59 $\varepsilon_5 = \varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3$

$$\begin{aligned} \varepsilon_5 \cdot \varepsilon_5 &= -1 \\ v(r', w') \cdot v(r, w) &= \frac{1}{4}(w^t w' - r^t r') \\ &+ \frac{1}{2}v(-j(r)r' + j(w)w', -j(w)r' - j(r)w') - \frac{1}{4}(w^t r' + r^t w')\varepsilon_5 \end{aligned}$$

the bracket on the Lie algebra:

$$\begin{aligned} [v(r, w), v(r', w')] &= v(r, w) \cdot v(r', w') - v(r', w') \cdot v(r, w) \\ [v(r, w), v(r', w')] &= v(j(r)r' - j(w)w', j(w)r' + j(r)w') \\ \varepsilon_5 \cdot v(r, w) &= v(r, w) \cdot \varepsilon_5 = v(r, -w) \end{aligned}$$

With signature (1,3) :

$$\begin{aligned} v(r, w) \cdot v(r', w') &= \frac{1}{4}(w^t w' - r^t r') - \frac{1}{2}v(-j(r)r' + j(w)w', j(w)r' + j(r)w') \\ &- \frac{1}{4}(w^t r' + r^t w')\varepsilon_5 \\ [v(r, w), v(r', w')] &= -v(j(r)r' - j(w)w', j(w)r' + j(r)w') \\ \varepsilon_5 \cdot v(r, w) &= v(r, w) \cdot \varepsilon_5 = v(w, r) \end{aligned}$$

Expression of elements of the spin group

Theorem 60 *The elements of the Spin groups read (see annex) in both signatures :*

$$s = a + v(r, w) + b\varepsilon_5 \quad (4.10)$$

$a, (w^j, r^j)_{j=1}^3, b$ are real scalar which are related. We have the necessary identities :

$$a^2 - b^2 = 1 + \frac{1}{4}(w^t w - r^t r) \quad (4.11)$$

$$ab = -\frac{1}{4}r^t w \quad (4.12)$$

The inverse is :

$$(a + v(r, w) + b\varepsilon_5)^{-1} = a - v(r, w) + b\varepsilon_5 \quad (4.13)$$

The exponential is not surjective on $so(3, 1)$ or $T_1 Spin(3, 1)$: for each $v(r, w) \in T_1 Spin(3, 1)$ there are two elements $\pm \exp v(r, w) \in Spin(3, 1)$:

$\exp t v(R, W) = \pm \sigma_w(t) \cdot \sigma_r(t)$ with opposite sign :

$$\begin{aligned} \sigma_w(t) &= \sqrt{1 + \frac{1}{4}W^t W \sinh^2 \frac{1}{2}t\sqrt{W^t W}} + \sinh \frac{1}{2}t\sqrt{W^t W}v(0, W) \\ \sigma_r(t) &= \sqrt{1 - \frac{1}{4}R^t R \sin^2 t\frac{1}{2}\sqrt{R^t R}} + \sin t\frac{1}{2}\sqrt{R^t R}v(R, 0) \end{aligned}$$

And we have the identity (Maths.1768) :

$$\forall v(r, w) \in T_1 Spin(3, 1) : \quad (4.14)$$

$$\exp \mathbf{Ad}_g v(r, w) = \mathbf{Ad}_g \exp v(r, w) = g \cdot \exp v(r, w) \cdot g^{-1} \quad (4.15)$$

The product s.s' reads :

$$(a + v(r, w) + b\varepsilon_5) \cdot (a' + v(r', w') + b'\varepsilon_5) = a'' + v(r'', w'') + b''\varepsilon_5$$

with :

$$a'' = aa' - b'b + \frac{1}{4}(w^t w' - r^t r')$$

$$b'' = ab' + ba' - \frac{1}{4}(w^t r' + r^t w')$$

and in $Spin(3, 1)$:

$$r'' = \frac{1}{2}(j(r)r' - j(w)w') + a'r + ar' - b'w - bw'$$

$$w'' = \frac{1}{2}(j(w)r' + j(r)w') + a'w + aw' + b'r + br'$$

and in $Spin(1, 3)$:

$$r'' = \frac{1}{2}(j(r)r' - j(w)w') + a'r + ar' + b'w + bw'$$

$$w'' = -\frac{1}{2}(j(w)r' + j(r)w') + a'w + aw' + b'r + br'$$

4.2.2 Scalar product and Norm

Scalar product on the Clifford algebra

There is a scalar product on $Cl(F, \langle \rangle)$ defined by :

$$\langle u_{i_1} \cdot u_{i_2} \cdot \dots \cdot u_{i_n}, v_{j_1} \cdot v_{j_2} \cdot \dots \cdot v_{j_n} \rangle = \langle u_{i_1}, v_{j_1} \rangle \langle u_{i_2}, v_{j_2} \rangle \dots \langle u_{i_n}, v_{j_n} \rangle$$

It does not depend on the choice of a basis, and any orthonormal basis defined as above is orthonormal :

$$\langle \varepsilon_{i_1} \cdot \varepsilon_{i_2} \cdot \dots \cdot \varepsilon_{i_n}, \varepsilon_{j_1} \cdot \varepsilon_{j_2} \cdot \dots \cdot \varepsilon_{j_n} \rangle = \eta_{i_1 j_1} \dots \eta_{i_n j_n} \epsilon(i_1, \dots, i_n, j_1, \dots, j_n) \text{ the latter term is the signature of the permutation } (i_1, \dots, i_n, j_1, \dots, j_n)$$

This scalar product on $Cl(3, 1), Cl(3, 1)$ has the signature (8, 8) : it is non degenerate but neither definite positive or negative. It is invariant by \mathbf{Ad} .

$$\forall w, w' \in Cl(F, \langle \rangle) : \langle \mathbf{Ad}_s w, \mathbf{Ad}_s w' \rangle_{Cl(E, \langle \rangle)} = \langle w, w' \rangle_{Cl(E, \langle \rangle)} \quad (4.16)$$

$(Cl(3, 1), \mathbf{Ad})$ is a unitary representation of $Spin(3, 1)$ and $(Cl(1, 3), \mathbf{Ad})$ a unitary representation of $Spin(1, 3)$.

It reads for elements of $T_1 Spin(3, 1)$:

$$\langle v(r, w), v(r', w') \rangle_{Cl} = \frac{1}{4}(r^t r' - w^t w') \quad (4.17)$$

Norm on the Lie algebra

Theorem 61 *There is a norm on the Lie algebra $T_1 Spin(3, 1)$ which is preserved by the adjoint map*

i) There is a bilinear symmetric form on $T_1Spin(3, 1)$, called the Killing form B which, in the basis above, has the same expression as in $so(3, 1)$ (Maths.1669) :

$$B(v(r, w), v(r', w')) = 4(r^t r' - w^t w') = 16 \langle v(r, w), v(r', w') \rangle_{Cl} \quad (4.18)$$

It is invariant by **Ad** as the scalar product.

ii) The Lie algebras $so(3, 1), T_1Spin(3, 1)$ are semi-simple. They have a Cartan involution, that is a map : $\theta : T_1Spin(3, 1) \rightarrow T_1Spin(3, 1)$ such that : $\theta^2 = Id$ and the bilinear form $B_\theta(X, Y) = -B(X, \theta Y)$ is definite positive.

θ has two eigen values ± 1 , by taking L_0 the eigen space for -1 and P_0 the eigen space for +1, the Lie algebra is the sum of two, orthogonal (for B so for $\langle \rangle_{Cl}$), vector subspaces (called a Cartan decomposition Maths.1742) such that :

$$T_1Spin(3, 1) = L_0 \oplus P_0$$

$$L_0 \text{ is a Lie subalgebra, } [L_0, L_0] \subset L_0, [L_0, P_0] \subset P_0, [P_0, P_0] \subset L_0$$

The Killing form and this decomposition do not depend on the choice of a basis in $T_1Spin(3, 1)$ (these are general properties of Lie algebras).

The Cartan involutions are defined in $T_1Spin(3, 1)$ by :

$$\theta(\varepsilon_0)(X) = \varepsilon_0 \cdot X \cdot \varepsilon_0$$

with any vector $\varepsilon_0 : \varepsilon_0 \cdot \varepsilon_0 = -1$. We have one involution for each orthonormal basis.

It is easy to see that for the basis built with ε_0 :

$$\forall a = 1, 2, 3 : \varepsilon_0 \cdot \vec{\kappa}_a \cdot \varepsilon_0 = -\vec{\kappa}_a$$

$$\forall a = 4, 5, 6 : \varepsilon_0 \cdot \vec{\kappa}_a \cdot \varepsilon_0 = \vec{\kappa}_a$$

$$\theta(\varepsilon_0)v(r, w) = v(-r, w)$$

$$B(v(r, w), \theta(\varepsilon_0)v(r, w)) = -4(w^t w + r^t r')$$

Thus :

$$L_0 = \{v(r, 0), r \in \mathbb{R}^3\} = \{X \in T_1Spin(3, 1) : \theta(\varepsilon_0)(X) = -X\}$$

$$P_0 = \{v(0, w), w \in \mathbb{R}^3\} = \{X \in T_1Spin(3, 1) : \theta(\varepsilon_0)(X) = X\}$$

iii) Let us check that the decomposition does not depend on the choice of a basis. In a change of basis on $T_1Spin(3, 1)$:

$$\varepsilon_i \rightarrow \tilde{\varepsilon}_i = Ad_s \varepsilon_i$$

$$\vec{\kappa}_a \rightarrow \widetilde{\vec{\kappa}}_a = Ad_s \vec{\kappa}_a$$

$$\theta(\tilde{\varepsilon}_0)\left(\widetilde{\vec{\kappa}}_a\right) = Ad_s \varepsilon_0 \cdot Ad_s \vec{\kappa}_a \cdot Ad_s \varepsilon_0 = Ad_s(\varepsilon_0 \cdot \vec{\kappa}_a \cdot \varepsilon_0) = Ad_s(\theta(\varepsilon_0)(\vec{\kappa}_a))$$

So :

$$\forall a = 1, 2, 3 : \theta(\tilde{\varepsilon}_0)\left(\widetilde{\vec{\kappa}}_a\right) = -Ad_s(\vec{\kappa}_a) = -\widetilde{\vec{\kappa}}_a$$

$$\forall a = 4, 5, 6 : \theta(\tilde{\varepsilon}_0)\left(\widetilde{\vec{\kappa}}_a\right) = Ad_s(\vec{\kappa}_a) = \widetilde{\vec{\kappa}}_a$$

$X \in L_0$:

$$X = \sum_{a=1}^3 X^a \vec{\kappa}_a = \sum_{a=1}^3 \tilde{X}^a \widetilde{\vec{\kappa}}_a \Rightarrow \theta(\tilde{\varepsilon}_0)(X) = -\sum_{a=1}^3 \tilde{X}^a \widetilde{\vec{\kappa}}_a = -X$$

$X \in P_0$:

$$X = \sum_{a=4}^6 X^a \vec{\kappa}_a = \sum_{a=4}^6 \tilde{X}^a \widetilde{\vec{\kappa}}_a \Rightarrow \theta(\tilde{\varepsilon}_0)(X) = \sum_{a=4}^6 \tilde{X}^a \widetilde{\vec{\kappa}}_a = X$$

iv) Because the vector spaces L_0, P_0 do not depend on the choice of an orthonormal basis we can define uniquely the projections :

$$\begin{aligned}\pi_L : T_1 Spin(3, 1) &\rightarrow L_0 :: \pi_L(X) = \frac{1}{2}(X - \theta(\varepsilon_0)(X)) = \frac{1}{2}(X - \varepsilon_0 \cdot X \cdot \varepsilon_0) \\ \pi_P : T_1 Spin(3, 1) &\rightarrow P_0 :: \pi_P(X) = \frac{1}{2}(X + \theta(\varepsilon_0)(X)) = \frac{1}{2}(X + \varepsilon_0 \cdot X \cdot \varepsilon_0) \\ X &= \pi_L(X) + \pi_P(X)\end{aligned}$$

L_0, P_0 are orthogonal for the scalar product, thus :

$$\langle X, X \rangle_{Cl} = \langle \pi_L(X), \pi_L(X) \rangle_{Cl} + \langle \pi_P(X), \pi_P(X) \rangle_{Cl}$$

The bilinear form :

$$\begin{aligned}\langle X, \pi_L X - \pi_P X \rangle_{Cl} &= \langle \pi_L(X) + \pi_P(X), \pi_L X - \pi_P X \rangle_{Cl} \\ &= \langle \pi_L(X), \pi_L(X) \rangle_{Cl} - \langle \pi_P(X), \pi_P(X) \rangle_{Cl}\end{aligned}$$

is definite positive and invariant by the adjoint map.

$$\|X\| = \sqrt{\langle X, \pi_L X - \pi_P X \rangle_{Cl}} \text{ is a norm on } T_1 Spin(3, 1)$$

$$\begin{aligned}\|X + X'\|^2 &= \langle X + X', \pi_L(X + X') - \pi_P(X + X') \rangle_{Cl} \\ &= \langle \pi_L(X + X'), \pi_L(X + X') \rangle_{Cl} - \langle \pi_P(X + X'), \pi_P(X + X') \rangle_{Cl} \\ \langle \pi_L(X + X'), \pi_L(X + X') \rangle_{Cl} &\leq \langle \pi_L(X), \pi_L(X) \rangle_{Cl} + \langle \pi_L(X'), \pi_L(X') \rangle_{Cl} \\ - \langle \pi_P(X + X'), \pi_P(X + X') \rangle_{Cl} &\leq - \langle \pi_P(X), \pi_P(X) \rangle_{Cl} - \langle \pi_P(X'), \pi_P(X') \rangle_{Cl}\end{aligned}$$

It reads :

$$\|v(r, w)\| = \frac{1}{2} \sqrt{r^t r + w^t w} = \frac{1}{2} \sqrt{\langle \pi_L X - \pi_P X, \pi_L X - \pi_P X \rangle_{Cl}} \quad (4.19)$$

and does not depend on a basis on $T_1 Spin(3, 1)$ and is invariant by the adjoint map.

4.2.3 Symmetry breakdown

Clifford algebra $Cl(3)$

The elements of $SO(3, 1)$ are the product of spatial rotations (represented by $\exp J(r)$) and boosts, linked to the speed and represented by $\exp K(w)$. We have similarly a decomposition of the elements of $Spin(3, 1)$. But to understand this topic, from both a mathematical and a physical point of view, we need to distinguish the abstract algebraic structure and the sets on which the structures have been defined.

From a vector space $(F, \langle \rangle)$ one can built only one Clifford algebra with the structure $Cl(3, 1)$: as a set $Cl(3, 1)$ must comprise all the vectors of F. But from any vector subspace of F one can built Clifford algebras : their algebraic structure depends on the dimension of the vector space, and on the signature of the metric induced on the vector subspace. To have a Clifford algebra structure $Cl(3)$ on F one needs a 3 dimensional vector subspace on which the scalar product is definite positive, so it cannot include any vector such that $\langle u, u \rangle < 0$ (and conversely for the signature $(1, 3)$: the scalar product must be definite negative). The subsets of F which are a 3 dimensional vector subspace and do not contain any vector such that $\langle u, u \rangle < 0$ are no unique. So we have different subsets of $Cl(3, 1)$ with the structure of a Clifford algebra $Cl(3)$, all isomorphic but which do not contain the same vectors. Because the Spin Groups are built from elements of the Clifford algebra, we have similarly isomorphic Spin groups $Spin(3)$, but with different elements.

The simplest way to deal with these issues is to fix an orthonormal basis. Any orthonormal basis of F contains one vector such that $\langle \varepsilon_i, \varepsilon_i \rangle = -1$ (or $+1$ with the signature $(1, 3)$). If we exclude this vector we can generate a vector subspace and then a Clifford algebra $Cl(3)$. So the identification of a specific set with the structure of $Cl(3)$ sums up to single out such a vector, that we will denote as ε_0 .

Decomposition of the elements of the Spin group

Theorem 62 *For a given vector ε_0 any element $s \in Spin(3, 1)$ can be written uniquely (up to sign) : $s = \epsilon (a_w + v(0, w)) \cdot \epsilon (a_w + v(r, 0))$ with $\epsilon = \pm 1$*

Let us choose an orthonormal basis $(\varepsilon_i)_{i=0}^3$. Then there is a unique vector subspace F^\perp orthogonal to ε_0 , where the scalar product is definite positive, and from $(F^\perp, \langle \rangle)$ one can build a Clifford algebra with structure $Cl(3)$. Its spin group has the structure $Spin(3)$ which has for Lie algebra $T_1 Spin(3)$.

In the basis :

an element of $T_1 Spin(3)$ reads : $X = v(r, 0) = \sum_{a=1}^3 r^a \overrightarrow{\kappa}_a$ and is associated to a 3×3 matrix $j(r)$

any element of $Spin(3)$ reads : $s_r = a_r + v(r, 0)$ with $a_r^2 = 1 - \frac{1}{4}r^t r$

$Spin(3)$ has 2 connected components. The connected component of the identity is comprised of elements with $a_r > 0$.

$Spin(3), T_1 Spin(3)$ are subsets of $Cl(3, 1)$: they are comprised of specific combinations of vectors of F .

The **quotient space** $Spin(3, 1)/Spin(3)$ is not a group but a 3 dimensional manifold, called a homogeneous space (Maths.22.4.3). It is characterized by the equivalence relation :

$$\forall s, s' \in Spin(3, 1) : s \sim s' \Leftrightarrow \exists s_r \in Spin(3) : s' = s \cdot s_r$$

The projection : $\pi_w : Spin(3, 1) \rightarrow Spin(3, 1)/Spin(3)$ is a submersion, its derivative $\pi'_w(s)$ is surjective.

In each class of $Spin(3, 1)/Spin(3)$ there are only two elements (which belong to $Spin(3, 1)$) which can be written as : $s_w = a_w + v(0, w)$ and they have opposite sign : $\pm s_w$ belong to the same class (see Annex for the proofs).

Any element $s \in Spin(3, 1)$ can be written : $s = s_w \cdot s_r$ with $s_w \in Spin(3, 1)/Spin(3), s_r \in Spin(3)$:

$$s = \epsilon (a_w + v(0, w)) \cdot \epsilon (a_w + v(r, 0)) \quad \text{with } \epsilon = \pm 1 \quad (4.20)$$

For each s there are two decompositions, unique up to sign. $\pm s_w$ belong to the class of equivalence of s in $Spin(3, 1)/Spin(3)$. They are specific representatives of the projection of s on the homogeneous space. *The decomposition depends on the choice of ε_0 .*

Theorem 63 *$Spin(3, 1)$ is a principal fiber bundle $Spin(3, 1)(Spin(3, 1)/Spin(3), Spin(3), \pi_w)$.*

and there is a left action of $Spin(3, 1)$ on $Spin(3, 1)/Spin(3)$:

$\lambda : Spin(3, 1) \times Spin(3, 1)/Spin(3) \rightarrow Spin(3, 1)/Spin(3)$:

$\lambda(s, s_w) = \pi_w(s \cdot s_w)$

Proof. Indeed we have :

$\lambda(s, s_w) = \pi_w(s \cdot s_w) \Leftrightarrow \exists s_r : s \cdot s_w = \pi_w(s \cdot s_w) \cdot s_r \Leftrightarrow \lambda(s, s_w) = s \cdot s_w \cdot s_r^{-1}$

$\lambda(s, \lambda(s', s_w)) = \lambda(s, s' \cdot s_w \cdot s_r'^{-1}) = s \cdot s' \cdot s_w \cdot s_r'^{-1} \cdot s_r^{-1}$

$\lambda(s \cdot s', s_w) = s \cdot s' \cdot s \cdot s_w \cdot s_r''$

$s \cdot s' \cdot s \cdot s_w \cdot s_r'' \sim s \cdot s' \cdot s_w \cdot s_r'^{-1} \cdot s_r^{-1}$ ■

The subset of $Spin(3, 1)$ of the elements s_r which commute with ε_0 :

$\mathbf{Ad}_{s_r} \varepsilon_0 = s_r \cdot \varepsilon_0 \cdot s_r^{-1} = \varepsilon_0$

read :

$$s_r = \epsilon \left(\sqrt{1 - \frac{1}{4} r^t r} + v(r, 0) \right) \text{ with } \epsilon = \pm 1, r^t r \leq 4 \quad (4.21)$$

They belong to the realization of $Spin(3)$ which is specified by ε_0 . They are generated by vectors belonging to the subspace spanned by the vectors $(\varepsilon_i)_{i=1}^3$. They have a special physical meaning : they are the spatial rotations for an observer *with a velocity in the direction of ε_0 . The vector r , with 3 components, has the same meaning as in Galilean Geometry.* This is just a different representation of the Lie algebra $o(3)$.

The elements of $Spin(3, 1)/Spin(3)$ are coordinated by w , and the matrix $[K(w)]$ corresponds to a gauge transformation for an observer moving with a spatial speed \vec{v} parallel to w , without spatial rotation.

Similarly the elements of the Lie algebra $T_1Spin(3)$ are characterized by

$X \in T_1Spin(3) \Leftrightarrow X \cdot \varepsilon_0 - \varepsilon_0 \cdot X = 0 \Leftrightarrow w = 0$

In the tangent space $T_m M$ of the manifold M all rotations (given by $Spin(3, 1)$) are on the same footing. But, because of our assumptions about the motion of observers (along time like lines), any observer introduces a breakdown of symmetry : some rotations are privileged. Indeed the spatial rotations are special, in that they are the ones for which the axis belongs to the physical space.

To sum up :

$(\varepsilon_i)_{i=1}^3$ are the generators in $Cl(F, \langle \rangle)$ of a set with the structure of $Spin(3)$ iff $\langle \varepsilon_i, \varepsilon_i \rangle = +1$ (then the 4th vector $\langle \varepsilon_0, \varepsilon_0 \rangle = -1$)

Similarly any element of the Lie algebra $T_1Spin(3, 1)$ read:

$X = v(r, w) = \sum_{a=1}^3 r^a \vec{\kappa}_a + w^a \vec{\kappa}_{a+3}$

and $\sum_{a=1}^3 r^a \vec{\kappa}_a$ identifies a specific Lie algebra $T_1Spin(3)$ if $\langle \varepsilon_i, \varepsilon_i \rangle = +1$.

In the following of this book we will always assume that the orthonormal basis is such that ε_0 is the 4th vector : $\langle \varepsilon_0, \varepsilon_0 \rangle = -1$ with signature (3,1) and $\langle \varepsilon_0, \varepsilon_0 \rangle = +1$ with signature (1,3).

4.2.4 Representation of Clifford algebras

Complexification of real Clifford algebras

A real Clifford algebra can be complexified, using the same procedure as for any finite dimensional vector space. The complexified of both $Cl(3, 1)$ and $Cl(1, 3)$ is $Cl(\mathbb{C}, 4)$, the Clifford algebra on \mathbb{C}^4 with the bilinear *symmetric* form of signature $(+ + + +)$.

There are real algebra morphisms (injective but not surjective) from the real Clifford algebras to $Cl(\mathbb{C}, 4)$.

Let $\tilde{C} : (\mathbb{R}^4, \langle \cdot, \cdot \rangle_L) \rightarrow Cl(\mathbb{C}, 4)$ be the real linear map defined by : $\tilde{C}(\varepsilon_j) = \varepsilon_j, j = 1, 2, 3; \tilde{C}(\varepsilon_0) = i\varepsilon_0$ and the scalar product $\langle \cdot, \cdot \rangle_L$ with signature $(- + + +)$.

It is easy to check that : $\tilde{C}(u) \cdot \tilde{C}(v) + \tilde{C}(v) \cdot \tilde{C}(u) = 2\langle u, v \rangle_L$ so, by the universal property of Clifford algebras, there is a unique real algebra morphism $C : Cl(3, 1) \rightarrow Cl(\mathbb{C}, 4)$ such that $\tilde{C} = C \circ j$ where j is the canonical injection $(\mathbb{R}^4, \langle \cdot, \cdot \rangle_L) \rightarrow Cl(3, 1)$. We will denote for simplicity $\tilde{C} = C$. The image $C(Cl(3, 1))$ is a real subalgebra of $Cl(\mathbb{C}, 4)$.

Similarly with $\tilde{C}'(\varepsilon_j) = i\varepsilon_j, j = 1, 2, 3; \tilde{C}'(\varepsilon_0) = \varepsilon_0$ we have a real algebra morphism $C' : Cl(1, 3) \rightarrow Cl(\mathbb{C}, 4)$ and $C'(Cl(1, 3))$ is a real subalgebra of $Cl(\mathbb{C}, 4)$. Moreover $C'(\varepsilon_j) = -i\eta_{jj}C(\varepsilon_j)$ (η always correspond to the signature $- + + +$).

Algebraic and geometric representations

An *algebraic* representation of a Clifford algebra is a map γ which associates to each element w of the Clifford algebra a matrix $[\gamma(w)]$ and such that γ is a isomorphism of algebra : all the operations in the Clifford algebra (multiplication by a scalar, sum, Clifford product) are reproduced on the matrices. A representation is fully defined by the family of matrices $(\gamma_i)_{i=0}^3$ representing each vector $(\varepsilon_i)_{i=0}^3$ of an orthonormal basis. The choice of these matrices is not unique : the only condition is that $[\gamma_i][\gamma_j] + [\gamma_j][\gamma_i] = 2\eta_{ij}[I]$ and any family of matrices deduced by conjugation with a fixed matrix gives an equivalent algebraic representation. An element of the Clifford algebra is then represented by a linear combination of generators :

$$\gamma(w) = \gamma\left(\sum_{\{i_1 \dots i_r\}} a^{i_1 \dots i_r} \varepsilon_{i_1} \cdot \dots \cdot \varepsilon_{i_r}\right) = \sum_{\{i_1 \dots i_r\}} a^{i_1 \dots i_r} \gamma_{i_1 \dots i_r}$$

A *geometric* representation (E, γ) of a Clifford algebra is an isomorphism $\gamma : Cl \rightarrow L(E; E)$ in which $[\gamma(w)]$ is the matrix of an endomorphism of E , represented in some basis. From an algebraic representation one can deduce a geometric representation, and they are equivalent up to the choice of a basis.

We look for a geometric representation : the quantity S that we are looking for is represented, not by γ matrices, but by vectors S of the space E , which are called spinors. Higher orders spinors are tensorial products of vectors of E .

A Clifford algebra has, up to isomorphism, a unique faithful algebraic irreducible representation in an algebra of matrices (γ is a bijection). As can be expected the representations depend on the signature :

For $Cl(3, 1)$ this is $\mathbb{R}(4)$ the 4×4 real matrices (the corresponding spinors are the Majorana spinors)

For $Cl(1, 3)$ this is $H(2)$ the 2×2 matrices with quaternionic elements

In both cases an element of the Clifford algebra is characterized by $2^4 = 16$ real parameters.

The geometry of the universe is based upon real structures. Thus we should consider representations of $Cl(3, 1)$ or $Cl(1, 3)$, which raises the issue of the signature. However it happens, from experience, that the vector space E must be complex.

The irreducible representation of $Cl(\mathbb{C}, 4)$ is by 4×4 matrices on complex numbers which must meet the condition : $\gamma_j \gamma_k + \gamma_k \gamma_j = 2\delta_{jk} I_4$.

If (E, γ) is a complex representation of $Cl(\mathbb{C}, 4)$ then $(E, \gamma \circ C)$ is a real geometric representation of $Cl(3, 1)$ on the complex vector space E : the map $\gamma \circ C : Cl(3, 1) \rightarrow L(E; E)$ is a real morphism of algebras, and the maps $\gamma \circ C(w)$ are complex linear. The matrices of the real representation are $i\gamma_0, \gamma_j, j = 1, 2, 3, i\gamma_0$. Similarly $(E, \gamma \circ C')$ is a real geometric representation of $Cl(1, 3)$ with matrices $\gamma_0, i\gamma_j, j = 1, 2, 3$.

Using this trick we see that we are fortunate, in that we have the same representation (E, γ) for both signatures, and a complex vector space E . Moreover it is easy to specify the representation through additional features of E (such as chirality as we will see). A spinor has 8 real components (vs 16 real components for elements of the real Clifford algebras) thus a spinor carries more information than a simple vector of \mathbb{R}^4 and this solves part of the issue of the number of parameters needed to represent the motion (both translation and rotation).

Chirality

Any Clifford algebra Cl is the direct sum of one Clifford subalgebra Cl_0 comprised of elements which are the sum of products of an even number of vectors, and a vector subspace Cl_1 comprised of elements which are the sum of products of an odd number of vectors. Moreover some Clifford algebras present a specific feature : they are the direct sum of two subalgebras which can be seen as algebras of left handed and right handed elements. This property depends on the existence of an element ϖ such that $\varpi \cdot \varpi = 1$. This element exists in any complex algebra, but not in $Cl(1, 3), Cl(3, 1)$. As chirality is a defining feature of particles, this is an additional argument for using $Cl(\mathbb{C}, 4)$.

In this subsection and the next one the vectors $(\varepsilon_j)_{j=0}^3$ denote the canonical basis of \mathbb{C}^4 endowed with the bilinear symmetric form of signature $+++ +$. In $Cl(\mathbb{C}, 4)$ the special element is : $\varpi = \pm \varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3 \in Spin(\mathbb{C}, 4)$. Thus there is a choice and we will use : $\varpi = \varepsilon_5 = \varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3$.

The Clifford algebra splits in two subalgebras :

$$Cl(\mathbb{C}, 4) = Cl^R(\mathbb{C}, 4) \oplus Cl^L(\mathbb{C}, 4) :$$

$$Cl^R(\mathbb{C}, 4) = \{w \in Cl(\mathbb{C}, 4) : \varepsilon_5 \cdot w = w\},$$

$$Cl^L(\mathbb{C}, 4) = \{w \in Cl(\mathbb{C}, 4) : \varepsilon_5 \cdot w = -w\}$$

and any element of $Cl(\mathbb{C}, 4)$ can be uniquely written as : $w = w_R + w_L$

The projections from $Cl(\mathbb{C}, 4)$ on each subalgebra are the maps

$$p_R = \frac{1}{2}(1 + \varepsilon_5), p_L = \frac{1}{2}(1 - \varepsilon_5) :$$

$$w_R = p_R \cdot w, w_L = p_L \cdot w$$

$$p_R \cdot p_L = p_L \cdot p_R = 0, p_R^2 = p_R, p_L^2 = p_L, p_R + p_L = 1$$

We have similarly : $E = E^R \oplus E^L$ with

$$E^R = \gamma_R(E), E^L = \gamma_L(E), \gamma_R = \gamma(p_R), \gamma_L = \gamma(p_L) \Rightarrow \gamma(\varepsilon_5) = \gamma_R - \gamma_L$$

$$u \in E : u = u_R + u_L :$$

$$u_R = \gamma_R(u) = \frac{1}{2}(u + \gamma(\varepsilon_5)u) ;$$

$$u_L = \gamma_L(u) = \frac{1}{2}(u - \gamma(\varepsilon_5)u)$$

For any homogeneous element $w = v_1 \cdot v_2 \dots \cdot v_k, v_j \in \mathbb{C}^4$ we have $\varepsilon_5 \cdot w = (-1)^k w \cdot \varepsilon_5$

$$\forall w \in Cl(\mathbb{C}, 4), u \in E :$$

$$\gamma_R(\gamma(w)u_R) = \frac{1}{2}(1 + (-1)^k)\gamma(w)u_R$$

$$k \text{ even} : \gamma_R(\gamma(w)u_R) = \gamma(w)u_R$$

$$k \text{ odd} : \gamma_R(\gamma(w)u_R) = 0$$

For k even : $\gamma(w)$ preserves both E^R, E^L (as vector subspaces)

For k odd : $\gamma(w)$ exchanges E^R, E^L

In particular the elements of the images $C(Spin(3, 1))$ and $C'(Spin(1, 3))$ by γ preserve both E^R, E^L . So we have reducible representations of these groups.

The choice of the representation γ

An algebraic representation is defined by the choice of its generators γ_i , and any set of generators conjugate by a fixed matrix gives an equivalent representation. We can specify the generators by the choice of a basis $(e_i)_{i=1}^4$ of E . The previous result leads to a natural choice : take $(e_i)_{i=1}^2$ as basis of E^R and $(e_i)_{i=3}^4$ as basis of E^L , then :

$$\gamma_R = \begin{bmatrix} I_2 & 0 \\ 0 & 0 \end{bmatrix}, \gamma_L = \begin{bmatrix} 0 & 0 \\ 0 & I_2 \end{bmatrix}, \gamma_5 = \gamma(\varepsilon_5) = \gamma_R - \gamma_L = \begin{bmatrix} I_2 & 0 \\ 0 & -I_2 \end{bmatrix}$$

Denote : $\gamma_j = \begin{bmatrix} A_j & B_j \\ C_j & D_j \end{bmatrix}$ with four 2×2 complex matrices.

ε_5 belongs to the Spin group $Spin(\mathbb{C}, 4)$, commutes with any element of $Cl_0(\mathbb{C}, 4)$ and anticommutes with any vector, thus $\gamma_5 \gamma_j = -\gamma_j \gamma_5$ which imposes the condition :

$$\begin{bmatrix} A_j & -B_j \\ C_j & -D_j \end{bmatrix} = - \begin{bmatrix} A_j & B_j \\ -C_j & -D_j \end{bmatrix} \Rightarrow \gamma_j = \begin{bmatrix} 0 & B_j \\ C_j & 0 \end{bmatrix}$$

The defining relations : $\gamma_j \gamma_k + \gamma_k \gamma_j = 2\delta_{jk}I_4$ lead to :

$$\begin{bmatrix} B_j C_k + B_k C_j & 0 \\ 0 & C_j B_k + C_k B_j \end{bmatrix} = 2\delta_{jk}I_4$$

$$j \neq k : B_j C_k + B_k C_j = C_j B_k + C_k B_j = 0$$

$$j = k : B_j C_j = C_j B_j = I_2 \Leftrightarrow C_j = B_j^{-1}$$

thus $(\gamma_i)_{i=0}^3$ is fully defined by a set $(B_i)_{i=0}^3$ of 2×2 complex matrices

$$\gamma_j = \begin{bmatrix} 0 & B_j \\ B_j^{-1} & 0 \end{bmatrix}$$

meeting : $j \neq k : B_j B_k^{-1} + B_k B_j^{-1} = B_j^{-1} B_k + B_k^{-1} B_j = 0$

which reads :

$$B_j B_k^{-1} = - (B_j B_k^{-1})^{-1} \Leftrightarrow (B_j B_k^{-1})^2 = -I_2$$

$$B_j^{-1} B_k = - (B_j^{-1} B_k)^{-1} \Leftrightarrow (B_j^{-1} B_k)^2 = -I_2$$

Let us define : $k = 1, 2, 3 : M_k = -i B_k B_0^{-1}$

The matrices $(M_k)_{k=1}^3$ are such that :

$$M_k^2 = - (B_j B_0^{-1})^2 = -I_2$$

$$M_j M_k + M_k M_j = -B_j B_0^{-1} B_k B_0^{-1} - B_k B_0^{-1} B_j B_0^{-1}$$

$$= - (-B_j B_k^{-1} B_0 - B_k B_j^{-1} B_0) B_0^{-1}$$

$$= B_j B_k^{-1} + B_k B_j^{-1} = 0$$

that is $k = 1, 2, 3 : M_j M_k + M_k M_j = 2\delta_{jk} I_2$

Moreover : $\gamma_5 = \gamma_0 \gamma_1 \gamma_2 \gamma_3 \Rightarrow$

$$B_0 B_1^{-1} B_2 B_3^{-1} = I_2$$

$$B_0^{-1} B_1 B_2^{-1} B_3 = -I_2$$

with $B_k = i M_k B_0, B_k^{-1} = -i B_0^{-1} M_k^{-1}$

$$B_0 (-i B_0^{-1} M_1^{-1}) (i M_2 B_0) (-i B_0^{-1} M_3^{-1}) = I_2 = -i M_1^{-1} M_2 M_3^{-1}$$

$$B_0^{-1} (i M_1 B_0) (-i B_0^{-1} M_2^{-1}) (i M_3 B_0) = -I_2 = i B_0^{-1} M_1 M_2^{-1} M_3 B_0$$

which reads :

$$i M_2 = -M_1 M_3 = M_3 M_1$$

$$-M_1^{-1} M_3^{-1} = i M_2^{-1} \Leftrightarrow i M_2 = M_3 M_1$$

$$M_2 M_3 + M_3 M_2 = 0 = i M_1 M_3 M_3 + M_3 M_2 \Leftrightarrow i M_1 = -M_3 M_2 = M_2 M_3$$

$$M_1 M_2 + M_2 M_1 = 0 = i M_3 M_2 M_2 + M_2 M_1 \Rightarrow i M_3 = -M_2 M_1 = M_1 M_2$$

The set of 3 matrices $(M_k)_{k=1}^3$ has the multiplication table :

$$\begin{bmatrix} 1 \setminus 2 & M_1 & M_2 & M_3 \\ M_1 & I & i M_3 & -i M_2 \\ M_2 & -i M_3 & I & i M_1 \\ M_3 & i M_2 & -i M_1 & I \end{bmatrix}$$

which is the same as the set of Pauli's matrices :

$$\sigma_1 = \begin{bmatrix} 0 & 1 \\ 1 & 0 \end{bmatrix}; \sigma_2 = \begin{bmatrix} 0 & -i \\ i & 0 \end{bmatrix}; \sigma_3 = \begin{bmatrix} 1 & 0 \\ 0 & -1 \end{bmatrix}; \sigma_0 = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} \quad (4.22)$$

$$\sigma_i^2 = \sigma_0; \text{ For } j \neq k : \sigma_j \sigma_k = \epsilon(j, k, l) i \sigma_l \quad (4.23)$$

Notation 64 $\epsilon(j, k, l)$ = the signature of the permutation of the three different integers i, j, k , 0 if two integers are equal

There is still some freedom in the choice of the γ_i matrices by the choice of B_0 and the simplest is : $B_0 = -iI_2 \Rightarrow B_k = \sigma_k$

Moreover, because scalars belong to Clifford algebras, one must have the identity matrix I_4 and $\gamma(z) = zI_4$

Thus :

$$\gamma_0 = \begin{bmatrix} 0 & -i\sigma_0 \\ i\sigma_0 & 0 \end{bmatrix}; \gamma_1 = \begin{bmatrix} 0 & \sigma_1 \\ \sigma_1 & 0 \end{bmatrix}; \gamma_2 = \begin{bmatrix} 0 & \sigma_2 \\ \sigma_2 & 0 \end{bmatrix}; \gamma_3 = \begin{bmatrix} 0 & \sigma_3 \\ \sigma_3 & 0 \end{bmatrix}; \quad (4.24)$$

The matrices γ_j are then unitary and Hermitian :

$$\gamma_j = \gamma_j^* = \gamma_j^{-1} \quad (4.25)$$

which is extremely convenient.

We will use the following (see the annex for more formulas) :

Notation 65 $j = 1, 2, 3 : \tilde{\gamma}_j = \begin{bmatrix} \sigma_j & 0 \\ 0 & \sigma_j \end{bmatrix}$

$$j \neq k, l = 1, 2, 3 : \gamma_j \gamma_k = -\gamma_k \gamma_j = i\epsilon(j, k, l) \tilde{\gamma}_l$$

$$j = 1, 2, 3 : \gamma_j \gamma_0 = -\gamma_0 \gamma_j = i \begin{bmatrix} \sigma_j & 0 \\ 0 & -\sigma_j \end{bmatrix} = i\gamma_5 \tilde{\gamma}_j$$

Notice that the choice of the matrices is done in $Cl(\mathbb{C}, 4)$, so it is independent of the choice of signature. However we have the representations of the real algebras by the matrices $\gamma C(\varepsilon_j)$ and $\gamma C'(\varepsilon_j)$

$$Cl(3, 1) : \gamma C(\varepsilon_j) = \gamma_j, j = 1, 2, 3; \gamma C(\varepsilon_0) = i\gamma_0; \gamma C(\varepsilon_5) = i\gamma_5 \quad (4.26)$$

$$Cl(1, 3) : \gamma C'(\varepsilon_j) = i\gamma_j, j = 1, 2, 3; \gamma C'(\varepsilon_0) = \gamma_0; \gamma C'(\varepsilon_5) = \gamma_5 \quad (4.27)$$

The representation that we have chosen here is not unique and others, equivalent, would hold. However from my point of view this is the most convenient because of the nice properties of the γ matrices. The choice of $\varpi = -\varepsilon_5 = -\varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3$ would have lead to take $\tilde{\gamma}_j = -\gamma_j$. In the Standard Model we have a representation of $Cl(1, 3)$ by the matrices : $\tilde{\gamma}_0 = i\gamma_0, \tilde{\gamma}_j = \gamma_j, j = 1, 2, 3$ and $\tilde{\gamma}_5 = -i\tilde{\gamma}_0\tilde{\gamma}_1\tilde{\gamma}_2\tilde{\gamma}_3$

Expression of the matrices for the Lie algebra and the Spin groups

The matrices $\gamma C(v(r, w)), \gamma C'(v(r, w))$ are of constant use.

In $\text{Cl}(\mathbf{R},3,1)$:

$$\begin{aligned} \gamma C(v(r, w)) &= -i\frac{1}{2} \sum_{a=1}^3 (w^a \gamma_a \gamma_0 + r^a \tilde{\gamma}_a) \\ &= \frac{1}{2} \sum_{a=1}^3 \begin{bmatrix} (w_a - ir_a) \sigma_a & 0 \\ 0 & -(w_a + ir_a) \sigma_a \end{bmatrix} \end{aligned} \quad (4.28)$$

In $\text{Cl}(\mathbf{R},1,3)$:

$$\begin{aligned} \gamma C'(v(r, w)) &= -i\frac{1}{2} \sum_{a=1}^3 (w^a \gamma_a \gamma_0 - r^a \tilde{\gamma}_a) \\ &= \frac{1}{2} \sum_{a=1}^3 \begin{bmatrix} (w_a + ir_a) \sigma_a & 0 \\ 0 & -(w_a - ir_a) \sigma_a \end{bmatrix} \end{aligned} \quad (4.29)$$

so one goes from one signature to the other by changing the sign of r .

The 2×2 matrices $\frac{1}{2} \sum_{a=1}^3 (w_a - ir_a) \sigma_a$ and $\frac{1}{2} \sum_{a=1}^3 (w_a + ir_a) \sigma_a$ belong to $SU(2)$

The elements of the spin groups are represented by the matrices :

In $\text{Cl}(\mathbf{R},3,1)$:

$$\gamma C(a + v(r, w) + b\varepsilon_5) = aI - i\frac{1}{2} \sum_{a=1}^3 (w_a \gamma_a \gamma_0 + r_a \tilde{\gamma}_a) + b\gamma_5 \quad (4.30)$$

In $\text{Cl}(\mathbf{R},1,3)$:

$$\gamma C'(a + v(r, w) + b\varepsilon_5) = aI - i\frac{1}{2} \sum_{a=1}^3 (w_a \gamma_a \gamma_0 - r_a \tilde{\gamma}_a) + b\gamma_5 \quad (4.31)$$

4.2.5 Scalar product of Spinors

We need a scalar product on E , preserved by a gauge transformation, that is by both $Spin(3,1)$ and $Spin(1,3)$.

Theorem 66 *The only scalar products on E , preserved by $\{\gamma C(\sigma), \sigma \in Spin(3,1)\}$*

are $G = \begin{bmatrix} 0 & k\sigma_0 \\ \bar{k}\sigma_0 & 0 \end{bmatrix}$ with $k \in \mathbb{C}$

Proof. It is represented in the basis of E by a 4×4 Hermitian matrix G such that : $G = G^*$

$$\forall s \in Spin(3,1) : [\gamma \circ C(s)]^* G [\gamma \circ C(s)] = G$$

$$\text{or } \forall s \in Spin(1,3) : [\gamma \circ C'(s)]^* G [\gamma \circ C'(s)] = G$$

$$[\gamma \circ C(s)]^* G = G [\gamma \circ C(s)]^{-1} = G [\gamma \circ C(s^{-1})]$$

$$\gamma C(s) = \begin{bmatrix} (a+ib)\sigma_0 + \frac{1}{2} \sum_{a=1}^3 (w_a - ir_a) \sigma_a & 0 \\ 0 & (a-ib)\sigma_0 - \frac{1}{2} \sum_{a=1}^3 (w_a + ir_a) \sigma_a \end{bmatrix}$$

$$\gamma C(s)^* = \begin{bmatrix} (a-ib)\sigma_0 + \frac{1}{2} \sum_{a=1}^3 (w_a + ir_a) \sigma_a & 0 \\ 0 & (a+ib)\sigma_0 - \frac{1}{2} \sum_{a=1}^3 (w_a - ir_a) \sigma_a \end{bmatrix}$$

$$G = \begin{bmatrix} A & B \\ B^* & C \end{bmatrix}, \text{ with } A = A^*, C = C^*$$

$$[\gamma \circ C(s)]^* G$$

$$\begin{aligned}
&= \begin{bmatrix} (a-ib)A + \frac{1}{2} \sum_{g=1}^3 (w_a + ir_a) \sigma_a A & (a-ib)B + \frac{1}{2} \sum_{g=1}^3 (w_a + ir_a) \sigma_a B \\ (a+ib)B^* - \frac{1}{2} \sum_{a=1}^3 (w_a - ir_a) \sigma_a B^* & (a+ib)C - \frac{1}{2} \sum_{a=1}^3 (w_a - ir_a) \sigma_a C \end{bmatrix} \\
G [\gamma \circ C (s^{-1})] \\
&= \begin{bmatrix} (a+ib)A - \frac{1}{2} \sum_{g=1}^3 (w_a - ir_a) A \sigma_a & (a-ib)B + \frac{1}{2} \sum_{g=1}^3 (w_a + ir_a) B \sigma_a \\ (a+ib)B^* - \frac{1}{2} \sum_{a=1}^3 (w_a - ir_a) B^* \sigma_a & (a-ib)C + \frac{1}{2} \sum_{a=1}^3 (w_a + ir_a) C \sigma_a \end{bmatrix} \\
&(a-ib)A + \frac{1}{2} \sum_{a=1}^3 (w_a + ir_a) \sigma_a A = (a+ib)A - \frac{1}{2} \sum_{a=1}^3 (w_a - ir_a) A \sigma_a \\
&(a-ib)B + \frac{1}{2} \sum_{a=1}^3 (w_a + ir_a) \sigma_a B = (a-ib)B + \frac{1}{2} \sum_{a=1}^3 (w_a + ir_a) B \sigma_a \\
&(a+ib)B^* - \frac{1}{2} \sum_{a=1}^3 (w_a - ir_a) \sigma_a B^* = (a+ib)B^* - \frac{1}{2} \sum_{a=1}^3 (w_a - ir_a) B^* \sigma_a \\
&(a+ib)C - \frac{1}{2} \sum_{a=1}^3 (w_a - ir_a) \sigma_a C = (a-ib)C + \frac{1}{2} \sum_{a=1}^3 (w_a + ir_a) C \sigma_a \\
&(w_a + ir_a) \sigma_a B = \sum_{a=1}^3 (w_a + ir_a) B \sigma_a \\
&\sum_{a=1}^3 (w_a - ir_a) \sigma_a B^* = \sum_{a=1}^3 (w_a - ir_a) B^* \sigma_a \\
2ibA \\
&= \frac{1}{2} \sum_{a=1}^3 (w_a + ir_a) \sigma_a A + (w_a - ir_a) A \sigma_a \\
&= \frac{1}{2} \sum_{a=1}^3 w_a (\sigma_a A + A \sigma_a) + ir_a ((\sigma_a A - A \sigma_a)) \\
2ibC \\
&= \frac{1}{2} \sum_{a=1}^3 (w_a + ir_a) C \sigma_a + (w_a - ir_a) \sigma_a C \\
&= \frac{1}{2} \sum_{a=1}^3 w_a (\sigma_a C + C \sigma_a) + ir_a ((C \sigma_a - \sigma_a C)) \\
&\text{By taking the adjoint on the two last equations :} \\
&-2ibA \\
&= \frac{1}{2} \sum_{a=1}^3 w_a (A \sigma_a + \sigma_a A) - ir_a ((A \sigma_a - \sigma_a A)) \\
&= -\frac{1}{2} \sum_{a=1}^3 w_a (\sigma_a A + A \sigma_a) + ir_a ((\sigma_a A - A \sigma_a)) \Rightarrow A = 0 \\
&-2ibC \\
&= \frac{1}{2} \sum_{a=1}^3 w_a (\sigma_a C + C \sigma_a) - ir_a ((C \sigma_a - \sigma_a C)) \\
&= -\frac{1}{2} \sum_{a=1}^3 w_a (\sigma_a C + C \sigma_a) + ir_a ((C \sigma_a - \sigma_a C)) \Rightarrow C = 0 \\
&\text{We are left with :} \\
&\forall w, r : \sum_{a=1}^3 (w_a + ir_a) \sigma_a B = \sum_{a=1}^3 (w_a + ir_a) B \sigma_a \\
&\text{which implies that B commutes with all the Dirac matrices, which happens} \\
&\text{only for the scalar matrices : } B = k \sigma_0. \\
G = \begin{bmatrix} 0 & k \sigma_0 \\ \bar{k} \sigma_0 & 0 \end{bmatrix} \blacksquare
\end{aligned}$$

The scalar product will never be definite positive, so we can take $k = -i$ that is $G = \gamma_0$. And it is easy to check that it works also for the signature (1,3).

Any vector of E reads :

$$u = \sum_{i=1}^4 u^i e_i = u_R + u_L \text{ with } u_R = \sum_{i=1}^2 u^i e_i, u_L = \sum_{i=3}^4 u^i e_i$$

The scalar product of two vectors u, v of E is then:

$$\left\langle \sum_{i=1}^4 u^i e_i, \sum_{i=1}^4 v^i e_i \right\rangle_E = [u]^* [\gamma_0] [v] = i (u_L^* v_R - u_R^* v_L) \quad (4.32)$$

It is not definite positive. It is preserved both by $Spin(3,1)$ and $Spin(1,3)$. It is definite positive on E_R and definite negative on E_L .

The basis $(e_i)_{i=1}^4$ of E is not orthonormal : $\langle e_j, e_k \rangle = i$

$$\begin{bmatrix} 0 & 0 & -1 & 0 \\ 0 & 0 & 0 & -1 \\ 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \end{bmatrix}$$

4.2.6 Norm on the space E of spinors

The scalar product on E is not definite, but we can define a norm, as we have done for the Lie algebra. We have introduced previously the two chiral operators p_R, p_L on $Cl(\mathbb{C}, 4)$. They do not belong to the images $\gamma C(Cl(3, 1)), \gamma C(Cl(1, 3))$ however they define two subspaces $E = E_R \oplus E_L$ and the elements of the images $C(Spin(3, 1))$ and $C'(Spin(1, 3))$ by γ preserve both E^R, E^L . So we have operators $\gamma_R : E \rightarrow E_R, \gamma_L : E \rightarrow E_L$ such that :

$$\gamma_R = \gamma_R \cdot \gamma_R; \gamma_L = \gamma_L \cdot \gamma_L$$

$$\gamma_R + \gamma_L = Id$$

$$\forall \sigma \in Spin(3, 1), Spin(1, 3) : \gamma C(\sigma) \circ \gamma_R = \gamma_R \circ \gamma C(\sigma); \gamma C(\sigma) \circ \gamma_L = \gamma_L \circ \gamma C(\sigma)$$

$\gamma_R = \gamma(p_R), \gamma_L = \gamma(p_L)$ are complex linear maps, as images of the complex linear maps p_R, p_L by the complex linear map γ . So they preserve any real structure on E : $\gamma_R(\text{Re } u + i \text{Im } u) = \gamma(p_R)(\text{Re } u + i \text{Im } u) = \gamma(p_R) \text{Re } u + i \gamma(p_R)(\text{Im } u)$ and $\gamma_R = \overline{\gamma}_R, \gamma_L = \overline{\gamma}_L$ (Maths.357).

In the basis $(e_i)_{i=1}^4$:

$$\gamma_R = \begin{bmatrix} \sigma_0 & 0 \\ 0 & 0 \end{bmatrix} = \gamma_R^*$$

$$\gamma_L = \begin{bmatrix} 0 & 0 \\ 0 & \sigma_0 \end{bmatrix} = \gamma_L^*$$

There is no definite scalar product on E , but there is a norm, that is a map :

$$\| \cdot \| : E \times E \rightarrow \mathbb{R}$$

such that :

$$\|S\| \geq 0$$

$$\|S\| = 0 \Rightarrow S = 0$$

$$\|kS\| = |k| \|S\|$$

$$\|S + S'\| \leq \|S\| + \|S'\|$$

Theorem 67 *The vector space E is a normed vector space with the norm, invariant by $Spin(3, 1), Spin(1, 3)$:*

$$\|S\|_E = \sqrt{\langle \gamma_R S, \gamma_R S \rangle_E - \langle \gamma_L S, \gamma_L S \rangle_E}$$

Proof. E_R, E_L are two 2 dimensional complex vector spaces, they can be endowed with a norm which is invariant by $Spin(3, 1), Spin(1, 3)$:

$$S \in E_R : \|S\|_{E_R}^2 = \langle S, S \rangle_E$$

$$S \in E_L : \|S\|_{E_L}^2 = -\langle S, S \rangle_E$$

The norm are invariant by γ_R, γ_L :

$$S \in E_R \Leftrightarrow \exists S' \in E : S = \gamma_R(S')$$

$$\begin{aligned} \|\gamma_R(S)\|_{E_R} &= \|\gamma_R^2(S')\|_{E_R} = \|\gamma_R(S')\|_{E_R} = \|S\|_{E_R} \\ \text{Define : } \|\cdot\|_E : E \times E &\rightarrow \mathbb{R} :: \|S\|_E = \sqrt{\|\gamma_R S\|_{E_R}^2 + \|\gamma_L S\|_{E_L}^2} \\ \|S\|_E = 0 &\Rightarrow \|\gamma_R S\|_{E_R} = 0; \|\gamma_L S\|_{E_L} = 0 \Rightarrow \gamma_R S = 0; \gamma_L S = 0 \Rightarrow \\ (\gamma_R + \gamma_L)[S] &= 0 = S \\ \|S + S'\|_E^2 &= \|\gamma_R(S + S')\|_{E_R}^2 + \|\gamma_L(S + S')\|_{E_L}^2 = \|S + S'\|_{E_R}^2 + \|S + S'\|_{E_L}^2 \\ &\leq \|S\|_{E_R}^2 + \|S'\|_{E_R}^2 + \|S\|_{E_L}^2 + \|S'\|_{E_L}^2 = \|S\|_E^2 + \|S'\|_E^2 \\ \text{This norm is invariant by } &Spin(3, 1), Spin(1, 3): \\ \|\gamma C(\sigma) S\|_E &= \sqrt{\|\gamma C(\sigma) \gamma_R S\|_{E_R}^2 + \|\gamma C(\sigma) \gamma_L S\|_{E_L}^2} \\ &= \sqrt{\|\gamma_R \gamma C(\sigma) S\|_{E_R}^2 + \|\gamma_L \gamma C(\sigma) S\|_{E_L}^2} \\ &= \sqrt{\|\gamma C(\sigma) S\|_{E_R}^2 + \|\gamma C(\sigma) S\|_{E_L}^2} = \sqrt{\|S\|_{E_R}^2 + \|S\|_{E_L}^2} = \|S\|_E \quad \blacksquare \end{aligned}$$

4.3 THE SPINOR MODEL OF KINEMATICS

We have now the mathematical tools to enter the representation of kinematics of material bodies in General Relativity. First we will make some adjustment to the fiber bundles used so far to represent the geometry, to account for the introduction of the Spin group.

4.3.1 Description of the fiber bundles

The geometric fiber bundles

The geometric model is similar to the previous one, with the replacement of $SO(3, 1)$ by the Spin group.

Definition 68 *The **principal bundle** $P_G(M, Spin_0(3, 1), \pi_G)$ has for fiber the connected component of the unity of the Spin group, for trivialization the map :*

$$\begin{aligned} \varphi_G : M \times Spin_0(3, 1) &\rightarrow P_G :: p = \varphi_G(m, s). \\ \text{The **standard gauge** used by observers is } \mathbf{p}(m) &= \varphi_G(m, \mathbf{1}) \end{aligned}$$

A section $\sigma \in \mathfrak{X}(P_G)$ is defined by a map: $\sigma : M \rightarrow Spin(3, 1)$ such that : $\sigma(m) = \varphi_G(m, \sigma(m))$ and in a change of gauge :

$$\mathbf{p}(m) = \varphi_G(m, \mathbf{1}) \rightarrow \tilde{\mathbf{p}}(m) = \mathbf{p}(m) \cdot \chi(m)^{-1} : \quad (4.33)$$

$$\sigma(m) = \varphi_G(m, \sigma(m)) = \tilde{\varphi}_G(m, \chi(m) \cdot \sigma(m)) \quad (4.34)$$

Definition 69 *The **vectors on the fiber bundle** TM are represented in the associated vector bundle $P_G[\mathbb{R}^4, \mathbf{Ad}]$ defined through the holonomic orthonormal basis :*

$$\varepsilon_i(m) = (\mathbf{p}(m), \varepsilon_i)$$

So $\varepsilon_0(m) = (\mathbf{p}(m), \varepsilon_0)$ is the 4th vector both in the Clifford algebra and in the tangent space $T_m M$. It corresponds to the velocity of the observer.

With the equivalence relation : $(\mathbf{p}(m), v) \sim (\varphi_G(m, g), \mathbf{Ad}_{g^{-1}} v)$

In a change of gauge on P_G the holonomic basis becomes :

$$\mathbf{p}(m) = \varphi_G(m, \mathbf{1}) \rightarrow \tilde{\mathbf{p}}(m) = \mathbf{p}(m) \cdot \chi(m)^{-1} : \quad (4.35)$$

$$\varepsilon_i(m) = (\mathbf{p}(m), \varepsilon_i) \rightarrow \tilde{\varepsilon}_i(m) = \mathbf{Ad}_{\chi(m)^{-1}} \varepsilon_i(m) \quad (4.36)$$

For a given observer $\varepsilon_0(p_o(t)) = \frac{1}{c} \frac{dp_o}{dt}$ is fixed along his world line.

The Lorentz scalar product on \mathbb{R}^4 is preserved by \mathbf{Ad} thus it can be extended to $P_G[\mathbb{R}^4, \mathbf{Ad}]$.

Definition 70 *The **Clifford bundle** $Cl(TM)$ is the associated vector bundle $P_G[Cl(3, 1), \mathbf{Ad}]$ defined through the basis $(\varepsilon_i(m))_{i=0}^3$.*

In a change of gauge on P_G the elements of $\text{Cl}(m)$ transforms as :

$$\mathbf{p}(m) = \varphi_G(m, 1) \rightarrow \tilde{\mathbf{p}}(m) = \mathbf{p}(m) \cdot \chi(m)^{-1} : \quad (4.37)$$

$$\sigma(m) = \varphi_P(m, \sigma(m)) = \tilde{\varphi}_P(m, \chi(m) \cdot \sigma(m)) \quad (4.38)$$

$$\mathbf{w}(m) = (\mathbf{p}(m), w) \rightarrow \mathbf{Ad}_{\chi(m)^{-1}} \mathbf{w}(m) \quad (4.39)$$

The kinematic bundle

In addition to the previous bundles we define the associated bundle in which the spinors live :

Proposition 71 *The relativist momentum of particles are represented by Spinors, which are, at each point of the world line of the particle, vectors of the associated vector bundle $P_G[E, \gamma C]$. They are measured by observers in the standard gauge defined through the holonomic basis : $\mathbf{e}_i(m) = (\mathbf{p}(m), e_i)$*

With the equivalence relation : $(\mathbf{p}(m), S) \sim (\varphi_G(m, g), \gamma C(g^{-1})S)$ so that in a change of gauge the holonomic basis becomes :

$$\mathbf{p}(m) = \varphi_G(m, 1) \rightarrow \tilde{\mathbf{p}}(m) = \mathbf{p}(m) \cdot \chi(m)^{-1} : \quad (4.40)$$

$$\mathbf{e}_i(m) = (\mathbf{p}(m), e_i) \rightarrow \tilde{\mathbf{e}}_i(m) = \gamma C(\chi(m)^{-1}) \mathbf{e}_i(m) \quad (4.41)$$

and the components of a section $S \in \mathfrak{X}(P_G[E, \gamma C])$ change as :

$$\mathbf{p}(m) = \varphi_G(m, 1) \rightarrow \tilde{\mathbf{p}}(m) = \mathbf{p}(m) \cdot \chi(m)^{-1} : \quad (4.42)$$

$$\mathbf{S}(m) = (\mathbf{p}(m), S_m) = \left(\mathbf{p}(m) \cdot \chi(m)^{-1}, \gamma C(\chi(m)) S_m \right) \quad (4.43)$$

From a mathematical point of view the holonomic basis $(\varepsilon_i(m))_{i=0}^3$, $(\mathbf{e}_i(m))_{i=1}^4$, are defined through the same standard gauge $\mathbf{p}(m)$ chosen by the observer. This gauge is arbitrary. For the tetrad the vectors $\varepsilon_i(m)$ can be measured in the holonomic basis of any chart through P . We have nothing similar for $\mathbf{e}_i(m)$, and actually the vectors e_i of E themselves are abstract. However we will see in the following how the basis $\mathbf{e}_i(m)$ used by an observer can be related to physical phenomena (inertial observers).

The scalar product on E is preserved by γC thus it can be extended to $P_G[E, \gamma C]$ and to the space of sections $\mathfrak{X}(P_G[E, \gamma C])$ by :

$$\langle \mathbf{S}, \mathbf{S}' \rangle = \int_{\Omega} \langle \mathbf{S}(m), \mathbf{S}'(m) \rangle_E \varpi_4(m)$$

Moreover we have the following :

Theorem 72 *The set of integrable sections :*

$$L^1(\mathfrak{X}(P_G[E, \gamma C]), \varpi_4) = \left\{ \int_{\Omega} \|S\| \varpi_4 < \infty \right\}$$

with the norm on E is a separable, infinite dimensional Fréchet space.

Proof. Consider the vector space : $\mathfrak{X}(P_G[E, \gamma C])$ endowed with the norm :

$$\|\mathbf{S}\| = \int_{\Omega} \|S(m)\|_E \varpi_4(m) \text{ and the norm}$$

$$\|S(m)\|_E = \sqrt{\langle \gamma_R S, \gamma_R S \rangle_E - \langle \gamma_L S, \gamma_L S \rangle_E}$$

Restrict this space to $L^1(M, P_G[E, \gamma C], \varpi_4)$

$$= \{ \mathbf{S} \in \mathfrak{X}(P_G[E, \gamma C]) \mid \int_{\Omega} \|S(m)\|_E \varpi_4(m) < \infty \}$$

This is a Fréchet space (Maths.2276). Moreover it is separable, because Ω is relatively compact and the smooth compactly supported maps are a countable basis in L^1 (see Lieb).

Because the norm is invariant by the Spin group this space does not depend on the choice of trivialization. ■

The result still holds if we impose that the sections are differentiable.

Fundamental symmetry breakdown

The observer uses the frame $(O, (\varepsilon_i)_{i=0}^3)$ to measure the components of vectors of TM, and the holonomic maps $(\mathbf{e}_i(m))_{i=0}^3$ to measure the spinors. The breakdown, specific to each observer, comes from the distinction of his present, and is materialized in his standard basis by the vector $\varepsilon_0(m)$. This choice leads to a split of the Spin group between the spatial rotations, represented by $Spin(3)$, and the homogeneous space $Spin(3,1)/Spin(3)$.

We have an associated fiber bundle :

$$P_W = P_G[Spin(3,1)/Spin(3), \lambda] :$$

$$(\mathbf{p}(m), s_w) = (\varphi_G(m, 1), s_w) \sim (\varphi_G(m, s), \lambda(s^{-1}, s_w))$$

with the left action :

$$\lambda : Spin(3,1) \times Spin(3,1)/Spin(3) \rightarrow Spin(3,1)/Spin(3) : \lambda(s, s_w) = \pi_w(s \cdot s_w)$$

On the manifold P_G there is a structure of principal fiber bundle

$$P_R(P_W, Spin(3), \pi_R) \text{ with trivialization :}$$

$$\varphi_R : P_W \times Spin(3) \rightarrow P_G ::$$

$$\varphi_R((\mathbf{p}(m), s_w), s_r) = \varphi_G(m, s_w \cdot s_r) = \varphi_R((\varphi_G(m, s), \lambda(s^{-1}, s_w)), s_r)$$

As the latest trivialization shows, for a given s , s_r depends on s_w in that it is a part of $s \in Spin(3,1)$.

It sums up to define the local basis in two steps : first by choosing s_w second by choosing s_r .

Any section $\sigma \in \mathfrak{X}(P_G)$ can be decomposed, for a given vector field ε_0 and a fixed $\epsilon = \pm 1$, in two sections :

$$\epsilon \sigma_w \in \mathfrak{X}(P_W), \epsilon \sigma_r \in \mathfrak{X}(P_R) \text{ with } \sigma(m) = \epsilon \sigma_w(m) \cdot \epsilon \sigma_r(m)$$

The set of vectors of $T_m M$ used to build $Spin(3)$ is defined by $\varepsilon_0(m)$.

Our objective is now to understand the relation between sections

$\mathbf{S} \in \mathbf{P}_G[E, \gamma C]$ of the vector bundle and the motion of a particle, and to precise the possible values of \mathbf{S} . We will proceed in several steps. First we will focus on the trajectories and we will show that they can be matched with sections of P_W .

4.3.2 Trajectories and the Spin Group

A change of observer is a change of gauge $\mathbf{p}(m) = \varphi_G(m, \mathbf{1})$ (or trivialization) on the principal bundle $P_G(M, Spin_0(3, 1), \pi_G)$. We see now how the formulas given previously can be expressed in the formalism of Clifford algebras.

Theorem 73 Any section $\sigma \in \mathfrak{X}(P_G)$ defines, for any positive function $f \in C_\infty(\Omega; \mathbb{R}_+)$ and observer, two vector fields $V \in \mathfrak{X}(P_G[\mathbb{R}^4, \mathbf{Ad}])$ by :

$$V(m) = f(m) \mathbf{Ad}_{\sigma(m)} \varepsilon_0(m) = f(m) ((2a_w^2 - 1) \varepsilon_0 + \epsilon a_w (w^1 \varepsilon_1 + w^2 \varepsilon_2 + w^3 \varepsilon_3)) \quad (4.44)$$

where $\sigma_w(m) = \epsilon(a_w + v(0, w))$ is the projection of σ on P_W along ε_0

Then V is time like, future oriented and $\langle V(m), V(m) \rangle = -f^2(m)$ and is invariant in a change of gauge on P_G

Conversely, for any time like, future oriented vector field $V \in \mathfrak{X}(P_G[\mathbb{R}^4, \mathbf{Ad}])$ there are two sections $\sigma_w \in \mathfrak{X}(P_G)$ such that :

$$\frac{V}{\sqrt{-\langle V, V \rangle}} = u = \mathbf{Ad}_{\sigma_w(m)} \varepsilon_0(m) : \sigma_w = \epsilon \left(\sqrt{\frac{1}{2}(u_0 + 1)} + \frac{1}{\sqrt{\frac{1}{2}(u_0 + 1)}} v(0, u_i) \right)$$

Proof. i) $\sigma(m) = \epsilon \sigma_w(m) \cdot \epsilon \sigma_r(m)$

$\sigma_w(m) = \epsilon(a_w + v(0, w))$ so let be $a_w > 0$ which defines ϵ

ii) Define

$u \in \mathfrak{X}(P_G[\mathbb{R}^4, \mathbf{Ad}]) : u(m) = \mathbf{Ad}_{\sigma_w(m)} \varepsilon_0(m) = (\mathbf{p}(m), \mathbf{Ad}_{\sigma_w} \varepsilon_0) = (\mathbf{p}(m), u)$

$$\begin{aligned} u &= \sigma_w \cdot \varepsilon_0 \cdot \sigma_w^{-1} = (\epsilon a_w + v(0, w)) \cdot \varepsilon_0 \cdot (\epsilon a_w - v(0, w)) \\ &= (\epsilon a_w + v(0, w)) \cdot (\epsilon a_w \varepsilon_0 - \varepsilon_0 \cdot v(0, w)) \\ &= a_w^2 \varepsilon_0 + \epsilon a_w (-\varepsilon_0 \cdot v(0, w) + v(0, w) \cdot \varepsilon_0) - v(0, w) \cdot \varepsilon_0 \cdot v(0, w) \\ &= a_w^2 \varepsilon_0 + \frac{1}{2} \epsilon a_w (-\varepsilon_0 \cdot \varepsilon_0 \cdot (w^1 \varepsilon_1 + w^2 \varepsilon_2 + w^3 \varepsilon_3) - (w^1 \varepsilon_1 + w^2 \varepsilon_2 + w^3 \varepsilon_3) \cdot \varepsilon_0 \cdot \varepsilon_0) \\ &\quad - \frac{1}{4} \varepsilon_0 \cdot (w^1 \varepsilon_1 + w^2 \varepsilon_2 + w^3 \varepsilon_3) \cdot \varepsilon_0 \cdot \varepsilon_0 \cdot (w^1 \varepsilon_1 + w^2 \varepsilon_2 + w^3 \varepsilon_3) \\ &= a_w^2 \varepsilon_0 + \frac{1}{2} \epsilon a_w (w + w) + \frac{1}{4} \varepsilon_0 \cdot w \cdot w \\ &= a_w^2 \varepsilon_0 + a_w \epsilon w + \frac{1}{4} \varepsilon_0 \cdot \langle w, w \rangle \\ &= (a_w^2 + \frac{1}{4} w^t w) \varepsilon_0 + \epsilon a_w w \end{aligned}$$

$$u = (2a_w^2 - 1) \varepsilon_0 + \epsilon a_w (w^1 \varepsilon_1 + w^2 \varepsilon_2 + w^3 \varepsilon_3)$$

$$\text{iii) } \langle u, \varepsilon_0 \rangle = -(2a_w^2 - 1) = 1 - 2(1 + \frac{1}{4} w^t w) = -1 - \frac{1}{2} w^t w < 0$$

$$\langle u, u \rangle = a_w^2 w^t w - (2a_w^2 - 1)^2 = a_w^2 (4(a_w^2 - 1)) - (2a_w^2 - 1)^2 = -1$$

iv) $V = f(m) u \Rightarrow$

$$\langle V, \varepsilon_0 \rangle = f(m) \langle u, \varepsilon_0 \rangle < 0$$

$$\langle V, V \rangle = -f^2 < 0$$

v) In a change of gauge on P_G :

$$\mathbf{p}(m) = \varphi_G(m, \mathbf{1}) \rightarrow \tilde{\mathbf{p}}(m) = \mathbf{p}(m) \cdot \chi(m)^{-1} :$$

$$\sigma(m) \rightarrow \tilde{\sigma}(m) = \chi(m)^{-1} \cdot \sigma(m)$$

$$u(m) = (\mathbf{p}(m), \mathbf{Ad}_{\sigma_w} \varepsilon_0) \sim (\mathbf{p}(m) \cdot \chi(m)^{-1}, \mathbf{Ad}_\chi \mathbf{Ad}_{\sigma_w} \varepsilon_0) = (\tilde{\mathbf{p}}(m), \mathbf{Ad}_{\tilde{\sigma}_w} \varepsilon_0) =$$

$\tilde{u}(m)$

vi) Let $u = \frac{V}{\sqrt{-\langle V, V \rangle}}$ then u is time like, $\langle u, u \rangle = -1, \langle u, \varepsilon_0 \rangle < 0$

$$\begin{aligned}
u &= ((2a_w^2 - 1) \varepsilon_0 + \epsilon a_w (w^1 \varepsilon_1 + w^2 \varepsilon_2 + w^3 \varepsilon_3)) = \sum_{i=0}^3 u_i \varepsilon_i \\
i=1,2,3 : w_i &= \epsilon u_i / a_w \\
w^t w &= \frac{\sum_{i=1}^3 u_i^2}{a_w^2} \\
\langle u, u \rangle &= -1 = \sum_{i=1}^3 u_i^2 - u_0^2 \\
u_0^2 &= 1 + \sum_{i=1}^3 u_i^2 \\
\langle u, \varepsilon_0 \rangle &= u_0 < 0 \\
u_0 &= -\sqrt{1 + \sum_{i=1}^3 u_i^2} < -1 \\
1 + \frac{1}{4} w^t w &= 1 + \frac{1}{4} \frac{\sum_{i=1}^3 u_i^2}{a_w^2} = \frac{1+4 \sum_{i=1}^3 u_i^2}{4a_w^2} = \frac{1+4(u_0^2-1)}{4a_w^2} = \frac{-3+4u_0^2}{4a_w^2} > \frac{-3+4}{4a_w^2} = \frac{1}{a_w^2} \\
\text{So we can define :} & \\
u_0 &= 2a_w^2 - 1 \\
a_w &= \epsilon \sqrt{\frac{1}{2}(u_0 + 1)} \\
w_i &= u_i / a_w = \epsilon u_i / \sqrt{\frac{1}{2}(u_0 + 1)} \\
\sigma_w &= \epsilon \left(\sqrt{\frac{1}{2}(u_0 + 1)} + \frac{1}{\sqrt{\frac{1}{2}(u_0 + 1)}} v(0, u_i) \right) = a_w + v(0, w) \\
\text{with } w_i &= \frac{u_i}{\sqrt{\frac{1}{2}(u_0 + 1)}} = \frac{V_i}{\sqrt{\frac{1}{2} \left(\frac{V_0}{\sqrt{-\langle V, V \rangle}} + 1 \right)}} \frac{1}{\sqrt{-\langle V, V \rangle}} = \frac{V_i}{\sqrt{-\frac{1}{2} (\langle V, V \rangle - V_0 \sqrt{-\langle V, V \rangle})}}
\end{aligned}$$

■

If we take $f(m) = c$ any section $\sigma \in \mathfrak{X}(P_G)$ defines two fields of world lines, with opposite spatial speed :

$$u = \frac{dp}{d\tau} = \frac{1}{\sqrt{1 - \frac{\|\vec{v}\|^2}{c^2}}} (\vec{v} + c\varepsilon_0(m)) = c \left((2a_w^2 - 1) \varepsilon_0 + \epsilon a_w \sum_{i=1}^3 w_i \varepsilon_i \right) \quad (4.45)$$

If we take $f(m) = \frac{c}{2a_w^2 - 1}$ any section $\sigma \in \mathfrak{X}(P_G)$ defines two fields of trajectories, with opposite spatial speed :

$$V = \frac{dp}{dt} = \vec{v} + c\varepsilon_0(m) = c \left(\varepsilon_0 + \epsilon \frac{a_w}{2a_w^2 - 1} \sum_{i=1}^3 w_i \varepsilon_i \right) \quad (4.46)$$

$$\begin{aligned}
\vec{v} &= \epsilon c \frac{a_w}{2a_w^2 - 1} \sum_{i=1}^3 w_i \varepsilon_i \\
a_w &= \sqrt{\frac{1}{2} \left(1 + \sqrt{1 + \frac{1}{4(1 - \frac{v^2}{c^2})}} \right)} \simeq 0.7 + \frac{1}{16} \frac{v^2}{c^2} \\
w &= \frac{\sqrt{1 + \frac{1}{4(1 - \frac{v^2}{c^2})}}}{\sqrt{1 + \sqrt{1 + \frac{1}{4(1 - \frac{v^2}{c^2})}}}} \frac{\vec{v}}{c} \simeq 0.8 \frac{\vec{v}}{c}
\end{aligned}$$

Remarks :

- i) All this is defined with respect to an observer, who fixes $\varepsilon_0(m)$
- ii) $V \in \mathfrak{X}(P_G[\mathbb{R}^4, \mathbf{Ad}])$ so can be equivalently defined as a section of TM :

$$V = \sum_{\alpha=0}^3 V^\alpha \partial \xi_\alpha = \sum_{\alpha=0}^3 \sum_{i=0}^3 V^i P_i^\alpha \partial \xi_\alpha$$

iii) If V is past oriented ($u_0 < 0$) or null ($\langle V, V \rangle = 0$) there is no solution :
 $2a_w^2 - 1 = \frac{1}{2}(u_0 - 1) < -\frac{1}{2} \Rightarrow a_w^2 < 1$ and $a_w^2 \neq 1 + \frac{1}{4}w^t w$

This gives a strong physical meaning to the representation of world lines by section of P_W .

iv) Any map $\sigma : \mathbb{R} \rightarrow P_G$ is projected on M as a curve, which is not necessarily time like or defines a world line.

4.3.3 Spatial spinor

A section $\sigma \in \mathfrak{X}(P_G)$ defines at each point an element of $Spin(3, 1)$, which can uniquely (up to sign) be decomposed in $\sigma = \epsilon\sigma_w \cdot \epsilon\sigma_r$ with respect to a given observer.

The first component $\sigma_w = \pm(a_w + v(0, w)) \in Spin(3, 1)$ defines, in the standard basis of the observer, a vector field of world lines, and a trajectory. w aligned in the direction of the spatial speed or the opposite.

The second part $\sigma_r = a_r + v(0, r)$ belonging to one of the two connected components of $Spin(3)$ (according to the sign of a_r , it is + for the component of the identity) leaves invariant $\varepsilon_0(m)$ and defines a spatial rotation, in the hyperplane orthogonal to $\varepsilon_0(m)$.

So with a single σ we have both a translational motion (along a worldline) and a spatial rotation, at the same point.

Notice that this is not a rotation around a point (like an orbit), but a *rotation at a point*. The action \mathbf{Ad}_{σ_r} on any vector of the tangent space $T_m M$ rotates the vector, but leaves invariant $\varepsilon_0(m)$, so this is an action on the physical space $\Omega_3(t)$. In the Spinor $S = \gamma C(\sigma(m)) S_0$ this action is done on vectors of E , and not on vector of $T_m M$.

The decomposition $\sigma = \epsilon\sigma_w \cdot \epsilon\sigma_r$ and the identification of the vectors of $Spin(3)$ request a choice of $\varepsilon_0(m)$: it is observer-dependant.

Our basic hypothesis is that the spinor S is a kinematic characteristic of the state of the particle : it represents the relativist momentum. With a section $\sigma \in \mathfrak{X}(P_G)$ and fixed vector $S_0 \in E$ we have a representation of the momenta, both translational and rotational.

Definition 74 We can then define the *spatial spinor* as :

$$\mathbf{S}_r(t) = \gamma C(\sigma_w^{-1}) \mathbf{S}(t) = \gamma C(\sigma_r(t)) S_0 \quad (4.47)$$

The spatial spinor is the representation of a rotational momentum. For a given trajectory and observer there are two possible, opposite, values of the spatial spinor : $S_r(t) = \pm \gamma C(\sigma_r(t)) S_0$. In all cases $S(t) = \gamma C(\sigma(t)) S_0$: the total spinor stays the same, the distinction between σ_w, σ_r and the opposite values is the consequence of the breakdown of symmetry induced by the observer. The sign \pm is related to a trajectory (the orientation of w with respect to the spatial speed) so one can speak of **spin up or down** with respect to the trajectory. This feature is entirely linked to the Relativist picture, and has

nothing to do with QM. The name spin is used freely in Quantum Physics, and this is sometimes confusing. The usual spin, as rotational momentum, is the Spatial Spinor and we will give a precise definition later. And to be clear I will call the present feature (spin up or down) **Relativist Spin** which takes the values of $\epsilon = \pm 1$.

If we assume that the spatial spinor is, by itself, an intrinsic feature of the particle, then one must assume that the map : $\sigma_r : \mathbb{R} \rightarrow P_R$ is continuous, thus σ_r must belong and stay in one of the two connected components of P_R . Normally the decomposition $\sigma = \epsilon\sigma_w \cdot \epsilon\sigma_r$ is continuous, and the passage to the opposite sign is, for the spatial spinor, a discontinuity, and also for the relativist spin.

The issue now is to precise what can be S_0 , that we will call **inertial spinor**. The 4 dimensional relativist momentum $P = \mu u$, which is a geometric quantity, has a constant scalar product : $\langle P, P \rangle = -\mu^2 c^2$ where μ is, by definition, the mass at rest. The scalar product $\langle S, S \rangle = \langle S_0, S_0 \rangle$ is preserved on the world line, so we will look at vectors S such that : $S \neq 0 \Rightarrow \langle S_0, S_0 \rangle \neq 0$.

4.3.4 Inertial spinor

We have seen that the unique scalar product on E is non degenerate, but not definite positive. So it is logical to require that S_0 belongs to some vector subspace E_0 of E , over which the scalar product is definite, either positive or negative. Moreover a change of spatial frame should change only S_r , thus E_0 should be invariant under the action of $Spin(3)$.

So there should be some vector subspace E_0 of E such that :

- it is invariant by $\gamma C(\sigma_r)$ for $\sigma_r \in Spin(3) : \forall S_0 \in E_0, s_r \in Spin(3) : \gamma C(s_r) S_0 \in E_0$

- on which the scalar product is either definite positive or definite negative : $\forall S_0 \in E_0 : \langle S_0, S_0 \rangle_E = 0 \Rightarrow S_0 = 0$

Theorem 75 *The only vector subspace of E invariant by γC on $Spin(3)$ and over which the scalar product is definite*

- positive is $E_0 = \left\{ S = \begin{bmatrix} S_R \\ S_L \end{bmatrix} = \begin{bmatrix} v \\ iv \end{bmatrix}, v \in \mathbb{C}^2 \right\}$

- negative is $E'_0 = \left\{ S = \begin{bmatrix} S_R \\ S_L \end{bmatrix} = \begin{bmatrix} v \\ -iv \end{bmatrix}, v \in \mathbb{C}^2 \right\}$

Proof. i) The scalar product on E (which does not depend on the signature) reads :

$$u = \begin{bmatrix} u_R \\ u_L \end{bmatrix} \in E : \begin{bmatrix} u_R^* & u_L^* \end{bmatrix} \begin{bmatrix} 0 & -i\sigma_0 \\ i\sigma_0 & 0 \end{bmatrix} \begin{bmatrix} u_R \\ u_L \end{bmatrix} = i(u_L^* u_R - u_R^* u_L) = i(\bar{u}_L^t u_R - (\bar{u}_R^t u_L))$$

$$u_L = v_L + iw_L \text{ with } v_L, w_L \in \mathbb{R}^2$$

$$u_R = v_R + iw_R \text{ with } v_R, w_R \in \mathbb{R}^2$$

$$\langle S, S \rangle = i((v_L^t - iw_L^t)(v_R + iw_R) - (v_R^t - iw_R^t)(v_L + iw_L))$$

$$= i(v_L^t v_R + i v_L^t w_R - i w_L^t v_R + w_L^t w_R - v_R^t v_L - i v_R^t w_L + i w_R^t v_L - w_R^t w_L)$$

$$= -2(v_L^t w_R - w_L^t v_R)$$

$$\langle S, S \rangle = 0 \Leftrightarrow v_L^t w_R = w_L^t v_R$$

So it is definite for any u such that :

$$v_L = -\epsilon w_R, w_L = \epsilon v_R \Rightarrow$$

$$u_L = -\epsilon w_R + \epsilon i v_R = \epsilon i (v_R + i w_R) = \epsilon i u_R$$

$$u = \begin{bmatrix} u_R \\ \epsilon i u_R \end{bmatrix}$$

$$\langle S, S \rangle = 2\epsilon (w_R^t w_R + v_R^t v_R) = 2\epsilon u_R^* u_R$$

It is definite positive for $\epsilon = +1$ and definite negative for $\epsilon = -1$

ii) The vector subspace must be invariant by $\gamma C(s_r)$. Which is equivalent to $S_L = \epsilon i S_R$

For any $S_0 \in E_0, E'_0, s \in Spin(3, 1)$

$$\gamma C(a + v(r, w) + b\epsilon_5) S_0 =$$

$$\begin{bmatrix} (a + ib)\sigma_0 + \frac{1}{2} \sum_a (w_a - ir_a)\sigma_a & 0 \\ 0 & (a - ib)\sigma_0 - \frac{1}{2} \sum_a (w_a + ir_a)\sigma_a \end{bmatrix} \begin{bmatrix} v \\ \epsilon i v \end{bmatrix}$$

$$= \begin{bmatrix} S_R \\ S_L \end{bmatrix}$$

$$S_R = \left((a + ib)\sigma_0 + \frac{1}{2} \sum_{a=1}^3 (w_a - ir_a)\sigma_a \right) v$$

$$S_L = \epsilon \left((a - ib)\sigma_0 - \frac{1}{2} \sum_{a=1}^3 (w_a + ir_a)\sigma_a \right) i v$$

and $S \in E_0 \Leftrightarrow S_L = \epsilon i S_R$

$$\Leftrightarrow \epsilon i \left((a - ib)\sigma_0 - \frac{1}{2} \sum_{a=1}^3 (w_a + ir_a)\sigma_a \right) v$$

$$= \epsilon i \left((a + ib)\sigma_0 + \frac{1}{2} \sum_{a=1}^3 (w_a - ir_a)\sigma_a \right) v$$

$$\Leftrightarrow \left(-ib\sigma_0 - \frac{1}{2} \sum_{a=1}^3 w_a \sigma_a \right) v = \left(ib\sigma_0 + \frac{1}{2} \sum_{a=1}^3 w_a \sigma_a \right) v$$

$$\Leftrightarrow \left(ib\sigma_0 + \frac{1}{2} \sum_{a=1}^3 w_a \sigma_a \right) v = 0$$

This condition is met for $w = 0$ that is $s \in Spin(3)$.

iii) It is easy to see that the result does not depend on the signature:

$$\gamma C'(s) \begin{bmatrix} v \\ \epsilon i v \end{bmatrix} = \begin{bmatrix} S_R \\ S_L \end{bmatrix} \Rightarrow$$

$$S_R = \left((a + b)\sigma_0 + \frac{1}{2} \sum_{a=1}^3 (w_a + ir_a)\sigma_a \right) v$$

$$S_L = \epsilon i \left((a - b)\sigma_0 + \frac{1}{2} \sum_{a=1}^3 (-w_a + ir_a)\sigma_a \right) v$$

$$\left(2b\sigma_0 + \sum_{a=1}^3 w_a \sigma_a \right) v = 0 \quad \blacksquare$$

So particles have both a left S_L and a right S_R part, which are linked but not equal. We have one of the known features of particles : chirality.

Because $E_0 \cap E'_0 = \{0\}$, $(E_0, \gamma C), (E'_0, \gamma C)$ are two, non equivalent, irreducible representations of $Spin(3)$. So they can be seen as corresponding to two kinds of particles according to ϵ .

The inertial spinor is defined, from the components of the two complex vectors of S_R , by 4 real scalars.

The quantity :

$$\langle S_0, S_0 \rangle_E = \epsilon 2S_R^* S_R \quad (4.48)$$

(with the same meaning of ϵ as above) is a scalar, which is conserved along the trajectory. Similarly to the 4 vector $P = \mu u$, we can define a mass at rest M_p by : $\langle S_0, S_0 \rangle = M_p^2 c^4$, but because $\langle S_0, S_0 \rangle$ can be negative we have to consider :

$$\langle S_0, S_0 \rangle = \epsilon M_p^2 c^4 \quad (4.49)$$

where ϵ is a characteristic of the particle. We retrieve a celebrated Dirac's result from his equation. So we define :

Definition 76 particles are such that $S_L = iS_R$.

Their mass is $M_p = \frac{1}{c^2} \sqrt{\langle S_0, S_0 \rangle_E} = \frac{1}{c^2} \sqrt{2S_R^* S_R}$

antiparticles are such that $S_L = -iS_R$

Their mass is $M_p = \frac{1}{c^2} \sqrt{-\langle S_0, S_0 \rangle_E} = \frac{1}{c^2} \sqrt{-2S_R^* S_R}$

Do antiparticles have negative mass ? The idea of a negative mass is still controversial. Dirac considered that antiparticles move backwards in time and indeed a negative mass combined with the first Newton's law seems to have this effect. But here the world line of the particle is defined by σ_w , and there is no doubt about the behavior of an antiparticle : it moves towards the future. The mass at rest M_p is somewhat conventional, the defining relation is $\langle S_0, S_0 \rangle = \epsilon M_p^2 c^2$ so we can choose any sign for M_p , and it seems more appropriate to take $M_p > 0$ both for particles and antiparticles.

Then S_R reads :

$$S_R = \frac{M_p}{c^2 \sqrt{2}} \begin{bmatrix} a \\ b \end{bmatrix} \text{ and } 1 = (|a|^2 + |b|^2)$$

It is customary to represent the polarization of the plane wave of an electric field by two complex quantities (the Jones vector) :

$$E_x = E_{0x} e^{i\alpha_x}$$

$$E_y = E_{0y} e^{i\alpha_y}$$

where (E_{0x}, E_{0y}) are the components of a vector E_0 along the axes x,y.

So we can write similarly :

$$S_R = \frac{M_p c^2}{\sqrt{2}} \begin{bmatrix} e^{i\alpha_1} \cos \alpha_0 \\ e^{i\alpha_2} \sin \alpha_0 \end{bmatrix} \quad (4.50)$$

Particles :

$$S_0 = \frac{M_p c^2}{\sqrt{2}} \begin{bmatrix} e^{i\alpha_1} \cos \alpha_0 \\ e^{i\alpha_2} \sin \alpha_0 \\ i e^{i\alpha_1} \cos \alpha_0 \\ i e^{i\alpha_2} \sin \alpha_0 \end{bmatrix}$$

Antiparticles :

$$S_0 = \frac{M_p c^2}{\sqrt{2}} \begin{bmatrix} e^{i\alpha_1} \cos \alpha_0 \\ e^{i\alpha_2} \sin \alpha_0 \\ -ie^{i\alpha_1} \cos \alpha_0 \\ -ie^{i\alpha_2} \sin \alpha_0 \end{bmatrix}$$

To each particle corresponds an antiparticle with the same mass. And particles show polarization characteristics similar to waves. This is how the inertial spinor is seen in Quantum Physics : the 3 real variables $\alpha_0, \alpha_1, \alpha_2$ define a polarization of the particle when it behaves as a wave. Each kind of elementary particles is characterized by a vector of S_0 , and we will see that it is associated to its charge with respect to the EM field, and the 3 variables $\alpha_0, \alpha_1, \alpha_2$ define the magnetic moment. But for neutral material bodies usually its value can be taken equal to one of the vectors of the orthonormal basis :

$$\mathcal{E}_1 = \frac{1}{\sqrt{2}} (e_1 + \epsilon i e_3), \mathcal{E}_2 = \frac{1}{\sqrt{2}} (e_2 + \epsilon i e_4)$$

The space E_0, E'_0 are orthogonal, so :

$$\forall S_0 \in E_0, S'_0 \in E'_0 : \langle \gamma C(\sigma) S_0, \gamma C(\sigma) S'_0 \rangle = 0$$

The definition of E_0, E'_0 does not depend on the observer.

The spatial spinor S_r belongs to E_0, E'_0 by construct. The action of $Spin(3)$ on E_0, E'_0 is proper, continuous and free, thus (Maths.1793) the orbits have a unique structure of manifold of dimension : $\dim E_0 - \dim Spin(3) = 1$. For any value $S_0 \in E_0, E'_0$ when $\sigma_r \in Spin(3)$ then $S_r = \gamma C(\sigma_r) S_0$ stays on a curve on E_0, E'_0 , and conversely each vector S of E_0, E'_0 belongs to a unique such curve. $Spin(3)$ is compact, so this curve is compact, $Spin(3)$ has two connected components, so the curve is formed of two, compact, connected components.

The demonstration above is actually the equivalent - expressed in the formalism of fiber bundles and spinors - of the classic Wigner's classification of particles (see for instance Weinberg), done through the analysis of equivariance of the relativist momentum by the Poincaré's group. We could, in the same way, consider also the null spinors (assimilated to bosons), which is the vector subspace of E :

$$\langle S, S \rangle = 0 \Leftrightarrow v_L^t w_R = w_L^t v_R$$

However we will see another way to deal with bosons.

4.3.5 Space and time reversal

Space reversal and time reversal are changes of basis in \mathbb{R}^4 , so they depend on the basis.

A change of orthonormal basis in \mathbb{R}^4 is represented by an orthogonal matrix, and in the Clifford algebra by the action \mathbf{Ad}_s for some element s of the Pin group (it is not necessarily represented by an element of the connected component of the Spin group) :

$$w \rightarrow \tilde{w} = \mathbf{Ad}_s w = s \cdot w \cdot s^{-1}$$

The impact on a representation is :

$$\gamma(w) \rightarrow \gamma(\tilde{w}) = [\gamma(s)] [\gamma(w)] [\gamma(s)^{-1}]$$

In a change of basis in E represented by a matrix Q the components of a vector $u \in E$ change according to : $[u] \rightarrow [\tilde{u}] = Q^{-1}[u]$ and the matrices γ representing endomorphisms change as : $\gamma \rightarrow \tilde{\gamma} = Q^{-1}\gamma Q$. So the change of basis in \mathbb{R}^4 corresponds to a change of basis represented by the matrix $Q = [\gamma(s)]^{-1}$ in E , and the components of a vector S of E change as : $[u] \rightarrow [\tilde{u}] = [\gamma(s)][u]$.

Time reversal

Time reversal is the operation :

$$u = u^0\varepsilon_0 + u^1\varepsilon_1 + u^2\varepsilon_2 + u^3\varepsilon_3 \rightarrow -u^0\varepsilon_0 + u^1\varepsilon_1 + u^2\varepsilon_2 + u^3\varepsilon_3$$

corresponding to $s = \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3$, with $s^{-1} = \varepsilon_3 \cdot \varepsilon_2 \cdot \varepsilon_1$ in $Cl(3, 1)$, $s^{-1} = \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3$ in $Cl(1, 3)$

$Cl(3, 1)$:

$$[\gamma C(\varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3)] = \gamma_3 \gamma_2 \gamma_1 = i \begin{bmatrix} 0 & \sigma_0 \\ \sigma_0 & 0 \end{bmatrix}$$

$$i \begin{bmatrix} 0 & \sigma_0 \\ \sigma_0 & 0 \end{bmatrix} \begin{bmatrix} v \\ \epsilon iv \end{bmatrix} = \begin{bmatrix} -\epsilon v \\ iv \end{bmatrix} = \begin{bmatrix} v' \\ -\epsilon iv' \end{bmatrix}$$

$Cl(1, 3)$:

$$[\gamma C(\varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3)] = -i\gamma_1\gamma_2\gamma_3 = \begin{bmatrix} 0 & \sigma_0 \\ \sigma_0 & 0 \end{bmatrix}$$

$$\begin{bmatrix} 0 & \sigma_0 \\ \sigma_0 & 0 \end{bmatrix} \begin{bmatrix} v \\ \epsilon iv \end{bmatrix} = \begin{bmatrix} \epsilon iv \\ v \end{bmatrix} = \begin{bmatrix} v' \\ -\epsilon iv' \end{bmatrix}$$

So with both signatures particles and antiparticles are exchanged.

Space reversal :

Space reversal is the operation :

$$u = u^0\varepsilon_0 + u^1\varepsilon_1 + u^2\varepsilon_2 + u^3\varepsilon_3 \rightarrow u^0\varepsilon_0 - u^1\varepsilon_1 - u^2\varepsilon_2 - u^3\varepsilon_3$$

corresponding to $s = \varepsilon_0$, $s^{-1} = -\varepsilon_0$ in $Cl(3, 1)$, $s^{-1} = \varepsilon_0$ in $Cl(1, 3)$

$Cl(3, 1)$:

$$[\gamma C(\varepsilon_0)] = i\gamma_0 = \begin{bmatrix} 0 & \sigma_0 \\ -\sigma_0 & 0 \end{bmatrix}$$

$$\begin{bmatrix} 0 & \sigma_0 \\ -\sigma_0 & 0 \end{bmatrix} \begin{bmatrix} v \\ \epsilon iv \end{bmatrix} = \begin{bmatrix} \epsilon iv \\ -v \end{bmatrix} = \begin{bmatrix} v' \\ \epsilon iv' \end{bmatrix}$$

$Cl(1, 3)$:

$$[\gamma C(\varepsilon_0)] = \gamma_0 = \begin{bmatrix} 0 & -i\sigma_0 \\ i\sigma_0 & 0 \end{bmatrix}$$

$$\begin{bmatrix} 0 & -i\sigma_0 \\ i\sigma_0 & 0 \end{bmatrix} \begin{bmatrix} v \\ \epsilon iv \end{bmatrix} = \begin{bmatrix} \epsilon v \\ iv \end{bmatrix} = \begin{bmatrix} v' \\ \epsilon iv' \end{bmatrix}$$

So with both signatures particle and antiparticles stay in the same category.

These results are consistent with what is checked in Particles Physics, and the Standard Model. However the latter does not consider both signatures. Here we see that this feature does not allow to distinguish one signature as more physical than the other.

4.3.6 Total Spinor

The relation $S_L = \epsilon i S_R$ does not hold any more at the level of the total spinor, however we have still $\langle S, S \rangle = \langle S_0, S_0 \rangle_E$ which is positive for particles, and negative for anti-particles, so the distinction holds. The total spinor $S = \gamma C(\sigma) S_0 = \gamma C(\sigma_w) S_r$ belongs to a subset \hat{E}_0 of E larger than E_0 .

Definition 77 $\hat{E}_0 = \{ \gamma C(\sigma_w) S_0, \sigma_w \in Spin(3,1) / Spin(3), S_0 \in E_0 \}$
 $= \left\{ a_w S_0 - i \frac{1}{2} \sum_{a=1}^3 w_a \gamma_a \gamma_0 S_0, S_0 \in E_0 \right\}$

with a similar set \hat{E}'_0 for antiparticles.

The expression is :

$$\gamma C(a + v(r, w) + b\epsilon_5) \begin{bmatrix} S_R \\ \epsilon i S_R \end{bmatrix} = \begin{bmatrix} ((a+b) + \frac{1}{2} \sum_a (w_a - ir_a) \sigma_a) S_R \\ \epsilon i ((a-b) - \frac{1}{2} \sum_a (w_a + ir_a) \sigma_a) S_R \end{bmatrix}$$

thus with $r = 0$ we have :

$$\gamma C(a + v(0, w)) \begin{bmatrix} S_R \\ \epsilon i S_R \end{bmatrix} = \begin{bmatrix} (a + \frac{1}{2} \sum_a w_a \sigma_a) S_R \\ \epsilon i (a - \frac{1}{2} \sum_a w_a \sigma_a) S_R \end{bmatrix}$$

\hat{E}_0, \hat{E}'_0 are not vector spaces ($a_w = \epsilon \sqrt{1 + \frac{1}{4} w^t w}$) but real manifolds, embedded in E, with real dimension 7 (4 parameters for S_0 , 3 for w). By definition they are invariant by $Spin(3,1)$.

Moreover we have the following :

Theorem 78 For a given value of the inertial spinor S_0 , and a measured value $S \in \hat{E}_0$ of the spinor S, there is a unique element $\sigma \in Spin(3,1)$ such that $\gamma C(\sigma) S_0 = S$

Proof. i) The action of $Spin(3,1)$ on E_0, E'_0 is free :

$$\forall S_0 \in E_0, E'_0 : \gamma C(s) S_0 = S_0 \Leftrightarrow \sigma = 1$$

$$S_0 = \begin{bmatrix} v \\ \epsilon i v \end{bmatrix}$$

$$\gamma C(s) S_0 = S_0 \Leftrightarrow$$

$$S_R = \left((a + ib) \sigma_0 + \frac{1}{2} \sum_{a=1}^3 (w_a - ir_a) \sigma_a \right) v = v$$

$$S_L = \left((a - ib) \sigma_0 - \frac{1}{2} \sum_{a=1}^3 (w_a + ir_a) \sigma_a \right) \epsilon i v = \epsilon i v$$

$$\Rightarrow \left(2a\sigma_0 - i \sum_{a=1}^3 r_a \sigma_a \right) v = 2v$$

$$\left(2ib\sigma_0 + \sum_{a=1}^3 w_a \sigma_a \right) v = 0$$

$$\left(\sum_{a=1}^3 r_a \sigma_a \right) v = 2i(1-a)v$$

$$\left(\sum_{a=1}^3 w_a \sigma_a \right) v = -2ibv$$

$$\Rightarrow \sum_{a=1}^3 r_a v^* \sigma_a v = 2i(1-a)v^*v$$

$$\sum_{a=1}^3 w_a v^* \sigma_a v = -2ibv^*v$$

The scalars $v^* \sigma_a v$ are real because the Dirac matrices are Hermitian, as is $v^* v$, so

$$\Rightarrow b = 0, a = 1$$

$$\Rightarrow r = w = 0$$

and the only solution is $\sigma = 1$.

$$\text{ii) } \gamma C(\sigma) S_0 = \gamma C(\sigma') S_0 \Rightarrow S_0 = \gamma C(\sigma^{-1}) \gamma C(\sigma') S_0 \Rightarrow \sigma^{-1} \cdot \sigma' = 1 \quad \blacksquare$$

We can assume that S_0 depends only on the type of particle, then with the knowledge of S_0 the measure of the spinor S defines uniquely the motion (translation and rotation) with respect to the observer. As $S = \gamma C(\sigma) S_0$ and σ can itself be uniquely, up to spin, decomposed in $\sigma = \sigma_w \cdot \sigma_r$, we have the correspondence with the formulas in the transition between observers where $[\chi] = \exp[K(w)] \exp[J(r)]$: σ_w corresponds to the boost, and we have seen how it can be computed from the spatial speed, and σ_r corresponds to $\exp[J(r)]$, the vector r is in both cases a Lie algebra representative of a spatial rotation. So conversely, knowing S_0 the spinor S can be computed from familiar data.

4.4 SPINOR FIELDS

4.4.1 Definition

The great interest of Spinors is that they sum up the kinematics of a particle in one single, geometric quantity which has a value at any point in a fiber bundle. It is then possible to conceive fields of particles whose world lines are defined by the same vector field, which is an usual case in Physics.

We have defined previously the associated vector bundle $P_G [E, \gamma C]$, defined through the holonomic basis : $\mathbf{e}_i (m) = (\mathbf{p} (m), e_i)$ with the equivalence relation : $(\mathbf{p} (m), S) \sim (\mathbf{p} (m) \cdot g, \gamma C (g^{-1}) S)$.

Definition 79 We will call **Spinor field** a section $\mathbf{S} \in \mathfrak{X} (P_G [E, \gamma C])$ which represents the relativist momentum of a particle or an antiparticle, such that $\int_{\Omega} \|S (m)\| \varpi_4 < \infty$

Equivalently a spinor field, denoted $\mathfrak{X} (S_0)$ is defined by a vector $S_0 \in E_0$ and a section $\sigma \in \mathfrak{X} (P_G)$ such that $\mathbf{S} (m) = \gamma C (\sigma (m)) S_0$.

Let $\mathbf{S} \in \mathfrak{X} (P_G [E, \gamma C])$. Then $\langle S (m), S (m) \rangle = i (S_L^* S_R - S_R^* S_L) = y (m)$ defines a function on M. If \mathbf{S} represents the relativist momentum of a particle, then $\langle S (m), S (m) \rangle = \langle S_0, S_0 \rangle$. A necessary condition for a section of $P_G [E, \gamma C]$ to represent the relativist momentum of a particle is that $\langle S (m), S (m) \rangle = y$ has a fixed, positive, value. Then the set $E (y)$ of vectors $S_0 \in E_0$ such that $\exists \sigma \in \mathfrak{X} (P_G) : S = \gamma C (\sigma) S_0$ is given by :

$$E (y) = \left\{ S_0 = \begin{bmatrix} v \\ iv \end{bmatrix}, v^* v = \frac{1}{2} y, v \in \mathbb{C}^2 \right\}$$

And for a given vector $S_0 \in E (y)$, at each point m there is a unique $\sigma (m) \in Spin (3, 1)$ such that : $S (m) = \gamma C (\sigma (m)) S_0$. It defines a section $\sigma \in \mathfrak{X} (P_G)$, and at each point m , for each value ± 1 of the relativist spin and for a given observer, a vector field which is the tangent to the world line, and a spatial spin.

And similarly for antiparticles.

A spinor field represents particles which have the same inertial behavior. If, in a model, we have several particles, interacting with each others or with force fields, each particle can be assigned to a spinor field, which represents a general solution of the problem. One can also associate a density to spinor fields. All these topics will be seen in more details in the next chapters.

For elementary particles the vector S_0 is one of its fundamental characteristic, and we will see how this can be interpreted in the Standard Model.

For other material bodies S_0 is a kinematic characteristic which, for deformable solids, can be computed, as we will see below.

A Spinor field is defined without any reference to an observer : it has an intrinsic meaning, as it was expected for the momentum of a particle. And the decomposition in translational momentum on one hand, and rotational momentum on the other hand, is relative to each observer. So we have here a new, significant, feature of the relativist momentum.

Two comments :

i) A section of P_G can be seen as belonging to the Clifford bundle $Cl(M)$, and its components are (a, b, r, w) so we could hope to define the section through some fiber bundle with only (r, w) . Unfortunately the formulas for the components of an element of a $Cl(M)$ in a change of gauge are complicated and non linear. So, even if the use of r, w is useful, one cannot hope to define simply an consistently σ only through two vector fields $r(m), w(m)$.

ii) In a spinor field, a particle follows, with respect to an observer, a trajectory which is given by a vector field V deduced from a_w . So the value of the spinor at a point incorporates also the instantaneous velocity : at a given time $t + \delta t$ the particle is located at a point $\Phi_o(t + \delta t, x(t + \delta t))$ and $x(t + \delta t)$ is given by \vec{v} , which is part of V and defined by a_w , and thus by the value of the spinor at t .

We will see now how the usual spin enters the picture, but for this we need some more mathematics on group representations.

4.4.2 More on the theory of the representations of groups

Functional Representations

Functional representations are representations on vector spaces of functions or maps. Any locally compact topological group has at least one unitary faithful representation (usually infinite dimensional) of this kind, and they are common in Physics. The principles are the following (Maths.23.2.2).

Let H be a Banach vector space of maps $\varphi : E \rightarrow F$ from a topological space E to a vector space F , G a topological group with a continuous left action λ on $E : \lambda : G \times E \rightarrow E :: \lambda(g, x)$ such that $\lambda(g \cdot g', x) = \lambda(g, \lambda(g', x)), \lambda(1, x) = x$

Define the left action Λ of G on $H : \Lambda : G \times H \rightarrow H :: \Lambda(g, \varphi)(x) = \varphi(\lambda(g^{-1}, x))$

Then (H, Λ) is a representation of G . Thus G acts *on the argument* of φ .

If H is a Hilbert space and G has a Haar measure μ (a measure on G , all the groups that we will encounter have one Maths.22.5) then the representation is unitary with the scalar product :

$$\langle \varphi_1, \varphi_2 \rangle = \int_G \langle \Lambda(g, \varphi_1), \Lambda(g, \varphi_2) \rangle_H \mu(g)$$

If G is a Lie group and the maps of H and λ are differentiable (which implies that E is a manifold) then $(H, \Lambda'_g(1, \cdot))$ is a representation of the Lie algebra T_1G where $X \in T_1G$ acts by a differential operator :

$$\Lambda'_g(1, \varphi)(X)(x) = -\varphi'(x) \lambda'_g(1, x) X = \frac{d}{dt} \varphi(\lambda(\exp(-tX), x)) |_{t=0}$$

For a right action $\rho : E \times G \rightarrow E :: \rho(g, x)$ we have similar results, with

$$P : H \times G \rightarrow H :: P(\varphi, g)(x) = \varphi(\rho(x, g))$$

$$P'_g(\varphi, 1)(X)(x) = -\varphi'(x) \rho'_g(x, 1) X = \frac{d}{dt} \varphi(\rho(x, \exp(-tX))) |_{t=0}$$

H can be a vector space of sections on a vector bundle. In a functional representation each function is a vector of the representation, so it is usually infinite dimensional. However the representation can be finite dimensional, by taking polynomials as functions, but this is not always possible : the set of polynomials must be algebraically closed under the action of the group.

Isomorphisms of groups

Most of the groups that are encountered in Physics are related to the group $SL(\mathbb{C}, 2)$ of 2×2 complex matrices with determinant 1 (Maths.24).

Any matrix of the Lie algebra $sl(\mathbb{C}, 2)$ reads with $Z = (z_1, z_2, z_3) \in \mathbb{C}^3$

$$f(Z) = \begin{bmatrix} iz_3 & z_2 + iz_1 \\ -z_2 + iz_1 & -iz_3 \end{bmatrix} \Rightarrow \text{Tr} f(Z) = 0$$

which is equivalent to take as basis the Dirac matrices.

The exponential is not surjective on $sl(\mathbb{C}, 2)$ and any matrix of $SL(\mathbb{C}, 2)$ reads :

$$I \cosh D + \frac{\sinh D}{D} f(Z) \text{ with } D^2 = -\det f(Z) = -(z_1^2 + z_2^2 + z_3^2)$$

The group $SU(2)$ of 2×2 unitary matrices ($NN^* = I$) is a compact real subgroup of $SL(\mathbb{C}, 2)$. Its Lie algebra is comprised of matrices $f(r)$ with $r \in \mathbb{R}^3$.

The exponential is surjective on $SU(2)$: $\exp f(r) = I \cos \sqrt{r^t r} + \frac{\sin \sqrt{r^t r}}{\sqrt{r^t r}} f(r)$

$T_1 Spin(3, 1)$ is isomorphic to $sl(\mathbb{C}, 2)$ (Math.1959) : $v(r, w) \rightarrow f(r + iw)$

$Spin(3, 1)$ is isomorphic to $SL(\mathbb{C}, 2)$: $a + v(r, w) + b\varepsilon_5 \rightarrow \exp f(r + iw)$

$T_1 Spin(3)$ is isomorphic to $su(2)$: $v(r, 0) \rightarrow f(r)$ and $so(3)$: $v(r, 0) \rightarrow j(r)$

$Spin(3)$ is isomorphic to $SU(2)$:

$$a_r + v(r, 0) \rightarrow \exp f(r) = I \cos \sqrt{r^t r} + \frac{\sin \sqrt{r^t r}}{\sqrt{r^t r}} f(r)$$

Representations of Spin(3,1), Spin(3) and SO(3)

$SL(\mathbb{C}, 2)$ and $Spin(3, 1)$ have the same representations which are (up to equivalence) :

- a unique, non unitary, irreducible representation of dimension n (Maths.1953), which can be seen as the tensorial product of two finite dimensional representations ($P^j \otimes P^k, D_j \times D_k$) of $SU(2) \times SU(2)$ (see below).

- the only unitary representations are over spaces of complex functions : they are infinite dimensional and each irreducible representation is parametrized by 2 scalars $z \in \mathbb{Z}, k \in \mathbb{R}$ (Maths.1955).

$SU(2)$ as $Spin(3)$ are compact groups, so their unitary representations are reducible (Math.1960) in a sum of orthogonal, finite dimensional, unitary representations. The only irreducible, finite dimensional, unitary, representations, denoted (P^j, D^j) are on the space P^j of degree $2j$ homogeneous polynomials with 2 complex variables z_1, z_2 , where conventionally j is an integer or half an integer. P^j is $2j + 1$ dimensional and the elements of an orthonormal basis are denoted :

$$|j, m\rangle = \frac{1}{\sqrt{(j-m)!(j+m)!}} z_1^{j+m} z_2^{j-m} \text{ with } -j \leq m \leq +j. \text{ And } D^j \text{ is defined by}$$

:

$$g \in U(2) : D^j(g) P \left(\begin{bmatrix} z_1 \\ z_2 \end{bmatrix} \right) = P \left([g]^{-1} \begin{bmatrix} z_1 \\ z_2 \end{bmatrix} \right)$$

Thus the functions read : $\varphi(z_1, z_2) = \sum_{j \in \frac{1}{2}\mathbb{Z}} \sum_{m=-j}^{m=+j} \varphi^{jm} |j, m\rangle$ with complex constants φ^{jm}

It induces a representation (P^j, d^j) of the Lie algebras where d^j is a differential operator acting on the polynomials P :

$$X \in su(2) : d_j(X)(P)(z_1, z_2) = \frac{d}{dt} P \left([\exp(-tX)] \begin{bmatrix} z_1 \\ z_2 \end{bmatrix} \right) |_{t=0}$$

which gives, for polynomials, another polynomial.

$d_j(X)$ is a linear map on P^j , which is also linear with respect to X , thus it is convenient to define d_j by the action $d_j(\kappa_a)$ of a basis $(\kappa_a)_{a=1}^3$ of the Lie algebra and the three operators are denoted L_x, L_y, L_z . They are expressed in the orthonormal basis $|j, m\rangle$ by square $2j+1$ matrices (depending on the conventions to represent the Lie algebra). The usage is to denote $L_z |j, m\rangle = m |j, m\rangle$.

The only irreducible, unitary, representations of $SO(3)$ are given by (P^j, D^j) with j integer.

Casimir element

The **universal enveloping algebra** U of a Lie algebra is actually a vector space, built from *tensorial powers* of the Lie algebra, and whose basis is given by ordered products of elements of the basis $(\kappa_i)_{i \in I}$ of the Lie algebra (Maths.1692). Universal enveloping algebras are necessary when interacting systems are considered (such as in Chemistry), because their representation involve the tensorial product of the variables.

Any representation (E, f) of the Lie algebra can be extended to a representation (E, F) of its universal enveloping algebra (Maths.1891) where the action is :

$$F \left(\kappa_{i_1}^{n_1} \dots \kappa_{i_p}^{n_p} \right) = f(\kappa_{i_1})^{n_1} \circ \dots \circ f(\kappa_{i_p})^{n_p}$$

When the representation (E, f) comes from a functional representation, in the induced representation on U the action of F is represented by differential operators, of the same order than $n_1 + n_2 + \dots + n_p$.

In the representation of $T_1 Spin(3, 1)$ by matrices of $so(3, 1)$ the universal enveloping algebra is actually an algebra of matrices (see Annex) where the operator j plays a key role.

The **Casimir element** is a special element Ω of U, defined through the Killing form (Maths.1698). In an irreducible representation (E, f) of a semi simple Lie algebra, as $Spin(3, 1)$, the image of the Casimir element acts by a non zero fixed scalar $F(\Omega)u = ku$. In functional representations it acts by a differential operator of second order : $F(\Omega)\varphi(x) = D_2\varphi(x) = k\varphi(x)$: φ is an eigen vector of D_2 . As a consequence, if there is a scalar product on E : $\langle F(\Omega)u, F(\Omega)u \rangle = \langle ku, ku \rangle = k^2 \langle u, u \rangle$. If $(E_1, f_1), (E_2, f_2)$ are two equivalent representations of the same algebra A :

$$\begin{aligned} \exists \phi : E_1 &\rightarrow E_2 \text{ such that :} \\ \forall \kappa \in A : f_1(\kappa) &= \phi^{-1} \circ f_2(\kappa) \circ \phi \end{aligned}$$

$$\begin{aligned}
F_1 \left(\kappa_{i_1}^{n_1} \dots \kappa_{i_p}^{n_p} \right) &= f_1 \left(\kappa_{i_1} \right)^{n_1} \circ \dots \circ f_1 \left(\kappa_{i_p} \right)^{n_p} = \left(\phi^{-1} \circ f_2 \left(\kappa_{i_1} \right) \circ \phi \right)^{n_1} \circ \dots \circ \\
\left(\phi^{-1} \circ f_2 \left(\kappa_{i_p} \right) \circ \phi \right)^{n_p} &= \phi^{-n_1 - \dots - n_p} \circ F_2 \left(\kappa_{i_1}^{n_1} \dots \kappa_{i_p}^{n_p} \right) \circ \phi^{n_1 + \dots + n_p} \\
F_1 \left(\Omega \right) &= \phi^{-n_1 - \dots - n_p} \circ F_2 \left(\kappa_{i_1}^{n_1} \dots \kappa_{i_p}^{n_p} \right) \circ \phi^{n_1 + \dots + n_p} \left(u \right) = \phi^{-n_1 - \dots - n_p} \circ \left(k_2 \phi^{n_1 + \dots + n_p} \left(u \right) \right) = \\
k_2 u &= k_1 u
\end{aligned}$$

Thus the Casimir element acts with the same scalar in all equivalent representations.

The Killing form on $T_1 Spin(3, 1)$ is :

$$B(v(r, w), v(r', w')) = 4(w^t w' - r^t r')$$

thus the elements

$$\kappa_1 = -\frac{1}{8}\varepsilon_3 \cdot \varepsilon_2, \kappa_2 = -\frac{1}{8}\varepsilon_1 \cdot \varepsilon_3, \kappa_3 = -\frac{1}{8}\varepsilon_2 \cdot \varepsilon_1,$$

$$\kappa_4 = \frac{1}{8}\varepsilon_0 \cdot \varepsilon_1, \kappa_5 = \frac{1}{8}\varepsilon_0 \cdot \varepsilon_2, \kappa_6 = \frac{1}{8}\varepsilon_0 \cdot \varepsilon_3$$

constitute an orthonormal basis for B and the Casimir element of $U(T_1 Spin(3, 1))$

is :

$$\Omega = \left(\sum_{i=4}^6 (\kappa_i)^2 - \sum_{i=1}^3 (\kappa_i)^2 \right)$$

The action of the Casimir element in the representation $(E, \gamma C)$ of $Spin(3, 1)$

is :

$$F_E(\Omega)u = \left(\sum_{i=4}^6 (\gamma C(\kappa_i))^2 - \sum_{i=1}^3 (\gamma C(\kappa_i))^2 \right) u = \frac{3}{2}u$$

In the representation (P^j, d^j) of $T_1 Spin(3)$, if we denote $L_x = f(\kappa_1), L_y = f(\kappa_2), L_z = f(\kappa_3)$ with 3 arbitrary orthogonal axes :

$$F(\Omega)|j, m\rangle = L^2|j, m\rangle = (L_x^2 + L_y^2 + L_z^2)|j, m\rangle = j(j+1)|j, m\rangle$$

$$d^j(\kappa_i) \left(\sum_{m=-j}^{m=+j} X^m |j, m\rangle \right) = \sum_{m=-j}^{m=+j} X^m d^j(\kappa_i) |j, m\rangle$$

4.4.3 The Spin of a particle

Definition

The space $\mathfrak{X}(P_G[E, \gamma C])$ of sections is a functional representation of $Spin(3, 1)$ with the global action γC and the argument σ . The subspace $\mathfrak{X}(S_0)$ is invariant by the right or left global actions of $Spin(3, 1)$: $\gamma C(\sigma(m))S_0 \rightarrow \gamma C(s)\gamma C(\sigma(m))S_0$ or $\gamma C(\sigma(m))S_0 \rightarrow \gamma C(\sigma(m))\gamma C(s)S_0$. In particular it is invariant by the action of $Spin(3)$:

$$\rho : \mathfrak{X}(S_0) \times Spin(3) \rightarrow \rho(S(m), s_r) = \gamma C(\sigma(m) \cdot s_r) S_0$$

Moreover the value of $Y(m) = \langle S(m), S(m) \rangle$ is invariant by $Spin(3, 1)$.

The spinor fields $S \in \mathfrak{X}(S_0)$ can equivalently be defined by a couple (S_0, σ) where $\sigma \in \mathfrak{X}(P_G)$. For a given observer each $\sigma(m)$ has two decompositions : $\sigma(m) = \epsilon \sigma_w(m) \cdot \epsilon \sigma_r(m)$ so the couple (S_0, σ) defines precisely two Spatial Spinor fields : $S_r(m) = \gamma C(\sigma_r(m))S_0$.

Conversely one can define Spatial Spinor Fields by a couple : (S_0, σ_r) where $\sigma_r \in \mathfrak{X}(P_R)$ and they constitute a set $\mathfrak{X}_r(S_0)$ which is invariant by $Spin(3)$ (but not by $Spin(3, 1)$).

Let us denote : $\pi_\epsilon : \mathfrak{X}(S_0) \rightarrow \mathfrak{X}_r(S_0)$ the maps which associates, for a given observer, to each Spinor field the Spatial Spinor field with $\epsilon = \pm 1$

On the set : $\mathfrak{X}(S_0)$ we can define the equivalence relation :

$$S \sim S' \Leftrightarrow \pi_\epsilon(S) = \pi_\epsilon(S')$$

Each class of equivalence is the set of spinor fields which have, for the observer, the same kinematic behavior with regard to a rotation by $Spin(3)$. The value of $\pi_\epsilon(S)$ for a given spinor field is **the Spin of the particle**, in its usual meaning. So to any given spinor field corresponds, for an observer, two Spins, with the Spin up or down. And conversely for a given Spin there can be infinitely many spinor fields, defined by a section of the associated bundle $\sigma_w \in \mathfrak{X}(P_W)$.

The projection π_ϵ depends on the choice of a vector ϵ_0 , so the Spin depends on the observer and its measure depends on the spatial basis he has chosen. The spin can be seen as a rotational momentum.

Similarly we have the projection : $\pi_w : Spin(3,1) \rightarrow Spin(3,1)/Spin(3)$ and we can define the equivalence relation in $\mathfrak{X}(S_0)$:

$$S = \gamma C(\sigma) S_0 \sim S' = \gamma C(\sigma') S_0 \Leftrightarrow \pi_w(\sigma) = \pi_w(\sigma')$$

The class of equivalence represents the particles which have the same trajectories. And we define the **translational momentum** by :

$$S_w = \pi_w(\sigma) S_0$$

This is also a geometric quantity, invariant by $Spin(3,1)$, but observer dependant.

Quantization of the Spinor

Theorem 80 *The set $L^1(S_0) = L^1(M, P_G[E, \gamma C], \varpi_4) \cap \mathfrak{X}(S_0)$ of integrable spinor fields associated to a particle is characterized by 2 scalars : $k \in \mathbb{R}, z \in \mathbb{Z}$.*

The Spin, up or down, associated to each section by an observer is characterized by a scalar $j \in \frac{1}{2}\mathbb{N}$ and belongs to a $2j+1$ dimensional vector space : $S_r(m) = \sum_{p=-j}^{+j} S_r^p |j, p\rangle$ with the constant components S_r^p and an orthonormal basis $|j, p\rangle$

Proof. i) The space $L^1(M, P_G[E, \gamma C], \varpi_4)$ is a Fréchet space. The Theorem 2 applies and there are a Hilbert space H and an isometry $\Upsilon : L^1 \rightarrow H :: \psi = \Upsilon(S)$. Moreover $(L^1, \gamma C)$ is an infinitely dimensional representation of $Spin(3,1)$ (the scalar product, thus the norm, is invariant by $Spin(3,1)$, and L^1 is invariant by $Spin(3,1)$). We can apply the theorem 22. $(H, \hat{\gamma})$ is a unitary representation of $Spin(3,1)$ with $\hat{\gamma}(\sigma) = \Upsilon \circ \gamma C(\sigma) \circ \Upsilon^{-1}$.

ii) Consider the function : $Y : L^1 \rightarrow \mathbb{C} :: Y = \langle S, S \rangle$. For a given section, Y has a value at each point of M and Y is invariant by $Spin(3,1)$. We can implement the theorem 24: to each value y of Y is associated a class of equivalence in L^1 and in H .

If we fix $Y = \langle S_0, S_0 \rangle = Ct$ we have two subsets

$$L^1(S_0) = L^1 \cap \mathfrak{X}(S_0) \text{ in } L^1 \text{ and } H(S_0) \text{ in } H.$$

$H(S_0)$ is invariant by $\hat{\gamma}$

iii) The unitary representations of $Spin(3,1)$ read : $H = \oplus_{z,k} H_{z,k} \oplus H_c$ where $H_{z,k}$ are unitary irreducible representations, defined by the parameters

$z \in \mathbb{Z}, k \in \mathbb{R}$, and H_c does not contain any irreducible representation, so H_c is not invariant under the action of $Spin(3, 1)$ (Maths.1914).

As a consequence $H(S_0)$ is isomorphic to a subset of one of the irreducible representations $H_{z,k}$ and the spinor field is characterized by two scalars $k \in \mathbb{R}, z \in \mathbb{Z}$ linked to S_0 .

iv) In L^1 , for each section S and a given observer, the Spatial Spinor S_r is a representation of $Spin(3)$. Moreover for S_0, ϵ fixed it belongs to one of the irreducible representations of $Spin(3)$. It is isomorphic to one of the representations (P^j, D^j) with $j \in \frac{1}{2}\mathbb{N}$. These representations are finite dimensional, so S_r belongs to a $2j + 1$ dimensional vector space : $S_r(m) = \sum_{p=-j}^{+j} S_r^p |j, p\rangle$ with the constant components S_r^p . ■

Assume that we study a system comprising of unknown particles $p = 1 \dots N$. The modeling of their kinematic characteristics leads naturally to assume that these particles belong to some spinor field : $S_p \in \mathfrak{X}(P_G[E, \gamma C])$ with unknown inertial spinor. Because no value of S_0 is imposed we have a vector space and we can implement the theorem 2.

What the theorem above tells us is that the solutions must be found in maps : $S_p : \Omega \rightarrow E$ which can be sorted out :

- by the value of $\langle S_0, S_0 \rangle$, that is their mass
- by the value of some integer $z \in \mathbb{Z}$
- and their spin by a half integer $j \in \frac{1}{2}\mathbb{N}$

They correspond to particles which have the same behavior when submitted to a force field (they have the same world lines and spatial spinor for any observer). In other words the spinor is not the only characteristic which determines the behavior of a particle, and these others characteristics can be labeled by a signed integer. This is the starting point to the representation of charged particles that we will see in the next chapter.

For elementary particles it is experimentally seen that $j = \frac{1}{2}$, and this is the origin of the name "particles of spin $\frac{1}{2}$ ". For composite particles or nuclei the spin can be higher.

Even if the set $\mathfrak{X}(S_0)$ is not a vector space, it is a manifold which is embedded in a vector space, so that each of its points (a map S_p) can be written as a fixed linear combination of vectors of a basis. The vector space is always infinite dimensional for the translational momentum, but each spin belongs to a finite dimensional vector space, which is isomorphic to some $(P^j, D^j) : S_r(m) = \sum_{p=-j}^{+j} S_r^p |j, p\rangle$ where S_r^p are fixed scalars and $|j, p\rangle$ are, for a given system, fixed maps $|j, p\rangle : \Omega \rightarrow E_0$, images of vectors of the basis of P^j by some isometry. Each vector $|j, p\rangle$ is assimilated to a state of the particle, and j, p are the quantum numbers labeling the state. The maps $|j, p\rangle$ are not polynomials (as in P^j), they are used only to define the algebraic structure of the space $H(S_0)$, however they have an interpretation for models of atoms (see below). Under the action of $Spin(3)$ the vectors $S_r(m)$ transform according to the same matrices as in D^j :

$\gamma C(\sigma(m) \cdot s_r) S_0 = \sum_{p=-j}^{+j} S_r^p [D^j(f(s_r))] |j, p\rangle$ where $f(s_r)$ is the image of σ in $SU(2)$.

By itself the theorem does not provide a solution : a vector of a basis of the vector spaces is itself some map $E_i : \Omega \rightarrow E$. But it shows that the solution cannot take any value, even before we implement any physical law relating the fields and the kinematic characteristics. In a given system the solutions that appear follow the same pattern, whatever the initial conditions, or the value of the other variables (notably the fields).

There is one important difference in the behavior of the spin, according to the value of j . The Spin is invariant by a rotation by $Spin(3)$, and the scalars $\pm 1 \in Spin(3)$. The actions of $+s$ and $-s$ give opposite results. $Spin(3)$ is the double cover of $SO(3)$: to the same element g of $SO(3)$ are associated two elements $\pm s$ of $Spin(3)$. The representations (P^j, D_j) with $j \in \mathbb{N}$ are also representations of $SO(3)$. It implies that the vector spaces are invariant by $\pm s$. The fact that j is an integer means that *the particle has a physical specific symmetry* : the rotations $\pm s$ give the same result. And equivalently, if j is half an integer the rotations by $\pm s$ give opposite results.

Measure of the spatial spin of a particle

A particle has, whatever the scale, by definition, no internal structure, so it is impossible to observe its geometric rotation. However it has a spin, its spatial spinor S_r is a variable which can be represented in a finite dimensional space : S_r is an observable. The measure of the spatial spinor, similar to a rotational momentum, is done by observing the behavior of the particle when it is submitted to a force field which acts differently according to the value of the spinor. This is similar to the measure of the rotation of a perfectly symmetric ball by observing its trajectory when it is submitted to a dissymmetric initial impulsion (golfers will understand).

Most particles have a magnetic moment, linked to their spin. So the usual way to measure the latter is to submit the particles to a non homogeneous magnetic field. This is the Stern-Gerlach analyzer described in all handbooks, where particles have different trajectories according to their magnetic moment. MRI uses a method based on the same principle with oscillating fields whose variation is measured. The process can be modelled as follows.

The spinors of the particles are represented by some section $\mathcal{S} \in \mathfrak{X}(S_0)$. The device operates only on the spin : $S_r(m) = \gamma C(\sigma_r(m)) S_0$ and is parametrized by a spatial rotation $s_r \in Spin(3)$, and usually by a vector $\rho \in \mathbb{R}^3$, corresponding to a rotation s_r .

The first effect is a breakdown of symmetry : s_r has not the same impact for the particles with spin up or down. This manifests by two separate beams in the Stern-Gerlach experiment.

An observable $\Phi(S_r)$ of S_r is a projection on some finite dimensional vector space of maps. Because of the quantization, this vector spaces has for vectors $|j, p\rangle$ which are fixed maps, image of the vectors of basis of P^j which are eigen vectors of the observable. The action of the device can be modelled as an operator $L(\rho)$ acting on this space, and the matrices to go from one orientation ρ_1 to another ρ_2 are the same as in (P^j, d^j) . It reads :

$$L(\rho) \Phi(S_r) = \sum_{p=-j}^{+j} S_r^p [d^j(\rho)] |j, p\rangle$$

For a given beam we have a breakdown of the measures, corresponding to each of the states labelled by p.

Arbitrary axes x, y, z are chosen for the device, which provide 3 measures $L_x(S_r), L_y(S_r), L_z(S_r)$, such that $L_z(S_r) |j, m\rangle = m |j, m\rangle$.

The Casimir operator Ω is such that $L^2\Phi(S_r) = (L_x^2 + L_y^2 + L_z^2)(S_r) = j(j+1)\Phi(S_r)$

Atoms and electrons

QM has been developed from the study of atoms, with a basic model (Bohr's atom) in which electrons move around the nucleus. Even if this idea still holds, and this is how atoms are commonly viewed, it had been quickly obvious that a classic model does not work. However using what has been developed previously, we can have another representation.

Let us consider a system comprised of one electron moving around a nucleus. If we consider the atom as a particle, that is without considering its internal structure, its relativist momentum can be represented by a spinor S , and its rotational momentum by a spin S_r . The previous results hold and the spin can be represented in a finite dimensional vector space isomorphic to P_j . However j , which belongs to $\frac{1}{2}\mathbb{N}$, is not necessarily equal to $\frac{1}{2}$.

As noticed before, the polynomials P^j have no physical meaning. However in this case it is usual to provide one. By a purely mathematical computation it is possible to show that the representation (P^j, D^j) is equivalent to a representation on square integrable functions $f(x)$ on \mathbb{R}^3 , and from there on spherical harmonic polynomials (Maths.1958). It is then assumed that the arguments of the function $f(x)$ are related to the coordinates (in an euclidean frame) of the electron. This is a legacy of the first models of atoms. Actually there is no need for such an assumption to build a consistent model, which would be useless in the GR context, and the image of electrons rotating around a nucleus and spinning has no physical support.

For atoms with several electrons, the model must involve the tensorial products of each spinor. The previous representations of $SU(2)$ are then extended to the tensorial products of P^j , and their derivative to representations of the universal enveloping algebra. It is often possible to rearrange these representations, by combinations using Clebsch-Jordan coefficients (Maths.1960), and in this endeavour the spherical harmonic polynomials are useful because they provide many identities. This is one major application of QM in Chemistry.

The same kind of model is used for composite particles in Quantum Theory of Fields.

4.4.4 Material bodies and spinors

Representation of a material body by sections of P_G

We have seen in the previous chapter that a material body B can be defined, from a geometric point of view, by a vector field u whose integral curves are the world lines of its particles. Then the flow $\Phi_u(\tau, a)$ defines the body B itself at each proper time τ as a compact subset $\omega(\tau)$ of a 3 dimensional hypersurface. And there are privileged observers B for whom $\omega(0) \subset \Omega(0)$.

So a material body can be defined with respect to these observers, up to a constant ± 1 by a section σ_w of P_W or, up to a spatial spinor, by a section of P_G and a compact, space like hypersurface $\omega(0)$. Then σ_w provides u , Φ_u and $\omega(0)$ defines $\omega(\tau)$. The section σ_w can be seen as the general definition of B , which can be fitted to any initial conditions $\omega(0)$. This is the most efficient way to define geometrically a material body in physical models.

Spinors representing a solid

We have seen that the usual concepts of motion of a body over itself (usually a rotation of the body) cannot be easily represented in relativist geometry. This is the main motivation for the introduction of spinors, and any material body whose internal structure can be neglected (at the scale of the study) can be represented, from the kinematic point of view, by a spinor which accounts for its rotation (through the spin). If the location of the material body can be represented by a geometric point, then the kinematic representation of B is given by a map : $S_B : \mathbb{R} \rightarrow P_G[E, \gamma C]$, such that $S_B(t) = \gamma C(\sigma(t)) S_0$. We do not need more : S_B provides everything, including the rotational momentum. Thus, even if no internal structure or rotation of the body is assumed, eventually it can be accounted for.

However this representation assumes that S_0 is known. As in Classic Mechanics for the inertial tensor, the computation of the inertial spinor S_0 is, for a solid, a separate issue. It can be done through the aggregation of material points (particles) with a specific law giving the shape and the density of the body. And the inertial spinor is not necessarily constant : we can consider deformable solids. Actually we can define a *rigid solid as a material body such that S_0 is constant.*

Proposition 81 *A deformable solid body can be represented by a map :*

$$S : \mathbb{R} \rightarrow P_G[E, \gamma C] \text{ such that } \langle S(\tau), S(\tau) \rangle > 0 \text{ or } \langle S(\tau), S(\tau) \rangle < 0$$

A rigid solid body can be represented by a map :

$$S : \mathbb{R} \rightarrow P_G[E, \gamma C] :: S(\tau) = \gamma C(\sigma(\tau)) S_0 \text{ for a fixed } S_0 \in E_0 \text{ or } S_0 \in E'_0$$

where τ is the proper time of the body

To assume that the material points behave in a coherent way in a solid assumes that there are forces which assure this cohesion. And indeed a material body can be deformed or broken. So we can say that the fact, assumed and which can be checked, that a material body can be represented by a unique spinor incorporates the existence of these internal forces. And ultimately the break of a material body can result in several spinors. So in modelling the evolution of a material body we should include additional assumptions about the laws (which are similar to the phenomenological laws for deformable solids) for the change of S_0 . And in a discontinuous process add the laws which rules the splitting in different spinors.

Aggregating matter fields

With these definitions we can consider the task to compute the spinor that we will denote S_B , for a deformable solid, by aggregating material points. This is similar to the computation of the inertial tensor in Classic Mechanics : this is a specific endeavour, done in a separate model, using specific assumptions (about the shape, density, motion of the particles) and the result is then used in a more general model (for instance to compute the motion of different bodies). The single spinor corresponding to the whole body is assigned, in the more general model, to any point : all the material points have then the same location.

The first issue is the definition of the motion of the material points with respect to the body. We need a chart to do it, which is given by an observer B, such that at his proper time $t = 0$ the set $\omega(0)$ is in his present $\Omega(0)$. Then at any given time t the set of particles constituting the solid stays in his present. B uses his standard chart :

$\varphi_B(t, \eta^1, \eta^2, \eta^3) = \Phi_{\varepsilon_0}(t, x(\eta^1, \eta^2, \eta^3))$ where $x(\eta^1, \eta^2, \eta^3)$ is a chart on $\omega(0)$ and ε_0 his time like vector field

The particles follow the trajectories given by a vector field V and their location at t is $\Phi_V(t, a) = \Phi_{\varepsilon_0}(t, x(t)) = \varphi_B(t, x(t))$ with $x(0) = a$.

$\omega(t) = \{\Phi_V(t, a), a \in \omega(0)\} = \{\Phi_{\varepsilon_0}(t, x(t)), x(0) \in \omega(0)\}$ represents the location of the body at t and $\Omega = \{\Omega(t), t \in [0, T]\}$

The material points are represented by a section $S(t, x) = \gamma C(\sigma(t, x)) S_0 \in \mathfrak{X}(S_0)$. The choice of S_0 can be arbitrary. S is a geometric quantity which does not depend on a chart, however σ_w provides a vector field of world lines u for the material points with respect to ε_0 .

We assume that the observer defines a tetrad $(\varepsilon_i(m))$ from which the metric and the volume form are deduced in the usual way.

The density $\mu(t, x)$ is defined over Ω with respect to the volume form ϖ_4 . Because $\omega(t) \subset \Omega(t)$ the unitary, future oriented, normal to $\omega(t)$ is $\varepsilon_0 = \partial\xi_0$ and μ induces a density $\mu_3(t, x)$ over $\omega(t)$ with respect to the volume form ϖ_3 :

$\mu_3(t, x) \varpi_3(t, x) = i_V(\mu(t, x) \varpi_4(t, x))$ which is the flux of matter going through $\omega(t)$.

As noticed before the holonomic basis $e_i(m) = (\mathbf{p}(m), \varepsilon_i)$ of $P_G[E, \gamma C]$ is

arbitrary, in that there is no physical reference for the choice of the vectors e_i . We can assume that, for a deformable solid, there is a common basis associated to the chart Φ_{ε_0} .

$$S(t, x) = \gamma C(\sigma(t, x)) S_0$$

Then the integral :

$$\int_{\omega(t)} \gamma C(\sigma(t, x)) \mu_3(t, x) S_0 \varpi_3(t, x) \\ = \left[\gamma C \left(\int_{\omega(t)} (\sigma(t, x)) \mu_3(\tau, x) \varpi_3(t, x) \right) \right] S_0$$

is well defined on the fixed vector space \mathbb{E} .

$$\sigma(t, x) = a(t, x) + v(r(t, x), w(t, x)) + b(t, x) \varepsilon_5$$

with the identities :

$$a(t, x)^2 - b(t, x)^2 = 1 + \frac{1}{4}(w^t w - r^t r)$$

$$a(t, x) b(t, x) = -\frac{1}{4} w^t r$$

Denote

$$\hat{r}(t) = \int_{\omega(t)} \mu_3(x, t) r(t, x) \varpi_3(t, x), \hat{w}(t) = \int_{\omega(t)} \mu_3(x, t) w(t, x) \varpi_3(t, x)$$

$$\hat{a}(t) = \int_{\omega(t)} a(t, x) \mu_3(t, x) \varpi_3(t, x), \hat{b}(t) = \int_{\omega(t)} b(t, x) \mu_3(t, x) \varpi_3(t, x)$$

$$\int_{\omega(t)} \mu_3(x, t) \sigma(t, x) \varpi_3(t, x) = \hat{a}(t) + v(\hat{r}(t), \hat{w}(t)) + \varepsilon_5 \hat{b}(t)$$

We impose, for a deformable solid, that :

$$\exists N(t) \in \mathbb{R}, R(t), W(t) \in \mathbb{R}^3 :$$

$$\hat{a}(t) + v(\hat{r}(t), \hat{w}(t)) + \varepsilon_5 \hat{b}(t) = N(t) (A(t) + v(R(t), W(t)) + B(t) \varepsilon_5) \quad (4.51)$$

such that $\sigma_B(t) = A(t) + v(R(t), W(t)) + B(t) \varepsilon_5 \in Spin(3, 1)$ which requires :

$$A^2 - B^2 = 1 + \frac{1}{4}(W^t W - R^t R)$$

$$AB = -\frac{1}{4} W^t R$$

and implies :

$$R(t) = \frac{1}{N} \hat{r}(t), W(t) = \frac{1}{N} \hat{w}(t)$$

$$A(t) = \frac{1}{N} \hat{a}(t), B(t) = \frac{1}{N} \hat{b}(t)$$

Which sums up to the two conditions :

$$\hat{a}(t) \hat{b}(t) = -\frac{1}{4} \hat{w}^t \hat{r} \quad (4.52)$$

$$\hat{a}^2 - \hat{b}^2 = N^2 + \frac{1}{4} (\hat{w}^t \hat{w} - \hat{r}^t \hat{r}) \quad (4.53)$$

$$\Rightarrow N^2 = \hat{a}^2 - \hat{b}^2 - \frac{1}{4} (\hat{w}^t \hat{w} - \hat{r}^t \hat{r}) > 0$$

Then the Spinor of the body is : $S_B(t) = N(t) \gamma C(\sigma_B(t)) S_0$

The conditions can be seen as resulting from the forces which keep the cohesion of the body.

The mass of the solid is proportional to

$$\langle S_B(t), S_B(t) \rangle = N^2(t) \langle S_0, S_0 \rangle = \left(\hat{a}^2 - \hat{b}^2 - \frac{1}{4} (\hat{w}^t \hat{w} - \hat{r}^t \hat{r}) \right) \langle S_0, S_0 \rangle$$

and is not necessarily constant. So we may impose the additional condition

:

$$\frac{d}{dt} \left(\widehat{a}^2 - \widehat{b}^2 - \frac{1}{4} (\widehat{w}^t \widehat{w} - \widehat{r}^t \widehat{r}) \right) = 0 \Leftrightarrow \widehat{a} \frac{d\widehat{a}}{dt} - \widehat{b} \frac{d\widehat{b}}{dt} - \frac{1}{4} \left(\widehat{w}^t \frac{d\widehat{w}}{dt} - \widehat{r}^t \frac{d\widehat{r}}{dt} \right) = 0$$

See below continuity equation.

In this aggregation the section σ represents the individual motion of the constituting material points, with respect to a gauge attached to the solid. The element $\sigma_B(t)$ represents the average motion of these points with respect to the gauge of the observer B in the computation of σ_B . The motion of the solid itself, with respect to the gauge of an observer O (in a different, more general model), is represented by an element $\sigma_o \in P_G$. The total motion (solid + solid on itself) is defined by a change of gauge in P_G and the resulting spinor (as it would be used in a model representing the solid) is then :

$$S(t) = N(t) \gamma C(\sigma_o(t)) \gamma C(\sigma_B(t)) S_0$$

Which sums up to replace the fixed inertial spinor S_0 by the variable spinor $S_B(t) = N(t) \gamma C(\sigma_B(t)) S_0$.

The physical meaning of σ_o must be understood with respect to the way the solid is defined : for instance if $\sigma(t, x)$ represents a rotation around an axis, then σ_o will be a rotation of this axis. The vector $r \in \mathbb{R}^3$ in σ_r , which has no geometric meaning for a particle, gets one for a solid, similar to the usual.

Continuity equation

The conservation of the mass of the body means, for the observer B, that :

$$\mathcal{M}(t) = \int_{\omega(t)} \mu_3(t, x) \varpi_3 = Ct = \int_{\omega(t)} i_V(\mu \varpi_4)$$

Consider the manifold $\omega([t_1, t_2])$ with borders $\omega(t_1), \omega(t_2)$:

$$\mathcal{M}(t_2) - \mathcal{M}(t_1) = \int_{\partial\omega([t_1, t_2])} i_V(\mu \varpi_4) = \int_{\omega([t_1, t_2])} d(i_V \mu \varpi_4)$$

$$d(i_V \mu \varpi_4) = \mathcal{L}_V(\mu \varpi_4) - i_V d(\mu \varpi_4) = \mathcal{L}_V(\mu \varpi_4) - i_V(d\mu \wedge \varpi_4) - i_V \mu d\varpi_4 = \mathcal{L}_V(\mu \varpi_4)$$

$$\mathcal{E}(t_2) - \mathcal{E}(t_1) = \int_{\omega([t_1, t_2])} \mathcal{L}_V(\mu \varpi_4)$$

with the Lie derivative \mathcal{L} (Maths.1517,1587)

The conservation of the mass is equivalent to the condition $\mathcal{L}_V(\mu \varpi_4) = 0$.

$$\begin{aligned} & \mathcal{L}_V \mu \varpi_4 \\ &= \frac{d\mu}{dt} \varpi_4 + \mu \mathcal{L}_V \varpi_4 \\ &= \frac{d\mu}{dt} \varpi_4 + \mu (\operatorname{div} V) \varpi_4 \\ &= \frac{d\mu}{dt} + \mu (\operatorname{div} V) \varpi_4 \end{aligned}$$

and we retrieve the usual continuity equation :

$$\frac{d\mu}{dt} + \mu \operatorname{div} V = 0 \tag{4.54}$$

$$\text{Then } N(t) = \int_{\omega(t)} \mu_3(t, x) \varpi_3$$

If $\mu = Ct$ (incompressible solid) the condition becomes : $\operatorname{div} V = 0$

Symmetries of a solid

By symmetries we mean symmetries of the whole body B : under a geometric transformation the body looks the same for an observer. So they are transformations occurring in each $\omega(t)$ and for a privileged observer who can see the

whole body. It is equivalent to consider either the transformation of the body or the transformation of the observer (as long as he keeps the same vector field ε_0), thus symmetries can be represented as a global change of observer with an element $s(t)$ belonging to a subgroup of $Spin(3)$. And B has a symmetry if the section $S(t, x) = \gamma C(\sigma(t, x)) S_0 \in \mathfrak{X}(S_0)$ is such that $\sigma(t, x) = \sigma_w(t, x) \cdot s(t)$. Then :

$$\begin{aligned} \sigma(t, x) &= \sigma_w(t, x) \cdot s(t) \\ &= (a_w(t, x) + v(0, w(t, x))) \cdot (a_r(t) + v(r(t), 0)) \\ &= a_w a_r + v(a_w r, \frac{1}{2} j(w) r + a_r w) - \frac{1}{4} (w^t r) \varepsilon_5 \\ &\text{and } \sigma_B(t) = N(t) \sigma_{Bw}(t) \cdot s(t) \\ &\text{with } \sigma_{Bw}(t) = \frac{1}{N(t)} \int_{\omega(t)} \sigma_w(t, x) \mu_3(t, x) \varpi_3(t, x) \text{ and the condition } \widehat{a}^2 - \\ &\widehat{b}^2 = N^2 + \frac{1}{4} (\widehat{w}^t \widehat{w} - \widehat{r}^t \widehat{r}). \end{aligned}$$

This approach can be useful in Astrophysics, where trajectories of stars systems or galaxies are studied. The spinor S_B can account for the rotational momentum of the bodies, which is significant and contributes to the total kinetic energy of the system.

4.4.5 Relativist Momentum, Spin and Energy

To sum up :

The Spinor S is the relativist momentum of the particle. Encompassing both the translational and the rotational motions, it is intrinsic and does not depend on the choice of a basis, and its measure $S = \gamma C(\sigma) S_0$ depends on the observer. This is the equivalent of the 4 vector $P = \mu u$, but it adds the rotational component.

The quantity $M_p = \frac{1}{c^2} \sqrt{|\langle S_0, S_0 \rangle_E|} = \frac{1}{c^2} \sqrt{|\langle S, S \rangle_E|}$ is the mass at rest of the particle. It does not depend on the motion, the gauge or the chart.

The decomposition of the relativist momentum by an observer shows :

- the relativist momentum $\epsilon = \pm 1$, spin up or down, with respect to the trajectory

- the translational momentum $\gamma C(\sigma_w) S_0$ with $\sigma_w = a_w + v(0, w)$. With respect to a basis the vector w can be identified with a spatial speed \vec{v} by the formulas 51. And the usual translational momentum is : $\vec{p}_r = M_p \frac{\vec{v}}{\sqrt{1 - \frac{\|\vec{v}\|^2}{c^2}}}$

- the spin $\gamma C(\sigma_r) S_0$ with $\sigma_r = a_r + v(r, 0)$. In a basis the vector r can be assimilated with a rotation of axis $\vec{r} = \sum_{i=1}^3 r_i \varepsilon_i(m)$ and rotational speed $\sqrt{\langle \vec{r}, \vec{r} \rangle_3}$. But this assimilation is formal. A particle, by definition, has no internal structure, so it should look the same after any rotation around its position. The Spin is not a geometric rotation, and because of that in the measure of the Spin the choice of the axis x, y, z does not matter, even if the measures change as if it was a rotation. The value of the spin is expressed in units, the SI unit of spin is the joule-second, as the classical angular momentum.

In practice, however, for elementary particles it is written as a multiple of the reduced Planck constant \hbar .

The **kinematic energy of a particle** can be defined as $\frac{1}{i} \langle S, \frac{dS}{dt} \rangle$ (the $\frac{1}{i}$ factor accounts for the fact that $\langle S, \frac{dS}{dt} \rangle$ is imaginary because $\frac{d}{dt} \langle S, S \rangle = 0$). Its measure depends on the observer, through the choice of t . The intrinsic quantity would be $\frac{1}{i} \langle S, \frac{dS}{d\tau} \rangle$ with the proper time τ , and $\frac{1}{i} \langle S, \frac{dS}{dt} \rangle = \left(\sqrt{1 - \frac{\|\vec{v}\|^2}{c^2}} \right) \frac{1}{i} \langle S, \frac{dS}{d\tau} \rangle$. It can be expressed with (r, w) and their derivatives³, as we will see in the following, and the result is :

$$\frac{1}{i} \langle S, \frac{dS}{dt} \rangle = k^t X$$

where

X is the vector : $X = [A_r(r)] \frac{dr}{d\tau} + [A_w(r, w)] \frac{dw}{d\tau}$, with $[A_r(r)]$, $[A_w(r, w)]$ matrices depending on r, w ,

and k is a vector, similar to the inertia tensor, which encompasses both the translational and the rotational motions, and acts also as a magnetic moment.

Its value is :

$$k = -\epsilon \frac{1}{4} M_p^2 c^4 \begin{bmatrix} (\sin 2\alpha_0) \cos(\alpha_2 - \alpha_1) \\ (\sin 2\alpha_0) \sin(\alpha_2 - \alpha_1) \\ 2 \cos 2\alpha_0 \end{bmatrix}$$

with

$$S_R = \frac{M_p c^2}{\sqrt{2}} \begin{bmatrix} e^{i\alpha_1} \cos \alpha_0 \\ e^{i\alpha_2} \sin \alpha_0 \end{bmatrix} \text{ and } \epsilon = -1 \text{ for antiparticles.}$$

The representation of the kinematic characteristics of a deformable solid has been seen before. For an object which has an internal structure this is done by tensorial products of spinors.

³We can expect that the kinematic energy has the dimension $[\text{Mass}][\text{Speed}]^2$. And we must remember that r, w have themselves the dimension of speed. So the definition is consistent.

Chapter 5

FORCE FIELDS

The concept of fields has appeared in the XIX^o century, in the wake of the electromagnetism theory, to replace the picture of action at a distance between particles. In the following by force field we mean one of the forces which interact with particles : the strong interaction, the weak and the electromagnetic forces combined in an electroweak interaction, gravitation being in one league by itself.

A force field is one object of Physics, which has distinctive properties :

i) Because particles are localized, a field must be able to act anywhere, that is to be present everywhere. So the first feature of force fields, as opposed to particles, is that, a priori, they are defined all over the universe, even if their action can decrease quickly with the distance.

ii) A force field propagates : the value of the field depends on the location, this propagation occurs when there is no particle, thus it is assumed that it results from the interaction of the force fields with themselves.

iii) Force fields interact with particles, which are themselves seen as the source of the fields. This interaction depend on charges which are carried by the particles.

iv) The interactions, of the fields with themselves or with particles are, in continuous processes, represented in the lagrangian according to the Principle of Least Action.

v) In some cases the force fields can act in discontinuous processes, in which they can be represented as particles (bosons and gravitons).

Thus we need a representation of the charges and of the fields. We will start with a short presentation of the Standard Model, as this is the most comprehensive picture of the force fields.

5.1 THE STANDARD MODEL

In the Standard Model there are 4 force fields which interact with particles (the gravitational field is not included) :

- the electromagnetic field (EM)
- the weak interactions
- the strong interactions
- the Higgs field

and two classes of elementary particles, fermions and bosons¹, in distinct families.

They are the main topic of the Quantum Theory of Fields (QTF) and their representation is summarized in the Standard Model.

5.1.1 Fermions and bosons

Fermions

The matter particles, that we will call fermions, are organized in 3 generations, comprised each of 2 leptons and 2 quarks :

- First generation : quarks up and down; leptons : electron, neutrino.
- Second generation : quarks charm and strange; leptons : muon, muon neutrino
- Third generation : quarks top and bottom; leptons : tau and tau neutrino

Their stability decreases with each generation, the first generation constitutes the usual matter. Each type of particle is called a flavor.

Fermions interact with the force fields, according to their charge, which are :

- color (strong interactions) : each type of quark can have one of 3 different colors (blue, green, red) and they are the only fermions which interact with the strong field
- hypercharge (electroweak interaction) : all fermions have an hypercharge (-2,-1,0,1,2) and interact with the weak field
- electric charge (electromagnetic interactions) : except the neutrinos all fermions have an electric charge and interact with the electromagnetic field.

All fermions have a weak isospin T_3 , equal to $\pm 1/2$ and there is a relation between the isospin, the electric charge Q and the hypercharge Y :

$$Y = 2(Q - T_3)$$

The total sum of weak isospin is conserved in interactions.

Each fermion (as it seems also true for the neutrinos) has a mass and so interacts also with the gravitational field. These kinematic properties are rep-

¹Actually the words fermions and bosons are also used for particles, which are not necessarily elementary, that follow the Fermi or the Bose rules in statistics related to many interacting particles. Here we are concerned only with elementary particles. So fermions mean elementary fermions and bosons elementary bosons.

resented in the Standard Model by a spinor with 4 components², and in weak and strong interactions the left and right components interact differently with the fields (this is the chirality effect noticed previously).

Each fermion has an associated antiparticle, which is represented by conjugation of the particle. In the process the charge changes (color becomes anticolor which are different, hypercharge takes the opposite sign), left handed spinors are exchanged with right handed spinors, but the mass is the same.

Elementary particles can be combined together to give other particles, which have mass, spin, charge,... and behave as a single particle. Quarks cannot be observed individually and group together to form a meson (a quark and anti-quark) or a baryon (3 quarks) : a proton is composed of 3 quarks udd and a neutron of 3 quarks uud . A particle can also transformed itself into another one, it can also disintegrate in other particles, and conversely particles can be created in discontinuous process, notably through collisions. The weak interaction is the only field which can change the flavor in a spontaneous, discontinuous, process, and is responsible for natural radioactivity.

Bosons

Besides the fermions, the Standard Model involves other objects, called bosons, linked to the force fields, which share some of the characteristics of particles. They are :

- 8 gluons linked with the strong interactions : they have no electric charge but each of them carries a color and an anticolor, and are massless. They are their own antiparticles.
- 3 bosons W^j linked with the electroweak field, which carry weak hypercharge and have a mass.
- 1 boson B, specific to the electromagnetic field, which carries a hypercharge and a mass.
- 1 Higgs boson, which has two bonded components, is its own antiparticle and has a mass but no charge or color

The bosons W, B combine to give the photon, the neutral boson Z and the charged bosons W^\pm . The photon and Z are their antiparticle, W^\pm are the antiparticle of each other. So in the Standard Model photons are not elementary particles (at least when electroweak interactions are considered).

5.1.2 The group representation

To put some order in the zoo of the many particles which were discovered a natural starting point is QM : since states of particles can be represented in Hilbert space, it seems logical to assign to each (truly) elementary fermion a vector of a basis of this Hilbert space F. Then the combinations which appear are represented by vectors ϕ , which are linear combinations (or in some case tensorial

²Because the right and left part are related, usually only one of them is used in computations, and we have two components Weyl spinors.

products) of these basis vectors, and the process of creation / annihilation are transitions between given states, following probability laws. The fact that there are three distinct generations of fermions, which interact together and appear in distinctive patterns, leads to the idea that they correspond to different representations of a group U . Indeed the representations of compact groups can be decomposed in sum of finite dimensional irreducible representations, thus one can have in the same way one group and several distinct but related Hilbert spaces. The problem was then to identify both the group U , and its representations. A given group has not always a representation of a given dimension, and representations can be combined together. Experiments lead to the choice of the direct product $SU(3) \times SU(2) \times U(1)$ as the group, and to precise the representations (whose definition is technical and complicated). Actually the range and the strength of the force fields are different : the range is very short for the strong and weak interactions, infinite for the electromagnetic field, moreover all fermions interact with the weak force and, except for the neutrinos, with the electromagnetic field. So this leads to associate more specifically a group to each force field :

- $SU(3)$ for the strong force
 - $U(1) \times SU(2)$ for the electroweak force (when the weak force is involved, the electromagnetic field is necessarily involved)
 - $U(1)$ for the electromagnetic force
- and to consider three layers : $U(1)$, $U(1) \times SU(2)$, $U(1) \times SU(2) \times SU(3)$ according to the forces that are involved in a problem.

On the other hand it was necessary to find a representation of the force fields, if possible which fits with the representation of the fermions. The first satisfying expression of the Maxwell's laws is relativist and leads to the introduction of the potential \hat{A} , which is a 1-form, and of the strength of the field \mathcal{F} , which is a two-form, to replace the electric and magnetic fields. It was soon shown that the Maxwell's equations can be expressed elegantly in the fiber bundle formalism, with the group $U(1)$. In the attempt to give a covariant (in the SR context) expression of the Schrodinger's equation including the electromagnetic field it was seen that this formalism was necessary. Later Yang and Mills introduced the same formalism for the weak interactions, which was extended to the strong interactions, and it became commonly accepted in what is called the gauge theories. The key object in this representation is a connection, coming from a potential, acting on a vector bundle, where ϕ lives, which corresponds to the representation of the group U .

5.1.3 The Standard Model

The Standard Model is a version of the Yang-Mills model, adapted to the Special Relativity geometry :

- i) Each of the groups or product of groups defines a principal bundle over the Minkovski affine space (which is \mathbb{R}^4 with the Lorentz metric).

ii) The physical characteristics (the charges) of the particles are vectors ϕ of a vector bundle associated to a principal bundle modelled on U .

iii) The state of the particles is then represented in a tensorial bundle, combining the spinor S (for the kinematic characteristics) and the physical characteristics ϕ .

iv) The masses are defined separately, because it is necessary to distinguish the proper mass and an apparent mass resulting from the screening by virtual particles.

v) Linear combination of these fermions give resonances which have usually a very short life. Stable elementary particles (such as the proton and the neutron) are bound states of elementary particles, represented as tensorial combinations of these fermions.

vi) The fields are represented by principal connections, which act on the vector bundles through ϕ . The Higgs field is represented through a complex valued function. The electroweak field acts differently on the chiral parts of fermions.

vii) The lagrangian is built from scalar products and the Dirac's operator.

viii) The bosons correspond to vectors of bases of the Lie algebras of each of the groups : the 8 gluons to $su(3)$, the 3 bosons W^j to $su(2)$, 1 boson B to $u(1)$.

5.1.4 The issues

The Standard Model does not sum up all of QTF, which encompasses many other aspects of the interactions between fields and particles. However there are several open issues in the Standard Model.

1. The Standard Model, built in the Special Relativity geometry, ignores gravitation. Considering the discrepancy between the forces at play, this is not really a problem for a model dedicated to the study of elementary particles. QTF is rooted in the Poincaré's algebra, and the localized state vectors, so it has no tool to handle trajectories, which are a key component of differential geometry.

2. The Higgs boson, celebrated recently, raises almost as many questions as it gives answers. It has been introduced in what can be considered as a patch, needed to solve the issue of masses for fermions and bosons. The Dirac's operator, as it is used for the fermions, does not give a definite positive scalar product and is null (and so their mass) whenever the particles are chiral. And as for the bosons, the equivariance in a change of gauge forbids the explicit introduction of the potential, which is assumed to be their correct representation, in the lagrangian. The Higgs boson solves these problems, but at the cost of many additional parameters, and the introduction of a fifth force which it should carry.

3. From a semi-classic lagrangian, actually most of the practical implementation of the Standard Model relies on particles to particles interactions, detailed by Feynmann's diagram and computed through perturbative methods. In many ways the actual QTF is, under a sophisticated guise, a return to the action at

a distance, the bosons and other virtual particles making the link. Of course this is consistent with a discrete representation of the world, but also difficult to conciliate with the matter fields representing particles.

4. The range of the weak and strong interactions is not well understood. Formally it is represented by the introduction of a Yukawa potential (which appears as a “constant coupling” in the Standard Model), proportional to $\frac{1}{r} \exp(-km)$ which implies that if the mass m of the carrier boson is not null the range decreases quickly with the distance r . Practically, as far as the system which is studied is limited to few particles, this is not a big issue.

5. We could wish to incorporate the three groups in a single one, meanwhile encompassing the gravitational field and explaining the hierarchy between the forces. This is the main topic of the Great Unification Theories (GUT) (see Seibatu for a review of the subject). The latest, undergone by Garrett Lisi, invokes the exceptional Lie group E_8 . Its sheer size (it has 248 dimensions) enables to account for everything, but also requires the introduction of as many parameters.

An option, which has been studied by Trayling and Lisi, would be to start, not from Lie groups, but from Clifford algebras as we have done for the Spinors. The real dimension of $SU(3) \times SU(2) \times U(1)$ is $12 = 8 + 3 + 1$ which implies to involve at least a Clifford algebra (dimension 2^n) on a four dimensional vector space and it makes sense to look at its complexified. The groups would then be Spin subgroups of the Clifford algebra. We have the following isomorphisms :

$$U(1) \sim Spin(\mathbb{R}, 2)$$

$$SU(2) \sim Spin(\mathbb{R}, 3)$$

but there is no simple isomorphism for $SU(3)$.

Albeit all together they are part of $Cl(\mathbb{R}, 10)$.

In the next sections we will see how the states of particles, force fields, including gravitation, and their interactions can be represented, in the geometrical context of GR. In the next chapter we will review the requirements that these representations impose to Lagrangians and continuous models. Two kinds of continuous models, simplified but similar to the Standard Model, will then be studied. They do not pretend to replace the Standard Model, but to help to understand the mechanisms at play, notably the motivation to use the mathematical tools in the representation of physical phenomena. So we will not insist on the many technical details of the Standard Model, heavily loaded with historical notations, and keep the formalism to a minimum.

5.2 STATES OF PARTICLES

We have seen that spinor fields can be characterized, beyond the inertial spinor, by an integer, which defines families of particles with similar behavior. Particles can then be differentiated, in addition to their kinematic characteristics summarized in the spinor, by a charge which accounts for their interaction with force fields. A particle can be seen as a system in itself. Its state is then a combination of its kinematic characteristics, represented by the spinor, and of its charge, which represents its interaction with the force fields. Using the description of elementary particles given by the Standard Model, it is then possible to set up a representation of elementary particles. From there the representation can be extended to particles and matter fields.

5.2.1 The space of representation of the states

The Law of Equivalence

We can follow some guidelines :

i) For any particle there are intrinsic characteristics ψ_0 , which do not change with the fields or the motion. If we assume that ψ belongs to a vector space V , then there is a set of vectors $\{\psi_{0p}\}_{p=1}^N$ such that ψ_{0p} characterizes a family of particles which have the same behavior.

ii) Motion is one of the features of the state of particles. It is represented by the action of $Spin(3, 1)$ on the space V , as we have done in the previous chapter.

iii) The intrinsic kinematic characteristics of particles are represented in the vector spaces E_0, E'_0 : each family of particles is associated to one vector of these spaces. Particles and anti-particles are distinguished by their inertial spinor.

iv) In the Newton's law of gravitation $F = G \frac{MM'}{r^2}$ and his law of Mechanics : $F = \mu\gamma$ the scalars M, μ represent respectively the gravitational charge and the inertial mass, and there is no reason why they should be equal. However this fact has been verified with great accuracy (two bodies fall in the vacuum at the same speed). This has lead Einstein to state the fundamental Law of Equivalence "Gravitational charge and inertial mass are identical". From which he built the Theory of General Relativity. This Law leads us to take as gravitational charge of particles the inertial spinor S_0

Proposition 82 *The Gravitational charge of a particle is represented by its inertial spinor S_0*

So, if we stay only with the gravitational field, the space E and the representation $(E, \gamma C)$ suffice to represent the state of particles. The kinematic characteristics of particles of the same flavor (quarks, leptons) are not differentiated according to their other charges. So we have $\psi_{0p} = S_{0p}$.

In the previous chapter we assumed that :

- there is, along the world line of a particle, a privileged frame $\varphi_G(m, \sigma(m))$ such that the spinor of the particle is $(\varphi_G(m, \sigma(m)), S_0)$ with $S_0 = Ct$

- the observer measures the spinor $S(m)$ in his gauge : $\varphi_G(m,1)$ and $(\varphi_G(m,\sigma(m)), S_0) \sim (\varphi_G(m,1), \gamma C(\sigma(m)) S_0) = (\varphi_G(m,1), S(m))$ thus : $S(m) = \gamma C(\sigma(m)) S_0$

We have now to consider an interpretation which is mathematically equivalent, but physically different :

- the observer measures the spinor $S(m)$ with $(\varphi_G(m,1), \gamma C(\sigma(m)) S_0)$
- in presence of gravity this spinor is equivalent to : $(\varphi_G(m,\sigma(m)), S_0)$

The privileged gauge (for the particle) is provided by the gravitational field. And the action of the motion, that is of the inertial forces, is equivalent to the action of gravity on the state of the particle, which is the meaning of the Law of Equivalence. This is a key point to understand the gravitational and the other fields : *particles have intrinsic properties, that they keep all over their travel on their world lines but, because of the existence of the field, their measure by an observer is distinct from this intrinsic value.* This leads to see the fields as the value of the element of the group ($\sigma \in Spin(3,1)$ for the gravitational field) but, as we will see, the action of the field goes through a special derivative because it manifests itself in the motion of the particle.

Representation of the charges for the other fields

For the other fields :

i) Bosons give the structure of the fields, in accordance with the dimension of the groups : 8 for the strong force ($SU(3)$ dimension 8), 3 for the weak force ($SU(2)$ dimension 3), 1 for the electromagnetic force ($U(1)$ dimension 1). In QTF the action of fields is represented by operators acting on V , in the representation of the Lie algebra of the groups. Because the exponential is surjective on compact groups it sums up to associate the fields to an action of the groups on V .

ii) The action depends on the charges - accounting for the possible combinations of charges, there all together 24 kinds of fermions - but also on the inertial spinors : particles and antiparticles do not behave the same way, and weak forces act differently according to the left or right chiral parts

Assuming that V is a vector space, and the actions of the fields are linear, the solution is to take V as the tensorial product $V = E \otimes F$ where F is a vector space such that (F, ϱ) is a representation of the group U corresponding to the forces other than gravity ($U = SU(3) \times SU(2) \times U(1)$ in the Standard Model).

That we sum up by :

Proposition 83 *There is a compact, connected, real Lie group U which characterizes the force fields other than gravitation.*

There is a n dimensional complex vector space F , endowed with a scalar product denoted $\langle \rangle_F$ and (F, ϱ) is a unitary representation of U

The states of particles are vectors of the tensorial product $E \otimes F$

*The intrinsic characteristics of each type of particles are represented by a tensor $\psi_0 \in E \otimes F$, that we call a **fundamental state**, and all particles sharing the same characteristics behave identically under the actions of all the fields.*

Notation 84 $(f_i)_{i=1}^n$ is a basis of F . We will assume that it is orthonormal.

$(\vec{\theta}_a)_{a=1}^m$ is a basis of the Lie algebra T_1U

$[\theta_a]$ is the matrix of $\rho'(1)$ $(\vec{\theta}_a)$ expressed in the basis $(f_i)_{i=1}^n$.

As a consequence :

i) Because (F, ρ) is a unitary representation, the scalar product is preserved by ρ : $\langle \rho(g)\phi, \rho(g)\phi' \rangle_F = \langle \phi, \phi' \rangle_F$

ii) $(F, \rho'(1))$ is a representation of the Lie algebra T_1U

iii) The derivative $\rho'(1)$ is anti-unitary and the matrices $[\rho'(1)\vec{\theta}_a] = [\theta_a]$ are anti-hermitian :

$$[\theta_a] = -[\theta_a]^* \quad (5.1)$$

F must be a complex vector space to account for the electromagnetic field. F is actually organized as different representations of the group U , and the representation is not irreducible, to account for the generations effect. Composite particles (such as the proton or the neutron) are represented by tensorial product of vectors of $E \otimes F$.

A basis of $E \otimes F$ is $(e_i \otimes f_j)_{i=0..3}^{j=1..n}$

The state of a particle is expressed as a tensor :

$\psi = \sum_{i=1}^4 \sum_{j=1}^n \psi^{ij} e_i \otimes f_j$ that we will usually denote in the matrix form : $[\psi]$ with 4 lines and n rows.

which reads :

$$\psi = \sum_{j=1}^n \left(\sum_{i=1}^4 \psi^{ij} e_i \right) \otimes f_j = \sum_{j=1}^n S^j \otimes f_j \text{ where } S^j \in E$$

So, when gravity alone is involved, the particles such as $\sum_{i=1}^4 \psi_0^{ij} e_i = S_0^j$ have the same behavior and can be seen as n particles, differentiated by their inertial spinor, and thus by their mass. At an elementary level the different values of the inertial spinors characterize the kinematics of each elementary particle.

The experimental fact that the action of the force fields depends also of the spinor part implies that the tensor is not necessarily decomposable (it cannot be written as the tensorial product of two vectors). However we will see that one can attribute a charge to a particle, but it is not expressed as a scalar quantity. There is no natural unit for the charges (except, for historical reasons, for the electric charge), and, indeed, what could be the unit for the colors of the strong force ? The set \mathfrak{F} of existing vectors ψ_0 is just an organized map of all the known combinations of spinors and charges. The formalism with the group representation is built on the experimental facts, but it does not answer the question : why is it so ?

The direct product group $Spin(3,1) \times U$ has an action denoted ϑ on $E \otimes F$
 $\vartheta : Spin(3,1) \times U \rightarrow \mathcal{L}(E \otimes F; E \otimes F)$
 defined by linear extension of γC and ρ :

$\vartheta(\sigma, \varkappa)(\psi) = \sum_{i,k=1}^4 \sum_{j,l=1}^n [\gamma C(\sigma)]_k^i [\varrho(\varkappa)]_l^j \psi^{kl} e_i \otimes f_j$
that we will denote in matrices :

Notation 85

$$\vartheta : Spin(3, 1) \times U \rightarrow \mathcal{L}(E \otimes F; E \otimes F) :: \vartheta(\sigma, \varkappa)[\psi] = [\gamma C(\sigma)][\psi][\varrho(\varkappa)] \quad (5.2)$$

One can extend the action of the Spin group to the action of the Clifford algebra. We define the action ϑ of $Cl(\mathbb{R}, 3, 1) \times U$ on $E \otimes F$ by the unique linear extension of :

$$\begin{aligned} \vartheta : Cl(\mathbb{R}, 3, 1) \times U &\rightarrow \mathcal{L}(E \otimes F; E \otimes F) :: \\ \vartheta(s, g)(S \otimes \phi) &= \gamma C(s)(S) \otimes \varrho(g)(\phi) \end{aligned}$$

to all tensors on $E \otimes F$

This is a morphism from $Cl(\mathbb{R}, 3, 1)$ on $L(E \otimes F; E \otimes F)$: ϑ is linear and preserves the Clifford product

Proof. $\vartheta(s \cdot s', gg')(S \otimes \phi) = \gamma C(s \cdot s')(S) \otimes \varrho(gg')(\phi)$
 $= \gamma C(s) \circ \gamma C(s')(S) \otimes \chi(g) \circ \varrho(g')(\phi)$
 $= \gamma C(s)(\gamma C(s')(S)) \otimes \varrho(g)(\varrho(g')(\phi))$
 $= (\gamma C(s) \otimes \varrho(g))(\gamma C(s')(S) \otimes \varrho(g')(\phi))$
 $= \gamma C(s) \otimes \varrho(g)(\gamma C(s') \otimes \varrho(g')(S \otimes \phi))$
 $= (\gamma C(s) \otimes \varrho(g)) \circ (\gamma C(s') \otimes \varrho(g'))(S \otimes \phi)$
 $= (\vartheta(s, g) \circ \vartheta(s', g'))(S \otimes \phi)$
 $\vartheta(1, 1)(S \otimes \phi) = \gamma C(1)(S) \otimes \varrho(1)(\phi) = S \otimes \phi \quad \blacksquare$
 $\vartheta(\sigma, 1)\psi = \gamma C(\sigma)\psi = \sum_{jkl} [\gamma C(\sigma)]_k^j \psi^{kl} e_j \otimes f_l$
 So the map ϑ defines a representation of $Cl(\mathbb{R}, 3, 1) \times U$ on $E \otimes F$.

Scalar product on the space $E \otimes F$

The scalar product on $E \otimes F$ is necessarily defined as :

$$\langle \psi, \psi' \rangle = \sum_{ijq} [\gamma_0]_k^i \delta_{jq} \bar{\psi}^{ij} \psi'^{kq} = \sum_{ijk} [\gamma_0]_k^i \bar{\psi}^{ij} \psi'^{kj} = Tr([\psi]^* [\gamma_0] [\psi'])$$

because the basis $(f_j)_{j=1}^n$ is orthonormal.

$$\langle \psi, \psi' \rangle = Tr([\psi]^* [\gamma_0] [\psi']) \quad (5.3)$$

Theorem 86 *The scalar product on $E \otimes F$ is preserved by ϑ :*

$$\langle \vartheta(\sigma, \varkappa)\psi, \vartheta(\sigma, \varkappa)\psi' \rangle = \langle \psi, \psi' \rangle$$

Proof. $\tilde{\psi}^{ij} = \sum_{k=1}^4 \sum_{l=1}^n [\gamma C(\sigma)]_k^i [\varrho(\varkappa)]_l^j \psi^{kl}$
 $\langle \tilde{\psi}, \tilde{\psi}' \rangle = \sum [\gamma_0]_k^i \overline{[\gamma C(\sigma)]_p^i} \overline{[\varrho(\varkappa)]_q^j} \tilde{\psi}^{pq} [\gamma C(\sigma)]_r^k [\varrho(\varkappa)]_s^j \psi'^{rs}$
 $= \sum ([\gamma C(\sigma)]^* [\gamma_0] [\gamma C(\sigma)])_r^p ([\varrho(\varkappa)]^* [\varrho(\varkappa)])_s^q \tilde{\psi}^{pq} \psi'^{rs}$
 $= \sum [\gamma_0]_r^p \tilde{\psi}^{pq} \psi'^{rs} \quad \blacksquare$

The scalar product is not definite, positive or negative, on $E \otimes F$. However there is a norm $\| \cdot \|_E$ on the space E and a norm on the space F , the latter defined by the scalar product. They define a norm on $E \otimes F$ by taking $\|e_i \otimes f_j\| =$

$\|e_i\|_E \|f_j\|_F$. Moreover this norm is invariant by ϑ . So that $E \otimes F$ is a Banach vector space.

Physical states of elementary particles

$(E \otimes F, \vartheta)$ is a unitary representation of the Lie group $Spin(3, 1) \otimes U$.

For any $\psi \in E \otimes F$ the set $\{\vartheta(\sigma, \varkappa) \psi, (\sigma, \varkappa) \in Spin(3, 1) \otimes U\}$ is the orbit of ψ . The orbits are the set of states corresponding to the same type of particles.

The relation of equivalence $\psi \sim \psi' \Leftrightarrow \exists (\sigma, \varkappa) \in Spin(3, 1) \otimes U : \psi' = \vartheta(\sigma, \varkappa) \psi$ defines a partition of $E \otimes F$ corresponding to the orbits. And each class of equivalence can be identified with a fundamental state ψ_0 .

All particles of the same type ψ_0 have the same behavior with the same fields \varkappa : so for ψ_0, \varkappa fixed, σ then ψ are fixed uniquely

The measures of fields is done by measuring the motion σ of known particles ψ_0 subjected to fields \varkappa : so from ψ, ψ_0 and σ one can compute a unique value \varkappa of the field.

Which sums up to, if \mathfrak{F} is the set of possible states of elementary particles :

Proposition 87 *The action of $Spin(3, 1) \times U$ on \mathfrak{F} is free and faithful : $\forall \psi \in \mathfrak{F} : \vartheta(\sigma, \varkappa) \psi = \psi \Leftrightarrow (\sigma, \varkappa) = (1, 1)$*

Then $\vartheta(\sigma, \varkappa) \psi = \vartheta(\sigma', \varkappa') \psi \Leftrightarrow (\sigma, \varkappa) = (\sigma', \varkappa')$

We had seen that this is the case for the spinor. This is extended to the states of particles.

The orbits are not vector subspaces :

Theorem 88 *For any fundamental state ψ_0 , the orbit $(E \otimes F)(\psi_0)$ of ψ_0 is a real finite dimensional Riemannian manifold, embedded in $E \otimes F$*

Proof. $Spin(3, 1)$ and U are real Lie groups, thus manifolds, take a chart in each

The vector spaces tangent at any point to the manifold are subspaces of the vector space $E \otimes F$

The metric on the tangent bundle is given by the scalar product, which is definite, positive or negative. ■

Particles and antiparticles

Chirality was introduced in the choice of the representation $(E, \gamma C)$ because it is significant in the behavior with the force fields. We will similarly distinguish in the matrix $[\psi]$ a right part, with the first 2 rows, and a left part, with the last 2 rows, so that in matrix form $[\psi] = \begin{bmatrix} \psi_R \\ \psi_L \end{bmatrix}$. In QTF this is called a Dirac's spinor, and ψ_R, ψ_L are Weyl's spinors.

The difference between particles and antiparticles was based on the sign of the scalar product $\langle S_0, S_0 \rangle$. So it is legitimate to discriminate particles and

antiparticles in a similar way. More precisely, we look for the subsets of $E \otimes F$ such that :

- i) the scalar product is definite either positive or negative : $\langle \psi_0, \psi_0 \rangle = 0 \Rightarrow \psi_0 = 0$
- ii) this is still true whenever ψ_0 is the tensorial product $\psi_0 = S_0 \otimes F_0$
- iii) the populations of antiparticles and particles are preserved by space reversal, and exchanged by time reversal, as we know that this is still true for particles in the Standard Model.

Theorem 89 *The only vector subspaces of $E \otimes F$ which meet these conditions are such that $\psi_L = \epsilon i \psi_R$ with $\epsilon = \pm 1$*

Proof. i) $\langle \psi, \psi \rangle = Tr([\psi]^* [\gamma_0] [\psi]) = i Tr(-\psi_R^* \psi_L + \psi_L^* \psi_R)$
 $Tr(\psi_L^* \psi_R) = Tr(\psi_L^* \psi_R)^t = Tr(\psi_R^t \overline{\psi_L}) = \overline{Tr(\psi_R^* \psi_L)}$
 Thus : $Tr(-\psi_R^* \psi_L + \psi_L^* \psi_R) = \overline{Tr(\psi_R^* \psi_L)} - Tr(\psi_R^* \psi_L)$
 $= -2i \text{Im} Tr(\psi_R^* \psi_L) \in i\mathbb{R}$
 and $\langle \psi, \psi \rangle = 2 \text{Im} Tr(\psi_R^* \psi_L) \in \mathbb{R}$

For $\psi = S \otimes F$ the matrix $[\psi]$ reads : $[\psi] = [S] [F]^t = \begin{bmatrix} S_R F^t \\ S_L F^t \end{bmatrix}$

and $\langle \psi, \psi \rangle = 2 \text{Im} Tr(\overline{[F]} [S_R]^* [S_L] [F]^t) = 2 \text{Im} [S_R]^* [S_L] Tr(\overline{[F]} [F]^t)$

It will be non degenerate iff : $S_L = \epsilon i S_R$ as seen previously and so we can generalize to $\psi_L = \epsilon i \psi_R$:

$\langle \psi, \psi \rangle = 2 \text{Im} Tr(\epsilon i \psi_R^* \psi_R) = 2\epsilon Tr(\psi_R^* \psi_R)$

ii) Time reversal is an operator on $E \otimes F$, represented by the matrix (see the section Spinor Model above) :

$T = \begin{bmatrix} 0 & i\sigma_0 \\ i\sigma_0 & 0 \end{bmatrix}$ with signature (3,1)

$T \begin{bmatrix} \psi_R \\ \epsilon i \psi_R \end{bmatrix} = \begin{bmatrix} 0 & i\sigma_0 \\ i\sigma_0 & 0 \end{bmatrix} \begin{bmatrix} \psi_R \\ \epsilon i \psi_R \end{bmatrix} = \begin{bmatrix} -\epsilon \psi_R \\ i \psi_R \end{bmatrix} = \begin{bmatrix} -\epsilon \psi_R \\ -\epsilon i (-\epsilon \psi_R) \end{bmatrix}$

$T = \begin{bmatrix} 0 & \sigma_0 \\ \sigma_0 & 0 \end{bmatrix}$ with signature (1,3)

$T \begin{bmatrix} \psi_R \\ \epsilon i \psi_R \end{bmatrix} = \begin{bmatrix} 0 & \sigma_0 \\ \sigma_0 & 0 \end{bmatrix} \begin{bmatrix} \psi_R \\ \epsilon i \psi_R \end{bmatrix} = \begin{bmatrix} i \epsilon \psi_R \\ \psi_R \end{bmatrix} = \begin{bmatrix} i \epsilon \psi_R \\ -\epsilon i (i \epsilon \psi_R) \end{bmatrix}$

iii) Space reversal is an operator on $E \otimes F$, represented by the matrix :

$S = i\gamma_0 = \begin{bmatrix} 0 & \sigma_0 \\ -\sigma_0 & 0 \end{bmatrix}$ with signature (3,1)

$S \begin{bmatrix} \psi_R \\ \epsilon i \psi_R \end{bmatrix} = \begin{bmatrix} 0 & \sigma_0 \\ -\sigma_0 & 0 \end{bmatrix} \begin{bmatrix} \psi_R \\ \epsilon i \psi_R \end{bmatrix} = \begin{bmatrix} i \epsilon \psi_R \\ -\psi_R \end{bmatrix} = \begin{bmatrix} i \epsilon \psi_R \\ \epsilon i (i \epsilon \psi_R) \end{bmatrix}$

$S = \begin{bmatrix} 0 & -i\sigma_0 \\ i\sigma_0 & 0 \end{bmatrix}$ with signature (1,3)

$S \begin{bmatrix} \psi_R \\ \epsilon i \psi_R \end{bmatrix} = \begin{bmatrix} 0 & -i\sigma_0 \\ i\sigma_0 & 0 \end{bmatrix} \begin{bmatrix} \psi_R \\ \epsilon i \psi_R \end{bmatrix} = \begin{bmatrix} \epsilon \psi_R \\ i \psi_R \end{bmatrix} = \begin{bmatrix} \epsilon \psi_R \\ \epsilon i (i \epsilon \psi_R) \end{bmatrix} \blacksquare$

And we can state :

Proposition 90 *The intrinsic characteristics ψ_0 of particles (fermions) are such that :*

$\psi_L = i\psi_R$ for particles, their mass M_p is such that

$$\langle \psi_0, \psi_0 \rangle = 2Tr(\psi_R^* \psi_R) = M_p^2 c^4$$

$\psi_L = -i\psi_R$ for antiparticles, their mass is

$$\langle \psi_0, \psi_0 \rangle = -2Tr(\psi_R^* \psi_R) = -M_p^2 c^4$$

To each fermion is associated an antiparticle which has the same mass.

As ϑ preserves the scalar product : $\langle \vartheta(\sigma, \varkappa) \psi_0, \vartheta(\sigma, \varkappa) \psi_0 \rangle = \langle \psi_0, \psi_0 \rangle$ the scalar product is definite positive or negative on the sets :

$(E \otimes F)(\psi_0) = \{\vartheta(\sigma, \varkappa) \psi_0, \sigma \in Spin(3, 1), \varkappa \in U\}$ for a fixed ψ_0 such that $\psi_L = \epsilon i \psi_R$

But these sets are not vector spaces. The expressions are :

$$[\psi] = [\gamma C(\sigma)] [\psi_0] [\varrho(\varkappa)] \text{ and } \sigma = \sigma_w \cdot \sigma_r$$

$$\gamma C(a + v(r, w) + b\varepsilon_5) \begin{bmatrix} \psi_{0R} \\ \epsilon i \psi_{0R} \end{bmatrix} [\varrho(\varkappa)]$$

$$= \begin{bmatrix} ((a+b) + \frac{1}{2} \sum_a (w_a - ir_a) \sigma_a) \psi_{0R} [\varrho(\varkappa)] \\ \epsilon i ((a-b) - \frac{1}{2} \sum_a (w_a + ir_a) \sigma_a) \psi_{0R} [\varrho(\varkappa)] \end{bmatrix}$$

$\gamma C(\sigma_r)$ preserves E_0 , and similarly the chiral relation $\psi_L = i\psi_R$

5.2.2 The Electromagnetic field (EM)

In the Standard Model the Electromagnetic field (EM) is represented by the group $U(1)$, the set of complex numbers with module 1 ($uu^* = 1$). It is a compact abelian group. Its irreducible representations are unidimensional, that is multiple of a given vector.

For any given arbitrary vector f there are 3 possible irreducible non equivalent representations :

- the standard one : $(F, \varrho) : \varrho(e^{i\phi}) f = e^{i\phi} f$ and $F = \{e^{i\phi} f, \phi \in \mathbb{R}\}$

- the contragredient : $(F, \bar{\varrho}) : \bar{\varrho}(e^{i\phi}) f = e^{-i\phi} f$ and $F = \{e^{i\phi} f, \phi \in \mathbb{R}\}$
(Maths.23.1.2)

- the trivial representation : $(F, \varrho) : \varrho(e^{i\phi}) f = f$ and $F = \{f\}$

The standard representation corresponds to negative charge, the contragredient representation to positive charge and the trivial one to neutral charge (the only known are the neutrinos). The choice positive / negative is arbitrary.

The EM field interacts similarly with the left and right part of a spinor, so the space of states of the particles is the sum of tensorial products : $S \otimes f$. The theory can be fully expressed this way. However it is legitimate to choose the vectors f in E , which is a 4 dimensional complex vector space. For elementary particles, then :

i) Let $\{S_p \in E_0, p = 1 \dots N\}$ be N vectors representing inertial spinors of particles. Then for each of them their states are represented by $\{e^{i\phi} S_p, \phi \in \mathbb{R}\}$ with the standard representation;

ii) Let $\{S'_p \in E'_0, p = 1 \dots N\}$ be N vectors representing inertial spinors of the antiparticles associated to S_p . Because E_0, E'_0 are orthogonal these vectors

are orthogonal to the S_p . Then for each of them their states are represented by $\{e^{i\phi}S'_p, \phi \in \mathbb{R}\}$ with the contragredient representation;

iii) Neutral particles $\{S_q \in E_0 \oplus E'_0, q = 1 \dots N'\}$ correspond to the trivial representation : their states is just one vector $S \in E_0$ or E'_0 .

Particles have a charge opposite to their anti-particle. Notice that the maps $\varrho, \bar{\varrho}$ are distinct from $\gamma C(\sigma)$ for any $\sigma \in Spin, T_1Spin$, so the state of a particle can change only by the action of $U(1)$.

A basis of E , for elementary particles, is then $\{(S_p, S'_p)_{p=1}^N, (S_q)_{q=1}^{N'}\}$. Each vector defines the mass, the charge and the type of the particle. In this picture there is no unit for the electric charge.

In the case of the EM field the structure brought by the charges is then built in the space E_0 or E'_0 . The state of an elementary particle is then :

$$\psi = e^{i\phi} S \quad (5.4)$$

where ϕ is a phase factor which must be considered as variable for charged particles, and $\phi = 0$ for neutral particles. As we will see actually the phase can usually be ignored (as the value \varkappa for the other field) : particles whose states differ by a phase factor have the same behavior, with regard to the electromagnetic field and they have the same mass : $\langle (\exp i\phi) S_0, (\exp i\phi') S_0 \rangle = \langle S_0, S_0 \rangle$ and the same EM charge.

This is the origin of the introduction of rays in QM. Two particles such that their states differ by a phase factor $e^{i\phi}$ behave the same way, for the gravitational field or the EM field, so they can be deemed representing the same state.

5.2.3 Composite material bodies

The picture that we have just drawn corresponds to elementary particles : the set $\mathfrak{F} \subset E \otimes F$ of possible states comes from the known combinations of charges and kinematic characteristics of the particles which are deemed elementary. But elementary particles are assumed to be the building blocks of material bodies, and they constitute more or less stable composites which can themselves be considered as particles. Several solutions can be considered to represent these composite bodies.

i) Whenever the weak or strong interactions are involved the state of composite bodies is represented by tensors in $\otimes_n \mathfrak{F}$, according to the theorem 28 of QM. In particular they are the only forces which can change the flavor of a particle.

ii) When only the EM field is involved the states of elementary particles can be represented in E . Mathematically the tensorial product of non equivalent representations is well defined. The action of $U(1)$ on the tensor $S_1 \otimes S_2$ is $e^{i\phi(\epsilon_1 + \epsilon_2)} S_1 \otimes S_2$ with $\epsilon_k = \pm 1$ depending on the representation. However a basis of the tensorial product is comprised of tensorial products of all the vectors of the basis, which would not have the same behavior under the action of $U(1)$ (the basis of E has positive, negative and neutral particles). So actually the only combinations which are acceptable are made of particles of the same kind

(positive, negative or neutral) and the action is then $e^{iq\phi\epsilon}$ where q is the number of particles, where q can be a positive, a negative or a null integer. As a consequence :

- when only the EM field is involved composite particles are comprised of particles with the same type of charge (this does not hold when the weak and strong interactions are considered)

- the electric charge of particles must be an integer multiple of an elementary charge.

Such tensorial products of spinors can be used for nuclei, atoms or molecules. The associated EM charge is an integer multiple of the elementary charge. Of course this does not matter for neutral particles but, as they do not interact with the EM field, we are actually in the next case.

iii) When only the gravitational field is involved the states of particles or material bodies can be represented either by a tensorial product of spinors, or by a single spinor in E .

Material bodies can be represented by a single spinor. There is no scale limit for these solutions, however the representation of a composite body by a single spinor assumes that the body meets the conditions stated in the previous chapter for deformable solids. The interactions between its particles are such that an equilibrium is reached, which preserves its cohesion. And the laws which rule its deformation account for these interactions. As noticed before the aggregation of particles in a deformable solid is done in a separate model, accounting for the specific characteristics of the body. It provides a single spinor which is an approximation of the collection of particles whose state can always be represented by tensors. Such spinors can be considered with the EM field, because they are the combination of an integer multiple of particles with the same charge, the inertial spinor of such composite must account for the total charge, expressed in an integer multiple of the elementary charge. For instance a nucleus can be represented by a single spinor with an EM charge.

Classic Mechanics provides efficient and simpler tools and the use of spinors would be just pedantic in common problems. However there is a domain where spinors could be useful : Astrophysics. It cannot ignore General Relativity, gravity is the main force at work, the common presentation of GR is awfully complicated and does not allow to account (at least in a simple way) for the rotation of stars or star systems, and there are some disturbing facts (nowadays linked to the existence of a dark matter) that are not explained, but about which spinors could provide some answer. In this case the introduction of density besides spinors is the natural path.

In the following we will keep the definition of particles of the previous sections : they are material bodies whose state can be represented by a single vector $\psi \in E \otimes F$, which corresponds either to an elementary particle or a material body whose structure is fully represented in ψ . And we will assume, at least formally, that the state of the particle belong to $E \otimes F$, even if it can be represented by a single spinor in E .

5.2.4 The fiber bundle representation

The action ϑ of the groups gives the value of ψ for any fundamental state ψ_0 :

$$\psi : (E \otimes F) \times (Spin(3,1) \times U) \rightarrow \psi = \vartheta(\sigma, \varkappa) \psi_0$$

The discrepancy between the fundamental state ψ_0 and the measured value ψ is assumed to come from the existence of the fields. So we are lead to assume that there is a principal bundle $Q(M, Spin(3,1) \times U, \pi_U)$ with fiber $Spin(3,1) \times U$, and the fields are represented by a section of this principal bundle : $\mathbf{q}_f \in \mathfrak{X}(Q) :: \mathbf{q}_f(m) = \varphi_Q(m, (\sigma(m), \varkappa(m)))$.

The observer is assumed to use a standard gauge : $\mathbf{q} \in \mathfrak{X}(Q) :: \mathbf{q}(m) = \varphi_Q(m, (1, 1))$.

Then the state of the particle is represented by a vector of the associated vector bundle $Q[E \otimes F, \vartheta]$ with fiber $E \otimes F$.

In the presence of fields the state of the particle is $(\mathbf{q}_f(m), \psi_0) = (\varphi_Q(m, \vartheta(\sigma(m), \varkappa(m))), \psi_0)$ which is equivalent to :

$$(\varphi_Q(m, \vartheta(1, 1)), \vartheta(\sigma(m), \varkappa(m)) \psi_0) = (\mathbf{q}(m), \vartheta(\sigma(m), \varkappa(m)) \psi_0)$$

and the observer measures $\psi(m) = \vartheta(\sigma(m), \varkappa(m)) \psi_0$ in his gauge $\varphi_Q(m, \vartheta(1, 1))$.

The *relativist momentum of the particle is the vector ψ* , belonging to the associated vector bundle $Q[E \otimes F, \vartheta]$. This is a geometric quantity, which is intrinsic to the particle and does not depend on a gauge. It adds the charges to the spinor S.

The measure of the relativist momentum depends on the observer. It represents the momentum of the particle. We will reserve the name State to this measure : *the state is the relativist momentum as measured by an observer*.

Notice that, as a consequence of this representation, the conservation of the characteristics ψ_0 of the particle entails that of its charge and mass during its motion. It is built in the formalism. And, meanwhile spinor and charge are entangled in the tensorial product $E \otimes F$, the gravitational field and the other fields keep their originality : Q has for fiber $Spin(3,1) \times U$ and not $Spin(3,1) \otimes U$.

That we sum up in :

Proposition 91 *There is a principal bundle $Q(M, Spin(3,1) \times U, \pi_U)$ with trivialization $\varphi_Q(m, (\sigma, \varkappa))$ and the fields are represented by sections $\mathbf{q}_f \in \mathfrak{X}(Q)$ of the principal bundle.*

The relativist momentum of particles is represented as vectors of the associated bundle $Q[E \otimes F, \vartheta]$

The observers measure the state of the particles in a standard gauge $\mathbf{q}(m) = \varphi_Q(m, (1, 1)) \in \mathfrak{X}(Q)$ and the measured states of particles in this gauge are $\psi(m) = \vartheta(\sigma(m), \varkappa(m)) \psi_0$

$Q[E \otimes F, \vartheta]$ has for trivialization :

$$(\varphi_Q(m, (1, 1)), \psi) \sim (\varphi_Q(m, (s^{-1}, g^{-1})), \vartheta(s, g) \psi)$$

and holonomic basis:

$$(\mathbf{e}_i(m) \otimes \mathbf{f}_j(m))_{i=0..3}^{j=1..n} = (\varphi_Q(m, (1, 1)), e_i \otimes f_j)$$

$$\psi(m) = \sum_{i=1}^4 \sum_{j=1}^n [\gamma C(\sigma(m))]_k^i [\varrho(\varkappa(m))]_l^j \psi_0^{kl}(m) \mathbf{e}_i(m) \otimes \mathbf{f}_j(m) \quad (5.5)$$

$$\text{in matrix form : } [\psi]_{4 \times n} = [\gamma C(\sigma)] [\psi] [\varrho(\varkappa)] \quad (5.6)$$

A change of trivialization with a section $\chi(m) \in \mathfrak{X}(Q)$ induces a change of gauge :

$$\mathbf{q}(m) = \varphi_Q(m, (1, 1)) \rightarrow \tilde{\mathbf{q}}(m) = \tilde{\varphi}_Q(m, (1, 1)) = \mathbf{q}(m) \cdot \chi(m)^{-1} \quad (5.7)$$

and a section $(\sigma, \varkappa) \in \mathfrak{X}(Q)$ takes a new expression :

$$(\sigma(m), \varkappa(m)) = \varphi_Q(m, (\sigma(m), \varkappa(m))) = \tilde{\varphi}_Q(m, \chi(m) \cdot (\sigma(m), \varkappa(m))) \quad (5.8)$$

It entails a change of holonomic basis in $E \otimes F$:

$$\mathbf{q}(m) = \varphi_Q(m, 1) \rightarrow \tilde{\mathbf{q}}(m) = \mathbf{q}(m) \cdot \chi(m)^{-1} :$$

$$\mathbf{e}_i(m) \otimes \mathbf{f}_j(m) = (\mathbf{p}(m), e_i \otimes f_j) \quad (5.9)$$

$$\rightarrow \tilde{\mathbf{e}}_i(m) \otimes \tilde{\mathbf{f}}_j(m) = \vartheta \left(\chi(m)^{-1} \right) (\mathbf{e}_i(m) \otimes \mathbf{f}_j(m)) \quad (5.10)$$

Sections $\psi \in \mathfrak{X}(Q[E \otimes F, \vartheta])$ have a new expression :

$$\mathbf{q}(m) = \varphi_Q(m, 1) \rightarrow \tilde{\mathbf{q}}(m) = \mathbf{q}(m) \cdot \chi(m)^{-1} :$$

$$[\psi(m)] \rightarrow [\tilde{\psi}(m)] = \vartheta(\chi(m)) [\psi(m)] = [\gamma C(s)] [\psi] [\varrho(g)] \quad (5.11)$$

$$\tilde{\psi}^{ij} = \sum_{k=1}^4 \sum_{l=1}^n [\gamma C(s)]_k^i [\varrho(g)]_l^j \psi^{kl}$$

The scalar product on $E \otimes F$ extends pointwise to $Q[E \otimes F, \vartheta]$:

$$\langle \psi(m), \psi'(m) \rangle = \text{Tr}([\psi(m)]^* [\gamma_0] [\psi'(m)])$$

It is preserved by ϑ .

The state of a particle along its world line is then represented by a path on the vector bundle :

$$\psi(\tau) = \vartheta(\tau) \psi_0 \text{ with } \vartheta(\tau) = \gamma C(\sigma(\tau), \rho(\varkappa(\tau))) \text{ and } \psi_0 \in \widehat{E}_0 \otimes F$$

$$\langle \psi, \psi \rangle = \langle \psi_0, \psi_0 \rangle = Ct \Leftrightarrow \text{Tr}([\psi]^* [\gamma_0] [\psi]) = \text{Tr}([\psi_0]^* [\gamma_0] [\psi_0]) \quad (5.12)$$

We will use the following bundles, which can be seen as restrictions of the previous ones :

By restriction to $\sigma = 1$ the principal bundle $Q(M, Spin(3, 1) \times U, \pi_U)$ is a principal bundle with fiber U , that we denote P_U with trivialization $\varphi_U(m, \varkappa)$.

A change of trivialization with a section $\chi(m) \in \mathfrak{X}(P_U)$ induces a change of gauge :

$$\mathbf{p}_U(m) = \varphi_U(m, 1) \rightarrow \tilde{\mathbf{p}}_U(m) = \tilde{\varphi}_U(m, 1) = \mathbf{p}_U(m) \cdot \chi(m)^{-1} \quad (5.13)$$

and a section $\varkappa \in \mathfrak{X}(P_U)$ takes a new expression :

$$\varkappa(m) = \varphi_U(m, \varkappa(m)) = \tilde{\varphi}_U(m, \chi(m) \cdot \varkappa(m)) \quad (5.14)$$

The principal fiber bundle P_U leads to the associated fiber bundle $P_U[F, \varrho]$ with holonomic basis : $\mathbf{f}_j(m) = (\mathbf{p}_U(m), f_j)$

A change of trivialization with a section $\chi(m) \in \mathfrak{X}(P_U)$ induces :
a change of holonomic basis in $E \otimes F$:

$$\mathbf{p}_U(m) = \varphi_{P_U}(m, 1) \rightarrow \tilde{\mathbf{p}}_U(m) = \mathbf{p}_U(m) \cdot \chi(m)^{-1} : \quad (5.15)$$

$$\mathbf{f}_j(m) = (\mathbf{p}(m), f_j) \rightarrow \tilde{\mathbf{f}}_j(m) = \varrho\left(\chi(m)^{-1}\right)(\mathbf{f}_j(m)) \quad (5.16)$$

a change in the expression of sections $\phi \in \mathfrak{X}(P_U[F, \varrho])$:

$$\mathbf{q}(m) = \varphi_U(m, 1) \rightarrow \tilde{\mathbf{q}}(m) = \mathbf{q}(m) \cdot \chi(m)^{-1} :$$

$$\mathbf{p}_U(m) = \varphi_{P_U}(m, 1) \rightarrow \tilde{\mathbf{p}}_U(m) = \mathbf{p}_U(m) \cdot \chi(m)^{-1} : \quad (5.17)$$

$$\phi(m) \rightarrow \tilde{\phi}(m) = \varrho(\chi(m))\phi(m) \quad (5.18)$$

5.2.5 Matter fields

The quantity ψ sums up everything (motion, kinematic, charge) about the particle. When particles are considered in a model they are naturally represented by their relativist momentum, which is a tensor ψ whose value can be measured at each point of its trajectory. So the most natural way to represent the particle is by a map : $\psi : [0, T] \rightarrow Q[E \otimes F, \vartheta]$ which can be parametrized either by the proper time or the time of the observer.

It is usual to consider models involving particles of the same type, submitted to similar conditions in a given area. Then, because they have the same behavior, one can assume that their trajectories can be represented by a unique vector field, and it is natural to represent their relativist momentum as sections of the fiber bundle $\mathfrak{X}(Q[E \otimes F, \vartheta])$.

Definition

Definition 92 A *matter field* is a section $\psi \in \mathfrak{X}(Q[E \otimes F, \vartheta])$ which, at each point, represents the relativist momentum of the same particle (or antiparticle). More precisely we will assume :

$$\begin{aligned} \exists (\sigma, \varkappa) \in \mathfrak{X}(Q), \exists \psi_0 \in E \otimes F : \psi_L = \epsilon i \psi_R :: \psi(m) = \vartheta(\sigma, \varkappa) \psi_0 \\ \int_{\Omega} \|\psi(m)\| \varpi_4(m) < \infty \end{aligned}$$

Notation 93 $\mathfrak{X}(M)$ is the set of matter fields, $\mathfrak{X}(\psi_0)$ the set of matter fields corresponding to $\psi_0 \in E \otimes F$.

For a matter field representing an elementary particle $\psi_0 \in \mathfrak{F}$. For a composite, as said before, the assumption that it can be represented as a particle implies then the existence of such a fundamental state, with a variable inertial spinor for a deformable solid (but there is still some fixed S_0).

A necessary condition to be a matter field is : $\langle \psi(m), \psi(m) \rangle = Ct$.

The set of matter fields is a subset of $\mathfrak{X}(Q[E \otimes F, \vartheta])$. This is not a vector space.

Mass, Spin and Charge of a matter field

We can proceed as for spinor fields.

The space $\mathfrak{X}(Q[E \otimes F, \vartheta])$ is a functional representation of $Spin(3, 1) \times U$ with the global action ϑ . The subset $\mathfrak{X}(\psi_0)$ is invariant by the right or left global actions of $Spin(3, 1) \times U$. Moreover the value of $Y(m) = \langle \psi(m), \psi(m) \rangle$ is invariant by $Spin(3, 1) \times U$.

The **mass of the particle** is defined as : $M_p = \frac{1}{c^2} \sqrt{|\langle \psi, \psi \rangle|}$ for $\psi \in \mathfrak{X}(\psi_0)$

The matter fields $\psi \in \mathfrak{X}(\psi_0)$ can equivalently be defined by a couple $(\psi_0, \sigma \times g)$ where $(\sigma \times g) \in \mathfrak{X}(Q)$. The representation is faithful : for given values of $\psi_0, \psi(m)$ there is a unique couple $(\sigma(m) \times g(m))$ and thus a unique $\sigma(m)$.

For a given observer $\sigma(m)$ admits two decompositions $\sigma(m) = \epsilon \sigma_w(m) \cdot \sigma_r(m)$. ϵ defines the relativist spin of the particle.

Define on $\mathfrak{X}(\psi_0)$ the equivalence relation :

$$\psi \sim \psi' \Leftrightarrow \forall m \in M : \sigma_r(m) = \sigma'_r(m)$$

Each class of equivalence is invariant by $Spin(3)$. The common value of $\psi = \gamma C(\sigma_r) \psi_0$ is **the Spin of the particle**, in his usual meaning. So to any given matter field correspond two Spins, with the Spin up or down.

Define on $\mathfrak{X}(\psi_0)$ the equivalence relation :

$$\psi \sim \psi' \Leftrightarrow \forall m \in M : \sigma_w(m) = \sigma'_w(m)$$

Each class of equivalence defines with the observer the same trajectories. The common value of $\psi = \gamma C(\sigma_w) \psi_0$ is **the translational momentum of the particle**, in his usual meaning.

Define on $\mathfrak{X}(\psi_0)$ the equivalence relation :

$$\psi \sim \psi' \Leftrightarrow \forall m \in M : g(m) = g'(m)$$

Each class of equivalence is invariant by U . The common value of $\psi = \varrho(g(m)) \psi_0$ is **the charge of the particle**. So the charge is not expressed by a scalar.

Quantization of the Spin and Charge

The set of sections $\mathfrak{X}(\psi_0)$ is not a vector space. However there is a norm on $E \otimes F$ invariant by ϑ . The space :

$L^1 = L^1(M, Q[E \otimes F, \vartheta]) = \{\psi \in \mathfrak{X}(Q[E \otimes F, \vartheta]) : \int_{\Omega} \|\psi(m)\| \varpi_4(m) < \infty\}$
is a separable Fréchet vector space. And we have the following :

Theorem 94 *The set $L^1(\psi_0) = L^1(M, Q[E \otimes U, \gamma C], \varpi_4) \cap \mathfrak{X}(\psi_0)$ of integrable spinor fields associated to a particle is characterized by 2 scalars : $k \in \mathbb{R}, z \in \mathbb{Z}$.*

The Spin, up or down, associated to each section is characterized by a scalar $j \in \frac{1}{2}\mathbb{N}$ and belongs to a $2j+1$ dimensional vector space isomorphic to (P^j, D^j)

The Charge of a matter field is characterized by a scalar and belongs to a finite dimensional vector space which is invariant by U

If the section ψ is continuous then the evaluation map : $\mathcal{E}(m) : L^1(\psi_0) \rightarrow E \otimes F :: \mathcal{E}(m)\psi = \psi(m)$ is continuous

Proof. i) For the first part the proof is the same as 80.

ii) For the charge, we add to the variable $\psi \in L^1(M, Q[E \otimes U, \gamma C], \varpi_4)$ the quantity $Z = \varrho(g(m))\psi_0$. For each value of Z we have a subset of the Hilbert space which is invariant by $\widehat{U} = \Upsilon \circ U \circ \Upsilon^{-1}$, so corresponds to an irreducible representation of U . U is compact, so Z belongs to one of the irreducible representation, which is a finite dimensional vector space characterized by scalars, and characteristic of ψ_0 .

iii) The space of continuous, compactly supported maps is dense in $L^1(M, E \otimes F, \varpi_4)$ (Maths.2292)

Let be ψ_n such a sequence converging to ψ in L^2

$(\psi - \psi_n, \psi - \psi_n)(m)$ is continuous, ≥ 0 on the open Ω so there are

$$A_n = \min_{m \in \Omega} (\psi - \psi_n, \psi - \psi_n)(m)$$

$$\int_{\Omega} A_n \varpi_4 \leq \int_{\Omega} (\psi - \psi_n, \psi - \psi_n) \varpi_4$$

$$\Rightarrow A_n \rightarrow 0$$

$$\Rightarrow \psi_n(m) \rightarrow \psi(m) \quad \blacksquare$$

Experiments show that the Spin of an elementary particle is $j = \frac{1}{2}$.

The scalar z in the representation of $\text{Spin}(3,1)$ corresponds to the charge of the particle. In each irreducible representation of U we can choose an orthonormal basis and the collection of these vectors is a basis of F , and z can label the irreducible representations. The Charge of a matter field is, as the Spin, a class of equivalence, which is represented as a section of $\mathfrak{X}(Q[E \otimes F, \vartheta])$ but the value of ψ itself is not simply a combination of the Spin and the charge. However the class of equivalence, and thus the charge, is constant.

When only the EM and gravitational fields are involved, then $E \otimes F$ is reduced to E and $z \in \mathbb{Z}$ is the charge of the particle, expressed in a suitable unit.

A matter field can be considered as a collection over Ω of tests particles, or as **virtual particles**. In a model where the relativist momentum of particles

are involved, a section of $\mathfrak{X}(\psi_0)$ is a general solution of a set of differential equations, whose solutions depend on the initial conditions.

Density of particles

A section $\psi \in \mathfrak{X}(Q[E \otimes F, \vartheta])$ or a map $\psi : \mathbb{R} \rightarrow Q[E \otimes F, \vartheta]$ such that $\langle \psi, \psi \rangle$ is not constant but has a constant sign can, formally, represent the state of a material body, however it is clear that the aggregation of particles has a physical meaning only for spinors (valued in E).

We can similarly consider a matter field with density μ . It has been introduced in the previous chapter to address the aggregation of particles in a solid body, but here we will consider the case of beams of particles, that is of a collection of particles of the same type which follow trajectories which do not cross. Their state can be represented by a matter field (with a constant ψ_0) and the density is then the number of particles by unit of volume. A solid body is characterized by the existence of a proper time, where the particles are localized on the space like hypersurface. A beam of particles will be defined as a collection of particles, continuously present over a period $[0, T]$ of an observer, so the time is measured with respect to an observer.

The chart which is used is the chart of the observer $\varphi_o(t, x), \varepsilon_0(m) = \partial\xi_0 = c\partial t$ and the volume is $\varpi_4 = \det[P'] d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3$.

The density is the number of particles by unit of 4 dimensional volume : on any area $\omega \subset \Omega$ this number is : $\int_\omega \mu(m) \varpi_4(m)$.

$$\mu(m) = \sqrt{\frac{\langle \psi(m), \psi(m) \rangle}{\langle \psi_0, \psi_0 \rangle}}$$

The velocity of the particles, with respect to the observer, is :

$$V = \frac{d\psi}{dt} = \vec{v} + c\varepsilon_0(m) = c \left((2a_w^2 - 1) \varepsilon_0 + \epsilon \frac{a_w}{2a_w^2 - 1} \sum_{i=1}^3 w_i \varepsilon_i \right)$$

The number of particles is constant, and the continuity equation is expressed by the Lie derivative of the volume form $\mu\varpi_4$ along the vector field $V : \mathcal{L}_V \mu\varpi_4 = 0$ (Maths.1524).

$$\mathcal{L}_V \mu\varpi_4 = 0 = \mu \mathcal{L}_V \varpi_4 + \varpi_4 \mathcal{L}_V \mu = \mu (\operatorname{div} V) \varpi_4 + \varpi_4 (\mu' (V)) = \mu (\operatorname{div} V) \varpi_4 + \frac{d\mu}{dV} \varpi_4$$

$$\mu \operatorname{div} V + \frac{d\mu}{dt} = 0 \quad (5.19)$$

Which is similar to the classic continuity equation in a fluid.

$$\begin{aligned} \operatorname{div} V &= \frac{1}{\det P'} \sum_{\alpha=0}^3 \partial_\alpha (V^\alpha \det P') \\ &= \frac{1}{\det P'} \sum_{\alpha=0}^3 V^\alpha \partial_\alpha (\det P') + \sum_{\alpha=0}^3 \partial_\alpha (V^\alpha) = \frac{1}{\det P'} \frac{d \det P'}{dt} + \sum_{\alpha=0}^3 \partial_\alpha (V^\alpha) \end{aligned}$$

$$\text{with } \mu(m) = \sqrt{\frac{\langle \psi, \psi \rangle}{\langle \psi_0, \psi_0 \rangle}}$$

$$\sqrt{|\langle \psi, \psi \rangle|} \left(\frac{1}{\det P'} \frac{d \det P'}{dt} + \sum_{\alpha=0}^3 \partial_\alpha (V^\alpha) \right) + \frac{1}{\sqrt{|\langle \psi, \psi \rangle|}} \frac{1}{2} \frac{d|\langle \psi, \psi \rangle|}{dt} = 0$$

$$\frac{1}{\det P'} \frac{d \det P'}{dt} + \sum_{\alpha=0}^3 \partial_\alpha (V^\alpha) + \frac{1}{2|\langle \psi, \psi \rangle|} \frac{d|\langle \psi, \psi \rangle|}{dt} = 0$$

$$\sum_{\alpha=0}^3 \partial_\alpha (V^\alpha) = -\frac{d}{dt} \ln \left(\sqrt{|\langle \psi, \psi \rangle|} \det P' \right)$$

Wave function

Usually one has a collection of particles of different types observed in a domain Ω , the goal of the experiment is to know the type and the motion of the particles. The states of the particles are represented by a unique section : $\psi \in L^1(M, Q[E \otimes F, \vartheta])$ and a primary observable is a linear map $\Phi : L^1(M, Q[E \otimes F, \vartheta]) \rightarrow V :: \Phi(\psi) = Y$ where V is a finite dimensional vector space. The observable can address some features of the particles only (such as the nature of the particles, their spin or charge,...).

There is a Hilbert space H associated to $L^1(M, Q[E \otimes F, \vartheta])$. This is an infinite dimensional, normed and separable vector space, and $E \otimes F$ is finite dimensional. The evaluation map $\mathcal{E}(m) : L^1(\psi_0) \rightarrow E \otimes F :: \mathcal{E}(m)\psi = \psi(m)$ is continuous. To Φ is associated the self adjoint operator $\hat{\Phi} = \Upsilon \circ \Phi \circ \Upsilon^{-1}$ on H .

We can apply the theorem 19. For any state ψ of the system there is a function : $W : M \times E \otimes F \rightarrow \mathbb{R}$ such that $W(m, Y) = \Pr(\Phi(\psi)(m) = y|\psi)$ is the probability that the measure of the value of the observable $\Phi(\psi)$ of ψ at m is y . It is given by :

$$\Pr(\Phi(\psi)(m) = y|\psi) = \frac{1}{\|\Upsilon(\psi)\|_H^2} \int_{Y \in \varpi(m, y)} \left\| \hat{\Phi}(\Upsilon(Y)) \right\|_H^2 \pi(Y) = W(m, y)$$

This can be seen as a density of probability, corresponding to the square of a wave function.

Of particular interest is the observable $\Phi(\psi) = \langle \psi, \psi \rangle$ which can be seen as the identification of the particles. The choice of the observable cannot be seen any longer as random. However one can assume that the choice of the point m is random. L^1 is partitioned in subsets $L^1(\psi_0)$ and any section ψ can be written as : $\psi(m) = \sum_j \varpi_j(m) \psi_j(m)$ where $\psi_j \in L^1(\psi_{0j})$ and $\varpi_j(m)$ is the characteristic function of the domain of ψ_j . Then the probability : $\Pr(\langle \psi(m), \psi(m) \rangle = \langle j, j \rangle | \psi) = \left(\int_{\Omega} \varpi_4 \right)^{-1} \int_{\Omega} \varpi_j \varpi_4$

5.3 CONNECTIONS

We know how the field changes the representation of the states of particles. But their action is not static and limited to ϑ . The propagation of the fields as well as their interaction with particles are, as usually, seen through infinitesimal changes and differential equations. So we need a mathematical way to define the derivative of $\mathbf{q}_f(m) = \varphi_Q(m, (\sigma(m), \varkappa(m)))$. There can be many different derivatives. Because of the anisotropy of the universe, the value of the derivative will depend on the direction on M , represented by a vector, so we are looking for a map : $M \rightarrow \Lambda_1(T_1Spin(3,1) \times T_1U)$, that is a one form valued in the Lie Algebra. This derivative is the **covariant derivative**. The action of a field on a particle, usually represented by a force $\vec{F} = \frac{d\vec{p}}{dt}$ which is the derivative of the momentum, is replaced by the covariant derivative $\nabla_V \psi$ of the relativist momentum along the direction V .

Covariant derivatives are built from more general mathematical objects, called connections. It will be necessary to refer to them, so it is better to introduce them from the start. This can be done in two complementary ways, geometric or through differential operators, and we need to remind some more mathematics.

5.3.1 Connections in Mathematics

Geometric connections

A fiber bundle $P(M, V, \pi)$ is a manifold, and its tangent space is split in two parts, related to its two manifolds components. By differentiation of the trivialization :

$$\varphi : M \times V \rightarrow P :: p = \varphi(m, u)$$

$$\varphi' : T_m M \times T_u V \rightarrow T_p P :: v_p = \varphi'_m(m, g) v_m + \varphi'_u(m, g) v_u$$

$\pi(p) = m \Rightarrow \pi'(p) v_p = v_m$ and the vector subspace $V_p P = \{\pi'(p) v_p = 0\}$ of $T_p P$ called the vertical space does not depend on the trivialization. It is isomorphic to the tangent space of V .

Our purpose is to look for a way to define a derivative of p , and the decomposition of the vector v_p shows that it requires two components : one linked to a motion in M , and another to a change in V . However, even if $\pi'(p) v_p = v_m$, this is not sufficient to define a decomposition which would be independent on the choice of a trivialization. A **connection** is just this tool : it is a projection of v_p on the vertical space $V_p P$. It is a one form on P valued in the vertical bundle VP . So it enables us to distinguish in a variation of p what can be imputed to a change of m and what can be imputed to a change of u . A section of P depends only on m : $\mathbf{p}(m) = \varphi(m, u(m))$ so by differentiation with respect to m this is a map from TM to TP and the value of a connection at each $\mathbf{p}(m)$ is a one form over M , valued in VP , called the **covariant derivative**. So it meets our purpose. Moreover because the vertical space is isomorphic to the tangent space on V , the value of the connection can be expressed in a simpler vector space.

All that has been said above holds for any kind of fiber bundle (Maths.27), but the connection takes different forms according to the kind of fiber bundle.

The covariant derivative issued from a linear connection on a vector bundle $P(M, V, \pi)$ reads:

$$\nabla X = \sum_{\alpha=0}^3 \sum_{a=1}^m \left(\partial_{\alpha} X^i(m) + \Gamma_{\alpha i}^j(m) X^j(m) \right) e_i \otimes d\xi^{\alpha}$$

where $\Gamma_{\alpha i}^j(m)$ is the Christoffel symbol of the connection. Its meaning is clear : the covariant derivative adds a linear combination of X to the derivative of X , this is the simplest form for the definition of a derivative on a fiber bundle. Readers who are familiar with GR are used to Christoffel symbols, and their definition through the metric. We will see how it works. All the connections that we will see depend on connections on principal bundles : a connection on a principal bundle leads to the definition of a covariant derivative on any associated bundle. Connections on associated vector bundles have the same form as above.

The second way to define the derivative is through differential operators, but for this we need to say a bit about r-jets.

r-jet extensions

In Differential Geometry one tries to avoid as much as possible the coordinates expressions. But when one deals with derivatives of orders higher than one this becomes impossible. It is always difficult to deal with partial derivatives, and notably to define the set in which they belong. The r-jet formalism provides a convenient solution. See Maths.26 for more.

For any r differentiable map $f \in C_r(M; N)$ between manifolds, the partial derivatives $\frac{\partial^s f}{\partial \xi^{\alpha_1} \dots \partial \xi^{\alpha_s}}$ are s symmetric linear maps from the tangent space $T_m M$ to the tangent space $T_p N$. Their expression in holonomic bases is a set of components $f_{\alpha_1 \dots \alpha_s}^i$ symmetric in the indices $\alpha_1, \dots, \alpha_s$. It defines fully the map $\frac{\partial^s f}{\partial \xi^{\alpha_1} \dots \partial \xi^{\alpha_s}}$ at m , and the set

$(p, f(p), f_{\alpha_1 \dots \alpha_s}^i, \alpha_k = 1.. \dim M, i = 1.. \dim N)$ is denoted $J_m^s(M, N)_{f(m)}$. Conversely one can conceive the same set of scalars and points of M,N, without any reference to the map f . It defines the class of equivalence of all maps which have the same derivatives up to order r, and is called the r-jet prolongation of $C_r(M; N)$. When one forgets the origin m and target $f(m)$ the set $J^s(M, N)$ is a vector space, and an affine space if one forgets only $f(m)$.

The formalism can be extended to fiber bundles, by replacing the maps f with sections on a fiber bundle. The **r-jet extension** $J^r P$ of the fiber bundle $P(M, V, \pi)$ is the r-jet prolongation of its sections. This is a closed manifold of the vector space $J^r(M, V)$. A map which associates to each $m \in M$ a set of values $Z = (m, z_{\alpha_1 \dots \alpha_s}^i, i = 1..n, s = 0, \dots, r)$ (called coordinates of the jet) in $J^r P$ is a **r-jet**, denoted j_m^r . A key point is that *there is a priori no relation between the $z_{\alpha_1 \dots \alpha_s}^i(m)$* : they do not correspond necessarily to the derivatives of the same map. But conversely a given section S of P provides a r-jet, denoted $J^r S$. A change of trivialization on the fiber bundle entails relations, usually complicated, between the elements $z_{\alpha_1 \dots \alpha_s}^i$, so they are not totally arbitrary.

The r-jet prolongation of a principal bundle is a principal bundle (but with a more complicated group), and the r-jet prolongation of an associated vector bundle is an associated vector bundle. Its elements have the coordinated expressions : $Z = (z_{\alpha_1 \dots \alpha_s}^i, i = 1 \dots n, s = 0, \dots, r)$. The index i refers to the component in the vector space V , the indices $\alpha_1, \dots, \alpha_s$ to the partial derivatives with respect to the coordinates in M .

The principal application of the r-jet formalism is in Differential Equations and Differential Operators.

Covariant derivatives as differential operators

A r differential operator is a base preserving morphism $D : \mathfrak{X}(J^r E_1) \rightarrow \mathfrak{X}(E_2)$ between two fiber bundles (Maths.32). It maps fiberwise $Z(m)$ in $J^r E_1$ to $Y(m)$ in E_2 . It is local : its computation involves only the values at m , and provides a result at m . By itself D does not involve any differentiation (it is defined for any section of the r-jet bundle $J^r E_1$). Combined with the map : $J^r : \mathfrak{X}(E_1) \rightarrow \mathfrak{X}(J^r E_1)$, $D \circ J^r$ maps sections on E_1 , to sections on E_2 . This distinction is useful : differential equations are equations between sections of $J^r E_1$ (the partial derivatives are, a priori, independent variables) and solutions are sections such that the components of the r-jet are deduced from the same map by differentiation. The most general mathematical definition of a set of differential equations of order r between variables defined on some fiber bundle E is that it is a closed subbundle of $J^r E$. The solutions are sections of E such that their r-jet extensions belong to this closed subbundle.

A linear r-differential operator is a linear, base preserving morphism, between two vector bundles (associated or not to a principal bundle, this does not matter here) : $E_1(M, V_1, \pi_1), E_2(M, V_2, \pi_2)$. The coordinates of a section $Z \in \mathfrak{X}(J^r E_1)$ read : $Z = (m, z_{\alpha_1 \dots \alpha_s}^i, i = 1 \dots n, s = 0, \dots, r)$ and DZ reads :

$$DZ = \sum_{s=0}^r \sum_{\alpha_i=1}^m \sum_{i=1}^n \sum_{j=1}^p A(m)_{i}^{j, \alpha_1 \dots \alpha_s} z_{\alpha_1 \dots \alpha_s}^i(m) e_{2j}(m)$$

with a basis $(e_{2j}(m))_{j=1}^p$ of E_2 , scalars $A(m)_{i}^{j, \alpha_1 \dots \alpha_s}$, and for a section $Z \in \mathfrak{X}(E_1) : z_{\alpha_1 \dots \alpha_s}^i(m) = \frac{\partial^s z^i}{\partial \xi^{\alpha_1} \dots \partial \xi^{\alpha_s}}$

In this framework it is easy to study the properties of Differential Operators such as action on distributions, adjoint of an operator, symbol, Fourier transform... We are concerned here with 1st order operators on a principal bundle.

A covariant derivative can then be considered as a differential operator (Maths.32.2.7) and indeed this is the way it is introduced in many books. For a vector bundle :

$$DZ = \sum_{\alpha=1}^m \sum_{i=1}^n \left(z_{\alpha}^i + \sum_{j=1}^p \Gamma_{\alpha i}^j z^j(m) \right) e_{2j}(m) \otimes d\xi^{\alpha}$$

and similarly for an associated vector bundle.

The Differential Operator approach is useful when one considers higher order derivatives (one can define higher order covariant derivative)

A general result

Using these tools we can prove the following theorem, for which there has been many demonstrations in specific cases :

Theorem 95 *Whenever a physical system is represented by a model with variables which are smooth sections $\mathbf{Z} \in \mathfrak{X}(E)$ of a vector bundle $E(\Omega, V, \pi)$ with fiber V on an open subset Ω of the manifold M , and their derivatives up to the order r , and the solutions are given by a set of differential equations, the system is determinist and the solutions depend linearly of the initial conditions.*

Proof. Using the previous notations, the initial conditions are defined in the fiber $\pi^{-1}(\Omega_3(0))$ of E over $\Omega_3(0)$. The set of differential equations is defined by a subbundle of $J^r E$, that is (Maths.2641) :

- a submanifold Ω_0 of Ω
- a vector bundle $E_0(\Omega_0, V_0, \pi_0)$ such that $\pi|_{\Omega_0} = \pi_0$,
- and there is a vector bundle morphism $F : E_0 \rightarrow J^r E$, that is a map such that :

$$\forall m \in \Omega_0 : F(\pi_0^{-1}(m)) \in \mathcal{L}(\pi_0^{-1}(m); \pi^{-1}(f(m)))$$

where : $f : \Omega_0 \rightarrow \Omega :: f \circ \pi_0 = \pi \circ F$

As E_0 is closed in $J^r E$ and π is an open map, Ω_0 is a closed submanifold of Ω , this is obviously $\Omega_3(0)$ and E_0 is the set of initial conditions. To each set of initial conditions is associated (by F) a section of $J^r E$. f is not necessarily surjective, but it is continuous and for $t' > t$, the state $S(t)$ of the system is in the fiber $\pi^{-1}(f(\Omega_3(0)) \cap S(t'))$ over $f(\Omega_3(0)) \cap \Omega_3(t')$, and is a continuously linear function of the initial conditions (by F). Thus the problem is well posed : the solutions depend continuously on the initial conditions, and moreover the relation is linear if we have a vector bundle. ■

This is in accordance with the usual assumption that, at least at an elementary level, the state of the system at a time t depends only on its value at some initial time, and not on its state at the intermediary times $0 < t' < t$. This characterizes continuous processes : they are defined by linear differential equations.

5.3.2 Connection for the force fields other than Gravity

A vector v_p on the tangent bundle $T_p P_U$ of the principal bundle $P_U(M, U, \pi_U)$ reads :

$$v_p = \varphi'_m(m, g) v_m + \varphi'_z(m, g) v_z = \sum_{\alpha=0}^3 v_m^\alpha \partial m_\alpha + \sum_{a=1}^m v_g^a \partial g_a$$

The vertical bundle VP_U is comprised of vectors $\varphi'_{Ug}(m, g) v_g$. It is isomorphic to the Lie algebra, and it is convenient to express v_u with respect to $T_1 U : v_g = L'_g(1) \theta$ where $L'_g(1)$ is the derivative of the left translation : $L : G \times G \rightarrow G :: L(g, h) = g \cdot h$ and $\theta \in T_1 U$ (Maths.1759).

$\zeta(\theta)(p) = \varphi'_{Ug}(m, g) L'_g(1) \theta = \sum_{a=1}^m v_a^u \partial g_a$ is called a **fundamental vector field** (they have many properties (Maths.1795)), in particular the map ζ does not depend on the trivialization.

Because the right action is the main characteristic of principal bundles, one distinguishes **principal connections**, which are equivariant by the right action. Then their value for any gauge of P_U can be defined through its value for $p = \varphi_U(m, 1)$. These are the only ones that we will use. Then a projection of $T_p P_U$ on $V_p P_U$ can be seen as a map : $T_p P_U \rightarrow T_1 U$. A principal connection $\dot{\mathbf{A}}$ on a principal bundle is uniquely defined by a family of maps $\dot{\mathbf{A}}$, called the **potential** of the connection :

$$\dot{\mathbf{A}} \in \Lambda_1(M; T_1 U) : TM \rightarrow T_1 U :: \dot{\mathbf{A}}(m) = \sum_{\alpha=0}^3 \sum_{a=1}^m \dot{A}_\alpha^a(m) \vec{\theta}_a \otimes d\xi^\alpha \quad (5.20)$$

such that in a change of trivialization on P_U :

$$\mathbf{p}_U(m) \rightarrow \tilde{\mathbf{p}}_U(m) = \tilde{\varphi}_U(m, 1) = \mathbf{p}_U(m) \cdot \chi(m)^{-1} : \quad (5.21)$$

$$\dot{\mathbf{A}}(m) \rightarrow \tilde{\dot{\mathbf{A}}}(m) = Ad_\chi \left(\dot{\mathbf{A}}(m) - L'_{\chi^{-1}}(\chi) \chi'(m) \right) \quad (5.22)$$

This is an affine law, which involves the derivative $\chi'(m)$ of the change of gauge, and this feature is at the origin of many specificities (and complications, such as the Higgs boson...). Notice that the potential is valued in the fixed vector space $T_1 U$.

Then the value of the connection $\dot{\mathbf{A}}$ is for $\varphi_U(m, g) = p$

$$\dot{\mathbf{A}}(p) (v^\alpha \partial m_\alpha + \zeta(v_g)(p)) = \zeta \left(v_g + Ad_{g^{-1}} \dot{\mathbf{A}}(m) v_m \right) (p) \quad (5.23)$$

The covariant derivative of a section $\mathbf{p}_f = \varphi_U(m, \varkappa(m)) \in \mathfrak{X}(P_U)$ (which could represent the field) is then :

$$\nabla^U \mathbf{p}_f = (L'_{\varkappa^{-1}} \varkappa)'(\varkappa'(m)) + \sum_{\alpha=0}^3 Ad_{\varkappa^{-1}} \dot{A}_\alpha(m) d\xi^\alpha \in \Lambda_1(M, T_1 U) \quad (5.24)$$

and for the holonomic gauge : $\mathbf{p}_U = \varphi_U(m, 1) : \nabla^U \mathbf{p}_U = \sum_{\alpha=0}^3 \dot{A}_\alpha(m) d\xi^\alpha$

With this connection on P_U it is possible to define a linear connection and a covariant derivative ∇^F , 1 form on M acting on sections $\phi(m) = \sum_{j=1}^n \phi^j(m) \mathbf{f}_j(m)$ of the associated vector bundle $P_U[F, \varrho]$ (Maths.27.4.2) :

$$\nabla^F \phi = \sum_{\alpha=0}^3 \left(\partial_\alpha \phi^i + \sum_{j=1}^n [\dot{A}_\alpha]_j^i \phi^j \right) \mathbf{f}_i(m) \otimes d\xi^\alpha \in \Lambda_1(M, P_U[F, \varrho]) \quad (5.25)$$

with the

Notation 96 $[\dot{A}_\alpha] = \sum_{a=1}^m \dot{A}_\alpha^a [\theta_a]$ is a $n \times n$ matrix representing $\varrho'(1) \dot{A}_\alpha \in \mathcal{L}(F; F)$

and $[\dot{A}_\alpha]_j^i$ has the same meaning as the Christoffel symbol of a linear connection.

A covariant derivative, when acting on a vector field $u \in TM$, becomes a section of the vector bundle $P_U[F, \rho]$, and transforms as such in a change of trivialization, so we have a map : $\mathfrak{X}(P_U[F, \rho]) \times \mathfrak{X}(TM) \rightarrow \mathfrak{X}(P_U[F, \rho])$. It meets our goal, and it can be proven that this is the only way to achieve it.

Practically this is the potential which represents the field. There has been some questions about the physical meaning of the potential. However some experiments such as Aharonov-Bohm's shows that, at least for the electromagnetic field, the potential is more than a simple formalism.

In QTF, because the groups are comprised of matrices with complex coefficients, and the elements of the Lie algebra T_1U are operators in the Hilbert spaces, it is usual to introduce the imaginary i everywhere, and to consider the complexified of the Lie algebra T_1U . However it is clear that the potential \dot{A}_α belongs to the real algebra, so it is a real quantity. And there are as many force carriers bosons (12) as the dimension of U

The electromagnetic field

The Lie algebra of $U(1)$ is \mathbb{R} . So the potential \dot{A} of the connection is a real valued one form on M : $\dot{A} = \sum_{\alpha=0}^3 \dot{A}_\alpha d\xi^\alpha \in \Lambda_1(M; \mathbb{R})$ which is usually represented as a vector field and not a form.

The action of $U(1)$ depends on the representation, thus on the charge of the particle :

- negative charge : $\varrho(\mathcal{X})\psi = (\exp i\mathcal{X})\psi$
- positive charge : $\varrho(\mathcal{X})\psi = (\exp(-i\mathcal{X}))\psi$
- neutral : $\varrho(\mathcal{X})\psi = \psi$

The matrix $[\theta_a] = \varrho'(1)\theta_a = i\theta_a$ so the covariant derivative reads :

$$\nabla_\alpha^F \psi = \partial_\alpha \psi + qi\dot{A}_\alpha \psi$$

where q is the charge (for composite bodies) expressed as a signed integer multiple of the negative elementary charge, and $q = 0$ for neutral particles.

5.3.3 The connection of the gravitational field

Potential

The principles are similar. The vertical bundle VP_G of the principal bundle $P_G(M, Spin(3, 1), \pi_G)$ is isomorphic to the Lie algebra $T_1Spin(3, 1)$.

The potential G of a principal connection \mathbf{G} on P_G is a map : $G \in \Lambda_1(M; T_1Spin(3, 1))$.

Using the Clifford algebra to represent the Lie algebra, G reads :

$$G \in \Lambda_1(M; T_1 Spin(3, 1)) : TM \rightarrow T_1 Spin(3, 1) :: \quad (5.26)$$

$$G(m) = \sum_{\alpha=0}^3 v(G_{r\alpha}(m), G_{w\alpha}(m)) d\xi^\alpha \quad (5.27)$$

$G_{r\alpha}(m), G_{w\alpha}(m)$ are two vectors $\in \mathbb{R}^3$. So the *gravitational field has a transversal ($G_{w\alpha}$) and a rotational ($G_{r\alpha}$) component*. This is the unavoidable consequence of the gauge group.

We will use also the equivalent notation :

$$G(m) = \sum_{a=1}^6 \sum_{\alpha=0}^3 G_\alpha^a(m) \vec{\kappa}_a \otimes d\xi^\alpha \quad (5.28)$$

$G_r(m) = \sum_{\alpha=0}^3 v(G_{r\alpha}(m), 0) d\xi^\alpha$ is a map $G \in \Lambda_1(M; T_1 Spin(3)) : TM \rightarrow T_1 Spin(3)$

In a change of gauge the potential transforms by an affine map :

$$\mathbf{p}(m) \rightarrow \tilde{\mathbf{p}}(m) = \mathbf{p}(m) \cdot \chi(m)^{-1} : \quad (5.29)$$

$$G(m) \rightarrow \tilde{G}(m) = \mathbf{Ad}_\chi \left(G(m) - L'_{\chi^{-1}}(\chi) \chi'(m) \right) \quad (5.30)$$

$$\mathbf{Ad}_\chi v(G_{r\alpha} - X_{r\alpha}, G_{w\alpha} - X_{w\alpha}) = v(\tilde{G}_r(m), \tilde{G}_w(m))$$

(see Annex for the values of X,Y)

We introduce the convenient notation that will be used in the following :

Notation 97 $v(\hat{G}_r(\tau), \hat{G}_w(\tau))$ is the value of the potential of the gravitational field along the integral curve $m(\tau) = \Phi_V(\tau, x)$ of any vector field V

$$v(\hat{G}_r(\tau), \hat{G}_w(\tau)) = \sum_{\alpha=0}^3 V^\alpha v(G_{r\alpha}(\tau), G_{w\alpha}(\tau))$$

$$\text{with : } V^\alpha = \sum_{i=0}^3 P_i^\alpha V^i$$

$$v(\hat{G}_r(\tau), \hat{G}_w(\tau)) = v([G_r][P][V]^a, [G_w][P][V]^a)$$

$$v(\hat{G}_r(\tau), \hat{G}_w(\tau)) \text{ can be extended to } v(G_r(m), G_w(m)) \text{ on } \Omega \text{ (Maths.1467).}$$

$$\text{And similarly : } \hat{A} = \sum_{\alpha=0}^3 V^\alpha \hat{A}_\alpha^a \vec{\theta}_a$$

There are several covariant derivatives deduced from this connection.

Covariant derivative on P_G

The connection acts on sections of the principal bundle, and the covariant derivative of $\sigma = \varphi_G(m, \sigma(m)) \in \mathfrak{X}(P_G)$ is (see general formula in the previous section) :

$$\nabla^G : \mathfrak{X}(P_G) \rightarrow \Lambda_1(M; T_1Spin) :: \quad (5.31)$$

$$\nabla^G \sigma = \sigma^{-1} \cdot \sigma' + \mathbf{Ad}_{\sigma^{-1}} G = \left(\sum_{\alpha=0}^3 \sigma^{-1} \cdot \partial_\alpha \sigma + \mathbf{Ad}_{\sigma^{-1}} G_\alpha \right) d\xi^\alpha \quad (5.32)$$

The covariant derivative is invariant in a change of gauge.

Proof. $\varphi_G(m, g) = \tilde{\varphi}_G(m, \chi(m)g)$

$$G(m) \rightarrow \tilde{G}(m) = Ad_\chi \left(G(m) - L'_{\chi^{-1}}(\chi) \chi'(m) \right)$$

$$\sigma(m) \rightarrow \tilde{\sigma}(m) = \chi(m) \cdot \sigma(m)$$

$$\nabla^G \sigma \rightarrow \widetilde{\nabla^G \sigma} = \tilde{\sigma}^{-1} \cdot \tilde{\sigma}' + \mathbf{Ad}_{\tilde{\sigma}^{-1}} \tilde{G}$$

$$= \sigma^{-1} \cdot \chi^{-1} \cdot (\chi' \cdot \sigma + \chi \cdot \sigma') + \mathbf{Ad}_{\sigma^{-1}} \mathbf{Ad}_{\chi^{-1}} (Ad_\chi (G - \chi^{-1} \cdot \chi'))$$

$$= \sigma^{-1} \cdot \chi^{-1} \cdot \chi' \cdot \sigma + \sigma^{-1} \cdot \chi^{-1} \cdot \chi \cdot \sigma' + \mathbf{Ad}_{\sigma^{-1}} (G - \chi^{-1} \cdot \chi')$$

$$= \sigma^{-1} \cdot \chi^{-1} \cdot \chi' \cdot \sigma + \sigma^{-1} \cdot \sigma' + \mathbf{Ad}_{\sigma^{-1}} G - \sigma^{-1} \cdot \chi^{-1} \cdot \chi' \cdot \sigma$$

$$= \sigma^{-1} \cdot \sigma' + \mathbf{Ad}_{\sigma^{-1}} G = \nabla^G \sigma \quad \blacksquare$$

The explicit formula is the following :

$$\nabla_\alpha^G \sigma = \sigma^{-1} \cdot \partial_\alpha \sigma + \mathbf{Ad}_{\sigma^{-1}} v(G_{r\alpha}, G_{w\alpha}) = v(X_\alpha, Y_\alpha) \quad (5.33)$$

with :

$$\sigma = a + v(r, w) + b\varepsilon_5$$

$$\sigma^{-1} \cdot \partial_\alpha \sigma =$$

$$v(-r\partial_\alpha a + a\partial_\alpha r + w\partial_\alpha b - b\partial_\alpha w + \frac{1}{2}(-j(r)\partial_\alpha r + j(w)\partial_\alpha w),$$

$$-w\partial_\alpha a + a\partial_\alpha w - r\partial_\alpha b + b\partial_\alpha r - \frac{1}{2}(j(w)\partial_\alpha r + j(r)\partial_\alpha w)$$

$$X_\alpha = -r\partial_\alpha a + a\partial_\alpha r + w\partial_\alpha b - b\partial_\alpha w + \frac{1}{2}(-j(r)\partial_\alpha r + j(w)\partial_\alpha w)$$

$$+ \left[(a+b)^2 - ab + \frac{1}{4}r^t r + (a+b)j(r) + \frac{1}{4}(j(w)j(r) - 2j(w)j(w) + 3j(r)j(r)) \right] G_{r\alpha}$$

$$+ \left[ab + \frac{1}{4}r^t r + (b-a)j(w) - \frac{1}{4}(3j(w)j(r) + 2j(r)j(w) - j(r)j(r)) \right] G_{w\alpha}$$

$$Y_\alpha = -w\partial_\alpha a + a\partial_\alpha w - r\partial_\alpha b + b\partial_\alpha r - \frac{1}{2}(j(w)\partial_\alpha r + j(r)\partial_\alpha w)$$

$$+ \left[\frac{1}{4}w^t w - ab + (a+b)j(w) + \frac{1}{4}(j(w)j(w) + 2j(w)j(r) + 3j(r)j(w)) \right] G_{r\alpha}$$

$$+ \left[(a-b)^2 - \frac{1}{4}w^t w - ab + (a-b)j(r) - \frac{1}{4}(3j(w)j(w) - j(r)j(w) - 2j(r)j(r)) \right] G_{w\alpha}$$

For the signature (1,3) change the sign of r.

Covariant derivative for spinors

The covariant derivative reads for a section $\mathbf{S} \in \mathfrak{X}(P_G [E, \gamma C])$:

$$\nabla^S S = \sum_{\alpha=0}^3 (\partial_\alpha S + \gamma C(G_\alpha) S) d\xi^\alpha = \sum_{\alpha=0}^3 (\partial_\alpha S + \gamma C(v(G_{r\alpha}, G_{w\alpha})) S) d\xi^\alpha \quad (5.34)$$

With the signature (3,1) :

$$\gamma C(v(G_{r\alpha}, G_{w\alpha})) = -i\frac{1}{2} \sum_{a=1}^3 (G_{w\alpha} \gamma_a \gamma_0 + G_{r\alpha} \tilde{\gamma}_a) \quad (5.35)$$

With the signature (1,3) :

$$\gamma C'(v(G_{r\alpha}, G_{w\alpha})) = -i\frac{1}{2} \sum_{a=1}^3 (G_{w\alpha} \gamma_a \gamma_0 - G_{r\alpha} \tilde{\gamma}_a) \quad (5.36)$$

So we go from the signature (3,1) to (1,3) by a change of the sign of $G_{r\alpha}$.

G_α being valued in $T_1Spin(3,1)$ and γC being a representation of the Clifford algebra the expression makes sense. Its coordinates expression is with right and left chiral parts:

$$\nabla^S S = \sum_{\alpha=0}^3 \left[\begin{array}{c} \partial_\alpha S_R + \frac{1}{2} \sum_{a=1}^3 (G_{w\alpha}^a - iG_{r\alpha}^a) \sigma_a S_R \\ \partial_\alpha S_L - \frac{1}{2} \sum_{a=1}^3 (G_{w\alpha}^a + iG_{r\alpha}^a) \sigma_a S_L \end{array} \right] d\xi^\alpha$$

It preserves the chirality.

In a change of gauge :

$$\mathbf{p}(m) \rightarrow \tilde{\mathbf{p}}(m) = \mathbf{p}(m) \cdot \chi(m)^{-1}$$

a section on $\mathfrak{X}(P_G[E, \gamma C])$ transforms as :

$$\mathbf{S}(m) = (\mathbf{p}(m), S(m)) = (\tilde{\mathbf{p}}(m), \tilde{S}(m)) \sim (\mathbf{p}(m), \gamma C(\chi(m)^{-1}) \tilde{S}(m))$$

$$\Rightarrow \tilde{S}(m) = \gamma C(\chi(m)) S(m)$$

The covariant derivative transforms as a section of $P_G[E, \gamma C]$:

Proof. $\nabla^S S \rightarrow \widetilde{\nabla^S S} = \sum_{\alpha=0}^3 (\partial_\alpha \tilde{S} + \gamma C(\tilde{G}_\alpha) \tilde{S}) d\xi^\alpha$

$$\begin{aligned} &= \sum_{\alpha=0}^3 (\gamma C(\partial_\alpha \chi) S + \gamma C(\chi) \partial_\alpha S + \gamma C(Ad_\chi(G - \chi^{-1} \partial_\alpha \chi)) \gamma C(\chi) S) d\xi^\alpha \\ &= \sum_{\alpha=0}^3 (\gamma C(\partial_\alpha \chi) S + \gamma C(\chi) \partial_\alpha S + \gamma C(\chi(G - \chi^{-1} \partial_\alpha \chi) \chi^{-1}) \gamma C(\chi) S) d\xi^\alpha \\ &= \sum_{\alpha=0}^3 (\gamma C(\partial_\alpha \chi) S + \gamma C(\chi) \partial_\alpha S + \gamma C(\chi) \gamma C(G) S - \gamma C(\partial_\alpha \chi) S) d\xi^\alpha \\ &= \sum_{\alpha=0}^3 \gamma C(\chi) (\partial_\alpha S + \gamma C(G) S) d\xi^\alpha = \gamma C(\chi) \nabla^S S \quad \blacksquare \end{aligned}$$

so the operator reads: $\nabla^S : \mathfrak{X}(P_G[E, \gamma C]) \rightarrow \wedge_1(M; \mathfrak{X}(P_G[E, \gamma C]))$

Covariant derivatives for vector fields on M

The connection on P_G induces a linear connection ∇^M on the associated vector bundle $P_G[\mathbb{R}^4, \mathbf{Ad}]$, which is TM with orthonormal bases, with Christoffel symbols :

$$\Gamma_M(m) = (\mathbf{Ad}_s)'_{s=1}(G(m))$$

with the product of vectors in $Cl(m)$:

$$v = \sum_{j=0}^3 v^j \varepsilon_j(m) \rightarrow$$

$$\sum_{i,j=0}^3 [\Gamma_G(m)]_i^j v^i \varepsilon_j(m) = v(G_{r\alpha}, G_{w\alpha}) \cdot v - v \cdot v(G_{r\alpha}, G_{w\alpha})$$

It is then more convenient to use the representation of $T_1Spin(3,1)$ by matrices of $so(3,1)$:

$$[\Gamma_{M\alpha}] = \sum_{a=1}^6 G_\alpha^a [\kappa_a] = \begin{bmatrix} 0 & G_{w\alpha}^1 & G_{w\alpha}^2 & G_{w\alpha}^3 \\ G_{w\alpha}^1 & 0 & -G_{r\alpha}^3 & G_{r\alpha}^2 \\ G_{w\alpha}^2 & G_{r\alpha}^3 & 0 & -G_{r\alpha}^1 \\ G_{w\alpha}^3 & -G_{r\alpha}^2 & G_{r\alpha}^1 & 0 \end{bmatrix}$$

In a change of gauge :

$$G(m) \rightarrow \tilde{G}(m) = Ad_\chi \left(G(m) - L'_{\chi^{-1}}(\chi) \chi'(m) \right)$$

$$[\tilde{\Gamma}_{M\alpha}] = [h(s)] ([\Gamma_{M\alpha}] - [h(s^{-1})] [h(s')])$$

The covariant derivative of a section $V \in \mathfrak{X}(P_G[\mathbb{R}^4, \mathbf{Ad}])$ is then :

$$\nabla^M V = \sum_{\alpha i=0}^3 \left(\partial_\alpha V^i + \sum_{j=0}^3 [\Gamma_{M\alpha}(m)]_j^i V^j \right) \varepsilon_i(m) \otimes d\xi^\alpha \quad (5.37)$$

For any vector field $W : \nabla_W^M : \mathfrak{X}(TM) \rightarrow \mathfrak{X}(TM)$ is a linear map which preserves the scalar product of vectors (Maths.2205):

$$\langle \nabla_W^M U, \nabla_W^M V \rangle = \langle U, V \rangle$$

The isomorphism $so(3,1) \rightarrow T_1 Spin(3,1) \subset Cl(\mathbb{R}, 3, 1)$ reads :

$$[J(r) + K(w)] \rightarrow v(r, w) = \frac{1}{4} \sum_{i=0}^3 ([J(r) + K(w)] [\eta])_j^i \varepsilon_i \cdot \varepsilon_j$$

thus in matrix form the Christoffel coefficient of the connection on $P_G[E, \gamma C]$

reads (Maths.9.2.4) :

$$\begin{aligned} [\Gamma_\alpha(m)] &= [\gamma C(G_\alpha)] \\ &= \frac{1}{4} \sum_{ijpq=0}^3 ([J(G_{r\alpha}) + K(G_{w\alpha})] [\eta])_j^i ([\gamma C(\varepsilon_i)] [\gamma C(\varepsilon_j)])_q^p \varepsilon_p(m) \otimes \varepsilon_q(m) \end{aligned}$$

But on the other hand the Christoffel coefficient of the connection on $P_G[\mathbb{R}^4, \mathbf{Ad}]$

is:

$$[\Gamma_{M\alpha}(m)] = \sum_{ij=0}^3 [K(G_{w\alpha}) + J(G_{r\alpha})]_j^i \varepsilon_i(m) \otimes \varepsilon_j(m)$$

thus :

$$[\Gamma_\alpha(m)] = \frac{1}{4} \sum_{ij=0}^3 ([\Gamma_{M\alpha}(m)] [\eta])_j^i [\gamma C(\varepsilon_i)] [\gamma C(\varepsilon_j)]$$

$P_G[E, \gamma C]$ is a spin bundle, and we have the identity between the derivatives

:

$$\forall V \in \mathfrak{X}(P_G[\mathbb{R}^4, \mathbf{Ad}]), S \in \mathfrak{X}(P_G[E, \gamma C]) :$$

$$\nabla(\gamma C(V)S) = \gamma C(\nabla^M V)S + \gamma C(V)\nabla S$$

which makes of \mathbf{G} a Clifford connection (Maths.2207).

5.3.4 Geodesics

There are several definitions of Geodesics, which, in different formulations, mean the curves of minimum length between two points. In Euclidean Geometry they are straight lines, in GR they are usually curves, and they play an important role because free particles move along geodesics. Moreover there is a unique geodesic passing through a point with a given tangent vector.

A connection enables to define the **parallel transport** of a vector (or a basis) along a curve (or a vector field).

Let C be a curve defined by a path $p : \mathbb{R} \rightarrow M : p(\tau)$ with $p(0) = a$, and a vector $v \in T_a M$. The parallel transported vector is given by a map :

$$V : \mathbb{R} \rightarrow T_{p(\tau)} M : V(\tau) \text{ such that } : \nabla_{\frac{dp}{d\tau}}^M V(\tau) = 0, V(0) = v$$

thus we have the differential equation with $V(\tau) = \sum_{i=0}^3 V^i(\tau) \varepsilon_i(p(\tau))$

$$\nabla_{\frac{dp}{d\tau}}^M V(\tau) = \sum_{\alpha i=0}^3 \left(\partial_\alpha V^i + \sum_{j=0}^3 \Gamma_M(p(\tau))_{\alpha j}^i V^j \right) \left(\frac{dp}{d\tau} \right)^\alpha \varepsilon_i(p(\tau)) = 0$$

$$\frac{dV^i}{d\tau} + \sum_{\alpha j=0}^3 \Gamma_M(p(\tau))_{\alpha j}^i V^j \left(\frac{dp}{d\tau} \right)^\alpha = 0$$

A **geodesic** is a path such that its tangent is parallel transported by the connection :

$$p : \mathbb{R} \rightarrow M : p(\tau) \text{ with } p(0) = a$$

$$V(\tau) = \frac{dp}{d\tau} = \sum_{i=0}^3 V^i(\tau) \varepsilon_i(p(\tau)) = \sum_{k\alpha=0}^3 V^k(\tau) P_k^\alpha(p(\tau)) \partial \xi_\alpha$$

$$\frac{dV^i}{d\tau} + \sum_{\alpha j k=0}^3 \Gamma_M(p(\tau))_{\alpha j}^i V^j(\tau) V^k(\tau) P_k^\alpha(p(\tau)) = 0$$

or in matrix form :

$$\left[\frac{dV}{d\tau} \right] + \sum_\alpha ([\Gamma_{M\alpha}] [V]) ([P'] [V])^\alpha = 0$$

The scalar product $\langle V, V \rangle$ is constant :

$$\frac{d}{d\tau} \langle V, V \rangle = \frac{d}{d\tau} \left([V]^t [\eta] [V] \right)$$

$$= - \sum_\alpha ([P'] [V])^\alpha [V]^t [\Gamma_{M\alpha}]^t [\eta] [V] - \sum_\alpha ([P'] [V])^\alpha [V]^t [\eta] ([\Gamma_{M\alpha}] [V])$$

$$= - \sum_\alpha ([P'] [V])^\alpha [V]^t \left([\Gamma_{M\alpha}]^t [\eta] + [\eta] [\Gamma_{M\alpha}] \right) [V] = 0$$

A **field of geodesics** is a vector field U such that it is parallel transported along its integral curves $p(\tau) = \Phi_U(\tau, x)$.

As $\langle U, U \rangle$ is constant, for a time like geodesic field we can take $\langle U, U \rangle = -1$ and, for a given observer, associate a section $\sigma_w \in P_W$:

$$U = \left((2a_w^2 - 1) \varepsilon_0(m) + a_w \sum_{j=1}^3 w_j \varepsilon_j(m) \right)$$

$$\text{and } \mathbf{U}(m) = \mathbf{Ad}_{\sigma_w} \varepsilon_0(m)$$

The formalism of vector bundles enables us to give a useful description of these geodesics, through the value of σ_w with respect to G .

Theorem 98 For a given observer geodesics are represented by sections $\sigma_w \in \mathfrak{X}(P_G)$ such that $\nabla_U^G \sigma_w \in T_1 \text{Spin}(3)$.

They are solutions of the differential equation :

$$a_w \frac{dw}{d\tau} =$$

$$\left((1 - a_w^2) a_w^2 + a_w j(w) - \frac{1}{4} a_w^2 j(w) j(w) \right) \widehat{G}_r + \left(3 - 4a_w^2 + \frac{1}{2} j(w) j(w) \right) \widehat{G}_w$$

where $v \left(\widehat{G}_r, \widehat{G}_w \right)$ is the value of the potential of the gravitational field along the geodesic

Proof. i) In the standard basis and with the Clifford algebra formalism :

$$\nabla_V^M U = \frac{dU}{d\tau} + \sum_{\alpha=0}^3 (V^\alpha (v(G_{r\alpha}, G_{w\alpha}) \cdot U - U \cdot v(G_{r\alpha}, G_{w\alpha})))$$

$$= \frac{d}{d\tau} \mathbf{Ad}_{\sigma_w} \varepsilon_0 + v \left(\widehat{G}_r, \widehat{G}_w \right) \cdot \mathbf{Ad}_{\sigma_w} \varepsilon_0 - \mathbf{Ad}_{\sigma_w} \varepsilon_0 \cdot v \left(\widehat{G}_r, \widehat{G}_w \right)$$

with $V^\alpha = \sum_i P_i^\alpha U^i$, $\widehat{G}_r = \sum_{\alpha=0}^3 G_{r\alpha} V^\alpha$; $\widehat{G}_w = \sum_{\alpha=0}^3 G_{w\alpha} V^\alpha$, $\mathbf{U}(m) = \mathbf{Ad}_{\sigma_w} \varepsilon_0(m)$

$$\begin{aligned}
\frac{d}{d\tau} \mathbf{Ad}_{\sigma_w} \varepsilon_0 &= \frac{d\sigma_w}{d\tau} \cdot \varepsilon_0 \cdot \sigma_w^{-1} - \sigma_w \cdot \varepsilon_0 \cdot \sigma_w^{-1} \cdot \frac{d\sigma_w}{d\tau} \cdot \sigma_w^{-1} \\
&= \left(\sigma_w \cdot \nabla_U^G \sigma_w - v \left(\widehat{G}_r, \widehat{G}_w \right) \cdot \sigma_w \right) \cdot \varepsilon_0 \cdot \sigma_w^{-1} - \sigma_w \cdot \varepsilon_0 \cdot \sigma_w^{-1} \cdot \left(\sigma_w \cdot \nabla_U^G \sigma_w - v \left(\widehat{G}_r, \widehat{G}_w \right) \cdot \sigma_w \right) \cdot \sigma_w^{-1} \\
&= \sigma_w \cdot \nabla_U^G \sigma_w \cdot \varepsilon_0 \cdot \sigma_w^{-1} - v \left(\widehat{G}_r, \widehat{G}_w \right) \cdot \sigma_w \cdot \varepsilon_0 \cdot \sigma_w^{-1} - \sigma_w \cdot \varepsilon_0 \cdot \nabla_U^G \sigma_w \cdot \sigma_w^{-1} + \\
&\sigma_w \cdot \varepsilon_0 \cdot \sigma_w^{-1} \cdot v \left(\widehat{G}_r, \widehat{G}_w \right)
\end{aligned}$$

$$\begin{aligned}
\text{with } \frac{d\sigma_w}{d\tau} &= \sigma_w \cdot \nabla_U^G \sigma_w - v \left(\widehat{G}_r, \widehat{G}_w \right) \cdot \sigma_w \\
\nabla_U^M U &= \sigma_w \cdot \nabla_U^G \sigma_w \cdot \varepsilon_0 \cdot \sigma_w^{-1} - \sigma_w \cdot \varepsilon_0 \cdot \nabla_U^G \sigma_w \cdot \sigma_w^{-1} \\
&- v \left(\widehat{G}_r, \widehat{G}_w \right) \cdot \sigma_w \cdot \varepsilon_0 \cdot \sigma_w^{-1} + \sigma_w \cdot \varepsilon_0 \cdot \sigma_w^{-1} \cdot v \left(\widehat{G}_r, \widehat{G}_w \right) \\
&+ v \left(\widehat{G}_r, \widehat{G}_w \right) \cdot \sigma_w \cdot \varepsilon_0 \cdot \sigma_w^{-1} - \sigma_w \cdot \varepsilon_0 \cdot \sigma_w^{-1} \cdot v \left(\widehat{G}_r, \widehat{G}_w \right) \\
&= \sigma_w \cdot \left(\nabla_U^G \sigma_w \cdot \varepsilon_0 - \varepsilon_0 \cdot \nabla_U^G \sigma_w \right) \cdot \sigma_w^{-1}
\end{aligned}$$

So, with the covariant derivative on the principal bundle P_G . We have a geodesic iff :

$$\nabla_U^G \sigma_w \cdot \varepsilon_0 - \varepsilon_0 \cdot \nabla_U^G \sigma_w = 0$$

that is iff $\nabla_u^G \sigma_w$ commutes with ε_0 .

For any element $v(r, w)$ of $T_1 Spin(3, 1)$ we have the identity :

$$v(r, w) \cdot \varepsilon_0 - \varepsilon_0 \cdot v(r, w) = w \text{ (see Annex for the proof)}$$

$$\text{So : } v(r, w) \in T_1 Spin(3) \Leftrightarrow v(r, w) \cdot \varepsilon_0 - \varepsilon_0 \cdot v(r, w) = 0 \Leftrightarrow w = 0$$

And the geodesics are represented by sections such that $\nabla_U^G \sigma_w \in T_1 Spin(3)$.

ii) The sections

$\nabla_\alpha^G \sigma_w = \sigma_w^{-1} \cdot (\partial_\alpha \sigma_w + v(G_{r\alpha}, G_{w\alpha}) \cdot \sigma_w)$ has been computed :

$$\begin{aligned}
\nabla_\alpha^G \sigma_w &= v \left(-\frac{1}{2} j(w) \partial_\alpha w + \left(a_w^2 + \frac{1}{4} - 2j(w)j(w) \right) G_{r\alpha} + a_w j(w) G_{w\alpha}, \right. \\
&\left. -w \partial_\alpha a_w + a_w \partial_\alpha w + \left(a_w^2 - 1 - a_w j(w) + \frac{1}{4} j(w)j(w) \right) G_{r\alpha} + \left(1 - \frac{3}{4} j(w)j(w) \right) G_{w\alpha} \right)
\end{aligned}$$

So geodesic fields are associated to the sections such that :

$$\begin{aligned}
\sum_\alpha V^\alpha \left(-w \partial_\alpha a_w + a_w \partial_\alpha w + \left(a_w^2 - 1 - a_w j(w) + \frac{1}{4} j(w)j(w) \right) G_{r\alpha} \right. \\
\left. + \left(1 - \frac{3}{4} j(w)j(w) \right) G_{w\alpha} \right) = 0
\end{aligned}$$

$$w \frac{da_w}{d\tau} - a_w \frac{dw}{d\tau} = \left(a_w^2 - 1 - a_w j(w) + \frac{1}{4} j(w)j(w) \right) \widehat{G}_r + \left(1 - \frac{3}{4} j(w)j(w) \right) \widehat{G}_w$$

By left multiplication with w^t :

$$w^t w \frac{da_w}{d\tau} - a_w w^t \frac{dw}{d\tau} = \left(a_w^2 - 1 \right) w^t \widehat{G}_r + w^t \widehat{G}_w$$

$$a_w^2 = 1 + \frac{1}{4} w^t w \Rightarrow \left(w^t \frac{dw}{dt} \right) = 4a_w \frac{da_w}{d\tau}$$

$$w^t w \frac{da_w}{d\tau} - 4a_w^2 \frac{da_w}{d\tau} = \left(a_w^2 - 1 \right) w^t \widehat{G}_r + w^t \widehat{G}_w$$

$$4 \left(a_w^2 - 1 - a_w^2 \right) \frac{da_w}{d\tau} = \left(a_w^2 - 1 \right) w^t \widehat{G}_r + w^t \widehat{G}_w$$

$$\frac{da_w}{d\tau} = \frac{1}{4} \left(1 - a_w^2 \right) w^t \widehat{G}_r - \frac{1}{4} w^t \widehat{G}_w$$

$$\frac{1}{4} \left(1 - a_w^2 \right) w w^t \widehat{G}_r - \frac{1}{4} w w^t \widehat{G}_w - a_w \frac{dw}{d\tau}$$

$$= \left(a_w^2 - 1 - a_w j(w) + \frac{1}{4} j(w)j(w) \right) \widehat{G}_r + \left(1 - \frac{3}{4} j(w)j(w) \right) \widehat{G}_w$$

$$\left(- \left(a_w^2 - 1 \right) a_w^2 + a_w j(w) - \frac{1}{4} a_w^2 j(w)j(w) \right) \widehat{G}_r$$

$$- \left(4 \left(a_w^2 - 1 \right) + 1 - \frac{1}{2} j(w)j(w) \right) \widehat{G}_w = a_w \frac{dw}{d\tau} \blacksquare$$

The interpretation of this result is clear : along the geodesic the forces, gravitational or inertial, are such that they induce spatial rotations, without change in the direction of the motion.

There is nothing equivalent for the null curves, such that their tangent vector u has a null scalar product : $\langle u, u \rangle = 0$. But the definition of the flow of a vector field, which does not involve the metric, still holds.

5.3.5 The Levi-Civita connection

In Differential Geometry one defines affine connections (Maths.1537), which are bilinear operators acting on vector fields (sections of the tangent bundle) $\nabla \in \mathcal{L}^2(\mathfrak{X}(TM), \mathfrak{X}(TM); \mathfrak{X}(TM))$ such that:

$$\begin{aligned} \forall f \in C_1(M; \mathbb{R}) : \\ \nabla_f V W = f \nabla_V W \\ \nabla_V f W = f \nabla_V W + (i_V df) W \end{aligned}$$

They read in holonomic basis of a chart :

$$\nabla_\alpha V = \sum_\beta (\partial_\beta V^\alpha + \sum_\gamma \Gamma_{\beta\gamma}^\alpha V^\gamma) \partial \xi^\beta \otimes d\xi_\alpha$$

with Christoffel symbols $\Gamma_{\beta\gamma}^\alpha(m)$ which change in a change of chart in a complicated way. So an affine connection is a covariant derivative, defined in the tangent bundle, and acting on sections of the tangent bundle, which are vector fields, or tensors. There can be many different affine connections.

An affine connection is said to be symmetric if $\Gamma_{\beta\gamma}^\alpha = \Gamma_{\gamma\beta}^\alpha$

When there is a metric (Riemannian or not) defined by a tensor g on a manifold, an affine connection is said to be metric if $\nabla_\alpha g = 0$: it preserves the scalar product of two vectors. With a metric one can define a unique, metric, symmetric connection, called the Levi-Civita connection. It reads (Maths.1626) :

$$\Gamma_{\beta\gamma}^\alpha = \frac{1}{2} \sum_\eta g^{\alpha\eta} (\partial_\beta g_{\gamma\eta} + \partial_\gamma g_{\beta\eta} - \partial_\eta g_{\beta\gamma})$$

And this has been the bread and butter of workers on GR for decenniums, in a formalism where the metric is at the core of the model.

With a principal bundle, and a principal connection, one can define covariant derivatives in any associated vector bundle, including of course the tangent bundle to M . And it has all the properties of the usual covariant derivative of affine connections. Connections on fiber bundles are a more general tool than usual affine connections which are strictly limited to the tangent bundle. We have seen that the connection \mathbf{G} on P_G induces a linear connection on $P_G[\mathbb{R}^4, \mathbf{Ad}]$, which is nothing more than TM with an orthonormal basis, and a covariant derivative ∇^M with Christoffel symbol Γ_M . By translating the orthonormal basis $(\varepsilon_i)_{i=0}^3$ into the holonomic basis $(\partial \xi_\alpha)_{\alpha=0}^3$ of any chart using the tetrad, a straightforward computation (Maths.2005) gives the Christoffel coefficients $\widehat{\Gamma}_{\alpha\beta}^\gamma$ of the affine connection Γ_M , expressed in the basis of the chart :

$$\widehat{\Gamma}_{\alpha\beta}^\gamma = P_i^\gamma (\partial_\alpha P_\beta^i + \Gamma_{M\alpha j}^i P_\beta^j)$$

In matrix form :

$$\widehat{\Gamma}_{\alpha\beta}^\gamma = \left[\widehat{\Gamma}_\alpha^\gamma \right]_\beta, \Gamma_{M\alpha j}^i = [\Gamma_{M\alpha}]_j^i,$$

$$[\Gamma_{M\alpha}] = \sum_{a=1}^6 G_{a\alpha} [\kappa_a]$$

$$\left[\widehat{\Gamma}_\alpha \right] = [P] ([\partial_\alpha P'] + [\Gamma_{M\alpha}] [P']) \Leftrightarrow [\Gamma_{M\alpha}] = \left([P'] \left[\widehat{\Gamma}_\alpha \right] - [\partial_\alpha P'] \right) [P]$$

with :

$$a=1,2,3 : [\kappa_a]_q^p = \sum_{bc=1}^3 \epsilon(a,b,c) \delta_c^p \delta_q^b$$

$$a=4,5,6 : ([\kappa_a]_q^p)^p = \delta_0^p \delta_q^{a-3} + \delta_{a-3}^p \delta_q^0$$

$$[\widehat{\Gamma}_\alpha]_\beta^\gamma = ([P] [\partial_\alpha P'])_\beta^\gamma$$

$$+ \sum_{a=1}^3 \left(\sum_{bc=1}^3 \epsilon(a,b,c) G_{r\alpha}^a [P]_c^\gamma [P']_\beta^b + G_{w\alpha}^a \left([P]_0^\gamma [P']_\beta^a + [P]_a^\gamma [P']_\beta^0 \right) \right)$$

Any affine connection deduced this way from a principal connection is necessarily metric, but it is not necessarily symmetric.

To sum up :

- affine connections are defined in the strict framework of the tangent bundle, and the Levi-Civita connection is one of these connections, with specific properties (it is metric and symmetric); the covariant derivative which is deduced acts only on vectors fields (or tensors) of the tangent bundle.

- connections on principal bundle define connections on any associated vector bundle and act on sections of these bundles. So one can compute a covariant derivative acting on vectors fields of the tangent bundle, which is necessarily metric but not necessarily symmetric.

So, using the formalism of fiber bundles we do not miss anything, we can get the usual results, but in a more elegant and simple way. One can require from the principal connection \mathbf{G} on P_G that the induced connection on TM is symmetric, which will then be identical to the Levi-Civita connection. This requests :

$$\forall \alpha, \beta, \gamma :$$

$$[\widehat{\Gamma}_\alpha]_\beta^\gamma = ([P] ([\partial_\alpha P'] + [\Gamma_{M\alpha}] [P']))_\beta^\gamma = [\widehat{\Gamma}_\beta]_\alpha^\gamma = ([P] ([\partial_\beta P'] + [\Gamma_{M\beta}] [P']))_\alpha^\gamma$$

which has no obvious meaning for Γ_M .

Actually the Levi-Civita connection is traditionally used because it is the natural mathematical choice when one starts from the metric. Moreover it is assumed that the gravitational field (whose action goes through the connection) acts symmetrically, in the meaning that it has no torsion (or no torque). But actually this assumption has not been verified (which is difficult), and different theories have been proposed, notably by Einstein and Cartan, which consider connections with torsion, that is connections other than the Levi-Civita connection. However, when starting from the metric, they lead mostly to more complicated computations, in what is already a dreadful endeavour. In the fiber bundle framework there is no such problem and actually it would be the requirement of symmetry, always possible at any point, which would introduce a complication. Moreover the introduction of spinors and the distinction of the components G_r, G_w of the connection, are a more efficient way to deal with rotation and torque so it is justified that we keep the more general connection. An additional argument is that the Levi-Civita connection does not make any distinction between the bases, which can be induced by any chart. But, as we have seen, there is always a privileged chart, that of the observer, and the use of an orthogonal basis, in the fiber bundle formalism, is a useful reminder of this feature.

5.3.6 The total connection

Definition

The covariant derivative is computed from the action of the Lie algebra of the group $Spin(3, 1) \times U$, that is by derivation of the action $\vartheta(\sigma, \varkappa)\psi$. And by combining the previous result we have :

Proposition 99 *There are on Q a principal connection defined by the potentials*

$$G \in \Lambda_1(M; T_1 Spin(3, 1)) : TM \rightarrow T_1 Spin(3, 1) ::$$

$$G(m) = \sum_{\alpha=0}^3 v(G_{r\alpha}(m), G_{w\alpha}(m)) d\xi^\alpha$$

$$\dot{A} \in \Lambda_1(M; T_1 U) : TM \rightarrow T_1 U :: \dot{A}(m) = \sum_{\alpha=0}^3 \sum_{a=1}^m \dot{A}_\alpha^a(m) \theta_a \otimes dm^\alpha$$

The action of the fields on the state of a particle is given by the covariant derivative.

It reads for $\psi \in \mathfrak{X}(Q[E \otimes F, \vartheta])$:

$$\nabla_\alpha \psi = \sum_{\alpha=0}^3 \left(\partial_\alpha \psi^{ij} + \sum_{a=1}^6 [\gamma C(G_\alpha)] [\psi] + \sum_{a=1}^m \dot{A}_\alpha^a [\psi] [\theta_a] \right)^{ij} \mathbf{e}_i(m) \otimes \mathbf{f}_i(m) \otimes d\xi^\alpha$$

In matrix form :

$$\nabla_\alpha \psi \in \Lambda_1(M, Q[E \otimes F, \vartheta]) : [\nabla_\alpha \psi] = \sum_{\alpha=0}^3 [\partial_\alpha \psi] + [\gamma C(G_\alpha)] [\psi] + [\psi] \left[\dot{A}_\alpha \right] \quad (5.38)$$

$$\text{with } [\gamma C(G_\alpha)] = [\gamma C(v(G_{r\alpha}, G_{w\alpha}))]$$

$$[\nabla_\alpha \psi] = [\partial_\alpha \psi] + \sum_{a=1}^m \dot{A}_\alpha^a [\psi] [\theta_a] - \frac{i}{2} \left(\sum_{a=1}^3 G_{w\alpha}^a [\gamma_a] [\gamma_0] + G_{r\alpha} [\tilde{\gamma}_a] \right) [\psi] \quad (5.39)$$

The action, that is the 4 dimensional force exercised on a particle with velocity u by the fields represented by G, \dot{A} is then given by :

$\psi \rightarrow \nabla_u \psi = \frac{d\psi}{d\tau} + [\gamma C(\hat{G})] [\psi] + [\psi] [\hat{A}]$ which is the equivalent of the Newton's law $F = \frac{dP}{dt} = u \nabla P$ with the operator $\nabla = \partial_\alpha$

It depends linearly on the potentials, and on the state of the particle and its derivative. Even if the path has been long, this is the simplest expression that one can imagine. The inertial or gravitational forces depend only on the inertial spinor, and the other forces on the charges.

Systemic energy of a particle

We have an important result :

Theorem 100 *The scalar product $\langle \psi, \nabla_\alpha \psi \rangle$ is purely imaginary : $\langle \psi, \nabla_\alpha \psi \rangle = i \text{Im} \langle \psi, \nabla_\alpha \psi \rangle$*

Proof. $[\nabla_\alpha \psi] = [\partial_\alpha \psi] + [\psi] \left[\dot{A}_\alpha \right] - \frac{i}{2} \sum_{a=1}^3 G_{w\alpha}^a \gamma_a \gamma_0 [\psi] + G_{r\alpha} \tilde{\gamma}_a [\psi]$

$$\begin{aligned} \langle \psi, \nabla_\alpha \psi \rangle &= Tr [\psi]^* \gamma_0 [\partial_\alpha \psi] + Tr [\psi]^* \gamma_0 [\psi] \left[\dot{A}_\alpha \right] \\ &\quad - \frac{i}{2} \sum_{a=1}^3 G_{w\alpha}^a Tr [\psi]^* \gamma_0 \gamma_a \gamma_0 [\psi] + G_{r\alpha} Tr [\psi]^* \gamma_0 \tilde{\gamma}_a [\psi] \\ &= Tr [\psi]^* \gamma_0 [\partial_\alpha \psi] + Tr [\psi]^* \gamma_0 [\psi] \left[\dot{A}_\alpha \right] - \frac{i}{2} \sum_{a=1}^3 \left(-G_{w\alpha}^a Tr [\psi]^* \gamma_a [\psi] + G_{r\alpha} Tr [\psi]^* \gamma_0 \tilde{\gamma}_a [\psi] \right) \\ &\quad Tr [\psi]^* \gamma_a [\psi], Tr [\psi]^* \gamma_0 \tilde{\gamma}_a [\psi] \text{ are real, } Tr [\psi]^* \gamma_0 [\psi] [\theta_a], Tr [\psi]^* \gamma_0 [\partial_\alpha \psi] \text{ are} \\ &\text{imaginary :} \\ \overline{Tr [\psi]^* \gamma_a [\psi]} &= Tr ([\psi]^* \gamma_a [\psi])^* = Tr [\psi]^* \gamma_a [\psi] \\ \overline{Tr [\psi]^* \gamma_0 \tilde{\gamma}_a [\psi]} &= Tr ([\psi]^* \gamma_0 \tilde{\gamma}_a [\psi])^* = Tr [\psi]^* (\gamma_0 \tilde{\gamma}_a)^* [\psi] = Tr [\psi]^* \gamma_0 \tilde{\gamma}_a [\psi] \\ \overline{Tr [\psi]^* \gamma_0 [\psi] [\theta_a]} &= Tr ([\psi]^* \gamma_0 [\psi] [\theta_a])^* = Tr [\theta_a]^* [\psi]^* [\gamma_0] [\psi] \\ &= -Tr [\theta_a] [\psi]^* [\gamma_0] [\psi] = -Tr [\psi]^* [\gamma_0] [\psi] [\theta_a] \\ \langle \psi, \psi \rangle &= \langle \psi_0, \psi_0 \rangle \Rightarrow \langle \psi, \partial_\alpha \psi \rangle + \langle \partial_\alpha \psi, \psi \rangle = 0 \quad \blacksquare \end{aligned}$$

$$\text{Im} \langle \psi, \nabla_\alpha \psi \rangle = \frac{1}{i} \left(\langle \psi, \partial_\alpha \psi \rangle + \left\langle \psi, [\psi] \left[\dot{A}_\alpha \right] \right\rangle + \langle \psi, [\gamma C (G_\alpha)] \psi \rangle \right) \quad (5.40)$$

Along the trajectory, defined by a vector ε_0 and the time t of an observer :

$$\text{Im} \langle \psi, \nabla_V \psi \rangle = \frac{1}{i} \left(\left\langle \psi, \frac{d\psi}{dt} \right\rangle + \left\langle \psi, [\psi] \left[\hat{A} \right] \right\rangle + \left\langle \psi, \gamma C \left(\hat{G} \right) \psi \right\rangle \right)$$

$\frac{1}{i} \left\langle \psi, [\psi] \left[\hat{A} \right] \right\rangle + \left\langle \psi, \gamma C \left(\hat{G} \right) \psi \right\rangle$ can be seen as the **potential energy** of the particle in the fields along its trajectory

$\frac{1}{i} \left\langle \psi, \frac{d\psi}{dt} \right\rangle$ can be seen as the **kinetic energy** of the particle with respect to the observer

$\text{Im} \langle \psi, \nabla_V \psi \rangle$ can be seen as the **systemic energy** of the particle, as part of a system which includes the fields. The energy depends on the choice of an observer : V is defined by σ_w with respect to ε_0 .

Inertial tensor

We can give a more precise expression of the covariant derivative, using the fundamental state ψ_0 :

$$\begin{aligned} &\partial_\alpha ([\gamma C (\sigma)] [\psi_0] [\varrho (\varkappa)]) \\ &= [\gamma C (\partial_\alpha \sigma)] [\psi_0] [\varrho (\varkappa)] + [\gamma C (\sigma)] [\psi_0] [\varrho' (\varkappa) \partial_\alpha \varkappa] \\ &[\varrho' (\varkappa) \partial_\alpha \varkappa] = [\varrho (\varkappa) \varrho' (1) (L'_{\varkappa^{-1}} \varkappa) \partial_\alpha \varkappa] = [\varrho (\varkappa)] \sum_{a=1}^m \partial_\alpha \varkappa^a [\theta_a] = [\varrho (\varkappa)] [\partial_\alpha \varkappa] \\ &\text{(Maths.1900)} \\ &\nabla_\alpha \psi = ([\gamma C (\partial_\alpha \sigma)] [\psi_0] [\varrho (\varkappa)]) + [\gamma C (\sigma)] [\psi_0] [\varrho (\varkappa)] [\partial_\alpha \varkappa] \\ &\quad + [\gamma C (G_\alpha)] [\gamma C (\sigma)] [\psi_0] [\varrho (\varkappa)] + [\gamma C (\sigma)] [\psi_0] [\varrho (\varkappa)] \left[\dot{A}_\alpha \right] \\ &= [\gamma C (\sigma)] \{ [\gamma C (\sigma^{-1} \cdot \partial_\alpha \sigma)] [\psi_0] + [\gamma C (\sigma^{-1} \cdot G_\alpha \cdot \sigma)] [\psi_0] + [\psi_0] [\varrho (\varkappa)] [\partial_\alpha \varkappa] [\varrho (\varkappa)]^{-1} + \\ &\quad [\psi_0] [\varrho (\varkappa)] \left[\dot{A}_\alpha \right] [\varrho (\varkappa)]^{-1} \} [\varrho (\varkappa)] \\ &\nabla_\alpha \psi = [\gamma C (\sigma)] \{ [\gamma C (\sigma^{-1} \cdot \partial_\alpha \sigma + Ad_{\sigma^{-1}} G_\alpha)] [\psi_0] + [\psi_0] \left(Ad_\varkappa [\partial_\alpha \varkappa] + \left[\dot{A}_\alpha \right] \right) \} [\varrho (\varkappa)] \\ &\sigma^{-1} \partial_\alpha \sigma + Ad_{\sigma^{-1}} G_\alpha \in T_1 Spin(3, 1), Ad_\varkappa \left(\partial_\alpha \varkappa + \dot{A}_\alpha \right) \in T_1 U \end{aligned}$$

We see that the action of the fields is expressed by $\partial_\alpha \varkappa + \dot{A}_\alpha \in T_1 U$, so actually, *as far as the covariant derivative is involved, the potential is defined up to $\partial_\alpha \varkappa$, and one can take $\partial_\alpha \varkappa = 0$.*

$$\nabla_\alpha \psi = \vartheta(\sigma, \varkappa) \left([\gamma C (\sigma^{-1} \partial_\alpha \sigma + Ad_{\sigma^{-1}} G_\alpha)] [\psi_0] + [\psi_0] [Ad_\varkappa (\dot{A}_\alpha)] \right) \quad (5.41)$$

By combination with the previous result :

$$\begin{aligned} & \text{Im} \langle \psi, \nabla_\alpha \psi \rangle \\ &= \frac{1}{i} \left\langle \vartheta(\sigma, \varkappa) \psi_0, \vartheta(\sigma, \varkappa) \left([\gamma C (\sigma^{-1} \partial_\alpha \sigma + Ad_{\sigma^{-1}} G_\alpha)] [\psi_0] + [\psi_0] [Ad_\varkappa (\dot{A}_\alpha)] \right) \right\rangle \\ &= \frac{1}{i} \left\langle \psi_0, [\gamma C (\sigma^{-1} \partial_\alpha \sigma + Ad_{\sigma^{-1}} G_\alpha)] [\psi_0] + [\psi_0] [Ad_\varkappa (\dot{A}_\alpha)] \right\rangle \\ &= \frac{1}{i} \left\langle \psi_0, [\gamma C (\sigma^{-1} \partial_\alpha \sigma + Ad_{\sigma^{-1}} G_\alpha)] [\psi_0] \right\rangle + \frac{1}{i} \left\langle \psi_0, [\psi_0] [Ad_\varkappa (\dot{A}_\alpha)] \right\rangle \end{aligned}$$

Let us denote $[\psi_0] = \begin{bmatrix} \psi_R \\ \epsilon i \psi_R \end{bmatrix}$ with ψ_R a $2 \times n$ matrix and $\epsilon = +1$ for particles, and $\epsilon = -1$ for antiparticles.

$$(\sigma^{-1} \partial_\alpha \sigma + Ad_{\sigma^{-1}} G_\alpha) = v(X_\alpha, Y_\alpha) \in T_1 Spin(3, 1)$$

$$\gamma C(v(X_\alpha, Y_\alpha)) = \frac{1}{2} \begin{bmatrix} \sum_{a=1}^3 (Y_\alpha^a - i X_\alpha^a) \sigma_a & 0 \\ 0 & -\sum_{a=1}^3 (Y_\alpha^a + i X_\alpha^a) \sigma_a \end{bmatrix}$$

$$\gamma C(v(X_\alpha, Y_\alpha)) [\psi_0] = \frac{1}{2} \begin{bmatrix} \sum_{a=1}^3 (Y_\alpha^a - i X_\alpha^a) \sigma_a \psi_R \\ -\epsilon i \sum_{a=1}^3 (Y_\alpha^a + i X_\alpha^a) \sigma_a \psi_R \end{bmatrix}$$

$$\langle \psi_0, \gamma C(\sigma^{-1} \partial_\alpha \sigma + Ad_{\sigma^{-1}} G_\alpha) \psi_0 \rangle$$

$$= \frac{1}{2} Tr \left(\psi_R^* \left(-\epsilon \sum_{a=1}^3 (Y_\alpha^a + i X_\alpha^a) \sigma_a \psi_R \right) - \epsilon i \psi_R^* \left(i \sum_{a=1}^3 (Y_\alpha^a - i X_\alpha^a) \sigma_a \psi_R \right) \right)$$

$$= \epsilon \frac{1}{2} \left(-\sum_{a=1}^3 (Y_\alpha^a + i X_\alpha^a) Tr(\psi_R^* \sigma_a \psi_R) + \sum_{a=1}^3 (Y_\alpha^a - i X_\alpha^a) Tr(\psi_R^* \sigma_a \psi_R) \right)$$

$$= \frac{1}{2} \epsilon \sum_{a=1}^3 \left(-(Y_\alpha^a + i X_\alpha^a) + (Y_\alpha^a - i X_\alpha^a) \right) Tr(\psi_R^* \sigma_a \psi_R)$$

$$= -\epsilon i \sum_{a=1}^3 X_\alpha^a Tr(\psi_R^* \sigma_a \psi_R)$$

$$= i k^t X_\alpha$$

where

$$k = -\epsilon (Tr(\psi_R^* \sigma_a \psi_R))_{a=1}^3 \quad (5.42)$$

are fixed, real scalars.

Thus :

$$\langle \psi_0, \gamma C(\sigma^{-1} \partial_\alpha \sigma + Ad_{\sigma^{-1}} G_\alpha) \psi_0 \rangle = i k^t X_\alpha \quad (5.43)$$

with both vectors k, X_α in \mathbb{R}^3 and :

$$\sum_{\alpha=0}^3 V^\alpha X_\alpha = \hat{X}$$

The systemic energy of the particle has the simple expression :

$$\frac{1}{i} \langle \psi, \nabla_V \psi \rangle = k^t \hat{X} + \frac{1}{i} \left\langle \psi_0, [\psi_0] [Ad_\varkappa \hat{A}] \right\rangle \quad (5.44)$$

For Spinors, with : $\langle S_0, S_0 \rangle = \epsilon M_p^2 c^4$

$$\psi_R = \frac{M_p c^2}{\sqrt{2}} \begin{bmatrix} e^{i\alpha_1} \cos \alpha_0 \\ e^{i\alpha_2} \sin \alpha_0 \end{bmatrix} = \begin{bmatrix} x_1 + iy_1 \\ x_2 + iy_2 \end{bmatrix}$$

$$k = -\epsilon \frac{1}{2} M_p^2 c^4 k_0 = -\epsilon \frac{1}{2} M_p^2 c^4 \begin{bmatrix} (\sin 2\alpha_0) \cos(\alpha_2 - \alpha_1) \\ (\sin 2\alpha_0) \sin(\alpha_2 - \alpha_1) \\ \cos 2\alpha_0 \end{bmatrix}; k_0^t k_0 = 1$$

The vector k is similar to the inertia tensor, but here this is a vector, and it encompasses both the translational and the rotational motions. And of course it holds for a particle without internal structure but can be computed for a solid body using the aggregation method.

Using $\sigma = \sigma_w \cdot \sigma_r$ one can compute X_α (see Annex). It reads :

$$X_\alpha = [A_r(r)] \partial_\alpha r + [A_w(r, w)] \partial_\alpha w + [D_r(r, w)] G_{r\alpha} + [D_w(r, w)] G_{w\alpha} \quad (5.45)$$

$$[A_r(r)] = \left[\frac{1}{a_r} + \frac{2-5a_r^2}{2} j(r) + \frac{5a_r^2+1}{4a_r} j(r) j(r) \right]$$

$$[A_w(r, w)] = [B_r(r)] [B_w(w)] + [C_r(r)] [C_w(w)]$$

$$[D_r(r, w)] = [B_r(r)] [B_{wr}(w)] + [C_r(r)] [C_{wr}(w)]$$

$$[D_w(r, w)] = [B_r(r)] [B_{ww}(w)] + [C_r(r)] [C_{ww}(w)]$$

with :

$$[B_r(r)] = \left[1 - a_r j(r) + \frac{3}{4} j(r) j(r) \right]$$

$$[B_w(w)] = \frac{1}{2} j(w)$$

$$[C_r(r)] = \left[1 - a_r^2 + \frac{1}{4} j(r) j(r) \right]$$

$$[C_w(w)] = \frac{1}{a_w} \left[1 - \frac{1}{4} j(w) j(w) \right]$$

$$[B_{wr}(w)] = \left[a_w^2 - \frac{1}{2} j(w) j(w) \right]$$

$$[B_{ww}(w)] = a_w [j(w)]$$

$$[C_{wr}(w)] = \left[a_w^2 - 1 - a_w j(w) + \frac{1}{4} j(w) j(w) \right]$$

$$[C_{ww}(w)] = \left[1 - \frac{3}{4} j(w) j(w) \right]$$

$$a_r^2 = 1 - \frac{1}{4} r^t r$$

$$a_w^2 = 1 - \frac{1}{4} w^t w$$

$$\text{Thus : } \widehat{X} = \sum_{\alpha=0}^3 V^\alpha X_\alpha$$

$$\widehat{X} = [A_r] \frac{dr}{dt} + [A_w] \frac{dw}{dt} + [D_r] \widehat{G}_r + [D_w] \widehat{G}_w \quad (5.46)$$

So \widehat{X} and the systemic energy of the particle can be expressed as a linear function of the first order derivatives $\frac{dr}{dt}, \frac{dw}{dt}$. The matrices $[A_r], [B_r], [B_w] \dots$ can be expressed as polynomials of $j(r), j(r) j(r), j(w), j(w) j(w),$

$j(r) j(w), j(w) j(r)$, with many nice algebraic properties.

The kinematic energy of the particle is : $k^t \left([A_r] \frac{dr}{dt} + [A_w] \frac{dw}{dt} \right)$.

The electromagnetic field

The electric charge is represented in the inertial spinor. Elementary particles correspond to negative charge, and antiparticles to positive charge. If only the EM field and gravity are present, $\psi = S$ the total covariant derivative reads :

$$\nabla_\alpha \psi = \partial_\alpha \psi + [\gamma C(G_\alpha)] \psi + iq [\dot{A}_\alpha] \psi$$

$$\langle S_0, S_0 \rangle = \epsilon M_p^2 c^4$$

$$k = -\epsilon \frac{1}{2} M_p^2 c^4 k_0$$

and the systemic energy is :

$$\frac{1}{i} \langle \psi, \nabla_V \psi \rangle = k^t \hat{X} + q \hat{A} \langle S_0, S_0 \rangle = -\epsilon \frac{1}{2} M_p^2 c^4 k_0^t \hat{X} + q \epsilon \hat{A} M_p^2 c^4$$

$$\frac{1}{i} \langle S, \nabla_V S \rangle = \epsilon M_p^2 c^4 \left(\frac{1}{2} k_0^t \hat{X} + q \dot{A}_\alpha \right) \quad (5.47)$$

5.3.7 The inertial observer

The states of the particles and the fields are linked, so to measure one we have to know the other : to measure a charge one uses a known field, and to measure a field one uses a known particle. This process requires actually two measures, involving the motion of the particle, it is done locally and is represented by the standard gauges : $\mathbf{p}_G(m) = \varphi_G(m, 1)$, $\mathbf{p}_U(m) = \varphi_U(m, 1)$ and the related holonomic bases $\mathbf{e}_i(m) = (\mathbf{p}_G(m), e_i)$, $\mathbf{f}_j(m) = (\mathbf{p}_U(m), f_j)$. The measures are done with respect to the standards (represented by 1), which are arbitrary. For this reason the gauges and the holonomic bases are not sections, just a specific choice done by the observer. This is consistent with the principle of locality (the measures are done locally) and the free will of the observer (he is not submitted himself to the laws of the system).

However one can consider another kind of gauges, such that they do not change on the travel of an observer on his world line. This seems more physical, and in accordance with the common understanding of inertial observers : they keep their gauges, in which they proceed to their measures, “constant”. However the implementation then must account for the existence of external fields : the observer keeps his free will (which is asserted by the fact that he adjusts his gauge), but he is also submitted to the action of the fields, and notably to the gravitational field.

The action of the fields is given by the covariant derivative of sections, so *the gauges of these observers are sections*, and they change with the fields according to :

$$p_G \in \mathfrak{X}(P_G) : p_G(m) = \varphi_G(m, \sigma(m)) : \nabla_{\varepsilon_0}^G p_G = 0$$

$$p_U \in \mathfrak{X}(P_U) : p_U(m) = \varphi_U(m, g(m)) : \nabla_{\varepsilon_0}^U p_U = 0$$

The world line of the observer is defined by $\varepsilon_0(m)$. The bases of the associated vector bundles $P_G[\mathbb{R}^4, \mathbf{Ad}]$, $P_G[E, \gamma C]$, $P_U[F, \varrho]$ change :

$$\varepsilon_i(m) \rightarrow \tilde{\varepsilon}_i(m) = Ad_{\sigma^{-1}} \varepsilon_i(m)$$

$$e_i(m) \rightarrow \tilde{e}_i(m) = \gamma C(\sigma) e_i(m)$$

$$f_i(m) \rightarrow \tilde{f}_i(m) = \varrho(g) f_i(m)$$

$$\begin{aligned}\nabla_{\varepsilon_0}^G p_G = 0 &\Leftrightarrow \sigma^{-1} \cdot \frac{d\sigma}{dt} + \mathbf{Ad}_{\sigma^{-1}} v \left(\widehat{G}_r, \widehat{G}_w \right) = 0 \Leftrightarrow \frac{d\sigma}{dt} = -v \left(\widehat{G}_r, \widehat{G}_w \right) \cdot \sigma \\ \nabla_{\varepsilon_0}^U p_U = 0 &\Leftrightarrow L'_{g^{-1}} g \left(\frac{dg}{dt} \right) + Ad_{g^{-1}} \widehat{A} = 0 \Leftrightarrow \frac{dg}{dt} = -R'_g 1 \left(\widehat{A} \right)\end{aligned}$$

So the inertial gauges depend on the existing fields along the trajectory. In the absence of any field one retrieves : $\sigma = Ct, g = Ct$ and the standard gauge. This is the usual meaning of inertial observers. This is obvious for p_U : the measure should avoid the interference of external fields, but more subtle for p_G . As for now one cannot escape gravity, inertial observers should follow trajectories such that their motion (through the inertial forces) balance the gravitational field. As it changes slowly with location it will require a constant acceleration, contrary to the common understanding of the inertial observer³.

A great consideration is given, both in Newtonian Mechanics and Relativity, to the inertial observers. The main motivation is that they allow the use of “fixed” frames in affine space, and we have seen that this is the only way to do it in SR. The formalism of fiber bundle seems abstract, but avoids this issue : the concept of standard gauge relies on the fact that the measures of the states of particles or of fields require the comparison with a known quantity (the standard), in similar conditions. And actually, as often in Physics, it is quite impossible to enshrine in the theories the conditions of the experiments. Doing this would bring more confusion than rigor.

The same remark applies to a procedure common in electromagnetism or linearized gravity. When facing a complicated mathematical relation it is tempting to reduce it to a simpler form by calling to what is called gauge freedom. Actually this procedure uses the fact that the same quantity is expressed in different forms according to the gauge (in the fiber bundle definition). So one can replace one by another, which is equivalent, and better looking. To have any meaning this procedure shall follow the requirements of the change of gauge, clearly stated in the fiber bundle formalism. But we have to keep in mind that a change of gauge has a physical meaning, and an implication on the observer who does the measures. A change of gauge can be physically unacceptable by the constraints which would be imposed to the observer, and any experimental proof which would ignore these requirements in its protocols would be non valid.

³An observer in the International Space Station can be considered as inertial, but obviously he is submitted to an acceleration which balances Earth gravity.

5.4 THE PROPAGATION OF FIELDS

We have noticed that Relativity obliges to dissociate the abstract representation of the reality in its entirety, from the representation of the reality which can be scientifically accessible to an observer : this is the true motivation for the use of fiber bundles. It is more necessary when addressing the propagation of fields. Fields are assumed to be defined everywhere, this is one of their key properties. Have the fields a defined value everywhere, including in the future ? What is the true meaning of their propagation, in a 4 dimensional Universe ? We have to remind that the manifold M is just a container, and its physical content is not frozen, it changes. So the propagation of a field is no more than the variation of its value along any future oriented path. But if this image is clear for particles, which have a definite path, it is less so for fields. Are there privileged paths for their propagation ? There is no simple answer to this question. We need to refer to one of our basic assumptions : the representation of the physical world depends on the observer. The observer has a privileged path, given by his velocity and represented by the vector $\varepsilon_0(m)$. This vector plays a fundamental role in the understanding of the propagation of fields. For any observer his own vector $\varepsilon_0(m)$ is the direction of the propagation of the fields. There is no other way to define it, practically, that is in a way which is accessible to measures. So *the propagation of fields is observer dependant*. As a consequence we will see that the field propagates at the speed of light⁴, whatever the observer (theorem 105). But this is not without problems, and one cannot conceive that the propagation of the fields is smooth : it is not instantaneous, and must conciliate the ever changing interactions with particles, and it would be a surprise if it could be done without discontinuities. This will be the topic of one the following chapters, with the introduction of bosons.

For the time being we will focus on the continuous propagation of fields, in the vacuum. The variation of the field is measured by a derivative, and as the fields are characterized by their potential, we need a derivative of the potentials G, \dot{A} . This is the strength of the field, a 2 form \mathcal{F} on M , which is logical because the potentials are themselves one form. However \mathcal{F} as well as the potential is valued in the Lie algebra. If we want to underline the role played by the observer, both in the definition of ε_0 and of holonomic bases in the Lie algebras, it is necessary to put in concordance the mathematical definitions. From this point of view the potential is not the good choice : it is only a map, which changes according to a strange rule in a change of gauge. So the key role in the propagation of the fields will be given to \mathcal{F} .

5.4.1 The strength of the connection

The potentials are one form over M , so their derivative will be a two form. There are several mathematical objects which can be considered, related to the curvature of the connection (Maths.27.1.4), but for principal connections the

⁴In the vacuum. The propagation of light in any medium is a process which involves the interaction of the field with the particles of the medium.

strength of the connection is the most pertinent. This is a 2-form \mathcal{F} on M valued in the Lie algebra (Maths.2194), which can be seen as the exterior covariant derivative of the potential and is a good estimate of its rate of change.

Gravitational field

The strength of the connection is a two form on M valued in the Lie algebra $T_1Spin(3,1)$ which reads with the basis $(\vec{\kappa}_a)_{a=1}^6$:

$$\mathcal{F}_G = \sum_{a=1}^6 \left(dG^a + \sum_{\alpha\beta=0}^3 [G_\alpha, G_\beta]^a d\xi^\alpha \wedge d\xi^\beta \right) \otimes \vec{\kappa}_a$$

$$\mathcal{F}_G = \sum_{a=1}^6 \sum_{\alpha,\beta=0}^3 \mathcal{F}_{G_{\alpha\beta}}^a d\xi^\alpha \wedge d\xi^\beta \otimes \vec{\kappa}_a \quad (5.48)$$

$$= \sum_{a=1}^6 \sum_{\alpha,\beta=0}^3 (\partial_\alpha G_\beta^a - \partial_\beta G_\alpha^a + [G_\alpha, G_\beta]^a) d\xi^\alpha \wedge d\xi^\beta \otimes \vec{\kappa}_a \quad (5.49)$$

where d is the exterior differential on TM and $[\]$ is the bracket in $T_1Spin(3,1)$.⁵

Or equivalently :

$$\mathcal{F}_G = d \left(\sum_{\alpha=0}^3 v(G_{r\alpha}, G_{w\alpha}) d\xi^\alpha \right) + \sum_{\alpha\beta=0}^3 [v(G_{r\alpha}, G_{w\alpha}), v(G_{r\beta}, G_{w\beta})] d\xi^\alpha \wedge d\xi^\beta$$

We can distinguish the two parts, $\mathcal{F}_r, \mathcal{F}_w$:

$$\mathcal{F}_G = \sum_{\alpha,\beta} v(\mathcal{F}_{r\alpha\beta}, \mathcal{F}_{w\alpha\beta}) d\xi^\alpha \wedge d\xi^\beta \quad (5.50)$$

and we have :

$$a = 1,2,3 : \mathcal{F}_{G_{\alpha\beta}}^a = \mathcal{F}_{r\alpha\beta}^a$$

$$a = 4,5,6 : \mathcal{F}_{G_{\alpha\beta}}^a = \mathcal{F}_{w\alpha\beta}^a$$

with the signature (3,1) :

$$[v(r, w), v(r', w')] = v(j(r) r' - j(w) w', j(w) r' + j(r) w')$$

$$[G_\alpha, G_\beta] = [v(G_{r\alpha}, G_{w\alpha}), v(G_{r\beta}, G_{w\beta})]$$

$$= v(j(G_{r\alpha}) G_{r\beta} - j(G_{w\alpha}) G_{w\beta}, j(G_{w\alpha}) G_{r\beta} + j(G_{r\alpha}) G_{w\beta})$$

$$\mathcal{F}_{r\alpha\beta} = v \left(\frac{\partial G_{r\beta}}{\partial \xi^\alpha} - \frac{\partial G_{r\alpha}}{\partial \xi^\beta} + j(G_{r\alpha}) G_{r\beta} - j(G_{w\alpha}) G_{w\beta}, 0 \right) \quad (5.51)$$

$$\mathcal{F}_{w\alpha\beta} = v \left(0, \frac{\partial G_{w\beta}}{\partial \xi^\alpha} - \frac{\partial G_{w\alpha}}{\partial \xi^\beta} + j(G_{w\alpha}) G_{r\beta} + j(G_{r\alpha}) G_{w\beta} \right) \quad (5.52)$$

With the signature (1,3):

$$[v(r, w), v(r', w')] = -v(j(r) r' - j(w) w', j(w) r' + j(r) w')$$

⁵The notations and conventions for r forms vary according to the authors and if the indices are ordered or not. On this see Maths.1525,1529.

$$\begin{aligned}
& [G_\alpha, G_\beta] \\
& = - (v (j (G_{r\alpha}) G_{r\beta} - j (G_{w\alpha}) G_{w\beta}), j (G_{w\alpha}) G_{r\beta} + j (G_{r\alpha}) G_{w\beta})) \\
\mathcal{F}_{r\alpha\beta} & = -v \left(\frac{\partial G_{r\beta}}{\partial \xi^\alpha} - \frac{\partial G_{r\alpha}}{\partial \xi^\beta} + j (G_{r\alpha}) G_{r\beta} - j (G_{w\alpha}) G_{w\beta}, 0 \right) \quad (5.53)
\end{aligned}$$

$$\mathcal{F}_{w\alpha\beta} = -v \left(0, \frac{\partial G_{w\beta}}{\partial \xi^\alpha} - \frac{\partial G_{w\alpha}}{\partial \xi^\beta} + j (G_{w\alpha}) G_{r\beta} + j (G_{r\alpha}) G_{w\beta} \right) \quad (5.54)$$

Notice that the indices α, β are not ordered, that it involves only the principal bundle, and not the associated vector bundles, and is valued in a fixed vector space.

\mathcal{F}_G can be written in the matrix form of $so(3, 1)$, using $\vec{\kappa}_a \rightarrow [\kappa_a]$:

$$[\mathcal{F}_{G\alpha\beta}] = \sum_{a=1}^6 \mathcal{F}_{G\alpha\beta}^a [\kappa_a] = [K (\mathcal{F}_{w\alpha\beta})] + [J (\mathcal{F}_{r\alpha\beta})] \quad (5.55)$$

$$[\mathcal{F}_{\alpha\beta}] = \begin{bmatrix} 0 & \mathcal{F}_{w\alpha\beta}^1 & \mathcal{F}_{w\alpha\beta}^2 & \mathcal{F}_{w\alpha\beta}^3 \\ \mathcal{F}_{w\alpha\beta}^1 & 0 & -\mathcal{F}_{r\alpha\beta}^3 & \mathcal{F}_{r\alpha\beta}^2 \\ \mathcal{F}_{w\alpha\beta}^2 & \mathcal{F}_{r\alpha\beta}^3 & 0 & -\mathcal{F}_{\alpha\beta}^1 \\ \mathcal{F}_{w\alpha\beta}^3 & -\mathcal{F}_{r\alpha\beta}^2 & \mathcal{F}_{r\alpha\beta}^1 & 0 \end{bmatrix} = \begin{bmatrix} 0 & \mathcal{F}_{w\alpha\beta} \\ \mathcal{F}_{w\alpha\beta} & j (\mathcal{F}_{r\alpha\beta}) \end{bmatrix}$$

which underlines the rotational feature of the component \mathcal{F}_r , and the transversal aspect of the component \mathcal{F}_w .

With :

$$\begin{aligned}
\Gamma_{M\alpha} & = [K (G_{w\alpha})] + [J (G_{r\alpha})] = \begin{bmatrix} 0 & G_{w\alpha}^1 & G_{w\alpha}^2 & G_{w\alpha}^3 \\ G_{w\alpha}^1 & 0 & -G_{r\alpha}^3 & G_{r\alpha}^2 \\ G_{w\alpha}^2 & G_{r\alpha}^3 & 0 & -G_{\alpha}^1 \\ G_{w\alpha}^3 & -G_{r\alpha}^2 & G_{r\alpha}^1 & 0 \end{bmatrix} \\
& = \begin{bmatrix} 0 & G_{w\alpha} \\ G_{w\alpha} & j (G_{r\alpha}) \end{bmatrix}
\end{aligned}$$

$$[\mathcal{F}_{\alpha\beta}] = [\partial_\alpha \Gamma_{M\beta}] - [\partial_\beta \Gamma_{M\alpha}] + [\Gamma_{M\alpha}] [\Gamma_{M\beta}] - [\Gamma_{M\beta}] [\Gamma_{M\alpha}] \quad (5.56)$$

Other fields

The strength \mathcal{F}_A of the connection is a 2 form on M valued in the Lie algebra T_1U , which reads, with non ordered indices α, β :

$$\mathcal{F}_A = d \left(\sum_{\alpha=0}^3 \dot{A}_\alpha d\xi^\alpha \right) + \sum_{\alpha\beta} [\dot{A}_\alpha, \dot{A}_\beta] d\xi^\alpha \wedge d\xi^\beta$$

where d is the exterior differential on TM and $[\]$ is the bracket in T_1U

$$\mathcal{F}_A = \sum_{\alpha, \beta} (\mathcal{F}_{A\alpha\beta}^a d\xi^\alpha \wedge d\xi^\beta) \otimes \vec{\theta}_a \in \Lambda_2 (M; T_1U) \quad (5.57)$$

and in components :

$$\mathcal{F}_{A\alpha\beta}^a = \frac{\partial \dot{A}_\beta^a}{\partial \xi^\alpha} - \frac{\partial \dot{A}_\alpha^a}{\partial \xi^\beta} + [\dot{A}_\alpha, \dot{A}_\beta]^a \quad (5.58)$$

Notice that the indices α, β are not ordered, that it involves only the principal bundle, and not the associated vector bundles, and is valued in a fixed vector space.

Adjoint bundle

The strength of the connection is a map valued in the Lie algebra, that is a fixed vector space. We have seen above how the potential changes in a change of gauge. In a change of gauge on the principal bundle the strength changes as :

$$\mathbf{p}_U(m) = \varphi_{P_U}(m, 1) \rightarrow \tilde{\mathbf{p}}_U(m) = \mathbf{p}_U(m) \cdot \varkappa(m)^{-1} : \quad (5.59)$$

$$\mathcal{F}_{A\alpha\beta} \rightarrow \tilde{\mathcal{F}}_{A\alpha\beta}(m) = Ad_{\varkappa(m)} \mathcal{F}_{A\alpha\beta} \quad (5.60)$$

and we have similarly :

$$\mathbf{p}_G(m) = \varphi_G(m, 1) \rightarrow \tilde{\mathbf{p}}_G(m) = \mathbf{p}_G(m) \cdot s(m)^{-1} : \quad (5.61)$$

$$\mathcal{F}_{G\alpha\beta} \rightarrow \tilde{\mathcal{F}}_{G\alpha\beta}(m) = \mathbf{Ad}_{s(m)} \mathcal{F}_{G\alpha\beta} \quad (5.62)$$

$$v \left(\tilde{\mathcal{F}}_{r\alpha\beta}, \tilde{\mathcal{F}}_{w\alpha\beta} \right) = \mathbf{Ad}_{s(m)} v \left(\mathcal{F}_{r\alpha\beta}, \mathcal{F}_{w\alpha\beta} \right)$$

This feature allows to consider the strength as sections of the **adjoint bundles**, which are defined as the associated vector bundles

$P_G [T_1 Spin(3, 1), \mathbf{Ad}]$, $P_U [T_1 U, Ad]$, using the representation of the groups on their Lie algebra through the adjoint map (Ad on $T_1 Spin(3, 1)$ is identical to \mathbf{Ad}). This gives a more geometrical meaning to the concept, and we will see that these relations are crucial in the definition of the lagrangian.

Electromagnetic field

The strength of the electromagnetic field is a 2 form valued in \mathbb{R} : $\mathcal{F}_A \in \Lambda_2(M; \mathbb{R})$.

Because the Lie algebra is abelian the bracket is null and : $\mathcal{F}_A = d\dot{A}$ which gives the first Maxwell's law : $d\mathcal{F}_A = 0$.

In a change of gauge : $\mathcal{F}_{A\alpha\beta} \rightarrow \tilde{\mathcal{F}}_{A\alpha\beta}(m) = Ad_{\varkappa(m)} \mathcal{F}_{A\alpha\beta} = \mathcal{F}_{A\alpha\beta}$. The strength of the EM field is invariant in a change of gauge.

5.4.2 Scalar curvature

The strength of the potential is a general object, which is related to the different concepts of curvature used in the theory of connections. It involves only the principal bundle, and not the associated vector bundles. However in GR another definition of curvature is commonly used, and it is necessary to see how these concepts are related.

Riemann curvature of a principal connection

A connection on a principal bundle leads to the definition of a quantity, called **Riemann curvature**, on any associated vector bundle, through a path which involves exterior covariant derivatives (Maths.2203). The result is a two-form on M, valued in the endomorphisms on the vector space, which in the case of $P_G [\mathbb{R}^4, \mathbf{Ad}]$ is expressed by the 4 order tensor :

$$R = \sum_{\{\alpha\beta\}ij} R_{\alpha\beta j}^i d\xi^\alpha \wedge d\xi^\beta \otimes \varepsilon_i(m) \otimes \varepsilon^j(m) \text{ with the dual basis } \varepsilon^j(m)$$

$$R_{\alpha\beta j}^i = \partial_\alpha \Gamma_{M\beta j}^i - \partial_\beta \Gamma_{M\alpha j}^i + \sum_{k=0}^3 \left(\Gamma_{M\alpha k}^i \Gamma_{M\beta j}^k - \Gamma_{M\beta k}^i \Gamma_{M\alpha j}^k \right)$$

In matrix form (see above) :

$$[R_{\alpha\beta}]_j^i = ([\partial_\alpha \Gamma_{M\beta}] - [\partial_\beta \Gamma_{M\alpha}] + [\Gamma_{M\alpha}] [\Gamma_{M\beta}] - [\Gamma_{M\beta}] [\Gamma_{M\alpha}])_j^i = [\mathcal{F}_{\alpha\beta}]_j^i$$

$$R = \sum_{\{\alpha\beta\}ij} [\mathcal{F}_{\alpha\beta}]_j^i d\xi^\alpha \wedge d\xi^\beta \otimes \varepsilon_i(m) \otimes \varepsilon^j(m)$$

The Riemann curvature is the image of the strength of the field on $P_G [\mathbb{R}^4, \mathbf{Ad}]$. This is the same quantity, but in the representation of $T_1 Spin(3, 1)$ in the matrix algebra $so(3, 1)$.

By construct this quantity is covariant (in a change of chart on M) and equivariant (in a change of gauge on P_G) :

In a change of gauge :

$$\mathbf{p}_G(m) = \varphi_G(m, 1) \rightarrow \tilde{\mathbf{p}}_G(m) = \mathbf{p}_G(m) \cdot s(m)^{-1} :$$

$$[\mathcal{F}_{\alpha\beta}]_j^i \rightarrow \sum_{kl=0}^3 [h(s)]_k^i [\mathcal{F}_{\alpha\beta}]_l^k [h(s^{-1})]_j^l$$

with $[h(s)]$ the matrix of $SO(3, 1)$ associated to $s \in Spin(3, 1)$

$$R \rightarrow \tilde{R} = \sum_{\{\alpha\beta\}ij} [\tilde{\mathcal{F}}_{\alpha\beta}]_j^i d\xi^\alpha \wedge d\xi^\beta \otimes \tilde{\varepsilon}_i(m) \otimes \tilde{\varepsilon}^j(m)$$

$$= \sum_{\{\alpha\beta\}ij} [h(s)]_k^i [\mathcal{F}_{\alpha\beta}]_l^k [h(s^{-1})]_j^l d\xi^\alpha \wedge d\xi^\beta \otimes [h(s^{-1})]_i^p \varepsilon_p(m) \otimes [h(s)]_q^j \varepsilon^q(m)$$

$$= \sum_{\{\alpha\beta\}ij} [\mathcal{F}_{\alpha\beta}]_q^p d\xi^\alpha \wedge d\xi^\beta \otimes \varepsilon_p(m) \otimes \varepsilon^q(m)$$

so $\tilde{R} = R$

It can be expressed in the holonomic basis of any chart on M using the tetrad

:

$$\varepsilon_i(m) = \sum_{\gamma=0}^3 P_i^\gamma \partial \xi_\gamma$$

$$\varepsilon^j(m) = \sum_{\eta=0}^3 P_\eta^j d\xi^\eta$$

$$R = \sum_{\{\alpha\beta\}\gamma\eta} ([P] [\mathcal{F}_{G\alpha\beta}] [P'])_\eta^\gamma d\xi^\alpha \wedge d\xi^\beta \otimes \partial \xi_\gamma \otimes d\xi^\eta$$

So we have the steps :

Principal connection $\mathbf{G} \rightarrow$ Riemann curvature R on $P_G [\mathbb{R}^4, \mathbf{Ad}] \rightarrow$ Riemann curvature R on TM in any chart

and the Riemann curvature R on TM is the same object as the strength of the connection \mathcal{F} , but expressed in matrix form in any holonomic basis of a chart.

Riemann tensor of an affine connection

With a common affine connection $\hat{\Gamma}_\alpha$ on TM one can also define similarly a **Riemann tensor** (Maths.1543) :

$$\hat{R} = \sum_{\{\alpha\beta\}} \sum_{\gamma\eta} \hat{R}_{\alpha\beta\eta}^\gamma d\xi^\alpha \wedge d\xi^\beta \otimes \partial \xi_\gamma \otimes d\xi^\eta$$

which, expressed in matrix form with : $\widehat{R}_{\alpha\beta\eta}^\gamma = \left[\widehat{R}_{\alpha\beta} \right]_\eta^\gamma$, reads :

$$\left[\widehat{R}_{\alpha\beta} \right] = \left[\partial_\alpha \widehat{\Gamma}_\beta \right] - \left[\partial_\beta \widehat{\Gamma}_\alpha \right] + \left[\widehat{\Gamma}_\alpha \right] \left[\widehat{\Gamma}_\beta \right] - \left[\widehat{\Gamma}_\beta \right] \left[\widehat{\Gamma}_\alpha \right]$$

$$\widehat{R} = \sum_{\{\alpha\beta\}} \sum_{\gamma\eta} \widehat{R}_{\alpha\beta\eta}^\gamma d\xi^\alpha \wedge d\xi^\beta \otimes \partial \xi_\gamma \otimes d\xi^\eta$$

When we take as affine connection the one which is deduced from \mathbf{G} :

$$\widehat{\Gamma}_{\alpha\beta}^\gamma = \left[\widehat{\Gamma}_\alpha \right]_\beta^\gamma = ([P] ([\partial_\alpha P'] + [\Gamma_{M\alpha}] [P']))_\beta^\gamma$$

we get the same result :

$$\left[\widehat{R}_{\alpha\beta} \right] = [R_{\alpha\beta}] = [P] [\mathcal{F}_{G\alpha\beta}] [P'] \Leftrightarrow [\mathcal{F}_{G\alpha\beta}] = [P'] [R_{\alpha\beta}] [P] \quad (5.63)$$

Proof. $\left[\widehat{R}_{\alpha\beta} \right]$

$$\begin{aligned} &= [\partial_\alpha P] [\partial_\beta P'] + [\partial_\alpha P] [\Gamma_{M\beta}] [P'] + [P] \left[\partial_{\beta\alpha}^2 P' \right] + [P] [\partial_\alpha \Gamma_{M\beta}] [P'] \\ &+ [P] [\Gamma_{M\beta}] [\partial_\alpha P'] - [\partial_\beta P] [\partial_\alpha P'] - [\partial_\beta P] [\Gamma_{M\alpha}] [P'] - [P] \left[\partial_{\alpha\beta}^2 P' \right] \\ &- [P] [\partial_\beta \Gamma_{M\alpha}] [P'] - [P] [\Gamma_{M\alpha}] [\partial_\beta P'] + [P] [\partial_\alpha P'] [P] [\partial_\beta P'] \\ &+ [P] [\Gamma_{M\alpha}] [P'] [P] [\partial_\beta P'] + [P] [\partial_\alpha P'] [P] [\Gamma_{M\beta}] [P'] \\ &+ [P] [\Gamma_{M\alpha}] [P'] [P] [\Gamma_{M\beta}] [P'] - [P] [\partial_\beta P'] [P] [\partial_\alpha P'] \\ &- [P] [\Gamma_{M\beta}] [P'] [P] [\partial_\alpha P'] - [P] [\partial_\beta P'] [P] [\Gamma_{M\alpha}] [P'] \\ &- [P] [\Gamma_{M\beta}] [P'] [P] [\Gamma_{M\alpha}] [P'] \\ &= + [P] ([\partial_\alpha \Gamma_{M\beta}] - [\partial_\beta \Gamma_{M\alpha}] + [\Gamma_{G\alpha}] [\Gamma_{M\beta}] - [\Gamma_{M\beta}] [\Gamma_{M\alpha}]) [P'] \\ &+ [\partial_\alpha P] [\partial_\beta P'] - [\partial_\beta P] [\partial_\alpha P'] + [P] [\partial_\alpha P'] [P] [\partial_\beta P'] \\ &- [P] [\partial_\beta P'] [P] [\partial_\alpha P'] + [\partial_\alpha P] [\Gamma_{M\beta}] [P'] - [\partial_\beta P] [\Gamma_{M\alpha}] [P'] \\ &+ [P] [\Gamma_{M\beta}] [\partial_\alpha P'] - [P] [\Gamma_{M\alpha}] [\partial_\beta P'] + [P] [\Gamma_{M\alpha}] [\partial_\beta P'] \\ &- [P] [\Gamma_{M\beta}] [\partial_\alpha P'] + [P] [\partial_\alpha P'] [P] [\Gamma_{M\beta}] [P'] - [P] [\partial_\beta P'] [P] [\Gamma_{M\alpha}] [P'] \\ &= [P] [\mathcal{F}_{G\alpha\beta}] [P'] + [\partial_\alpha P] [\partial_\beta P'] - [\partial_\beta P] [\partial_\alpha P'] \\ &- [\partial_\alpha P] [P'] [P] [\partial_\beta P'] + [\partial_\beta P] [P'] [P] [\partial_\alpha P'] + [\partial_\alpha P] [\Gamma_{M\beta}] [P'] \\ &- [\partial_\beta P] [\Gamma_{M\alpha}] [P'] + [P] [\Gamma_{M\beta}] [\partial_\alpha P'] - [P] [\Gamma_{M\alpha}] [\partial_\beta P'] + [P] [\Gamma_{M\alpha}] [\partial_\beta P'] \\ &- [P] [\Gamma_{M\beta}] [\partial_\alpha P'] - [\partial_\alpha P] [P'] [P] [\Gamma_{M\beta}] [P'] + [\partial_\beta P] [P'] [P] [\Gamma_{M\alpha}] [P'] \\ &= [P] [\mathcal{F}_{G\alpha\beta}] [P'] \\ &\text{with } [P] [\partial_\alpha P'] + [\partial_\alpha P] [P'] = 0 \quad \blacksquare \end{aligned}$$

So the Riemann tensor is the Riemann curvature of the principal connection, expressed in the holonomic basis of a chart, and it is the same object as the strength of the connection.

The Riemann tensor can be computed with any affine connection, as well as with any principal connection. In the usual RG formalism the Riemann tensor is computed with a special connection : the Levy-Civita connection.

The Riemann tensor is antisymmetric, in the meaning :

$$\begin{aligned} \widehat{R}_{\alpha\beta\gamma\eta} &= -\widehat{R}_{\alpha\beta\eta\gamma} \text{ with } \widehat{R}_{\alpha\beta\gamma\eta} = \sum_\lambda \widehat{R}_{\alpha\beta\gamma}^\lambda g_{\lambda\eta} \\ [\mathcal{F}_{G\alpha\beta}] &\in so(3,1) \text{ so } [\eta] [\mathcal{F}_{G\alpha\beta}] + [\mathcal{F}_{G\alpha\beta}]^t [\eta] = 0 \text{ and} \\ \widehat{R}_{\alpha\beta\gamma\eta} &= \sum_\lambda \widehat{R}_{\alpha\beta\gamma}^\lambda g_{\lambda\eta} = \sum_\lambda ([P] [\mathcal{F}_{G\alpha\beta}] [P'])_\gamma^\lambda g_{\lambda\eta} = \left([P']^t [\eta] [\mathcal{F}_{G\alpha\beta}] [P'] \right)_\gamma^\eta \\ &= \left(\left([P']^t [\eta] [\mathcal{F}_{G\alpha\beta}] [P'] \right)^t \right)_\eta^\gamma = \left([P']^t [\mathcal{F}_{G\alpha\beta}]^t [\eta] [P'] \right)_\eta^\gamma \end{aligned}$$

$$= - \left([P']^t [\eta] [\mathcal{F}_{G\alpha\beta}] [P'] \right)_\eta^\gamma = - \widehat{R}_{\alpha\beta\eta\gamma}$$

Thus this symmetry is not specific to the Lévi-Civita connection as it is usually assumed (Wald p.39).

Ricci tensor and scalar curvature

The Riemann tensor \widehat{R} , coming from any connection, is a 2 form but can be expressed as an antisymmetric tensor with non ordered indices :

$$\widehat{R} = \sum_{\alpha\beta\gamma\eta} \left[\widehat{R}_{\alpha\beta} \right]_\eta^\gamma d\xi^\alpha \otimes d\xi^\beta \otimes \partial\xi_\gamma \otimes d\xi^\eta$$

and we can contract the covariant index α, β or η with the contravariant index γ . The result does not depend on a basis : it is covariant (Maths.385).

The different solutions give :

$$\begin{aligned} \alpha &: \sum_{\beta\eta} \left(\sum_\alpha \left[\widehat{R}_{\alpha\beta} \right]_\eta^\alpha \right) d\xi^\beta \otimes d\xi^\eta \\ \beta &: \sum_{\alpha\eta} \left(\sum_\beta \left[\widehat{R}_{\alpha\beta} \right]_\eta^\beta \right) d\xi^\alpha \otimes d\xi^\eta \\ \eta &: \sum_{\alpha\eta} \left(\sum_\gamma \left[\widehat{R}_{\alpha\beta} \right]_\gamma^\eta \right) d\xi^\alpha \otimes d\xi^\beta \end{aligned}$$

The last solution has no interest because :

$$Tr ([P] [\mathcal{F}_{G\alpha\beta}] [P']) = Tr ([\mathcal{F}_{G\alpha\beta}] [P'] [P]) = Tr ([\mathcal{F}_{G\alpha\beta}]) = 0$$

The first two read :

$$\begin{aligned} \sum_{\beta\gamma} [P]_k^\alpha [\mathcal{F}_{G\alpha\beta}]_l^k [P']_\eta^l [P]_i^\beta \varepsilon^i \otimes [P]_j^\eta \varepsilon^j &= \sum_{\beta\gamma} [P]_k^\alpha [\mathcal{F}_{G\alpha\beta}]_j^k [P]_i^\beta \varepsilon^i \otimes \varepsilon^j = \\ \sum_{\alpha\beta j} ([P] [\mathcal{F}_{G\alpha\beta}])_j^\alpha d\xi^\beta \otimes \varepsilon^j & \\ \sum_{\alpha\gamma} [P]_k^\beta [\mathcal{F}_{G\alpha\beta}]_l^k [P']_\eta^l [P]_i^\alpha \varepsilon^i \otimes [P]_j^\eta \varepsilon^j &= \sum_{\alpha\gamma} [P]_k^\beta [\mathcal{F}_{G\alpha\beta}]_j^k [P]_i^\alpha \varepsilon^i \otimes \varepsilon^j = \\ \sum_{\beta\gamma} ([P] [\mathcal{F}_{G\alpha\beta}])_j^\beta d\xi^\alpha \otimes \varepsilon^j & \end{aligned}$$

The **Ricci tensor** is the contraction on the two indices γ, β of \widehat{R} :

$$Ric = \sum_{\alpha\eta} Ric_{\alpha\eta} d\xi^\alpha \otimes d\xi^\eta = \sum_{\alpha\eta} \left(\sum_\beta \left[\widehat{R}_{\alpha\beta} \right]_\eta^\beta \right) d\xi^\alpha \otimes d\xi^\eta$$

This is a tensor, from which one can compute another tensor by lowering the last index:

$$\sum_\lambda g^{\eta\lambda} Ric_{\alpha\eta} d\xi^\alpha \otimes d\xi^\eta = \sum_{\alpha\lambda} Ric_\alpha^\lambda d\xi^\alpha \otimes \partial\xi_\lambda$$

and the contraction (called the trace of this tensor) provides the **scalar curvature** :

$$\mathbf{R} = \sum_\alpha Ric_\alpha^\alpha = \sum_{\alpha\beta\eta} g^{\alpha\eta} \left[\widehat{R}_{\alpha\beta} \right]_\eta^\beta$$

The same procedure applied to the contraction on the two indices γ, α of \widehat{R} gives the opposite scalar :

$$\mathbf{R} = \sum_{\alpha\beta\eta} g^{\beta\eta} \left[\widehat{R}_{\alpha\beta} \right]_\eta^\alpha = - \sum_{\alpha\beta\eta} g^{\alpha\eta} \left[\widehat{R}_{\beta\alpha} \right]_\eta^\beta = - \sum_{\alpha\beta\eta} g^{\alpha\eta} \left[\widehat{R}_{\alpha\beta} \right]_\eta^\beta$$

This manipulation is mathematically valid, and provides a unique scalar, which does not depend on a chart, and can be used in a lagrangian. However its physical justification (see Wald) is weak.

In the usual GR formalism the scalar curvature is computed with the Riemann tensor deduced from the Levy-Civita connection but, as we can see, it can be computed with any principal connection. Starting from $[P] [\mathcal{F}_{G\alpha\beta}] [P'] = [\widehat{\mathcal{F}}_{G\alpha\beta}]$ one gets :

$$\begin{aligned} \mathbf{R} &= \sum_{\alpha\beta\eta} g^{\alpha\eta} ([P] [\mathcal{F}_{G\alpha\beta}] [P'])_{\eta}^{\beta} \text{ and with } [g]^{-1} = [P] [\eta] [P]^t \\ \mathbf{R} &= \sum_{\alpha\beta\eta} ([P] [\mathcal{F}_{G\alpha\beta}] [P'])_{\eta}^{\beta} [g^{-1}]_{\alpha}^{\eta} = \sum_{\alpha\beta} ([P] [\mathcal{F}_{G\alpha\beta}] [P'] [P] [\eta] [P]^t)_{\alpha}^{\beta} \\ \mathbf{R} &= \sum_{\alpha\beta} ([P] [\mathcal{F}_{G\alpha\beta}] [\eta] [P]^t)_{\alpha}^{\beta} \end{aligned} \quad (5.64)$$

The computation of the scalar curvature gives ::

$$\begin{aligned} \mathbf{R} &= \sum_{a=1}^6 \sum_{\alpha\beta} \mathcal{F}_{G\alpha\beta}^a ([P] [\kappa_a] [\eta] [P]^t)_{\alpha}^{\beta} \\ \text{For } a = 1,2,3 : [\kappa_a] [\eta] &= [\kappa_a] \\ [\kappa_a]_q^p &= \sum_{bc=1}^3 \epsilon(a,b,c) \delta_c^p \delta_q^b \\ ([P] [\kappa_a] [\eta] [P]^t)_{\alpha}^{\beta} &= \sum_{p,q=1}^3 \epsilon(a,q,p) [P]_p^{\beta} [P]_q^{\alpha} \\ \sum_{a=1}^3 \sum_{\alpha\beta} \mathcal{F}_{G\alpha\beta}^a ([P] [\kappa_a] [\eta] [P]^t)_{\alpha}^{\beta} &= \sum_{a=1}^3 \sum_{\alpha\beta} \mathcal{F}_{r\alpha\beta}^a \sum_{p,q=1}^3 \epsilon(a,q,p) [P]_p^{\beta} [P]_q^{\alpha} \\ \text{For } a = 4,5,6 : \\ ([\kappa_a] [\eta])_q^p &= \delta_0^p \delta_q^{a-3} - \delta_{a-3}^p \delta_q^0 \\ ([P] [\kappa_a] [\eta] [P]^t)_{\alpha}^{\beta} &= [P]_0^{\beta} [P]_{a-3}^{\alpha} - [P]_{a-3}^{\beta} [P]_0^{\alpha} \\ \sum_{a=4}^6 \sum_{\alpha\beta} \mathcal{F}_{G\alpha\beta}^a ([P] [\kappa_a] [\eta] [P]^t)_{\alpha}^{\beta} &= \sum_{a=1}^3 \sum_{\alpha\beta} \mathcal{F}_{w\alpha\beta}^a ([P]_0^{\beta} [P]_a^{\alpha} - [P]_a^{\beta} [P]_0^{\alpha}) \end{aligned}$$

Thus :

$$\mathbf{R} = \sum_{a=1}^3 \sum_{\alpha\beta} \left(\mathcal{F}_{w\alpha\beta}^a ([P]_0^{\beta} [P]_a^{\alpha} - [P]_a^{\beta} [P]_0^{\alpha}) + \mathcal{F}_{r\alpha\beta}^a \sum_{p,q=1}^3 \epsilon(a,q,p) [P]_p^{\beta} [P]_q^{\alpha} \right) \quad (5.65)$$

which reads in a geometric way :

$$\begin{aligned} \mathbf{R} &= \sum_{a=1}^3 \left(\mathcal{F}_w^a (\varepsilon_a, \varepsilon_0) + \sum_{p,q=1}^3 \epsilon(a,q,p) \mathcal{F}_r^a (\varepsilon_q, \varepsilon_p) \right) \\ \text{In the standard chart with : } [P]_0^{\alpha} &= \delta_0^{\alpha} \\ \mathbf{R} &= \sum_{a=1}^3 \sum_{\alpha\beta} \left(\mathcal{F}_{w\alpha\beta}^a (\delta_0^{\beta} [P]_a^{\alpha} - [P]_a^{\beta} \delta_0^{\alpha}) + \mathcal{F}_{r\alpha\beta}^a \sum_{p,q=1}^3 \epsilon(a,q,p) [P]_p^{\beta} [P]_q^{\alpha} \right) \\ &= 2 \sum_{\alpha < \beta} \left\{ \sum_{a=1}^3 \mathcal{F}_{G\alpha\beta}^a \sum_{p,q=1}^3 \epsilon(a,q,p) [P]_p^{\beta} [P]_q^{\alpha} \right. \\ &\quad \left. + \sum_{a=4}^6 \mathcal{F}_{G\alpha\beta}^a ([P]_0^{\beta} [P]_{a-3}^{\alpha} - [P]_{a-3}^{\beta} [P]_0^{\alpha}) \right\} \end{aligned}$$

This expression has two important features :

- the scalar curvature is linear with respect to the strength of the field.
In the implementation of the Principle of Least Action it provides equations which are linear with respect to \mathcal{F}_G , which is a big improvement from the usual computations.

- it shows that *the scalar curvature has a transversal component and a rotational component*. This happens for any scalar curvature, but is just masked in the usual expression through the metric. This feature is not without significance, as it is related to the distinction between the space and the time Universe.

To sum up with the fiber bundle and connections formalism it is possible to compute, more easily, a scalar curvature which has the usual meaning. And by imposing symmetry to the affine connection we get exactly the same quantity. However, as we have seen before, the symmetry of the connection has no obvious physical meaning, and similarly for the scalar curvature. So, in the following, we will stay with the strength of the connection, which gives a good representation of the propagation of the field, and puts gravitation in the same footing as the other fields.

5.4.3 Polarization of the fields

Polarization of light is a concept familiar to Physicists, and can be easily observed. It has a deeper meaning, related to the fact that the vector $\partial\xi_0$ is the only privileged direction for the propagation of the field with respect to the observer.

With singling out $d\xi^0$, any two form read :

$$\mathcal{F} = \mathcal{F}_R + \mathcal{F}_W$$

with :

$$\mathcal{F}_R = 2 (\mathcal{F}_{32}d\xi^3 \wedge d\xi^2 + \mathcal{F}_{13}d\xi^1 \wedge d\xi^3 + \mathcal{F}_{21}d\xi^2 \wedge d\xi^1)$$

$$\mathcal{F}_W = 2 (\mathcal{F}_{01}d\xi^0 \wedge d\xi^1 + \mathcal{F}_{02}d\xi^0 \wedge d\xi^2 + \mathcal{F}_{03}d\xi^0 \wedge d\xi^3)$$

(the 2 accounting for the symmetric part)

The action of these two forms on vectors is :

$$\mathcal{F}_W \left(\sum_{\alpha=0}^3 u^\alpha \partial\xi_\alpha \right) = -2u^0 (\mathcal{F}_{01}d\xi^1 + \mathcal{F}_{02}d\xi^2 + \mathcal{F}_{03}d\xi^3)$$

$$+2 (\mathcal{F}_{01}u^1 + \mathcal{F}_{02}u^2 + \mathcal{F}_{03}u^3) d\xi^0$$

$$\mathcal{F}_R (u^0 \partial\xi_0) = 0$$

$$\mathcal{F}_R \left(\sum_{\alpha=0}^3 u^\alpha \partial\xi_\alpha \right) = - \sum_{\alpha=1}^3 [j (\mathcal{F}_{32}d\xi^1 + \mathcal{F}_{13}d\xi^2 + \mathcal{F}_{21}d\xi^3) u]_\alpha d\xi^\alpha$$

so the action of \mathcal{F}_R^a can be seen as a rotation in the physical space by the vector \mathcal{F}_R .

This decomposition seems a bit formal, however for the EM field this is exactly the decomposition in electric field (\mathcal{F}_W) and magnetic field (\mathcal{F}_R). It characterizes the **polarization of the field** in the direction given by the vector $\varepsilon_0 = \partial\xi_0$ which characterizes the observer. And we see that it can be extended to any field, separately for each component \mathcal{F}^a . The decomposition into the components $\mathcal{F}_R, \mathcal{F}_W$ depends on the chart. For instance the decomposition of the EM field in electric and magnetic field depends on the observer. But we can go further in exploring this decomposition.

5.4.4 The Relativist Momentum of Fields

\mathcal{F} is formally defined as a 2 form, that is with respect to any chart, in one hand, and is valued in the adjoint bundle, that is with respect to a gauge provided by P_U or P_G on the other hand. The privileged direction $\partial\xi_0 = \varepsilon_0$ is masked in the anonymity of the labels in a banalized chart. But there is no such banal chart in Relativity : a chart is always linked to an observer, and each observer has a privileged orientation, that of its future. To put the matter right it is necessary to express \mathcal{F} in a basis which varies with the observer in a consistent and clear way. It is possible to do this, first by expressing the two forms in the orthonormal basis, then by expressing it in a unique vector bundle.

From the holonomic basis of a chart to the orthonormal basis

Any scalar two form on M can be expressed in the dual basis $(\varepsilon^i(m))_{i=0}^3$:

$$\mathcal{F} = \sum_{\alpha\beta=0}^3 \mathcal{F}_{\alpha\beta} d\xi^\alpha \wedge d\xi^\beta = \sum_{ij=0}^3 F_{ij} \varepsilon^i(m) \wedge \varepsilon^j(m)$$

This is a classic change of basis with :

$$\varepsilon^i(m) = \sum_{\alpha=0}^3 P_\alpha^i d\xi^\alpha \Leftrightarrow d\xi^\alpha = \sum_{i=0}^3 P_i^\alpha \varepsilon^i(m)$$

$$F_{ij} = \sum_{\alpha\beta=0}^3 \mathcal{F}_{\alpha\beta} P_i^\alpha P_j^\beta \Leftrightarrow \mathcal{F}_{\alpha\beta} = \sum_{ij=0}^3 F_{ij} P_\alpha^i P_\beta^j \quad (5.66)$$

In the Special Relativity context the components F_{ij} are the components of the field, measured in an inertial orthonormal frame, so all usual formulas in this context can be transposed in curved space time by using the frame $(\varepsilon^i(m))_{i=0}^3$.

The dual Clifford bundle

The Clifford bundle $Cl(TM)$ is defined through the basis $(\varepsilon_i(m))_{i=0}^3$. In a change of gauge on P_G :

$$\mathbf{p}(m) = \varphi_G(m, 1) \rightarrow \tilde{\mathbf{p}}(m) = \mathbf{p}(m) \cdot \chi(m)^{-1}$$

The holonomic basis of $P_G[\mathbb{R}^4, \mathbf{Ad}]$ changes as :

$$\varepsilon_i(m) = (\mathbf{p}(m), \varepsilon_i) \rightarrow \tilde{\varepsilon}_i(m) = \chi(m)^{-1} \cdot \varepsilon_i(m) \cdot \chi(m)$$

and the Clifford product of vectors as :

$$\varepsilon_{i_1}(m) \cdot \varepsilon_{i_2}(m) \cdot \dots \cdot \varepsilon_{i_p}(m) \rightarrow \tilde{\varepsilon}_{i_1}(m) \cdot \dots \cdot \tilde{\varepsilon}_{i_p}(m) = \chi(m)^{-1} \cdot \varepsilon_{i_1}(m) \cdot \varepsilon_{i_2}(m) \cdot \dots \cdot \varepsilon_{i_p}(m) \cdot \chi(m)$$

that is :

$$\tilde{\varepsilon}_{i_1}(m) \cdot \dots \cdot \tilde{\varepsilon}_{i_p}(m) = \mathbf{Ad}_{\chi(m)^{-1}} \varepsilon_{i_1}(m) \cdot \dots \cdot \varepsilon_{i_p}(m)$$

and the components of an element of $Cl(TM)$ in the holonomic basis $(\varepsilon_i(m))_{i=0}^3$ as

$$\tilde{w}^{i_1 \dots i_p} = [\mathbf{Ad}]_{j_1}^{i_1} \dots [\mathbf{Ad}]_{j_p}^{i_p} w^{j_1 \dots j_p}$$

$$\text{so } Cl(TM) \equiv P_G[Cl(3,1), \mathbf{Ad}]$$

The Clifford algebra built on covectors : ε^i in \mathbb{R}^4 with the bisymmetric linear form of same signature is isomorphic to the Clifford algebras that we have denoted $Cl(3,1)$ or $Cl(1,3)$: this is just the replacement of ε_i by the

dual ε^i . Notice that $\varepsilon^i(\varepsilon_j) = \delta_j^i$ and not η_{ij} . We will denote them $Cl(3,1)^*$ or $Cl(1,3)^*$.

The Clifford bundle $Cl(TM^*)$ is the associated vector bundle defined through the basis $(\varepsilon^i(m))_{i=0}^3$. In a change of gauge on P_G the holonomic basis changes as :

$$\varepsilon^i(m) = (\mathbf{p}(m), \varepsilon^i) \rightarrow \tilde{\varepsilon}^i(m) = \chi(m) \cdot \varepsilon^i(m) \cdot \chi(m)^{-1}$$

the elements of $Cl(TM^*)$ transform as :

$$\varepsilon^{i_1}(m) \dots \varepsilon^{i_p}(m) \rightarrow \tilde{\varepsilon}^{i_1}(m) \dots \tilde{\varepsilon}^{i_p}(m) = \mathbf{Ad}_{\chi(m)} \varepsilon^{i_1}(m) \dots \varepsilon^{i_p}(m)$$

and the components of an element of $Cl(TM^*)$ in the holonomic basis $(\varepsilon^i(m))_{i=0}^3$ as

$$\tilde{w}_{i_1 \dots i_p} = [\mathbf{Ad}]_{i_1}^{j_1} \dots [\mathbf{Ad}]_{i_p}^{j_p} w_{j_1 \dots j_p}$$

So we will denote this associated vector bundle $P_G [Cl(3,1)^*, \mathbf{Ad}^{-1}]$.

There are a Spin group $Spin(3,1)^*$ and its Spin algebra $T_1Spin(3,1)^*$ defined in $P_G [Cl(3,1)^*, \mathbf{Ad}^{-1}]$ and we denote the basis of $T_1Spin(3,1)^*$ ⁶ :

$$\begin{aligned} \vec{\kappa}^1 &= \frac{1}{2} \varepsilon^3 \cdot \varepsilon^2, \vec{\kappa}^2 = \frac{1}{2} \varepsilon^1 \cdot \varepsilon^3, \vec{\kappa}^3 = \frac{1}{2} \varepsilon^2 \cdot \varepsilon^1, \\ \vec{\kappa}^4 &= \frac{1}{2} \varepsilon^0 \cdot \varepsilon^1, \vec{\kappa}^5 = \frac{1}{2} \varepsilon^0 \cdot \varepsilon^2, \vec{\kappa}^6 = \frac{1}{2} \varepsilon^0 \cdot \varepsilon^2 \end{aligned}$$

2 forms expressed in the Clifford bundle

The algebra $\oplus_{r=0}^4 \Lambda_r TM^*$ of forms on M is isomorphic (as vector space, the isomorphism does not extend to the product) to the Clifford algebra $Cl(TM^*)$. This isomorphism is defined through any orthonormal basis.

$$j : \Lambda TM^* \rightarrow Cl(TM^*) : j(\varepsilon^{i_1} \wedge \dots \wedge \varepsilon^{i_p}) = \varepsilon^{i_1} \dots \varepsilon^{i_p} \quad (5.67)$$

Then we can associate, by the isomorphism j the 2 form $\sum_{ij=0}^3 F_{ij} \varepsilon^i(m) \wedge \varepsilon^j(m)$ to the element of the dual Clifford bundle $Cl(TM^*)$:

$$j(\mathcal{F}) = \sum_{ij=0}^3 F_{ij} \varepsilon^i(m) \cdot \varepsilon^j(m) \quad (5.68)$$

It reads :

$$j(\mathcal{F}) = 2(F_{32} \varepsilon^3 \cdot \varepsilon^2 + F_{13} \varepsilon^1 \cdot \varepsilon^3 + F_{21} \varepsilon^2 \cdot \varepsilon^1 + F_{01} \varepsilon^0 \cdot \varepsilon^1 + F_{02} \varepsilon^0 \cdot \varepsilon^2 + F_{03} \varepsilon^0 \cdot \varepsilon^3)$$

With the decomposition :

$$\mathcal{F}_R = 2(F_{32} \varepsilon^3 \wedge \varepsilon^2 + F_{13} \varepsilon^1 \wedge \varepsilon^3 + F_{21} \varepsilon^2 \wedge \varepsilon^1)$$

$$\mathcal{F}_W = 2(F_{01} \varepsilon^0 \wedge \varepsilon^1 + F_{02} \varepsilon^0 \wedge \varepsilon^2 + F_{03} \varepsilon^0 \wedge \varepsilon^3)$$

and $v^*(F_R, F_W)$

$$= \frac{1}{2}(F_{32} \varepsilon^3 \cdot \varepsilon^2 + F_{13} \varepsilon^1 \cdot \varepsilon^3 + F_{21} \varepsilon^2 \cdot \varepsilon^1 + F_{01} \varepsilon^0 \cdot \varepsilon^1 + F_{02} \varepsilon^0 \cdot \varepsilon^2 + F_{03} \varepsilon^0 \cdot \varepsilon^3)$$

$$j(\mathcal{F}) = 4v^*((F_{32}, F_{13}, F_{21}), (F_{01}, F_{02}, F_{03})) = 4v^*(F_R, F_W) \quad (5.69)$$

In the basis $\vec{\kappa}^a : j(\mathcal{F}) = \sum_{a=1}^6 [j(\mathcal{F})]_a \vec{\kappa}^a$

⁶Notice that this is the basis of the subset $T_1Spin(3,1)^*$ of $Cl(M^*)$ and not the basis of the vector space dual of $T_1Spin(3,1)$.

That we can sum in the formula :

$$\{[j(\mathcal{F})]_a\}_{a=1}^6 = 4 \sum_{\alpha\beta=0}^3 \mathcal{F}_{\alpha\beta} \left\{ P_3^\alpha P_2^\beta, P_1^\alpha P_3^\beta, P_2^\alpha P_1^\beta, P_0^\alpha P_1^\beta, P_0^\alpha P_2^\beta, P_0^\alpha P_3^\beta \right\}$$

The isomorphism uses the bases of $T_m M^*$ and $P_G [\mathbb{R}^4, \mathbf{Ad}]$. We need to check how $j(\mathcal{F})$ behaves in a change of gauge or chart.

In a change of chart on M, the variables change as :

$$P_i^\alpha \rightarrow \tilde{P}_i^\alpha = \sum_\lambda J_\lambda^\alpha P_i^\lambda$$

$$\mathcal{F}_{\alpha\beta} \rightarrow \tilde{\mathcal{F}}_{\alpha\beta} = \sum_{\gamma\eta} K_\alpha^\gamma K_\beta^\eta \mathcal{F}_{\gamma\eta}$$

$$\text{with the jacobian : } J = \left[J_\beta^\alpha \right] = \left[\frac{\partial \tilde{\xi}^\alpha}{\partial \xi^\beta} \right] \text{ and } K = J^{-1}$$

$$\widetilde{[j(\mathcal{F})]_a} = \sum_{\alpha\beta=0}^3 \tilde{\mathcal{F}}_{\alpha\beta} \left\{ \tilde{P}_3^\alpha \tilde{P}_2^\beta, \tilde{P}_1^\alpha \tilde{P}_3^\beta, \tilde{P}_2^\alpha \tilde{P}_1^\beta, \tilde{P}_0^\alpha \tilde{P}_1^\beta, \tilde{P}_0^\alpha \tilde{P}_2^\beta, \tilde{P}_0^\alpha \tilde{P}_3^\beta \right\}$$

$$\widetilde{[j(\mathcal{F})]_a} = \sum_{\alpha\beta=0}^3 \sum_{\gamma\eta} K_\alpha^\gamma K_\beta^\eta \mathcal{F}_{\gamma\eta} \sum_{\lambda\mu} J_\lambda^\alpha J_\mu^\beta \left\{ P_3^\lambda P_2^\mu, P_1^\lambda P_3^\mu, P_2^\lambda P_1^\mu, P_0^\lambda P_1^\mu, P_0^\lambda P_2^\mu, P_0^\lambda P_3^\mu \right\}$$

$$\widetilde{[j(\mathcal{F})]_a} = \sum_{\lambda\mu=0}^3 \mathcal{F}_{\lambda\mu} \left\{ P_3^\lambda P_2^\mu, P_1^\lambda P_3^\mu, P_2^\lambda P_1^\mu, P_0^\lambda P_1^\mu, P_0^\lambda P_2^\mu, P_0^\lambda P_3^\mu \right\}$$

$$\widetilde{[j(\mathcal{F})]_a} = [j(\mathcal{F})]_a$$

So $j(\mathcal{F})$ is invariant in a change of chart on M.

Strength of the fields in the Clifford algebras

The fields are valued in the Lie algebras. The corresponding quantities read :

$$F_{Aij}^a = \sum_{\alpha\beta=0}^3 \mathcal{F}_{A\alpha\beta}^a P_i^\alpha P_j^\beta$$

$$j(\mathcal{F}_A) = 2 \sum_{a=1}^m$$

$$\left(F_{A32}^a \varepsilon^3 \cdot \varepsilon^2 + F_{A13}^a \varepsilon^1 \cdot \varepsilon^3 + F_{A21}^a \varepsilon^2 \cdot \varepsilon^1 + F_{A01}^a \varepsilon^0 \cdot \varepsilon^1 + F_{A02}^a \varepsilon^0 \cdot \varepsilon^2 + F_{A03}^a \varepsilon^0 \cdot \varepsilon^3 \right) \otimes \vec{\theta}_a$$

Expressed with the $m \times 6$ matrix $[j(\mathcal{F}_A^a)]$:

$$[j(\mathcal{F}_A)]_1^a = 4F_{A32}^a; [j(\mathcal{F}_A)]_2^a = 4F_{A13}^a; [j(\mathcal{F}_A)]_3^a = 4F_{A21}^a; [j(\mathcal{F}_A)]_4^a = 4F_{A01}^a; [j(\mathcal{F}_A)]_5^a = 4F_{A02}^a; [j(\mathcal{F}_A)]_6^a = 4F_{A03}^a$$

$$j(\mathcal{F}_A)(m) = \sum_{a=1}^m \sum_{b=1}^6 [j(\mathcal{F}_A^a)]_b^a \vec{\kappa}^b(m) \otimes \vec{\theta}_a(m) \quad (5.70)$$

Similarly for the gravitational field :

$$j(\mathcal{F}_G) = \sum_{a=1}^6 \sum_{\alpha\beta=0}^3 \mathcal{F}_{G\alpha\beta}^a P_i^\alpha P_j^\beta \varepsilon^i \otimes \varepsilon^j \otimes \vec{\kappa}_a$$

$$= \sum_{a=1}^6 \sum_{\alpha\beta=0}^3 v(\mathcal{F}_{r\alpha\beta}, \mathcal{F}_{w\alpha\beta}) P_i^\alpha P_j^\beta \varepsilon^i \otimes \varepsilon^j$$

$$j(\mathcal{F}_G)(m) = \sum_{a,b=1}^6 [j(\mathcal{F}_G)]_b^a \vec{\kappa}^b(m) \otimes \vec{\kappa}_a(m) \quad (5.71)$$

$$F_{rij}^a = \sum_{\alpha\beta=0}^3 \mathcal{F}_{r\alpha\beta}^a P_i^\alpha P_j^\beta$$

$$F_{wij}^a = \sum_{\alpha\beta=0}^3 \mathcal{F}_{w\alpha\beta}^a P_i^\alpha P_j^\beta$$

$$j(\mathcal{F}_G) = 2 \left\{ v(F_{r32}, F_{w32}) \varepsilon^3 \cdot \varepsilon^2 + v(F_{r13}, F_{w13}) \varepsilon^1 \cdot \varepsilon^3 + v(F_{r21}, F_{w21}) \varepsilon^2 \cdot \varepsilon^1 + v(F_{r01}, F_{w01}) \varepsilon^0 \cdot \varepsilon^1 + v(F_{r02}, F_{w02}) \varepsilon^0 \cdot \varepsilon^2 + v(F_{r03}, F_{w03}) \varepsilon^0 \cdot \varepsilon^3 \right\}$$

Expressed with the 6×6 matrix $[j(\mathcal{F}_G^a)]$
 $a=1,2,3 : [j(\mathcal{F}_G)]_1^a = 4F_{r32}^a; [j(\mathcal{F}_G)]_2^a = 4F_{r13}^a; [j(\mathcal{F}_G)]_3^a = 4F_{r21}^a; [j(\mathcal{F}_G)]_4^a = 4F_{r01}^a; [j(\mathcal{F}_G)]_5^a = 4F_{r02}^a; [j(\mathcal{F}_G)]_6^a = 4F_{r03}^a$
 $a=4,5,6 : [j(\mathcal{F}_G)]_1^a = 4F_{w32}^a; [j(\mathcal{F}_G)]_2^a = 4F_{w13}^a; [j(\mathcal{F}_G)]_3^a = 4F_{w21}^a; [j(\mathcal{F}_G)]_4^a = 4F_{w01}^a; [j(\mathcal{F}_G)]_5^a = 4F_{w02}^a; [j(\mathcal{F}_G)]_6^a = 4F_{w03}^a$

In a change of gauge on the principal bundle $P_G : \mathbf{p}_G(m) = \varphi_G(m, 1) \rightarrow \tilde{\mathbf{p}}_G(m) = \mathbf{p}_G(m) \cdot \chi(m)^{-1}$ the holonomic basis becomes with $\chi(m) \in Spin(3, 1)$
 $\vec{\kappa}^b \rightarrow \widetilde{\vec{\kappa}}^b = \mathbf{Ad}_\chi \vec{\kappa}^b$ notice that this is a 6×6 matrix for $\mathbf{Ad}_\chi : T_1 Spin(3, 1) \rightarrow T_1 Spin(3, 1)$

$$\begin{aligned} j(\mathcal{F}_A) &= \sum_{a=1}^m \sum_{b=1}^6 [j(\mathcal{F}_A)]_b^a \vec{\kappa}^b \otimes \vec{\theta}_a = \sum_{a=1}^m \sum_{b=1}^6 [\widetilde{j(\mathcal{F}_A)}]_b^a \widetilde{\vec{\kappa}}^b \otimes \widetilde{\vec{\theta}}_a \\ &= \sum_{a=1}^m \sum_{b=1}^6 [j(\mathcal{F}_A)]_b^a [\mathbf{Ad}_\chi]_c^b \vec{\kappa}^c \otimes \vec{\theta}_a \\ [j(\mathcal{F}_A)]_c^a &= \sum_{b=1}^6 [j(\mathcal{F}_A)]_b^a [\mathbf{Ad}_\chi]_c^b \\ [\widetilde{j(\mathcal{F}_A)}] &= [j(\mathcal{F}_A)] [\mathbf{Ad}_{\chi^{-1}}] \end{aligned} \quad (5.72)$$

For the gravitational field :

$$\begin{aligned} \vec{\kappa}^b &\rightarrow \mathbf{Ad}_\chi \vec{\kappa}^b \\ \vec{\kappa}_a &\rightarrow \mathbf{Ad}_{\chi^{-1}} \vec{\kappa}_a \\ j(\mathcal{F}_G)(m) &= \sum_{a,b=1}^6 [j(\mathcal{F}_G)]_b^a \vec{\kappa}^b \otimes \vec{\kappa}_a = \sum_{a,b=1}^6 [\widetilde{j(\mathcal{F}_G)}]_b^a \widetilde{\vec{\kappa}}^b \otimes \widetilde{\vec{\kappa}}_a \\ &= \sum_{a,b=1}^6 [j(\mathcal{F}_G)]_b^a \mathbf{Ad}_\chi \vec{\kappa}^b \otimes \mathbf{Ad}_{\chi^{-1}} \vec{\kappa}_a \\ &= \sum_{a,bcd=1}^6 [j(\mathcal{F}_G)]_b^a [\mathbf{Ad}_\chi]_d^b \vec{\kappa}^d \otimes [\mathbf{Ad}_{\chi^{-1}}]_a^c \vec{\kappa}_c \\ [j(\mathcal{F}_G)] &= [\mathbf{Ad}_{\chi^{-1}}] [j(\mathcal{F}_G)] [\mathbf{Ad}_\chi] \\ [\widetilde{j(\mathcal{F}_G)}] &= [\mathbf{Ad}_\chi] [j(\mathcal{F}_G)] [\mathbf{Ad}_{\chi^{-1}}] \end{aligned} \quad (5.73)$$

In a change of gauge on $P_U : \mathbf{p}_U(m) = \varphi_U(m, 1) \rightarrow \tilde{\mathbf{p}}_U(m) = \mathbf{p}_U(m) \cdot \chi(m)^{-1}$

$j(\mathcal{F}_G)$ does not change.

$$\begin{aligned} \vec{\theta}_a(m) &\rightarrow \widetilde{\vec{\theta}}_a(m) = \mathbf{Ad}_{\chi^{-1}} \vec{\theta}_a \\ j(\mathcal{F}_A) &= 4 \sum_{a=1}^m \sum_{b=1}^6 [j(\mathcal{F}_A)]_b^a \vec{\kappa}^b \otimes \vec{\theta}_a = \sum_{a=1}^m \sum_{b=1}^6 [\widetilde{j(\mathcal{F}_A)}]_b^a \widetilde{\vec{\kappa}}^b \otimes \widetilde{\vec{\theta}}_a \\ &= \sum_{ac=1}^m \sum_{b=1}^6 [j(\mathcal{F}_A)]_b^a \vec{\kappa}^b \otimes [\mathbf{Ad}_{\chi^{-1}}]_a^c \vec{\theta}_c \\ [\widetilde{j(\mathcal{F}_A)}] &= [\mathbf{Ad}_\chi] [j(\mathcal{F}_A)] \end{aligned} \quad (5.74)$$

$(T_1 Spin(3, 1)^* \otimes T_1 Spin(3, 1), \mathbf{Ad}^{-1} \times \mathbf{Ad})$ is a representation of $Spin(3, 1)$

$P_G [T_1 Spin(3, 1)^* \otimes T_1 Spin(3, 1), \mathbf{Ad}^{-1} \times \mathbf{Ad}]$ is an associated vector bundle

The fields are defined everywhere, and we can associate to \mathcal{F}_G the section $j(\mathcal{F}_G)$ of the associated bundle :

$$(\mathbf{p}(m), j(\mathcal{F}_G)) \sim (\mathbf{p}(m) \cdot \chi^{-1}, [\mathbf{Ad}_\chi] j(\mathcal{F}_G) [\mathbf{Ad}_{\chi^{-1}}])$$

$(T_1 Spin(3, 1)^* \otimes T_1 U, \mathbf{Ad}^{-1} \times Ad)$ is a representation of $Spin(3, 1) \times U$

$Q [T_1 Spin(3, 1)^* \otimes T_1 U, \mathbf{Ad}^{-1} \times Ad]$ is an associated vector bundle

The fields are defined everywhere, and we can associate to \mathcal{F}_A the section $j(\mathcal{F}_A)$ of the associated bundle :

$$(\mathbf{q}(m), j(\mathcal{F}_A)) \sim (\mathbf{q}(m) \cdot (\chi^{-1}, g^{-1}), [Ad_g] j(\mathcal{F}_A) [\mathbf{Ad}_{\chi^{-1}}])$$

$j(\mathcal{F}_G), j(\mathcal{F}_A)$ are deemed tensors : they have an intrinsic existence, independent of any chart or gauge, and their measure changes, in a change of gauge on P_G, Q , in a clear and unified way, with respect to the observer. They involve only the strength of the field, as expected, and we will see that the potentials cannot figure explicitly in a lagrangian, so this is consistent. And we state :

Proposition 101 *The relativist momentum density of the fields is represented as sections of the associated vector bundles :*

For the gravitational field :

$$j(\mathcal{F}_G) \in \mathfrak{X}(P_G [T_1 Spin(3, 1)^* \otimes T_1 Spin(3, 1), \mathbf{Ad}^{-1} \times \mathbf{Ad}])$$

For the other fields : $j(\mathcal{F}_A) \in \mathfrak{X}(Q [T_1 Spin(3, 1)^* \otimes T_1 U, \mathbf{Ad}^{-1} \times Ad])$

For the EM field $T_1 U = \mathbb{R}, T_1 Spin(3, 1)^* \otimes \mathbb{R} \sim T_1 Spin(3, 1)^*$ so the relativist momentum :

$$j(\mathcal{F}_{EM}) \in P_G [T_1 Spin(3, 1)^*, \mathbf{Ad}^{-1}]$$

$$j(\mathcal{F}_{EM}) = \sum_{a=1}^6 F_a \vec{k}^a = 2 \sum_{i,j=0}^3 \sum_{\alpha,\beta=0}^3 \mathcal{F}_{\alpha\beta} P_i^\alpha P_j^\beta \varepsilon^i \cdot \varepsilon^j$$

5.4.5 Energy of the fields

It is intuitive that Force fields carry energy. Because fields are present everywhere, it will be a density, valued at any point and related to the volume form ϖ_4 . It should be a quantity which changes in a consistent way with the observer. And, as for particles, it would be logical if it involves a scalar product of the relativist momentum. We could do it from the expression above, but it is more illuminating to proceed from the expression in any chart, as it leads to introduce useful tools. There are two vector spaces involved : the tangent space of 2 forms with the holonomic basis of a chart, and the Lie algebra. So we need to address successively both spaces.

Scalar product of forms over M

There is a scalar product G_r on the space $\Lambda_r M$ of scalar r forms computed with the metric g (Maths.1611). This is a bilinear symmetric form, which does not depend on a chart, is non degenerate and definite positive if g is Riemannian.

The **Hodge dual** $*\lambda$ of a r form λ is a $4 - r$ form (Maths.1613) such that :

$$\forall \mu \in \Lambda_r(M) : \mu \wedge *\lambda = G_r(\mu, \lambda) \varpi_4$$

The Hodge dual $*\mathcal{F}$ of a scalar 2-form $\mathcal{F} \in \Lambda_2 M$ is a 2 form whose expression, with the Lorentz metric, is simple when a specific ordering is used, which have been used before about the polarization of the fields :

$$\begin{aligned}\mathcal{F} &= \mathcal{F}_R + \mathcal{F}_W \text{ with} \\ \mathcal{F}_R &= 2 (\mathcal{F}_{32} d\xi^3 \wedge d\xi^2 + \mathcal{F}_{13} d\xi^1 \wedge d\xi^3 + \mathcal{F}_{21} d\xi^2 \wedge d\xi^1) \\ \mathcal{F}_W &= 2 (\mathcal{F}_{01} d\xi^0 \wedge d\xi^1 + \mathcal{F}_{02} d\xi^0 \wedge d\xi^2 + \mathcal{F}_{03} d\xi^0 \wedge d\xi^3) \\ &\text{(the 2 accounting for the symmetric part)}\end{aligned}$$

By raising the indices with g we get :

$$\begin{aligned}\mathcal{F}^{\alpha\beta} &= \sum_{\lambda\mu} g^{\alpha\lambda} g^{\beta\mu} \mathcal{F}_{\lambda\mu} \\ *\mathcal{F} &= *\mathcal{F}_R + *\mathcal{F}_W \\ *\mathcal{F}_R &= 2 (\mathcal{F}^{01} d\xi^3 \wedge d\xi^2 + \mathcal{F}^{02} d\xi^1 \wedge d\xi^3 + \mathcal{F}^{03} d\xi^2 \wedge d\xi^1) \det P' \\ *\mathcal{F}_W &= 2 (\mathcal{F}^{32} d\xi^0 \wedge d\xi^1 + \mathcal{F}^{13} d\xi^0 \wedge d\xi^2 + \mathcal{F}^{21} d\xi^0 \wedge d\xi^3) \det P'\end{aligned}$$

thus the components of the parts are exchanged and the indices are raised with the metric g . Notice that the Hodge dual is a 2 form : even if the notation uses raised indexes, they refer to the basis $d\xi^\alpha \wedge d\xi^\beta$.

Take any two scalar 2 forms \mathcal{F}, K and their decomposition as above, a straightforward computation gives :

$$\begin{aligned}*\mathcal{F}_W \wedge K_W &= 0 \\ *\mathcal{F}_W \wedge K_R &= -4 (\mathcal{F}^{32} K_{32} + \mathcal{F}^{13} K_{13} + \mathcal{F}^{21} K_{21}) \varpi_4 \\ *\mathcal{F}_R \wedge K_W &= -4 (\mathcal{F}^{01} K_{01} + \mathcal{F}^{02} K_{02} + \mathcal{F}^{03} K_{03}) \varpi_4 \\ *\mathcal{F}_R \wedge K_R &= 0 \\ G_2(\mathcal{F}_W, K_W) &= G_2(\mathcal{F}_R, K_R) = 0 \\ G_2(\mathcal{F}_W, K_R) &= -4 (\mathcal{F}^{32} K_{32} + \mathcal{F}^{13} K_{13} + \mathcal{F}^{21} K_{21}) \\ G_2(\mathcal{F}_R, K_W) &= -4 (\mathcal{F}^{01} K_{01} + \mathcal{F}^{02} K_{02} + \mathcal{F}^{03} K_{03}) \\ \text{From there, because } G_2 &\text{ is bilinear :} \\ G_2(\mathcal{F}, K) &= G_2(\mathcal{F}_W + \mathcal{F}_R, K_R + K_W) \\ &= -4 (\mathcal{F}^{32} K_{32} + \mathcal{F}^{13} K_{13} + \mathcal{F}^{21} K_{21} + \mathcal{F}^{01} K_{01} + \mathcal{F}^{02} K_{02} + \mathcal{F}^{03} K_{03})\end{aligned}$$

$$G_2(\mathcal{F}, K) = -2 \sum_{\alpha\beta} \mathcal{F}^{\alpha\beta} K_{\alpha\beta} \quad (5.75)$$

$$\begin{aligned}\text{These quantities can be expressed in the orthonormal basis } &(\varepsilon^i(m))_{i=0}^3. \\ G_2(\mathcal{F}, K) &= -2 \sum_{\alpha\beta} \mathcal{F}^{\alpha\beta} K_{\alpha\beta} = -2 \sum_{\alpha\beta} g^{\alpha\lambda} g^{\beta\mu} \sum_{ijpq=0}^3 F_{ij} P_\lambda^i P_\mu^j K_{pq} P_\alpha^p P_\beta^q \\ &= -2 \sum_{ijpq} \eta^{pi} \eta^{qj} F_{ij} K_{pq} = -2 \sum_{ij=0}^3 \eta^{ii} \eta^{jj} F_{ij} K_{ij} \\ &= -2 \left(\sum_{j=1}^3 \eta^{00} \eta^{jj} F_{0j} K_{0j} + \sum_{i=0}^3 \eta^{ii} \eta^{00} F_{i0} K_{i0} + \sum_{ij=1}^3 \eta^{ii} \eta^{jj} F_{ij} K_{ij} \right) \\ &= -2 \left(\sum_{ij=1}^3 F_{ij} K_{ij} - 2 \sum_{j=1}^3 F_{0j} K_{0j} \right)\end{aligned}$$

$$G_2(\mathcal{F}, \mathcal{F}) = 4 \left(\sum_{j=1}^3 (F_{0j})^2 - \sum_{\{ij\}=1}^3 (F_{ij})^2 \right) \quad (5.76)$$

G_2 is not definite positive or negative. The decomposition emphasizes the role played by the component along ε_0 : the spatial components (F_{ij}) and the temporal components (F_{0j}) contribute to the scalar product with definite opposite signs.

This scalar product can easily be expressed for $J(\mathcal{F})$ using the scalar product on the Clifford algebra $Cl(3,1)^*$:

$$\begin{aligned} \langle v(r, w), v(r', w') \rangle_{Cl(3,1)^*} &= \frac{1}{4} (r^t r' - w^t w') \\ J(\mathcal{F}) &= 4 \sum_{a=1}^6 \sum_{b=1}^m v^* ((F_{32}, F_{13}, F_{21}), (F_{01}, F_{02}, F_{03})) \in T_1 Spin(3,1)^* \\ \langle J(\mathcal{F}), J(\mathcal{F}) \rangle_{Cl(3,1)^*} &= \frac{1}{4} 16 \left((F_{32})^2 + (F_{13})^2 + (F_{21})^2 - \left((F_{01})^2 + (F_{02})^2 + (F_{03})^2 \right) \right) \\ \langle J(\mathcal{F}), J(\mathcal{F}) \rangle_{Cl(3,1)^*} &= 4 (F_R^t F_R - F_W^t F_W) = -G_2(\mathcal{F}, \mathcal{F}) \end{aligned}$$

$$\langle J(\mathcal{F}), J(\mathcal{F}) \rangle_{Cl(3,1)^*} = 4 \left(\sum_{\{ij\}=1}^3 (F_{ij})^2 - \sum_{j=1}^3 (F_{0j})^2 \right) = -G_2(\mathcal{F}, \mathcal{F}) \quad (5.77)$$

So, up to the sign, we have the same quantity.

Scalar products on the Lie algebras

The strength can be seen as a section of the associated vector bundles $P_G [T_1 Spin(3,1), Ad], P_U [T_1 U, Ad]$ and then the scalar product must be preserved by the adjoint map Ad . There are not too many possibilities. It can be shown that, for simple groups of matrices, the only scalar products on their Lie algebra which are invariant by the adjoint map are of the kind : $\langle [X], [Y] \rangle = k Tr ([X]^* [Y])$ which sums up, in our case, to use the Killing form. This is a bilinear form (Maths.1609) which is preserved by any automorphism of the Lie algebra (thus in any representation). However it is negative definite if and only if the group is compact and semi-simple (Maths.1847).

Scalar product for the gravitational field : We have seen in the previous Chapter that the scalar product on $T_1 Spin(3,1)$, induced by the scalar product on the Clifford algebra, is up to a constant the Killing form :

$$\begin{aligned} \langle v(r, w), v(r', w') \rangle_{Cl(3,1)} &= \frac{1}{4} (r^t r' - w^t w') \\ a=1,2,3 : \mathcal{F}_{G\alpha\beta}^a &= \mathcal{F}_{r\alpha\beta}^a \\ a=4,5,6 : \mathcal{F}_{G\alpha\beta}^a &= \mathcal{F}_{w\alpha\beta}^a \\ \langle \mathcal{F}_{G\alpha\beta}(m), \mathcal{F}'_{G\lambda\mu}(m) \rangle_{Cl} &= \left\langle v(\mathcal{F}_{r\alpha\beta}, \mathcal{F}_{w\alpha\beta}), v(\mathcal{F}'_{r\lambda\mu}, \mathcal{F}'_{w\lambda\mu}) \right\rangle_{Cl} \\ &= \frac{1}{4} \left(\mathcal{F}_{r\alpha\beta}^t \mathcal{F}'_{r\lambda\mu} - \mathcal{F}_{w\alpha\beta}^t \mathcal{F}'_{w\lambda\mu} \right) \\ \langle \mathcal{F}_{G\alpha\beta}(m), \mathcal{F}'_{G\lambda\mu}(m) \rangle_{Cl} &= \frac{1}{4} \left(\sum_{a=1}^3 \mathcal{F}_{G\alpha\beta}^a \mathcal{F}'_{G\lambda\mu} - \sum_{a=4}^6 \mathcal{F}_{G\alpha\beta}^a \mathcal{F}'_{G\lambda\mu} \right) \quad (5.78) \end{aligned}$$

The result does not depend on the signature. This scalar product is invariant in a change of gauge, non degenerate but not definite positive.

Scalar product for the other fields

The group U is assumed to be compact and connected. If U is semi-simple, its Killing form, which is invariant by the adjoint map, is then definite negative, and we can define a definite positive scalar product, invariant in a change of gauge, on its Lie algebra. This is the case for $SU(2)$ and $SU(3)$ but not for $U(1)$, however the Lie algebra of $U(1)$ is \mathbb{R} and there is an obvious definite positive scalar product. As T_1U is a real vector space the scalar product is a *bilinear symmetric* form.

So we will assume that :

Proposition 102 *There is a definite positive scalar product on the Lie algebra T_1U , defined by a bilinear symmetric form preserved by the adjoint map, that we will denote $\langle \rangle_{T_1U}$. The basis $(\vec{\theta}_a)_{a=1}^m$ of T_1U is orthonormal for this scalar product.*

Notice that it is different from the scalar product on F (which defines the charges), which is Hermitian. In the standard model, because several groups are involved, three different constants are used, called the “gauge coupling”. Here we consider only one group, and we can take the basis $(\vec{\theta}_a)_{a=1}^m$ as orthonormal for the scalar product.

The scalar product between sections \mathcal{F}_A of $\Lambda_2(M; T_1U)$ is then defined, pointwise, as

$$\langle \mathcal{F}_{A\alpha\beta}(m), \mathcal{F}'_{A\lambda\mu}(m) \rangle_{T_1U} = \sum_{a=1}^m \mathcal{F}_{A\alpha\beta}^a(m) \mathcal{F}'_{A\lambda\mu}{}^a(m) \quad (5.79)$$

Scalar product for the strength of the fields

We have to combine both scalar products.

The scalar product on $T_1Spin(3, 1)^* \otimes T_1Spin(3, 1)$ is expressed by :

$$\begin{aligned} & \left\langle \sum_{ab=1}^6 [X]_a^b \kappa^a \otimes \kappa_b, \sum_{cd=1}^6 [Y]_c^d \kappa^c \otimes \kappa_d \right\rangle \\ &= \left\langle \sum_{b=1}^6 \left(\sum_{a=1}^6 [X]_a^b \kappa^a \right) \otimes \kappa_b, \sum_{d=1}^6 \left(\sum_{c=1}^6 [Y]_c^d \kappa^c \right) \otimes \kappa_d \right\rangle \\ &= \frac{1}{4} \left(\sum_{b=1}^3 \left\langle \sum_{a=1}^6 [X]_a^b \kappa^a, \sum_{c=1}^6 [Y]_c^b \kappa^c \right\rangle - \sum_{b=4}^6 \left\langle \sum_{a=1}^6 [X]_a^b \kappa^a, \sum_{c=1}^6 [Y]_c^b \kappa^c \right\rangle \right) \\ &= \frac{1}{4} \sum_{b=1}^3 \frac{1}{4} \left(\sum_{a=1}^3 [X]_a^b [Y]_a^b - \sum_{a=4}^6 [X]_a^b [Y]_a^b \right) \\ &\quad - \frac{1}{4} \sum_{b=4}^6 \left(\frac{1}{4} \left(\sum_{a=1}^3 [X]_a^b [Y]_a^b - \sum_{a=4}^6 [X]_a^b [Y]_a^b \right) \right) \\ &= \frac{1}{16} \sum_{a=1}^3 \left(\sum_{b=1}^3 [X]_a^b [Y]_a^b - \sum_{b=4}^6 [X]_a^b [Y]_a^b \right) \\ &\quad - \sum_{a=4}^6 \left(\sum_{b=1}^3 [X]_a^b [Y]_a^b - \sum_{b=4}^6 [X]_a^b [Y]_a^b \right) \\ & \langle j(\mathcal{F}_G), j(\mathcal{F}_G) \rangle \\ &= \sum_{b=1}^3 \left([F_{r32}^b]^2 + [F_{r13}^b]^2 + [F_{r21}^b]^2 \right) - \left([F_{w32}^b]^2 + [F_{w13}^b]^2 + [F_{w21}^b]^2 \right) \end{aligned}$$

$$\begin{aligned}
& - \left([F_{r01}^b]^2 + [F_{r02}^b]^2 + [F_{r03}^b]^2 \right) + \left([F_{w01}^b]^2 + [F_{w02}^b]^2 + [F_{w03}^b]^2 \right) \\
& = \sum_{b=1}^3 \left([F_{r32}^b]^2 + [F_{r13}^b]^2 + [F_{r21}^b]^2 \right) - \left([F_{r01}^b]^2 + [F_{r02}^b]^2 + [F_{r03}^b]^2 \right) \\
& - \left([F_{w32}^b]^2 + [F_{w13}^b]^2 + [F_{w21}^b]^2 \right) - \left([F_{w01}^b]^2 + [F_{w02}^b]^2 + [F_{w03}^b]^2 \right) \\
& = \sum_{b=1}^3 -\frac{1}{4} G_2 (\mathcal{F}_r^b, \mathcal{F}_r^b) + \frac{1}{4} G_2 (\mathcal{F}_w^b, \mathcal{F}_w^b) \\
& = \sum_{b=1}^3 -\frac{1}{4} \left(-2 \sum_{\alpha\beta} \mathcal{F}_r^{b\alpha\beta} \mathcal{F}_{r\alpha\beta}^b \right) + \frac{1}{4} \left(-2 \sum_{\alpha\beta} \mathcal{F}_w^{b\alpha\beta} \mathcal{F}_{w\alpha\beta}^b \right) \\
& = \frac{1}{2} \sum_{\alpha\beta=0}^3 \left((\mathcal{F}_r^{\alpha\beta})^t \mathcal{F}_{r\alpha\beta} - (\mathcal{F}_w^{\alpha\beta})^t \mathcal{F}_{w\alpha\beta} \right) \\
\langle j(\mathcal{F}_G), j(\mathcal{F}_G) \rangle & = \frac{1}{2} \sum_{\alpha\beta=0}^3 \left((\mathcal{F}_r^{\alpha\beta})^t \mathcal{F}_{r\alpha\beta} - (\mathcal{F}_w^{\alpha\beta})^t \mathcal{F}_{w\alpha\beta} \right) \quad (5.80)
\end{aligned}$$

For the other fields, pointwise :

$$\begin{aligned}
& \left\langle \sum_{b=1}^m \sum_{a=1}^6 [X]_a^b \kappa^a \otimes \theta_b, \sum_{d=1}^m \sum_{c=1}^6 [Y]_c^d \kappa^c \otimes \theta_d \right\rangle \\
& = \left\langle \sum_{b=1}^m \left(\sum_{a=1}^6 [X]_a^b \kappa^a \right) \otimes \theta_b, \sum_{d=1}^m \left(\sum_{c=1}^6 [Y]_c^d \kappa^c \right) \otimes \theta_d \right\rangle \\
& = \sum_{b=1}^m \left\langle \sum_{a=1}^6 [X]_a^b \kappa^a, \sum_{c=1}^6 [Y]_c^b \kappa^c \right\rangle \\
& = \sum_{b=1}^m \frac{1}{4} \left(\sum_{a=1}^3 [X]_a^b [Y]_a^b - \sum_{a=4}^6 [X]_a^b [Y]_a^b \right)
\end{aligned}$$

$$\begin{aligned}
& \langle j(\mathcal{F}_A), j(\mathcal{F}_A) \rangle \\
& = 4 \sum_{b=1}^m \left(\sum_{a=1}^3 [F_{A32}^b]^2 + [F_{A13}^b]^2 + [F_{A21}^b]^2 - \left([F_{A01}^b]^2 + [F_{A02}^b]^2 + [F_{A03}^b]^2 \right) \right) \\
& = 4 \sum_{b=1}^m -\frac{1}{4} G (\mathcal{F}_A^b, \mathcal{F}_A^b) \\
& = -2 \sum_{b=1}^m \sum_{\alpha\beta} \mathcal{F}_A^{b\alpha\beta} \mathcal{F}_{A\alpha\beta}^b \\
& = -2 \sum_{\alpha\beta} \left(\mathcal{F}_A^{\alpha\beta} \right)^t \mathcal{F}_{A\alpha\beta}
\end{aligned}$$

$$\langle j(\mathcal{F}_A), j(\mathcal{F}_A) \rangle = -2 \sum_{\alpha\beta=0}^3 \sum_{a=1}^m \mathcal{F}_A^{a\alpha\beta} \mathcal{F}_{A\alpha\beta}^a = -2 \sum_{\alpha\beta=0}^3 \left(\mathcal{F}_A^{\alpha\beta} \right)^t \mathcal{F}_{A\alpha\beta} \quad (5.81)$$

These scalar products are invariant in a change of charts or in a change of gauge. However their expression depends on the decomposition between the transversal and rotational components, characteristic of the observer, as can be seen in the orthonormal basis. They do not depend on the signature of the metric but they are not definite positive. All these quantities are, of course, estimated up to constants depending on the units.

Energy of the fields

And we can state :

Proposition 103 *The energy density of the fields is the scalar product $\langle j(\mathcal{F}_G), j(\mathcal{F}_G) \rangle, \langle j(\mathcal{F}_A), j(\mathcal{F}_A) \rangle$*

The question of the energy of the gravitational field has been a topic of discussion for years and it is generally accepted that it has no precise value (see Wald). We have here another, consistent answer : there is a density of relativist momentum and energy, which have precise definitions, and whose measure depends on the observer. Notice that these quantities can be computed in the usual frame work : it then involves the Riemann tensor, it is simply more obvious in the fiber bundle presentation. Moreover the gravitational field has a distinctive feature : the energy of the gravitational field has two components which have opposite signs. This can explain the fact that the gravitational field, as we measure it, is exceptionally weak : the electromagnetic force is some 39 orders of magnitude greater than the force of gravity.

The energy of the fields does not depend on the chart or the observer, it is similar to $\langle \psi_0, \psi_0 \rangle$ for particles. This quantity is not necessarily positive. However we will see that, incorporated in a lagrangian along with the definition of the energy for particles, it provides the expected solutions. We must also remember that the similar scalar product $\langle S_0, S_0 \rangle$ can be negative for antiparticles.

In Thermodynamics, only the variation of the energy of a system has a physical meaning, and it was a source of discussion when Relativity introduced the idea of an absolute energy for particles with $c^2 \langle P, P \rangle$. This absolute energy is linked to the existence of a fundamental state (here represented by ψ_0). And one can wonder if something similar does exist for the fields. Indeed QTF considers the “energy of the vacuum”. This vacuum is intended as an area where there is no particle or boson (so where is no field), but it has quite an animated life with virtual particles. So it is worth to consider the existence of a fundamental state for the fields.

The quantity for which it could be defined must be geometric, so it would apply to $j(\mathcal{F}_G), j(\mathcal{F}_A)$, which would change with the action $\mathbf{Ad}^{-1} \times \mathbf{Ad}, \mathbf{Ad}^{-1} \times \mathbf{Ad}$ of the groups. But then we would have to introduce explicitly the value $s(m), g(m)$ of elements of $Spin(3, 1)$ and U :

$$\begin{aligned} j(\mathcal{F}_G)(m) &= [\mathbf{Ad}_{s(m)}] [j(\mathcal{F}_G)_0] [\mathbf{Ad}_{s^{-1}(m)}] \\ j(\mathcal{F}_A)(m) &= [Ad_{g(m)}] [j(\mathcal{F}_G)_0] [\mathbf{Ad}_{s^{-1}(m)}] \end{aligned}$$

Formally this is not much trouble for $Spin(3, 1)$ but less so for U , at least for the fields other than the EM field. Physically we would expect that similarly :

$$\psi(m) = \vartheta(s(m), g(m)) \psi_0$$

Then the key variables of any model would be sections $s \in P_G, g \in P_U$. A study in this picture can be done with the model presented in the next chapters, but we will stick to a more conventional venue.

Identity

We have a useful property which is more general, and holds for all the fields:

Theorem 104 *On the Lie algebra T_1U of a Lie group U , endowed with a symmetric scalar product $\langle \rangle_{T_1U}$ which is preserved by the adjoint map :*

$$\forall X, Y, Z \in T_1U : \langle X, [Y, Z] \rangle = \langle [X, Y], Z \rangle \quad (5.82)$$

Proof. $\forall g \in U : \langle Ad_g X, Ad_g Y \rangle = \langle X, Y \rangle$

take the derivative with respect to g at $g = 1$ for $Z \in T_1U$:

$$(Ad_g X)'(Z) = ad(Z)(X) = [Z, X]$$

$$\langle [Z, X], Y \rangle + \langle X, [Z, Y] \rangle = 0 \Leftrightarrow \langle X, [Y, Z] \rangle = \langle [Z, X], Y \rangle$$

exchange X, Z :

$$\Rightarrow \langle Z, [Y, X] \rangle = \langle [X, Z], Y \rangle = -\langle [Z, X], Y \rangle = -\langle X, [Y, Z] \rangle = -\langle Z, [X, Y] \rangle$$

■

5.4.6 Norm on the spaces of the relativist momentum of the fields

As usual a norm is required for the application of QM theorems.

For the gravitational field :

$$j(\mathcal{F}_G) \in \mathfrak{X}(P_G [T_1Spin(3, 1)^* \otimes T_1Spin(3, 1), \mathbf{Ad}^{-1} \times \mathbf{Ad}])$$

$$\text{For the other fields : } j(\mathcal{F}_A) \in \mathfrak{X}(Q [T_1Spin(3, 1)^* \otimes T_1U, \mathbf{Ad}^{-1} \times Ad])$$

With the Cartan decomposition in $Cl(3, 1)^* \otimes Cl(3, 1)$:

$$T_1Spin(3, 1)^* \otimes T_1Spin(3, 1) = (L_0^* \oplus P_0^*) \otimes (L_0 \oplus P_0) = (L_0^* \otimes L_0) \oplus (L_0^* \otimes P_0) \oplus (P_0^* \otimes L_0) \oplus (P_0^* \otimes P_0)$$

The projections :

$$\pi_{ll} : T_1Spin(3, 1)^* \otimes T_1Spin(3, 1) \rightarrow L_0^* \otimes L_0$$

$$\pi_{lp} : T_1Spin(3, 1)^* \otimes T_1Spin(3, 1) \rightarrow L_0^* \otimes P_0, \dots$$

are well defined

$$\forall T \in T_1Spin(3, 1)^* \otimes T_1Spin(3, 1) = \pi_{ll}(T) + \pi_{lp}(T) + \pi_{pl}(T) + \pi_{pp}(T)$$

The scalar product is extended to the tensorial product by :

$$\forall X, X' \in l_0, Y, Y' \in p_0 : \langle X \otimes Y, X' \otimes Y' \rangle = \langle X, X' \rangle_{l_0} \langle Y, Y' \rangle_{p_0}$$

and similarly for the others. The norm on $T_1Spin(3, 1)^* \otimes T_1Spin(3, 1)$ is then :

$$\|T\| = \sqrt{\langle \pi_{ll}(T), \pi_{ll}(T) \rangle + \langle \pi_{lp}(T), \pi_{lp}(T) \rangle + \langle \pi_{pl}(T), \pi_{pl}(T) \rangle + \langle \pi_{pp}(T), \pi_{pp}(T) \rangle}$$

We define the norm on the space $\Lambda_2(M; T_1Spin(3, 1))$ by :

$$\|\mathcal{F}_G\| = \|j(\mathcal{F}_G)\|$$

Similarly for \mathcal{F}_A :

$$T_1Spin(3, 1) \otimes T_1U = (l_0^* \oplus p_0^*) \otimes T_1U = (l_0^* \otimes T_1U) \oplus (p_0^* \otimes T_1U)$$

$$\forall T \in T_1Spin(3, 1) \otimes T_1U = \pi_l(T) + \pi_p(T)$$

$$\forall X, X' \in l_0, Y, Y' \in T_1U : \langle X \otimes Y, X' \otimes Y' \rangle = \langle X, X' \rangle_{l_0} \langle Y, Y' \rangle_{T_1U}$$

$$\|T\| = \sqrt{\langle \pi_l(T), \pi_l(T) \rangle + \langle \pi_p(T), \pi_p(T) \rangle}$$

$$\|\mathcal{F}_A\| = \|j(\mathcal{F}_A)\|$$

These norms are invariant in a change of chart or in a change of gauge.

With these norms the sets :

$$\begin{aligned} L_G^1 &= L^1(M, \mathfrak{X}(P_G [T_1 Spin(3, 1)^* \otimes T_1 Spin(3, 1), \mathbf{Ad}^{-1} \times \mathbf{Ad}], \varpi_4)) = \\ &= \{j(\mathcal{F}_G) \in \mathfrak{X}(P_G [T_1 Spin(3, 1)^* \otimes T_1 Spin(3, 1), \mathbf{Ad}^{-1} \times \mathbf{Ad}]) : \int_{\Omega} \|\mathcal{F}_G\| \varpi_4 < \infty\} \\ L_A^1 &= L^1(M, \mathfrak{X}(Q [T_1 Spin(3, 1)^* \otimes T_1 U, \mathbf{Ad}^{-1} \times Ad]), \varpi_4) = \\ &= \{j(\mathcal{F}_A) \in \mathfrak{X}(Q [T_1 Spin(3, 1)^* \otimes T_1 U, \mathbf{Ad}^{-1} \times Ad]) : \int_{\Omega} \|\mathcal{F}_A\| \varpi_4 < \infty\} \end{aligned}$$

are separable Fréchet spaces.

5.4.7 Structure of the force fields

Force Fields present several specific features, which can be easily identified in the representation with fiber bundles.

Quantization of the force fields

We consider as variables sections of the vector bundle as defined above. We can implement the theorem 22. There are Hilbert spaces H_G, H_A , isometries Υ_G, Υ_A , which associate a vector to each map $j(\mathcal{F}_G), j(\mathcal{F}_A)$:

$$\begin{aligned} \Upsilon_G &: L_G^1 \rightarrow H_G \\ \Upsilon_A &: L_A^1 \rightarrow H_A \\ \widehat{U}_G &= \Upsilon_G \circ \mathbf{Ad}^{-1} \times \mathbf{Ad} \circ \Upsilon_G^{-1} \\ \widehat{U}_A &= \Upsilon_A \circ \mathbf{Ad}^{-1} \times Ad \circ \Upsilon_A^{-1} \\ (H_G, \widehat{U}_G) &\text{ is a unitary representation of } Spin(3, 1), \\ (H_A, \widehat{U}_A) &\text{ is a unitary representation of } Spin(3, 1) \times U \end{aligned}$$

Moreover we can also implement the theorem 24 by adding the variables $\langle j(\mathcal{F}_G), j(\mathcal{F}_G) \rangle = Y_G, \langle j(\mathcal{F}_A), j(\mathcal{F}_A) \rangle = Y_A$. To each level of energy of the fields in the system is associated a subset $H_G(Y_G), H_A(Y_A)$ invariant by $\widehat{U}_G, \widehat{U}_A$, that is belonging to one of the irreducible representations of $(H_G, \widehat{U}_G), (H_A, \widehat{U}_A)$.

The only unitary representations of $Spin(3, 1)$ are infinite dimensional, and parametrized by two scalars $z \in \mathbb{Z}, k \in \mathbb{R}$.

U is a compact group so its unitary representations are the sum of finite dimensional unitary irreducible representations.

As a consequence, for a given level of energy, the gravitational field is characterized by two scalars $z \in \mathbb{Z}, k \in \mathbb{R}$. The real k can be linked to Y_G and we will see below that z is linked to the spin of the field.

For the other fields we have the two scalars $z \in \mathbb{Z}, k \in \mathbb{R}$ and a scalar d related to the representation of U. The representation (F, ϱ) is reducible for particles, and similarly the force fields split in the components corresponding to the different kinds of interactions. The scalar k can be linked to Y_A and d to the type of field.

Spin of the Field

Let us consider any section $j(\mathcal{F}_A) \in \mathfrak{X}(Q[T_1Spin(3,1)^* \otimes T_1U, \mathbf{Ad}^{-1} \times Ad])$. Its decomposition in :

$$j(\mathcal{F}_A) = j(\mathcal{F}_{Ar}) + j(\mathcal{F}_{Aw})$$

with :

$$j(\mathcal{F}_{Ar}) = 2 \sum_{a=1}^m (F_{A32}^a \varepsilon^3 \cdot \varepsilon^2 + F_{A13}^a \varepsilon^1 \cdot \varepsilon^3 + F_{A21}^a \varepsilon^2 \cdot \varepsilon^1) \otimes \vec{\theta}_a(m)$$

$$j(\mathcal{F}_{Aw}) = 2 \sum_{a=1}^m (F_{A01}^a \varepsilon^0 \cdot \varepsilon^1 + F_{A02}^a \varepsilon^0 \cdot \varepsilon^2 + F_{A03}^a \varepsilon^0 \cdot \varepsilon^3) \otimes \vec{\theta}_a(m)$$

depends only on the choice of ε^0 , that is of ε_0 . For a given ε^0 the set $\{\gamma^C(\sigma_r)j(\mathcal{F}_{Ar}), \sigma_r \in SO(3)\}$ is invariant by any rotation of the spatial basis, which respects the future orientation (whence the restriction to $SO(3)$). But we can go further.

There is a projection $\pi : T_1Spin(3,1) \rightarrow T_1Spin(3)$ within $Cl(3,1)$ which does not depend on the choice of a basis, as long as $\langle \varepsilon_0, \varepsilon_0 \rangle = -1$ (see the Chapter Clifford Algebras). So we can define similarly the projection :

$$\pi : T_1Spin(3,1)^* \otimes T_1U \rightarrow T_1Spin(3)^* \otimes T_1U :: \pi(j(\mathcal{F}_A)) = j(\mathcal{F}_{Ar})$$

and the relation of equivalence on $\mathfrak{X}(Q[T_1Spin(3,1)^* \otimes T_1U, \mathbf{Ad}^{-1} \times Ad])$:

$$j(\mathcal{F}_A) \sim j(\mathcal{F}'_A) \Leftrightarrow \pi(j(\mathcal{F}_A)) = \pi(j(\mathcal{F}'_A))$$

For the EM field $j(\mathcal{F}_{Ar})$ corresponds to the magnetic component. Actually $j(\mathcal{F}_{Ar})$ is similar to the spin of a particle, and we will call it the **Spin of the field**. It can be represented by a section

$j(\mathcal{F}_{Ar}) \in \mathfrak{X}(Q[T_1Spin(3)^* \otimes T_1U, \mathbf{Ad}^{-1} \times Ad])$, that is by a geometric object which does not depend on a spatial gauge (ε_0 must stay constant), whose measure changes with the action of $SO(3)$. Moreover $SO(3)$ is a compact group so $j(\mathcal{F}_{Ar})$ is quantized : for a level of energy of the field $j(\mathcal{F}_{Ar})$ belongs to a $2j + 1$ dimensional vector space, isomorphic to (P^j, D^j) with $j \in \mathbb{N}$. *The field has an integer Spin.*

Curvature of the Gravitational field

With the Cartan decomposition in $Cl(3,1)^* \otimes Cl(3,1)$:

$$T_1Spin(3,1)^* \otimes T_1Spin(3,1) = (L_0^* \oplus P_0^*) \otimes (L_0 \oplus P_0) = (L_0^* \otimes L_0) \oplus (L_0^* \otimes P_0) \oplus (P_0^* \otimes L_0) \oplus (P_0^* \otimes P_0)$$

The projection :

$\pi : T_1Spin(3,1)^* \otimes T_1Spin(3,1) \rightarrow L_0^* \otimes L_0$ is well defined and $L_0^* \otimes L_0$ has the structure $T_1Spin(3)^* \otimes T_1Spin(3)$

$$\begin{aligned} \pi(j(\mathcal{F}_G)) &= \sum_{a,b=1}^3 [j(\mathcal{F}_G)]_b^a \vec{\kappa}^b(m) \otimes \vec{\kappa}_a(m) \\ &= 4 \sum_{a=1}^3 (F_{r32}^a \vec{\kappa}^1 + F_{r13}^a \vec{\kappa}^2 + F_{r21}^a \vec{\kappa}^3) \otimes \vec{\kappa}_a \end{aligned}$$

$$\pi(j(\mathcal{F}_G)) = 4 \sum_{a=1}^3 \sum_{\alpha\beta=0}^3 \mathcal{F}_{r\alpha\beta}^a \left(P_3^\alpha P_2^\beta \vec{\kappa}^1 + P_3^\alpha P_2^\beta \vec{\kappa}^2 + P_3^\alpha P_2^\beta \vec{\kappa}^3 \right) \otimes \vec{\kappa}_a \quad (5.83)$$

We can define as above a relation of equivalence on

$\mathfrak{X}(P_G[T_1Spin(3,1)^* \otimes T_1Spin(3,1), \mathbf{Ad}^{-1} \times \mathbf{Ad}])$:

$$j(\mathcal{F}_G) \sim j(\mathcal{F}'_G) \Leftrightarrow \pi(j(\mathcal{F}_G)) = \pi(j(\mathcal{F}'_G))$$

and the class of equivalence of $j(\mathcal{F}_G)$ does not depend on a spatial basis. As above $\pi(j(\mathcal{F}_G))$ can be defined as the **spin of the gravitational field**. It is quantized : $\pi(j(\mathcal{F}_G))$ belongs to a $2j + 1$ dimensional vector space, isomorphic to (P^j, D^j) with $j \in \mathbb{N}$. *The gravitational field has an integer Spin*. The spin depends only on the components $\mathcal{F}_{r\alpha\beta}^a$.

An elementary representation of the gravitational field, on a given $\Omega_3(t)$, involves only $\pi(j(\mathcal{F}_G))$: a tensor defined in an euclidean, spatial basis. The vacuum, on a hypersurface $\Omega_3(t)$, is invariant by the action of $SO(3)$. For t fixed the maps $\pi(j(\mathcal{F}_G))(\varphi_o(t, x))$ are sum of harmonic polynomials, and clearly one can identify the arguments of the functions, at least locally, with coordinates of points in $\Omega_3(t)$. Harmonic polynomials have been introduced originally as solutions of the gravitational field, and so we are lead to identify $\pi(j(\mathcal{F}_G))$ with the spatial component of the gravitational field, as it is usually understood. Which gives a physical meaning to the decomposition of \mathcal{F}_G in $\mathcal{F}_r, \mathcal{F}_w$.

The tensor $j(\mathcal{F}_G)$ reads, with the dual basis

$$j(\mathcal{F}_G) = 4\left\{\sum_{a=1}^6 F_{G32}^a \vec{\kappa}^1 \otimes \vec{\kappa}_a + F_{G13}^a \vec{\kappa}^2 \otimes \vec{\kappa}_a + F_{G21}^a \vec{\kappa}^3 \otimes \vec{\kappa}_a + F_{G01}^a \vec{\kappa}^4 \otimes \vec{\kappa}_a + F_{G02}^a \vec{\kappa}^5 \otimes \vec{\kappa}_a + F_{G03}^a \vec{\kappa}^6 \otimes \vec{\kappa}_a\right\}$$

Thus we can contract the tensor, and the scalar :

$$\begin{aligned} R' &= 4(F_{G32}^1 + F_{G13}^2 + F_{G21}^3 + F_{G01}^4 + F_{G02}^5 + F_{G03}^6) \\ &= 4\sum_{\alpha\beta=0}^3 \mathcal{F}_{r\alpha\beta}^1 P_3^\alpha P_2^\beta + \mathcal{F}_{r\alpha\beta}^2 P_3^\alpha P_2^\beta + \mathcal{F}_{r\alpha\beta}^3 P_3^\alpha P_2^\beta + \mathcal{F}_{w\alpha\beta}^1 P_0^\alpha P_1^\beta + \mathcal{F}_{w\alpha\beta}^2 P_0^\alpha P_2^\beta + \\ &\quad \mathcal{F}_{w\alpha\beta}^3 P_0^\alpha P_3^\beta \end{aligned}$$

$$R' = 2\left(-\sum_{aij=1}^3 \sum_{\alpha\beta} \epsilon(a, i, j) \mathcal{F}_{r\alpha\beta}^a P_i^\alpha P_j^\beta + 2\sum_{a=1}^3 \sum_{\alpha=1}^3 \mathcal{F}_{w0\alpha}^a P_a^\alpha\right)$$

is invariant in a change of gauge or chart. And we retrieve the usual scalar curvature, up to a constant :

$$\begin{aligned} \mathbf{R} &= \sum_{a=1}^3 2\sum_{\alpha=1}^3 \mathcal{F}_{w0\alpha}^a [P]_a^\alpha + \sum_{\alpha\beta} \mathcal{F}_{r\alpha\beta}^a \sum_{p,q=1}^3 \epsilon(a, q, p) [P]_p^\beta [P]_q^\alpha \\ R' &= -2R \end{aligned}$$

From the quantization of $j(\mathcal{F}_G)$ one can deduce that *the scalar curvature should be related to the level of energy and to the spin*.

The curvature is comprised of a translational and a rotational part. The latter is fully defined by the spin :

$$R'_R = 4(F_{G32}^1 + F_{G13}^2 + F_{G21}^3)$$

It is invariant by a change of spatial gauge, and its value is quantized.

The decomposition of \mathcal{F}_G in its transversal and spatial components seems physically significant. \mathcal{F}_G is equivalent to the Riemann tensor, for which the decomposition is less obvious. The physical meaning of the torsion has perhaps been wrongly understood. So this remark suggests that the right variable to study the gravitational field is (G_r, G_w) with 24 components, and not the Levi-Civita connection.

Symmetry of the force field

It is common to assume that a field is symmetric : a geometric transformation in the space $\Omega(t)$ of an observer is such that the field looks the same. It is

equivalent to consider a global change of observer given by an element $s(t) \in Spin(3)$.

For the gravitational field :

$$[j(\mathcal{F}_G)] = [\mathbf{Ad}_s] [j(\mathcal{F}_G)] [\mathbf{Ad}_{s^{-1}}] = [j(\mathcal{F}_G)] \Leftrightarrow [\mathbf{Ad}_s] [j(\mathcal{F}_G)] = [j(\mathcal{F}_G)] [\mathbf{Ad}_s]$$

with the 6×6 matrix :

$$[j(\mathcal{F}_G)] = \begin{bmatrix} F_r^r & F_w^r \\ F_r^w & F_w^w \end{bmatrix}$$

and $s(t) = a_r(t) + v(r(t), 0)$ it sums up to :

$$F_r^r = \left([1 + a_r j(r) + \frac{3}{4} j(r) j(r)] F_r^r + [1 - a_r^2 + \frac{1}{4} j(r) j(r)] F_w^r \right) [1 - a_r j(r) + \frac{3}{4} j(r) j(r)]$$

$$F_r^w = [a_r^2 + a_r j(r) + \frac{1}{2} j(r) j(r)] F_r^w [1 - a_r j(r) + \frac{3}{4} j(r) j(r)] =$$

$$F_w^r = \left\{ \left([1 + a_r j(r) + \frac{3}{4} j(r) j(r)] F_r^r + [1 - a_r^2 + \frac{1}{4} j(r) j(r)] F_w^r \right) [1 - a_r^2 + \frac{1}{4} j(r) j(r)] \right. \\ \left. + \left([1 + a_r j(r) + \frac{3}{4} j(r) j(r)] F_w^r + [1 - a_r^2 + \frac{1}{4} j(r) j(r)] F_r^w \right) [a_r^2 - a_r j(r) + \frac{1}{2} j(r) j(r)] \right\}$$

$$F_w^w = [a_r^2 + a_r j(r) + \frac{1}{2} j(r) j(r)] F_r^w [1 - a_r^2 + \frac{1}{4} j(r) j(r)]$$

$$+ [a_r^2 + a_r j(r) + \frac{1}{2} j(r) j(r)] F_w^w [a_r^2 - a_r j(r) + \frac{1}{2} j(r) j(r)] =$$

The set of polynomials of 3×3 matrices $A + B_j(r) + C_j(r) j(r)$ is a commutative algebra, so the solution is :

$$[j(\mathcal{F}_G)] = \begin{bmatrix} [A_r^r + B_r^r j(r) + C_r^r j(r) j(r)] & [A_w^r + B_w^r j(r) + C_w^r j(r) j(r)] \\ [A_r^w + B_r^w j(r) + C_r^w j(r) j(r)] & [A_w^w + B_w^w j(r) + C_w^w j(r) j(r)] \end{bmatrix}$$

and the equations give 3 potential solutions :

i) $r = 0$: this is the spherical solution

ii) $a_r^2 = \frac{3}{5}; A_r^w = 0;$

$$\left[-\frac{2}{5} - \frac{1}{4} j(r) j(r) \right] F_w^r + \left[\frac{2}{5} + \frac{1}{4} j(r) j(r) \right] (F_r^r + \frac{3}{5} F_w^w) = 0$$

iii) $F_r^w = 0;$

$$[a_r^2 - 1 + (\frac{5}{4} a_r^2 - 1) j(r) j(r)] F_w^r + [1 - a_r^2 + \frac{1}{4} j(r) j(r)] (F_r^r + a_r^2 F_w^w) = 0$$

For the other fields :

$$[j(\mathcal{F}_A)] = [j(\mathcal{F}_A)] [\mathbf{Ad}_{\chi^{-1}}] = [j(\mathcal{F}_A)]$$

with the $m \times 6$ matrix :

$$[j(\mathcal{F}_G)] = [F_r \quad F_w]$$

$$F_r [1 - a_r j(r) + \frac{3}{4} j(r) j(r)] = F_r$$

$$F_r [\frac{1}{4} r^t r + \frac{1}{4} j(r) j(r)] + F_w [a_r^2 - a_r j(r) + \frac{1}{2} j(r) j(r)] = F_w$$

For the EM field the only solution is :

$[F_r \quad F_w] = \lambda(t) [r(t) \quad r(t)]$ with a scalar function $\lambda(t)$, that is plane waves with a wave vector $r(t)$.

Chern-Weil Theory

There is a, complementary, approach to the quantization of fields. On a given principal bundle not any principal connections can exist. The conditions, which involve \mathcal{F} , are the topic of the Chern-Weil theory (Maths.27.4.5). They apply also for the passage from one strength \mathcal{F} to another one \mathcal{F}' , so to the evolution of \mathcal{F} , for instance in a discontinuous process. The rules are a bit abstract, but

depend only on the manifold M and the group U (or $Spin(3, 1)$). This leads to a topological oriented study of the fields.

We could consider the definition of the Charge of a field, as we have done for particles. However we do not have the ingredient which was crucial : a fundamental state $j(\mathcal{F}_A)_0$.

Chapter 6

LAGRANGIAN

The Principle of Least Action states that for any system there is some quantity which is stationary when the system is at its equilibrium. For a system which is represented by quantitative variables and their derivatives, defined in the context of a vector bundle $E(M, V, \pi)$ the involved quantity can be expressed in a general way as a functional, that is a map :

$$\ell : \mathfrak{X}(J^r E) \rightarrow \mathbb{R}$$

which acts on sections Z of the r -jet prolongation of E , represented by their coordinates $Z = (z_{\alpha_1 \dots \alpha_s}^i, i = 1 \dots n, s = 0, \dots, r)$. And the stationarity is understood as a local extremum : around the equilibrium Z_0 it is impossible to change $\ell(Z_0)$ by an infinitesimal variation δZ of Z . It is natural to assume that there is some function $L : \mathfrak{X}(J^r E) \rightarrow \mathbb{C}$ over which ℓ acts and, if this action is linear, then the functional can be expressed as an integral (Maths.2342), in this case:

$$\begin{aligned} \ell &= \int_{\Omega} \mathcal{L}(Z(m)) d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3 = \\ &= \int_{\Omega} L(Z(m)) \det P' d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3 = \int_{\Omega} L(Z(m)) \varpi_4 \end{aligned}$$

L is then a real scalar function called the scalar lagrangian.

In Mathematics a lagrangian is usually defined as a $\dim M$ form on M , and their theory, as well as the Variational Calculus, is extensive but also rather technical (see Maths.34.2). When only vector bundles are considered, as this is usually the case in Physics, the problems are more simple and one can use an extension of the theory of distributions (or generalized functions) which is similar to the functional derivatives well known by the physicists. We will see its application in the next chapter, and for the time being, we will study the specification of the scalar lagrangian L .

This is a crucial step, because L sums up much of the Physics of the model. The specification of a lagrangian is an art in itself and many variants have been proposed. The Standard Model is built around a complicated lagrangian (see Wikipedia Standard Model for its expression) which is the result of many attempts and patches to find a solution which fits the results of experiments. It is useful to remind, at this step, that one of the criteria in the choice and validation of a scientific theory is efficiency. Physicists must be demanding about their basic concepts, upon which everything is built, but, as they proceed to more

specific problems, they can relax a bit. There is no Theory or a unique Model of Everything, which would be suited to all problems. The framework that we have exposed provides several tools, which can be selected according to the problems at hand. So we continue in the same spirit, and, fortunately, in the choice of the right lagrangian there are logical rules, coming essentially from the Principle of Relativity : the solution should be equivariant in a change of observer, which entails that the lagrangian itself, which is a scalar function, should be invariant. This condition provides strong guidelines in its specification, that we will see now. The methods that we expose are general, but as we have done so far, they are more easily understood when implemented on an example, and we will use the variables and representations which have been developed in the previous chapters.

6.1 THE SPECIFICATION OF THE LAGRANGIAN

6.1.1 General issues

Which variables ?

We have to decide which are the variables that enter the lagrangian and the order of their derivatives. The lagrangian is a function

$L(z_{\alpha_1 \dots \alpha_s}^i, i = 1 \dots n, s = 0, \dots, r)$ and this is under this form that a solution is found. However usually the variables appear as composite expressions : for instance the derivatives of the potential $\partial_\alpha \dot{A}$ appear not as such but in the strength \mathcal{F} . So this is always a two steps process, from the composite variables to the variables and their derivatives.

We will limit ourselves to the variables which have been introduced previously, as they give a comprehensive picture of the problems. The key variables are : ψ for the state of the particle, considering that the particle belongs to some matter field, G the potential of the gravitational field, \dot{A} the potential of the other fields, and the tetrad P which, in the fiber bundle model is a variable as the others and defines the metric g . We can add a density μ , which is a scalar function on M .

All these variables are maps defined on a bounded area Ω of M , and valued in various vector bundles, so expressed in components in the relevant holonomic frames. The use of the formalism of fiber bundle enables us to study the most general problem with 4 variables only, and the consideration of matter fields (which are no more than general definition of motion and states) to address the issue of the trajectories and of rotation. Of course, whenever only gravitation is involved, ψ is reduced to the spinor S , but we will keep the most general framework.

The model is based on first order derivatives : the covariant derivative is at its core, and this is a first order operator. The strength \mathcal{F} is of first order with respect to the potentials. So in the lagrangian it is legitimate to stay at : $\partial_\alpha \psi, \partial_\alpha G_\beta, \partial_\alpha \dot{A}_\beta, \partial_\alpha P$.

Equivariance and Covariance

A lagrangian is assumed to sum up the physics of a system, so locations, velocity and all quantities are expressed in a unique chart and frame, with the flexibility offered by a change of gauge. We can assume that the chart is the standard chart φ_o of an observer. The equilibrium of the Principle of Least Action is the equilibrium for this observer. But the notion of equilibrium, by itself, suggests that it does not depend on the observer. So the lagrangian and the solutions should be, not invariant, by equivariant in a change of observer. The equilibrium is not expressed by the same figures, but it is still an equilibrium and one can go from one set of data to another by using mathematical relations deduced from the respective disposition of the observers.

In any model based on manifolds (and I remind that an affine space is a manifold, so this applies also in Galilean Geometry) a lagrangian, as any other

mathematical relation, should stay the same in a change of chart. This condition is usually called covariance.

In a model based on fiber bundles there is an additional condition : the expressions must change according to the rules in a change of gauge. This condition is usually called equivariance, but it has the same meaning.

Covariance and equivariance are expressed as conditions that any quantity, and of course the lagrangian, must meet. These conditions are also a way to deal with uncertainty which comes for the choice of some variables. For instance the orthonormal basis $(\varepsilon_i(m))$ is defined (and the tetrad with it) up to a $SO(3,1)$ matrix. The equivariance relations account for this fact.

Equivariance is usually expressed as Noether's currents (from the Mathematician Emmy Noether) and presented as the consequence of symmetries in the model. Of course if there are additional, physical symmetries, they can be accounted for in the same way. But the Noether's currents are the genuine expression of the freedom of gauge.

Once we have checked that our lagrangian (and more generally any quantity) is compliant with equivariance and covariance, of course we can exercise our freedom of gauge by choosing one specific gauge. This is how Gauge Freedom is usually introduced in Physics (in Electromagnetism we have the Gauss gauge, the Coulomb gauge,...). The goal is to simplify an expression by imposing some relations between variables. This is legitimate but, as noticed before, one must be aware that it has practical implications on the observer himself who must actually use this gauge in the collection of his data.

Time

The Principle of Locality leads naturally to express all quantities related to particles with respect to their proper time. But, whenever the propagation of the fields or several particles are considered, the state of the system must be related to a unique time, which is the time of an observer (who is arbitrary). This is necessary to have a common definition of the area of integration in the action.

The proper time of a particle and the time of the observer are related. The basic relations are (with the notations used previously) :

between the proper time τ of a particle and the time t of an observer :

$$\frac{d\tau}{dt} = \sqrt{1 - \frac{\|\vec{v}\|^2}{c^2}} = \frac{1}{c} \sqrt{-\langle V, V \rangle} = \frac{c}{u^0} = \frac{1}{2a_w^2 - 1}$$

between the velocity u of a particle and the speed V as measured by an observer :

$$u = \frac{dp}{d\tau} = V \frac{c}{\sqrt{-\langle V, V \rangle}} = \frac{1}{\sqrt{1 - \frac{\|\vec{v}\|^2}{c^2}}} (\vec{v} + c\varepsilon_0(m))$$

$$= c \left((2a_w^2 - 1) \varepsilon_0 + \epsilon a_w \sum_{i=1}^3 w_i \varepsilon_i \right)$$

$$V = \frac{dp}{dt} = u \frac{\sqrt{-\langle V, V \rangle}}{c} = \vec{v} + c\varepsilon_0(m) = c \left(\varepsilon_0 + \epsilon \frac{a_w}{2a_w^2 - 1} \sum_{i=1}^3 w_i \varepsilon_i \right)$$

Whenever particles are represented as matter fields these relations can be fully expressed with the σ_w component of the spinor.

We see in the expressions above that the distinction between proper time and time of the observer, by breaking the fundamental symmetry, requires to distinguish the relativist spin (represented by $\epsilon = \pm 1$). It is assumed that its value is a continuous variable : a change along the world line is a discontinuous process,

The distinction between proper time and time of the observer is usually ignored in QTF, in spite of its obvious significance. Some attempts have been made to confront this issue, which is linked, in Quantum Physics, to the speed of propagation of the perturbation of a wave function (see Schnaid).

Fundamental state

The assumption of the existence of a fundamental state ψ_0 is at the core of the theory. For elementary particles it is fixed, and given by the type of the particle. For a composite body, the assumption that the state can be represented by a element of $E \otimes F$ implies that there is some way to define a fundamental state : there are internal forces which keep the cohesion of the body, and the definition of the body itself provides a way to compute (in a separate model) some fundamental state.

For a solid body the computation given previously provides :

- for a rigid solid a constant inertial spinor S_0 , which is simply the basis for a spinor $S(t) = \gamma C(\sigma(t)) S_0$ along the trajectory
- for a deformable solid an inertial spinor S_B which varies along the world line : $S_B(\tau(t)) = N_B(\tau(t)) \gamma C(\sigma_B(\tau(t))) S_0$ so that $S(t) = \gamma C(\sigma(t)) S_B(\tau(t))$.

Moreover the introduction of a density (assimilated to a number of particles by unit of volume) provides some flexibility.

Partial derivatives and covariant derivatives

To implement the rules of Variational Calculus the partial derivatives $\partial_\alpha \psi$,

$\partial_\alpha \dot{A}, \partial_\alpha G, \dots$ are required. However other quantities, and notably the lagrangian, can be expressed in using the covariant derivative ∇ or the strength \mathcal{F} , which have a more physical meaning. The question is then : is it legitimate to express $\partial_\alpha G, \partial_\alpha \dot{A}$ only through the strength \mathcal{F} , and $\partial_\alpha \psi$ through the covariant derivative ? And we will see that the answer is definitively positive.

However there is an issue with the use of the covariant derivative. The interaction of the force fields with particles goes through $\nabla_\alpha \psi$ but the covariant derivative is a 1 form, and we need to choose a vector along which to take this derivative.

Because particles move along world lines, the most natural choice is to take as vector the velocity u of the particle, or, using the rules above, the speed V . The resulting quantities are proportional :

$$\nabla_V \psi = \frac{\sqrt{-(V,V)}}{c} \nabla_u \psi$$

The difference between the two quantities has a physical meaning : $\nabla_V \psi$ is the covariant derivative as measured by the observer, $\nabla_u \psi$ is an intrinsic quantity, which does not depend on the observer. However we will use the chart

of an observer, with its time t , and u is related to σ_w in the standard basis of an observer, thus the choice of $\nabla_V \psi$ is more legitimate.

The value of the potentials along the trajectory have been denoted

$$\widehat{G} = \sum_{\alpha=0}^3 V^\alpha G_\alpha$$

$$\widehat{A} = \sum_{\alpha=0}^3 V^\alpha \dot{A}_\alpha$$

V is a variable of the model, it is related to the motion that we should compute, for a given ψ_0 and for an observer by :

$$[\psi] = [\gamma C(\sigma_w \cdot \sigma_r)] [\psi_0] [\varrho(\varkappa)]$$

$$V = \frac{dp}{dt} = u \frac{\sqrt{-\langle V, V \rangle}}{c} = \vec{v} + c\varepsilon_0(m) = c \left(\varepsilon_0 + \epsilon \frac{a_w}{2a_w^2 - 1} \sum_{i=1}^3 w_i \varepsilon_i \right)$$

The choice of the key variables depends on the problem at hand :

i) Whenever the trajectory is known (for instance for bonded particles) V and then u, σ_w are known. But usually σ_r is not fixed and is a variable of the problem.

ii) For particles studied individually the trajectories are given (with respect to ε_0 for an observer) by a map : $\sigma : [0, T] \rightarrow Spin(3, 1)$ which is a variable. There is a “state equation” which is an ODE in ψ with \widehat{G}, \widehat{A} as parameters, and key variables maps $r(t), w(t)$ on one hand, and “fields equations” which relate the value of the strength $\mathcal{F}_G, \mathcal{F}_A$ to the motion of the particles, on the other hand. The combination of these equations and their fitting with the initial conditions provide the trajectories.

iii) For particles represented by matter fields, the trajectories are given by a map : $\sigma : \Omega \rightarrow Spin(3, 1)$ which is a variable. Usually there is also a density, with the continuity equation. The solutions are computed as in ii).

In QTF the solution which is commonly chosen is different, this is the Dirac’s operator, celebrated because it is mathematically clever, but has serious drawbacks.

Dirac operator

The Dirac operator is a differential operator, and no longer a 1-form on M , defined from the covariant derivative, which does not require the choice of a vector : so it “absorbs” the α of the covariant derivative. Actually this is required in the Standard Model because the world lines are not explicit, but the Dirac’s operator can be defined in a very large context (Maths.32.28), including GR, and in our formalism its meaning is more obvious. The general Dirac operator D is weakly elliptic, and D^2 is a scalar operator (Maths.2495,2496).

The mechanism is the following :

i) using the isomorphism between TM and the dual bundle TM^* provided by the metric g , to each covector $\omega = \sum_{\alpha=0}^3 \omega_\alpha d\xi^\alpha$ one can associate a vector : $\omega^* = \sum_{\alpha\beta=0}^3 g^{\alpha\beta} \omega_\alpha \partial\xi_\beta$

ii) vectors $v = \sum_{\alpha=0}^3 v^\alpha \partial\xi^\alpha$ of TM can be seen as elements of the Clifford bundle $Cl(M)$ and as such acts on $\mathbf{e}_p(m) \otimes \mathbf{f}_q(m)$ by :

$$v = \sum_{\alpha j=0}^3 v^\alpha P_\alpha^{!j} \varepsilon_j(m) \text{ in the orthogonal frame}$$

ε_j acts on $\mathbf{e}_p(m) \otimes \mathbf{f}_q(m)$ by γC :

$$(\mathbf{q}(m), \mathbf{e}_p(m) \otimes \mathbf{f}_q(m))$$

$$\rightarrow \left(\mathbf{q}(m), \sum_{\alpha,j=0}^3 v^\alpha P_\alpha'^j ([\gamma C(\varepsilon_j)] \mathbf{e}_p(m)) \otimes \mathbf{f}_q(m) \right)$$

iii) thus there is an action of TM^* on $\mathbf{e}_p(m) \otimes \mathbf{f}_q(m)$ with $v = \omega^*$

$$(\mathbf{q}(m), \gamma C(\omega^*) \psi(m))$$

$$= \left(\mathbf{p}(m), \sum_{\alpha,\beta,j=0}^3 g^{\alpha\beta} \omega_\alpha P_\beta'^j ([\gamma C(\varepsilon_j)] \mathbf{e}_p(m)) \otimes \mathbf{f}_q(m) \right)$$

and as the tetrad defines the metric g :

$$\sum_\beta g^{\alpha\beta} P_\beta'^j = \sum_{\beta,kl} \eta^{kl} P_k^\alpha P_l^\beta P_\beta'^j = \sum_k \eta^{kj} P_k^\alpha$$

$$\sum_{\alpha,\beta,j=0}^3 g^{\alpha\beta} \omega_\alpha P_\beta'^j [\gamma C(\varepsilon_j)] \mathbf{e}_p(m) \otimes \mathbf{f}_q(m) = \sum_{\alpha,\beta=0}^3 g^{\alpha\beta} \omega_\alpha [\gamma C(\partial\xi_\beta)] \mathbf{e}_p(m) \otimes$$

$$\mathbf{f}_q(m) = \sum_{\alpha=0}^3 \omega_\alpha [\gamma C(d\xi^\alpha)] \mathbf{e}_p(m) \otimes \mathbf{f}_q(m)$$

iv) the covariant derivative is a one form on M so one can take $\varpi = \nabla_\alpha$ and the Dirac operator is :

$$D : \mathfrak{X}(J^1Q[E \otimes F, \vartheta]) \rightarrow \mathfrak{X}(J^1Q[E \otimes F, \vartheta]) :: D\psi = \sum_{\alpha=0}^3 [\gamma C(d\xi^\alpha)] [\nabla_\alpha \psi] \quad (6.1)$$

$$D\psi = \sum_{\alpha=0}^3 [P_i^\alpha [\gamma C(\varepsilon^i)] [\nabla_\alpha \psi]$$

$$\varepsilon^i(\varepsilon_j) = \delta_j^i \Rightarrow \gamma C(\varepsilon^i) = \gamma C(\varepsilon_i)^{-1}$$

$$D\psi = \sum_{\alpha=0}^3 [P_a^\alpha [\gamma C(\varepsilon_a)] [\nabla_\alpha \psi]$$

So the Dirac operator can be seen as the trace of the covariant operator, which averages the action of the covariant derivative along the directions $\alpha = 0 \dots 3$ which are put on the same footing. This is mathematically convenient, and consistent with the notion of undifferentiated matter field, but this has no real physical justification : it is clear that one direction is privileged on the world line.

We have seen that $\langle \psi, \nabla_\alpha \psi \rangle = i \operatorname{Im} \langle \psi, \nabla_\alpha \psi \rangle$ which is convenient to define the energy of the particle in the system. But the Dirac's operator exchanges the chirality. The scalar product $\langle \psi, D\psi \rangle$ is not necessarily a real quantity and, with the matrices γ used in QTF, can be null, which is one of the reasons for the introduction of the Higgs boson (see Schücker).

Hamiltonian

In Classic Mechanics the time t is totally independent from the other geometric coordinates, so the most natural formulation of the Principle of Least Action takes the form (Maths.2606) :

$$\ell(Z) = \int_0^T L(t, q^i, y^i) dt$$

where y^i stands for $\frac{dq^i}{dt}$ in the 1-jet formalism, and the change of variable with the conjugate momenta :

$$p^i = \frac{\partial L}{\partial q^i}$$

$$H = \sum_{i=1}^n p^i y^i - L$$

leads to the Hamilton equations :

$$\frac{dq^i}{dt} = \frac{\partial H}{\partial p^i}, \quad \frac{dp^i}{dt} = -\frac{\partial H}{\partial q^i}$$

which are the translation of the Euler-Lagrange equations with the new variables.

In QM the operator in the Schrödinger equation is assumed to be the Hamiltonian : $i\hbar \frac{\partial \psi}{\partial t} = H\psi$ and this has been an issue at the origin of Quantum Physics, because of the specific role played by the time, which seemed to be inconsistent with the covariance required by Relativity. After many attempts it has lead to the path integral formalism, which uses the lagrangian and is viewed as compatible both with Relativity and QM.

However, as we noticed above, even if in a relativist lagrangian the coordinates are masked by a chart, it is not true that the coordinate time is banalized. To study, in a consistent manner, any system, we need a single time, and this is necessarily the time of an observer. We have to check that the formulation of the lagrangian is consistent with the Principle of Relativity : the equilibrium must be an equilibrium for any observer, but the definition of the system itself is observer-dependant. This is obvious with the foliation : the geometric area Ω of the Universe encompassed by the system during its evolution is not the same as the one of another observer. The covariance must be assured in any change of chart which respects this foliation, but that does not mean that the time itself is not specific. The Hamiltonian formulation is certainly not appropriate in the relativist context, but for many other reasons (for instance the Maxwell's equations, and more generally the concept of fields are not compatible with the Galilean Geometry) than the distinction of a privileged time.

Conservation of Momentum and Energy

In a system comprised of interacting particles and field, each object has its own momentum and energy, which is not necessarily conserved. So to define a conservation of momentum and energy of the system we should find a way to aggregate these components, and it would be with respect to an observer who defines the time of the system.

Usually the lagrangian is built around the energy, which is a scalar function L and can be summed with respect to a volume form ϖ_4 . The energy of the system can then be defined as :

$$\mathcal{E} = \int_{\Omega} L(z^i, z_{\alpha}^i) \varpi_4$$

and the conservation of energy for the observer means that ¹:

$$\mathcal{E}(t) = \int_{\Omega(t)} L(z^i, z_{\alpha}^i) \varpi_3 = Ct = \int_{\Omega(t)} i_{\varepsilon_0} (L(z^i, z_{\alpha}^i) \varpi_4)$$

Consider the manifold $\Omega([t_1, t_2])$ with borders $\Omega(t_1), \Omega(t_2)$:

$$\mathcal{E}(t_2) - \mathcal{E}(t_1) = \int_{\partial\Omega([t_1, t_2])} i_{\varepsilon_0} (L\varpi_4) = \int_{\Omega([t_1, t_2])} d(i_{\varepsilon_0} L\varpi_4)$$

$$d(i_{\varepsilon_0} L\varpi_4) = \mathcal{L}_{\varepsilon_0} (L\varpi_4) - i_{\varepsilon_0} d(L\varpi_4)$$

$$= (\mathcal{L}_{\varepsilon_0} L) \varpi_4 + L \mathcal{L}_{\varepsilon_0} \varpi_4 - i_{\varepsilon_0} (dL \wedge \varpi_4) - i_{\varepsilon_0} L d\varpi_4$$

¹Notice the difference with a similar computation done for material bodies : material bodies are characterized by a unique vector field V , but in a general system the unique reference is ε_0 .

$$\begin{aligned}
&= L'(\varepsilon_0) \varpi_4 + L(\operatorname{div} \varepsilon_0) \varpi_4 - i_{\varepsilon_0}(dL \wedge \varpi_4) \\
&= \operatorname{div}(L\varepsilon_0) \varpi_4 \\
\mathcal{E}(t_2) - \mathcal{E}(t_1) &= \int_{\Omega(t_1, t_2)} \operatorname{div}(L\varepsilon_0) \varpi_4 \\
&\text{(Maths.1517,1587)}
\end{aligned}$$

The conservation of energy for the observer imposes a specific condition : $\operatorname{div}(L\varepsilon_0)$ which is not necessarily met for a stationary solution. However we will see that it is met for usual models.

As for the momenta the situation is worse, because they are not defined in the same bundles.

So, at least for continuous models which are based on lagrangians, the conservation of energy and momentum is not a key principle.

6.1.2 Specification of a General Lagrangian

We will use the precise notation :

L denotes the scalar lagrangian $L(z^i, z_\alpha^i)$ function of the variables z^i , expressed by the components in the gauge of the observer, and their partial derivatives which, in the jets bundle formalism, are considered as independent variables z_α^i .

$$\mathcal{L} = L(z^i, z_\alpha^i) (\det P')$$

$$L\varpi_4 = L(z^i, z_\alpha^i) (\det P') d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3 \text{ is the 4-form}$$

$\frac{\partial \mathcal{L}}{\partial z}$ to denote the usual partial derivative with respect to the variable z

$\frac{d\mathcal{L}}{dz}$ to denote the total derivative with respect to the variable z , meaning accounting for the composite expressions in which it is an argument.

We will illustrate how to compute the rules of equivariance and covariance for a general lagrangian, using the variables that we have defined previously, expressed by their coordinates : $\psi^{ij}, G_\alpha^a, \dot{A}_\alpha^a, P_i^\alpha, \partial_\beta \psi^{ij}, \partial_\beta G_\alpha^a, \partial_\beta \dot{A}_\alpha^a, \partial_\beta P_i^\alpha, V^\alpha$.

So in this section :

$$L(\psi^{ij}, G_\alpha^a, \dot{A}_\alpha^a, P_i^\alpha, \partial_\beta \psi^{ij}, \partial_\beta G_\alpha^a, \partial_\beta \dot{A}_\alpha^a, \partial_\beta P_i^\alpha, V^\alpha) \text{ in an action such as : } \int_\Omega L \mu \det P' d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3$$

All variables are represented by their coordinates in relevant bases, by real or complex scalars. L is not supposed to be holomorphic, so the real and imaginary part of the variables $\psi^{ij}, \partial_\alpha \psi^{ij}$ must appear explicitly. We will use the convenient notation for complex variables z and their conjugates \bar{z} , by introducing the holomorphic complex valued functions :

$$\frac{\partial L}{\partial z} = \frac{1}{2} \left(\frac{\partial L}{\partial \operatorname{Re} z} + \frac{1}{i} \frac{\partial L}{\partial \operatorname{Im} z} \right); \quad \frac{\partial L}{\partial \bar{z}} = \frac{1}{2} \left(\frac{\partial L}{\partial \operatorname{Re} z} - \frac{1}{i} \frac{\partial L}{\partial \operatorname{Im} z} \right) \quad (6.2)$$

\Leftrightarrow

$$\begin{aligned}
\frac{\partial L}{\partial \operatorname{Re} z} &= \frac{\partial L}{\partial z} + \frac{\partial L}{\partial \bar{z}} \\
\frac{\partial L}{\partial \operatorname{Im} z} &= i \left(\frac{\partial L}{\partial z} - \frac{\partial L}{\partial \bar{z}} \right)
\end{aligned}$$

The partial derivatives $\frac{\partial L}{\partial \operatorname{Re} z}, \frac{\partial L}{\partial \operatorname{Im} z}$ are real valued functions, so $\frac{\partial L}{\partial \bar{z}} = \overline{\frac{\partial L}{\partial z}}$. And we have the identities for any complex valued function u :

$$\frac{\partial L}{\partial \operatorname{Re} z} \operatorname{Re} u + \frac{\partial L}{\partial \operatorname{Im} z} \operatorname{Im} u = 2 \operatorname{Re} \frac{\partial L}{\partial z} u; -\frac{\partial L}{\partial \operatorname{Re} z} \operatorname{Im} u + \frac{\partial L}{\partial \operatorname{Im} z} \operatorname{Re} u = -2 \operatorname{Im} \frac{\partial L}{\partial z} u \quad (6.3)$$

To find a solution we need the explicit presence of the variables and their partial derivatives. But as our goal is to precise the specification of L, we can, without loss of generality, make the replacements :

$$\begin{aligned} \partial_\alpha \psi^{ij} &\rightarrow \nabla_\alpha \psi^{ij} = \partial_\alpha \psi^{ij} + \sum_{k=1}^4 \sum_{a=1}^6 [\gamma C(G_\alpha^a)]_k^i \psi^{kj} + \sum_{k=1}^n \psi^{ik} [\dot{A}_\alpha]_j^k \\ \partial_\beta G_\alpha^a &\rightarrow \mathcal{F}_{G\alpha\beta}^a = \partial_\alpha G_\beta^a - \partial_\beta G_\alpha^a + [G_\alpha, G_\beta]^a \text{ and } F_{G\alpha\beta} = \partial_\alpha G_\beta^a + \partial_\beta G_\alpha^a \\ \partial_\beta \dot{A}_\alpha^a &\rightarrow \mathcal{F}_{A\alpha\beta}^a = \partial_\alpha \dot{A}_\beta^a - \partial_\beta \dot{A}_\alpha^a + [\dot{A}_\alpha, \dot{A}_\beta]^a \text{ and } F_{G\alpha\beta} = \partial_\alpha \dot{A}_\beta^a + \partial_\beta \dot{A}_\alpha^a \end{aligned}$$

And the lagrangian is then a function :

$$\mathcal{L} \left(\psi^{ij}, G_\alpha^a, \dot{A}_\alpha^a, P_i^\alpha, \nabla_\alpha \psi^{ij}, \mathcal{F}_{G\alpha\beta}^a, F_{G\alpha\beta}^a, \mathcal{F}_{A\alpha\beta}^a, F_{A\alpha\beta}^a, \partial_\beta P_i^\alpha, u^\alpha \right)$$

Most of the variables above are defined up to some transformation : for instance the components of the tetrad are defined up to a matrix of $SO(3,1)$. The function L should be intrinsic, meaning invariant by :

- a change of gauge in the principal bundles P_G, P_U and their associated bundles
- a change of chart in the manifold M

The operations below will give the relations which must exist consequently between the variables and the partial derivatives of L, and some precious information about the presence or the absence of some variables.

6.1.3 Equivariance in a change of gauge

The mechanism is exposed below to find the relations between variables, but it is also at the foundation of Noether's currents.

One parameter group of change of gauge

One parameter groups of change of trivialization on a principal bundle are defined by sections of their adjoint bundle (Maths.2070) :

$$\kappa \in \mathfrak{X}(P_G [T_1 \operatorname{Spin}(3,1), \mathbf{Ad}])$$

$$\theta \in \mathfrak{X}(P_U [T_1 U, \operatorname{Ad}])$$

$\kappa = v(\kappa_r, \kappa_w), \theta$ are maps from M to the Lie algebras. At each point m , for a given value of a scalar parameter τ , the exponential on the Lie algebra defines an element of the groups at m (Maths.22.2.6) :

$$\exp : \mathbb{R} \times T_1 \operatorname{Spin}(3,1) \rightarrow \operatorname{Spin}(3,1) :: \exp(\tau \kappa(m))$$

$$\exp : \mathbb{R} \times T_1 U \rightarrow U :: \exp(\tau \theta(m))$$

The exponential on $T_1 Spin(3, 1)$ is expressed by (see Annex) :

$$\begin{aligned} \exp t\kappa &= \exp \tau v(\kappa_r, \kappa_w) = \sigma_w(\tau, \kappa_w) \cdot \sigma_r(\tau, \kappa_r) \\ \sigma_w(\tau, \kappa_w) &= a_w(\tau, \kappa_w) + \sinh \frac{1}{2} \tau \sqrt{\kappa_w^t \kappa_w} v(0, \kappa_w) \\ a_w(\tau, \kappa_w) &= \sqrt{1 + \frac{1}{4} \left(\kappa_w^t \kappa_w \sinh^2 \frac{1}{2} \tau \sqrt{\kappa_w^t \kappa_w} \right)} \\ \sigma_r(\tau, \kappa_r) &= a_r(\tau, \kappa_w) + \sin t \frac{1}{2} \sqrt{\kappa_r^t \kappa_r} v(\kappa_r, 0) \\ a_r(\tau, \kappa_w) &= \sqrt{1 - \frac{1}{4} \kappa_r^t \kappa_r \sin^2 t \frac{1}{2} \sqrt{\kappa_r^t \kappa_r}} \end{aligned}$$

It is actually multivalued (because of the double cover) so we assume that one of the value has been chosen (for instance $a > 0$). This does not matter here.

By definition the derivative of these exponential for $\tau = 0$ gives back the elements of the Lie algebras :

$$\begin{aligned} \frac{d}{d\tau} \exp(\tau\kappa(m))|_{\tau=0} &= \kappa(m) \\ \frac{d}{d\tau} \exp(\tau\theta(m))|_{\tau=0} &= \theta(m) \end{aligned}$$

With the change of gauge :

$$\begin{aligned} \mathbf{p}_G(m) &\rightarrow \tilde{\mathbf{p}}_G(m, \tau) \cdot \exp(-\tau\kappa(m)) \\ \mathbf{p}_U(m) &\rightarrow \tilde{\mathbf{p}}_U(m) \cdot \exp(-\tau\theta(m)) \end{aligned}$$

The components of the variables become :

$P_i^\alpha \rightarrow \tilde{P}_i^\alpha(m, \tau) = \sum_{j=0}^3 [h(\exp(-\tau\kappa))]_i^j P_j^\alpha$ where $[h]$ is the $SO(3, 1)$ corresponding matrix

$$\psi^{ij} \rightarrow \tilde{\psi}^{ij}(m, \tau) = \sum_{k=1}^4 \sum_{l=1}^n [\gamma C(\exp(\tau\kappa))]_k^i [\varrho(\exp(\tau\theta))]_l^j \psi^{kl}$$

$$G_\alpha(m) \rightarrow \tilde{G}_\alpha(m) = \mathbf{Ad}_{\exp \tau\kappa} (G_\alpha - \exp(-\tau\kappa) (\exp \tau\kappa)' \tau \partial_\alpha \kappa)$$

$$\dot{A}_\alpha \rightarrow \tilde{\dot{A}}_\alpha(m, \tau) = \mathbf{Ad}_{\exp \tau\theta} (\dot{A}_\alpha - \exp(-\tau\theta) \exp(\tau\theta)' \tau \partial_\alpha \theta)$$

$$\nabla_\alpha \psi \rightarrow \widetilde{\nabla_\alpha \psi}^{ij}(m, \tau) = \sum_{k=1}^4 \sum_{l=1}^n [\gamma C(\exp(\tau\kappa))]_k^i [\varrho(\exp(\tau\theta))]_l^j \nabla_\alpha \psi^{kl}$$

All these expressions depend on m , as well as $\kappa(m), \theta(m)$, so they can be differentiated with respect to the coordinates of m to get :

$$\partial_\beta P_i^\alpha \rightarrow \tilde{\partial}_\beta P_i^\alpha(m, \tau) = \sum_{j=0}^3 \left([h(\exp(-\tau\kappa))]' \partial_\beta \kappa \right)_i^j P_j^\alpha + [h(\exp(-\tau\kappa))]_i^j \partial_\beta P_i^\alpha$$

$$\partial_\beta \tilde{G}_\alpha(m, \tau)$$

$$= [(\exp -\tau\kappa) (\exp \tau\kappa)' \tau \partial_\beta \kappa, G_\alpha - \tau \partial_\alpha \kappa]$$

$$+ \mathbf{Ad}_{\exp \tau\kappa} \{ \partial_\beta G_\alpha -$$

$$\{ (\exp -\tau\kappa)' \tau \partial_\beta \kappa \circ (\exp \tau\kappa)' \tau \partial_\alpha \kappa$$

$$+ \exp(-\tau\kappa) \circ (\exp \tau\kappa)'' (\tau \partial_\beta \kappa, \tau \partial_\alpha \kappa) + \exp(-\tau\kappa) \circ \exp(\tau\kappa)' \tau \partial_{\alpha\beta}^2 \kappa \}$$

$$\partial_\beta \tilde{\dot{A}}_\alpha(m, \tau) = [(\exp -\tau\theta) (\exp \tau\theta)' \tau \partial_\beta \theta, \dot{A}_\alpha - \tau \partial_\alpha \theta]$$

$$+ \mathbf{Ad}_{\exp \tau\theta} (\partial_\beta \dot{A}_\alpha - (\exp(-\tau\theta))' \tau \partial_\beta \theta \circ (\exp \tau\theta)' \tau \partial_\alpha \theta)$$

$$+ \exp(-\tau\theta) \circ (\exp \tau\theta)'' (\tau \partial_\beta \theta, \tau \partial_\alpha \theta) + \exp(-\tau\theta) \circ \exp(\tau\theta)' \tau \partial_{\alpha\beta}^2 \theta)$$

$$\mathcal{F}_{G\alpha\beta}^a \rightarrow \tilde{\mathcal{F}}_{G\alpha\beta}(\tau) = \mathbf{Ad}_{\exp \tau\kappa} \mathcal{F}_{G\alpha\beta}$$

$$\mathcal{F}_{A\alpha\beta}^a \rightarrow \tilde{\mathcal{F}}_{A\alpha\beta}(\tau) = \mathbf{Ad}_{\exp \tau\theta} \mathcal{F}_{A\alpha\beta}$$

$$F_{G\alpha\beta} \rightarrow \mathbf{Ad}_{\exp \tau\kappa} F_{G\alpha\beta}$$

$$+ [(\exp -\tau\kappa) (\exp \tau\kappa)' \tau \partial_\beta \kappa, G_\alpha - t \partial_\alpha \kappa] + [(\exp -\tau\kappa) (\exp \tau\kappa)' \tau \partial_\alpha \kappa, G_\beta - \tau \partial_\beta \kappa]$$

$$- \mathbf{Ad}_{\exp \tau\kappa} ((\exp -\tau\kappa)' \tau \partial_\beta \kappa \circ (\exp \tau\kappa)' \tau \partial_\alpha \kappa + \exp(-\tau\kappa) \circ (\exp \tau\kappa)'' (\tau \partial_\beta \kappa, \tau \partial_\alpha \kappa)$$

$$+ \exp(-\tau\kappa) \circ \exp(\tau\kappa)' \tau \partial_{\alpha\beta}^2 \kappa)$$

$$\begin{aligned}
& -Ad_{\exp \tau \kappa}((\exp -\tau \kappa)' \tau \partial_\alpha \kappa \circ (\exp \tau \kappa)' \tau \partial_\beta \kappa + \exp(-\tau \kappa) \circ (\exp \tau \kappa)'' (\tau \partial_\alpha \kappa, \tau \partial_\beta \kappa) \\
& + \exp(-\tau \kappa) \circ \exp(\tau \kappa)' \tau \partial_{\alpha\beta}^2 \kappa) \\
& F_{A\alpha\beta} \rightarrow Ad_{\exp(\tau\theta) F_{A\alpha\beta}} \\
& + \left[(\exp -\tau\theta) (\exp \tau\theta)' \tau \partial_\beta \theta, \dot{A}_\alpha - \tau \partial_\alpha \theta \right] + \left[(\exp -\tau\theta) (\exp \tau\theta)' \tau \partial_\alpha \theta, \dot{A}_\beta - \tau \partial_\beta \theta \right] \\
& - Ad_{\exp \tau\theta}(\exp(-\tau\theta)' \tau \partial_\beta \theta \circ (\exp \tau\theta)' \tau \partial_\alpha \theta + \exp(-\tau\theta) \circ (\exp \tau\theta)'' (\tau \partial_\beta \theta, \tau \partial_\alpha \theta) \\
& + \exp(-\tau\theta) \circ \exp(\tau\theta)' \tau \partial_{\alpha\beta}^2 \theta) \\
& - Ad_{\exp \tau\theta}(\exp(-\tau\theta)' \tau \partial_\alpha \theta \circ (\exp \tau\theta)' \tau \partial_\beta \theta + \exp(-\tau\theta) \circ (\exp \tau\theta)'' (\tau \partial_\alpha \theta, \tau \partial_\beta \theta) \\
& + \exp(-\tau\theta) \circ \exp(\tau\theta)' \tau \partial_{\alpha\beta}^2 \theta)
\end{aligned}$$

The vector V is defined in the holonomic basis $\partial \xi_\alpha$ so its components are not impacted.

The determinant $\det P'$ is invariant, because we have a change of orthonormal basis, so the scalar lagrangian L is invariant :

$$\begin{aligned}
L & \left(\psi^{ij}, G_\alpha^a, \dot{A}_\alpha, P_i^\alpha, \nabla_\beta \psi^{ij}, \mathcal{F}_{G\alpha\beta}, F_{G\alpha\beta}, \mathcal{F}_{A\alpha\beta}, F_{A\alpha\beta}, \partial_\beta P_i^\alpha, V^\alpha \right) \\
& = L \left(\widetilde{\psi}^{ij}, \widetilde{G}_\alpha^a, \widetilde{\dot{A}}_\alpha, \widetilde{P}_i^\alpha, \widetilde{\nabla}_\beta \widetilde{\psi}^{ij}, \widetilde{\mathcal{F}}_{G\alpha\beta}, \widetilde{F}_{G\alpha\beta}, \widetilde{\mathcal{F}}_{A\alpha\beta}, \widetilde{F}_{A\alpha\beta}, \widetilde{\partial}_\beta \widetilde{P}_i^\alpha, V^\alpha \right)
\end{aligned}$$

If we take the derivative of this identity for $\tau = 0$ we get an identity which must hold whatever $\kappa(m), \theta(m)$ and then identities between the partial derivatives of \mathcal{L} . From a mathematical point of view this derivative with respect to τ is the Lie derivative of the lagrangian along the vertical vector fields generated by the derivative $\frac{dz^i}{d\tau}|_{\tau=0}$ for each variable. These vector fields are the **Noether currents** (Maths.34.3.4). Here we will not explicit these currents, but simply deduce some compatibilities between the partial derivatives.

Equivariance on P_G

The computation for $\exp(\tau \kappa(m))$ gives :

$$\begin{aligned}
& \frac{d}{d\tau} \widetilde{P}_i^\alpha(m, \tau)|_{\tau=0} = -\sum_a \kappa^a ([P][\kappa_a])_i^\alpha \\
& \frac{d}{d\tau} \operatorname{Re} \widetilde{\psi}^{ij}(m, \tau)|_{\tau=0} \\
& = \sum_a \kappa^a \sum_{k=1}^4 \left(\operatorname{Re} \left([\gamma C(\kappa_a)]_k^i \right) \operatorname{Re} \psi^{kj} - \operatorname{Im} \left([\gamma C(\kappa_a)]_k^i \right) \operatorname{Im} \psi^{kj} \right) \\
& = \sum_a \kappa^a \operatorname{Re} \left([\gamma C(\kappa_a)] [\psi] \right)_j^i \\
& \frac{d}{d\tau} \operatorname{Im} \widetilde{\psi}^{ij}(m, \tau)|_{\tau=0} \\
& = \sum_a \kappa^a \sum_{k=1}^4 \left(\operatorname{Re} \left([\gamma C(\kappa_a)]_k^i \right) \operatorname{Im} \psi^{kj} + \operatorname{Im} \left([\gamma C(\kappa_a)]_k^i \right) \operatorname{Re} \psi^{kj} \right) \\
& = \sum_a \kappa^a \operatorname{Im} \left([\gamma C(\kappa_a)] [\psi] \right)_j^i \\
& \frac{d}{d\tau} \partial_\beta \widetilde{P}(m, t)_j^\alpha|_{\tau=0} = -\sum_a \kappa^a ([\partial_\beta P][\kappa_a])_j^\alpha + \partial_\beta \kappa^a ([P][\kappa_a])_j^\alpha \\
& \frac{d}{d\tau} \operatorname{Re} \widetilde{\nabla}_\alpha \psi^{ij}(m, \tau)|_{\tau=0} \\
& = \sum_a \kappa^a \sum_{k=1}^4 \left(\operatorname{Re} \left([\gamma C(\kappa_a)]_k^i \right) \operatorname{Re} \nabla_\alpha \psi^{kj} - \operatorname{Im} \left([\gamma C(\kappa_a)]_k^i \right) \operatorname{Im} \nabla_\alpha \psi^{kj} \right) \\
& = \sum_a \kappa^a \operatorname{Re} \left([\gamma C(\kappa_a)] [\nabla_\alpha \psi] \right)_j^i \\
& \frac{d}{d\tau} \operatorname{Im} \widetilde{\nabla}_\alpha \psi^{ij}(m, \tau)|_{\tau=0} \\
& = \sum_a \kappa^a \sum_{k=1}^4 \left(\operatorname{Re} \left([\gamma C(\kappa_a)]_k^i \right) \operatorname{Im} \nabla_\alpha \psi^{kj} + \operatorname{Im} \left([\gamma C(\kappa_a)]_k^i \right) \operatorname{Re} \nabla_\alpha \psi^{kj} \right)
\end{aligned}$$

$$\begin{aligned}
&= \sum_a \kappa^a \operatorname{Im}([\gamma C(\kappa_a)] [\nabla_\alpha \psi])_j^i \\
&\frac{d}{d\tau} \tilde{G}_\alpha^a(m) |_{\tau=0} = \sum_b \kappa^b [\vec{\kappa}_b, G_\alpha]^a - \partial_\alpha \kappa^a \\
&\frac{d}{d\tau} \partial_\beta \tilde{G}_\alpha^a(m, \tau) |_{\tau=0} = \sum_b \kappa^b [\vec{\kappa}_b, \partial_\beta G_\alpha]^a + \partial_\beta \kappa^b [\vec{\kappa}_b, G_\alpha]^a - \partial_{\alpha\beta} \kappa^a \\
&\frac{d}{d\tau} \tilde{\mathcal{F}}_{G\alpha\beta}^a(\tau) |_{\tau=0} = \sum_b \kappa^b [\vec{\kappa}_b, \mathcal{F}_{G\alpha\beta}]^a \\
&\frac{d}{d\tau} \tilde{F}_{G\alpha\beta}^a |_{\tau=0} = \sum_b \kappa^b [\vec{\kappa}_b, F_{G\alpha\beta}]^a + \partial_\beta \kappa^b [\vec{\kappa}_b, G_\alpha]^a + \partial_\alpha \kappa^b [\vec{\kappa}_b, G_\beta]^a - 2\partial_{\alpha\beta} \kappa^a
\end{aligned}$$

So we have the identity :

$$\forall \kappa_a, \partial_\beta \kappa^a, \partial_{\alpha\beta} \kappa^a :$$

$$\begin{aligned}
0 &= \\
&\sum_a \kappa^a \left\{ \sum_{ij} \frac{\partial L}{\partial \operatorname{Re} \psi^{i\bar{j}}} \operatorname{Re}([\gamma C(\kappa_a)] [\psi])_j^i + \frac{\partial L}{\partial \operatorname{Im} \psi^{i\bar{j}}} \operatorname{Im}([\gamma C(\kappa_a)] [\psi])_j^i \right. \\
&+ \sum_{\alpha ij} \frac{\partial L}{\partial \operatorname{Re} \nabla_\alpha \psi^{i\bar{j}}} \operatorname{Re}([\gamma C(\kappa_a)] [\nabla_\alpha \psi])_j^i + \frac{\partial L}{\partial \operatorname{Im} \nabla_\alpha \psi^{i\bar{j}}} \operatorname{Im}([\gamma C(\kappa_a)] [\nabla_\alpha \psi])_j^i \left. \right\} \\
&+ \sum_{i\alpha} \frac{\partial L}{\partial P_i^\alpha} (-\sum_a \kappa^a ([P] [\kappa_a])_i^\alpha) \\
&+ \sum_{i\alpha\beta} \frac{\partial \mathcal{L}}{\partial \partial_\beta P_i^\alpha} \left(-\sum_a \kappa^a ([\partial_\beta P] [\kappa_a])_j^\alpha + \partial_\beta \kappa^a ([P] [\kappa_a])_j^\alpha \right) \\
&+ \sum_{a\alpha} \frac{\partial L}{\partial G_\alpha^a} (\sum_b \kappa^b [\vec{\kappa}_b, G_\alpha]^a - \partial_\alpha \kappa^a) \\
&+ \sum_{a\alpha\beta} \frac{\partial L}{\partial \mathcal{F}_{G\alpha\beta}^a} (\sum_b \kappa^b [\vec{\kappa}_b, \mathcal{F}_{G\alpha\beta}]^a) \\
&+ \frac{\partial L}{\partial F_{G\alpha\beta}^a} (\sum_b \kappa^b [\vec{\kappa}_b, F_{G\alpha\beta}]^a + \partial_\beta \kappa^b [\vec{\kappa}_b, G_\alpha]^a + \partial_\alpha \kappa^b [\vec{\kappa}_b, G_\beta]^a - 2\partial_{\alpha\beta} \kappa^a)
\end{aligned}$$

With the component in $\partial_{\alpha\beta} \kappa^a$ we have immediately : $\forall a, \alpha, \beta : \frac{\partial L}{\partial F_{G\alpha\beta}^a} = 0$

With the component in $\partial_\alpha \kappa^a : \forall a, \alpha : \sum_{\beta i} \frac{\partial L}{\partial \partial_\alpha P_i^\beta} ([P] [\kappa_a])_i^\beta = -\frac{\partial L}{\partial G_\alpha^a}$

And we are left with :

$$\forall a = 1..6 :$$

$$\begin{aligned}
0 &= \\
&\sum_{ij} \frac{\partial L}{\partial \psi^{i\bar{j}}} ([\gamma C(\kappa_a)] [\psi])_j^i + \sum_{\alpha ij} \frac{\partial L}{\partial \nabla_\alpha \psi^{i\bar{j}}} ([\gamma C(\kappa_a)] [\nabla_\alpha \psi])_j^i \\
&- \sum_{i\alpha} \frac{\partial L}{\partial P_i^\alpha} ([P] [\kappa_a])_i^\alpha - \sum_{i\alpha\beta} \frac{\partial \mathcal{L}}{\partial \partial_\beta P_i^\alpha} ([\partial_\beta P] [\kappa_a])_j^\alpha \\
&+ \sum_{b\alpha} \frac{\partial L}{\partial G_\alpha^b} [\vec{\kappa}_a, G_\alpha]^b + \sum_{a\alpha\beta} \frac{\partial L}{\partial \mathcal{F}_{G\alpha\beta}^b} [\vec{\kappa}_a, \mathcal{F}_{G\alpha\beta}]^b
\end{aligned}$$

Moreover, by taking the derivative with respect to the initial variables we get :

$$\begin{aligned}
&\sum_{k=1}^4 [\gamma C(\exp(\tau\kappa(m)))_i^k] \frac{\partial \tilde{L}}{\partial \psi^{k\bar{j}}} = \frac{\partial L}{\partial \psi^{i\bar{j}}} \\
&\sum_{k=1}^4 [\gamma C(\exp(\tau\kappa(m)))_i^k] \frac{\partial \tilde{L}}{\partial \nabla_\alpha \psi^{k\bar{j}}} = \frac{\partial L}{\partial \nabla_\alpha \psi^{i\bar{j}}} \\
&\sum_j [h(\exp(-\tau\kappa(m)))_i^j] \frac{\partial \tilde{L}}{\partial P_j^\alpha} = \frac{\partial L}{\partial P_i^\alpha} \\
&L([Ad_{\exp \tau\kappa}]_b^a \mathcal{F}_{G\alpha\beta}^b) = L(\mathcal{F}_{G\alpha\beta}) \\
&\sum_b [Ad_{\exp \tau\kappa}]_a^b \frac{\partial \tilde{L}}{\partial \mathcal{F}_{G\alpha\beta}^b} = \frac{\partial L}{\partial \mathcal{F}_{G\alpha\beta}^a}
\end{aligned}$$

and other similar identities, which show that the partial derivatives are tensors, with respect to the vector bundles :

$$\sum_i \frac{\partial L}{\partial \psi^{i\bar{j}}} \mathbf{e}^i, \sum_i \frac{\partial L}{\partial \nabla_\alpha \psi^{i\bar{j}}} \mathbf{e}^i, \sum_i \frac{\partial L}{\partial P_i^\alpha} \varepsilon_i, \sum_i \frac{\partial L}{\partial \partial_\beta P_i^\alpha} \varepsilon_i, \sum_a \frac{\partial L}{\partial \mathcal{F}_{G\alpha\beta}^a} \vec{\kappa}^a$$

Equivariance on P_U

We have similarly :

$$\begin{aligned}
\frac{d}{d\tau} \widetilde{\psi}^{ij} (m, \tau) |_{\tau=0} &= \sum_{k=1}^n \sum_{a=1}^m \theta^a \psi^{ik} [\theta_a]_j^k \\
\frac{d}{d\tau} \operatorname{Re} \widetilde{\psi}^{ij} (m, \tau) |_{\tau=0} &= \sum_{a=1}^m \theta^a \operatorname{Re} (\psi [\theta_a])^{ij} \\
\frac{d}{d\tau} \operatorname{Im} \widetilde{\psi}^{ij} (m, \tau) |_{\tau=0} &= \sum_{a=1}^m \theta^a \operatorname{Im} (\psi [\theta_a])^{ij} \\
\frac{d}{d\tau} \widetilde{A}_\alpha^a (m, \tau) |_{\tau=0} &= \sum_{b=1}^m \theta^b \left[\vec{\theta}_b, \dot{A}_\alpha \right]^a - \partial_\alpha \theta^a \\
\frac{d}{d\tau} \operatorname{Re} \widetilde{\nabla_\alpha \psi}^{ij} (m, \tau) |_{\tau=0} &= \sum_{a=1}^m \theta^a \operatorname{Re} (\nabla_\alpha \psi [\theta_a])^{ij} \\
\frac{d}{d\tau} \operatorname{Im} \widetilde{\nabla_\alpha \psi}^{ij} (m, \tau) |_{\tau=0} &= \sum_{a=1}^m \theta^a \operatorname{Im} (\nabla_\alpha \psi [\theta_a])^{ij} \\
\frac{d}{d\tau} \partial_\beta \widetilde{A}_\alpha^a (m, \tau) |_{\tau=0} &= \sum_{b=1}^m \theta^b \left[\theta_b, \partial_\beta \dot{A}_\alpha \right]^a + \partial_\beta \theta^b \left[\theta_b, \dot{A}_\alpha \right]^a - \partial_{\alpha\beta} \theta^a \\
\frac{d}{d\tau} \widetilde{\mathcal{F}}_{A\alpha\beta} (\tau) |_{\tau=0} &= \sum_{b=1}^m \theta^b \left[\vec{\theta}_b, \mathcal{F}_{A\alpha\beta} \right]^a \\
\frac{d}{d\tau} \widetilde{F}_{A\alpha\beta} |_{\tau=0} &= \sum_{b=1}^m \theta^b \left[\theta_b, F_{A\alpha\beta} \right]^a + \partial_\beta \theta^b \left[\theta_b, \dot{A}_\alpha \right]^a + \partial_\alpha \theta^b \left[\theta_b, \dot{A}_\beta \right]^a - 2\partial_{\alpha\beta} \theta^a \\
\sum_{ij} \frac{\partial L}{\partial \operatorname{Re} \psi^{ij}} \sum_{a=1}^m \theta^a \operatorname{Re} (\psi [\theta_a])^{ij} &+ \frac{\partial L}{\partial \operatorname{Im} \psi^{ij}} \sum_{a=1}^m \theta^a \operatorname{Im} (\psi [\theta_a])^{ij} \\
+ \sum_{ij\alpha} \frac{\partial L}{\partial \operatorname{Re} \nabla_\alpha \psi^{ij}} \sum_{a=1}^m \theta^a \operatorname{Re} (\psi [\theta_a])^{ij} &+ \sum_{ij\alpha} \frac{\partial L}{\partial \operatorname{Im} \nabla_\alpha \psi^{ij}} \sum_{a=1}^m \theta^a \operatorname{Im} (\psi [\theta_a])^{ij} \\
+ \sum_{a\alpha} \frac{\partial L}{\partial \dot{A}_\alpha^a} \left(\sum_{b=1}^m \theta^b \left[\vec{\theta}_b, \dot{A}_\alpha \right]^a - \partial_\alpha \theta^a \right) &+ \sum_{a\alpha\beta} \frac{\partial L}{\partial \mathcal{F}_{A\alpha\beta}^a} \left(\sum_{b=1}^m \theta^b \left[\vec{\theta}_b, \mathcal{F}_{A\alpha\beta} \right]^a \right) \\
+ \sum_{a\alpha\beta} \frac{\partial L}{\partial F_{A\alpha\beta}^a} \left(\sum_{b=1}^m \theta^b \left[\theta_b, F_{A\alpha\beta} \right]^a + \partial_\beta \theta^b \left[\theta_b, \dot{A}_\alpha \right]^a \right. &+ \left. \partial_\alpha \theta^b \left[\theta_b, \dot{A}_\beta \right]^a - 2\partial_{\alpha\beta} \theta^a \right) \\
&= 0
\end{aligned}$$

Which implies :

$$\forall a, \alpha, \beta : \frac{\partial L}{\partial F_{A\alpha\beta}^a} = 0, \frac{\partial L}{\partial \dot{A}_\alpha^a} = 0$$

$\forall a = 1..m :$

$$\sum_{ij} \frac{\partial L}{\partial \psi^{ij}} (\psi [\theta_a])^{ij} + \sum_{ij\alpha} \frac{\partial L}{\partial \nabla_\alpha \psi^{ij}} (\nabla_\alpha \psi [\theta_a])^{ij} + \sum_{b\alpha\beta} \frac{\partial L}{\partial \mathcal{F}_{A\alpha\beta}^b} \left(\left[\vec{\theta}_a, \mathcal{F}_{A\alpha\beta} \right]^b \right) =$$

0

By taking the derivative with respect to the initial variables we check that the partial derivatives are tensors, with respect to the vector bundles : $\sum_i \frac{\partial L}{\partial \psi^{ij}} \mathbf{f}^j$, $\sum_i \frac{\partial L}{\partial \nabla_\alpha \psi^{ij}} \mathbf{f}^j$, $\sum_a \frac{\partial L}{\partial \mathcal{F}_{A\alpha\beta}^a} \vec{\theta}^a$ with $\vec{\theta}^a$ the basis vector of the dual of T_1U : $\vec{\theta}^a \left(\vec{\theta}_b \right) = \delta_b^a$

6.1.4 Covariance

In a change of charts on M with the jacobian : $J = \left[J_\beta^\alpha \right] = \left[\frac{\partial \tilde{\xi}^\alpha}{\partial \xi^\beta} \right]$ and $K = J^{-1}$

the 4-form on M which defines the action changes as :

$$L\mu \det [P] d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3 = \tilde{L}\tilde{\mu} \det [\tilde{P}] d\tilde{\xi}^0 \wedge d\tilde{\xi}^1 \wedge d\tilde{\xi}^2 \wedge d\tilde{\xi}^3$$

and because :

$$\tilde{\mu} \det [\tilde{P}] d\tilde{\xi}^0 \wedge d\tilde{\xi}^1 \wedge d\tilde{\xi}^2 \wedge d\tilde{\xi}^3 = \mu \det [P] d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3$$

the scalar lagrangian L should be invariant.

The variables change as :

ψ^{ij} do not change

The covariant derivatives are one form :

$$\nabla_\alpha \psi^{ij} \rightarrow \widetilde{\nabla}_\alpha \psi^{ij} = \sum_\beta K_\alpha^\beta \nabla_\beta \psi^{ij}$$

P, V are vectors, but their components are functions :

$$V^\alpha \rightarrow \widetilde{V}^\alpha = \sum_\gamma J_\gamma^\alpha u^\gamma$$

$$P_i^\alpha \rightarrow \widetilde{P}_i^\alpha = \sum_\gamma J_\gamma^\alpha P_i^\gamma$$

$$\widetilde{\partial}_\beta P_i^\alpha = \frac{\partial}{\partial \xi^\beta} \left(\sum_\gamma J_\gamma^\alpha (\xi) P_i^\gamma (\xi) \right) = \sum_\gamma \left(\frac{\partial}{\partial \xi^\beta} J_\gamma^\alpha (\xi) \right) P_i^\gamma (\xi) + J_\gamma^\alpha (\xi) \frac{\partial}{\partial \xi^\beta} P_i^\gamma (\xi)$$

$$\widetilde{\partial}_\beta P_i^\alpha = \sum_{\gamma\eta} (\partial_\eta J_\gamma^\alpha) K_\beta^\eta P_i^\gamma + \left((\partial_\eta P_i^\gamma) J_\gamma^\alpha K_\beta^\eta \right)$$

The potentials are 1-form :

$$G_\alpha^a \rightarrow \widetilde{G}_\alpha^a = \sum_\beta K_\alpha^\beta G_\beta^a$$

$$\dot{A}_\alpha^a \rightarrow \widetilde{\dot{A}}_\alpha^a = \sum_\beta K_\alpha^\beta \dot{A}_\beta^a$$

The strengths of the fields are 2-forms :

$$\mathcal{F}_{G\alpha\beta}^a \rightarrow \widetilde{\mathcal{F}}_{G\alpha\beta}^a = \sum_{\gamma\eta} K_\alpha^\gamma K_\beta^\eta \mathcal{F}_{G\gamma\eta}^a$$

$$\mathcal{F}_{A\alpha\beta}^a \rightarrow \widetilde{\mathcal{F}}_{A\alpha\beta}^a = \sum_{\gamma\eta} K_\alpha^\gamma K_\beta^\eta \mathcal{F}_{A\gamma\eta}^a$$

So we have the identity, for any change of chart :

$$\begin{aligned} & L \left(P_i^\alpha, \psi^{ij}, G_\alpha^a, \partial_\beta P_i^\alpha, \nabla_\alpha \psi^{ij}, \mathcal{F}_{A\alpha\beta}^a, \mathcal{F}_{G\alpha\beta}^a, V^\alpha \right) \\ &= \widetilde{L} \left(\widetilde{P}_i^\alpha, \widetilde{\psi}^{ij}, \widetilde{G}_\alpha^a, \widetilde{\partial}_\beta P_i^\alpha, \widetilde{\nabla}_\alpha \psi^{ij}, \widetilde{\mathcal{F}}_{A\alpha\beta}^a, \widetilde{\mathcal{F}}_{G\alpha\beta}^a, \widetilde{V}^\alpha \right) \\ &= L \left(\sum_\gamma J_\gamma^\alpha P_i^\gamma, \psi^{ij}, \sum_\beta K_\alpha^\beta G_\beta^a, \sum_{\gamma\eta} (\partial_\eta J_\gamma^\alpha) K_\beta^\eta P_i^\gamma + \left((\partial_\eta P_i^\gamma) J_\gamma^\alpha K_\beta^\eta \right), \right. \\ & \quad \left. \sum_\beta K_\alpha^\beta \nabla_\beta \psi^{ij}, \sum_{\gamma\eta} K_\alpha^\gamma K_\beta^\eta \mathcal{F}_{G\gamma\eta}^a, \sum_{\gamma\eta} K_\alpha^\gamma K_\beta^\eta \mathcal{F}_{A\gamma\eta}^a, \sum_\gamma J_\gamma^\alpha V^\gamma \right) \end{aligned}$$

If we take the derivative of this identity with respect to $(\partial_\eta J_\mu^\lambda)$:

$$0 = \sum_{i\alpha\beta} \frac{\partial L}{\partial \partial_\beta P_i^\alpha} \sum_{\gamma\eta} K_\beta^\eta P_i^\gamma \delta_\alpha^\lambda \delta_\gamma^\mu = \sum_{\alpha\beta\eta i} \frac{\partial L}{\partial \partial_\beta P_i^\alpha} P_i^\mu K_\beta^\eta$$

$$\text{take } J_\mu^\lambda = \delta_\mu^\lambda \Rightarrow K_\beta^\eta = \delta_\beta^\eta$$

$$\sum_i \frac{\partial L}{\partial \partial_\eta P_i^\lambda} P_i^\mu = 0$$

$$\forall \alpha, \beta, \gamma : \sum_i \frac{\partial L}{\partial \partial_\alpha P_i^\beta} P_i^\gamma = 0$$

$$\text{by product with } P_\gamma^{j\prime} \text{ and summation : } \forall \alpha, \beta, j : \frac{\partial L}{\partial \partial_\alpha P_j^\beta} = 0$$

and as we had :

$$\forall a, \alpha : \sum_{\beta i} \frac{\partial L}{\partial \partial_\alpha P_i^\beta} ([P] [\kappa_a])_i^\beta = -\frac{\partial L}{\partial G_\alpha^a} \Rightarrow \forall a, \alpha : \frac{\partial L}{\partial G_\alpha^a} = 0$$

The derivative with respect to J_μ^λ :

$$\begin{aligned} & \sum_{i\alpha} \frac{\partial L}{\partial P_i^\alpha} \sum_\gamma P_i^\gamma \delta_\alpha^\lambda \delta_\gamma^\mu + \sum_{i\alpha} \frac{\partial L}{\partial \text{Re} \nabla_\alpha \psi^{i\gamma}} \sum_\beta \left(\frac{\partial}{\partial J_\mu^\lambda} K_\alpha^\beta \right) \text{Re} \nabla_\beta \psi^{ij} \\ &+ \sum_{i\alpha} \frac{\partial L}{\partial \text{Im} \nabla_\alpha \psi^{i\gamma}} \sum_\beta \left(\frac{\partial}{\partial J_\mu^\lambda} K_\alpha^\beta \right) \text{Im} \nabla_\beta \psi^{ij} \\ &+ \sum_{a\alpha\beta} \frac{\partial L}{\partial \mathcal{F}_{G\alpha\beta}^a} \sum_{\gamma\eta} \left(\left(\frac{\partial}{\partial J_\mu^\lambda} K_\alpha^\gamma \right) K_\beta^\eta + K_\alpha^\gamma \frac{\partial}{\partial J_\mu^\lambda} K_\beta^\eta \right) \mathcal{F}_{G\gamma\eta}^a \end{aligned}$$

$$\begin{aligned}
& + \frac{\partial L}{\partial \mathcal{F}_{A\alpha\beta}^a} \sum_{\gamma\eta} \left(\left(\frac{\partial}{\partial J_\mu^\lambda} K_\alpha^\gamma \right) K_\beta^\eta + K_\alpha^\gamma \frac{\partial}{\partial J_\mu^\lambda} K_\beta^\eta \right) \mathcal{F}_{A\gamma\eta}^a + \frac{\partial L}{\partial V^\alpha} \sum_\gamma V^\gamma \delta_\alpha^\lambda \delta_\gamma^\mu = 0 \\
& \text{with } \frac{\partial}{\partial J_\mu^\lambda} K_\alpha^\beta = -K_\lambda^\beta K_\alpha^\mu \\
& \sum_{i\alpha} \frac{\partial L}{\partial P_i^\alpha} P_i^\mu + \sum_{ij\alpha} \frac{\partial L}{\partial \nabla_\alpha \psi^{ij}} \sum_\beta \left(-K_\lambda^\beta K_\alpha^\mu \right) \nabla_\beta \psi^{ij} \\
& + \sum_{a\alpha\beta} \frac{\partial L}{\partial \mathcal{F}_{G\alpha\beta}^a} \sum_{\gamma\eta} \left(\left((-K_\lambda^\gamma K_\alpha^\mu) \right) K_\beta^\eta + K_\alpha^\gamma \left(-K_\lambda^\eta K_\beta^\mu \right) \right) \mathcal{F}_{G\gamma\eta}^a \\
& + \frac{\partial L}{\partial \mathcal{F}_{A\alpha\beta}^a} \sum_{\gamma\eta} \left(\left((-K_\lambda^\gamma K_\alpha^\mu) \right) K_\beta^\eta + K_\alpha^\gamma \left(-K_\lambda^\eta K_\beta^\mu \right) \right) \mathcal{F}_{A\gamma\eta}^a + \frac{\partial L}{\partial V^\lambda} V^\mu = 0 \\
& \text{Let us take } J_\mu^\lambda = \delta_\mu^\lambda \Rightarrow K_\mu^\lambda = \delta_\mu^\lambda \\
& \sum_i \frac{\partial L}{\partial P_i^\lambda} P_i^\mu - \sum_{i\alpha} \frac{\partial L}{\partial \nabla_\mu \psi^{i\alpha}} \nabla_\lambda \psi^{ij} - \sum_{a\eta} \frac{\partial L}{\partial \mathcal{F}_{G\mu\eta}^a} \mathcal{F}_{G\lambda\eta}^a - \sum_{a\gamma} \frac{\partial L}{\partial \mathcal{F}_{G\gamma\mu}^a} \mathcal{F}_{G\gamma\lambda}^a - \sum_{a\eta} \frac{\partial L}{\partial \mathcal{F}_{A\mu\eta}^a} \mathcal{F}_{A\lambda\eta}^a - \\
& \sum_{a\gamma} \frac{\partial L}{\partial \mathcal{F}_{A\gamma\mu}^a} \mathcal{F}_{A\gamma\lambda}^a + \frac{\partial L}{\partial V^\lambda} V^\mu = 0 \\
& \text{that is :} \\
& \forall \alpha, \beta : \sum_{ij} \frac{\partial L}{\partial \nabla_\beta \psi^{ij}} \nabla_\alpha \psi^{ij} + 2 \sum_{a\gamma} \frac{\partial L}{\partial \mathcal{F}_{G\beta\gamma}^a} \mathcal{F}_{G\alpha\gamma}^a + \frac{\partial L}{\partial \mathcal{F}_{A\beta\gamma}^a} \mathcal{F}_{A\alpha\gamma}^a = \sum_i \frac{\partial L}{\partial P_i^\alpha} P_i^\beta + \\
& \frac{\partial L}{\partial V^\alpha} V^\beta
\end{aligned}$$

The derivative with respect to the initial variables gives the identities :

$$\begin{aligned}
\frac{\partial L}{\partial P_i^\alpha} &= \sum_\beta J_\alpha^\beta \frac{\partial L}{\partial \tilde{P}_i^\beta} \\
\frac{\partial L}{\partial u^\alpha} &= \sum_\beta J_\alpha^\beta \frac{\partial L}{\partial \tilde{u}^\beta} \\
\frac{\partial L}{\partial \nabla_\beta \psi^{ij}} &= \sum_\gamma K_\gamma^\beta \frac{\partial \tilde{L}}{\partial \nabla_\gamma \psi^{ij}}; \\
\frac{\partial L}{\partial \mathcal{F}_{A\alpha\beta}^a} &= \sum_{\gamma\eta} K_\gamma^\alpha K_\eta^\beta \frac{\partial L}{\partial \mathcal{F}_{A\gamma\eta}^a}; \quad \frac{\partial L}{\partial \mathcal{F}_{G\alpha\beta}^a} = \sum_{\gamma\eta} K_\gamma^\alpha K_\eta^\beta \frac{\partial L}{\partial \mathcal{F}_{G\gamma\eta}^a}
\end{aligned}$$

which show that the corresponding quantities are tensors.

6.1.5 Conclusion

i) The potentials \dot{A}, G , and the derivatives $\partial_\beta P_i^\alpha$ do not figure explicitly, the derivatives of the potential \dot{A}, G factor in the strength.

The lagrangian is a function of 6 variables only :

$$L = L(\psi, \nabla_\alpha \psi, P_i^\alpha, \mathcal{F}_{G\alpha\beta}, \mathcal{F}_{A\alpha\beta}, V^\alpha) \quad (6.4)$$

ii) The following quantities are tensors :

$$\begin{aligned}
\Pi_\psi &= \sum_{ij} \frac{\partial L}{\partial \psi^{ij}} \mathbf{e}^i \otimes \mathbf{f}^j \\
\Pi_\nabla &= \sum_\alpha \frac{\partial L}{\partial \nabla_\alpha \psi^{ij}} \partial \xi_\alpha \otimes \mathbf{e}^i \otimes \mathbf{f}^j \\
\Pi_P &= \sum_\alpha \frac{\partial L}{\partial P_i^\alpha} d\xi^\alpha \otimes \varepsilon_i; \\
\Pi_A &= \sum_{\alpha\beta} \frac{\partial L}{\partial \mathcal{F}_{A\alpha\beta}^a} \partial \xi_\alpha \wedge \partial \xi_\beta \otimes \vec{\theta}^a; \\
\Pi_G &= \sum_{\alpha\beta} \frac{\partial L}{\partial \mathcal{F}_{G\alpha\beta}^a} \partial \xi_\alpha \wedge \partial \xi_\beta \otimes \vec{\kappa}^a
\end{aligned}$$

and similarly $\sum_{\alpha\beta} v^* \left(\frac{\partial L}{\partial \mathcal{F}_{r\alpha\beta}}, \frac{\partial L}{\partial \mathcal{F}_{w\alpha\beta}} \right) \partial \xi_\alpha \wedge \partial \xi_\beta$

This result is useful because we can express cumbersome equations with simpler and more flexible geometric quantities. Notice that these quantities, when $\det[P']$ is added to L, are no longer covariant.

iii) We have the identities

$$\forall a = 1..6 : \Pi_\psi [\gamma C (\vec{\kappa}_a)] \psi + \sum_\alpha \Pi_{\nabla}^\alpha [\gamma C (\vec{\kappa}_a)] \nabla_\alpha \psi - \Pi_P [P] [\kappa_a] + \sum_{b\alpha\beta} \Pi_{Gb}^{\alpha\beta} [\vec{\kappa}_a, \mathcal{F}_{G\alpha\beta}]^b = 0$$

$$\forall a = 1..m : (\Pi_\psi \psi + \sum_\alpha \Pi_{\nabla}^\alpha \nabla_\alpha \psi) [\theta_a] + \sum_{b\alpha\beta} \Pi_{Ab}^{\alpha\beta} [\vec{\theta}_a, \mathcal{F}_{A\alpha\beta}]^b = 0$$

$$\forall \alpha, \beta : \Pi_{\nabla}^\beta \nabla_\alpha \psi + 2 \sum_{a\gamma} \Pi_{Ga}^{\beta\gamma} \mathcal{F}_{G\alpha\gamma}^a + \Pi_{Aa}^{\beta\gamma} \mathcal{F}_{A\alpha\gamma}^a - \sum_i \Pi_{P\alpha}^i P_i^\beta = \frac{\partial L}{\partial V^\alpha} V^\beta$$

These identities are minimal necessary conditions for the lagrangian : the calculations could be continued to higher derivatives. They do not depend on the signature. Whenever the lagrangian is expressed with the geometrical quantities, these identities are automatically satisfied.

The last identity, by multiplication with $P_i^{\beta k}$ and summation gives :

$$\forall \alpha, i : \Pi_{P\alpha}^i = \sum_{\beta pq} \Pi_{\nabla pq}^\beta P_\beta^i (\nabla_\alpha \psi)_q^p + 2 \sum_{a\gamma} \Pi_{Ga}^{\beta\gamma} P_\beta^i \mathcal{F}_{G\alpha\gamma}^a + \Pi_{Aa}^{\beta\gamma} P_\beta^i \mathcal{F}_{A\alpha\gamma}^a - \frac{\partial L}{\partial V^\alpha} V^\beta P_\beta^i \quad (6.5)$$

with the first identity we get :

$$\sum_{i\alpha} \Pi_{P\alpha}^i ([P] [\kappa_a])_i^\alpha = \sum_{ij} \Pi_{\psi ij} ([\gamma C (\kappa_a)] [\psi])_j^i + \sum_{\alpha ij} \Pi_{\nabla ij}^\alpha ([\gamma C (\kappa_a)] [\nabla_\alpha \psi])_j^i + \sum_{b\alpha\beta} \Pi_{Gb}^{\alpha\beta} [\vec{\kappa}_a, \mathcal{F}_{G\alpha\beta}]^b$$

$\forall a = 1..6 :$

$$\begin{aligned} & \sum_{ij} \Pi_{\psi ij} ([\gamma C (\kappa_a)] [\psi])_j^i + \sum_{\alpha ij} \Pi_{\nabla ij}^\alpha \left([\gamma C (\kappa_a)] [\nabla_\alpha \psi] - \sum_\beta ([P] [\kappa_a] [P'])_\alpha^\beta (\nabla_\beta \psi) \right)_j^i \\ & + \sum_{b\alpha\beta} \Pi_{Gb}^{\alpha\beta} \left([\vec{\kappa}_a, \mathcal{F}_{G\alpha\beta}]^b - 2 ([P] [\kappa_a] [P'])_\alpha^\gamma \mathcal{F}_{G\gamma\beta}^b \right) - 2 \Pi_{Ab}^{\alpha\beta} ([P] [\kappa_a] [P'])_\alpha^\gamma \mathcal{F}_{A\gamma\beta}^b = \\ & \sum_{i\alpha\beta} \frac{\partial L}{\partial V^\alpha} V^\beta ([P] [\kappa_a] [P'])_\beta^\alpha \end{aligned}$$

and with the Riemann tensor $[R_{\alpha\beta}] = [P] [\mathcal{F}_{G\alpha\beta}] [P'] = \sum_{a=1}^6 \mathcal{F}_{G\alpha\beta}^a [P] [\kappa_a] [P']$

$\forall \lambda, \mu = 0..3 :$

$$\begin{aligned} & \sum_{ij} \Pi_{\psi ij} ([\gamma C (\mathcal{F}_{G\lambda\mu})] [\psi])_j^i + \sum_{\alpha ij} \Pi_{\nabla ij}^\alpha \left([\gamma C (\mathcal{F}_{G\lambda\mu})] [\nabla_\alpha \psi] - \sum_\beta [R_{\lambda\mu}]_\alpha^\beta (\nabla_\beta \psi) \right)_j^i \\ & + \sum_{b\alpha\beta} \Pi_{Gb}^{\alpha\beta} \left([\mathcal{F}_{G\lambda\mu}, \mathcal{F}_{G\alpha\beta}]^b - 2 [R_{\lambda\mu}]_\alpha^\gamma \mathcal{F}_{G\gamma\beta}^b \right) - 2 \Pi_{Ab}^{\alpha\beta} [R_{\lambda\mu}]_\alpha^\gamma \mathcal{F}_{A\gamma\beta}^b = \sum_{\alpha\beta} \frac{\partial L}{\partial V^\alpha} V^\beta [R_{\lambda\mu}]_\beta^\alpha \end{aligned}$$

6.2 THE POINT PARTICLE ISSUE

A lagrangian must suit the case of particles alone, fields alone and interacting fields and particles. So it comprises a part for the fields, and another one for the particles and their interactions. So if we consider a population of N particles interacting with the fields the action is :

$$\int_{\Omega} L_1(P_i^\alpha, \mathcal{F}_{G\alpha\beta}, \mathcal{F}_{A\alpha\beta}) \varpi_4 + \sum_{p=1}^N \int_{\tau_p^1}^{\tau_p^2} L_2(\psi_p, \nabla_\alpha \psi_p, P_i^\alpha, V_p^\alpha) cd\tau_p$$

because the induced volume form on the world line is $cd\tau_p$

And this raises several issues, mathematical and physical, depending on the system considered.

6.2.1 Propagation of Fields

If we consider a system without any particle, focus on the fields and aim at knowing their propagation in Ω , the variables are just the components of the tetrad P , and the strength of the fields $\mathcal{F}_A, \mathcal{F}_G$. We have a unique integral over Ω and the Euler-Lagrange equations give general solutions which are matched to the initial conditions. A direct and simple answer can be found, at least for the variation with respect to t . In the usual models the propagation at the speed of light is postulated, and introduced separately in a linearized approximation. The classic examples are, in General Relativity (with the Levi-Civita connection) the Einstein equation :

$$Ric_{\alpha\beta} - \frac{1}{2}g_{\alpha\beta}(R + \Lambda) = 0$$

and the Maxwell equations :

$$\sum_{\alpha\beta} \partial_\alpha (\mathcal{F}^{\alpha\beta} \sqrt{|\det P^r|}) = 0$$

with the lagrangian : $L = \sum_{\alpha\beta} Gg^{\alpha\beta} Ric_{\alpha\beta} + \mu_0 \mathcal{F}_{\alpha\beta} \mathcal{F}^{\alpha\beta}$

6.2.2 Particles moving in known Fields

When the system is comprised of particles moving in known fields actually only the second part of the action is involved. Replacing the proper time of each particle with the time of the observer, using the relations above, we have a classic variational problem over the interval $[0, T]$ of the experiment.

If the fields induced by the particle are negligible we can expect a solution, but it will be at best expressed as general conditions that the trajectories must meet. The main example is the trajectory of free particles, that is particles which are not submitted to a field. With the simple lagrangian $L_1 = 1$ and the Levi-Civita connection one finds that the trajectory must be a geodesic, and there is a unique geodesic passing through any point m with a given tangent $V(0)$. But the equation does not give by itself the coordinates of the geodesic (which require the knowledge of G) or the value of the field. For the electromagnetic field, if we know the value of the field and we neglect the field induced by the particle, we get similarly a solution : $\nabla_u u = \mu_0 \frac{q}{mc} \sum_\alpha \mathcal{F}^{\alpha\beta} u_\beta$ with $u = \frac{c}{\sqrt{-\langle V, V \rangle}} V$

If we want to account for the field induced by the particle we have a problem. As the field propagates, we need to know the field out of the trajectory. It

could be computed by the more general model, and the results reintegrated in the single particle model. The resulting equation for the trajectory is known, for the electromagnetic field, as the “Lorentz-Dirac equation” (see Poisson and Quinn). The procedure is not simple, and there are doubts about the physical meaning of the equation itself.

6.2.3 Particles and Fields interacting

The fundamental issue is that the particles are not present everywhere, so even if we can represent the states of the particles by a matter field, that is a section of a vector bundle, we have to account for the actual presence of the particles : virtual particles do not interact. There are different solutions.

Common solutions

If the trajectories of the particles are known, a direct computation gives usually the field that they induce. This is useful for particles which are bonded (such as in condensed matter).

In QTF the introduction of matter fields in the lagrangian is in part formal, as most of the computations, notably when they address the problem of the creation / annihilation of particles, is done through Feynman’s diagram, which is a way to reintroduce the action at a distance between identified particles.

In the classical picture the practical solutions which have been implemented with the Principle of Least Action have many variants, but share the following assumptions :

- they assume that the particles follow some kind of continuous trajectories and keep their physical characteristics (this condition adds usually a separate constraint)
- the trajectory is the key variable, but the model gives up the concept of point particle, replaced by some form of density of particles.

These assumptions makes sense when we are close to the equilibrium, and we are concerned not by the behavior of each individual particle but by global results about distinguished populations, measured as cross sections over an hypersurface. They share many characteristics with the models used in fluid mechanics. In the usual QM interpretation the density of particles can be seen as a probability of presence, but these models are used in the classical picture, and actually the state of the particles is represented as sections of the vector bundle TM (with a constraint imposed by the mass), combined with a density function. So the density has a direct, classic interpretation.

The simplest solution is, assuming that the particles have the same physical characteristics, to take as key variable a density $\mu\varpi_4$. Then the application of the principle of least action with a 4 dimensional integral gives the equations relating the fields and the density of charge.

The classic examples are :

- the 2nd Maxwell equation in GR :

$$\nabla^\beta \mathcal{F}_{\beta\alpha} = -\mu_0 J_\alpha \Leftrightarrow \mu_0 J^\alpha \sqrt{-\det g} = \sum_\beta \partial_\beta (\mathcal{F}^{\alpha\beta} \sqrt{-\det g})$$

with the current : $J = \mu(m) qu$ and the lagrangian

$$L = \mu_0 \sum_{\alpha} \dot{A}_{\alpha} J^{\alpha} + \frac{1}{2} \sum_{\alpha\beta} \mathcal{F}_{\alpha\beta} \mathcal{F}^{\alpha\beta}$$

- the Einstein Equation in GR :

$$Ric_{\alpha\beta} - \frac{1}{2} (R + \Lambda) g_{\alpha\beta} = \frac{8\pi G}{\sqrt{c}} T_{\alpha\beta}$$

with the momentum energy tensor $T_{\alpha\beta} = \frac{\partial T}{\partial g^{\alpha\beta}} - \frac{1}{2} g_{\alpha\beta} T$

and the lagrangian $L = T(g, z^i, z_{\alpha}^i) + \frac{\sqrt{c}}{8\pi G} (R + \Lambda)$

The distribution of charges is defined independently, but it must meet a conservation law. In the examples above we must have :

$$\sum_{\alpha} \partial_{\alpha} J^{\alpha} = 0$$

$$\sum_{\alpha} \nabla^{\alpha} T_{\alpha\beta} = 0$$

The Einstein-Vlasov systems are also based on a distribution function $f(m, p)$ depending on the localization m and the linear momentum p , which must follow a conservation law, expressed as a differential equation (the Vlasov equation). The particles are generally assumed to have the same mass, so there is an additional constraint on the momentum as above. When only the gravitational field is considered the particles follow geodesics, to which the conservation law is adjusted. These systems have been extensively studied for plasmas and Astrophysics (see Andréasson).

This kind of model has been adjusted to Yang-Mills fields (Choquet-Bruhat) : the particles have different physical characteristics (similar to the vector ϕ seen previously), and must follow an additional conservation law given by $\nabla_V \phi = 0$ (the Wong equation).

In all these solutions the 4 dimensional action, with a lagrangian adapted to the fields considered, gives an equation relating the field and the distribution of charges.

So the situation is not satisfying. These difficulties have physical roots. The concept of field is aimed at removing the idea of action at a distance, but, as the example of the motion of a single particle in its own field shows, it seems difficult to circumvent the direct consideration of mutual interactions between particles, which needs to identify separately each of them.

However, from these classic examples, two results seem quite clear :

- the trajectories should belong to some family of curves, defined by the interactions

- the initial conditions, that is the beginning x of the curve and its initial tangent, should determine the curve in the family.

They are consistent with our description of the motions by sections of fiber bundles and matter fields. Moreover the Spinor formalism avoids the introduction of constraints on the state (or momentum) of the particles : the conservation law is satisfied by construct. Indeed the particles keep their intrinsic properties through ψ_0 .

Finite number of particles

For a finite, fixed, number of particles, with known fundamental state ψ_{0p} , the second integral of the action reads :

$$\begin{aligned} & \sum_{p=1}^N \int_{\tau_p^1}^{\tau_p^2} L_2((\psi_p, \nabla_{u_p} \psi_p, P_i^\alpha, V_p^\alpha)) d\tau_p \\ &= \sum_{p=1}^N \int_0^T L_2((\psi_p, \nabla_{u_p} \psi_p, P_i^\alpha)(t)) \sqrt{-\langle V_p, V_p \rangle} dt \end{aligned}$$

The states of the particles are represented by maps : $\psi_p : [0, T] \rightarrow E \otimes F$ with a fundamental state ψ_{0p} . The key variables are then $r_p(t), w_p(t)$ which are related to V_p , and the value of the potential along the world lines \widehat{G}, \widehat{A} .

There is an obvious mathematical problem : the fields and the particles are defined over domains which are manifolds of different dimensions, which precludes the usual method by Euler-Lagrange equations. It is common to put a Dirac's function in the second part, but this, naive, solution is just a formal way to rewrite the same integral without any added value.

If the model considers only a finite, fixed, number of particles, there is a rigorous mathematical solution, by functional derivatives, that we will see in the next chapter.

Density of particles

The system is a population of identical particles (they have the same fundamental state ψ_0) which follow trajectories without collisions (the trajectories do not cross) and are observed in an area Ω over the period $[0, T]$. It is then possible to represent the system by a matter field $\psi \in \mathfrak{X}(\psi_0)$ with a density μ with respect to the volume form ϖ_4 .

The key variables are :

- maps $r, w \in C_1(\Omega; \mathbb{R}^3)$ which define a section $\sigma \in \mathfrak{X}(P_G)$, and $\varkappa \in \mathfrak{X}(P_U)$
- the density $\mu \in C(\Omega; \mathbb{R}_+)$: $\mu(m) = \sqrt{\frac{\langle \psi, \psi \rangle}{\langle \psi_0, \psi_0 \rangle}}$
- the fields represented by the potentials (if there are external fields they are accounted for in addition to the fields generated by the particles).

The lagrangian for the interactions is then :

$$\int_{\Omega} L_2(\psi, \nabla_u \psi, P_i^\alpha, V^\alpha) \mu \varpi_4$$

With the conditions :

$$\psi = \mu(m) \vartheta(\sigma(m), \varkappa(m)) \psi_0$$

$$\mu \operatorname{div} V + \frac{d\mu}{dt} = 0$$

and the initial conditions, defined on $\Omega_3(0)$.

We will see an example in the next chapter.

As a specific case we have the aggregation of particles, in order to compute the inertial spinor of a deformable solid.

6.3 PERTURBATIVE LAGRANGIAN

In a perturbative approach, meaning close to the equilibrium, which are anyway the conditions in which the principle of least action applies, the lagrangian can be estimated by a development in Taylor series, meaning that each term is represented by polynomials. Because all the variables are derivatives at most of the second order and are vectorial, it is natural to look for scalar products.

6.3.1 Interactions Fields / Fields

It is generally assumed that there is no direct interaction gravitation / other fields (the deviation of light comes from the fact that the Universe, as seen in an inertial frame, is curved). So we have two distinct terms, which can involve only the strength of the field. They are two forms on M valued in the Lie algebra, which transform in a change of gauge by the adjoint map, thus the scalar product must be invariant by Ad .

We have such quantities, the density of energy of the field, defined by scalar products. So this is the obvious choice. However for the gravitational field there is the usual solution of the scalar curvature R which can be computed with our variables. It is invariant by a change of gauge or chart. The action with the scalar curvature is then $\int_{\Omega} R \varpi_4$, called the Hilbert action. Any scalar constant added to a lagrangian (with the volume form) leads to a lagrangian which is still covariant. So one can add a constant Λ to the curvature, the equation then becomes : $Ric_{\alpha\beta} - \frac{1}{2}g_{\alpha\beta}(R + \Lambda) = 0$. The **cosmological constant** Λ acts as a pressure, positive or negative, to impact the expansion of the Universe in Cosmological models. Its existence and value have been a hot topic, but it is nowadays generally acknowledged that, at least for cosmological models, it should be non null.

The models use traditionally the scalar curvature, with the Levi-Civita connection. The application of the principle of least action leads then in the vacuum to the Einstein equation : $Ric_{\alpha\beta} - \frac{1}{2}g_{\alpha\beta}R = 0$.

In our formalism this solution leads to linear equations : R is a linear function of \mathcal{F}_G , so it leads to much simpler computations than the usual method (and of course they provide the same solutions).

In all the, difficult, experimental verifications, the models are highly simplified, and to tell that the choice of R is validated by facts would be a bit excessive. We have seen that its computation, mathematically legitimate, has no real physical justification : the contraction of indices is actually similar to the procedure used to define the Dirac's operator.

It seems logical to use the same quantity for the gravitational field as for the other fields. This is the option that we will follow. It is more pedagogical, and opens the possibility to study a dissymmetric gravitational field.

So, to represent the interactions fields / fields we will take in a perturbative lagrangian (without cosmological constant) :

$$\int_{\Omega} \left(\sum_{\alpha\beta} C_G \left(\sum_{a=1}^3 \mathcal{F}_{G\alpha\beta}^a \mathcal{F}_G^{a\alpha\beta} - \sum_{a=4}^6 \mathcal{F}_{G\alpha\beta}^a \mathcal{F}_G^{a\alpha\beta} \right) + C_A \sum_{a=1}^m \mathcal{F}_{A\alpha\beta}^a \mathcal{F}_A^{a\alpha\beta} \right) \varpi_4(m) \quad (6.6)$$

where C_G, C_A are real constant scalars.

6.3.2 Interactions Particles /Fields

We have seen in the previous section that the derivative of the fields appear through the covariant derivative, and that it must be taken along the world line. We have to choose the scalar product. As the lagrangian involves only variables at the first order, the key quantity is the scalar product $\langle \psi, \nabla_{\alpha} \psi \rangle$.

We have already introduced the quantity :

$$\frac{1}{i} \langle \psi, \nabla_{\alpha} \psi \rangle = \text{Im} \langle \psi, \nabla_{\alpha} \psi \rangle = \frac{1}{i} \left(\langle \psi, \partial_{\alpha} \psi \rangle + \langle \psi, [\psi] [\dot{A}_{\alpha}] \rangle + \langle \psi, \gamma C(G_{\alpha}) \psi \rangle \right)$$

So $\frac{1}{i} \langle \psi, \nabla_V \psi \rangle$ can be seen as the energy of the particle in the system, as measured by the observer.

$$\frac{1}{i} \langle \psi, \nabla_V \psi \rangle = k^t \widehat{X} + \frac{1}{i} \langle \psi_0, [\psi_0] [Ad_{\ast} \widehat{A}] \rangle$$

where :

$$k = -\epsilon (Tr(\psi_R^* \sigma_a \psi_R))_{a=1}^3$$

and for spinors :

$$k = -\epsilon \frac{1}{2} M_p^2 c^4 k_0 = -\epsilon \frac{1}{2} M_p^2 c^4 \begin{bmatrix} (\sin 2\alpha_0) \cos(\alpha_2 - \alpha_1) \\ (\sin 2\alpha_0) \sin(\alpha_2 - \alpha_1) \\ \cos 2\alpha_0 \end{bmatrix}; k_0^t k_0 = 1$$

V depends on the vectors r, w by :

$$V = \frac{dp}{dt} = \vec{v} + c\varepsilon_0(m) = c \left(\varepsilon_0 + \epsilon \frac{a_w}{2a_w^2 - 1} \sum_{i=1}^3 w_i \varepsilon_i \right) = \sum_{i=0}^3 V^i \varepsilon_i = \sum_{\alpha=0}^3 V^{\alpha} \partial_{\xi_{\alpha}}$$

$$V^{\alpha} = \sum_{i=0}^3 P_i^{\alpha} V^i$$

$$\sqrt{-\langle V, V \rangle} = \frac{c}{2a_w^2 - 1}$$

and \widehat{X} depends on r, w and linearly on their first derivative by :

$$\widehat{X} = \sum_{\alpha=0}^3 V^{\alpha} X_{\alpha}$$

$$X_{\alpha} = [A_r(r)] \partial_{\alpha} r + [A_w(r, w)] \partial_{\alpha} w + [D_r(r, w)] G_{r\alpha} + [D_w(r, w)] G_{w\alpha}$$

with the matrix polynomials $A_r(r), \dots$ which can be expressed as 4nd order polynomials in the components r_a, w_a .

The part of the action representing the interactions is then with a density of particles :

$$\int_{\Omega} C_I \frac{1}{i} \langle \psi, \nabla_V \psi \rangle \mu \varpi_4 = \int_{\Omega} C_I \left(k^t \widehat{X} + \frac{1}{i} \langle \psi_0, [\psi_0] [Ad_{\ast} \widehat{A}] \rangle \right) \mu \varpi_4 \quad (6.7)$$

and for a collection of particles :

$$\sum_{p=1}^N \int_0^T C_I \left(k_p^t \hat{X}_p + \frac{1}{i} \langle \psi_{0p}, [\psi_{0p}] [Ad_{\mathcal{K}} \hat{A}_p] \rangle \right) \sqrt{-\langle V_p, V_p \rangle} dt \quad (6.8)$$

In the lagrangian used in the Standard Model ² we have a similar expression, with the Dirac operator and \hat{A} is identified with the bosons as force carriers (which requires the introduction of the Higgs boson).

There could be a part expressing the energy at rest of the particles. With the assumption of the existence of a fundamental state if takes the form $C_P \langle \psi_0, \psi_0 \rangle$. In a model with a density of particle this would just be the addition of a constant, similar to a cosmological constant.

²Of course the tools used in QTF are quite different (the key variables are located operators), but the comparison with the lagrangian of the Standard Model makes sense.

Chapter 7

CONTINUOUS MODELS

Continuous models represent systems where no discontinuous process occurs : the variables are mainly defined as differentiable sections of vector bundles, the trajectories do not cross, there is no creation or annihilation of particles and the bosons are absent. The application of the Principle of Least Action with a lagrangian provides usually a set of differential equations for the variables involved, and we have seen that they can be restricted to 6 sections on vector bundles. Then the theorem 95 tells us that the problem is well posed : the evolution is determinist, and the solutions depend linearly on the initial conditions. However the solutions are not necessarily unique, and the Principle does not tell us how to go from one equilibrium to another.

Continuous models correspond to an ideal situation, which are nevertheless useful to study the basic relations between the 6 variables. Using the results of the previous chapter, we will study 2 models : matter field with a density, individual particles. The main purpose is to show the computation methods, and introduce quantities which will be useful, such as currents and energy-momentum tensor.

Even if the Principle of Least action does not tell how to go from one equilibrium to the other, it provides a way to see what are the resistance that a given system opposes to a change in the equilibrium. This is done with the Energy-Momentum tensor, which can be seen as representing the inertia of the system.

The equations for the fields show that 2 special vector fields, the currents, one linked to the particles and the other to the fields, play a specific role and explain how the particles are the sources of the fields.

But at first we will see what the theorems of the QM Chapter about the evolution of a system (the equivalent of the Schrödinger equation) can tell. Without using the Principle of Least Action they give important results.

7.1 THE SCHRÖDINGER'S EQUATION

Continuous models are the natural picture for the Schrödinger's equation. The theorem 26 tells that, if the evolution of a system is quasi-determinist, then there is a unitary operator $\Theta(t)$ such that $X(t) = \Theta(t)X(0)$. To be implemented several conditions are required.

There can be other arguments in the variable X , but the time t must be unique for the system. So this is necessarily the time of an observer of Ω , who uses his standard chart $\varphi_o(t, x)$ and follows the system during an interval $[0, T]$. This choice is not neutral : it means that t is specific to an observer, but also that the result will hold for any observer with his proper time.

The second is about the determinism of the evolution. In a continuous model the Principle of Least Action provides solutions, which can be not unique, however going from one equilibrium to another requires an action on the system as it will be seen with the Energy-Momentum tensor. So, given initial conditions, and without external action, we can safely assume that the system is determinist.

The third is about the variables, which must belong to Fréchet spaces, thus for variables which are section of vector bundles for which there is at least a norm which is gauge invariant. So the only candidates are the state ψ of the particles and the strength of the fields $\mathcal{F}_A, \mathcal{F}_G$.

The theorem appears in two forms. In the stronger form the state of the system must be followed for $t \in \mathbb{R}$, which imposes conditions which can be met only in some cases.

7.1.1 Schrödinger equation for the particles

Whenever there is a fundamental state, ψ reads :

$$\psi(t) = \vartheta(\sigma(t), \varkappa(t)) \psi_0$$

for a particle on its trajectory. For $t = 0$: $\psi(0) = \vartheta(\sigma(0), \varkappa(0)) \psi_0$ with known values and

$$\psi(t) = \vartheta\left(\sigma(t) \cdot \sigma^{-1}(0), \varkappa(t) \cdot \varkappa(0)^{-1}\right) \psi(0)$$

$$\text{So : } \psi(t) = \Theta(t) \psi(0) \text{ with } \Theta(t) = \vartheta\left(\sigma(t) \cdot \sigma^{-1}(0), \varkappa(t) \cdot \varkappa(0)^{-1}\right)$$

In all common computations in QM, the wave function can be replaced by the state vector ψ , and the variables $r(t), w(t)$, replace the linear and rotational momentum operators. The lagrangian, notably under its perturbative form, replaces the Hamiltonian, with parameters the potential of the fields \widehat{G}, \widehat{A} along the trajectory. We have a clear mathematical framework in which the usual problems can be addressed.

With a matter field, and a density of particles :

$$\psi(m) = \vartheta(\sigma(m), \varkappa(m)) \psi_0$$

the matter field itself defines a vector field for the trajectories, with respect to an observer, thus we have natural maps :

$$[0, T] \rightarrow Q[E \otimes F, \vartheta] :: \psi(\Phi_V(t, x)) = \vartheta(\sigma(\Phi_V(t, x)), \varkappa(\Phi_V(t, x))) \psi_0$$

which represent the evolution of the state of a given particle (labeled by x on $\Omega_3(0)$) along its trajectory, and indeed the implementation of the Principle

of Least Action provides differential equations along the trajectory. So we still have :

$$\psi(\Phi_V(t, x)) = \Theta(t)(\Phi_V(0, x))$$

The stronger version of the theorem requires that ψ is defined all over \mathbb{R} : the evolution of the particle is in a continuous process, with no beginning or ending, and this assumption is crucial. An important special case is of bonded particles in a regular environment, such as a crystal : it can then be assumed that $\psi(m)$ is a periodic map over a lattice defined by the geometric structure of the medium. The observer is then defined with respect to this lattice (which sums up to choose a suitable chart of $\Omega(0)$). The value of the potentials is defined in this chart.

7.1.2 Schrödinger equation for the fields

We do not have a fundamental state for the fields, but we can implement the theorem, with an interesting result.

Theorem 105 *In the vacuum the fields propagate at the speed of light c for any observer. There is a continuous, unitary, map $\Theta(t)$ such that:*

$$\forall t \in [0, T] : \mathcal{F}(\varphi_o(t, x)) = \Theta(t)(\mathcal{F}(\varphi_o(0, x))) \quad (7.1)$$

Proof. $\mathcal{F}_A, \mathcal{F}_G$ take value in the Lie algebras, which are normed spaces.

We assume that the fields are observed on an area Ω on the time interval $[0, T]$ of an observer in his standard chart $\varphi_o(t, x) = m$

If we assume that $\mathcal{F}_A, \mathcal{F}_G$ are sections of $L^1(M, T_1Spin(3, 1), \varpi_4), L^1(M, T_1U, \varpi_4)$ which are separable Fréchet spaces, then the evaluations maps are continuous if we use the standard chart of the observer, because :

$$\begin{aligned} \mathcal{E}(t) : \mathcal{E}(t)\mathcal{F}_A &= \mathcal{F}_A(\varphi_o(t, x)) \\ \|\mathcal{F}\| &= \int_{\Omega} \|\mathcal{F}(m)\| \varpi_4(m) = \int_0^T c dt \int_{\Omega_3(t)} \|\mathcal{F}(\varphi_o(t, x))\| \varpi_3(x) \end{aligned}$$

Then we can implement the theorem 26.

There is a Hilbert space H and a continuous, unitary, map $\Theta(t)$ such that :

$$\begin{aligned} \mathcal{F}(\varphi_o(t, \cdot)) &\in H \\ \Theta(t) &= \mathcal{E}(t) \circ \mathcal{E}(0)^{-1} \\ \forall t \in [0, T] : \mathcal{F}(\varphi_o(t, x)) &= \Theta(t)(\mathcal{F}(\varphi_o(0, x))) \end{aligned}$$

So the value of the variable over $\Omega_3(t)$ is deduced from its value on $\Omega_3(0)$

The 4 dimensional distance between the points $\varphi_o(t + \theta, x)$ and $\varphi_o(t, x)$ for the observer is $c\theta$ so, in this meaning, for the observer, the field propagates at the speed c . ■

$\Theta(t)$ is a $m \times m$ Hermitian matrix in the orthonormal basis $(\vec{\theta}_a)_{a=1}^m$

The speed of propagation is the same for any observer, because it needs to be defined with respect to one unique time all over Ω , and it cannot be anything other than t . This is the genuine interpretation of the Michelson & Morley experiment : it can be explained by the variations of the length of the apparatus, but this is not necessary (even if it is a consequence of the

assumptions of the Relativist Geometry). The deeper fact is that to measure the speed of propagation of a field, we need to measure this propagation with respect to an observer, and all the observers are submitted to the same constraints and use similar charts.

For the EM field, we have $\mathcal{F}_A = d\dot{A} \in \Lambda_2(M; \mathbb{R})$ so that :

$$\begin{aligned} d\dot{A}(\varphi_o(t, x)) &= \Theta(t) \left(d\dot{A}(\varphi_o(0, x)) \right) \\ \Leftrightarrow \partial_\alpha \dot{A}(\varphi_o(t, x)) &= \Theta(t) \left(\partial_\alpha \dot{A}(\varphi_o(0, x)) \right) \end{aligned}$$

$$\dot{A}(\varphi_o(t, x)) = \Theta(t) \dot{A}(\varphi_o(0, x))$$

The same reasoning can be done with $j(\mathcal{F})$ so :

$j(\mathcal{F}(\varphi_o(t, x))) = \tilde{\Theta}(t) j(\mathcal{F}(\varphi_o(0, x)))$ for a unitary map on a Hilbert space which contains the Lie algebras, but not necessarily with the usual scalar product.

We have seen previously how the fields can be quantized, and the more interesting results come from the use of invariance by the compact groups U and $SO(3)$. The present results add a complement, which is directly linked to the 4th dimension t . If the propagation is continuous, then $\Theta(t)$ is continuous and the Spin of the fields $j(\mathcal{F}_R)$ stays in the same irreducible representation.

We could consider to use the Theorem 27. It requires the condition that for any fixed θ the variables $X(t)$ and $X(t + \theta)$ represent the same state of the system. If the conditions are met, then $\Theta(t) = \exp t\theta$.

But the theorem requires that the system is followed over all $t \in \mathbb{R}$. One can deal with an interval such that $[0, +\infty[$. A semi-one parameter group $U(t)$ (that is defined only for $t > 0$) can be extended to a parameter group (for any t) if : $\lim_{t \rightarrow 0^+} \|U(t) - Id\| = 0$ (Maths.1030). Continuity from the initial conditions are generally assumed, so we can consider that the theorem applies if

$$\int_0^\infty c dt \int_{\Omega_3(t)} \|\mathcal{F}(\varphi_o(t, x))\| \varpi_3(x) < \infty$$

This assumes that the field vanishes at infinity, or that it is enclosed in a box, and null outside the box, or is periodic (as in a crystal). This holds only for the fields other than gravitation.

If \mathcal{F}_A is assumed to be periodic over $[0, T]$, then a general theorem (Maths.2428) says that any periodic function of L^2 can be expressed as a Fourier series :

$$\Theta(t) = \sum_{n=-\infty}^{+\infty} \exp(2i\pi t n \nu) \Theta_n \text{ with the frequency } \nu = 1/T$$

$$\mathcal{F}_A(\varphi_o(t, x)) = \sum_{n=-\infty}^{+\infty} (\exp(2i\pi t n \nu) \Theta_n) (\mathcal{F}_A(\varphi_o(0, x)))$$

In the case of the EM field it reads :

$\mathcal{F}_{EM}(\varphi_o(t, x)) = \sum_{n=-\infty}^{+\infty} (\exp 2i\pi t n \nu) \mathcal{F}_{EM}(\varphi_o(0, x))$. This is the usual expression of the EM field as the sum of plane waves (on each $\Omega(t)$). Because \mathcal{F}_{EM} is real we have to take the real Fourier series with cos and sin. And this expression extends to the potential as seen above.

7.2 MODEL WITH A DENSITY OF PARTICLES

Whenever the action is represented by a single integral the solution of the variational calculus problem is given by the Euler-Lagrange equations.

7.2.1 Variational calculus with Euler-Lagrange Equations

Find a section Z for which the integral $\int_{\Omega} L(z^i, z_{\alpha}^i) \varpi_4$ is stationary is a classic problem of variational calculus, and the solution is given by the Euler-Lagrange equation, for each variable (Maths.34.3).

We will use the same notations as in the Previous Chapter :

L denotes the scalar lagrangian $L(z^i, z_{\alpha}^i)$ function of the variables z^i , expressed by the components in the gauge of the observer, and their partial derivatives which, in the jets bundle formalism, are considered as independent variables z_{α}^i .

$$\mathcal{L} = L(z^i, z_{\alpha}^i) (\det P')$$

$$L \varpi_4 = L(z^i, z_{\alpha}^i) (\det P') d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3 \text{ is the 4-form}$$

$\frac{\partial \mathcal{L}}{\partial z}$ to denote the usual partial derivative with respect to the variable z

$\frac{d \mathcal{L}}{dz}$ to denote the total derivative with respect to the variable z , meaning accounting for the composite expressions in which it is an argument.

For an action $\int_{\Omega} L(z^i, z_{\alpha}^i) \varpi_4$ where (z^i, z_{α}^i) is a 1-jet section of a vector bundle, the Euler-Lagrange equations read :

$$\frac{d(L \det P')}{dz^i} - \sum_{\beta} \frac{d}{d\xi^{\beta}} \frac{d(L \det P')}{dz_{\beta}^i} = 0 \quad (7.2)$$

where $\frac{d}{d\xi^{\beta}}$ is the derivative with respect to the coordinates in M . $\det [P']$ is necessary to account for ϖ_4 which involves P' .

We must compute separately the derivatives for the real and imaginary parts, and with respect to the z^i, z_{α}^i .

The equation involve the partial derivatives of each variable (not of composite maps) and the real and imaginary parts must be considered.

Whenever z^i is complex, we have two families of real valued equations :

$$\frac{\partial L \det P'}{\partial \operatorname{Re} z^i} - \sum_{\beta} \frac{d}{d\xi^{\beta}} \frac{\partial L \det P'}{\partial \operatorname{Re} z_{\beta}^i} = 0$$

$$\frac{\partial L \det P'}{\partial \operatorname{Im} z^i} - \sum_{\beta} \frac{d}{d\xi^{\beta}} \frac{\partial L \det P'}{\partial \operatorname{Im} z_{\beta}^i} = 0$$

and by defining the holomorphic complex valued functions :

$$\frac{\partial L \det P'}{\partial z^i} = \frac{\partial L \det P'}{\partial \operatorname{Re} z^i} + \frac{1}{i} \frac{\partial L \det P'}{\partial \operatorname{Im} z^i}$$

$$\frac{\partial L \det P'}{\partial z_{\beta}^i} = \frac{\partial L \det P'}{\partial \operatorname{Re} z_{\beta}^i} - \frac{1}{i} \frac{\partial L \det P'}{\partial \operatorname{Im} z_{\beta}^i}$$

the equations read :

$$\frac{\partial L \det P'}{\partial z^i} = \sum_{\beta} \frac{d}{d\xi^{\beta}} \left(\frac{\partial L \det P'}{\partial \operatorname{Re} z_{\beta}^i} + \frac{1}{i} \frac{\partial L \det P'}{\partial \operatorname{Im} z_{\beta}^i} \right) = \sum_{\beta} \frac{d}{d\xi^{\beta}} \left(\frac{\partial L \det P'}{\partial z_{\beta}^i} \right)$$

$$\begin{aligned} \frac{\partial L \det P'}{\partial z^i} &= \frac{\partial L \det P'}{\partial z^i} = \sum_{\beta} \frac{d}{d\xi^{\beta}} \left(\frac{\partial L \det P'}{\partial \operatorname{Re} z_{\beta}^i} - \frac{1}{i} \frac{\partial L \det P'}{\partial \operatorname{Im} z_{\beta}^i} \right) \\ &= \sum_{\beta} \frac{d}{d\xi^{\beta}} \left(\frac{\partial L \det P'}{\partial z_{\beta}^i} \right) = \sum_{\beta} \frac{d}{d\xi^{\beta}} \left(\frac{\partial L \det P'}{\partial z_{\beta}^i} \right) \\ \text{and we are left with the unique equation :} \\ \frac{\partial L \det P'}{\partial z^i} &= \sum_{\beta} \frac{d}{d\xi^{\beta}} \left(\frac{\partial L \det P'}{\partial z_{\beta}^i} \right) \end{aligned}$$

The equation holds pointwise for any $m \in \Omega$. However when one considers a point along a trajectory : $p(t) = m(\Phi_V(t, x))$ then the expressions like : $\sum_{\beta} V^{\beta} \frac{d}{d\xi^{\beta}} (X(p(t)))$ read : $\frac{dX}{dt}(p(t))$.

On the manifold M endowed with the volume form ϖ_4 the divergence of a vector field $X = \sum_{\alpha} X^{\alpha} \partial_{\xi_{\alpha}}$ is the function $\operatorname{div}(X) : \mathcal{L}_X \varpi_4 = \operatorname{div}(X) \varpi_4$ and its expression in coordinates is (Maths.17.2.4) :

$$\begin{aligned} \operatorname{div}(X) &= \frac{1}{\det P'} \sum_{\alpha} \partial_{\alpha} (X^{\alpha} \det P') \text{ which reads in the SR geometry : } \operatorname{div}(X) = \\ &= \sum_{\alpha} \partial_{\alpha} (X^{\alpha}) \\ \frac{dL}{dz^i} &\text{ is a vector : } Z_i = \frac{dL}{dz_{\beta}^i} \partial_{\xi_{\beta}} \text{ and } \frac{1}{\det P'} \sum_{\beta} \frac{d}{d\xi^{\beta}} \left(\frac{dL}{dz_{\beta}^i} \det P' \right) = \operatorname{div}(Z_i) \\ \frac{dL \det P'}{dz^i} \det P' + L \frac{d \det P'}{dz^i} &= \sum_{\beta} \frac{d}{d\xi^{\beta}} \left(\frac{\partial L \det P'}{\partial z_{\beta}^i} \right) \\ \frac{dL \det P'}{\partial dz^i} + L \frac{1}{\det P'} \frac{d \det P'}{dz^i} &= \operatorname{div}(Z_i) \\ \text{thus, when P' does not depend on } z^i &: \frac{dL \det P'}{dz^i} = \operatorname{div}(Z_i) \end{aligned}$$

We will do the computation with the perturbative lagrangian :

$$\int_{\Omega} \{ \sum_{\alpha\beta} C_G \langle \mathcal{F}_G, \mathcal{F}_G \rangle + C_A \langle \mathcal{F}_A, \mathcal{F}_A \rangle + C_I \mu \frac{1}{i} \langle \psi, \nabla_V \psi \rangle \} \varpi_4$$

For the state equation (ψ), which is specific, we will use a perturbative lagrangian with the EM field.

7.2.2 Equation for the Matter Field

The lagrangian

The state equation can be made explicit with a perturbative lagrangian where only the EM and gravitational field are present. The interaction term is then : $C_I \mu \frac{1}{i} \langle S, \nabla_V S \rangle = C_I \mu \left(k^t \hat{X} + \hat{A} \right)$. The variables are the vectors r, w which define V.

$$\begin{aligned} X_{\alpha} &= [A_r(r)] \partial_{\alpha} r + [A_w(r, w)] \partial_{\alpha} w + [D_r(r, w)] G_{r\alpha} + [D_w(r, w)] G_{w\alpha} \\ [A_r(r)] &= \left[\frac{1}{a_r} + \frac{2-5a_r^2}{2} j(r) + \frac{5a_r^2+1}{4a_r} j(r) j(r) \right] \\ [A_w(r, w)] &= [B_r(r)] [B_w(w)] + [C_r(r)] [C_w(w)] \\ [D_r(r, w)] &= [B_r(r)] [B_{wr}(w)] + [C_r(r)] [C_{wr}(w)] \\ [D_w(r, w)] &= [B_r(r)] [B_{ww}(w)] + [C_r(r)] [C_{ww}(w)] \\ [B_r(r)] &= \left[1 - a_r j(r) + \frac{3}{4} j(r) j(r) \right] \\ [B_w(w)] &= \frac{1}{2} j(w) \\ [C_r(r)] &= \left[1 - a_r^2 + \frac{1}{4} j(r) j(r) \right] \\ [C_w(w)] &= \frac{1}{a_w} \left[1 - \frac{1}{4} j(w) j(w) \right] \end{aligned}$$

$$\begin{aligned}
[B_{wr}(w)] &= [a_w^2 - \frac{1}{2}j(w)j(w)] \\
[B_{ww}(w)] &= a_w [j(w)] \\
[C_{wr}(w)] &= [a_w^2 - 1 - a_w j(w) + \frac{1}{4}j(w)j(w)] \\
[C_{ww}(w)] &= [1 - \frac{3}{4}j(w)j(w)] \\
a_r^2 &= 1 - \frac{1}{4}r^t r \\
a_w^2 &= 1 - \frac{1}{4}w^t w \\
\widehat{X} &= \sum_{\alpha=0}^3 V^\alpha X_\alpha
\end{aligned}$$

It is assumed that the density μ meets the continuity equation : $\frac{d\mu}{dt} + \mu \operatorname{div} V = 0$

Equations

The equations are :

$$\begin{aligned}
\forall a = 1, 2, 3 : \\
\frac{dL_I}{dr_a} &= \frac{1}{\det P'} \sum_{\alpha=0}^3 \frac{d}{d\xi^\alpha} \left(\frac{\partial L_I}{\partial \partial_\alpha r_a} \det P' \right) \\
\frac{dL_I}{dw_a} &= \frac{1}{\det P'} \sum_{\alpha=0}^3 \frac{d}{d\xi^\alpha} \left(\frac{\partial L_I}{\partial \partial_\alpha w_a} \det P' \right)
\end{aligned}$$

There are two solutions : (a_r, r, a_w, w) and $(-a_r, -r, -a_w, -w)$ corresponding to opposite relativist spatial spin.

Derivatives :

In order to get equations expressed with vectors r, w and not only their coordinates some work must be done involving the properties of the operator j which makes the matrices A_r, A_w, \dots

$$\begin{aligned}
j \in \mathcal{L}(\mathbb{R}^3; \mathcal{L}(\mathbb{R}^3; \mathbb{R}^3)) &\Rightarrow \frac{dj}{dr} = j' = j : (j(r))'(u) = j(u) \\
j^2 \in \mathcal{L}^2(\mathbb{R}^3, \mathbb{R}^3; \mathcal{L}(\mathbb{R}^3; \mathbb{R}^3)) &\Rightarrow (j(x)j(y))'(u, v) = j(u)j(y) + j(x)j(v)
\end{aligned}$$

For any matrix Y such that :

$$\begin{aligned}
Y(r) &= a + bj(r) + cj(r)j(r) \text{ with } a, b, c = Ct \\
\frac{\partial Y}{\partial r}(u) &= bj(u) + c(j(u)j(r) + j(r)j(u)) \text{ is a matrix}
\end{aligned}$$

$$\begin{aligned}
v^t \frac{\partial Y}{\partial r}(u) &= bv^t j(u) + cv^t (ru^t + ur^t - 2r^t u) \\
&= u^t [-bj(v) + c((v^t r) + vr^t - 2rv^t)] \\
&= u^t [-bj(v) + c((v^t r) + j(r)j(v) + v^t r - 2j(v)j(r) - 2v^t r)] \\
&= u^t [-bj(v) + c(j(r)j(v) - 2j(v)j(r))]
\end{aligned}$$

and we will denote :

$$[\delta Y(r, v)] = [-bj(v) + c(j(r)j(v) - 2j(v)j(r))] \quad (7.3)$$

$$v^t \frac{\partial Y}{\partial r}(u) = u^t [\delta Y(r, v)] \Leftrightarrow \left[\frac{\partial Y}{\partial r}(u) \right] v = [\delta Y(r, v)]^t u \quad (7.4)$$

In particular, by definition : $\frac{\partial Y}{\partial r_a} = \frac{\partial Y}{\partial r}(\varepsilon_a)$

$$v^t \left[\frac{\partial Y}{\partial r_a} \right] = \varepsilon_a^t [\delta Y(r, u)] \Leftrightarrow [\delta Y(r, u)]^t \varepsilon_a = \left[\frac{\partial Y}{\partial r_a} \right]^t v \quad (7.5)$$

Equation for r :

$$\begin{aligned} \frac{dL_I}{dr_a} &= \frac{d}{dr_a} \left(C_{I\mu} \left(k^t \widehat{X} + \widehat{A} \right) \right) = C_{I\mu} k^t \frac{d\widehat{X}}{dr_a} = C_{I\mu} k^t \left(\frac{\partial X_\alpha}{\partial a_r} \frac{\partial a_r}{\partial r_a} + \frac{\partial X_\alpha}{\partial r_a} \right) \\ &= C_{I\mu} \left(-\frac{1}{4a_r} k^t \frac{\partial \widehat{X}}{\partial a_r} r_a + k^t \frac{\partial \widehat{X}}{\partial r_a} \right) = C_{I\mu} \left(-\frac{1}{4a_r} (\varepsilon_a^t r) \left(k^t \frac{\partial \widehat{X}}{\partial a_r} \right) + k^t \frac{\partial \widehat{X}}{\partial r_a} \right) \\ k^t \frac{\partial \widehat{X}}{\partial r_a} &= k^t \left[\frac{\partial A_r}{\partial r_a} \right] \frac{dr}{dt} + k^t \left[\frac{\partial B_r}{\partial r_a} \right] \left([B_w] \frac{dw}{dt} + [B_{wr}] \widehat{G}_r + [B_{ww}] \widehat{G}_w \right) \\ &+ k^t \left[\frac{\partial C_r}{\partial r_a} \right] \left([C_w] \frac{dw}{dt} + [C_{wr}] \widehat{G}_r + [C_{ww}] \widehat{G}_w \right) \\ &= \varepsilon_a^t \{ [\delta A_r(r, k)] \frac{dr}{dt} + ([\delta B_r(r, k)] [B_w] + [\delta C_r(r, k)] [C_w]) \frac{dw}{dt} + [A_3] \widehat{G}_r + [A_4] \widehat{G}_w \} \\ &\text{with :} \\ [A_3] &= [\delta B_r(r, k)] [B_{wr}] + [\delta C_r(r, k)] [C_{wr}] \\ [A_4] &= [\delta B_r(r, k)] [B_{ww}] + [\delta C_r(r, k)] [C_{ww}] \\ \frac{dL_I}{dr_a} &= C_{I\mu} \varepsilon_a^t \left\{ -\frac{1}{4a_r} \left(k^t \frac{\partial \widehat{X}}{\partial a_r} \right) r + [\delta A_r(r, k)] \frac{dr}{dt} \right. \\ &+ ([\delta B_r(r, k)] [B_w] + [\delta C_r(r, k)] [C_w]) \frac{dw}{dt} \\ &+ [A_3] \widehat{G}_r + [A_4] \widehat{G}_w \} \\ \frac{\partial L_I}{\partial \alpha r_a} &= \frac{\partial}{\partial \alpha r_a} \left(C_{I\mu} \left(k^t \widehat{X} + \widehat{A} \right) \right) = C_{I\mu} k^t \frac{\partial \widehat{X}}{\partial \alpha r_a} = C_{I\mu} V^\alpha (k^t [A_r] \varepsilon_a) = \\ &C_{I\mu} V^\alpha \left(\varepsilon_a^t [A_r]^t k \right) \end{aligned}$$

The equation read :

$$\begin{aligned} &\mu \left\{ -\frac{1}{4a_r} \left(k^t \frac{\partial \widehat{X}}{\partial a_r} \right) r + [\delta A_r(r, k)] \frac{dr}{dt} \right. \\ &+ ([\delta B_r(r, k)] [B_w] + [\delta C_r(r, k)] [C_w]) \frac{dw}{dt} + [A_3] \widehat{G}_r + [A_4] \widehat{G}_w \} \\ &= \frac{1}{\det P'} \sum_{\alpha=0}^3 \frac{d}{d\xi^\alpha} \left(\mu V^\alpha ([A_r]^t k) \det P' \right) \\ &= \left([A_r]^t k \right) \sum_{\alpha=0}^3 V^\alpha \frac{d\mu}{d\xi^\alpha} + \left([A_r]^t k \right) \mu \frac{1}{\det P'} \sum_{\alpha=0}^3 \frac{d}{d\xi^\alpha} (V^\alpha \det P') \\ &+ \mu \sum_{\alpha=0}^3 V^\alpha \frac{d}{d\xi^\alpha} \left([A_r]^t k \right) \\ &= \left([A_r]^t k \right) \left(\frac{d\mu}{dt} + \mu \operatorname{div} V \right) + \mu \frac{d}{dt} \left([A_r]^t k \right) \end{aligned}$$

With the continuity equation we are left with :

$$\begin{aligned} &-\frac{1}{4a_r} \left(k^t \frac{\partial \widehat{X}}{\partial a_r} \right) r + [\delta A_r(r, k)] \frac{dr}{dt} + ([\delta B_r(r, k)] [B_w] + [\delta C_r(r, k)] [C_w]) \frac{dw}{dt} \\ &+ [A_3] \widehat{G}_r + [A_4] \widehat{G}_w = \frac{d}{dt} \left([A_r]^t k \right) \end{aligned}$$

By product with r^t :

$$\begin{aligned} &-\frac{1}{4a_r} \left(k^t \frac{\partial \widehat{X}}{\partial a_r} \right) = \frac{1}{4(1-a_r^2)} \left\{ r^t \frac{d}{dt} \left([A_r]^t k \right) \right. \\ &\left. - r^t \left([\delta A_r(r, k)] \frac{dr}{dt} + ([\delta B_r(r, k)] [B_w] + [\delta C_r(r, k)] [C_w]) \frac{dw}{dt} + [A_3] \widehat{G}_r + [A_4] \widehat{G}_w \right) \right\} \end{aligned}$$

and substitution, the equation reads :

$$\begin{aligned} &\left(1 - \frac{1}{4(1-a_r^2)} r r^t \right) \times \\ &\left([\delta A_r(r, k)] \frac{dr}{dt} + ([\delta B_r(r, k)] [B_w] + [\delta C_r(r, k)] [C_w]) \frac{dw}{dt} + [A_3] \widehat{G}_r + [A_4] \widehat{G}_w \right) \end{aligned}$$

$$= \left(1 - \frac{1}{4(1-a_r^2)} r r^t\right) \frac{d}{dt} \left([A_r]^t k\right)$$

with : $r r^t = j(r) j(r) + r^t r$

$$\left(1 - \frac{1}{4(1-a_r^2)} r r^t\right) = -\frac{1}{4(1-a_r^2)} j(r) j(r)$$

The only eigenvector with eigen value 0 of $j(r) j(r)$ is r itself. So we have the solution with some scalar function $K_r(t)$:

$$[\delta A_r(r, k)] \frac{dr}{dt} + ([\delta B_r(r, k)] [B_w] + [\delta C_r(r, k)] [C_w]) \frac{dw}{dt} + [A_3] \widehat{G}_r + [A_4] \widehat{G}_w - \frac{d}{dt} \left([A_r]^t k\right) = r K_r$$

$$\frac{d}{dt} \left([A_r]^t k\right)$$

$$= \frac{\partial [A_r]^t}{\partial a_r} k \frac{da_r}{dt} + \sum_{b=1}^3 \frac{\partial [A_r]^t}{\partial r_b} k \frac{dr_b}{dt}$$

$$= -\frac{1}{4a_r} \frac{\partial [A_r]^t}{\partial a_r} k \left(r^t \frac{dr}{dt}\right) + \left[\frac{\partial [A_r]^t}{\partial r} \left(\frac{dr}{dt}\right)\right] k$$

$$= -\frac{1}{4a_r} \frac{\partial [A_r]^t}{\partial a_r} (j(r) j(k) + k^t r) \frac{dr}{dt} + [\delta A_r^t(r, k)]^t \frac{dr}{dt}$$

$$= \left(-\frac{1}{4a_r} \frac{\partial [A_r]^t}{\partial a_r} (j(r) j(k) + k^t r) + [\delta A_r^t(r, k)]^t\right) \frac{dr}{dt}$$

The equation reads :

$$A_1(r) \frac{dr}{dt} + A_2(r, w) \frac{dw}{dt} + A_3(r, w) \widehat{G}_r + A_4(r, w) \widehat{G}_w = r K_r \quad (7.6)$$

With :

$$[A_1(r)] = [\delta A_r(r, k)] - [\delta A_r^t(r, k)]^t + \frac{1}{4a_r} \frac{\partial [A_r]^t}{\partial a_r} (j(r) j(k) + k^t r)$$

$$= \frac{1}{4a_r} \left\{ (k^t r) \left(-\frac{1}{a_r^2} + 5a_r j(r) + \frac{5a_r^2 - 1}{4a_r^2} j(r) j(r)\right) - 20a_r (1 - a_r^2) j(k) \right.$$

$$\left. + \frac{-3 + a_r^2 + 20a_r^4}{a_r^2} j(r) j(k) - 3(5a_r^2 + 1) j(k) j(r) \right\}$$

$$[A_2(r, w)] = [\delta B_r(r, k)] [B_w] + [\delta C_r(r, k)] [C_w]$$

$$= \frac{1}{2} a_r j(k) j(w) + \frac{1}{4} [j(r) j(k) - 2j(k) j(r)] \left[\frac{1}{a_w} + \frac{3}{2} j(w) - \frac{1}{4a_w} j(w) j(w)\right]$$

$$[A_3(r, w)] = [\delta B_r(r, k)] [B_{wr}] + [\delta C_r(r, k)] [C_{wr}]$$

$$= a_r j(k) [a_w^2 - \frac{1}{2} j(w) j(w)] + [j(r) j(k) - 2j(k) j(r)] \frac{1}{4a_w^2} [3a_w^4 + 1 - \frac{1}{4} (6a_w^2 + 1) j(w) j(w)]$$

$$[A_4(r, w)] = [\delta B_r(r, k)] [B_{ww}] + [\delta C_r(r, k)] [C_{ww}]$$

$$= a_r a_w j(k) j(w) + \frac{1}{4} [j(r) j(k) - 2j(k) j(r)] \left[1 + 3a_w j(w) - \frac{3}{4} j(w) j(w)\right]$$

One can check that if (a_r, r, a_w, w) is a solution, then $(-a_r, -r, -a_w, -w)$ is still a solution.

Equation for w :

$$\frac{dL_I}{d\partial_\alpha w_\alpha} = \frac{d}{d\partial_\alpha w_\alpha} \left(C_I \mu \sum_{\alpha=0}^3 V^\alpha \left(k^t X_\alpha + \dot{A}_\alpha\right)\right)$$

$$= C_I \mu V^\alpha \left(k^t \frac{dX_\alpha}{d\partial_\alpha w_\alpha}\right)$$

$$= C_I \mu V^\alpha k^t ([B_r] [B_w] + [C_r] [C_w]) \varepsilon_a$$

$$= \varepsilon_a^t C_I \mu V^\alpha ([B_r] [B_w] + [C_r] [C_w])^t k$$

$$\begin{aligned}
\frac{dL_I}{dw_a} &= \frac{d}{dw_a} \left(C_I \mu \sum_{\alpha=0}^3 V^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) \right) = C_I \mu \sum_{\alpha=0}^3 \frac{dV^\alpha}{dw_a} \left(k^t X_\alpha + \dot{A}_\alpha \right) + \\
&V^\alpha k^t \frac{dX_\alpha}{dw_a} \\
V &= \sum_{\alpha=0}^3 V^\alpha \partial \xi_\alpha \text{ is related to } w \text{ by } V^\alpha = \sum_{i=0}^3 V^i P_i^\alpha \text{ with : } V = \\
\sum_{i=0}^3 V^i \varepsilon_i &= c \left(\varepsilon_0 + \frac{a_w}{2a_w^2-1} \sum_{i=1}^3 w_i \varepsilon_i \right) \\
\frac{dV^\alpha}{dw_a} &= \sum_{i=0}^3 \frac{dV^i}{dw_a} P_i^\alpha \\
&= \sum_{i=1}^3 \left(w_i \frac{d}{dw_a} \left(\frac{ca_w}{2a_w^2-1} \right) + \frac{ca_w}{2a_w^2-1} \delta_i^a \right) P_i^\alpha \\
&= \sum_{i=1}^3 \left(-cw_i \frac{2a_w^2+1}{(2a_w^2-1)^2} \frac{1}{4a_w} w_a + \frac{ca_w}{2a_w^2-1} \delta_i^a \right) P_i^\alpha \\
&= -w_a \frac{2a_w^2+1}{(2a_w^2-1)^2} \frac{1}{4a_w} \sum_{i=1}^3 cw_i P_i^\alpha + \frac{ca_w}{2a_w^2-1} \sum_{i=1}^3 \delta_i^a P_i^\alpha \\
&= -(\varepsilon_a^t w) \frac{c(2a_w^2+1)}{4a_w(2a_w^2-1)^2} \sum_{i=1}^3 w_i P_i^\alpha + \frac{ca_w}{2a_w^2-1} \sum_{i=1}^3 P_i^\alpha (\varepsilon_a^t \varepsilon_i) \\
\sum_{\alpha=0}^3 V^\alpha k^t \frac{dX_\alpha}{dw_a} &= k^t \frac{d\widehat{X}}{dw_a} = k^t \left(\frac{d\widehat{X}}{da_w} \frac{da_w}{dw_a} + \frac{\partial \widehat{X}}{\partial w_a} \right) = k^t \left(\frac{d\widehat{X}}{da_w} \frac{1}{4a_w} w_a + \frac{\partial \widehat{X}}{\partial w_a} \right) = \\
(\varepsilon_a^t w) &\left(k^t \frac{d\widehat{X}}{da_w} \right) \frac{1}{4a_w} + k^t \frac{\partial \widehat{X}}{\partial w_a} \\
k^t \frac{\partial \widehat{X}}{\partial w_a} &= k^t [B_r] \left(\left[\frac{\partial B_w}{\partial w_a} \right] \frac{dw}{dt} + \left[\frac{\partial B_{wr}}{\partial w_a} \right] \widehat{G}_r + \left[\frac{\partial B_{ww}}{\partial w_a} \right] \widehat{G}_w \right) \\
&+ k^t [C_r] \left(\left[\frac{\partial C_w}{\partial w_a} \right] \frac{dw}{dt} + \left[\frac{\partial C_{wr}}{\partial w_a} \right] \widehat{G}_r + \left[\frac{\partial C_{ww}}{\partial w_a} \right] \widehat{G}_w \right) \\
&= \varepsilon_a^t \left\{ \left(\left[\delta B_w \left(w, [B_r]^t k \right) \right] + \left[\delta C_w \left(w, [C_r]^t k \right) \right] \right) \frac{dw}{dt} + [B_3] \widehat{G}_r + [B_4] \widehat{G}_w \right\} \\
\text{with :} & \\
[B_3] &= \left[\delta B_{wr} \left(w, [B_r]^t k \right) \right] + \left[\delta C_{wr} \left(w, [C_r]^t k \right) \right] \\
[B_4] &= \left(\left[\delta B_{ww} \left(w, [B_r]^t k \right) \right] + \left[\delta C_{ww} \left(w, [C_r]^t k \right) \right] \right) \\
\frac{dL_I}{dw_a} &= \varepsilon_a^t C_I \mu \sum_{\alpha=0}^3 \left(-w \frac{c(2a_w^2+1)}{4a_w(2a_w^2-1)^2} \sum_{i=1}^3 w_i P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) + \frac{ca_w}{2a_w^2-1} \sum_{i=1}^3 P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) \varepsilon_i \right) \\
&+ \varepsilon_a^t w \left(k^t \frac{d\widehat{X}}{da_w} \right) \frac{1}{4a_w} \\
&+ \varepsilon_a^t C_I \mu \left\{ \left(\left[\delta B_w \left(w, [B_r]^t k \right) \right] + \left[\delta C_w \left(w, [C_r]^t k \right) \right] \right) \frac{dw}{dt} + [B_3] \widehat{G}_r + [B_4] \widehat{G}_w \right\}
\end{aligned}$$

The equation reads :

$$\begin{aligned}
&\mu \sum_{\alpha=0}^3 \left(-w \frac{c(2a_w^2+1)}{4a_w(2a_w^2-1)^2} \sum_{i=1}^3 w_i P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) + \frac{ca_w}{2a_w^2-1} \sum_{i=1}^3 P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) \varepsilon_i \right) \\
&+ \mu w \left(k^t \frac{d\widehat{X}}{da_w} \right) \frac{1}{4a_w} \\
&+ \mu \left\{ \left(\left[\delta B_w \left(w, [B_r]^t k \right) \right] + \left[\delta C_w \left(w, [C_r]^t k \right) \right] \right) \frac{dw}{dt} + [B_3] \widehat{G}_r + [B_4] \widehat{G}_w \right\} \\
&= \frac{1}{\det P'} \sum_{\alpha=0}^3 \frac{d}{d\xi^\alpha} \left(\mu V^\alpha \left([B_r] [B_w] + [C_r] [C_w] \right)^t k \det P' \right) \\
&= \left([B_r] [B_w] + [C_r] [C_w] \right)^t k \left(\sum_{\alpha=0}^3 V^\alpha \frac{d\mu}{d\xi^\alpha} + \mu \frac{1}{\det P'} \sum_{\alpha=0}^3 \frac{d}{d\xi^\alpha} \left(V^\alpha \det P' \right) \right) \\
&+ \mu \sum_{\alpha=0}^3 V^\alpha \frac{d}{d\xi^\alpha} \left(\left([B_r] [B_w] + [C_r] [C_w] \right)^t k \right) \\
&= \left([B_r] [B_w] + [C_r] [C_w] \right)^t k \left(\frac{d\mu}{dt} + \mu \operatorname{div} V \right) + \mu \frac{d}{dt} \left(\left([B_r] [B_w] + [C_r] [C_w] \right)^t k \right)
\end{aligned}$$

$$= \mu \frac{d}{dt} \left(([B_r] [B_w] + [C_r] [C_w])^t k \right)$$

with the continuity equation, that is :

$$\begin{aligned} & w \left(\left(k^t \frac{d\hat{X}}{da_w} \right) \frac{1}{4a_w} - \frac{c(2a_w^2+1)}{4a_w(2a_w^2-1)^2} \sum_{i=1}^3 w_i P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) \right) \\ & + \frac{ca_w}{2a_w^2-1} \sum_{\alpha,i=1}^3 P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) \varepsilon_i \\ & + \left(\left[\delta B_w \left(w, [B_r]^t k \right) \right] + \left[\delta C_w \left(w, [C_r]^t k \right) \right] \right) \frac{dw}{dt} + [B_3] \hat{G}_r + [B_4] \hat{G}_w \\ & - \frac{d}{dt} \left(([B_r] [B_w] + [C_r] [C_w])^t k \right) \\ & = 0 \end{aligned}$$

With product by w^t :

$$\begin{aligned} & \left(\left(k^t \frac{d\hat{X}}{da_w} \right) \frac{1}{4a_w} - \frac{c(2a_w^2+1)}{4a_w(2a_w^2-1)^2} \sum_{i=1}^3 w_i P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) \right) \\ & = \frac{1}{4(a_w^2-1)} w^t \left\{ -\frac{ca_w}{2a_w^2-1} \sum_{\alpha,i=1}^3 P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) \varepsilon_i \right. \\ & \left. - \left(\left[\delta B_w \left(w, [B_r]^t k \right) \right] + \left[\delta C_w \left(w, [C_r]^t k \right) \right] \right) \frac{dw}{dt} + [B_3] \hat{G}_r + [B_4] \hat{G}_w + \frac{d}{dt} \left(([B_r] [B_w] + [C_r] [C_w])^t k \right) \right\} \end{aligned}$$

and substitution as above :

$$\begin{aligned} & \left(1 - \frac{1}{4(a_w^2-1)} w w^t \right) \times \\ & \left\{ \frac{ca_w}{2a_w^2-1} \sum_{\alpha,i=1}^3 P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) \varepsilon_i \right. \\ & \left. + \left(\left[\delta B_w \left(w, [B_r]^t k \right) \right] + \left[\delta C_w \left(w, [C_r]^t k \right) \right] \right) \frac{dw}{dt} + [B_3] \hat{G}_r + [B_4] \hat{G}_w - \frac{d}{dt} \left(([B_r] [B_w] + [C_r] [C_w])^t k \right) \right\} = \\ & 0 \\ & 1 - \frac{1}{4(a_w^2-1)} w w^t = -\frac{1}{4(a_w^2-1)} j(w) j(w) \end{aligned}$$

Thus :

$$\begin{aligned} & \left(\left[\delta B_w \left(w, [B_r]^t k \right) \right] + \left[\delta C_w \left(w, [C_r]^t k \right) \right] \right) \frac{dw}{dt} + [B_3] \hat{G}_r + [B_4] \hat{G}_w - \frac{d}{dt} \left(([B_r] [B_w] + [C_r] [C_w])^t k \right) \\ & = -\frac{a_w}{2a_w^2-1} \sum_{\alpha,i=1}^3 P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) \varepsilon_i + w K_w \end{aligned}$$

with some scalar function $K_w(t)$

$$\begin{aligned} & \frac{d}{dt} \left(([B_r] [B_w] + [C_r] [C_w])^t k \right) \\ & = \frac{dB_w^t}{dt} B_r^t k + B_w^t \frac{dB_r^t}{dt} k + \frac{dC_w^t}{dt} C_r^t k + C_w^t \frac{dC_r^t}{dt} k \\ & = \left(\frac{1}{4a_w} \frac{\partial B_w^t}{\partial a_w} B_r^t k \left(w^t \frac{dw}{dt} \right) + \left[\frac{\partial B_w^t}{\partial w} \left(\frac{dw}{dt} \right) \right] B_r^t k \right) + B_w^t \left(-\frac{1}{4a_r} \frac{\partial B_r^t}{\partial a_r} k \left(r^t \frac{dr}{dt} \right) + \left[\frac{\partial B_r^t}{\partial r} \left(\frac{dr}{dt} \right) \right] k \right) \\ & + \left(\frac{1}{4a_w} \frac{\partial C_w^t}{\partial a_w} C_r^t k \left(w^t \frac{dw}{dt} \right) + \left[\frac{\partial C_w^t}{\partial w} \left(\frac{dw}{dt} \right) \right] C_r^t k \right) + C_w^t \left(-\frac{1}{4a_r} \frac{\partial C_r^t}{\partial a_r} k \left(r^t \frac{dr}{dt} \right) + \left[\frac{\partial C_r^t}{\partial r} \left(\frac{dr}{dt} \right) \right] k \right) \\ & = \left(\frac{1}{4a_w} \frac{\partial B_w^t}{\partial a_w} B_r^t \left(j(w) j(k) + (k^t w) \right) \frac{dw}{dt} + \left[\delta B_w^t \left(w, B_r^t k \right) \right]^t \left(\frac{dw}{dt} \right) \right) \\ & + B_w^t \left(-\frac{1}{4a_r} \frac{\partial B_r^t}{\partial a_r} \left(j(r) j(k) + (k^t r) \right) \frac{dr}{dt} + \left[\delta B_r^t \left(r, k \right) \right]^t \frac{dr}{dt} \right) \\ & + \left(\frac{1}{4a_w} \frac{\partial C_w^t}{\partial a_w} C_r^t \left(j(w) j(k) + (k^t w) \right) \frac{dw}{dt} + \left[\delta C_w^t \left(w, C_r^t k \right) \right]^t \left(\frac{dw}{dt} \right) \right) \\ & + C_w^t \left(-\frac{1}{4a_r} \frac{\partial C_r^t}{\partial a_r} \left(j(r) j(k) + (k^t r) \right) \frac{dr}{dt} + \left[\delta C_r^t \left(r, k \right) \right]^t \frac{dr}{dt} \right) \\ & = \left(\frac{1}{4a_w} \left(\frac{\partial B_w^t}{\partial a_w} B_r^t + \frac{\partial C_w^t}{\partial a_w} C_r^t \right) \left(j(w) j(k) + (k^t w) \right) + \left[\delta B_w^t \left(w, B_r^t k \right) \right]^t + \left[\delta C_w^t \left(w, C_r^t k \right) \right]^t \right) \frac{dw}{dt} \\ & + \left(-\frac{1}{4a_r} \left(B_w^t \frac{\partial B_r^t}{\partial a_r} + C_w^t \frac{\partial C_r^t}{\partial a_r} \right) \left(j(r) j(k) + (k^t r) \right) + B_w^t \left[\delta B_r^t \left(r, k \right) \right]^t + C_w^t \left[\delta C_r^t \left(r, k \right) \right]^t \right) \frac{dr}{dt} \end{aligned}$$

The equation reads :

$$B_1(r, w) \frac{dr}{dt} + B_2(r, w) \frac{dw}{dt} + B_3(r, w) \widehat{G}_r + B_4(r, w) \widehat{G}_w = -\frac{ca_w}{2a_w^2 - 1} \sum_{\alpha, i=1}^3 P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) \varepsilon_i + w K_w \quad (7.7)$$

Using the identities :

$$j(j(x)y) = j(x)j(y) - j(y)j(x)$$

$$j(j(x)j(x)y) = (y^t x)j(x) - (x^t x)j(y)$$

one can compute the matrices and one gets :

$$\begin{aligned} [B_1(r, w)] &= \frac{1}{4a_r} \left(\frac{\partial B_r}{\partial a_r} B_w + \frac{\partial C_r}{\partial a_r} C_w \right)^t (j(r)j(k) + (k^t r)) - ([\delta B_r^t(r, k)] B_w + [\delta C_r^t(r, k)] C_w)^t \\ &= \frac{1}{2a_w} \left[-(k^t r) + 3j(r)j(k) - \frac{1}{2}j(k)j(r) \right] \\ &\quad + j(w) \frac{1}{8a_r} \left[(4a_r^2 - 6a_r j(r) - j(r)j(r))j(k) - (k^t r)j(r) + 3a_r j(k)j(r) \right] \\ &\quad + \frac{1}{8a_w} j(w)j(w) \left[(k^t r) - 7j(r)j(k) + \frac{1}{2}j(k)j(r) \right] \\ [B_2(r, w)] &= \left[\delta B_w(w, [B_r]^t k) \right] - [\delta B_w^t(w, B_r^t k)]^t \\ &\quad + \left[\delta C_w(w, [C_r]^t k) \right] - [\delta C_w^t(w, C_r^t k)]^t \\ &\quad - \frac{1}{4a_w} \left(B_r \frac{\partial B_w}{\partial a_w} + C_r \frac{\partial C_w}{\partial a_w} \right)^t (j(w)j(k) + (k^t w)) \\ &= \frac{3}{16a_w} (k^t r) (j(r)j(w) - j(w)j(r)) + \\ &\quad \frac{1}{a_w^3} \left[1 - \frac{1}{4}j(w)j(w) \right] \left[1 - a_r^2 + \frac{1}{4}j(r)j(r) \right] (j(w)j(k) + (k^t w)) \\ [B_3(r, w)] &= \left[\delta B_{wr}(w, [B_r]^t k) \right] + \left[\delta C_{wr}(w, [C_r]^t k) \right] \\ &= -\frac{a_w}{4} (k^t r) j(r) - j(w) \left[\frac{1}{2} (1 - 3a_r^2) j(k) - (k^t r) \frac{5}{16} j(r) + \frac{a_r}{2} (j(k)j(r) - j(r)j(k)) \right] \\ &\quad - \left[(-2 + 3a_r^2) j(k) + (k^t r) \frac{5}{8} j(r) + a_r (j(r)j(k) - j(k)j(r)) \right] j(w) \\ [B_4(r, w)] &= \left[\delta B_{ww}(w, [B_r]^t k) \right] + \left[\delta C_{ww}(w, [C_r]^t k) \right] \\ &= a_w \left[(3a_r^2 - 2) j(k) + \frac{3}{4} (k^t r) j(r) - a_r (j(r)j(k) - j(k)j(r)) \right] \\ &\quad - \frac{3}{16} (k^t r) (2j(r)j(w) - j(w)j(r)) \end{aligned}$$

One can check that if (a_r, r, a_w, w) is a solution, then $(-a_r, -r, -a_w, -w)$ is still a solution.

We will see below that on shell $k^t \widehat{X} + \widehat{A} = 0$

7.2.3 Equations for the gravitational field

We will use the lagrangian with the interaction term in its general form $C_I \mu^{\frac{1}{2}} \langle \psi, \nabla_V \psi \rangle$

The equations are :

$\forall a = 1..6, \alpha = 0..3 :$

$$\frac{d(L \det P')}{dG_\alpha^a} = \sum_\beta \frac{d}{d\xi^\beta} \frac{d(L \det P')}{d\delta_\beta G_\alpha^a}$$

Derivatives

$$\begin{aligned}
\text{i) } \frac{dL}{dG_\alpha^a} &= C_I \mu \frac{1}{i} \frac{\partial}{\partial G_\alpha^a} \langle \psi, \nabla_V \psi \rangle + C_G \frac{\partial}{\partial G_\alpha^a} \sum_{\lambda\mu} \langle \mathcal{F}_G^{\lambda\mu}, \mathcal{F}_{G\lambda\mu} \rangle \\
\text{ii) } \frac{\partial}{\partial G_\alpha^a} \langle \psi, \nabla_V \psi \rangle &= V^\alpha \frac{\partial}{\partial G_\alpha^a} \langle \psi, \nabla_\alpha \psi \rangle = V^\alpha \left\langle \psi, \frac{\partial}{\partial G_\alpha^a} \nabla_\alpha \psi \right\rangle = V^\alpha \langle \psi, [\gamma C(\vec{\kappa}_a)] [\psi] \rangle \\
\text{iii) } \frac{\partial}{\partial G_\alpha^a} \langle \mathcal{F}_G, \mathcal{F}_G \rangle &= \frac{\partial}{\partial G_\alpha^a} \left(\sum_{b=1}^3 \sum_{prqs} (\mathcal{F}_{r pq}^b g^{pr} g^{qs} \mathcal{F}_{rrs}^b - \mathcal{F}_{wpq}^b g^{pr} g^{qs} \mathcal{F}_{wrs}^b) \right) \\
&= \sum_{b=1}^3 \sum_{prqs} \left(\frac{\partial}{\partial G_\alpha^a} \mathcal{F}_{r pq}^b \right) g^{pr} g^{qs} \mathcal{F}_{rrs}^b - \left(\frac{\partial}{\partial G_\alpha^a} \mathcal{F}_{wpq}^b \right) g^{pr} g^{qs} \mathcal{F}_{wrs}^b \\
&+ \sum_{b=1}^3 \sum_{prqs} \mathcal{F}_{r pq}^b g^{pr} g^{qs} \left(\frac{\partial}{\partial G_\alpha^a} \mathcal{F}_{rrs}^b \right) - \mathcal{F}_{wpq}^b g^{pr} g^{qs} \left(\frac{\partial}{\partial G_\alpha^a} \mathcal{F}_{wrs}^b \right) \\
&= 2 \sum_{b=1}^3 \sum_{\lambda\mu} \left(\frac{\partial}{\partial G_\alpha^a} \mathcal{F}_{r\lambda\mu}^b \right) \mathcal{F}_r^{b\lambda\mu} - \left(\frac{\partial}{\partial G_\alpha^a} \mathcal{F}_{w\lambda\mu}^b \right) \mathcal{F}_w^{b\lambda\mu} \\
\frac{\partial}{\partial G_\alpha^a} \mathcal{F}_{r\lambda\mu}^b &= \frac{\partial}{\partial G_\alpha^a} [j(G_{r\lambda}) G_{r\mu} - j(G_{w\lambda}) G_{w\mu}]^b = \frac{\partial}{\partial G_\alpha^a} \sum_{p,q=1}^3 \epsilon(b,p,q) [G_{r\lambda}^p G_{r\mu}^q - G_{w\lambda}^p G_{w\mu}^q] \\
\frac{\partial}{\partial G_\alpha^a} \mathcal{F}_{w\lambda\mu}^b &= \frac{\partial}{\partial G_\alpha^a} [j(G_{w\lambda}) G_{r\mu} + j(G_{r\lambda}) G_{w\mu}]^b = \frac{\partial}{\partial G_\alpha^a} \sum_{p,q=1}^3 \epsilon(b,p,q) [G_{w\lambda}^p G_{r\mu}^q + G_{r\lambda}^p G_{w\mu}^q] \\
a = 1, 2, 3 : \\
\frac{\partial}{\partial G_\alpha^a} \mathcal{F}_{r\lambda\mu}^b &= \sum_{p,q=1}^3 \epsilon(b,p,q) (\delta_\alpha^\lambda \delta_a^p G_{r\mu}^q + \delta_\alpha^\mu \delta_a^q G_{r\lambda}^p) \\
&= \epsilon(b,a,c) (\delta_\alpha^c G_{r\mu}^c - \delta_\alpha^c G_{r\lambda}^c) \\
\frac{\partial}{\partial G_\alpha^a} \mathcal{F}_{w\lambda\mu}^b &= \sum_{p,q=1}^3 \epsilon(b,p,q) [\delta_\alpha^\mu \delta_a^q G_{w\lambda}^p + \delta_\alpha^\lambda \delta_a^p G_{w\mu}^q] \\
&= \epsilon(b,a,c) (-\delta_\alpha^c G_{w\lambda}^c + \delta_\alpha^c G_{w\mu}^c) \\
\frac{\partial}{\partial G_\alpha^a} \langle \mathcal{F}_G, \mathcal{F}_G \rangle &= -4 [j(\mathcal{F}_r^{\alpha\lambda}) G_{r\lambda} - j(\mathcal{F}_w^{\alpha\lambda}) G_{w\lambda}]^a \\
a = 4, 5, 6 : \\
\frac{\partial}{\partial G_\alpha^a} \mathcal{F}_{r\lambda\mu}^b &= \sum_{p,q=1}^3 \epsilon(b,p,q) (-\delta_\alpha^\lambda \delta_a^p G_{w\mu}^q - \delta_\alpha^\mu \delta_a^q G_{w\lambda}^p) \\
&= \sum_{c=1}^3 \epsilon(b,a,c) (-\delta_\alpha^c G_{w\mu}^c + \delta_\alpha^c G_{w\lambda}^c) \\
\frac{\partial}{\partial G_\alpha^a} \mathcal{F}_{w\lambda\mu}^b &= \sum_{p,q=1}^3 \epsilon(b,p,q) [\delta_a^p \delta_\alpha^\lambda G_{r\mu}^q + \delta_a^q \delta_\alpha^\mu G_{r\lambda}^p] \\
&= \epsilon(b,a,c) (\delta_\alpha^c G_{r\mu}^c - \delta_\alpha^c G_{r\lambda}^c) \\
\frac{\partial}{\partial G_\alpha^a} \langle \mathcal{F}_G, \mathcal{F}_G \rangle &= -4 \sum_{\lambda=0}^3 [j(\mathcal{F}_r^{\alpha\lambda}) G_{w\lambda} + j(\mathcal{F}_w^{\alpha\lambda}) G_{r\lambda}]^a \\
\text{Using the formula in the Clifford algebra :} \\
\frac{\partial}{\partial G_\alpha^a} \langle \mathcal{F}_G, \mathcal{F}_G \rangle &= -4 [v^* ([j(\mathcal{F}_r^{\alpha\lambda}) G_{r\lambda} - j(\mathcal{F}_w^{\alpha\lambda}) G_{w\lambda}], [j(\mathcal{F}_r^{\alpha\lambda}) G_{w\lambda} + j(\mathcal{F}_w^{\alpha\lambda}) G_{r\lambda}])]^a \\
\frac{\partial}{\partial G_\alpha^a} \langle \mathcal{F}_G, \mathcal{F}_G \rangle &= -4 \sum_{\lambda=0}^3 [\mathcal{F}_G^{\alpha\lambda}, G_\lambda]^a \\
\text{iv) } \frac{dL}{dG_\alpha^a} &= C_I \mu \frac{1}{i} V^\alpha \langle \psi, [\gamma C(\vec{\kappa}_a)] [\psi] \rangle - 4C_G \sum_{\lambda=0}^3 [\mathcal{F}_G^{\alpha\lambda}, G_\lambda]^a \\
\text{v) } \frac{dL}{d\beta G_\alpha^a} &= 2 \sum_{b=1}^3 \sum_{\lambda\mu} \left(\frac{\partial}{\partial \beta G_\alpha^a} \mathcal{F}_{r\lambda\mu}^b \right) \mathcal{F}_r^{b\lambda\mu} - \left(\frac{\partial}{\partial \beta G_\alpha^a} \mathcal{F}_{w\lambda\mu}^b \right) \mathcal{F}_w^{b\lambda\mu} \\
a = 1, 2, 3 : \\
\frac{dL}{d\beta G_\alpha^a} &= 2C_G (\mathcal{F}_r^{b\beta\alpha} - \mathcal{F}_r^{b\alpha\beta}) = 4C_G \mathcal{F}_r^{a\beta\alpha} \\
a = 4, 5, 6 : \\
\frac{dL}{d\beta G_\alpha^a} &= -2C_G (\mathcal{F}_w^{b\beta\alpha} - \mathcal{F}_w^{b\alpha\beta}) = -4C_G \mathcal{F}_w^{a\beta\alpha}
\end{aligned}$$

Equations :

$$\forall \alpha = 0, \dots, 3$$

$$\forall a = 1, 2, 3$$

$$C_I \mu \frac{1}{i} V^\alpha \langle \psi, [\gamma C(\vec{\kappa}_a)] [\psi] \rangle - 4C_G \sum_{\lambda=0}^3 [\mathcal{F}_G^{\alpha\lambda}, G_\lambda]^a = \frac{1}{\det P'} 4C_G \sum_\beta \frac{d}{d\xi^\beta} (\mathcal{F}_r^{a\beta\alpha} \det P')$$

$$\forall a = 4, 5, 6$$

$$C_I \mu \frac{1}{i} V^\alpha \langle \psi, [\gamma C(\vec{\kappa}_a)] [\psi] \rangle - 4C_G \sum_{\lambda=0}^3 [\mathcal{F}_G^{\alpha\lambda}, G_\lambda]^a = -\frac{1}{\det P'} 4C_G \sum_\beta \frac{d}{d\xi^\beta} (\mathcal{F}_w^{a\beta\alpha} \det P')$$

That is :

$$\forall \alpha = 0, \dots, 3, a = 1, \dots, 6 : \sum_{\lambda=0}^3 [\mathcal{F}_G^{\alpha\lambda}, G_\lambda]^a - \frac{1}{\det P'} \sum_{\beta} \frac{d}{d\xi^\beta} (\mathcal{F}_r^{a\alpha\beta} \det P') = J_a^\alpha \quad (7.8)$$

with :

$$\forall a = 1, 2, 3 : J_a^\alpha = \frac{C_I}{4C_G} \mu \frac{1}{i} V^\alpha \langle \psi, [\gamma C(\vec{\kappa}_a)] [\psi] \rangle \quad (7.9)$$

$$\forall a = 4, 5, 6 : J_a^\alpha = -\frac{C_I}{4C_G} \mu \frac{1}{i} V^\alpha \langle \psi, [\gamma C(\vec{\kappa}_a)] [\psi] \rangle \quad (7.10)$$

7.2.4 Equation for the other fields

We will keep the general expression.

Derivatives :

$$\begin{aligned} \text{i) } \frac{dL}{d\dot{A}_\alpha^a} &= C_I \mu \frac{1}{i} \left\langle \psi, \frac{\partial}{\partial \dot{A}_\alpha^a} \nabla_V \psi \right\rangle + C_G \frac{\partial}{\partial \dot{A}_\alpha^a} \langle \mathcal{F}_A, \mathcal{F}_A \rangle \\ &\left\langle \psi, \frac{\partial}{\partial \dot{A}_\alpha^a} \nabla_V \psi \right\rangle = V^\alpha \left\langle \psi, \frac{\partial}{\partial \dot{A}_\alpha^a} [\theta_a] [\psi] \right\rangle \\ &\frac{\partial}{\partial \dot{A}_\alpha^a} \langle \mathcal{F}_A, \mathcal{F}_A \rangle \\ &= \frac{\partial}{\partial \dot{A}_\alpha^a} \sum_{b=1}^3 \sum_{pqrs} (\mathcal{F}_{Apq}^b g^{pr} g^{qs} \mathcal{F}_{rrs}^b) \\ &= 2 \sum_{b=1}^3 \sum_{\lambda\mu} \left(\frac{\partial}{\partial \dot{A}_\alpha^a} \mathcal{F}_{A\lambda\mu}^b \right) \mathcal{F}_A^{b\lambda\mu} \\ &= 2 \sum_{b=1}^3 \sum_{\lambda\mu} \left(\frac{\partial}{\partial \dot{A}_\alpha^a} [\dot{A}_\lambda, \dot{A}_\mu]^b \right) \mathcal{F}_A^{b\lambda\mu} \\ &= 2 \sum_{b=1}^3 \sum_{\lambda=0}^3 \left[\vec{\theta}_a, \dot{A}_\lambda \right]^b \mathcal{F}_A^{b\alpha\lambda} + \left(\frac{\partial}{\partial \dot{A}_\alpha^a} [\dot{A}_\lambda, \vec{\theta}_a]^b \right) \mathcal{F}_A^{b\lambda\alpha} \\ &= 4 \sum_{b=1}^3 \sum_{\lambda=0}^3 \left[\vec{\theta}_a, \dot{A}_\lambda \right]^b \mathcal{F}_A^{b\alpha\lambda} \\ &= 4 \sum_{\lambda=0}^3 \left\langle \left[\vec{\theta}_a, \dot{A}_\lambda \right], \mathcal{F}_A^{\alpha\lambda} \right\rangle \\ &= 4 \sum_{\lambda=0}^3 \left\langle \vec{\theta}_a, [\dot{A}_\lambda, \mathcal{F}_A^{\alpha\lambda}] \right\rangle \\ &= 4 \sum_{\lambda=0}^3 \left[\dot{A}_\lambda, \mathcal{F}_A^{\alpha\lambda} \right]^a \end{aligned}$$

Using : $\forall X, Y, Z \in T_1 U : \langle X, [Y, Z] \rangle = \langle [X, Y], Z \rangle$ and the fact that the basis is orthonormal.

$$\begin{aligned} \frac{dL}{d\dot{A}_\alpha^a} &= C_I \mu \frac{1}{i} V^\alpha \left\langle \psi, \frac{\partial}{\partial \dot{A}_\alpha^a} [\theta_a] [\psi] \right\rangle - 4C_A \sum_{\lambda=0}^3 \left[\mathcal{F}_A^{\alpha\lambda}, \dot{A}_\lambda \right]^a \\ \text{ii) } \frac{dL}{d\partial_\beta \dot{A}_\alpha^a} &= 2C_A \sum_{b=1}^3 \sum_{\lambda\mu} \left(\frac{\partial}{\partial \dot{A}_\alpha^a} \mathcal{F}_{A\lambda\mu}^b \right) \mathcal{F}_A^{b\lambda\mu} \\ \frac{d\mathcal{F}_{A\lambda\mu}^b}{d\partial_\beta \dot{A}_\alpha^a} &= \frac{d}{d\partial_\beta \dot{A}_\alpha^a} \left(\partial_\lambda \dot{A}_\mu^b - \partial_\mu \dot{A}_\lambda^b \right) \\ \frac{dL}{d\partial_\beta \dot{A}_\alpha^a} &= 4C_A \mathcal{F}_A^{a\beta\alpha} \end{aligned}$$

Equation

The equation reads :

$$\forall \alpha = 0 \dots 3, \forall a = 1, \dots m$$

$$C_I \mu \frac{1}{i} V^\alpha \left\langle \psi, \frac{\partial}{\partial \dot{A}_\alpha} [\theta_a] [\psi] \right\rangle - 4C_A \sum_{\lambda=0}^3 \left[\mathcal{F}_A^{\alpha\lambda}, \dot{A}_\lambda \right]^a = \frac{1}{\det P'} \sum_\beta \frac{d}{d\xi^\beta} \left(4C_A \mathcal{F}_A^{b\beta\alpha} \right)$$

That is :

$$\forall a = 1 \dots m, \alpha = 0 \dots 3 : \sum_\beta \left[\mathcal{F}_A^{\alpha\beta}, \dot{A}_\beta \right]^a - \frac{1}{\det P'} \frac{d}{d\xi^\beta} \left(\mathcal{F}_A^{a\alpha\beta} \det P' \right) = J_A^{a\alpha} \quad (7.11)$$

with

$$J_A^{a\alpha} = \frac{C_I}{4C_A} \mu V^\alpha \sum_{a=1}^m \left\langle \psi, [\psi] \frac{1}{i} [\theta_a] \right\rangle \quad (7.12)$$

7.2.5 Equation for the tetrad

The equations are :

$$\forall \alpha, i : \frac{dL \det P'}{dP_i^\alpha} = \sum_\beta \frac{d}{d\xi^\beta} \left(\frac{\partial L \det P'}{\partial \theta_\beta P_i^\alpha} \right) = 0$$

The derivative of the determinant is (Maths.490) :

$$\frac{\partial \det P'}{\partial P_i^\alpha} = - \left(\frac{1}{\det P} \right)^2 \frac{\partial \det P}{\partial P_i^\alpha} = - \left(\frac{1}{\det P} \right)^2 P_\alpha^i \det P = -P_\alpha^i \det P'$$

So the equations read :

$$\frac{dL}{dP_i^\alpha} \det P' - L (\det P') P_\alpha^i = 0$$

By product with P_i^β and summation on i :

$$\forall \alpha, \beta = 0 \dots 3 : \sum_i \frac{dL}{dP_i^\alpha} P_i^\beta - L \delta_\beta^\alpha = 0 \quad (7.13)$$

Derivatives

For the part related to the fields :

$$\frac{dL_1}{dP_i^\alpha} = \sum_{\rho\theta\lambda\mu} \frac{\partial}{\partial P_i^\alpha} \left((g^{\lambda\rho} g^{\mu\theta}) (C_G \langle \mathcal{F}_{G\lambda\mu}, \mathcal{F}_{G\rho\theta} \rangle + C_A \langle \mathcal{F}_{A\lambda\mu}, \mathcal{F}_{A\rho\theta} \rangle) \right)$$

$$\begin{aligned} \frac{\partial}{\partial P_i^\alpha} (g^{\lambda\rho} g^{\mu\theta}) &= \sum_{pqjk} \eta^{jk} \eta^{pq} \frac{\partial}{\partial P_i^\alpha} P_j^\lambda P_k^\rho P_p^\mu P_q^\theta \\ &= \sum_{pqjk} \eta^{ij} (\delta_\alpha^\lambda g^{\mu\theta} P_j^\rho + \delta_\alpha^\rho g^{\mu\theta} P_j^\lambda + \delta_\alpha^\mu g^{\lambda\rho} P_j^\theta + g^{\lambda\rho} \delta_\alpha^\theta P_j^\mu) \end{aligned}$$

$$\frac{dL_1}{dP_i^\alpha} = 4 \sum_{j\lambda\mu\theta} \eta^{ij} g^{\mu\lambda} P_j^\theta (C_G \langle \mathcal{F}_{G\alpha\mu}, \mathcal{F}_{G\theta\lambda} \rangle + C_A \langle \mathcal{F}_{A\alpha\mu}, \mathcal{F}_{A\theta\lambda} \rangle)$$

For the part related to the interactions we will keep the general expression :

$$\frac{d}{dP_i^\alpha} \left(\sum_{\alpha, i=0}^3 C_I \mu \frac{1}{i} V^i P_i^\alpha \langle \psi, \nabla_\alpha \psi \rangle \right) = C_I \mu \frac{1}{i} V^i \langle \psi, \nabla_\alpha \psi \rangle$$

Equations :

$$\begin{aligned}
& 4 \sum_{ij\lambda\mu} \eta^{ij} g^{\gamma\mu} P_j^\lambda P_i^\beta (C_G \langle \mathcal{F}_{G\alpha\gamma}, \mathcal{F}_{G\lambda\mu} \rangle + C_A \langle \mathcal{F}_{A\alpha\gamma}, \mathcal{F}_{A\lambda\mu} \rangle) \\
& + C_I \mu \frac{1}{i} \sum_i V^i P_i^\beta \langle \psi, \nabla_\alpha \psi \rangle - L \delta_\beta^\alpha = 0 \\
& 4 \sum_{ij\lambda\mu} g^{\gamma\mu} g^{\lambda\beta} (C_G \langle \mathcal{F}_{G\alpha\gamma}, \mathcal{F}_{G\lambda\mu} \rangle + C_A \langle \mathcal{F}_{A\alpha\gamma}, \mathcal{F}_{A\lambda\mu} \rangle) \\
& + C_I \mu \frac{1}{i} V^\beta \langle \psi, \nabla_\alpha \psi \rangle - L \delta_\beta^\alpha = 0 \\
& 4 \sum_\gamma \left(C_G \langle \mathcal{F}_{G\alpha\gamma}, \mathcal{F}_G^{\beta\gamma} \rangle + C_A \langle \mathcal{F}_{A\alpha\gamma}, \mathcal{F}_A^{\beta\gamma} \rangle \right) \\
& + C_I \mu \frac{1}{i} V^\beta \langle \psi, \nabla_\alpha \psi \rangle - L \delta_\beta^\alpha = 0
\end{aligned}$$

The equation reads :

$$\forall \alpha, \beta = 0 \dots 3 : C_I \mu \frac{1}{i} V^\beta \langle \psi, \nabla_\alpha \psi \rangle + 4 \sum_{\gamma=0}^3 \left(C_G \langle \mathcal{F}_{G\alpha\gamma}, \mathcal{F}_G^{\beta\gamma} \rangle + C_A \langle \mathcal{F}_{A\alpha\gamma}, \mathcal{F}_A^{\beta\gamma} \rangle \right) \tag{7.14}$$

$$= \delta_\beta^\alpha \sum_{\lambda\mu} \left(C_G \langle \mathcal{F}_{G\lambda\mu}, \mathcal{F}_G^{\lambda\mu} \rangle + C_A \langle \mathcal{F}_{A\lambda\mu}, \mathcal{F}_A^{\lambda\mu} \rangle + C_I \mu \frac{1}{i} \langle \psi, \nabla_V \psi \rangle \right) \tag{7.15}$$

By taking $\alpha = \beta$ and summing :

$$\begin{aligned}
& C_I \mu \frac{1}{i} \langle \psi, \nabla_V \psi \rangle + 4 \sum_{\alpha\gamma} (C_G \langle \mathcal{F}_{G\alpha\gamma}, \mathcal{F}_G^{\alpha\gamma} \rangle + C_A \langle \mathcal{F}_{A\alpha\gamma}, \mathcal{F}_A^{\alpha\gamma} \rangle) - 4L = 0 \\
& \text{with } L = \sum_{\lambda\mu} C_G \langle \mathcal{F}_{G\lambda\mu}, \mathcal{F}_G^{\lambda\mu} \rangle + C_A \langle \mathcal{F}_{A\lambda\mu}, \mathcal{F}_A^{\lambda\mu} \rangle + C_I \frac{1}{i} \langle \psi, \nabla_V \psi \rangle \mu \\
& - 3C_I \mu \frac{1}{i} \langle \psi, \nabla_V \psi \rangle = 0 \\
& \Rightarrow
\end{aligned}$$

$$\langle \psi, \nabla_V \psi \rangle = 0 \tag{7.16}$$

As a consequence, in the general case :

$$\frac{1}{i} \langle \psi, \nabla_V \psi \rangle = k^t \widehat{X} + \frac{1}{i} \left\langle \psi_0, [\psi_0] \left[Ad_{z\alpha} \widehat{A} \right] \right\rangle = 0$$

and with the EM field :

$$k^t \widehat{X} + \widehat{A} = 0 \tag{7.17}$$

We have all together a set of 7 first order differential equations for r, w , which defines the motion on the trajectory, with respect to the standard chart of an observer.

$$\begin{aligned}
& A_1(r) \frac{dr}{dt} + A_2(r, w) \frac{dw}{dt} + A_3(r, w) \widehat{G}_r + A_4(r, w) \widehat{G}_w = rK_r \\
& B_1(r, w) \frac{dr}{dt} + B_2(r, w) \frac{dw}{dt} + B_3(r, w) \widehat{G}_r + B_4(r, w) \widehat{G}_w \\
& = \frac{caw}{2a_w^2 - 1} \sum_{\alpha, i=1}^3 P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) \varepsilon_i + wK_w \\
& k^t [A_r(r)] \frac{dr}{dt} + k^t [A_w(r, w)] \frac{dw}{dt} + k^t [D_r(r, w)] \widehat{G}_r + k^t [D_w(r, w)] \widehat{G}_w = 0
\end{aligned}$$

These general solutions must be adjusted to the fields equations to find the trajectory. The density is deduced from the continuity equation.

We will see in the following what can be deduced from the equations related to the fields.

7.2.6 Deformable solid

One can compute the single spinor representing a deformable solid by aggregating particles. In the chart of an observer B the solid is defined by :

a space like compact hypersurface $\omega(0)$,

a section $S(t, x) = \gamma C(\sigma(t, x)) S_0 \in \mathfrak{X}(S_0)$, which provides, with respect to the observer a field V of trajectories,

a density μ which follows a continuity equation.

The spinor $S_B(t)$ of the solid is defined, in this model, by integration of σ :

$$S_B(t) = N(t) \gamma C(\sigma_B(t)) S_0$$

with

$$\sigma_B(t) = A(t) + v(R(t), W(t)) + B(t) \varepsilon_5$$

$$R(t) = \frac{1}{N} \widehat{r}(t), W(t) = \frac{1}{N} \widehat{w}(t)$$

$$A(t) = \frac{1}{N} \widehat{a}(t), B(t) = \frac{1}{N} \widehat{b}(\tau t)$$

$$N(t) = \int_{\omega(t)} \mu(t, x) \varpi_3(t, x)$$

$$\sigma(t, x) = a(t, x) + v(r(t, x), w(t, x)) + b(t, x) \varepsilon_5$$

$$\widehat{r}(t) = \int_{\omega(t)} \mu(t, x) r(t, x) \varpi_3(t, x), \widehat{w}(t) = \int_{\omega(0)} \mu(t, x) w(t, x) \varpi_3(t, x)$$

$$\widehat{a}(t) = \int_{\omega(0)} a(t, x) \mu(t, x) \varpi_3(t, x), \widehat{b}(t) = \int_{\omega(0)} b(t, x) \mu(\tau, x) \varpi_3(t, x)$$

which requires that the two conditions :

$$\widehat{a}(t) \widehat{b}(t) = -\frac{1}{4} \widehat{w}^t \widehat{r}$$

$$\widehat{a}^2 - \widehat{b}^2 = N^2 + \frac{1}{4} (\widehat{w}^t \widehat{w} - \widehat{r}^t \widehat{r})$$

$$\Rightarrow N^2 = \widehat{a}^2 - \widehat{b}^2 - \frac{1}{4} (\widehat{w}^t \widehat{w} - \widehat{r}^t \widehat{r}) > 0$$

are met, as the effect of the internal forces which keep together the material body.

It can be seen as a special case of the previous model, with the perturbative Lagrangian and a vector k deduced from S_0 . It can be used in different situations.

i) The external fields are given. Then the maps r, w, μ are deduced from the state equation and the continuity equations, with the parameters \widehat{G}, \widehat{A} , and adjustment to the initial conditions (of which $\omega(0)$). This is the study of the deformation of the body submitted to given forces.

The state and the continuity equations are ODE in $\frac{dr}{dt}, \frac{dw}{dt}$. If \widehat{G}, \widehat{A} do not depend on x (the fields are assumed to have the same value at any point of the material body) then the solutions r, w depend only on t and the initial conditions. The conditions above are then satisfied :

$a(t), b(t)$ are defined by $r(t), w(t)$ with the usual identities.

$$\widehat{r}(t) = r(t) N(t), \widehat{w}(\tau) = w(t) N(t), \widehat{a}(t) = a(t) N(t), \widehat{b}(t) = b(t) N(t)$$

$$\widehat{a}(\tau) \widehat{b}(\tau) = a(t) b(t) N(t)^2 = -\frac{1}{4} r(t)^t w(t) N(t)^2 = -\frac{1}{4} R(t)^t W(t)$$

$$\widehat{a}^2 - \widehat{b}^2 = \left(a(t)^2 - b(t)^2 \right) N(t)^2 = \left(1 + \frac{1}{4} \left(w(t)^t w(t) - r(t)^t r(t) \right) \right) N(t)^2 = N^2 + \frac{1}{4} (\widehat{w}^t \widehat{w} - \widehat{r}^t \widehat{r})$$

If \widehat{G}, \widehat{A} depend on x the conditions above are met if other fields are added, which represent the internal forces necessary to keep the cohesion of the body.

So actually the state equation can be used the other way around. The model is built explicitly on the assumption that the particles constituting the material body are represented by a matter field, so that its cohesion is kept. Then, for each value of the parameters $\widehat{G} = \widehat{G}_{ext} + \widehat{G}_{int}, \widehat{A} = \widehat{A}_{ext} + \widehat{A}_{int}$, the equation gives the deformation of the solid under external fields, and the value of the internal fields which are necessary to keep its cohesion. In particular a rigid body is such that S_B , and thus σ , is constant. As can be seen in the equation $\frac{dr}{dt} = 0, \frac{dw}{dt} = 0 \nRightarrow r = 0, w = 0$. The constant solutions depend on the value of the fields, whose total must be constant : the internal forces counterbalance the external fields. And the conditions above are limiting conditions for the internal forces.

ii) In the previous case the fields equations are ignored : the deformation of the body does not change the value of the fields. But in some cases this effect cannot be ignored and the full model is required. Then usually there is no external field : the system is self-contained (for instance it can be a star system). However we need to precise what is the observer. In the previous case ε_0 is arbitrary (it represents the trajectory of the whole body in a more general model) and the tetrad (and the metric) are given. With the full model the tetrad is deformed with respect to a chart (through P), and this is this chart which provides the external reference for the full motion. This chart is arbitrary : it will be the chart of a model in which the body will be inserted as a single spinor.

Of course such models are simplistic : collisions, with thermodynamics effects, should be accounted for. But it is of interest to know that the general formalism presented here can be used, quite easily, to model complex physical phenomena in the GR context.

7.3 MODEL WITH INDIVIDUAL PARTICLES

We consider a system of a fixed number N of particles $p=1\dots N$ interacting with the fields, represented as above :

- the state of each particle is represented by a map : $\psi_p : [0, T] \rightarrow Q [E \otimes F, \vartheta]$
- each of them occupies a location, which is a point m_p in M , and in the chart φ_o of an observer : $m_p(t) = \varphi_o(t, x_p(t)) \in \Omega_3(t)$ with

$$V_p(t) = \frac{dm_p}{dt} = \vec{v}_p + c\varepsilon_0(m_p(t))$$

- the fields are represented by their potential G_α, \dot{A}_α and their strength $\mathcal{F}_{G\alpha\beta}, \mathcal{F}_{A\alpha\beta}$

- the value of the potential of the fields at m_p is denoted as usual :

$$\widehat{A}_p^\alpha = \sum_{\alpha=0}^3 \dot{A}_\alpha^\alpha(m_p(t)) V_p^\alpha(t)$$

$$\widehat{G}_p^\alpha = \sum_{\alpha=0}^3 G_\alpha^\alpha(m_p(t)) V_p^\alpha(t)$$

If there is an external field it should be added as a parameter. Similarly if the trajectory is known, it should be incorporated in the model. What we have here is a collection of particles in equilibrium with their own field (for $N=1$ that is the ‘‘Lorentz-Dirac equation’’) and exterior fields.

We will study the equations with an action of the general form :

$$\int_\Omega \left(\sum_{\alpha\beta} C_G \left(\mathcal{F}_{r\alpha\beta}^t \mathcal{F}_r^{\alpha\beta} - \mathcal{F}_{w\alpha\beta}^t \mathcal{F}_w^{\alpha\beta} \right) + C_A \mathcal{F}_{A\alpha\beta}^t \mathcal{F}_A^{\alpha\beta} \right) \varpi_4 \\ + \sum_{p=1}^N \int_0^T \left(C_I \operatorname{Im} \langle \psi_p, \nabla_{V_p} \psi_p \rangle + C_P \sqrt{\epsilon_p \langle \psi_p, \psi_p \rangle} \right) \sqrt{-\langle V_p, V_p \rangle} dt$$

The particles are assumed to be of a type defined by a fundamental state $\psi_{0p}, p = 1\dots N$ so that :

$$\psi_p(t) = \vartheta(\sigma_p(t), \varkappa_p(t)) \psi_{0p}$$

There is no need for a density, but for a constant term : $C_P \sqrt{\epsilon_p \langle \psi_{0p}, \psi_{0p} \rangle}$ the square root accounting for the energy at rest : $M_p c^2$

We will use a detailed version of the interactions in the state equation (ψ) only, as it is only there that $r(t), w(t)$ are involved :

$$C_I \operatorname{Im} \langle \psi_p, \nabla_{V_p} \psi_p \rangle = C_I \sum_{\alpha=0}^3 V_{\alpha p} \left(k_p^t X_{\alpha p} + \dot{A}_\alpha \right) = C_I \left(k^t \widehat{X}_p + \widehat{A}_p \right)$$

7.3.1 Functional derivatives

We cannot use the Euler-Lagrange equations because the two integrals are of different order. Functional derivatives (derivative with respect to a function) are commonly used by physicists, but in an uncertain mathematical rigor. Actually their mathematical theory can be done in a very general context, using the extension of distributions on vector bundles (see Maths.30.3.2 and 34.1).

A functional : $\ell : J^r E \rightarrow \mathbb{R}$ defined on a normed subspace of sections $\mathfrak{X}(J^r E)$ of a vector bundle E has a functional derivative $\frac{\delta \ell}{\delta z}(Z_0)$ with respect to a section $Z \in \mathfrak{X}(E)$ in Z_0 if there is a distribution $\frac{\delta \ell}{\delta z}$ such that for any smooth, compactly supported $\delta Z \in \mathfrak{X}_{c\infty}(E)$:

$$\lim_{\|\delta Z\| \rightarrow 0} \left\| \ell(Z_0 + \delta Z) - \ell(Z_0) - \frac{\delta \ell}{\delta z}(Z_0) \delta Z \right\| = 0$$

Because Z and δZ are sections of E their r -jets extensions are computed by taking the partial derivatives. The key point in the definition is that only δZ ,

and not its derivatives, is involved. It is clear that the functional is stationary in Z_0 if $\frac{\delta \ell}{\delta z}(Z_0) = 0$.

When the functional is linear in Z then $\frac{\delta \ell}{\delta z} = \ell$.

When the functional is given by an integral : $\int_{\Omega} \lambda(J^r Z) d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3$ the functional derivative is the distribution :

$$\frac{\delta \ell}{\delta z}(\delta Z) = \int_{\Omega} \sum_{s=0}^r \sum_{\alpha_1 \dots \alpha_s} (-1)^s D_{\alpha_1 \dots \alpha_s} \frac{\partial \lambda}{\partial Z_{\alpha_1 \dots \alpha_s}} \delta Z d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3$$

so that we get back the Euler- Lagrange equations if all the functionals are integral similarly defined.

7.3.2 Equations for the particles

The variables ψ_p are involved in the last integral only :

$$\int_0^T \left(C_I \sum_{\alpha=0}^3 \left(k^t X_{\alpha p} + \dot{A}_{\alpha} \right) + C_P \sqrt{\epsilon_p \langle \psi_{0p}, \psi_{0p} \rangle} \right) \sqrt{-\langle V_p, V_p \rangle} dt$$

so the equations can be deduced from the Euler-Lagrange equations with the variables $r_p(t), w_p(t)$.

All the variables are valued at the location on their trajectories given by t , and specific to each particle. We will drop the index p .

We have :

$$\sqrt{-\langle V, V \rangle} = \frac{c}{2a_w^2 - 1}$$

There are two solutions : (a_r, r, a_w, w) and $(-a_r, -r, -a_w, -w)$ corresponding to opposite relativist spatial spin.

Equation for r :

The equations are :

$$\forall a = 1, 2, 3 : \sqrt{-\langle V, V \rangle} \frac{dL}{dr_a} = \frac{d}{dt} \left(\frac{dL}{d\frac{dr_a}{dt}} \sqrt{-\langle V, V \rangle} \right)$$

With :

$$\frac{dL_I}{dr_a} = \frac{d}{dr_a} \left(C_I \left(k^t \hat{X} + \hat{A} \right) \right) = C_I k^t \frac{d\hat{X}}{dr_a}$$

$$\frac{dL_I}{d\frac{dr_a}{dt}} = \frac{d}{d\frac{dr_a}{dt}} \left(C_I \left(k^t \hat{X} + \hat{A} \right) \right) = C_I k^t \frac{d\hat{X}}{d\frac{dr_a}{dt}} = C_I (k^t [A_r] \varepsilon_a)$$

they read :

$$C_I k^t \frac{d\hat{X}}{dr_a} = \frac{1}{\sqrt{-\langle V, V \rangle}} \frac{d}{dt} \left(C_I (k^t [A_r] \varepsilon_a) \sqrt{-\langle V, V \rangle} \right)$$

$$= C_I (k^t [A_r] \varepsilon_a) \frac{1}{\sqrt{-\langle V, V \rangle}} \frac{d}{dt} \left(\sqrt{-\langle V, V \rangle} \right) + \frac{d}{dt} (C_I (k^t [A_r] \varepsilon_a))$$

$$\frac{1}{\sqrt{-\langle V, V \rangle}} \frac{d}{dt} \left(\sqrt{-\langle V, V \rangle} \right) = -\frac{4a_w}{(2a_w^2 - 1)^2} \frac{da_w}{dt} (2a_w^2 - 1) = -\frac{4a_w}{2a_w^2 - 1} \frac{da_w}{dt} = -\frac{1}{2a_w^2 - 1} \left(w^t \frac{dw}{dt} \right)$$

$$k^t \frac{d\hat{X}}{dr_a} = -\frac{1}{2a_w^2 - 1} \left(w^t \frac{dw}{dt} \right) (k^t [A_r] \varepsilon_a) + \frac{d}{dt} (k^t [A_r] \varepsilon_a)$$

The quantities have been computed in the previous section.

$$k^t \frac{d\hat{X}}{dr_a} = -\frac{1}{4a_r} (\varepsilon_a^t r) \left(k^t \frac{\partial \hat{X}}{\partial a_r} \right)$$

$$+ \varepsilon_a^t \left\{ [\delta A_r(r, k)] \frac{dr}{dt} + [\delta B_r(r, k)] \left([B_w] \frac{dw}{dt} + [B_{wr}] \hat{G}_r + [B_{ww}] \hat{G}_w \right) \right.$$

$$\left. + [\delta C_r(r, k)] \left([C_w] \frac{dw}{dt} + [C_{wr}] \hat{G}_r + [C_{ww}] \hat{G}_w \right) \right\}$$

The equation reads :

$$\begin{aligned} & -\frac{1}{4a_r} (\varepsilon_a^t r) \left(k^t \frac{\partial \widehat{X}}{\partial a_r} \right) \\ & + \varepsilon_a^t \{ [\delta A_r (r, k)] \frac{dr}{dt} + ([\delta B_r (r, k)] [B_w] + [\delta C_r (r, k)] [C_w]) \frac{dw}{dt} \\ & + [A_3] \widehat{G}_r + [A_4] \widehat{G}_w \} \\ & = -\frac{1}{2a_w^2 - 1} (w^t \frac{dw}{dt}) \left(\varepsilon_a^t [A_r]^t k \right) + \frac{d}{dt} \left(\varepsilon_a^t [A_r]^t k \right) \end{aligned}$$

with :

$$\begin{aligned} [A_3] &= [\delta B_r (r, k)] [B_{wr}] + [\delta C_r (r, k)] [C_{wr}] \\ [A_4] &= [\delta B_r (r, k)] [B_{ww}] + [\delta C_r (r, k)] [C_{ww}] \end{aligned}$$

$$\begin{aligned} & -\frac{1}{4a_r} r \left(k^t \frac{\partial \widehat{X}}{\partial a_r} \right) + [\delta A_r (r, k)] \frac{dr}{dt} + ([\delta B_r (r, k)] [B_w] + [\delta C_r (r, k)] [C_w]) \frac{dw}{dt} + \\ & [A_3] \widehat{G}_r + [A_4] \widehat{G}_w \\ & = -\frac{1}{2a_w^2 - 1} (w^t \frac{dw}{dt}) [A_r]^t k + \frac{d}{dt} \left([A_r]^t k \right) \end{aligned}$$

By product with r^t :

$$\begin{aligned} & -\frac{1}{4a_r} r^t r \left(k^t \frac{\partial \widehat{X}}{\partial a_r} \right) \\ & = -r^t [\delta A_r (r, k)] \frac{dr}{dt} - r^t ([\delta B_r (r, k)] [B_w] + [\delta C_r (r, k)] [C_w]) \frac{dw}{dt} \\ & + [A_3] \widehat{G}_r + [A_4] \widehat{G}_w - \frac{1}{2a_w^2 - 1} (w^t \frac{dw}{dt}) r^t [A_r]^t k + r^t \frac{d}{dt} \left([A_r]^t k \right) \end{aligned}$$

and substitution as above :

$$\begin{aligned} & [\delta A_r (r, k)] \frac{dr}{dt} + ([\delta B_r (r, k)] [B_w] + [\delta C_r (r, k)] [C_w]) \frac{dw}{dt} + [A_3] \widehat{G}_r + [A_4] \widehat{G}_w \\ & + \frac{1}{2a_w^2 - 1} (w^t \frac{dw}{dt}) [A_r]^t k - \frac{d}{dt} \left([A_r]^t k \right) = rK_r \\ & \frac{d}{dt} \left([A_r]^t k \right) = \left(-\frac{1}{4a_r} \frac{\partial [A_r]^t}{\partial a_r} (j(r) j(k) + k^t r) + [\delta A_r^t (r, k)]^t \right) \frac{dr}{dt} \\ & \frac{1}{2a_w^2 - 1} (w^t \frac{dw}{dt}) [A_r]^t k = \frac{1}{2a_w^2 - 1} [A_r]^t k w^t \frac{dw}{dt} = \frac{1}{2a_w^2 - 1} [A_r]^t (j(w) j(k) + k^t w) \frac{dw}{dt} \end{aligned}$$

The equations sums up to :

$$\begin{aligned} & \left([\delta A_r (r, k)] - [\delta A_r^t (r, k)]^t + \frac{1}{4a_r} \frac{\partial [A_r]^t}{\partial a_r} (j(r) j(k) + k^t r) \right) \frac{dr}{dt} \\ & + \left([\delta B_r (r, k)] [B_w] + [\delta C_r (r, k)] [C_w] + \frac{1}{2a_w^2 - 1} [A_r]^t (j(w) j(k) + k^t w) \right) \frac{dw}{dt} \\ & + [A_3] \widehat{G}_r + [A_4] \widehat{G}_w = rK_r \end{aligned}$$

$$A_1 (r) \frac{dr}{dt} + A_2 (r, w) \frac{dw}{dt} + A_3 (r, w) \widehat{G}_r + A_4 (r, w) \widehat{G}_w = rK_r \quad (7.18)$$

The coefficients are the same as in the previous section, except A_2 :

$$\begin{aligned} [A_2 (r, w)] &= ([\delta B_r (r, k)] [B_w] + [\delta C_r (r, k)] [C_w]) + \frac{1}{2a_w^2 - 1} [A_r]^t (j(w) j(k) + k^t w) \\ &= \frac{1}{2} a_r j(k) j(w) + \frac{1}{4} (j(r) j(k) - 2j(k) j(r)) \left[\frac{1}{a_w} + \frac{3}{2} j(w) - \frac{1}{4a_w} j(w) j(w) \right] \\ &+ \frac{1}{2a_w^2 - 1} \left[\frac{1}{a_r} - \frac{2-5a_r^2}{2} j(r) + \frac{5a_r^2+1}{4a_r} j(r) j(r) \right] (j(w) j(k) + k^t w) \end{aligned}$$

Equation for w :

We have similarly the equation :

$$\forall a = 1, 2, 3 : \sqrt{-\langle V, V \rangle} \frac{dL}{dw_a} = \frac{d}{dt} \left(\frac{dL}{d\frac{dw_a}{dt}} \sqrt{-\langle V, V \rangle} \right)$$

with :

$$\begin{aligned} \frac{dL_I}{dw_a} &= \frac{d}{dw_a} \left(\sum_{\alpha=0}^3 V^\alpha C_I \left(k^t X_\alpha + \dot{A}_\alpha \right) \right) = C_I k^t \left(\sum_{\alpha=0}^3 \frac{dV^\alpha}{dw_a} \left(k^t X_\alpha + \dot{A}_\alpha \right) + V^\alpha \frac{dX_\alpha}{dw_a} \right) \\ &= \varepsilon_a^t \sum_{\alpha=0}^3 \left(w \frac{c(2a_w^2+1)}{4a_w(2a_w^2-1)^2} \sum_{i=1}^3 w_i P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) + \frac{ca_w}{2a_w^2-1} \sum_{i=1}^3 P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) \varepsilon_i \right) \\ &\quad \sum_{\alpha=0}^3 V^\alpha k^t \frac{dX_\alpha}{dw_a} \\ &= (\varepsilon_a^t w) \left(k^t \frac{d\hat{X}}{da_w} \right) \frac{1}{4a_w} + \varepsilon_a^t \left(\left[\delta B_w \left(w, [B_r]^t k \right) \right] + \left[\delta C_w \left(w, [C_r]^t k \right) \right] \right) \frac{dw}{dt} \\ &\quad + \varepsilon_a^t \left([B_3] \hat{G}_r + [B_4] \hat{G}_w \right) \\ \frac{dL}{d\frac{dw_a}{dt}} &= \frac{d}{d\frac{dw_a}{dt}} \left(C_I \left(k^t \hat{X} + \hat{A} \right) \right) = C_I k^t \frac{d\hat{X}}{d\frac{dw_a}{dt}} = C_I k^t \left([B_r] [B_w] + [C_r] [C_w] \right) \varepsilon_a \\ &= C_I \varepsilon_a^t \left([B_r] [B_w] + [C_r] [C_w] \right)^t k \end{aligned}$$

The equation is :

$$\begin{aligned} &w \left(\frac{2a_w^2+1}{(2a_w^2-1)^2} \frac{1}{4a_w} \sum_{\alpha=0}^3 \sum_{i=1}^3 cw_i P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) + \left(k^t \frac{d\hat{X}}{da_w} \right) \frac{1}{4a_w} \right) \\ &+ \frac{ca_w}{2a_w^2-1} \sum_{\alpha,i=1}^3 P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) \varepsilon_i \\ &+ \left[\delta B_w \left(w, [B_r]^t k \right) \right] + \left[\delta C_w \left(w, [C_r]^t k \right) \right] \frac{dw}{dt} + \left([B_3] \hat{G}_r + [B_4] \hat{G}_w \right) \\ &= \frac{1}{\sqrt{-\langle V, V \rangle}} \frac{d}{dt} \left(\left([B_r] [B_w] + [C_r] [C_w] \right)^t k \sqrt{-\langle V, V \rangle} \right) \\ &= \left([B_r] [B_w] + [C_r] [C_w] \right)^t k \frac{1}{\sqrt{-\langle V, V \rangle}} \frac{d}{dt} \left(\sqrt{-\langle V, V \rangle} \right) + \frac{d}{dt} \left(\left([B_r] [B_w] + [C_r] [C_w] \right)^t k \right) \\ &= -\frac{1}{2a_w^2-1} \left([B_r] [B_w] + [C_r] [C_w] \right)^t k \left(w^t \frac{dw}{dt} \right) + \frac{d}{dt} \left(\left([B_r] [B_w] + [C_r] [C_w] \right)^t k \right) \end{aligned}$$

By product with w^t :

$$\begin{aligned} &\frac{2a_w^2+1}{(2a_w^2-1)^2} \frac{1}{4a_w} \sum_{\alpha=0}^3 \sum_{i=1}^3 cw_i P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) + \left(k^t \frac{d\hat{X}}{da_w} \right) \frac{1}{4a_w} \\ &= \frac{1}{4(a_w^2-1)} w^t \left\{ -c \frac{a_w}{2a_w^2-1} \sum_{b,\alpha=1}^3 P_b^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) \varepsilon_b \right. \\ &\quad \left. - \left[\delta B_w \left(w, [B_r]^t k \right) \right] + \left[\delta C_w \left(w, [C_r]^t k \right) \right] \frac{dw}{dt} - \left([B_3] \hat{G}_r + [B_4] \hat{G}_w \right) \right. \\ &\quad \left. - \frac{1}{2a_w^2-1} \left([B_r] [B_w] + [C_r] [C_w] \right)^t k \left(w^t \frac{dw}{dt} \right) + \frac{d}{dt} \left(\left([B_r] [B_w] + [C_r] [C_w] \right)^t k \right) \right\} \end{aligned}$$

and substitution :

$$\begin{aligned} &\left(1 - \frac{1}{4(a_w^2-1)} w w^t \right) \times c \frac{a_w}{2a_w^2-1} \sum_{i,\alpha=1}^3 P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) \varepsilon_i \\ &= \left(1 - \frac{1}{4(a_w^2-1)} w w^t \right) \times \\ &\quad \left\{ - \left(\left[\delta B_w \left(w, [B_r]^t k \right) \right] + \left[\delta C_w \left(w, [C_r]^t k \right) \right] \right) \frac{dw}{dt} - [B_3] \hat{G}_r - [B_4] \hat{G}_w \right. \\ &\quad \left. - \frac{1}{2a_w^2-1} \left([B_r] [B_w] + [C_r] [C_w] \right)^t k \left(w^t \frac{dw}{dt} \right) + \frac{d}{dt} \left(\left([B_r] [B_w] + [C_r] [C_w] \right)^t k \right) \right\} \end{aligned}$$

The equation becomes :

$$\left(\left[\delta B_w \left(w, [B_r]^t k \right) \right] + \left[\delta C_w \left(w, [C_r]^t k \right) \right] \right) \frac{dw}{dt} + [B_3] \hat{G}_r + [B_4] \hat{G}_w$$

$$\begin{aligned}
& + \frac{1}{2a_w^2-1} ([B_r][B_w] + [C_r][C_w])^t k \left(w^t \frac{dw}{dt} \right) - \frac{d}{dt} \left(([B_r][B_w] + [C_r][C_w])^t k \right) \\
& = -c \frac{a_w}{2a_w^2-1} \sum_{i,\alpha=1}^3 P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) \varepsilon_i + w K_w \\
& \frac{1}{2a_w^2-1} ([B_r][B_w] + [C_r][C_w])^t k \left(w^t \frac{dw}{dt} \right) = \frac{1}{2a_w^2-1} ([B_r][B_w] + [C_r][C_w])^t (j(w)j(k) + k^t w) \frac{dw}{dt} \\
& \frac{d}{dt} \left(([B_r][B_w] + [C_r][C_w])^t k \right) \\
& = \left(\frac{1}{4a_w} \left(\frac{\partial B_w^t}{\partial a_w} B_r^t + \frac{\partial C_w^t}{\partial a_w} C_r^t \right) (j(w)j(k) + (k^t w)) + [\delta B_w^t(w, B_r^t k)]^t + [\delta C_w^t(w, C_r^t k)]^t \right) \frac{dw}{dt} \\
& + \left(-\frac{1}{4a_r} \left(B_w^t \frac{\partial B_r^t}{\partial a_r} + C_w^t \frac{\partial C_r^t}{\partial a_r} \right) (j(r)j(k) + (k^t r)) + B_w^t [\delta B_r(r, k)]^t + C_w^t [\delta C_r(r, k)]^t \right) \frac{dr}{dt} \\
& \text{The equations sums to :}
\end{aligned}$$

$$B_1(r, w) \frac{dr}{dt} + B_2(r, w) \frac{dw}{dt} + B_3(r, w) \widehat{G}_r + B_4(r, w) \widehat{G}_w = c \frac{a_w}{2a_w^2-1} \sum_{i,\alpha=1}^3 P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) \varepsilon_i + w K_w \quad (7.19)$$

The coefficients have the same value as in the previous section.

7.3.3 Equation for the fields

We use the more general expression for the interactions $C_I \frac{1}{i} \langle \psi, \nabla_V \psi \rangle$.

Let us consider a variation $\delta \dot{A}_\alpha^a$ of \dot{A}_α^a , given by a compactly supported map (so it has a value anywhere).

The functional derivative of the first integral is (Maths.2601), with $d\xi = d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3$:

$$\begin{aligned}
& \frac{\delta}{\delta \dot{A}_\alpha^a} \left(\int_\Omega C_A \mathcal{F}_{A\alpha\beta}^t \mathcal{F}_A^{\alpha\beta} \varpi_4 \right) \left(\delta \dot{A}_\alpha^a \right) \\
& = \int_\Omega C_A \left(\frac{\partial}{\partial \dot{A}_\alpha^a} \langle \mathcal{F}_{A\alpha\beta}, \mathcal{F}_A^{\alpha\beta} \rangle \det P' - \sum_\beta \frac{d}{d\xi^\beta} \frac{\partial}{\partial \dot{A}_\alpha^a} \left(\langle \mathcal{F}_{A\alpha\beta}, \mathcal{F}_A^{\alpha\beta} \rangle \det P' \right) \right) \left(\delta \dot{A}_\alpha^a \right) d\xi \\
& = \int_\Omega C_A \left(4 \sum_\beta \left[\dot{A}_\beta, \mathcal{F}_A^{\alpha\beta} \right]^a \det P' - \sum_\beta \frac{d}{d\xi^\beta} \left(-4 C_A \mathcal{F}_A^{\alpha\beta} \det P' \right) \right) \left(\delta \dot{A}_\alpha^a \right) d\xi \\
& = 4 C_A \int_\Omega \left(\sum_\beta \left[\dot{A}_\beta, \mathcal{F}_A^{\alpha\beta} \right]^a + \frac{1}{\det P'} \sum_\beta \frac{d}{d\xi^\beta} \left(\mathcal{F}_A^{\alpha\beta} \det P' \right) \right) \left(\delta \dot{A}_\alpha^a \right) \varpi_4
\end{aligned}$$

For the simple integral a direct computation gives the functional integral :

$$\begin{aligned}
& \frac{\delta}{\delta \dot{A}_\alpha^a} \left(\sum_{p=1}^N \int_0^T C_I \frac{1}{i} \langle \psi_p, \nabla_{V_p} \psi_p \rangle \sqrt{-\langle V_p, V_p \rangle} dt \right) \left(\delta \dot{A}_\alpha^a \right) \\
& = C_I \frac{1}{i} \sum_{p=1}^N \int_0^T \langle \psi_p, \psi_p V_p^\alpha \delta \dot{A}_\alpha^a(m_p) [\theta_a] \rangle \sqrt{-\langle V_p, V_p \rangle} dt \\
& = C_I \frac{1}{i} \sum_{p=1}^N \int_0^T V_p^\alpha \delta \dot{A}_\alpha^a(m_p) \langle \psi_p, \psi_p [\theta_a] \rangle \sqrt{-\langle V_p, V_p \rangle} dt
\end{aligned}$$

The equation reads :

$$\begin{aligned}
& \forall \delta \dot{A}_\alpha^a : \\
& 4 C_A \int_\Omega \left(\sum_\beta \left[\dot{A}_\beta, \mathcal{F}_A^{\alpha\beta} \right]^a + \frac{1}{\det P'} \sum_\beta \frac{d}{d\xi^\beta} \left(\mathcal{F}_A^{\alpha\beta} \det P' \right) \right) \left(\delta \dot{A}_\alpha^a \right) \varpi_4 \\
& + \sum_{p=1}^N \int_0^T C_I \frac{1}{i} \left(V_p^\alpha \delta \dot{A}_\alpha^a(m_p) \right) \langle \psi_p, [\psi_p] [\theta_a] \rangle \sqrt{-\langle V_p, V_p \rangle} dt = 0
\end{aligned}$$

The equation holds for any compactly smooth $\delta \dot{A}_\alpha^a$. Take $\delta \dot{A}_\alpha^a$ null outside a small tube ∂C_p enclosing the trajectory of each particle. By shrinking ∂C_p

the first integral converges to the integral along the trajectory, with its volume form $\sqrt{-\langle V_p, V_p \rangle} dt$:

$$\begin{aligned} & 4C_A \int_{\Omega} \left(\sum_{\beta} \left[\dot{A}_{\beta}, \mathcal{F}_A^{\alpha\beta} \right]^a + \frac{1}{\det P'} \sum_{\beta} \frac{d}{d\xi^{\beta}} \left(\mathcal{F}_A^{a\alpha\beta} \det P' \right) \right) \left(\delta \dot{A}_{\alpha}^a \right) \varpi_4 \\ & \rightarrow 4C_A \int_0^T \left(\sum_{\beta} \left[\dot{A}_{\beta}, \mathcal{F}_A^{\alpha\beta} \right]^a + \frac{1}{\det P'} \sum_{\beta} \frac{d}{d\xi^{\beta}} \left(\mathcal{F}_A^{a\alpha\beta} \det P' \right) \right) \left(\delta \dot{A}_{\alpha}^a (m_p(t)) \right) \sqrt{-\langle V_p, V_p \rangle} dt \end{aligned}$$

and the equation reads :

$$\forall \delta \dot{A}_{\alpha}^a :$$

$$\begin{aligned} & 4C_A \int_0^T \left(\sum_{\beta} \left[\dot{A}_{\beta}, \mathcal{F}_A^{\alpha\beta} \right]^{a'} + \frac{1}{\det P'} \sum_{\beta} \frac{d}{d\xi^{\beta}} \left(\mathcal{F}_A^{a\alpha\beta} \det P' \right) \right) \left(\delta \dot{A}_{\alpha}^a (m_p(t)) \right) \sqrt{-\langle V_p, V_p \rangle} dt \\ & + \int_0^T C_I \frac{1}{i} \left(\delta \dot{A}_{\alpha}^a (m_p(t)) \right) V_p^{\alpha} \langle \psi_p, [\psi_p] [\theta_a] \rangle \sqrt{-\langle V_p, V_p \rangle} dt = 0 \\ \forall a, \alpha : & 4C_A \sum_{\beta} \left[\dot{A}_{\beta}, \mathcal{F}_A^{\alpha\beta} \right]^a + C_I \frac{1}{i} V_p^{\alpha} \langle \psi_p, [\psi_p] [\theta_a] \rangle \sqrt{-\langle V_p, V_p \rangle} \\ & + 4C_A \frac{1}{\det P'} \sum_{\beta} \frac{d}{d\xi^{\beta}} \left(\mathcal{F}_A^{\alpha\beta} \det P' \right) = 0 \end{aligned}$$

$$\forall \alpha, a, p : \sum_{\beta} \left[\mathcal{F}_A^{\alpha\beta}, \dot{A}_{\beta} \right]^a - \frac{1}{\det P'} \sum_{\beta} \frac{d}{d\xi^{\beta}} \left(\mathcal{F}_A^{\alpha\beta} \det P' \right) = J_{Ap}^{a\alpha} \quad (7.20)$$

with :

$$J_{Ap}^{a\alpha} = \frac{C_I}{4C_A} \sqrt{-\langle V_p, V_p \rangle} V_p^{\alpha} \sum_{a=1}^m \left\langle \psi_p, [\psi_p] \frac{1}{i} [\theta_a] \right\rangle \quad (7.21)$$

So we have the same equation as in the first model and μ disappears. We have similarly :

$$\forall \alpha, a, p : \sum_{\beta} \left[\mathcal{F}_G^{\alpha\beta}, G_{\beta} \right]^a - \frac{1}{\det P'} \sum_{\beta} \frac{d}{d\xi^{\beta}} \left(\mathcal{F}_G^{a\alpha\beta} \det P' \right) = J_{Gp}^{a\alpha} \quad (7.22)$$

$$a = 1, 2, 3 : J_{Gp}^{a\alpha} = \frac{C_I}{4C_G} \sqrt{-\langle V_p, V_p \rangle} V_p^{\alpha} \langle \psi_p, \gamma C(\vec{\kappa}_a) \psi_p \rangle \quad (7.23)$$

$$a = 4, 5, 6 : J_{Gp}^{a\alpha} = -\frac{C_I}{4C_G} \sqrt{-\langle V_p, V_p \rangle} V_p^{\alpha} \langle \psi_p, \gamma C(\vec{\kappa}_a) \psi_p \rangle \quad (7.24)$$

These equations holds only on the trajectories : $m = m_p(t)$

7.3.4 Tetrad equation

We have to compute the functional derivative on both integrals.

The functional derivative reads for the first :

$$\begin{aligned} & \frac{\delta}{\delta P_i^{\alpha}} \left(\int_{\Omega} \sum_{\lambda\mu} \left(C_G \langle \mathcal{F}_{G\lambda\mu}, \mathcal{F}_G^{\lambda\mu} \rangle + C_A \langle \mathcal{F}_{A\lambda\mu}, \mathcal{F}_A^{\lambda\mu} \rangle \right) \varpi_4 \right) (\delta P_i^{\alpha}) \\ & = 4 \sum_{j\gamma\lambda\mu} \int_{\Omega} \eta^{ij} P_j^{\lambda} g^{\mu\gamma} \left(C_G \langle \mathcal{F}_{G\alpha\mu}, \mathcal{F}_{G\lambda\gamma} \rangle + C_A \langle \mathcal{F}_{A\alpha\mu}, \mathcal{F}_{A\lambda\gamma} \rangle \right) (\delta P_i^{\alpha}) \varpi_4 \\ & - \int_{\Omega} \left(\sum_{\lambda\mu} \left(C_G \langle \mathcal{F}_{G\lambda\mu}, \mathcal{F}_G^{\lambda\mu} \rangle + C_A \langle \mathcal{F}_{A\lambda\mu}, \mathcal{F}_A^{\lambda\mu} \rangle \right) \right) (P_i^{\alpha}) (\delta P_i^{\alpha}) \varpi_4 \end{aligned}$$

(to account for the derivative with respect to $\det P'$)

For the part related to the interactions V is defined by $V^\alpha = \sum_{i=0}^3 P_i^\alpha V^i$ and $\sqrt{-\langle V_p, V_p \rangle} = \frac{c}{2a_{w_p}^2 - 1}$

$$\begin{aligned} C_I \frac{1}{i} \langle \psi, \nabla_V \psi \rangle &= \sum_{\alpha, i=0}^3 C_I \frac{1}{i} V^i P_i^\alpha \langle \psi, \nabla_\alpha \psi \rangle \\ \frac{\delta}{\delta P_i^\alpha} \int_0^T \left(C_I \frac{1}{i} \langle \psi_p, \nabla_{V_p} \psi_p \rangle + C_P \sqrt{\epsilon_p \langle \psi_{0p}, \psi_{0p} \rangle} \right) \sqrt{-\langle V_p, V_p \rangle} dt \\ &= C_I \frac{1}{i} \int_0^T V^i \langle \psi_p, \nabla_\alpha \psi_p \rangle \sqrt{-\langle V_p, V_p \rangle} dt \end{aligned}$$

Thus :

$$\begin{aligned} \delta \mathcal{L} &= 4 \sum_{j\gamma\lambda\mu} \int_\Omega \eta^{ij} P_j^\lambda g^{\mu\gamma} (C_G \langle \mathcal{F}_{G\alpha\mu}, \mathcal{F}_{G\lambda\gamma} \rangle + C_A \langle \mathcal{F}_{A\alpha\mu}, \mathcal{F}_{A\lambda\gamma} \rangle) (\delta P_i^\alpha) \varpi_4 \\ &\quad - \int_\Omega \left(\sum_{\lambda\mu} \left(C_G \langle \mathcal{F}_{G\lambda\mu}, \mathcal{F}_G^{\lambda\mu} \rangle + C_A \langle \mathcal{F}_{A\lambda\mu}, \mathcal{F}_A^{\lambda\mu} \rangle \right) \right) (P_i^\alpha) (\delta P_i^\alpha) \varpi_4 \\ &\quad + \sum_{p=1}^N C_I \frac{1}{i} \int_0^T V^i \langle \psi_p, \nabla_\alpha \psi_p \rangle (\delta P_i^\alpha) \sqrt{-\langle V_p, V_p \rangle} dt \end{aligned}$$

And the equation $\frac{\delta \mathcal{L}}{\delta P_i^\alpha} (\delta P_i^\alpha) = 0$ reads, for the solutions :

$\forall \delta P_i^\alpha :$

$$\begin{aligned} 4 \sum_{j\gamma\lambda\mu} \int_\Omega \eta^{ij} P_j^\lambda g^{\mu\gamma} (C_G \langle \mathcal{F}_{G\alpha\mu}, \mathcal{F}_{G\lambda\gamma} \rangle + C_A \langle \mathcal{F}_{A\alpha\mu}, \mathcal{F}_{A\lambda\gamma} \rangle) (\delta P_i^\alpha) \varpi_4 \\ - \int_\Omega \left(\sum_{\lambda\mu} \left(C_G \langle \mathcal{F}_{G\lambda\mu}, \mathcal{F}_G^{\lambda\mu} \rangle + C_A \langle \mathcal{F}_{A\lambda\mu}, \mathcal{F}_A^{\lambda\mu} \rangle \right) \right) (P_i^\alpha) (\delta P_i^\alpha) \varpi_4 \\ + \sum_{p=1}^N C_I \frac{1}{i} \int_0^T V^i \langle \psi_p, \nabla_\alpha \psi_p \rangle (\delta P_i^\alpha) \sqrt{-\langle V_p, V_p \rangle} dt = 0 \end{aligned}$$

With the same reasoning as above, for each particle :

$$\begin{aligned} 4 \sum_{j\gamma\lambda\mu} \int_0^T \eta^{ij} P_j^\lambda g^{\mu\gamma} (C_G \langle \mathcal{F}_{G\alpha\mu}, \mathcal{F}_{G\lambda\gamma} \rangle + C_A \langle \mathcal{F}_{A\alpha\mu}, \mathcal{F}_{A\lambda\gamma} \rangle) (\delta P_i^\alpha) \sqrt{-\langle V_p, V_p \rangle} dt \\ - \int_0^T \left(\sum_{\lambda\mu} \left(C_G \langle \mathcal{F}_{G\lambda\mu}, \mathcal{F}_G^{\lambda\mu} \rangle + C_A \langle \mathcal{F}_{A\lambda\mu}, \mathcal{F}_A^{\lambda\mu} \rangle \right) \right) (P_i^\alpha) (\delta P_i^\alpha) \sqrt{-\langle V_p, V_p \rangle} dt \\ + C_I \frac{1}{i} \int_0^T V^i \langle \psi_p, \nabla_\alpha \psi_p \rangle (\delta P_i^\alpha) \sqrt{-\langle V_p, V_p \rangle} dt = 0 \end{aligned}$$

$$\begin{aligned} 4 \sum_{j\gamma\lambda\mu} \eta^{ij} P_j^\lambda g^{\mu\gamma} (C_G \langle \mathcal{F}_{G\alpha\mu}, \mathcal{F}_{G\lambda\gamma} \rangle + C_A \langle \mathcal{F}_{A\alpha\mu}, \mathcal{F}_{A\lambda\gamma} \rangle) \\ - \left(\sum_{\lambda\mu} \left(C_G \langle \mathcal{F}_{G\lambda\mu}, \mathcal{F}_G^{\lambda\mu} \rangle + C_A \langle \mathcal{F}_{A\lambda\mu}, \mathcal{F}_A^{\lambda\mu} \rangle \right) \right) (P_i^\alpha) \\ + C_I \frac{1}{i} V^i \langle \psi_p, \nabla_\alpha \psi_p \rangle = 0 \end{aligned}$$

The equation, by product with P_i^β and summation on i gives :

$$\begin{aligned} 4 \sum_{\lambda\mu} g^{\lambda\beta} g^{\mu\gamma} (C_G \langle \mathcal{F}_{G\alpha\mu}, \mathcal{F}_{G\lambda\gamma} \rangle + C_A \langle \mathcal{F}_{A\alpha\mu}, \mathcal{F}_{A\lambda\gamma} \rangle) \\ - \delta_\alpha^\beta \left(\sum_{\lambda\mu} \left(C_G \langle \mathcal{F}_{G\lambda\mu}, \mathcal{F}_G^{\lambda\mu} \rangle + C_A \langle \mathcal{F}_{A\lambda\mu}, \mathcal{F}_A^{\lambda\mu} \rangle \right) \right) \\ + C_I \frac{1}{i} V^\beta \langle \psi_p, \nabla_\alpha \psi_p \rangle = 0 \end{aligned}$$

$$4 \sum_{\mu} \left(C_G \langle \mathcal{F}_{G\alpha\mu}, \mathcal{F}_G^{\beta\mu} \rangle + C_A \langle \mathcal{F}_{A\alpha\mu}, \mathcal{F}_A^{\beta\mu} \rangle \right) + C_I \frac{1}{i} V^\beta \langle \psi_p, \nabla_\alpha \psi_p \rangle \quad (7.25)$$

$$= \delta_\alpha^\beta \sum_{\lambda\mu} \left(C_G \langle \mathcal{F}_{G\lambda\mu}, \mathcal{F}_G^{\lambda\mu} \rangle + C_A \langle \mathcal{F}_{A\lambda\mu}, \mathcal{F}_A^{\lambda\mu} \rangle \right) \quad (7.26)$$

With $\alpha = \beta$ and summing :

$$\begin{aligned} 4 \sum_{\alpha\lambda} (C_G \langle \mathcal{F}_{G\alpha\lambda}, \mathcal{F}_G^{\alpha\lambda} \rangle + C_A \langle \mathcal{F}_{A\alpha\lambda}, \mathcal{F}_A^{\alpha\lambda} \rangle) + C_I \frac{1}{i} \langle \psi_p, \nabla_{V_p} \psi_p \rangle \\ - 4 \sum_{\lambda\mu} \left(C_G \langle \mathcal{F}_{G\lambda\mu}, \mathcal{F}_G^{\lambda\mu} \rangle + C_A \langle \mathcal{F}_{A\lambda\mu}, \mathcal{F}_A^{\lambda\mu} \rangle \right) \\ + 4 C_I \frac{1}{i} \langle \psi_p, \nabla_{V_p} \psi_p \rangle = 0 \end{aligned}$$

⇒

$$\langle \psi_p, \nabla_V \psi_p \rangle = 0 \quad (7.27)$$

and as a consequence :

$$k^t \widehat{X}_p + \widehat{A}_p = 0 \quad (7.28)$$

The equations are (up to the density μ) identical or similar to the previous model. However here they hold on the trajectories : the coordinates of $m(t)$ are in the chart of the observer : $m = \varphi_0(t, x(\xi^1, \xi^2, \xi^3))$ and $x(t)$ is related to t . So the equations are differential equations with variables $r(t), w(t), \varkappa(t), \widehat{G}(t), \widehat{A}(t)$. We cannot tell anything outside the trajectories. But the previous model holds pointwise in the vacuum so can be used to compute G and \dot{A} , and the solutions must be adjusted, as well as with the initial conditions.

7.3.5 Particle submitted to an external field

The equations hold, but we have to add the external fields \dot{A}_{ext}, G_{ext} to the field due to the particle itself. The equations read :

$$\begin{aligned} A_1(r) \frac{dr}{dt} + A_2(r, w) \frac{dw}{dt} + A_3(r, w) \widehat{G}_r + A_4(r, w) \widehat{G}_w &= rK_r \\ B_1(r, w) \frac{dr}{dt} + B_2(r, w) \frac{dw}{dt} + B_3(r, w) \widehat{G}_r + B_4(r, w) \widehat{G}_w \\ &= \frac{ca_w}{2a_w^2 - 1} \sum_{\alpha, i=1}^3 P_i^\alpha \left(k^t X_\alpha + \dot{A}_\alpha \right) \varepsilon_i + wK_w \\ k^t [A_r(r)] \frac{dr}{dt} + k^t [A_w(r, w)] \frac{dw}{dt} + k^t [D_r(r, w)] \widehat{G}_r + k^t [D_w(r, w)] \widehat{G}_w &= 0 \end{aligned}$$

They are impressive but, using the many algebraic properties of the operator j , they can be solved. As an example, if σ_w is fixed (for instance the particle is bonded) $w = 0$. Without gravitation, the equations are :

$$\begin{aligned} \frac{1}{a_r} k^t \left[1 + \left(1 - \frac{5}{2} a_r^2 \right) j(r) + \frac{1}{4} (5a_r^2 + 1) j(r) j(r) \right] \frac{dr}{dt} + \widehat{A} &= 0 \\ \frac{1}{4a_r} \{ (k^t r) \left(-\frac{1}{a_r^2} + 5a_r j(r) + \frac{5a_r^2 - 1}{4a_r^2} j(r) j(r) \right) \right. \\ \left. - 20a_r (1 - a_r^2) j(k) + \frac{-3 + a_r^2 + 20a_r^4}{a_r^2} j(r) j(k) - 3(5a_r^2 + 1) j(k) j(r) \right] \frac{dr}{dt} &= rK_r \end{aligned}$$

We look for solutions such that : $r = \lambda(t) k$ with a scalar function λ .

The first equation gives :

$$\begin{aligned} \frac{1}{a_r} \|k\|^2 \frac{d\lambda}{dt} + \widehat{A} &= 0 \\ \frac{\|k\|^2}{\sqrt{1 - \frac{1}{4}\lambda^2 \|k\|^2}} \frac{d\lambda}{dt} &= -\dot{A} \end{aligned}$$

With : $\lambda = \frac{2}{\|k\|} \sin \omega(t)$ we get : $\varpi' = -\frac{1}{2\|k\|} \dot{A}$

The second equation becomes :

$$\begin{aligned} -\frac{1}{4a_r^3} \|k\|^2 \frac{d\lambda}{dt} k &= kK_r \\ K_r &= -\frac{1}{4(\cos \omega)^{3/2}} \|k\|^2 \frac{2}{\|k\|} \varpi' \cos \varpi = -\frac{\|k\| \varpi'}{2\sqrt{|\cos \omega|}} \end{aligned}$$

The motion is a rotation with a speed $-\frac{1}{2\|k\|}\dot{A}$. This can be seen as a precession with gyromagnetic ratio $\gamma = -\frac{1}{2\|k\|} \sim -\frac{e\hbar}{2m}$. The vector k can be seen as a magnetic moment.

These models are useful : after all they hold for any kind of particles, submitted to any field. They can represent beams of particles used in experiments. They are similar to the Standard Model, with GR and gravitation added, and a much simpler formalism. They hold whenever there is no particle, that is in the vacuum. And they can be used for simple problems such as the magnetic moment. We will use them in the following to introduce some physical concepts, and mathematical tools. But we have to keep in mind that these models represent ideal physical cases : no collision, no discontinuity, no change in the number or the characteristics of the particles. By analogy with fluid mechanics they represent steady flows. But, as real flows are more complicated, but also more interesting, with vortex, Particle Mechanics must consider models allowing for collisions or creation / annihilation of particles.

We will address these issues in the next chapter, but we need more tools before that.

7.4 CURRENTS

The Noether currents are usually introduced through the equivariance of the Lagrange equations, by computing the effects of a change of gauge or chart on the lagrangian. This is exactly what we have done before, and from that we have deduced some basic rules for the specification of the lagrangian, and identities which must be satisfied by the partial derivatives. Whenever the lagrangian is defined from geometric quantities these identities are met, and the Noether currents do not appear this way. But we have a more interesting, and more intuitive, view of the currents, from the equations that we have computed with the perturbative lagrangian.

7.4.1 Definition

The equations for the force fields (usually called the equations of motion) are :

$$\begin{aligned} \forall \alpha = 0 \dots 3 : \\ \forall a = 1 \dots 6 : \sum_{\beta} \left[\mathcal{F}_G^{\alpha\beta}, G_{\beta} \right]^a - \frac{1}{\det P'} \sum_{\beta} \partial_{\beta} \left(\mathcal{F}_G^{a\alpha\beta} \det P' \right) &= J_G^{a\alpha} \\ \forall a = 1 \dots m : \sum_{\beta} \left[\mathcal{F}_A^{\alpha\beta}, \dot{A}_{\beta} \right]^a - \frac{1}{\det P'} \sum_{\beta} \partial_{\beta} \left(\mathcal{F}_A^{a\alpha\beta} \det P' \right) &= J_A^{a\alpha} \end{aligned}$$

$J_G^{a\alpha}, J_A^{a\alpha}$ are components of tensors :

$$J_A = \frac{C_I}{4C_A} \mu \sum_{a=1}^m \left\langle \psi, [\psi] \frac{1}{i} [\theta_a] \right\rangle V \otimes \vec{\theta}^a \in TM \otimes T_1 U^* \quad (7.29)$$

$$a = 1, 2, 3 : J_G = \frac{C_I}{4C_G} \mu \frac{1}{i} \langle \psi, \gamma C(\vec{\kappa}_a) \psi \rangle V \otimes \vec{\kappa}^a \in TM \otimes T_1 Spin(3, 1)^* \quad (7.30)$$

$$a = 4, 5, 6 : J_G = -\frac{C_I}{4C_G} \mu \frac{1}{i} \langle \psi, \gamma C(\vec{\kappa}_a) \psi \rangle V \otimes \vec{\kappa}^a \in TM \otimes T_1 Spin(3, 1)^* \quad (7.31)$$

Similarly are tensors the quantities :

$$\phi_G = \sum_{\beta} \left[\mathcal{F}_G^{\alpha\beta}, G_{\beta} \right] \otimes \partial \xi_{\alpha} \in T_1 Spin(3, 1)^* \otimes TM \quad (7.32)$$

$$\phi_A = \sum_{\beta} \left[\mathcal{F}_A^{\alpha\beta}, \dot{A}_{\beta} \right] \otimes \partial \xi_{\alpha} \in T_1 U^* \otimes TM \quad (7.33)$$

because $\sum_{\alpha\beta} \frac{\partial L}{\partial \mathcal{F}_A^{\alpha\beta}} \partial \xi_{\alpha} \wedge \partial \xi_{\beta} \otimes \vec{\theta}^a, \sum_{\alpha\beta} \frac{\partial L}{\partial \mathcal{F}_G^{\alpha\beta}} \partial \xi_{\alpha} \wedge \partial \xi_{\beta} \otimes \vec{\kappa}^a$ are tensors (see lagrangian).

In the second model the quantities are the same, but the density μ is replaced by $\sqrt{-\langle V_p, V_p \rangle}$ and they hold for each particle.

A computation (see Annex) gives :

$$a = 1, 2, 3 : J_G^a = \frac{C_I}{4C_G} \mu k^t [D_r(r, w)] \varepsilon_a V^{\alpha}$$

$$a = 4, 5, 6 : J_G^a = -\frac{C_I}{4C_G} \mu k^t [D_w(r, w)] \varepsilon_a V^\alpha$$

$$J_G = \frac{C_I}{4C_G} \mu V \otimes v^* (k^t [D_r(r, w)], -k^t [D_w(r, w)]) \quad (7.34)$$

where $[D_r(r, w)]$, $[D_w(r, w)]$ and k have the same definition as in the total connection.

$$\begin{aligned} \phi_G &= \sum_\beta \left[\mathcal{F}_G^{\alpha\beta}, G_\beta \right] \otimes \partial\xi_\alpha \\ &= \sum_{\beta\lambda\mu} g^{\lambda\alpha} g^{\beta\mu} [\mathcal{F}_{G\lambda\mu}, G_\beta] \otimes \partial\xi_\alpha \\ &= \sum_{\beta\lambda\mu} g^{\lambda\alpha} g^{\beta\mu} [v(\mathcal{F}_{r\lambda\mu}, \mathcal{F}_{w\lambda\mu}), v(G_{r\beta}, G_{w\beta})] \otimes \partial\xi_\alpha \\ &= \sum_{\beta\lambda\mu} g^{\lambda\alpha} g^{\beta\mu} v(j(\mathcal{F}_{r\lambda\mu}) G_{r\beta} - j(\mathcal{F}_{w\lambda\mu}) G_{w\beta}, j(\mathcal{F}_{w\lambda\mu}) G_{r\beta} + j(\mathcal{F}_{r\lambda\mu}) G_{w\beta}) \otimes \partial\xi_\alpha \end{aligned}$$

$$\phi_G = \sum_\beta v^* (j(\mathcal{F}_r^{\alpha\beta}) G_{r\beta} - j(\mathcal{F}_w^{\alpha\beta}) G_{w\beta}, j(\mathcal{F}_w^{\alpha\beta}) G_{r\beta} + j(\mathcal{F}_r^{\alpha\beta}) G_{w\beta}) \otimes \partial\xi_\alpha \quad (7.35)$$

For the EM field :

$$J_A = \frac{C_I}{4C_A} \mu \langle \psi_0, \psi_0 \rangle V$$

J_G, J_A, ϕ_G, ϕ_A are vector fields over TM, valued in the Lie algebras, called the **currents** associated to the fields labeled by a . There is one collection of currents associated to each particle, and *their support is the trajectory of the particles*. For individual particles the equality is met at each point of the trajectory, thus we can say that the current ϕ follows the same trajectory as the particle.

There is one vector field for each value of a , 12 in the Standard Model for the other fields and 6 for gravitation. This latter result does not come from the choice of a connection different from the Levi-Civita connection (\mathcal{F} corresponds to the Riemann tensor) but from the choice of the scalar product over the scalar curvature. In many ways this is more in accordance with the law of equivalence, and in a GUT we would have strictly the same equations.

In a time reversal, given by the matrix

$$T = \begin{bmatrix} 0 & i\sigma_0 \\ i\sigma_0 & 0 \end{bmatrix}$$

particles are exchanged with antiparticles, and the scalar products in the currents take the opposite sign, so we have opposite currents.

The interaction term in the lagrangian reads :

$$\begin{aligned} C_I \mu \operatorname{Im} \langle \psi, \nabla_V \psi \rangle &= C_I \mu \frac{1}{i} \langle \psi, \nabla_V \psi \rangle \\ &= C_I \mu \frac{1}{i} \left\langle \psi, \frac{d\psi}{dt} \right\rangle + 4C_A \sum_{a=1}^m \dot{A}_\alpha^a J_A^a + 4C_G \sum_{a=1}^3 G_{r\alpha}^a J_G^a - G_{w\alpha}^a J_G^{a+3} \end{aligned}$$

so we have an expression using the currents, which is more familiar.

7.4.2 Main theorem

We have introduced previously the scalar product of forms on M, and the Hodge dual. The Hodge dual $*\mathcal{F}^a$ of a 2-form is a 2 form. The exterior differential $d(*\mathcal{F}^a)$ is a 3 form, which reads :

$d(*\mathcal{F}^a) = \sum_{\beta=0}^3 (-1)^\alpha \partial_\beta (\mathcal{F}^{a\alpha\beta} \det P') d\xi^0 \wedge \dots \widehat{d\xi^\alpha} \dots \wedge d\xi^3$ where $\widehat{}$ means that the vector is skipped ¹.

The equations can be written :

$$\begin{aligned} \forall a, \alpha : \sum_\beta \left[\mathcal{F}_A^{\alpha\beta}, \dot{A}_\beta \right]^a - J_A^{a\alpha} &= \frac{1}{\det P'} \sum_\beta \partial_\beta \left(\mathcal{F}_A^{a\alpha\beta} \det P' \right) \\ \forall a : \sum_{\alpha=0}^3 (-1)^\alpha \left(\sum_\beta \left[\mathcal{F}_A^{\alpha\beta}, \dot{A}_\beta \right]^a - J_A^{a\alpha} \right) \det P' d\xi^0 \wedge \dots \widehat{d\xi^\alpha} \dots \wedge d\xi^3 \\ &= \sum_{\alpha=0}^3 (-1)^\alpha \sum_\beta \partial_\beta \left(\mathcal{F}_A^{a\alpha\beta} \det P' \right) d\xi^0 \wedge \dots \widehat{d\xi^\alpha} \dots \wedge d\xi^3 \\ &= d(*\mathcal{F}_A^a) \end{aligned}$$

Moreover :

$$\sum_{\alpha=0}^3 (-1)^\alpha \left(\sum_\beta \left[\mathcal{F}_G^{\alpha\beta}, G_\beta \right]^a \right) (\det P') d\xi^0 \wedge \dots \widehat{d\xi^\alpha} \dots \wedge d\xi^3 = i_{\phi_G} \varpi_4$$

$$\sum_{\alpha=0}^3 (-1)^\alpha J_G^{a\alpha} (\det P') d\xi^0 \wedge \dots \widehat{d\xi^\alpha} \dots \wedge d\xi^3 = i_{J_G} \varpi_4$$

are the values of the 4 form ϖ_4 for the vector ϕ_A, J_A .

Thus we can write the equations in the geometric form :

$$i_{J_G} \varpi_4 = i_{\phi_G} \varpi_4 + d \circ (*\mathcal{F}_G^a)$$

$$i_{J_A} \varpi_4 = i_{\phi_A} \varpi_4 + d \circ (*\mathcal{F}_A^a)$$

$i_{B_G} \varpi_4, i_{J_G} \varpi_4, i_{B_A} \varpi_4, i_{J_A} \varpi_4$ can be interpreted as the densities of the currents ϕ_G, J_G, ϕ_A, J_A and because :

$$d \circ d \circ (*\mathcal{F}_G^a) = 0, d \circ d \circ (*\mathcal{F}_A^a) = 0$$

$$d(i_{J_G} \varpi_4) = d(i_{\phi_G} \varpi_4); d(i_{J_A} \varpi_4) = d(i_{\phi_A} \varpi_4)$$

The flow of the vector $\phi_G - J_G$ through a hypersurface ∂C can be defined as $\int_{\partial C} i_{\phi_G} \varpi_4 - i_{J_G} \varpi_4$, and if C is a manifold with boundary : $\int_C d \circ (i_{\phi_G} \varpi_4 - i_{J_G} \varpi_4) = \int_{\partial C} (i_{\phi_G} \varpi_4 - i_{J_G} \varpi_4) = 0 \Leftrightarrow \int_{\partial C} i_{\phi_G} \varpi_4 = \int_{\partial C} i_{J_G} \varpi_4$

We can also express the currents by the corresponding 1 form J^*, ϕ^* , by raising the indexes with g , and proceed to the computations (Maths.1613) :

$$J = \sum_\alpha J^\alpha \partial \xi_\alpha \rightarrow J^* = \sum_{\lambda\alpha} g_{\alpha\lambda} J^\lambda d\xi^\alpha = \sum_\alpha J_\alpha^* d\xi^\alpha$$

$$J^* \rightarrow *J^* = \sum_{\alpha,\beta=0}^3 (-1)^\alpha g^{\alpha\beta} J_\beta^* (\det P') d\xi^0 \wedge \dots \widehat{d\xi^\alpha} \dots \wedge d\xi^3$$

$$= \sum_{\alpha,\beta=0}^3 (-1)^\alpha g^{\alpha\beta} g_{\alpha\beta} J^\lambda (\det P') d\xi^0 \wedge \dots \widehat{d\xi^\alpha} \dots \wedge d\xi^3$$

$$= \sum_{\alpha=0}^3 (-1)^\alpha J^\alpha (\det P') d\xi^0 \wedge \dots \widehat{d\xi^\alpha} \dots \wedge d\xi^3$$

$$i_{J_G} \varpi_4 = *J^*$$

and similarly

$$\phi_G^* = \sum_{\alpha\beta\lambda} \left[g_{\alpha\lambda} \mathcal{F}_G^{\lambda\beta}, G_\beta \right]^a d\xi^\alpha$$

$$*\phi_G^* = \sum_{\alpha=0}^3 (-1)^\alpha \left(\sum_\beta \left[\mathcal{F}_G^{\alpha\beta}, G_\beta \right]^a \right) (\det P') d\xi^0 \wedge \dots \widehat{d\xi^\alpha} \dots \wedge d\xi^3 = i_{\phi_G} \varpi_4$$

The equations read :

$$*J_G^* = *\phi_G^* + d \circ (*\mathcal{F}_G^a)$$

¹Beware. The exponent is α and not $\alpha - 1$ because the vectors are labelled 0,1,2,3 and not 1,2,3,4. A legacy of decennies of notation.

$$*J_A^* = *\phi_A^* + d \circ (*\mathcal{F}_A^a)$$

Using : $**J_G^* = -J_G^*$, $**\phi_G^* = -\phi_G^*$ and $*d \circ (*\mathcal{F}_G) = -\delta\mathcal{F}_G$ where δ is the **codifferential** (Maths.32.3.1) the equations read equivalently :

$$J_G^* = \phi_G^* + \delta(\mathcal{F}_G); J_A^* = \phi_A^* + \delta(\mathcal{F}_A)$$

For the EM field

$\phi_A^* = 0$ because $U(1)$ is abelian

$J_A^* = \frac{C_I}{4C_A} \mu \langle \psi, \psi \rangle V^*$ because $\theta_a = i$

and the equation sums up to :

$$J_A^* = \delta(\mathcal{F}_A)$$

$$\forall \alpha, \alpha : -\frac{C_I}{4C_A} \mu \langle \psi_0, \psi_0 \rangle V^* = \frac{1}{\det P'} \sum_{\beta} \partial_{\beta} \left(\mathcal{F}_A^{\alpha\beta} \det P' \right)$$

and for a single particle :

$$-\frac{C_I}{4C_A} \langle \psi_{0p}, \psi_{0p} \rangle \sqrt{-\langle V_p, V_p \rangle} V_p^* = \frac{1}{\det P'} \sum_{\beta} \partial_{\beta} \left(\mathcal{F}_A^{\alpha\beta} \det P' \right)$$

The Maxwell's equation in GR reads for a single particle : $\delta\mathcal{F}_A = \mu_0 \sqrt{-\langle V, V \rangle} q V^*$ with a universal constant μ_0 and a charge q , so we are lead to take : $\frac{C_I}{4C_A} \langle \psi_0, \psi_0 \rangle = \mu_0 q$. The sign of q depends on the sign of $\langle \psi_0, \psi_0 \rangle = \langle S_0, S_0 \rangle$ as expected.

The practical solutions in electrodynamics rely on the knowledge of the propagation of fields, notably with Liénard-Wiechert retarded potentials.

The codifferential reduces the order of a form by one. It is in some way the inverse operator of the exterior differential d . So for the fields other than EM :

$$\delta J_G^* = \frac{1}{\det P'} \sum_{\alpha\beta} \partial_{\alpha} \left(g^{\alpha\beta} J_{\beta G}^* \det P' \right) = \frac{1}{\det P'} \sum_{\alpha} \partial_{\alpha} \left(J_G^{\alpha} \det P' \right)$$

$$\delta \phi_G^* = \frac{1}{\det P'} \sum_{\alpha\beta} \partial_{\alpha} \left(g^{\alpha\beta} \left[\mathcal{F}_G^{\alpha\beta}, G_{\beta} \right]_{\alpha}^* \det P' \right) = \frac{1}{\det P'} \sum_{\alpha\beta} \partial_{\alpha} \left(\left[\mathcal{F}_G^{\alpha\beta}, G_{\beta} \right] \det P' \right)$$

$$\delta J_A^* = \frac{1}{\det P'} \sum_{\alpha\beta} \partial_{\alpha} \left(g^{\alpha\beta} J_{\beta A}^* \det P' \right) = \frac{1}{\det P'} \sum_{\alpha} \partial_{\alpha} \left(J_A^{\alpha} \det P' \right)$$

are functions.

We can take similarly the codifferential of ϕ_A^* if it is a one form (that is for the fields other than EM) :

$$\delta \phi_A^* = \frac{1}{\det P'} \sum_{\alpha\beta} \partial_{\alpha} \left(g^{\alpha\beta} \left[\mathcal{F}_G^{\alpha\beta}, G_{\beta} \right]_{\alpha}^* \det P' \right) = \frac{1}{\det P'} \sum_{\alpha\beta} \partial_{\alpha} \left(\left[\mathcal{F}_G^{\alpha\beta}, G_{\beta} \right] \det P' \right)$$

The codifferential is such that : $\delta^2 = 0$

$$\Rightarrow \delta J_G^* = \delta \phi_G^*; \delta J_A^* = \delta \phi_A^*$$

and for the fields other than EM :

$$\sum_{\alpha} \partial_{\alpha} \left(J_G^{\alpha} \det P' \right) = \sum_{\alpha} \partial_{\alpha} \left(\left[\mathcal{F}_G^{\alpha\beta}, G_{\beta} \right] \det P' \right)$$

$$\sum_{\alpha} \partial_{\alpha} \left(J_A^{\alpha} \det P' \right) = \sum_{\alpha} \partial_{\alpha} \left(\left[\mathcal{F}_A^{\alpha\beta}, A_{\beta} \right] \det P' \right)$$

The codifferential is the adjoint of the exterior differential with respect to the scalar product of forms on TM (Maths.2498). So for any 1-form λ on TM :

$$\forall \lambda \in \Lambda_1 TM : G_1(\lambda, \delta(\mathcal{F}_G)) = G_1(d\lambda, \mathcal{F}_G)$$

$$G_1(\lambda, J_G^*) = G_1(\lambda, \phi_G^*) + G_1(\lambda, \delta(\mathcal{F}_G)) = G_1(\lambda, \phi_G^*) + G_1(d\lambda, \mathcal{F}_G)$$

thus if $d\lambda = 0$: $G_1(\lambda, J_G^*) = G_1(\lambda, \phi_G^*)$ which reads :

$$\forall \mu \in \Lambda_1(M) : \mu \wedge * \lambda = G_2(\mu, \lambda) \varpi_4$$

$$G_2(\lambda, J_G^*) \varpi_4 = G_2(J_G^*, \lambda) \varpi_4 = J_G^* \wedge * \lambda$$

$$J_G^* \wedge * \lambda = \phi_G^* \wedge * \lambda$$

Take : $\lambda = df$ with any function $f \in C_1(M; \mathbb{R})$:

$$\begin{aligned} *df &= \sum_{\alpha, \beta=0}^3 (-1)^\alpha g^{\alpha\beta} \partial_\beta f (\det P') d\xi^0 \wedge \dots \widehat{d\xi^\alpha} \dots \wedge d\xi^3 \\ &(\sum_\alpha J_\alpha^* d\xi^\alpha) \wedge \left(\sum_{\alpha, \beta=0}^3 (-1)^{\alpha+1} g^{\alpha\beta} \partial_\beta f (\det P') d\xi^0 \wedge \dots \widehat{d\xi^\alpha} \dots \wedge d\xi^3 \right) \\ &= (\sum_\alpha \phi_\alpha^* d\xi^\alpha) \wedge \left(\sum_{\alpha, \beta=0}^3 (-1)^{\alpha+1} g^{\alpha\beta} \partial_\beta f (\det P') d\xi^0 \wedge \dots \widehat{d\xi^\alpha} \dots \wedge d\xi^3 \right) \end{aligned}$$

$$\sum_{\alpha\beta} J_\alpha^* g^{\alpha\beta} \partial_\beta f = \sum_{\alpha\beta} \phi_\alpha^* g^{\alpha\beta} \partial_\beta f$$

$$\sum_\beta J^\beta \partial_\beta f = \sum_\beta \phi^\beta \partial_\beta f \Leftrightarrow f'(m) J = f'(m) \phi$$

$$\forall f \in C_1(M; \mathbb{R}) : f'(m) J_G = f'(m) \phi_G; f'(m) J_A = f'(m) \phi_A$$

Take $f(m) = \xi^\alpha$ with $\alpha = 0, \dots, 3$:

$$f'(m) J = J^\alpha = f'(m) \phi_G = \phi^\alpha$$

and we have the result :

$$J_A = \phi_A; J_G = \phi_G; \delta\mathcal{F}_A = 0; \delta\mathcal{F}_G = 0$$

As we had :

$\forall a, \alpha$:

$$\phi_A^{a\alpha} - J_A^{a\alpha} = \frac{1}{\det P'} \sum_\beta \partial_\beta \left(\mathcal{F}_A^{a\alpha\beta} \det P' \right) = 0$$

$$\phi_G^{a\alpha} - J_G^{a\alpha} = \frac{1}{\det P'} \sum_\beta \partial_\beta \left(\mathcal{F}_G^{a\alpha\beta} \det P' \right) = 0$$

The Laplacian is the differential operator : $\Delta = -(d\delta + \delta d)$ (Maths.2500).

Thus :

$$\Delta\mathcal{F}_A = -\delta d\mathcal{F}_A; \Delta\mathcal{F}_G = -\delta d\mathcal{F}_G$$

$$i_{J_A} \varpi_4 = i_{\phi_A} \varpi_4 + d \circ (*\mathcal{F}_A^a) \Rightarrow d \circ (*\mathcal{F}_A^a) = 0$$

Thus for all fields :

$$\Delta\mathcal{F} = -\delta d\mathcal{F}$$

For the EM field :

$$\mathcal{F}_{EM} = d\dot{A}$$

$$i_{J_G} \varpi_4 = i_{\phi_G} \varpi_4 + d \circ (*\mathcal{F}_G^a) \Rightarrow d \circ (*\mathcal{F}_G^a) = 0$$

$$J_A^* = \delta(\mathcal{F}_A) = \delta d\dot{A} = -\Delta\dot{A}$$

To sum up :

Theorem 106 For the gravitational field in a continuous model :

$$J_G = \phi_G \Leftrightarrow \frac{C_I}{4C_G} \mu V \otimes v^* (k^t [D_r(r, w)], -k^t [D_w(r)]) = \sum_\beta \left[\mathcal{F}_G^{\alpha\beta}, G_\beta^a \right] \otimes \partial\xi_\alpha \quad (7.36)$$

$$\forall a, \forall \alpha = 0..3 : \sum_{\beta=0}^3 \partial_\beta \left(\mathcal{F}_G^{a\alpha\beta} \det P' \right) = 0 \quad (7.37)$$

$$\Delta\mathcal{F}_G = -\delta d\mathcal{F}_G \quad (7.38)$$

For the EM field in a continuous model :

$$J_{EM}^* = \delta\mathcal{F}_{EM} \Leftrightarrow \forall a, \alpha : -\frac{C_I}{4C_A} \mu \langle \psi_0, \psi_0 \rangle V^* = \frac{1}{\det P'} \sum_\beta \partial_\beta \left(\mathcal{F}_{EM}^{a\alpha\beta} \det P' \right) \quad (7.39)$$

$$J_{EM}^* = \frac{C_I}{4C_A} \mu \langle \psi, \psi \rangle V^* = -\Delta \dot{A} \quad (7.40)$$

For the other fields :

$$J_A = \phi_A \Leftrightarrow \frac{C_I}{4C_A} \mu \sum_{a=1}^m \left\langle \psi, [\psi] \frac{1}{i} [\theta_a] \right\rangle V \otimes \vec{\theta}_a = \sum_{\beta} \left[\mathcal{F}_A^{\alpha\beta}, \dot{A}_\beta \right] \otimes \partial \xi_\alpha \quad (7.41)$$

$$\delta \mathcal{F}_A = 0 \Leftrightarrow \forall a, \forall \alpha = 0..3 : \sum_{\beta=0}^3 \partial_\beta \left(\mathcal{F}_A^{\alpha\beta} \det P' \right) = 0 \quad (7.42)$$

$$\Delta \mathcal{F}_A = -\delta d \mathcal{F}_A \quad (7.43)$$

In the SR context : $\alpha = 0, 1, 2, 3 : \det P' = -1$,

For the EM field $J_A^* = \frac{C_I}{4C_A} \mu \langle \psi, \psi \rangle V^* = -\Delta \dot{A} = \square \dot{A}$ with the d'alambertian \square and we have the model of propagation by waves (Maths.2503, 33.2.6).

For the gravitational field :

the equation $\sum_{\beta=0}^3 \partial_\beta \left(\mathcal{F}_G^{\alpha\beta} \det P' \right) = 0$ reads :

$$\begin{aligned} & \sum_{\lambda\mu\beta} Y_1^{\alpha\beta\lambda\mu} \partial_\beta \partial_\lambda G_{r\mu}^a + \sum_{\lambda\beta} \sum_{\mu} \left(-Y_{2\mu}^{\alpha\beta\lambda} + Y_{\mu 2}^{\alpha\lambda\beta} - Y_1^{\alpha\beta\lambda\mu} j(G_{r\mu}^a) \right) (\partial_\beta G_{r\lambda}^a) \\ & + \left(\sum_{\mu} Y_1^{\alpha\beta\lambda\mu} j(G_{w\mu}^a) \right) \partial_\beta G_{w\lambda}^a + \sum_{\mu} Y_{2\mu}^{\alpha\beta\lambda} \left(j(G_{r\lambda}^a) G_{r\beta}^a - j(G_{w\lambda}^a) G_{w\beta}^a \right) = \\ 0 & \sum_{\lambda\beta} Y_1^{\alpha\beta\lambda\mu} \partial_\beta \partial_\lambda G_{w\mu}^a + \sum_{\lambda\beta} \sum_{\mu} \left(-Y_{2\mu}^{\alpha\beta\lambda} + Y_{2\mu}^{\alpha\lambda\beta} - Y_1^{\alpha\beta\lambda\mu} j(G_{r\mu}^a) \right) \partial_\beta G_{w\lambda}^a \\ & - \left(\sum_{\mu} Y_1^{\alpha\beta\lambda\mu} j(G_{w\mu}^a) \right) \partial_\beta G_{r\lambda}^a + \sum_{\mu} Y_{2\mu}^{\alpha\beta\lambda} \left(j(G_{w\lambda}^a) G_{r\beta}^a + j(G_{r\lambda}^a) G_{w\beta}^a \right) = 0 \end{aligned}$$

with

$$Y_1^{\alpha\beta\lambda\mu} = g^{\alpha\lambda} g^{\beta\mu} - g^{\alpha\mu} g^{\beta\lambda}$$

$$Y_{2\mu}^{\alpha\beta\lambda} = \partial_\mu (g^{\alpha\lambda} g^{\beta\mu}) - g^{\alpha\lambda} g^{\beta\mu} \sum_{i\gamma=1}^3 P_i^\gamma \partial_\mu P_\gamma^i$$

the equation : $\sum_{\beta} \left[\mathcal{F}_G^{\alpha\beta}, G_\beta \right] \otimes \partial \xi_\alpha = J_G$ reads in the vacuum :

$\Rightarrow \forall \alpha$

$$\sum_{\beta} j(G_{r\beta}) \mathcal{F}_r^{\alpha\beta} = \sum_{\beta} j(G_{w\beta}) \mathcal{F}_w^{\alpha\beta}$$

$$\sum_{\beta} j(G_{r\beta}) \mathcal{F}_w^{\alpha\beta} = - \sum_{\beta} j(G_{w\beta}) \mathcal{F}_r^{\alpha\beta}$$

An observable for the field is a projection on a finite dimensional vector space. These vector spaces must meet the conditions above. The natural choice is then polynomials, for instance for the EM field the potential is estimated by harmonic polynomials.

7.4.3 Energy of the system

The lagrangian represents the energy of the system. In the first model, the energy density is :

$$\mathcal{E} = \sum_{\alpha\beta} C_G \left\langle \mathcal{F}_G^{\alpha\beta}, \mathcal{F}_{G\alpha\beta} \right\rangle + C_A \left\langle \mathcal{F}_A^{\alpha\beta}, \mathcal{F}_{A\alpha\beta} \right\rangle + C_I \mu \frac{1}{i} \langle \psi, \nabla_V \psi \rangle$$

and on shell the last term is null : the energy is stored in the fields.

We have for the fields :

$$\begin{aligned}
& \sum_{\alpha\beta} \langle \mathcal{F}_A^{\alpha\beta}, \mathcal{F}_{A\alpha\beta} \rangle = 2 \sum_{\{\alpha\beta\}} \langle \mathcal{F}_A^{\alpha\beta}, \mathcal{F}_{A\alpha\beta} \rangle \\
& = 2 \sum_{\{\alpha\beta\}} \langle \mathcal{F}_A^{\alpha\beta}, \partial_\alpha \dot{A}_\beta - \partial_\beta \dot{A}_\alpha + [\dot{A}_\alpha, \dot{A}_\beta] \rangle \\
& = 2 \sum_{\{\alpha\beta\}} \langle \mathcal{F}_A^{\alpha\beta}, \partial_\alpha \dot{A}_\beta - \partial_\beta \dot{A}_\alpha \rangle + 2 \sum_{\{\alpha\beta\}} \langle \mathcal{F}_A^{\alpha\beta}, [\dot{A}_\alpha, \dot{A}_\beta] \rangle \\
& 2 \sum_{\{\alpha\beta\}} \langle \mathcal{F}_A^{\alpha\beta}, [\dot{A}_\alpha, \dot{A}_\beta] \rangle = 2 \sum_{\alpha>\beta} \langle \mathcal{F}_A^{\alpha\beta}, [\dot{A}_\alpha, \dot{A}_\beta] \rangle \\
& = \sum_{\alpha>\beta} \langle \mathcal{F}_A^{\alpha\beta}, [\dot{A}_\alpha, \dot{A}_\beta] \rangle + \sum_{\beta>\alpha} \langle \mathcal{F}_A^{\beta\alpha}, [\dot{A}_\beta, \dot{A}_\alpha] \rangle \\
& = - \sum_{\alpha>\beta} \langle \mathcal{F}_A^{\alpha\beta}, [\dot{A}_\beta, \dot{A}_\alpha] \rangle + \sum_{\beta>\alpha} \langle \mathcal{F}_A^{\alpha\beta}, [\dot{A}_\beta, \dot{A}_\alpha] \rangle \\
& = - \sum_{\alpha>\beta} \langle [\mathcal{F}_A^{\alpha\beta}, \dot{A}_\beta], \dot{A}_\alpha \rangle - \sum_{\beta>\alpha} \langle [\mathcal{F}_A^{\alpha\beta}, \dot{A}_\beta], \dot{A}_\alpha \rangle \\
& = - \sum_{\alpha\beta} \langle [\mathcal{F}_A^{\alpha\beta}, \dot{A}_\beta], \dot{A}_\alpha \rangle \\
& = - \sum_\alpha \langle \phi_A^\alpha, \dot{A}_\alpha \rangle \\
& 2 \sum_{\{\alpha\beta\}} \langle \mathcal{F}_A^{\alpha\beta}, \partial_\alpha \dot{A}_\beta - \partial_\beta \dot{A}_\alpha \rangle = 2 \sum_{\alpha>\beta} \langle \mathcal{F}_A^{\alpha\beta}, \partial_\alpha \dot{A}_\beta - \partial_\beta \dot{A}_\alpha \rangle \\
& = \sum_{\alpha>\beta} \langle \mathcal{F}_A^{\alpha\beta}, \partial_\alpha \dot{A}_\beta \rangle - \sum_{\beta>\alpha} \langle \mathcal{F}_A^{\beta\alpha}, \partial_\alpha \dot{A}_\beta \rangle \\
& = 2 \sum_{\alpha\beta} \langle \mathcal{F}_A^{\alpha\beta}, \partial_\alpha \dot{A}_\beta \rangle
\end{aligned}$$

Thus the energy density of the fields can be written :

$$\mathcal{E}_{Fields} = - \sum_\alpha C_A \langle \phi_A^\alpha, \dot{A}_\alpha \rangle + C_G \langle \phi_G^\alpha, G_\alpha \rangle + 2 \sum_{\alpha\beta} C_G \langle \mathcal{F}_G^{\alpha\beta}, \partial_\alpha G_\beta \rangle + C_A \langle \mathcal{F}_A^{\alpha\beta}, \partial_\alpha \dot{A}_\beta \rangle$$

For the particles :

$$\begin{aligned}
& C_I \mu \frac{1}{i} V^\beta \langle \psi, \nabla_\alpha \psi \rangle \\
& = C_I \mu \frac{1}{i} V^\beta \left(\langle \psi, \partial_\alpha \psi \rangle + \sum_{a=1}^m \dot{A}_\alpha^a \langle \psi, [\psi] [\theta_a] \rangle + \sum_{a=1}^6 G_\alpha^a \langle \psi, \gamma C(\vec{\kappa}_\alpha) \psi \rangle \right) \\
& = C_I \mu \frac{1}{i} V^\beta \langle \psi, \partial_\alpha \psi \rangle + \sum_{a=1}^m \dot{A}_\alpha^a C_I \mu V^\beta \frac{1}{i} \langle \psi, [\psi] [\theta_a] \rangle + \sum_{a=1}^6 G_\alpha^a C_I \mu V^\beta \frac{1}{i} \langle \psi, \gamma C(\vec{\kappa}_\alpha) \psi \rangle \\
& = C_I \mu \frac{1}{i} V^\beta \langle \psi, \partial_\alpha \psi \rangle + 4 C_A \langle \dot{A}_\alpha, J_A^\beta \rangle + 4 C_G \langle G_\alpha, J_G^\beta \rangle
\end{aligned}$$

$$\begin{aligned}
\mathcal{E}_{Particles} & = C_I \mu \frac{1}{i} \langle \psi, \nabla_V \psi \rangle = C_I \mu \frac{1}{i} \left\langle \psi, \frac{d\psi}{dt} \right\rangle + 4 \sum_{\alpha=0}^3 \left(C_A \langle \dot{A}_\alpha, J_A^\alpha \rangle + 4 C_G \langle G_\alpha, J_G^\alpha \rangle \right) \\
& = C_I \mu \left(k^t X_\alpha + \frac{1}{i} \langle \psi, [\psi] [\dot{A}_\alpha] \rangle \right)
\end{aligned}$$

The energy density of the system reads :

$$\begin{aligned}
\mathcal{E} & = 2 \sum_{\alpha\beta} C_G \langle \mathcal{F}_G^{\alpha\beta}, \partial_\alpha G_\beta \rangle + C_A \langle \mathcal{F}_A^{\alpha\beta}, \partial_\alpha \dot{A}_\beta \rangle + C_I \mu \left(k^t X_\alpha + \frac{1}{i} \langle \psi, [\psi] [\dot{A}_\alpha] \rangle \right) - \\
& \sum_\alpha C_A \langle \phi_A^\alpha, \dot{A}_\alpha \rangle + C_G \langle \phi_G^\alpha, G_\alpha \rangle
\end{aligned}$$

On shell :

$$\begin{aligned}
& C_I \mu \frac{1}{i} \langle \psi, \nabla_V \psi \rangle = 0 \\
& = C_I \mu \left(k^t X_\alpha + \frac{1}{i} \langle \psi, [\psi] [\dot{A}_\alpha] \rangle \right) \\
& = C_I \mu \frac{1}{i} \left\langle \psi, \frac{d\psi}{dt} \right\rangle + 4 \sum_{\alpha=0}^3 \left(C_A \langle \dot{A}_\alpha, J_A^\alpha \rangle + 4 C_G \langle G_\alpha, J_G^\alpha \rangle \right)
\end{aligned}$$

The kinetic energy of the particles is the opposite of the energy stored in the currents :

$$\begin{aligned}
C_I \mu \frac{1}{i} \left\langle \psi, \frac{d\psi}{dt} \right\rangle &= -4 \sum_{\alpha=0}^3 \left(C_A \left\langle \dot{A}_\alpha, J_A^\alpha \right\rangle + 4C_G \left\langle G_\alpha, J_G^\alpha \right\rangle \right) \\
\text{and as : } J_A &= \phi_A, J_G = \phi_G \\
\sum_{\alpha=0}^3 \left(C_A \left\langle \dot{A}_\alpha, J_A^\alpha \right\rangle + C_G \left\langle G_\alpha, J_G^\alpha \right\rangle \right) \\
&= \sum_{\alpha} C_A \left\langle \phi_A^\alpha, \dot{A}_\alpha \right\rangle + C_G \left\langle \phi_G^\alpha, G_\alpha \right\rangle \\
&= -\frac{1}{4} C_I \mu \frac{1}{i} \left\langle \psi, \frac{d\psi}{dt} \right\rangle \\
&= -\frac{1}{4} C_I \mu k^t \left([A_r] \frac{dr}{dt} + [A_w] \frac{dw}{dt} \right)
\end{aligned}$$

So that, on shell, the energy density of the system reads :

$$\mathcal{E}_{shell} = 2 \sum_{\alpha\beta=0}^3 C_A \left\langle \mathcal{F}_A^{\alpha\beta}, \partial_\alpha \dot{A}_\beta \right\rangle + C_G \left\langle \mathcal{F}_G^{\alpha\beta}, \partial_\alpha G_\beta \right\rangle + \frac{1}{4} C_I \mu k^t \left([A_r] \frac{dr}{dt} + [A_w] \frac{dw}{dt} \right) \quad (7.44)$$

Notice that the rotational energy is part of the energy of the system.

7.4.4 Tetrad equation

The tetrad equation reads :

$$\begin{aligned}
\forall \alpha, \beta = 0 \dots 3 : & 4 \sum_{\gamma=0}^3 \left(C_G \left\langle \mathcal{F}_{G\alpha\gamma}, \mathcal{F}_G^{\beta\gamma} \right\rangle + C_A \left\langle \mathcal{F}_{A\alpha\gamma}, \mathcal{F}_A^{\beta\gamma} \right\rangle \right) + C_I \mu \frac{1}{i} V^\beta \left\langle \psi, \nabla_\alpha \psi \right\rangle \\
&= \delta_\beta^\alpha \sum_{\lambda\mu} \left(C_G \left\langle \mathcal{F}_{G\lambda\mu}, \mathcal{F}_G^{\lambda\mu} \right\rangle + C_A \left\langle \mathcal{F}_{A\lambda\mu}, \mathcal{F}_A^{\lambda\mu} \right\rangle \right) = \delta_\beta^\alpha \mathcal{E}_{shell} \\
\sum_{\gamma=0}^3 \left\langle \mathcal{F}_A^{\beta\gamma}, \mathcal{F}_{A\alpha\gamma} \right\rangle &= \sum_{\gamma=0}^3 \left\langle \mathcal{F}_A^{\beta\gamma}, \partial_\alpha \dot{A}_\gamma - \partial_\gamma \dot{A}_\alpha + [\dot{A}_\alpha, \dot{A}_\gamma] \right\rangle \\
&= \sum_{\gamma=0}^3 \left\langle \mathcal{F}_A^{\beta\gamma}, \partial_\alpha \dot{A}_\gamma - \partial_\gamma \dot{A}_\alpha \right\rangle + \sum_{\gamma=0}^3 \left\langle \mathcal{F}_A^{\beta\gamma}, [\dot{A}_\alpha, \dot{A}_\gamma] \right\rangle \\
&= \sum_{\gamma=0}^3 \left\langle \mathcal{F}_A^{\beta\gamma}, \partial_\alpha \dot{A}_\gamma - \partial_\gamma \dot{A}_\alpha \right\rangle - \sum_{\gamma=0}^3 \left\langle \mathcal{F}_A^{\beta\gamma}, [\dot{A}_\gamma, \dot{A}_\alpha] \right\rangle \\
&= \sum_{\gamma=0}^3 \left\langle \mathcal{F}_A^{\beta\gamma}, \partial_\alpha \dot{A}_\gamma - \partial_\gamma \dot{A}_\alpha \right\rangle - \sum_{\gamma=0}^3 \left\langle [\mathcal{F}_A^{\beta\gamma}, \dot{A}_\gamma], \dot{A}_\alpha \right\rangle \\
&= \sum_{\gamma=0}^3 \left\langle \mathcal{F}_A^{\beta\gamma}, \partial_\alpha \dot{A}_\gamma - \partial_\gamma \dot{A}_\alpha \right\rangle - \left\langle \phi_A^\beta, \dot{A}_\alpha \right\rangle \\
4 \sum_{\gamma=0}^3 \left(C_G \left\langle \mathcal{F}_{G\alpha\gamma}, \mathcal{F}_G^{\beta\gamma} \right\rangle + C_A \left\langle \mathcal{F}_{A\alpha\gamma}, \mathcal{F}_A^{\beta\gamma} \right\rangle \right) &+ C_I \mu \frac{1}{i} V^\beta \left\langle \psi, \nabla_\alpha \psi \right\rangle \\
&= 4 \sum_{\gamma=0}^3 C_A \left\langle \mathcal{F}_A^{\beta\gamma}, \partial_\alpha \dot{A}_\gamma - \partial_\gamma \dot{A}_\alpha \right\rangle + C_G \left\langle \mathcal{F}_G^{\beta\gamma}, \partial_\alpha G_\gamma - \partial_\gamma G_\alpha \right\rangle \\
&- 4C_A \left\langle \phi_A^\beta, \dot{A}_\alpha \right\rangle - 4C_A \left\langle \phi_G^\beta, G_\alpha \right\rangle \\
&+ C_I \mu \frac{1}{i} V^\beta \left\langle \psi, \partial_\alpha \psi \right\rangle + 4C_A \left\langle \dot{A}_\alpha, J_A^\beta \right\rangle + 4C_G \left\langle G_\alpha, J_G^\beta \right\rangle \\
&= C_I \mu \frac{1}{i} V^\beta \left\langle \psi, \partial_\alpha \psi \right\rangle + 4 \sum_{\gamma=0}^3 C_A \left\langle \mathcal{F}_A^{\beta\gamma}, \partial_\alpha \dot{A}_\gamma - \partial_\gamma \dot{A}_\alpha \right\rangle + C_G \left\langle \mathcal{F}_G^{\beta\gamma}, \partial_\alpha G_\gamma - \partial_\gamma G_\alpha \right\rangle \\
&= \delta_\beta^\alpha \left(2 \sum_{\alpha\beta=0}^3 C_A \left\langle \mathcal{F}_A^{\alpha\beta}, \partial_\alpha \dot{A}_\beta \right\rangle + C_G \left\langle \mathcal{F}_G^{\alpha\beta}, \partial_\alpha G_\beta \right\rangle + \frac{1}{4} C_I \mu \frac{1}{i} \left\langle \psi, \frac{d\psi}{dt} \right\rangle \right)
\end{aligned}$$

$$C_{I\mu} \frac{1}{i} V^\beta \langle \psi, \partial_\alpha \psi \rangle + 4 \sum_{\gamma=0}^3 C_A \langle \mathcal{F}_A^{\beta\gamma}, \partial_\alpha \dot{A}_\gamma - \partial_\gamma \dot{A}_\alpha \rangle + C_G \langle \mathcal{F}_G^{\beta\gamma}, \partial_\alpha G_\gamma - \partial_\gamma G_\alpha \rangle \quad (7.45)$$

$$= \delta_\beta^\alpha \left(\frac{1}{4} C_{I\mu} \frac{1}{i} \left\langle \psi, \frac{d\psi}{dt} \right\rangle + 2 \sum_{\lambda\mu=0}^3 C_A \langle \mathcal{F}_A^{\lambda\mu}, \partial_\lambda \dot{A}_\mu \rangle + C_G \langle \mathcal{F}_G^{\lambda\mu}, \partial_\lambda G_\mu \rangle \right) \quad (7.46)$$

It is the only equation which shows a relation between the gravitational field and the other fields. In the equations of the models, if exterior fields are present they must be added in the equations above. The gravitational field represents the action of the inertial forces, and if an exterior field \dot{A}_{ext} is added, then there must be a compensation of the gravitation field : the trajectories change and then inertial forces appear.

But this equation has another meaning.

7.4.5 Conservation of energy

We have seen (see Lagrangian) that the conservation of the energy of the system requires : $div(\varepsilon_0 L) = 0$ that is :

$$\begin{aligned} \sum_{\alpha=0}^3 \frac{d}{d\xi^\alpha} (\varepsilon_0^\alpha L \det P') &= \frac{d}{d\xi^0} (L \det P') \\ &= \frac{d}{dt} (L \det P') = \det P' \frac{dL}{dt} + L \frac{d}{dt} (\det P') \\ &= \det P' \frac{dL}{dt} - L \sum_{i\alpha} P_\alpha^i \frac{dP_\alpha^i}{dt} (\det P') \end{aligned}$$

that is :

$$\frac{dL}{dt} = L \sum_{i\alpha} P_\alpha^i \frac{dP_\alpha^i}{dt}$$

But the tetrad equation reads

$$\frac{dL}{dP_\alpha^i} - L P_\alpha^i = 0$$

and on shell the condition sums up to the identity :

$$\frac{dL}{dt} = \sum_{i\alpha} L P_\alpha^i \frac{dP_\alpha^i}{dt} = \sum_{i\alpha} \frac{dL}{dP_\alpha^i} \frac{dP_\alpha^i}{dt}$$

So the tetrad equation implies the conservation of energy. The result holds for any lagrangian.

7.5 ENERGY-MOMENTUM TENSOR

7.5.1 Definition

The concept of equilibrium is at the core of the Principle of Least Action. So, for any tentative change of the values of the variables, beyond the point of equilibrium, the system reacts by showing resistance against the change : this is the inertia of the system. It is better understood with the functional derivatives.

The idea applies foremost to models of the first kind, with a continuous distribution of particles and fields. Let us consider an action with integrals of the kind :

$$\begin{aligned} \ell(z^i, z_\alpha^i) &= \int_{\Omega} \mathcal{L}(z^i, z_\alpha^i, P_i^\alpha) \varpi_4 \\ &= \int L(z^i, z_\alpha^i, P_i^\alpha) (\det P') d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3 \end{aligned}$$

with the variables $Z = (z^i, z_\alpha^i, P_i^\alpha)$ sections of a 1st jet bundle $J^1 E$.

For a change $\delta Z = (\delta z^i, \delta z_\alpha^i, \delta P_i^\alpha)$ the change of the action is, at the second order :

$$\begin{aligned} \delta \ell &= \int \sum_{i\alpha} \left(\frac{\partial L}{\partial z^i} \delta z^i + \frac{\partial L}{\partial z_\alpha^i} \delta z_\alpha^i + \frac{\partial L}{\partial P_i^\alpha} \delta P_i^\alpha - L P_i^\alpha \delta P_i^\alpha \right) \det P' d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3 \\ &\text{with } \frac{\partial \det P'}{\partial P_i^\alpha} = -P_i^\alpha \det P' \\ \delta \ell &= \int \sum_{i\alpha} \left(\frac{\partial L}{\partial z^i} \delta z^i + \frac{\partial L}{\partial z_\alpha^i} \delta z_\alpha^i + \left(\frac{\partial L}{\partial P_i^\alpha} - L P_i^\alpha \right) \delta P_i^\alpha \right) \varpi_4 \end{aligned}$$

The variational derivative is, when δZ comes from a section (see the proof of Maths.2601), that is when : $\delta z_\alpha^i = \partial_\alpha \delta z^i$:

$$\begin{aligned} \frac{\delta \ell}{\delta z^i} (\delta z^i) &= \int \left(\frac{\partial \mathcal{L}}{\partial z^i} - \sum_\alpha \frac{d}{d\xi^\alpha} \left(\frac{\partial \mathcal{L}}{\partial z_\alpha^i} \right) \right) (\delta z^i) d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3 \\ &= \int \left(\frac{\partial L}{\partial z^i} \det P' - \sum_\alpha \frac{d}{d\xi^\alpha} \left(\frac{\partial L}{\partial z_\alpha^i} \det P' \right) \right) (\delta z^i) d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3 \\ &\Leftrightarrow \int \frac{\partial L}{\partial z^i} \delta z^i \varpi_4 = \frac{\delta \ell}{\delta z^i} (\delta z^i) + \int \left(\sum_\alpha \frac{d}{d\xi^\alpha} \left(\frac{\partial L}{\partial z_\alpha^i} \det P' \right) \right) \delta z^i d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3 \\ \frac{\delta \ell}{\delta P_i^\alpha} (\delta P_i^\alpha) &= \int \left(\frac{\partial L}{\partial P_i^\alpha} \det P' - P_i^\alpha \det P' \right) (\delta P_i^\alpha) d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3 \\ &= \int \left(\frac{\partial L}{\partial P_i^\alpha} - P_i^\alpha \right) (\delta P_i^\alpha) \varpi_4 \end{aligned}$$

so we can rewrite :

$$\delta \ell = \sum_i \frac{\delta \ell}{\delta z^i} \delta z^i + \sum_{i\alpha} \frac{\delta \ell}{\delta P_i^\alpha} \delta P_i^\alpha + \int \left(\sum_{i\alpha} \frac{d}{d\xi^\alpha} \left(\frac{\partial L}{\partial z_\alpha^i} \det P' \right) \right) \delta z^i d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3$$

By definition, on shell (at equilibrium) the first part is null, thus :

$$\begin{aligned} \delta \ell &= \int \left(\sum_\alpha \frac{d}{d\xi^\alpha} \left(\frac{\partial L}{\partial z_\alpha^i} \det P' \right) \right) \delta z^i d\xi^0 \wedge d\xi^1 \wedge d\xi^2 \wedge d\xi^3 \\ \delta Y &= \sum_i \delta z^i \sum_\alpha \frac{\partial L}{\partial z_\alpha^i} \partial \xi_\alpha \text{ is a vector field (see Lagrangian) and} \\ \sum_\alpha \frac{d}{d\xi^\alpha} \left(\sum_i \frac{\partial L}{\partial z_\alpha^i} \delta z^i (\det P') \right) &= (\det P') \operatorname{div} (\delta Y) \end{aligned}$$

$$\delta \ell = \int_{\Omega} \operatorname{div} (\delta Y) \varpi_4 \quad (7.47)$$

The variables z^i are sections of a fiber bundle. If we consider a translation : $\delta v = \sum_{\beta=0}^3 v^\alpha \delta \xi_\beta$:

$$\delta z^i = \sum_\beta \frac{\partial z^i}{\partial \xi^\beta} v^\beta \delta \xi_\beta = \sum_\beta v^\beta z_i^\beta \delta \xi_\beta$$

and :

$$\delta Y = \sum_i \left(\sum_\beta v^\beta z_i^\beta \delta \xi_\beta \right) \sum_\alpha \frac{\partial L}{\partial z_\alpha^i} \partial \xi_\alpha = \sum_{i\alpha\beta} \frac{\partial L}{\partial z_\alpha^i} z_i^\beta v^\beta \delta \xi_\beta$$

The quantity :

$$T = \sum_{i\alpha\beta} \frac{\partial L}{\partial z_\alpha^i} z_i^\beta \partial \xi_\alpha \otimes d\xi^\beta \quad (7.48)$$

is a tensor, and $\delta Y = T(\delta v)$

$$\delta \ell = \int_\Omega \text{div}(T(\delta v)) \varpi_4 \quad (7.49)$$

T is called the **energy-momentum tensor**. Its name and definition, closely related to the lagrangian, come from fluid mechanics. The lagrangian has the meaning of a density of energy for the whole system, $\text{div}(T(\delta v))$ has the meaning of a variation of this energy due to the action of 4 dimensional forces T acting along the direction δv . T can be interpreted as the resistance of the system against a change in the direction $\partial \xi_\alpha$, or equivalently as the inertia of the system, and it is opposite to the 4 dimensional force necessary to change the equilibrium. So it can be assimilated to the relativist momentum of the system.

7.5.2 Expression of the Energy Momentum tensor

With the more general lagrangian $\mathcal{L} = L(\psi, \nabla_\alpha \psi_p, P_i^\alpha, \mathcal{F}_{G\alpha\beta}, \mathcal{F}_{A\alpha\beta}, V^\alpha) \det P'$ the tensor T reads :

$$T = \sum_{\alpha\beta} \left\{ \sum_{ij} \frac{\partial L}{\partial \partial_\alpha \psi^{ij}} \partial_\beta \psi^{ij} + \sum_{a\gamma} \frac{\partial L}{\partial \partial_\alpha A_\gamma^a} \partial_\beta \dot{A}_\gamma^a + \sum_{a\gamma} \frac{\partial L}{\partial \partial_\alpha G_\gamma^a} \partial_\beta G_\gamma^a \right\} \partial \xi_\alpha \otimes d\xi^\beta$$

Notice that P_i^α, V^α do not appear. With :

$$\begin{aligned} \frac{dL}{d\partial \partial_\alpha \psi^{ij}} &= \Pi_{\nabla}^{\alpha}{}_{ij} \\ \frac{dL}{d\partial_\beta G_\alpha^a} &= -2\Pi_{G_a}^{\alpha\beta} \\ \frac{dL}{d\partial_\beta \dot{A}_\alpha^a} &= -2\Pi_{A_a}^{\alpha\beta} \end{aligned}$$

$$T = \sum_{\alpha\beta} \left(\sum_{ij} \Pi_{\nabla}^{\alpha}{}_{ij} \partial_\beta \psi^{ij} - 2 \sum_{a\gamma} \Pi_{A_a}^{\alpha\gamma} \partial_\beta \dot{A}_\gamma^a + \Pi_{G_a}^{\alpha\gamma} \partial_\beta G_\gamma^a \right) \partial \xi_\alpha \otimes d\xi^\beta \quad (7.50)$$

So that the momenta can be derived from the Energy-Momentum tensor, in a way which is usual in fluid mechanics : $\Pi_{\nabla}^{\alpha}{}_{ij} = \frac{\partial T}{\partial \partial_\beta \psi^{ij}}, \dots$

With the perturbative lagrangian in a model of the first type :

$$T_\beta^\alpha = C_I \mu V^\alpha k^t ([A_r] \partial_\beta r + [A_w] \partial_\beta w) + 4 \sum_\gamma C_A \langle \mathcal{F}_A^{\alpha\gamma}, \partial_\beta \dot{A}_\gamma \rangle + C_G \langle \mathcal{F}_G^{\alpha\gamma}, \partial_\beta G_\gamma \rangle \quad (7.51)$$

The trace of the tensor is :

$$Tr(T) = \sum_{\alpha=0}^3 T_{\alpha}^{\alpha} = C_I \mu k^t \left([A_r] \frac{dr}{dt} + [A_w] \frac{dw}{dt} \right) + 4 \sum_{\alpha\gamma} C_A \left\langle \mathcal{F}_A^{\alpha\gamma}, \partial_{\alpha} \dot{A}_{\gamma} \right\rangle + C_G \left\langle \mathcal{F}_G^{\alpha\gamma}, \partial_{\alpha} G_{\gamma} \right\rangle$$

The density of energy of the system is :

$$\mathcal{E} = 2 \sum_{\alpha\beta} C_G \left\langle \mathcal{F}_G^{\alpha\beta}, \partial_{\alpha} G_{\beta} \right\rangle + C_A \left\langle \mathcal{F}_A^{\alpha\beta}, \partial_{\alpha} \dot{A}_{\beta} \right\rangle + C_I \mu \left(k^t X_{\alpha} + \frac{1}{i} \left\langle \psi, [\psi] \left[\dot{A}_{\alpha} \right] \right\rangle \right) - \sum_{\alpha} C_A \left\langle \phi_A^{\alpha}, \dot{A}_{\alpha} \right\rangle + C_G \left\langle \phi_G^{\alpha}, G_{\alpha} \right\rangle$$

Thus :

$$\begin{aligned} & \sum_{\alpha\gamma} C_A \left\langle \mathcal{F}_A^{\alpha\gamma}, \partial_{\alpha} \dot{A}_{\gamma} \right\rangle + C_G \left\langle \mathcal{F}_G^{\alpha\gamma}, \partial_{\alpha} G_{\gamma} \right\rangle \\ &= \frac{1}{4} Tr(T) - \frac{1}{4} C_I \mu k^t \left([A_r] \frac{dr}{dt} + [A_w] \frac{dw}{dt} \right) \\ &= \frac{1}{2} \mathcal{E} - \frac{1}{2} C_I \mu \left(k^t \left([A_r] \frac{dr}{dt} + [A_w] \frac{dw}{dt} \right) + [D_r] \hat{G}_r + [D_w] \hat{G}_w + \frac{1}{i} \left\langle \psi, [\psi] \left[\dot{A}_{\alpha} \right] \right\rangle \right) + \\ & \frac{1}{2} \sum_{\alpha} C_A \left\langle \phi_A^{\alpha}, \dot{A}_{\alpha} \right\rangle + C_G \left\langle \phi_G^{\alpha}, G_{\alpha} \right\rangle \end{aligned}$$

$$\mathcal{E} = \frac{1}{2} Tr(T) + \frac{1}{2} C_I \mu k^t \left([A_r] \frac{dr}{dt} + [A_w] \frac{dw}{dt} \right) + C_I \mu \left(k^t \left([D_r] \hat{G}_r + [D_w] \hat{G}_w \right) + \frac{1}{i} \left\langle \psi, [\psi] \left[\dot{A}_{\alpha} \right] \right\rangle \right) + \sum_{\alpha} C_A \left\langle \phi_A^{\alpha}, \dot{A}_{\alpha} \right\rangle + C_G \left\langle \phi_G^{\alpha}, G_{\alpha} \right\rangle$$

and on shell :

$$\begin{aligned} & \sum_{\alpha\beta=0}^3 C_A \left\langle \mathcal{F}_A^{\alpha\beta}, \partial_{\alpha} \dot{A}_{\beta} \right\rangle + C_G \left\langle \mathcal{F}_G^{\alpha\beta}, \partial_{\alpha} G_{\beta} \right\rangle \\ &= \frac{1}{2} \mathcal{E} - \frac{1}{8} C_I \mu k^t \left([A_r] \frac{dr}{dt} + [A_w] \frac{dw}{dt} \right) = \\ &= \frac{1}{4} Tr(T) - \frac{1}{4} C_I \mu k^t \left([A_r] \frac{dr}{dt} + [A_w] \frac{dw}{dt} \right) \end{aligned}$$

$$\mathcal{E} = -\frac{1}{4} C_I \mu k^t \left([A_r] \frac{dr}{dt} + [A_w] \frac{dw}{dt} \right) + \frac{1}{2} Tr(T) \quad (7.52)$$

In the vacuum : $\mathcal{E} = \frac{1}{2} Tr(T)$.

The energy is stationary on shell, any change to the trajectory of the particles meets a resistance, due to the fields, equal in energy to 2 times the kinetic energy.

For a deformable solid which is not submitted to external fields we have seen that the interactions of the fields induced by the motion of the particles can be seen as the forces resulting from the deformation. So one can state similarly that for any deformation $(\delta r, \delta w)$ the energy-momentum tensor provides, in a continuous deformation at equilibrium, the value of the induced fields, or equivalently the deformation induced by external fields.

7.5.3 Momentum of the system

The Energy-Momentum tensor can be assimilated to the relativist momentum of the system. However, because this is a tensor, defined at each point in the chart of the observer, there is no way to sum it over the whole system (at least in the GR framework). The conservation at each point along the vector ε_0 would require : $\mathcal{L}_{\varepsilon_0} T = 0$.

On shell, $\forall \alpha, \beta$:

$$C_I \mu \frac{1}{i} V^{\alpha} \left\langle \psi, \partial_{\beta} \psi \right\rangle + 4 \sum_{\gamma=0}^3 C_A \left\langle \mathcal{F}_A^{\alpha\gamma}, \partial_{\beta} \dot{A}_{\gamma} - \partial_{\gamma} \dot{A}_{\beta} \right\rangle + C_G \left\langle \mathcal{F}_G^{\alpha\gamma}, \partial_{\beta} G_{\gamma} - \partial_{\gamma} G_{\beta} \right\rangle = \delta_{\beta}^{\alpha} \mathcal{E}_{shell}$$

$$\begin{aligned}
T &= \sum_{\alpha\beta} \left(\delta_{\beta}^{\alpha} \mathcal{E}_{shell} + 4 \sum_{\gamma=0}^3 C_A \langle \mathcal{F}_A^{\alpha\gamma}, \partial_{\gamma} \dot{A}_{\beta} \rangle + C_G \langle \mathcal{F}_G^{\alpha\gamma}, \partial_{\gamma} G_{\beta} \rangle \right) \partial \xi_{\alpha} \otimes d\xi^{\beta} \\
\mathcal{L}_{\varepsilon_0} T &= \mathcal{L}_{\varepsilon_0} \left(\sum_{\alpha\beta} \delta_{\beta}^{\alpha} \mathcal{E}_{shell} \partial \xi_{\alpha} \otimes d\xi^{\beta} \right) \\
&\quad + 4 \mathcal{L}_{\varepsilon_0} \sum_{\alpha\beta\gamma=0}^3 \left(C_A \langle \mathcal{F}_A^{\alpha\gamma}, \partial_{\gamma} \dot{A}_{\beta} \rangle + C_G \langle \mathcal{F}_G^{\alpha\gamma}, \partial_{\gamma} G_{\beta} \rangle \right) \partial \xi_{\alpha} \otimes d\xi^{\beta}
\end{aligned}$$

The first term corresponds to the conservation of energy, thus it is null. But there is no obvious reason for the second term to be null.

Chapter 8

DISCONTINUOUS MODELS

Most physical phenomena involve, at some step, processes which can be seen as discontinuous :

- collision of molecules or particles, elastic (without loss of energy) or not
- the photo-electric effect : interaction of an electromagnetic field with electrons of the atoms of a solid material
- disintegration of a nucleus, spontaneous or following collisions
- black body : interaction of the electromagnetic field with the electrons of a body, under thermal agitation

We have seen that, in a continuous model, the spatial spin does not change. So the change of its orientation (up or down) requires a discontinuous process.

The models used to represent these phenomena feature an explicit discontinuous operation. There are many variants.

The oldest are the kinetic models. They usually derive from a hydrodynamic model (similar to the continuous models), but add a “collision operator” to represent elastic collisions between particles : because there is no loss of energy, the pure spatial momenta are conserved, which brings a fundamental relation between the variables before and after a collision. The models are based upon a distribution function $f(m, p)$ of particles of linear momentum p which shall follow a conservation law, using the collision operator. So the distribution of charges is itself given by a specific equation. Then the 4 dimensional action, with a lagrangian adapted to the fields considered, gives an equation relating the field and the distribution of charges. Usually the particles are assumed to have the same physical characteristics (mass and charge), which imposes an additional condition on the linear momentum : $\langle p, p \rangle = mc^2$. The frequency of collisions is related to a thermodynamic variable similar to temperature. Such models have been extensively studied with gravitational fields only (Boltzman systems), notably in Astrophysics, and the electromagnetic field for plasmas (Vlasov-Maxwell systems).

In QTF these discontinuous processes are represented by Feynman diagrams, which describe all the possible interactions between fermions and bosons, and the probability of occurrence of a given interaction is computed by path-integrals.

An equilibrium is not necessarily the result of a continuous process but, in the physical world, no process is totally discontinuous : the discontinuity appears as a singular event, between periods of equilibrium. So the Principle of Least Action and the kind of models that we have seen are not useless, and we need to find a way to introduce discontinuities with respect to smooth variables. At an elementary level discontinuities can happen in collisions of particles, but also with fields in the vacuum. Particles and fields adjust to each other by their interactions, but these adjustments are not instantaneous, and, as the fields propagate, there are points in the vacuum where a smooth adjustment is not possible and a discontinuity appears. The discontinuities in fields are linked with bosons : they are seen as the carriers of force fields, in the meaning that discontinuities occur when they meet each other or particles. This is clear in all the classic experiments involving photons, such as the photo-electric effect. So we start with the discontinuities for fields.

8.1 BOSONS

8.1.1 Discontinuities for the fields

Wave propagations, wave packets, burst of fields, or solitons are continuous processes : they are solutions, sometimes very specific to initial conditions or to the nature of the medium in which the field propagates, of regular differential equations such as $\square A = 0$. Discontinuities are different, and can take at the macroscopic level the form of shock waves, but we need to precise what we mean for the fields.

The force fields have been represented so far by smooth maps over M . A discontinuity in the fields must be understood as :

- i) occurring as the consequence of some discontinuous process, occurring at a location $a \in M$ with the time a_t , which can be outside Ω
- ii) propagating in M , the propagation is a continuous process, which conserves some of the initial characteristics of the discontinuity
- iii) it has a geometric support. Phenomena discontinuous at any point are unphysical. The discontinuity occurs either in isolated points, or at the border $\partial\omega(t)$ of a closed domain $\omega(t) \subset \Omega_3(t)$ for $t > a_t$, subset of a submanifold of $\Omega(t)$ of dimension 1 (a curve), 2 (a surface) or 3 (a volume).

The propagation can then be represented by a differentiable map, starting at $a_t : f(t) : \partial\omega(a_t) \rightarrow \partial\omega(t)$ such that $\omega(t) = \{\varphi_M(t, f(t)(x)), x \in \omega(a_t)\}$ and $\frac{df}{dt} \neq 0$, then the map is bijective and a local diffeomorphism, so the dimension of $\varpi(t)$ must be preserved.

The field is then continuous in the interior of $\omega(t)$ (but with the effects of the discontinuity added) and discontinuous at the boundary $\partial\omega(t)$.

Our work has been so far done at the most elementary level, where it is possible to discard the structural relations which exist around a material point : we consider geometric points. If the discontinuity happens at a point, the conservation of the dimension of $\omega(t)$ makes that the only physically conceivable solution is that it propagates along a path, time like or null, future oriented, and meets each $\Omega(t)$ at a singular point. The other solutions are actually the consequence of discontinuous processes happening simultaneously at points which are spatially close. In models using a density it is clear that one can consider discontinuities represented by an area $\partial\omega(t)$ which is not limited to a point, but they involve other processes and require assumptions about the structure of the body or the fluid and their thermodynamics, that are not our topic. So we will focus on the case of a singular discontinuity occurring at some point and propagating along a curve. This implies that the discrepancy does not spread around its trajectory, moves as the field propagates, and keeps some of its initial characteristics along its trajectory.

A discontinuity of the field can appear either as a discontinuity of the potential or as a discontinuity of the strength of the field. The latter, at least for the EM field, has a physical meaning, for which one can easily conceive discontinuities. However its general definition, with the brackets of potentials, makes difficult a study of discontinuities. On the other hand the potential is a more

abstract quantity, and a discontinuous potential would preclude the existence of the strength of the field. So what we will study is a *discontinuity in the derivative of the potential*.

The measure of a discontinuity requires to compare the situations before and after the discontinuity. To do this we need other mathematical tools, they are similar to those used in the Lie derivative that we have met several times.

8.1.2 Lie derivative

Definitions

Let $f \in C_1(M; M)$ be a differentiable map on the manifold M . Its derivative $f'(m) : T_m M \rightarrow T_{f(m)} M$ is a continuous linear map between the vector spaces tangent to M at the points m and $f(m)$, bijective if f is a diffeomorphism. It enables to compare tensors in the two vector spaces.

The push-forward of a vector field is the operator $f_* : \mathfrak{X}(TM) \rightarrow \mathfrak{X}(TM) :: (f_* V)(f(m)) = f'(m) V(m)$

The pull-back of a one form field is the operator $f^* : \mathfrak{X}(TM^*) \rightarrow \mathfrak{X}(TM^*) :: (f^* \lambda)(m)(u_m) = \lambda(f(m))(f'(m) u_m)$

One can define similarly the push-forward of a form, and the pull-back of a vector if f is invertible, and extend the operations to tensors (Maths.16.1.5). f_*, f^* are linear operators, which are inverse of each other : $(f_*)^{-1} = (f^{-1})^*$, preserve the exterior product of forms and commute with the exterior differential.

Using these operators one can define a derivative of any tensor field at one point m . The idea is to use the flow $\Phi_V(\theta, m)$ of a given vector field V , which is itself a diffeomorphism, to take a tensor $T(\Phi_V(-\theta, m))$ at some point $\Phi_V(-\theta, m)$, push-forward this tensor at

$m = \Phi_V(0, m) = \Phi_V(\theta, \Phi_V(-\theta, m))$ using $\Phi_V(\theta, m)_*$. The tensor which is so transported belong to the same vector space as $T(\Phi_V(\theta, m))$ and we can compute their difference :

$$\Delta_L(\theta) = T(m) - \Phi_V(\theta, \Phi_V(-\theta, m))_* T(\Phi_V(-\theta, m))$$

Equivalently we can transport $T(\Phi_V(\theta, m))$ in m by $\Phi_V(-\theta, m)$ by pull back and compute :

$$\Delta_R(\theta) = \Phi_V(-\theta, \Phi_V(\theta, m))^* T(\Phi_V(\theta, m)) - T(m)$$

Because $(\Phi_V(\theta, m)_*)^{-1} = (\Phi_V(-\theta, m))^*$ we can use both operations in each case.

We have the following diagram with $m_L = \Phi_V(-\theta, m), m_R = \Phi_V(\theta, m)$

$$\begin{array}{ccccc}
 & & & \text{Left} & \\
 & & & \Phi_V(\theta, m_L)_* T(m_L) & \\
 TM & T(m_L) & \longrightarrow & \longrightarrow & T(m) \\
 & \uparrow & & & \uparrow \\
 & T & & & T \\
 & \uparrow & & \Phi_V(\theta, m_L) & \uparrow \\
 M : & m_L & \longrightarrow & \longrightarrow & m
 \end{array}$$

$$\begin{array}{ccccc}
 & & \text{Right} & & \\
 & & \Phi_V(-\theta, m_R)^* T(m_R) & & \\
 T(m) & \longleftarrow & & \longleftarrow & T(m_R) \\
 \uparrow & & & & \uparrow \\
 T & & & & T \\
 \uparrow & & \Phi_V(-\theta, m_R) & & \uparrow \\
 m & \longleftarrow & & \longleftarrow & m_R
 \end{array}$$

The Lie derivative of T at m is defined as :
 $\mathcal{L}_V T = \lim_{\theta \rightarrow 0^+} \frac{1}{\theta} \Delta_L(\theta) = \lim_{\theta \rightarrow 0^+} \frac{1}{\theta} \Delta_R(\theta) = \frac{\partial}{\partial \theta} \Phi_V(-\theta, m_R)^* T(m_R) |_{\theta=0}$
 $= -\frac{\partial}{\partial \theta} \Phi_V(\theta, m_L)_* T(m_L) |_{\theta=0}$

The Lie derivative is linear with respect to T and V, it preserves the type of tensors, antisymmetric tensors go to antisymmetric tensors, and commutes with the exterior differential (Maths.16.2.2).

Measure of discontinuities on the tangent bundle

We will use the Lie derivative to represent a discontinuity. Let us assume that there is a discontinuity of the derivative of a tensor T happening at some point m along a curve represented by the flow of a vector field V. If T has a derivative at the left and a derivative at the right, along the curve, then :

$$\begin{aligned}
 \mathcal{L}_V^L T &= \lim_{\theta \rightarrow 0^+} \frac{1}{\theta} \Delta_L(\theta) = \lim_{\theta \rightarrow 0^+} \frac{1}{\theta} (T(m) - \Phi_V(\theta, m_L)^* T(m_R)) \\
 \mathcal{L}_V^R T &= \lim_{\theta \rightarrow 0^+} \frac{1}{\theta} \Delta_R(\theta) = \lim_{\theta \rightarrow 0^+} \frac{1}{\theta} (\Phi_V(-\theta, m_R)^* T(m_R) - T(m))
 \end{aligned}$$

and we can define the discontinuity in the derivative by :

$$\begin{aligned}
 \Delta T(m) &= \lim_{\theta \rightarrow 0^+} \frac{1}{\theta} (\Delta_R(\theta) - \Delta_L(\theta)) \\
 &= \lim_{\theta \rightarrow 0^+} \frac{1}{\theta} (\Phi_V(-\theta, m_R)^* T(m_R) - \Phi_V(\theta, m_L)_* T(m_L))
 \end{aligned}$$

We have assumed that discontinuities happen along the curve at singular points. T is continuous in m, is left and right differentiable, but the derivatives are different. Because the discontinuity propagates toward the future its effects will be seen on the left.

$\Delta T(m)$ is a tensor of the same kind as T, and its definition is fully covariant : it does not involve a basis. The propagation of the discontinuity can be seen as the discrepancy between what would be the Lie derivative of a fully differentiable tensor $T(\Phi_V(\theta, m))$, and its actual value :

$$\mathcal{L}_V T + \Delta T(m) = \frac{\partial}{\partial \theta} \Phi_V(-\theta, m_R)^* T(m_R) |_{\theta=0} + \Delta T(m)$$

The propagation occurs along V, which is an essential part of the discontinuity. V must be a time like or null vector, future oriented : $\langle \varepsilon_0(m), V(m) \rangle < 0$ for any observer. The parameter θ is uniquely defined by V and the condition $\Phi_V(0, m) = m$. But because $\Phi_V(\theta + h, m) = \Phi_V(h, \Phi_V(\theta, m))$ and that push-forwards can be composed, we can take any origin for θ if we are consistent with the definition of $\Delta T(m)$.

Because all the formalism that we have developed so far is based upon continuous representations, it is then easy to treat discontinuities as the continuous propagation of discontinuities, *which comes in addition to the continuous variables*. But these discontinuities have a support : if T is defined everywhere, the discontinuity appears only along the curve. So we have a situation similar to

particles. *Discontinuities of continuous fields are represented as particles.* We have already met this kind of situation : the currents ϕ and J have the same support (the curve along which the particle moves) and for individual particles the current ϕ follows the same trajectory as the particle. So discontinuities can be seen as currents, similar to ϕ , without associated particle.

However in the mechanism above the tensor is transported unchanged : $\Phi_V(\theta, m)_*$ is a linear operator and $\Delta T(\Phi_V(\theta, m)) = \Phi_V(\theta, m)_* \Delta T(m)$ is just a copy of $\Delta T(m)$. This is the difference between the Lie derivative and the covariant derivative, which allows for changes along the transport. This is where the concept of gauge comes handy. As well as a particle keeps its intrinsic state along its world line, the discontinuity will not change in its own gauge, but it can change in the gauge of an observer. But of course it will depend on the vector bundle on which the tensor is defined, and this will bring good news.

Discontinuities on fiber bundles

We see how it works for $P_U(M, U, \pi_U)$, that is on the manifold P_U , and its tangent space TP_U , where live the quantities to be derived. A vector of $T_p P_U$ at $p = \varphi_U(m, g)$ is defined by a couple $(v_m, v_g) \in T_m M \times T_g U$.

$$v_p = \sum_{\alpha} v_m^{\alpha} \partial m_{\alpha} + \sum_a v_g^a \partial g_a \in T_p P_U$$

The vertical bundle VP_U is comprised of vectors $\varphi'_{Ug}(m, g)v_g$. It is isomorphic to the Lie algebra, and it is convenient to use the fundamental vectors : $\zeta(v_1)(p) = \sum_{a=1}^m [L'_g 1(v_1)]^a \partial g_a$ where v_1 is a vector of the Lie algebra.

We will need a vector field W on TP_U , but not any vector field will do, we need a vector field W which is projected as a vector field V on M , called a projectable vector field. It is such that $\pi'_U(p)W(p) = V(\pi_U(p))$, meaning that the components v_m^{α} depend on m only, and not g . It reads :

$$W(p) = \sum_{\alpha=0}^3 V^{\alpha}(m) \partial m_{\alpha} + \zeta(X(p))(p) \text{ where } X(p) \in T_1 U$$

The flow of a projectable vector field is a fibered morphism, in the meaning :

$$\forall p \in P_U : \pi_U(\Phi_W(\theta, p)) = \Phi_V(\theta, \pi(p))$$

and because $\pi'_U(p)W(p) = V(\pi_U(p))$ the integral curves of W on P_U and V on M have the same parameter θ .

A vertical vector field, such that $\pi'_U(p)W(p) = 0$, is projectable.

The Lie derivative of *sections* $\phi \in \mathfrak{X}(P_U)$ on P_U is then defined similarly to the scheme above, P_U replaces TM (Maths.2015). We define the right and left transports on P_U by transporting the point in M with the flow of V , then taking this value in P_U by ϕ , and transporting it by the flow of W , according to the diagram below. Because ϕ is a section its value depends only on m , and because W is projectable one comes back to the same fiber $\pi_U^{-1}(m)$.

$$\text{With } \Phi_V(-\theta, m) = m_L, \Phi_V(\theta, m) = m_R$$

$$\begin{array}{ccccc}
 & & \text{Left} & & \\
 & & \Phi_W(\phi(m_L), \theta)_* & & \\
 & & T(\phi(m_L)) & & \\
 T_p P_U: & T(\phi(m_L)) & \longrightarrow & \longrightarrow & T(\phi(m)) \\
 & \uparrow & & & \uparrow \\
 & T & & & T \\
 & \uparrow & & & \uparrow \\
 P_U: & \phi(m_L) & \longrightarrow & \Phi_W(\phi(m_L), \theta) & \longrightarrow & \phi(m) \\
 & \uparrow & & & \uparrow \\
 & \phi & & & \phi \\
 & \uparrow & & & \uparrow \\
 M: & m_L & \longrightarrow & \Phi_V(\theta, m_L) & \longrightarrow & m
 \end{array}$$

$$\begin{array}{ccccc}
 & & \text{Right} & & \\
 & & \Phi_W(\phi(m_R), -\theta)^* & & \\
 & & T(\phi(m_R)) & & \\
 T(\phi(m)) & \longleftarrow & \longleftarrow & \longleftarrow & T(\phi(m_R)) \\
 & \uparrow & & & \uparrow \\
 & T & & & T \\
 & \uparrow & & & \uparrow \\
 \phi(m) & \longleftarrow & \Phi_W(\phi(m_R), -\theta) & \longleftarrow & \phi(m_R) \\
 & \uparrow & & & \uparrow \\
 & \phi & & & \phi \\
 & \uparrow & & & \uparrow \\
 m & \longleftarrow & \Phi_V(-\theta, m_R) & \longleftarrow & m_R
 \end{array}$$

The Lie derivative of ϕ at m along W is then :

$$\mathcal{L}_W \phi(m) = \frac{\partial}{\partial \theta} \Phi_W(\phi(\Phi_V(-\theta, m)), \theta) |_{\theta=0}$$

and we can define the discontinuity in P_U as :

$$\Delta\phi(m) = \lim_{\theta \rightarrow 0} \frac{1}{\theta} (\Phi_W(\phi(\Phi_V(\theta, m)), -\theta) - \Phi_W(\phi(\Phi_V(-\theta, m)), \theta))$$

The Lie derivative $\mathcal{L}_W T$ of a section T of the tensorial bundle on TP_U (that is of a tensor which is defined at each point p in the tangent bundle TP_U) along a projectable vector field, can be defined similarly, if it is defined through a section on P_U : the value of T is given at a value of $p = \phi(m)$, and so by the point m . Because the Lie derivative is a natural operator (Maths.2226) the operation has the key features :

- $\mathcal{L}_W T$ is a tensor of the same type : same order, antisymmetric or symmetric
- $\mathcal{L}_W T$ belongs to the vertical bundle which projects on T
- if T transforms with a linear relation in a change of gauge (a natural operator) then $\mathcal{L}_W T$ transforms with the same relation.

And these features are shared with the computation of a discontinuity.

8.1.3 Representation of a discontinuity of the potential

We will prove the following :

Theorem 107 *A discontinuity of the derivative of the potential \dot{A} of a field, moving on a path with tangent V , can be represented as an antisymmetric tensor :*

$$\begin{aligned} J(\Delta\dot{A}) &= 4 \sum_{a=1}^m \sum_{i,j=0}^3 \left[\delta\dot{A} \right]_i^a V_j^* \varepsilon^i(m) \cdot \varepsilon^j(m) \otimes \vec{\theta}_a \\ &\in Q [T_1 Spin(3,1)^* \otimes T_1 U, \mathbf{Ad}^{-1} \times Ad] \\ \text{where :} \\ \delta\dot{A} &= \sum_{a=1}^m \sum_{i=0}^3 \left[\delta\dot{A} \right]_i^a \varepsilon^i(m) \otimes \vec{\theta}_a \\ V^* &= \sum_{i=0}^3 \sum_{\alpha\beta=0}^3 g_{\alpha\beta} V^\beta P_i^\alpha \varepsilon^i(m) \end{aligned}$$

The result holds for the gravitational field. This will be done in several steps, using P_U as example.

1st step : Discontinuities in the potential at one point

Lie derivative of the potential A principal connection on a principal bundle is a map \dot{A} from the tangent space of the fiber bundle to its vertical bundle. At the point $p \in P_U$:

$$\dot{A}(p) : T_p P_U \rightarrow V_p P_U : \dot{A}(p)(v^\alpha \partial m_\alpha + \zeta(v_g)(p)) = \zeta(v_g + Ad_{g^{-1}} \dot{A}(m) v_m)(p)$$

\dot{A} is the potential of the connection. This is a map : $\dot{A} : M \rightarrow T_1 U$ which depends only on m because the connection is principal.

The fundamental vector field $\zeta(p) : T_1 U \rightarrow V_p P_U$ as well as the vertical space do not depend on the trivialization. So the connection \dot{A} is a tensor field over P_U . To implement the method above we need a section of P_U to define the passage from $m \in M$ to $p \in P_U$. We take as gauge at each point $\mathbf{p}(m) = \varphi_U(m, 1)$ the gauge of the observer. Then $\dot{A}(\mathbf{p}(m)) = \zeta(\dot{A}(m) dm^\alpha)(\mathbf{p}(m))$

The discontinuity propagates along a vector V on M . W is the derivative of $\mathbf{p}(m)$ along V :

$$\mathbf{p}'(m)V = \varphi'_{Um}(m, 1)V(m) = \sum_{\alpha=0}^3 V^\alpha \partial m_\alpha$$

W is a projectable vector field which projects as $V : V = \pi'_P(W)$.

We can define the discontinuity of the tensor \dot{A} at $\mathbf{p}(a) = \varphi_U(a, 1) \in P_U$, along the projectable vector field W by :

$$\begin{aligned} \Delta_R \dot{A}(m, \theta) &= \Phi_W(\mathbf{p}(\Phi_V(\theta, m)), -\theta)^* \dot{A}(\mathbf{p}(\Phi_V(\theta, m))) - \dot{A}(\mathbf{p}(m)) \\ \Delta_L \dot{A}(m, \theta) &= \dot{A}(\mathbf{p}(m)) - \Phi_W(\mathbf{p}(\Phi_V(-\theta, m)), \theta)_* \dot{A}(\mathbf{p}(\Phi_V(-\theta, m))) \\ \text{and } \Delta \dot{A}(m) &= \lim_{\theta \rightarrow 0^+} \frac{1}{\theta} \left(\Delta_R \dot{A}(m, \theta) - \Delta_L \dot{A}(m, \theta) \right) \end{aligned}$$

$\dot{A}(\mathbf{p}(\Phi_V(\theta, m))) = \dot{A}(\mathbf{p}(m_R))$ reads :

$$\dot{A}(\mathbf{p}(m_R)) = \zeta \left(\sum_{\alpha} \dot{A}_\alpha(m_R) dm^\alpha \right) (\mathbf{p}(m_R))$$

The map $X(m, \theta) \rightarrow \zeta(X(m, \theta))(p)$ is linear and smooth,

$\lim_{\theta \rightarrow 0} \mathbf{p}(\Phi_V(\theta, m)) = \mathbf{p}(m)$ so :

$$\begin{aligned} \Delta \dot{A}(m) &= \zeta \left(\lim_{\theta \rightarrow 0^+} \frac{1}{\theta} \sum_{\alpha} \left(\dot{A}_\alpha(m_R) - \dot{A}_\alpha(m_L) \right) dm^\alpha \right) (\mathbf{p}(m)) \\ &= \zeta \left(\sum_{\alpha} \Delta \dot{A}_\alpha(m) dm^\alpha \right) (\mathbf{p}(m)) \end{aligned}$$

The discontinuity is expressed as a fundamental vector

$\zeta \left(\sum_{\alpha} \Delta \dot{A}_{\alpha}(m) dm^{\alpha} \right) (\mathbf{p}(m))$ where $\Delta \dot{A}(m) \in \Lambda_1(M; T_1U)$ is valued in the Lie algebra.

At this point we need to tell more about fundamental vectors denoted $\zeta(Y)(p)$ (see Maths.2066).

Fundamental vectors

The right action on P_U is the map :

$$\rho : P_U \times U \rightarrow P_U :: \rho(\varphi_U(m, g), g') = \varphi_U(m, g \cdot g')$$

It does not depend on the trivialization, and can be written : $\rho(p, g')$

Fundamental vectors are defined by the map :

$$\zeta : T_1U \rightarrow \mathfrak{X}(VP_U) :: \rho'_g(p, 1)(Y) = \varphi'_{Ug}(m, g)(L'_g 1)(Y) = \zeta(Y)(p)$$

This is a continuous linear map :

$$\zeta(\alpha Y + \beta Z)(p) = \alpha \zeta(Y)(p) + \beta \zeta(Z)(p)$$

which does not depend on the trivialization. With a fixed element Y of the Lie algebra T_1U the map ζ defines at any point p of P_U a vector on P_U belonging to the vertical bundle, that is $\zeta(Y)(p)$ is a vector field over the manifold P_U : its value depends on the point p of P_U , like a vector field over M depends on m , and its components in the holonomic basis $(\partial m, \partial g)$ given by a trivialization change according to the usual rules for vectors. Because it is vertical it reads in a holonomic basis at p :

$$\zeta(Y)(p) = \sum_a Y^a \varphi'_{Ug}(m, g)(L'_g 1) \vec{\theta}_a = \sum_a Y^a \partial \theta_a(p)$$

They have many properties, in particular *their value change with p as* :

$$\zeta(Y)(p \cdot g) = \rho'_p(p, g) \zeta(Ad_g Y)(p)$$

The map $\rho'_p(p, g) : T_p P_U \rightarrow T_{p \cdot g} P_U$ is linear and transports a basis of $T_p P_U$ in a basis of $T_{p \cdot g} P_U$, $\zeta(Y)(p \cdot g)$ and $\zeta(Y)(p)$ are vertical vectors, thus the rule is expressed as :

$$\begin{aligned} \partial \theta_a(p \cdot g) &= \rho'_p(p, g) \partial \theta_a(p) \\ \sum_a Y^a \partial \theta_a(p \cdot g) &= \sum_a [Ad_g]_b^a Y^b \partial \theta_a(p) \end{aligned}$$

Representation of the discontinuity in one point

The value of the discontinuity at $\mathbf{p}(m) = \varphi_U(m, 1)$ is $\zeta \left(\Delta \dot{A}(m) \right) (\mathbf{p}(m))$

The observer measures the discontinuity as the potential, in its standard gauge : $\mathbf{p}(m) = \varphi_U(m, 1)$, which is not a section, but just a particular point of P_U in the fiber over m . The value measured by the observer is $\zeta \left(\Delta \dot{A}(m) \right) (\mathbf{p}(m))$. It depends on m and is expressed in the basis of $T_p P_U$ at \mathbf{p} as a vertical vector :

$$\zeta \left(\Delta \dot{A}(m) \right) (\mathbf{p}(m)) = \sum_a \left(\Delta \dot{A}(m) \right)^a \partial \theta_a(\mathbf{p})$$

that is as a vector of T_1U , varying with m .

But it depends also on the trivialization, which entails that in a change of standard gauge the discontinuity is measured at another point :

$$\tilde{\mathbf{p}} = \tilde{\varphi}_U(m, 1) = \varphi_U(m, \chi(m)) = \mathbf{p} \cdot \chi(m) \text{ so the measure changes as :}$$

$$\zeta \left(\Delta \dot{A}(m) \right) (\tilde{\mathbf{p}}) = \zeta \left(\Delta \dot{A}(m) \right) (\mathbf{p} \cdot \chi(m))$$

$$\begin{aligned}
&= \rho'_p(\mathbf{p}, \chi(m)) \zeta \left(Ad_{\chi(m)} \Delta \dot{A}(m) \right) (\mathbf{p}) \\
&\text{in components :} \\
&\sum_a \left(\Delta \dot{A}(m) \right)^a \partial \theta_a(\mathbf{p}) \rightarrow \\
&\sum_a \left(\widetilde{\Delta \dot{A}(m)} \right)^a \partial \theta_a(\tilde{\mathbf{p}}) = \sum_{ab} [Ad_{\chi(m)}]_b^a \left(\Delta \dot{A}(m) \right)^b \partial \theta_a(\tilde{\mathbf{p}})
\end{aligned}$$

So the measure of the discontinuity changes as $Ad_{\chi(m)} \Delta \dot{A}(m)$. We have the same rule as for the strength of the field \mathcal{F}_A . *The discontinuity at m can be represented as a one form on M valued in the adjoint bundle $P_U [T_1 U, Ad]$.*

This result is important : the reason why a potential cannot explicitly be present in the lagrangian comes from its special rule in a change of gauge (see lagrangian), but this rule applies no longer to $\zeta \left(\Delta \dot{A}(m) \right) (p)$, which becomes a section of an associated vector bundle like the others. In the Standard Model bosons are represented by potentials, changing in the usual way, so that, to be part of a lagrangian, a patch is necessary, provided for by the Higgs Boson. This patch is no longer required if bosons are represented as discontinuities in the fields (the other reason to introduce the Higgs boson comes from the use of the Dirac's operator, and we have dealt with it).

2nd step : Propagation of the discontinuity

The discontinuity is seen as a change (the discontinuity is additive to the derivative of \dot{A}) in the value of a continuous potential, which propagates along a vector V , which is part of the definition of the discontinuity. It should propagate with the field itself, so we should have $\langle V, V \rangle = 0$. Vector fields such that $\langle V, V \rangle = 0$ (null vector fields) are not geodesics, and they cannot be defined by a point and a vector of origin. We have seen that fields should propagate at c , however the propagation of the fields other than the EM field is not well known, and we will reserve this point and assume only that V is a future oriented vector.

2 form valued in the Lie algebra

\dot{A} is a map :

$$\dot{A} : M \rightarrow TM^* \otimes T_1 U :: \dot{A}(m) = \sum_{a=1}^m \sum_{\alpha=0}^3 \dot{A}_\alpha^a(m) d\xi^\alpha \otimes \vec{\theta}_a$$

Its derivative is :

$$\dot{A}' : M \rightarrow TM^* \otimes TM^* \otimes T_1 U :: \dot{A}'(m) = \sum_{a=1}^m \sum_{\alpha\beta=0}^3 \partial_\beta \dot{A}_\alpha^a(m) d\xi^\beta \otimes d\xi^\alpha \otimes \vec{\theta}_a$$

The derivative of \dot{A} along the vector field V is expressed as :

$$\dot{A}'(m)(V) = \sum_{a=1}^m \sum_{\alpha\beta=0}^3 \partial_\beta \dot{A}_\alpha^a(m) V^\beta d\xi^\alpha \otimes \vec{\theta}_a$$

So we can write :

$$\Delta \dot{A}(m) = \sum_{a=1}^m \sum_{\alpha\beta=0}^3 \delta \dot{A}_\alpha^a(m) V_\beta^* d\xi^\alpha \otimes d\xi^\beta \otimes \vec{\theta}_a$$

where V^* is the dual of V : $V^* = \sum_{\lambda=0}^3 g_{\alpha\lambda} V^\lambda d\xi^\alpha$

The tensor $\Delta \dot{A}$ is not necessarily symmetric in α, β .

Fields have a density of energy which, at equilibrium and in the vacuum, reads : $\sum_{\alpha\beta=0}^3 \sum_{a=1}^m \mathcal{F}_A^{a\alpha\beta} \partial_\alpha \dot{A}_\beta$. The same quantity can be computed for the

discontinuities : $\sum_{\alpha\beta=0}^3 \sum_{a=1}^m \mathcal{F}_A^{a\alpha\beta} \Delta \dot{A}_{\alpha\beta}^a$. Physically it represents the discontinuity in the density of energy. By interchanging the indexes α, β we get :

$$\begin{aligned} \sum_{\alpha\beta=0}^3 \sum_{a=1}^m \mathcal{F}_A^{a\alpha\beta} \Delta \dot{A}_{\alpha\beta}^a &\rightarrow \sum_{\alpha\beta=0}^3 \sum_{a=1}^m \mathcal{F}_A^{a\beta\alpha} \Delta \dot{A}_{\beta\alpha}^a \\ &= - \sum_{\alpha\beta=0}^3 \sum_{a=1}^m \mathcal{F}_A^{a\alpha\beta} \Delta \dot{A}_{\beta\alpha}^a \end{aligned}$$

thus, necessarily, $\Delta \dot{A}_{\beta\alpha}^a = -\Delta \dot{A}_{\alpha\beta}^a$: it is an antisymmetric tensor, and we can write :

$$\Delta \dot{A} = \frac{1}{2} \left(\sum_{a=1}^m \sum_{\alpha=0}^3 \delta \dot{A}_{\alpha}^a d\xi^{\alpha} \otimes \vec{\theta}_a \right) \wedge \left(\sum_{\beta=0}^3 V_{\beta}^* d\xi^{\beta} \right) \quad (8.1)$$

$$\Delta \dot{A} = \sum_{a=1}^m \sum_{\{\alpha\beta\}=0}^3 \left(\delta \dot{A}_{\alpha}^a V_{\beta}^* - \delta \dot{A}_{\beta}^a V_{\alpha}^* \right) d\xi^{\alpha} \wedge d\xi^{\beta} \otimes \vec{\theta}_a$$

Representation in the Clifford algebra

$\Delta \dot{A}$ is a 2 form on M, valued in the Lie algebra, as the strength of the field. We can proceed as for the strength of the field and associate to $\Delta \dot{A}$ the element of $T_1 Spin(3, 1)^* \otimes T_1 U$:

$$\begin{aligned} j(\Delta \dot{A}) &= 2 \sum_{a=1}^m \{ (\delta \dot{A}_3^a V_2^* - \delta \dot{A}_2^a V_3^*) \vec{\kappa}^1 + (\delta \dot{A}_1^a V_3^* - \delta \dot{A}_3^a V_1^*) \vec{\kappa}^2 \\ &+ (\delta \dot{A}_2^a V_1^* - \delta \dot{A}_1^a V_2^*) \vec{\kappa}^3 + (\delta \dot{A}_0^a V_1^* - \delta \dot{A}_1^a V_0^*) \vec{\kappa}^4 + (\delta \dot{A}_0^a V_2^* - \delta \dot{A}_2^a V_0^*) \vec{\kappa}^5 \\ &+ (\delta \dot{A}_0^a V_3^* - \delta \dot{A}_3^a V_0^*) \vec{\kappa}^6 \} \otimes \vec{\theta}_a \in T_1 Spin(3, 1)^* \otimes T_1 U \end{aligned}$$

where the components V and $\delta \dot{A}^a$ are expressed in the basis $(\varepsilon^i(m))_{i=0}^3$

$[j(\Delta \dot{A})]$ is a $m \times 6$ matrix.

Replacing $\vec{\kappa}^a$ by their value in $\varepsilon^i \cdot \varepsilon^j$ a straightforward computation gives :

$$\begin{aligned} j(\Delta \dot{A}) &= \sum_{a=1}^m \{ (\delta \dot{A}_3^a V_2^* - \delta \dot{A}_2^a V_3^*) \varepsilon^3 \cdot \varepsilon^2 + (\delta \dot{A}_1^a V_3^* - \delta \dot{A}_3^a V_1^*) \varepsilon^1 \cdot \varepsilon^3 \\ &+ (\delta \dot{A}_2^a V_1^* - \delta \dot{A}_1^a V_2^*) \varepsilon^2 \cdot \varepsilon^1 + (\delta \dot{A}_0^a V_1^* - \delta \dot{A}_1^a V_0^*) \varepsilon^0 \cdot \varepsilon^1 \\ &+ (\delta \dot{A}_0^a V_2^* - \delta \dot{A}_2^a V_0^*) \varepsilon^0 \cdot \varepsilon^2 + (\delta \dot{A}_0^a V_3^* - \delta \dot{A}_3^a V_0^*) \varepsilon^0 \cdot \varepsilon^3 \} \otimes \vec{\theta}_a \\ &= \sum_{a=1}^m \sum_{i=0}^3 \delta \dot{A}_i^a \varepsilon^i \cdot \left(\sum_{j=0}^3 V_j^* \varepsilon^j - V_i^* \varepsilon^i \right) \otimes \vec{\theta}_a \\ &= \sum_{a=1}^m \left(\delta \dot{A}^a \cdot V^* - \sum_{i=0}^3 \delta \dot{A}_i^a V_i^* \varepsilon^i \cdot \varepsilon^i \right) \otimes \vec{\theta}_a \\ &= \sum_{a=1}^m \left(\delta \dot{A}^a \cdot V^* - \langle \delta \dot{A}^a, V^* \rangle \right) \otimes \vec{\theta}_a \\ &= \sum_{a=1}^m \left(\delta \dot{A}^a \cdot V^* - V^* \cdot \delta \dot{A}^a \right) \otimes \vec{\theta}_a \text{ with : } \delta \dot{A}^a \cdot V^* + V^* \cdot \delta \dot{A}^a = 2 \langle \delta \dot{A}^a, V^* \rangle \\ j(\Delta \dot{A}^a) &= \left(\delta \dot{A}^a \cdot V^* - V^* \cdot \delta \dot{A}^a \right) \otimes \vec{\theta}_a \quad (8.2) \end{aligned}$$

We can represent equivalently $j(\Delta \dot{A}^a)$ in the basis $\vec{\kappa}^a$ of $T_1 Spin(3, 1)^*$ or $\varepsilon^i \cdot \varepsilon^j$ of $Cl(3, 1)^*$.

In the basis $\vec{\kappa}^a$:

$$\begin{aligned} j(\Delta \dot{A}) &= \sum_{a=1}^m v^* ((\Delta F_{32}^a, \Delta F_{13}^a, \Delta F_{21}^a), (\Delta F_{01}^a, \Delta F_{02}^a, \Delta F_{03}^a)) \\ &= v^* (\Delta F_R^a, \Delta F_W^a) \otimes \vec{\theta}_a \in T_1 Spin(3, 1)^* \otimes T_1 U \end{aligned}$$

with

$$\begin{aligned}\Delta F_R^a &= 2 \left(\left(\delta \dot{A}_3^a V_2^* - \delta \dot{A}_2^a V_3^* \right), \left(\delta \dot{A}_1^a V_3^* - \delta \dot{A}_3^a V_1^* \right), \left(\delta \dot{A}_2^a V_1^* - \delta \dot{A}_1^a V_2^* \right) \right) \\ \Delta F_W^a &= 2 \left(\left(\delta \dot{A}_0^a V_1^* - \delta \dot{A}_1^a V_0^* \right), \left(\delta \dot{A}_0^a V_2^* - \delta \dot{A}_2^a V_0^* \right), \left(\delta \dot{A}_0^a V_3^* - \delta \dot{A}_3^a V_0^* \right) \right) \\ (\Delta F_R^a)^t \Delta F_W^a &= 0 \\ (\Delta F_R^a)^t \Delta F_R^a - (\Delta F_W^a)^t \Delta F_W^a &= \langle \delta \dot{A}^a, \delta \dot{A}^a \rangle \langle V, V \rangle - \langle \delta \dot{A}^a, V \rangle^2\end{aligned}$$

There should be a gauge $q(m, \vartheta(\sigma, g))$ depending on the location $\Phi_V(t, a)$ on the propagation in which the discontinuity has a constant value $B_0 \in T_1 Spin(3, 1)^* \otimes T_1 U$:

$$\begin{aligned}\left(q(m, \vartheta(1, 1)), j(\Delta \dot{A}) \right) &\sim \left(q(m, \vartheta(\sigma^{-1}, g^{-1})), Ad_g j(\Delta \dot{A}) \mathbf{Ad}_{\sigma^{-1}} \right) \\ &= \left(q(m, \vartheta(\sigma^{-1}, g^{-1})), B_0 \right)\end{aligned}$$

In this gauge the tensors $\delta \dot{A}, V^*$ do not change. We could expect that in this gauge the axis $\tilde{\varepsilon}_0$ would be in the direction of V , but this is impossible if $\langle V, V \rangle = 0 : V = k\tilde{\varepsilon}_0 \Rightarrow \langle V, V \rangle = -k^2$. This is an important difference between bosons and particles : they do not have a privileged gauge (at least if they propagate at the speed of light) and as a consequence the concept of relativist momentum loses its main purpose. Notice that this remark holds, whatever the formalism used to represent the bosons.

8.1.4 Properties of Bosons

Such discontinuities propagate on a curve, as a particle. They can be assimilated to the quasi-particles which, in QTF, are the carriers of forces fields, gravitons for the gravitational field, and bosons for the other fields. And we state :

Proposition 108 *Bosons are discontinuities of the derivative of the potential \dot{A} , which propagate along a path with velocity V , future oriented. They can be represented at each point of their trajectory as a tensor :*

$$j(\Delta \dot{A})(t) = \left(\delta \dot{A}^a \cdot V^* - V^* \cdot \delta \dot{A}^a \right) \otimes \vec{\theta}_a$$

$$\in Q [T_1 Spin(3, 1)^* \otimes T_1 U, \mathbf{Ad}^{-1} \times Ad]$$

where $\delta \dot{A} \in Q [\mathbb{R}^{4*} \otimes T_1 U, \mathbf{Ad}^{-1} \times Ad]$, V^* is the dual of V

Gravitons are discontinuities of the derivative of the potential G , which propagate along a path with velocity V , future oriented. They can be represented at each point of their trajectory as a tensor :

$$j(\Delta G)(t) = \left(\delta G^a \cdot V^* - V^* \cdot \delta G^a \right) \otimes \vec{\kappa}_a$$

$$\in P_G [T_1 Spin(3, 1)^* \otimes T_1 Spin(3, 1), \mathbf{Ad}^{-1} \times \mathbf{Ad}]$$

where $\delta G \in P_G [\mathbb{R}^{4*} \otimes T_1 Spin(3, 1), \mathbf{Ad}^{-1} \times \mathbf{Ad}]$, V^* is the dual of V

Equation of propagation

The trajectory of a discontinuity is represented by a map :

$$j(\Delta \dot{A}) : [0, T] \rightarrow Q [T_1 Spin(3, 1)^* \otimes T_1 U, \mathbf{Ad}^{-1} \times Ad] :$$

$$j(\Delta\dot{A})(\mathbf{q}(\Phi_V(t, a))) \\ = (\mathbf{q}(\Phi_V(t, a), 1), \sum_{a=1}^m (\delta\dot{A}^a(t) \cdot V^*(t) - V^*(t) \cdot \delta\dot{A}^a(t)) \otimes \vec{\theta}_a)$$

The set L_A^1 of maps $C([0, T]; Q[T_1Spin(3, 1)^* \otimes T_1U])$ such that

$\int_0^T \|j(\Delta\dot{A})(\mathbf{q}(\Phi_V(t, a)))\| dt < \infty$ is a Fréchet space, and a representation of $Spin(3, 1) \times U$.

The propagation of bosons is continuous, if it is also determinist then the theorem 25 applies. There is Hilbert space H , including $T_1Spin(3, 1)^* \otimes T_1U$, and a unitary operator $\Theta(t)$ such that :

$$j(\Delta\dot{A})(\mathbf{q}(\Phi_V(t, a))) = \Theta(t)j(\Delta\dot{A})(\mathbf{q}(\Phi_V(0, a))) \quad (8.3)$$

which is the propagation law for the discontinuity. This is a geometric equation, and it holds whatever the gauge. In a global change of gauge :

$$\mathbf{q}(m) = \varphi_Q(m, (1, 1)) \rightarrow \tilde{\mathbf{q}}(m) = \tilde{\varphi}_Q(m, (1, 1)) = \mathbf{q}(m) \cdot (\sigma^{-1}, g^{-1})$$

$$\mathbf{Ad}_{\sigma^{-1}j}(\Delta\dot{A})(\mathbf{q}(\Phi_V(t, a))) Ad_g = \Theta(t) \mathbf{Ad}_{\sigma^{-1}j}(\Delta\dot{A})(\mathbf{q}(\Phi_V(0, a))) Ad_g = \\ \mathbf{Ad}_{\sigma^{-1}}\Theta(t)j(\Delta\dot{A})(\mathbf{q}(\Phi_V(0, a))) Ad_g$$

$\Theta(t)$ commute with both $\mathbf{Ad}_{\sigma^{-1}}$ and Ad_g , in any global change of gauge. Thus it is necessarily an isomorphism by the Schur's lemna (Maths.1909).

The result holds for the gravitational field.

Spin of a boson

We can proceed as for particles and fields.

Using the projection (see Clifford Algebra) :

$$\pi_L : T_1Spin(3, 1)^* \rightarrow L_0^* :: \pi_L^*(X) = \frac{1}{2}(X - \varepsilon^0 \cdot X \cdot \varepsilon^0)$$

we define :

$$\pi : T_1Spin(3, 1)^* \otimes T_1U \rightarrow L_0^* \otimes T_1U ::$$

$$\pi(j(\Delta\dot{A})) = \frac{1}{2}(j(\Delta\dot{A}) - \varepsilon^0 \cdot j(\Delta\dot{A}) \cdot \varepsilon^0)$$

$$\pi(j(\Delta\dot{A})) = \sum_{a=1}^m \{(\delta\dot{A}_3^a V_2^* - \delta\dot{A}_2^a V_3^*) \varepsilon^3 \cdot \varepsilon^2$$

$$+ (\delta\dot{A}_1^a V_3^* - \delta\dot{A}_3^a V_1^*) \varepsilon^1 \cdot \varepsilon^3 + (\delta\dot{A}_2^a V_1^* - \delta\dot{A}_1^a V_2^*) \varepsilon^2 \cdot \varepsilon^1\} \otimes \vec{\theta}_a$$

and the relation of equivalence on L_A^1 :

$$j(\Delta\dot{A}) \sim j(\Delta\dot{A}') \Leftrightarrow \pi(j(\Delta\dot{A})) = \pi(j(\Delta\dot{A}'))$$

we define the **Spin of the boson** as $\pi(j(\Delta\dot{A}))$.

It can be represented by a map in L_A^1 , whose measure changes with the action of $SO(3)$. It is quantized and belongs to a $2j + 1$ dimensional vector space, isomorphic to (P^j, D^j) with $j \in \mathbb{N}$. *The boson has an integer Spin.*

The experiments show that actually bosons have a spin = 1.

Spin and Curvature of gravitons

For gravitons the situation is more interesting. A graviton reads :

$$\begin{aligned} j(\Delta G) &= \sum_{a=1}^6 (\delta G^a \cdot V^* - V^* \cdot \delta G^a) \otimes \vec{\kappa}_a \\ &\in P_G [T_1 Spin(3, 1)^* \otimes T_1 Spin(3, 1), \mathbf{Ad}^{-1} \times \mathbf{Ad}] \\ j(\Delta G) &= \\ &2\{(\delta G_3^a V_2^* - \delta G_2^a V_3^*) \vec{\kappa}^1 + (\delta G_1^a V_3^* - \delta G_3^a V_1^*) \vec{\kappa}^2 + (\delta G_2^a V_1^* - \delta G_1^a V_2^*) \vec{\kappa}^3 \\ &+ (\delta G_0^a V_1^* - \delta G_1^a V_0^*) \vec{\kappa}^4 + (\delta G_0^a V_2^* - \delta G_2^a V_0^*) \vec{\kappa}^5 + (\delta G_0^a V_3^* - \delta G_3^a V_0^*) \vec{\kappa}^6\} \otimes \\ &\vec{\kappa}_a \end{aligned}$$

Let for $a=1,2,3$: $\delta G^a = \delta G_r^a, \delta G^{a+3} = \delta G_w^a$
 $[\delta G_r^a] = [\delta G_{r0}^a \quad \delta G_{r1}^a \quad \delta G_{r2}^a \quad \delta G_{r3}^a]$ be a 1 row 4 columns matrix, labeled from 0 to 3,

$[\delta G_w^a] = [\delta G_{w0}^a \quad \delta G_{w1}^a \quad \delta G_{w2}^a \quad \delta G_{w3}^a]$ be a 1 row 4 columns matrix, labeled from 0 to 3,

$[V^*] = [V_0^* \quad V_1^* \quad V_2^* \quad V_3^*]$ be a 1 row 4 columns matrix, labeled from 0 to 3,

$[\Delta G_r^a] = [\delta G_r^a]^t [V^*] - [V^*]^t [\delta G_r^a]$ is a 4×4 matrix, with entries labeled from 0 to 3,

$$\begin{aligned} j(\Delta G) &= \sum_{i,j=0}^3 [\Delta G_r^1]_j^i \varepsilon^i \cdot \varepsilon^j \otimes \varepsilon_3 \cdot \varepsilon_2 + [\Delta G_r^2]_j^i \varepsilon^i \cdot \varepsilon^j \otimes \varepsilon_1 \cdot \varepsilon_3 \\ &+ [\Delta G_r^1]_j^i \varepsilon^i \cdot \varepsilon^j \otimes \varepsilon_2 \cdot \varepsilon_1 + [\Delta G_w^1]_j^i \varepsilon^i \cdot \varepsilon^j \otimes \varepsilon_0 \cdot \varepsilon_1 \\ &+ [\Delta G_w^2]_j^i \varepsilon^i \cdot \varepsilon^j \otimes \varepsilon_0 \cdot \varepsilon_2 + [\Delta G_w^3]_j^i \varepsilon^i \cdot \varepsilon^j \otimes \varepsilon_0 \cdot \varepsilon_3 \end{aligned}$$

We can proceed as above and single out the components which do not involve ε_0 :

$$\pi(j(\Delta G)) = \sum_{i,j=1}^3 [\Delta G_r^1]_j^i \varepsilon^i \cdot \varepsilon^j \otimes \varepsilon_3 \cdot \varepsilon_2 + [\Delta G_r^2]_j^i \varepsilon^i \cdot \varepsilon^j \otimes \varepsilon_1 \cdot \varepsilon_3 + [\Delta G_r^1]_j^i \varepsilon^i \cdot \varepsilon^j \otimes \varepsilon_2 \cdot \varepsilon_1$$

which can then be defined as the Spin of the Graviton. It can be represented by a map in L_G^1 , whose measure changes with the action of $SO(3)$. It is quantized and belongs to a $2j+1$ dimensional vector space, isomorphic to (P^j, D^j) with $j \in \mathbb{N}$. *The graviton has an integer Spin.* It would impact the same components of the gravitational field, that we have identified with the spin of the field, and as the usual spatial gravity tensor.

The contraction on the 2nd and 4th indices gives a tensor :

$$\begin{aligned} j(\Delta G)_1 &= \sum_{i=0}^3 [\Delta G_r^1]_2^i \varepsilon^i \otimes \varepsilon_3 + [\Delta G_r^2]_3^i \varepsilon^i \otimes \varepsilon_1 + [\Delta G_r^1]_1^i \varepsilon^i \otimes \varepsilon_2 \\ &+ \left(([\Delta G_w^1]_1^i + [\Delta G_w^2]_2^i + [\Delta G_w^3]_3^i) \varepsilon^i \right) \otimes \varepsilon^0 \end{aligned}$$

and a scalar :

$$R_1 = [\Delta G_r^1]_2^3 + [\Delta G_r^2]_3^1 + [\Delta G_r^1]_1^2 + [\Delta G_w^1]_1^0 + [\Delta G_w^2]_2^0 + [\Delta G_w^3]_3^0$$

The contraction on the 1st and 3d indices gives a tensor :

$$\begin{aligned} j(\Delta G)_2 &= \sum_{j=0}^3 [\Delta G_r^1]_j^3 \varepsilon^j \otimes \varepsilon_2 + [\Delta G_r^2]_j^1 \varepsilon^j \otimes \varepsilon_3 + [\Delta G_r^1]_j^2 \varepsilon^j \otimes \varepsilon_1 + [\Delta G_w^1]_j^0 \varepsilon^j \otimes \\ &\varepsilon_1 + [\Delta G_w^2]_j^0 \varepsilon^j \otimes \varepsilon_2 + [\Delta G_w^3]_j^0 \varepsilon^j \otimes \varepsilon_3 \\ j(\Delta G)_2 &= \sum_{j=0}^3 \left([\Delta G_r^1]_j^2 + [\Delta G_w^1]_j^0 \right) \varepsilon^j \otimes \varepsilon_1 \end{aligned}$$

$$+ \left([\Delta G_r^1]_j^3 + [\Delta G_w^2]_j^0 \right) \varepsilon^j \otimes \varepsilon_2 + \left([\Delta G_r^2]_j^1 + [\Delta G_w^3]_j^0 \right) \varepsilon^j \otimes \varepsilon_3$$

and a scalar :

$$R_2 = [\Delta G_r^1]_1^2 + [\Delta G_w^1]_1^0 + [\Delta G_r^1]_2^3 + [\Delta G_w^2]_2^0 + [\Delta G_r^2]_3^1 + [\Delta G_w^3]_3^0 = R_1$$

The contraction on all indices for j (ΔG) gives the curvature :

$$\delta R = [\Delta G_r^1]_2^3 + [\Delta G_r^2]_3^1 + [\Delta G_r^1]_1^2 + [\Delta G_w^1]_1^0 + [\Delta G_w^2]_2^0 + [\Delta G_w^3]_3^0 = R_1 = R_2$$

The curvature of the field is :

$$R' = 4 (F_{r32}^1 + F_{r13}^2 + F_{r21}^3 + F_{w01}^1 + F_{w02}^2 + F_{w03}^3)$$

so δR appears as a discontinuity of the scalar curvature. All these quantities are equivariant in a change of gauge.

In the absence of any experimental evidence of gravitons, it is difficult to assess the physical meaning of these tensors. However we can assume that the curvature δR is constant along the trajectory, and is one of the characteristic of the graviton.

Charge of a boson

The charges have been defined with respect to the vector space F . But we have seen that in a continuous model they appear in currents :

$$\phi_A = \sum_{\beta} \left[\mathcal{F}_A^{\alpha\beta}, \dot{A}_{\beta} \right] \vec{\theta}_a \otimes \partial \xi_{\alpha} \in T_1 U \otimes TM$$

The simplest expression of the current created by the boson is then :

$$\begin{aligned} \phi_B &= \sum_{a=1}^m \sum_{\beta, \alpha} \left[\mathcal{F}_A^{\alpha\beta}, \Delta \dot{A}(V)_{\beta} \right] \vec{\theta}_a \otimes \partial \xi_{\alpha} \\ \Delta \dot{A}(V) &= \sum_{a=1}^m \left(\langle V, V \rangle \delta \dot{A}^a - \delta \dot{A}^a(V) V_* \right) \otimes \vec{\theta}_a \end{aligned}$$

$$\phi_B = \sum_{a=1}^m \sum_{\beta, \alpha} \left(\langle V, V \rangle \left[\mathcal{F}_A^{\alpha\beta}, \delta \dot{A}_{\beta} \right]^a - V_{\beta}^* \left[\mathcal{F}_A^{\alpha\beta}, \delta \dot{A}(V) \right]^a \right) \vec{\theta}_a \otimes \partial \xi_{\alpha} \quad (8.4)$$

In the orthonormal basis :

$$\mathcal{F}_A^{\alpha\beta} = \sum_{ij=0}^3 P_i^{\alpha} P_j^{\beta} F^{aij}$$

$$V_{\beta}^* = \sum_p V^p \eta_{pp} P_{\beta}^p$$

$$\delta \dot{A}_{\beta} = \sum_q P_{\beta}^q \delta \dot{A}_q$$

$$\phi_B = \sum_{a=1}^m \sum_{ij=0}^3 \left(\langle V, V \rangle \left[F_A^{ij}, \delta \dot{A}_j \right]^a - \eta_{jj} V^j \left[F_A^{ij}, \delta \dot{A}(V) \right]^a \right) \vec{\theta}_a \otimes \varepsilon_i \quad (8.5)$$

The current created by a photon is null, because $U(1)$ is abelian.

If the tensor δA was decomposable : $\delta \dot{A} = \delta C \otimes V^*$ then $\delta \dot{A}(V) = \delta C \langle V, V \rangle$ and the current would be null :

$$\begin{aligned} \phi_B &= \\ \sum_{a=1}^m \sum_{ij=0}^3 \left(\langle V, V \rangle V_j \left[F_A^{ij}, \delta C \right]^a - \langle V, V \rangle V_j \left[F_A^{ij}, \delta C \right]^a \right) \vec{\theta}_a \otimes \varepsilon_i &= 0 \end{aligned}$$

Actually the charge of the boson is measured, as the charge of any particle, with respect to its behavior when submitted to a field. So we can say that two bosons have the same charge if there is $C(\Delta\dot{A}) \in T_1U$ such that :

$$C(\Delta\dot{A}) = C(\Delta\dot{A}') \Leftrightarrow \phi_B = C(\Delta\dot{A}) \otimes V, \phi'_B = C(\Delta\dot{A}') \otimes V'$$

That is : there is a projection π_C from $T_1U \otimes TM$ onto T_1U such that $\pi_C(\phi_B) = \pi_C(\phi'_B)$

The projection should follow the rules in a change of gauge which sum up in : $\forall g \in U : \pi_C \circ Ad_g = Ad_g \circ \pi_C$

Using the notations above, the trajectories of bosons can be represented by a map : $j(\Delta\dot{A}) : [0, T] \rightarrow Q[T_1Spin(3, 1)^* \otimes T_1U, \mathbf{Ad}^{-1} \times Ad]$ belonging to a Fréchet space. The quantization of $j(\Delta\dot{A})$, accounting for the compacity of U , leads to tell that $j(\Delta\dot{A})$ belongs to a vector space isomorphic to an irreducible, finite dimensional, representation of U . And actually, in the Standard Model, F itself is the sum of irreducible representations, accounting for the generations feature. Moreover if we add the variable $\pi_C(\phi_B)$, assuming that the projection does exist, then we have families of bosons, characterized by their charge. And indeed this is what is seen experimentally : there are 12 kinds of elementary bosons, which can be combined together and carry a specific charge encompassing several dimensions a of the fields (so the tensor $j(\Delta\dot{A})$ is not decomposable). And the photon (when only the EM field is considered) has no charge.

So we are lead to state, in our framework :

Proposition 109 *The currents associated to bosons are decomposable tensors, and there is a charge $C(\Delta\dot{A}) \in T_1U$, specific to each type of boson, such that :*

$$\phi_B = C(\Delta\dot{A}) \otimes V \quad (8.6)$$

The same idea could be implemented to gravitons, but $C(\Delta G)$ would not be a scalar, contrary to the mass. So it seems more speculative.

8.1.5 Relativist Momentum and Energy of bosons

The issues

$j(\Delta\dot{A})$ can be considered as the state of the boson, as measured by the observer.

This is a geometric quantity. The scalar product is

$$\begin{aligned} \langle j(\Delta\dot{A}), j(\Delta\dot{A}) \rangle &= -2 \sum_{\alpha\beta=0}^3 \sum_{a=1}^m \Delta\dot{A}^{a\alpha\beta} \Delta\dot{A}_{\alpha\beta}^a \\ &= -2 \sum_{a=1}^m \sum_{\alpha\beta=0}^3 (\delta\dot{A}^{a\alpha} V^\beta) (\delta\dot{A}_\alpha^a V_\beta^*) = -2 \langle V, V \rangle \sum_{a=1}^m g^{\alpha\beta} \delta\dot{A}_\alpha^a \delta\dot{A}_\beta^a \\ \langle j(\Delta\dot{A}), j(\Delta\dot{A}) \rangle &= -2 \langle V, V \rangle \sum_{a=1}^m \langle \delta\dot{A}^a, \delta\dot{A}^a \rangle \end{aligned} \quad (8.7)$$

$$\begin{aligned}
\langle j(\Delta G), j(\Delta G) \rangle &= \frac{1}{2} \sum_{\alpha\beta=0}^3 \left(\sum_{a=1}^3 \Delta G_r^{a\alpha\beta} \Delta G_{r\alpha\beta}^a - \Delta G_w^{a\alpha\beta} \Delta G_{w\alpha\beta}^a \right) \\
&= \frac{1}{2} \sum_{\alpha\beta=0}^3 \left(\sum_{a=1}^3 \delta G_r^{a\alpha} V^\beta \delta G_{r\alpha\beta}^a V_\beta^* - \delta G_w^{a\alpha} V^\beta \delta G_{w\alpha\beta}^a V_\beta^* \right) \\
&= \frac{1}{2} \langle V, V \rangle \sum_{\alpha\beta=0}^3 \left(\sum_{a=1}^3 \delta G_r^{a\alpha} \delta G_{r\alpha}^a - \delta G_w^{a\alpha} \delta G_{w\alpha}^a \right) \\
\langle j(\Delta G), j(\Delta G) \rangle &= \frac{1}{2} \langle V, V \rangle \sum_{\alpha\beta=0}^3 \left(\sum_{a=1}^3 \langle \delta G_r^a, \delta G_r^a \rangle - \langle \delta G_w^a, \delta G_w^a \rangle \right) \quad (8.8)
\end{aligned}$$

If the mass is $M_B^2 = \frac{1}{c^2} \langle j(\Delta \dot{A}), j(\Delta \dot{A}) \rangle$ we see that it is null if the boson travels at the speed of light, as expected.

The equivalent of the Relativist Momentum of a boson is $j(\Delta \dot{A})$. However there are several issues.

The first is that $j(\Delta \dot{A})$ is not valued in the same space $E \otimes F$ as the particles. One could conceive of a representation through the map :

$$D = \gamma C \otimes \varrho'(1) : T_1 Spin(3,1)^* \otimes T_1 U \rightarrow \mathcal{L}(E \otimes F; E \otimes F)$$

$$D \left(j(\Delta \dot{A}) \right) \psi =$$

$$\sum_{i=1}^4 \sum_{j=1}^n \sum_{a=1}^m \sum_{b=1}^6 \left[j(\Delta \dot{A}) \right]_b^a ([\gamma C(\vec{\kappa}^b)] [B] [\theta_a])_j^i e_i \otimes f_j$$

acting on a fixed tensor B but the action is not consistent with the rules on $Q[E \otimes F, \vartheta]$.

The second is that there is no privileged gauge along which the boson would be in a fundamental, constant, state.

The third is that $\langle j(\Delta \dot{A}), j(\Delta \dot{A}) \rangle$ is null for bosons travelling at the speed of light. This is consistent with a null mass for the photon, but problematic for the energy.

The only general, and sensible, definition of the energy of a boson is in relation with the underlying field, by :

$$\delta \mathcal{E} = -2 \sum_{a=1}^m \sum_{\alpha\beta=0}^3 \langle \mathcal{F}_A^{a\alpha\beta}, \Delta \dot{A}_{\alpha\beta}^a \rangle = \langle j(\mathcal{F}_A), j(\Delta \dot{A}) \rangle \quad (8.9)$$

where $\delta \mathcal{E}$ is a discrepancy with respect to the density of energy of the underlying field.

For the photon, which is by far and large the best known boson, a historic representation prevails, but which leaves the issue of the definition of energy a bit confusing. For the other bosons the definition of mass and energy, is complicated, and involves screening by virtual particles. So there is no unified and satisfying solution, which shows the limits of the similitude between bosons and particles.

The representation of photons

$$j(\Delta\dot{A}) = v^*(\Delta F_R, \Delta F_W)$$

where $\Delta F_R, \Delta F_W$ are two orthogonal vectors of \mathbb{R}^3

With :

$$V^* = v^* + c\varepsilon^0$$

$$\delta\dot{A} = \delta a + \delta\dot{A}_0\varepsilon^0$$

$$\Delta F_R^a = 2j(v^*)\delta a$$

$$\Delta F_W^a = 2\left(\delta\dot{A}_0^a v^* - c\delta a\right)$$

So it depends on 7 parameters.

ΔF_R^a is the spin. It is measured as for a particle. The result is represented as a vector S of \mathbb{R}^3 and its component S_z in the direction of the trajectory \vec{v} of the photon is the helicity. The 3 dimensional norm of S is normalized to $\sqrt{2\hbar}$.

The translational momentum is represented through \vec{v} . The momentum \vec{p} is linked to the energy and the frequency ν by : $\vec{p} = \frac{h\nu}{c^2}\vec{v}$. Which raises the question : what is ν ?

The Planck's law

It seems so natural to link the energy of a photon to a frequency, through the Planck's Law $E = h\nu$ that one forgets to ask for the origin of this frequency, so much that most often the frequency is defined by the energy. In order to try to explain this law, we have to scrutinize how, actually, the energy of a photon is measured.

Let be a EM field, observed in an area Ω , during a time interval $[0, T]$, without particles. The field is represented by a 2 form on M , valued in the Lie algebra \mathbb{R} , such that the norm $\|\mathcal{F}\|$ is integrable over Ω . The density of energy of the field is well defined by $\langle \mathcal{F}, \mathcal{F} \rangle$ and, using the results of 5.4.7 and 105 for each level of the energy $\int_{\Omega} \langle \mathcal{F}, \mathcal{F} \rangle \varpi_4 = Y$ there is a unique map \mathcal{F} .

The detection of a photon is done either when it interacts with a particle, or as a variation of the energy of the field. In the vacuum the photon appears as a singularity in the energy of the field, to which it is added. In the first chapter (Theorem 16) we have seen the requirements to scientifically accept a singularity, from a sample of points, over some time period. A discontinuity is acknowledged if the ratio $\frac{\rho_2(x)}{\rho(x)}$ is at least > 1 , where $\rho(x)$ is the frequency with which the value x is taken by the underlying quantity (here the energy density of the field $\mathcal{E} = \sum_{\alpha\beta} \langle \mathcal{F}_A^{\alpha\beta}, \mathcal{F}_{A\alpha\beta} \rangle$) and $\rho_2(x)$ is the frequency with which the value x is taken by the discrepancy (here $\delta\mathcal{E} = \sum_{\alpha\beta=0}^3 \langle \mathcal{F}_A^{\alpha\beta}, \Delta\dot{A}_{\alpha\beta}^a \rangle$). So this is not the absolute value of $\delta\mathcal{E}$ which matters, and not even its relative value $\delta\mathcal{E}/\mathcal{E}$, but the frequency with which this ratio is > 1 .

If the theorem 27 holds, then the field is decomposable in a Fourier series with frequency ν :

$$\mathcal{F}(\varphi_o(t, x)) = \sum_{n=-\infty}^{+\infty} (\exp 2i\pi t n \nu) \mathcal{F}(\varphi_o(0, x))$$

For a level of energy the frequency is fixed and the energy density \mathcal{E} is constant. So the quantity $\frac{\rho_2(x)}{\rho(x)}$ is proportional to the frequency ν . The higher the frequency, the greater the probability that a given discrepancy in the density of energy is acknowledged. Conversely, to a given $\delta\mathcal{E}$ is associated a minimum frequency. And actually it is quite impossible to detect a photon with fields of large wave lengths. In the EM field higher frequency means higher density of energy of the field, and then, because $\delta\mathcal{E} > \mathcal{E}$ photons with higher energy. This is the basic interpretation of the Planck's law : $E = h\nu$. The frequency ν is linked to the underlying field, for which the photon is a discontinuity. And, indeed, this is the only answer to the origin of the quantity ν itself : it must come from a field. Notice that this result is not related to the precision of the measure, or to the purity of the field. On the contrary, the probability to detect a singularity is greater for a perfectly monochromatic field.

Actually most of the experiments involve photons in direct interaction with particles. In these discontinuous processes the role of the underlying field is less obvious, however the detection of the photons, directly or indirectly (such as in the Black Body), relies on the measure of singularities in a continuous field, which is the motivation to keep the relation between energy and frequency.

Formally the reasoning can be extended to the other fields. But because they are either not well known, or difficult to measure, this kind of quantization has little interest, and bosons appear essentially through their direct interactions with particles.

Detection of Gravitons

All the previous results hold, at least formally, for the gravitational field. However the graviton, has never been observed, no more than gravitational waves. This just gives more interest to a theoretical approach, albeit speculative.

The effect on the underlying gravitational field is given by the scalar product :

$$\langle j(\mathcal{F}_G), j(\Delta G) \rangle = \frac{1}{2} \sum_{a=1}^3 j(\mathcal{F}_r^a)^t j(\Delta G^a) - j(\mathcal{F}_w^a)^t j(\Delta G^{a+3})$$

As seen in the theorem 105 , along the trajectory of the graviton $\Phi_V(t, a)$

$$j(\mathcal{F}_G^a)(\Phi_V(t, a)) = \Theta_r(t) j(\mathcal{F}_G^a)(\Phi_V(0, a))$$

The gravitational field does not vanish at infinity, and there is no reason why it could be decomposed in periodic plane waves, so the Planck's law does not hold any longer. However one can expect that $\Theta(t)$ vary very slowly. So the condition to detect a graviton is that

$$\langle j(\mathcal{F}_G), j(\Delta G) \rangle > \langle j(\mathcal{F}_G), j(\mathcal{F}_G) \rangle$$

which is a very strong condition. Moreover, if gravitons exist in flavors encompassing all dimensions a of the Lie algebra the total impact is the difference of two quantities. So these are reasons which can explain the non observation, so far, of both gravitational waves and gravitons.

If gravitons exist, the decomposition along the vector subspaces L_0, P_0 and

their dual leads to the spin $(L_0^* \otimes L_0)$ and 3 other tensors, which all involve ε_0 .

The spin of the boson would interfere only with the usual spatial gravity. The impact of the other components is less obvious. But the projection on $P_0^* \otimes P_0$:

$$[j(\Delta G)]_{P_0^* \otimes P_0} = \sum_{i=1}^3 [\Delta G_w^1]_0^i \varepsilon^i \cdot \varepsilon^0 \otimes \varepsilon_0 \cdot \varepsilon_1 + [\Delta G_w^2]_0^i \varepsilon^i \cdot \varepsilon^0 \otimes \varepsilon_0 \cdot \varepsilon_2 + [\Delta G_w^3]_0^i \varepsilon^i \cdot \varepsilon^0 \otimes \varepsilon_0 \cdot \varepsilon_3$$

would have a maximum impact on the foliation $\Omega_3(t)$, so practically it would cause a distortion of the measure of the time. As time is the physical quantity which is measured with the best precision, the detection of transversal gravitons should be easier, if they exist and the theory is right !

8.1.6 Spontaneous discontinuities

How can discontinuities appear spontaneously in the fields ? The models that we have studied show that the fields must satisfy differential equations at the equilibrium. The models represent systems of interacting fields and particles, which adjust to each other continuously : they are the equivalent of the reversible equilibrium of thermodynamics. To achieve this adjustment fields propagate, their speed is huge, but this propagation is not instantaneous. If some discontinuous process occurs, such as the action of an external field, or the introduction of new particles in the system, a new equilibrium is reached, the strength of the field \mathcal{F} takes a new value over Ω . But we can conceive that it is not always possible to go from one equilibrium to another by a continuous deformation of the field : discontinuities can be necessary to reach the adjustment. This is similar to what occurs with shock waves in fluids : when a strong variation of pressure happens at some point, the equilibrium cannot any longer be achieved by the smooth propagation of reversible adjustments, discontinuities appear along a front wave which propagates.

A similar process occurs in a collision between particles : each particle generates its own field, and an adjustment must be done when the particles approach each other. Depending on the velocities of the particles, this adjustment can be done smoothly or not.

The conditions for the apparition of discontinuities in systems are an important domain of research in Physics, which involves Topological Algebra (to which the Chern-Weil theory can be attached), in condensed matter (phases transitions) as well as systems of particles.

8.2 DISCONTINUITIES WITH PARTICLES AND FIELDS

Potentials cannot appear explicitly in Lagrangians because of their special behavior in a change of gauge. This obstruction does not hold any longer for bosons, with our definition. However there is another one : when they are considered as particles, their relativist momentum should be evaluated along their trajectories, but there is no measure on these curves, at least if we assume that they are null curves : $\langle V, V \rangle = 0$. So the model with a single particle cannot be extended to bosons. One can consider bosonic fields, similar to the matter fields. However the physical meaning of such models is questionable. Another issue is that bosons interact with particles but not as the fields do : the covariant derivative mechanism does not hold anymore. Actually bosons interact with particles during direct contact, either in collisions, or as the result of a collision or the disintegration of a particle. In this sections collisions will encompass all the discontinuous processes during which particles or bosons exist together at one point. It can be the consequence of the encounter of two objects, which exist before the collision, the splitting of the object in different parts or the emission of a boson by a particle.

There are several, complementary, methods to deal with discontinuous processes : the use of general conservations law for momentum and energy, the path integrals and the representation of interacting systems with their prolongation in Fock's space.

8.2.1 General rules in collisions

The cases vary but the solutions which are implemented follow some basic rules.

The first is the conservation of energy and momentum, which holds in discontinuous processes. They have both a clear definition for particles, with the relativist momentum, from which the translational and rotational momenta can be deduced. For bosons the energy is linked to the energy of the underlying field (which is always present, at least by the field for which the particle is the source) and, for photons to its frequency. In QTF their momentum is reduced to their translational momentum, measured by the spatial 3 components as it is assumed that they travel at the speed of light. The conservation is stated pointwise, where the collision occurs, and it should be met in any gauge.

There are some rules about the conservation of charge. They are well defined for particles, and for bosons through the currents. The rules for the conservation of charges depend on the fields which are involved.

When only the gravitational field is present, it sums up to the conservation of the mass at rest, which is a basic tenet of Classical Mechanics, generalized to the Relativist context.

When only the EM field is involved, the charge is incorporated in the spinor, and there is a clear law for the conservation of charge expressed in units. Moreover the bosons have no charge.

When the other fields are involved, the situation is more complicated. There are rules, more empirical than based on strict principles, such as the conservation of the sum of weak isospin or of the number of baryons.

A discontinuous process occurs between periods during which an equilibrium is assumed to have been reached. So one considers transitions from one state of equilibrium to another, which is represented as a **scattering** process in experiments using beams of particles, with a population of incoming particles in “in” states, and outgoing particles in “out” states. In an equilibrium the Principle of Least Action applies, and the states of the particles follow equations similar to one of the two models which have been studied in the previous chapter. Moreover the states of particles can be represented, either as matter fields for a population of particles, or for a single particle as maps belonging to Fréchet spaces. The states of the system “in” and “out” can then be computed as if they were equilibrium and one can consider the possible outcomes.

The states of the “in” particles can be represented by maps :

$$t \in]-\infty, 0[: \psi(t) = \vartheta(\sigma(t), \varkappa(t)) \psi_{0in}$$

and for the “out” particles by :

$$t \in]0, +\infty[: \psi(t) = \vartheta(\sigma(t), \varkappa(t)) \psi_{out}$$

where $(\sigma(t), \varkappa(t))$ are computed as in a continuous process.

The conservation of energy and momentum is met at each location, for all the particles or bosons involved. And the previous rules are implemented locally. If a photon strikes a particle, whose fundamental state does not change, the variation will be imparted to the boson (this is the Compton effect). Conversely if the state of the particle changes, following a collision with another particle for instance, the new equilibrium is reached with the emission of a boson (Black Body radiation). There is always an underlying field, at least the one created by the particle.

8.2.2 Path integrals

If we have different possible outcomes, the question arises to find which one will occur. This is the main purpose of the path integral theory.

As many others in Quantum Physics, the idea of path integral comes from Statistical Mechanics, and was proposed notably by Wiener.

If the evolution of the system meets the criteria of the Theorem 26 (the variables are maps depending on time and valued in a normed vector space and the process is determinist) there is an operator $\Theta(t)$ such that : $X(t) = \Theta(t) X(0)$. When in addition the variables $X(t)$ and $X(t + \theta)$ represents the same state, $\Theta(t) = \exp t\Theta$ with a constant operator. The exponential of an operator on a Banach space is a well known object in Mathematics (Maths.1025), so the law of evolution is simple when $\Theta(t)$ is constant which, has we have seen several times, requires fairly strong conditions. However, because discontinuities

are isolated points, at least at an elementary level, we can consider that between the transitions points Θ is constant. Then we have a succession of laws :

$$t \in [t_p, t_{p+1}[: X(t) = (\exp t\Theta_p) X(t_p)$$

and :

$$X(t) = (\exp(t - t_p)\Theta_p) (\exp(t_p - t_{p-1})\Theta_{p-1}) \dots (\exp t_1\Theta_p) X(0)$$

which are usually represented, starting from the derivative.

This is a generalization of the mathematical method to express the solution of the differential equation in \mathbb{R}^m : $\frac{dX}{dt} = \Theta(t) X(t)$:

$$X(t) = \lim_{n \rightarrow \infty} \left(\prod_{p=0}^n \exp(t_{p+1} - t_p)\Theta(t_p) \right) X(0) \text{ (Maths.2570)}$$

The Θ_p and the intermediary transition points are not known, but if we can attribute a probability to each transition, then we have a stochastic process (see Maths.11.4.4). The usual assumption is that the transitions are independent events, and the increment $(\Theta_{p+1} - \Theta_p)$ follow a fixed normal distribution law (a Wiener process). In this scheme all possible paths must be considered.

In QM the starting point is the Schrödinger equation, $i\hbar \frac{d\psi}{dt} = H\psi$, which has a similar meaning. However in a conventional QM interpretation there is no definite path (only the initial and the final states are considered) and furthermore, because of the singular role given to t , it seemed not compatible with Relativity. Dirac proposed the use of the lagrangian, and Feynman provided a full theory of path integrals, which is still one of the essential tools of QTF. The fundamental ideas, as expressed by Feynman, are that :

- to any physical event is associated a complex scalar ϕ , called an amplitude of probability,
- a physical process is represented by a path, in which several events occur successively
- the amplitude of probability of a process along a path is the sum of the amplitude of probability of each event
- the probability of occurrence of a process is the square of the module of the sum of the amplitudes of probability along any path which starts and ends as the initial and final states of the process (at least if there is no observation of any intermediate event).

The amplitude of probability of a given process is given by : $e^{\frac{i}{\hbar}S[z]}$ where $S[z]$ is the action, computed with the lagrangian :

$S[z] = \int_A^B L(z^i, z_{\alpha}^i \dots z_{\alpha_1 \dots \alpha_r}^i) dm$ evaluated from the r-jet extension of z , with a classical lagrangian. The total amplitude of probability to go from a state A to a state B is $\phi = \int e^{\frac{i}{\hbar}S[z]} Dz$ where Dz means that all the imaginable processes must be considered. Then the probability to go from A to B is $|\phi|^2$. So each path contributes equally to the amplitude of probability, but the probability itself is the square of the module of the second integral.

The QM wave function follows : $\psi(x, t) = \int_{-\infty}^{+\infty} \phi(x, t; \xi, \tau) \psi(\xi, \tau) d\xi d\tau$

If a process can be divided as : $A \rightarrow B \rightarrow C$ then

$\phi(A, C) = \phi(A, B) \phi(B, C)$ which is actually the idea of dividing the path in small time intervals.

It can be shown that in the classical limit ($\hbar \rightarrow 0$) and certain conditions the path integral is equivalent to the Principle of Least Action. With simplifications most of the usual results of QM can be retrieved.

Even if the literature emphasizes simple examples (such as the trajectory of a single particle), the path integral is used, with many variants, mainly to address the case of discontinuous processes in QTF, as this is the only general method known. It leads then to consider the multiple possibilities of collisions, emissions,... involving different kinds of particles or bosons, in paths called Feynman's diagrams.

The quantities which are involved are either force fields (gravitation is not considered), fermionic fields or bosonic fields. In the latter two cases a trajectory is computed as a path, but this is a bit awkward because if bosons follow a null curve the lagrangian is not defined as we have seen.

It is clear that this formalism is grounded in the philosophical point of view that all physical processes are discreet and random. One can subscribe or not to this vision, but it leads to some strange explanations. For instance all the paths must be considered, even when they involve unphysical behaviors for the particles (the virtual particle are not supposed to follow the usual laws of physics). An explanation which is not necessary : we have eventually a variational calculus, so the r-jets, in which the derivatives are independent variables, is the natural mathematical framework and we must consider all possible values for the variables, independently of their formal relations.

Beyond the simplest case, where it has little added value, the computation of path integrals is a dreadful mathematical endeavour. This is done essentially in a perturbative approach, where the lagrangian is simplified as we have done previously, so as to come back to quadratic expressions. The results are then developed in series of some scale constant. However it is full of mathematical inconsistencies, such as divergent integrals. The theory of path integrals is then essentially dedicated to find new computational methods or tricks, without few or no physical justification : renormalization, ghosts fields, Gladston bosons, Wick's rotation, BRST,...

8.2.3 Interacting micro-systems

The continuous model of type 1 was inspired by Fluid Mechanics, and the natural extension is Gas Mechanics, where a great number of particles interact together. Fortunately the framework used here fits well with the idea of interacting microsystems.

The single particle model does not provide the value of the fields out of the trajectory. However it shows that the interactions force fields / particles, in spite of being represented by maps which have supports of different dimension (a 4 manifold for force fields and a curve for fermions or bosons) lead to relations in which only the values in one point matter : $\psi(m), \hat{A}(m) = \sum_{\alpha=0}^3 \dot{A}_\alpha V^\alpha, \hat{G}(m) = \sum_{\alpha=0}^3 G_\alpha V^\alpha$. All these variables are geometric, so actually the location m does not matter : their value can be expressed in any system

of frames or coordinates. Moreover the state of the particle or the boson incorporates all the information about their motion, which is then defined without the need to exhibit components of vectors in a frame.

We can then consider micro systems, comprising one particle, the fields and possibly one boson, all located at the same point, and represented by the variables which enter the lagrangian : $\psi, \mathcal{F}_A, \mathcal{F}_G, \Delta\dot{A}, \Delta G$ and P in the GR context, albeit it is a bit of an over kill in these kinds of problems, the SR approximation suffices ¹. These variables are seen as force, fermionic or bosonic fields but, because their location does not matter, this is the value in the associated vector spaces which is considered (which is equivalent to take the gauge of the observer). If the conditions of the theorem 28 are met, then the state of a microsystem is represented in a Hilbert space H by a vector, which is the direct product of vectors representing each variable.

If we have a system comprised of N such microsystems, and if the bosons and fermions are of the same type, they have the same behavior, and are indistinguishable : we have a homogeneous system and we can apply the theorems 31 and 33. The interactions between the micro systems lead to the quantization of the states. This is done in several steps.

1. The states of the microsystems (encompassing all the variables

$\psi, \mathcal{F}_A, \mathcal{F}_G, \Delta\dot{A}, \Delta G$) are associated to a Hilbert space H , and the states of the system are associated to the tensorial product $\otimes_{n=1}^N H$ of the Hilbert space H associated to each microsystem. An equilibrium of the system corresponds to a vector subspace \mathbf{h} of $\otimes_{n=1}^N H$ which is defined by :

i) a class of conjugacy $\mathfrak{S}(\lambda)$ of the group of permutations $\mathfrak{S}(N)$, defined itself by a decomposition of N in p parts :

$$\lambda = \{0 \leq n_p \leq \dots \leq n_1 \leq N, n_1 + \dots + n_p = N\}.$$

ii) p distinct vectors $(\tilde{\varepsilon}_j)_{j=1}^p$ of a Hermitian basis of H which together define a vector space H_J

And \mathbf{h} is then either $\odot_{n_1} H_J \otimes \odot_{n_2} H_J \dots \otimes \odot_{n_p} H_J$ or $\wedge_{n_1} H_J \otimes \wedge_{n_2} H_J \dots \otimes \wedge_{n_p} H_J$

The state Ψ of the system is then : $\Psi = \sum_{(i_1 \dots i_n)} \Psi^{i_1 \dots i_n} \tilde{\varepsilon}_{i_1} \otimes \dots \otimes \tilde{\varepsilon}_{i_n}$ with an antisymmetric or a symmetric tensor.

2. We have global variables, which can be taken equivalently as the number of particles, or their charge, and the energy of the system. For each value of the global variables the state Ψ of the system belongs to one of the irreducible representations. The class of conjugacy λ and the vectors $(\tilde{\varepsilon}_j)_{j=1}^p$ are fixed.

3. At the level of each microsystem, each vector $\tilde{\varepsilon}_j \in H$ represents a definite state of a micro system, and the value of each variable of the micro-system is quantized. In a probabilist interpretation one can say that there are $(n_i)_{i=1}^p$ microsystems in the state $\tilde{\varepsilon}_{j_i}$. But of course one cannot say with certainty what is the state of a given microsystem.

The quantization of each microsystem means that the vector ϕ representing its state in H belongs to a finite dimensional vector space :

¹Notice that P is a vector (ε_i) so a geometric quantity, and is not submitted a priori to a constraint because g is defined from P .

$\phi = \sum_{(i_1, \dots, i_q)} \phi^{i_1 \dots i_q} |e_{i_1}, e_{i_2} \dots e_{i_q}\rangle$ where the vectors $|e_{i_1}, e_{i_2} \dots e_{i_q}\rangle$ correspond to the $\tilde{\varepsilon}_j$

The spin of the particle corresponds to one of the vectors e_j of the basis. The spin of a particle is always invariant by $Spin(3)$, the action of s and $-s$ give opposite results. But if the spin number j is an integer then the particle has a specific, physical symmetry, and its spin is invariant by $SO(3)$. We have seen that the force fields and the bosons have a spin number which is an integer. So we have a distinct behavior of the system, if its particles have a spin number which is an integer or half an integer. This property must be reflected in the states of the system.

If j is half an integer the representation of the system is by antisymmetric tensors to account for the antisymmetry by $Spin(3)$. As a consequence in each vector space $\wedge_n H_J$ the components of the tensors, expressed in any basis, which correspond to the diagonal are null :

$$\psi^{i_1 \dots i_n} = 0 \text{ for } i_1 = i_2 = \dots = i_n$$

The micro systems belonging to the same $\wedge_n H_J$ must be in different states. This is the Pauli's exclusion principle.

The particles whose spin number is half an integer are called fermions and are said to follow the Fermi-Dirac statistic.

The particles whose spin number is an integer are called bosons and are said to follow the Bose-Einstein statistic.

So the denominations fermions / bosons are here different from that we have used so far. All elementary particles are fermions, all discontinuities of the fields are bosons, but composite particles or atoms can be bosons if the spin number is an integer.

Notice that the exclusion principle does not apply to all the micro-systems. In a system there are usually different sets of microsystems, which corresponds to different subspaces $\wedge_n H_J$ and therefore micro-systems belonging to different subspaces can have the same spin, however each of these subspaces is distinguished by other global variables, such the energy (for instance the electrons are organized in bands of valence in an atom).

The physical manifestation of these mechanisms have been experimentally checked. But the definition of a homogeneous system must be met, which restricts them practically to systems of identical particles enclosed in a limited area.

8.2.4 Fock Spaces

In the model of interacting systems above, the number of particles was assumed to be constant. It is possible to suppress this limitation by using an extension of the tensorial product of Hilbert spaces.

The space of representation of the system is then the tensorial algebra, called a Fock space : $\mathcal{F} = \bigoplus_{k=0}^{\infty} (\otimes_k H)$ where H stands for the Hilbert space of the microsystems (Maths.1209). k can be 0 so scalars can be vectors of the Fock spaces.

A vector Ψ of $\mathcal{F}_n = \bigoplus_{k=0}^n (\otimes_k H)$ is given by $n + 1$ tensors :

$$(\psi^m, \psi^m \in \otimes^m H, m = 0 \dots n)$$

The “ground state” is the vector $(1, 0, 0, \dots)$ in the algebra.

Any operator on the Hilbert spaces can be extended to a linear continuous operator on the Fock space.

For each Fock space $\oplus_{k=1}^{\infty} (\otimes_k H)$ there is a number operator N , whose, dense, domain is :

$$D(N) = \left\{ \psi^m \in \otimes_m H, \sum_{k \geq 0} m^2 \|\psi^m\|^2 < \infty \right\}$$

$$N(\Psi) = (0, \psi^1, 2\psi^2, \dots, m\psi^m \dots)$$

N is self adjoint.

The annihilation operator cuts a tensor at its beginning :

$$a_m : H \rightarrow \mathcal{L}(\otimes_m H; \otimes_{m-1} H) ::$$

$$a_m(\psi)(\psi_1 \otimes \psi_2 \dots \otimes \psi_m) = \frac{1}{\sqrt{m}} \langle \psi, \psi_1 \rangle_H \psi_2 \otimes \psi_3 \dots \otimes \psi_m$$

The creation operator adds a vector to a tensor at its beginning :

$$a_m^* : H \rightarrow \mathcal{L}(\otimes_m H; \otimes_{m+1} H) ::$$

$$a_m^*(\psi)(\psi_1 \otimes \psi_2 \dots \otimes \psi_m) = \sqrt{m+1} \psi \otimes \psi_1 \otimes \psi_2 \otimes \psi_3 \dots \otimes \psi_m$$

a_m^* is the adjoint of a_m and a_m, a_m^* can be extended to the Fock space as a, a^* .

The physical meaning of these operators is clear from their names. They are the main tools to represent the crucial phenomenon in a discontinuous process, that is the variation of the number of particles.

The spaces of symmetric (called the Bose-Fock space) and antisymmetric (called the Fermi-Fock space) tensors in a Fock space have special properties. They are closed vector subspaces, so are themselves Hilbert spaces, with an adjusted scalar product. Any tensor of the Fock space can be projected on the Bose subspace (by P_+) or the Fermi space (by P_-) by symmetrization and antisymmetrization respectively, and P_+, P_- are orthogonal. The operator $\exp itN$ leaves both subspaces invariant. Any self-adjoint operator on the underlying Hilbert space has an essentially self adjoint prolongation on these subspaces (called its second quantification). However the creation and annihilation operators have extensions with specific commutation rules :

Canonical commutation rules (CCR) in the Bose space:

$$[a_+(u), a_+(v)] = [a_+^*(u), a_+^*(v)] = 0$$

$$[a_+(u), a_+^*(v)] = \langle u, v \rangle 1$$

Canonical anticommutation rules (CAR) in the Fermi space :

$$\{a_+(u), a_+(v)\} = \{a_-^*(u), a_-^*(v)\} = 0$$

$$\{a_+(u), a_+^*(v)\} = \langle u, v \rangle 1$$

where

$$[X, Y] = X \circ Y - Y \circ X$$

$$\{X, Y\} = X \circ Y + Y \circ X$$

These differences have important mathematical consequences. In the Fermi space the operators a_-, a_-^* have bounded (continuous) extensions. Any configuration of particles can be generated by the product of creation operators acting on the ground state. We have nothing equivalent for the bosons.

We will not pursue on these topics, which are exposed in many books, and requires more mathematical concepts.

Chapter 9

CONCLUSION

At the end of this book I hope that the reader has a better understanding of how Mathematical tools such as Group Representations, Clifford Algebras, Fiber Bundles, Connections can be used in Physics. They can be a bit abstract, but actually they are well fitted to address the issues of modern Physics, and quite efficient. I hope also to have brought some clarification on Quantum Mechanics, Relativity and gauge theories, as well as on concepts, such as the duality between particles and fields.

1. In the Second Chapter it has been proven that most of the axioms of QM come from the way models are expressed in Physics, and the following chapters have shown how the theorems which have been proven can be used, and support the usual computations. The theorems state precise guidelines and requirements for their validity, and these requirements, albeit expressed as Mathematical conditions, lead to a deeper investigation of the physical meaning of the quantities which are used. For a property, the fact to be geometric is not a simple formality : it means that this is an entity which exists beyond the measures which can be made, and that these measures vary according to precise rules. The role of the observer in the process of measurement is then clearly specified. There has been few examples of the use of observables, whose role is more central in models representing practical experiments, but their meaning should be clear.

2. Relativity, and particularly General Relativity, which is often seen as a difficult topic, can be understood if we accept to start from the beginning, from Geometry, the particularities of our Universe and if we accept to give up schemes and representations which have become too familiar. With the formalism of fiber bundles it is then easy to address very general topics without losing the mathematical rigor.

3. Clifford algebras are not new, but they appear really useful when one accepts fully the riches of their structure, without resorting to hybrid concepts such as quasi or axial vectors. With spinors the concept of matter fields becomes clear. In my opinion they are the only way to represent in a consistent and efficient way the motion and the kinematics properties of material bodies in the

GR context. So Spinors should be useful in Astrophysics, where gravitation is the only force involved and GR cannot be dismissed.

4. The use of connections to represent the force fields has become a standard in gauge theories. The strict usage of fiber bundles and spinors enables to put the gravitational field in the same framework, and it appears clearly that the traditional method based on the metric and the Levi-Civita connection imposes useless complications and miss some features which can be physically important, such that the decomposition in transversal and spatial components.

5. The two models presented were essentially an example of how the theory of Lagrangians, with its difficulties, can be used practically. They have given the opportunity to introduce in a consistent manner the concepts of energy-momentum tensor and currents. And we have given a strong mathematical backing to the functional derivatives calculus, based of an original theory of distributions on vector bundles introduced in the Mathematics Book.

6. The idea of bosons as discontinuities in the fields is more speculative. But it seems necessary to complement the concept of matter fields, which is quite clear, with an interpretation of the corpuscular nature of the fields. The presentation leaves some gaps, which are in part due to our limited knowledge of the propagation of weak and strong interactions, and to the imprecision in the definition of mass and energy, beyond the Higgs mechanism. Anyway the differences between continuous and discontinuous processes should be clear.

There are some new results in this book : QM, deformable solids in RG, spinors, bosons. They are worth to be extended, by filling the gaps, or simply using the methods which have been introduced. For instance many other theorems could be proven in QM, a true Mechanics of deformable solids could be built, with the addition of Thermodynamics concepts, the representation of bosons could be more firmly grounded by the consideration of a “state of the vacuum”. But from my point of view the one which is worth of the most efforts is gravitation. This is the most common and weakest of all force fields, but we are still unable to use it or understand it properly. The representation of the gravitational field by connections on one hand, and of the gravitational charges by spinors on the other hand, shows striking similarities with the EM field : indeed they are the only fields which have an infinite range, the EM charge can be incorporated in the gravitational charge, and their propagation equations are similar. This similitude has been remarked by many authors, Heaviside, Negut, Jefimenko, Tajmar, de Matos,...and it has been developed in a full Theory, which has sometimes be opposed to GR. We find here that these similitudes exist in the frame of a GR theory which allows for a more general connection and the use of the Riemann tensor, so it seems more promising to explore this avenue than to fight against GR. As we have seen the gravitational field shows in all its aspects two components. The “magnetic” component can be assimilated to the usual gravity : this is the one which acts in the 3 dimensional space. The “electric” component acts in the time dimension, and it seems logical to give it a cosmological interpretation : it would be the engine which moves matter on its world line. Both components have opposite effects, and there is no compelling reason that

it should always be attractive. The representation of the gravitational field in the bundle $P_G [T_1 Spin(3, 1)^* \otimes T_1 Spin(3, 1), \mathbf{Ad}^{-1} \times \mathbf{Ad}]$ enables to explore more efficiently the structure of the field than the usual Petrov-Pirani-Penrose classification based on the Weyl's tensor.

QM and Relativity have deeply transformed the way we do Physics. I am not the first to state the obvious. I hope that the reader understands better what has changed, and why.

We were used to an eternal, flat, infinite Universe (an idea which is after all not so obvious), with Relativity we had to accept that we could represent the Universe as a four dimensional, curved, structure, which integrates the time. Beyond the change of mathematical formalism, Relativity has also put limits to our capability to know the Universe. We are allowed to model it as we want, with an infinite extension, in space and time, but the only Universe that is accessible to our measures and experiments is specific to each observer : we have as many windows on the Reality that there are observers. We can dream the whole world, we can put in our models variables which are related to the past or the future, as if they were there, but the world that I can perceive is the world that I see from my window, and my neighbor uses another window. I can imagine what is beyond my window, but to get a comprehensive picture I need to patch together different visions.

With QM we have realized that we can model the reality, whatever the scale, with mathematical objects, but these objects exist only in the abstract world of Mathematics, they are some idealization that we use because they are efficient in our computations, but we can access reality only with cruder objects, finite samples and statistic estimations. The discrepancy between the measures, necessarily circumstantial and probabilist, and the real world does not mean that the real world is discreet and proceeds according to random behaviors, only that we have to acknowledge the difference between a representation and the reality. And conversely it does not preclude the use of the models, as long as we are aware of their specific place : it is not because we cannot measure simultaneously location and speed that their concepts are void.

Contrary to many, I am a realist, I believe that there is a unique real world outside, it can be understood, it is not ruled by strange and erratic behaviors, but I fully acknowledge that my world, the one in which I live, measure and experiment, is specific. This is one the multiple windows which are offered. Modern Physics, in a mischievous turn, has imposed the need to reintroduce the individual in Science, in the guise of the observer, and the discrepancy between imagination, which enables to see the whole as if it was there, and the limited possibility to keep it in check. The genuine feature of the human brain is that it can conceive of things that do not exist, that will never occur as we dreamed them. This is precious and Science would be impossible without it. To impart to reality our limitations or to limit our ambitions to what we can check are equally wrong. Actually the only way for a Scientist to keep his sanity in front of all the possible explanations which are provided is that to remember that there is one world : the one in which he lives. And that Science can progress

only through a shared knowledge.

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Chapter 10

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Appendix A

ANNEX

A.1 CLIFFORD ALGEBRAS

This annex gives proofs of some results presented in the core of the paper.

A.1.1 Products in the Clifford algebra

Many results are consequences of the computation of products in the Clifford algebra. The computations are straightforward but the results precious. In the following $\langle \varepsilon_0, \varepsilon_0 \rangle = -1$ with the signature (3,1) and +1 with the signature (1,3). The operator j is reminded in the Formulas at the end of this Annex.

Product $v(r, w) \cdot v(r', w')$

$$v(r, w) = \frac{1}{2} (w^1 \varepsilon_0 \cdot \varepsilon_1 + w^2 \varepsilon_0 \cdot \varepsilon_2 + w^3 \varepsilon_0 \cdot \varepsilon_3 + r^3 \varepsilon_2 \cdot \varepsilon_1 + r^2 \varepsilon_1 \cdot \varepsilon_3 + r^1 \varepsilon_3 \cdot \varepsilon_2)$$
$$v(r', w') = \frac{1}{2} (w'^1 \varepsilon_0 \cdot \varepsilon_1 + w'^2 \varepsilon_0 \cdot \varepsilon_2 + w'^3 \varepsilon_0 \cdot \varepsilon_3 + r'^3 \varepsilon_2 \cdot \varepsilon_1 + r'^2 \varepsilon_1 \cdot \varepsilon_3 + r'^1 \varepsilon_3 \cdot \varepsilon_2)$$

With signature (3,1) :

$$v(r, w) \cdot v(r', w') = \frac{1}{4} (w^t w' - r^t r') + \frac{1}{2} v(j(r) r' - j(w) w', j(w) r' + j(r) w') - \frac{1}{4} (w^t r' + r^t w') \varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3$$

From there the bracket on the Lie algebra :

$$[v(r, w), v(r', w')] = v(r, w) \cdot v(r', w') - v(r', w') \cdot v(r, w)$$

$$[v(r, w), v(r', w')] = v(j(r) r' - j(w) w', j(w) r' + j(r) w') \quad (\text{A.1})$$

With signature (1,3) :

$$v(r, w) \cdot v(r', w') = \frac{1}{4} (w^t w' - r^t r') - \frac{1}{2} v(-j(r) r' + j(w) w', j(w) r' + j(r) w') - \frac{1}{4} (w^t r' + r^t w') \varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3$$

From there the bracket on the Lie algebra :

$$[v(r, w), v(r', w')] = -v(j(r)r' - j(w)w', j(w)r' + j(r)w') \quad (\text{A.2})$$

In both signatures the basis of the Lie algebra is denoted :

$$\begin{aligned} \vec{\kappa}_1 &= \frac{1}{2}\varepsilon_3 \cdot \varepsilon_2, \\ \vec{\kappa}_2 &= \frac{1}{2}\varepsilon_1 \cdot \varepsilon_3, \\ \vec{\kappa}_3 &= \frac{1}{2}\varepsilon_2 \cdot \varepsilon_1, \\ \vec{\kappa}_4 &= \frac{1}{2}\varepsilon_0 \cdot \varepsilon_1, \\ \vec{\kappa}_5 &= \frac{1}{2}\varepsilon_0 \cdot \varepsilon_2, \\ \vec{\kappa}_6 &= \frac{1}{2}\varepsilon_0 \cdot \varepsilon_3 \\ a, b, c &= 1, 2, 3 \\ [\vec{\kappa}_a, \vec{\kappa}_b] &= \epsilon(c, a, b) \vec{\kappa}_c \\ [\vec{\kappa}_a, \vec{\kappa}_{3+b}] &= \epsilon(c, a, b) \vec{\kappa}_{3+c} \\ [\vec{\kappa}_{3+a}, \vec{\kappa}_{3+b}] &= \epsilon(c, a, b) \vec{\kappa}_{3+c} \end{aligned}$$

Product on $Spin(3, 1)$

Because they belong to $Cl_0(3, 1)$ the elements of $Spin(3, 1)$ can be written :

$$s = a + \frac{1}{2}(w^1\varepsilon_0 \cdot \varepsilon_1 + w^2\varepsilon_0 \cdot \varepsilon_2 + w^3\varepsilon_0 \cdot \varepsilon_3 + r^3\varepsilon_2 \cdot \varepsilon_1 + r^2\varepsilon_1 \cdot \varepsilon_3 + r^1\varepsilon_3 \cdot \varepsilon_2) + b\varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3$$

where $a, (w^j, r^j)_{j=1}^3, b$ are real scalar which are related. That we will write with

$$\varepsilon_5 = \varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3 \quad (\text{A.3})$$

$$s = a + v(r, w) + b\varepsilon_5 \quad (\text{A.4})$$

And similarly in $Cl(1, 3)$

$$s = a + v(r, w) + b\varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3$$

The product of two elements of the spin group expressed as :

$$s = a + v(r, w) + b\varepsilon_5$$

$$s' = a' + v(r', w') + b'\varepsilon_5$$

can be computed with the previous formulas.

i) With signature (3,1)

$$v(r, w)\varepsilon_5 = v(r, -w)$$

$$\varepsilon_5 v(r', w') = v(r', -w')$$

$$\varepsilon_5 \cdot \varepsilon_5 = -1$$

$$s \cdot s' = a'' + v(r'', w'') + b''\varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3$$

with :

$$a'' = aa' - b'b + \frac{1}{4}(w^t w' - r^t r')$$

$$b'' = ab' + ba' - \frac{1}{4}(w^t r' + r^t w')$$

$$r'' = \frac{1}{2}(j(r)r' - j(w)w') + a'r + ar' - b'w - bw'$$

$$w'' = \frac{1}{2}(j(w)r' + j(r)w') + a'w + aw' + b'r + br'$$

So we have in particular :

$$(a + v(0, w)) \cdot (a' + v(0, w')) = aa' + \frac{1}{4}w^t w' + v\left(-\frac{1}{2}(j(w)w', a'w + aw')\right)$$

$$(a + v(r, 0)) \cdot (a' + v(r', 0)) = aa' - \frac{1}{4}r^t r' + v\left(\frac{1}{2}j(r)r' + (a'r + ar'), 0\right)$$

$$(a_w + v(0, w)) \cdot (a_r + v(r, 0)) = a_w a_r + v(a_w r, a_r w) - \frac{1}{2}(w^t r) \varepsilon_5$$

ii) With signature (1,3)

$$s = a + v(r, w) + b\varepsilon_5$$

$$s' = a' + v(r', w') + b'\varepsilon_5$$

$$v(r, w) \cdot \varepsilon_5 = v(w, r)$$

$$\varepsilon_5 v(r', w') = v(w', r')$$

$$\varepsilon_5 \cdot \varepsilon_5 = -1$$

$$s \cdot s' = a'' + v(r'', w'') + b''\varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3$$

with :

$$a'' = aa' - b'b + \frac{1}{4}(w^t w' - r^t r')$$

$$b'' = ab' + ba' - \frac{1}{4}(w^t r' + r^t w')$$

$$r'' = \frac{1}{2}(j(r)r' - j(w)w') + a'r + ar' + b'w + bw'$$

$$w'' = -\frac{1}{2}(j(w)r' + j(r)w') + a'w + aw' + b'r + br'$$

A.1.2 Characterization of the elements of the Spin group

Inverse

The elements of $Spin(3, 1)$ are the product of an even number of vectors of norm ± 1 . Consequently we have :

$$s \cdot s^t = (v_1 \cdot \dots \cdot v_{2p}) \cdot (v_{2p} \cdot \dots \cdot v_1) = 1$$

The transposition is an involution on the Clifford algebra, thus :

$$(a + v(r, w) + b\varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3) \cdot (a + v(r, w)^t + b\varepsilon_3 \cdot \varepsilon_2 \cdot \varepsilon_1 \cdot \varepsilon_0) = 1$$

$$(a + v(r, w) + b\varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3) \cdot (a - v(r, w) + b\varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3) = 1$$

$$\Leftrightarrow (a + v(r, w) + b\varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3)^{-1} = (a - v(r, w) + b\varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3)$$

and we have the same result in $Cl(1, 3)$

$$(a + v(r, w) + b\varepsilon_5)^{-1} = a - v(r, w) + b\varepsilon_5 \quad (\text{A.5})$$

Relation between a, b, r, w

By a straightforward computation this identity gives the following relation between a, b, r, w :

1. With signature (3,1)

$$(a + v(r, w) + b\varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3) \cdot (a - v(r, w) + b\varepsilon_3 \cdot \varepsilon_2 \cdot \varepsilon_1 \cdot \varepsilon_0) = 1$$

$$= a'' + v(r'', w'') + b''\varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3$$

with :

$$a'' = a^2 - b^2 + \frac{1}{4}(-w^t w + r^t r) = 1$$

$$b'' = ab + ba - \frac{1}{4}(-w^t r - r^t w) = 0$$

$$r'' = \frac{1}{2}(-j(r)r + j(w)w) + ar - ar - bw + bw = 0$$

$$\begin{aligned}
w'' &= \frac{1}{2}(-j(w)r - j(r)w) + aw - aw + br - br = 0 \\
a^2 - b^2 &= 1 + \frac{1}{4}(w^t w - r^t r) \\
\text{So, for any element : } a + v(r, w) + b\varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3 \\
&\text{we have :}
\end{aligned}$$

$$a^2 - b^2 = 1 + \frac{1}{4}(w^t w - r^t r) \quad (\text{A.6})$$

$$ab = -\frac{1}{4}r^t w \quad (\text{A.7})$$

and we keep only 6 free parameters. a, b are defined from r, w , up to sign, with the conditions:

$$\begin{aligned}
\text{i) } r^t w \neq 0 : b &= -\frac{1}{4a}r^t w \\
a^2 &= \frac{1}{2} \left(\left(1 + \frac{1}{4}(w^t w - r^t r)\right) + \sqrt{\left(1 + \frac{1}{4}(w^t w - r^t r)\right)^2 + \frac{1}{4}(r^t w)^2} \right) \\
\text{ii) } r^t w = 0 : \\
(w^t w - r^t r) \geq -4 : a &= \epsilon \sqrt{1 + \frac{1}{4}(w^t w - r^t r)}; b = 0 \\
(w^t w - r^t r) \leq -4 : b &= \epsilon \sqrt{-\left(1 + \frac{1}{4}(w^t w - r^t r)\right)}; a = 0 \\
\text{So :} \\
\text{if } r = 0 \text{ then : } s &= \epsilon \sqrt{1 + \frac{1}{4}w^t w} + v(0, w) \\
\text{if } w = 0 \text{ then} \\
r^t r \leq 4 : s &= \epsilon \sqrt{1 - \frac{1}{4}r^t r} + v(r, 0) \\
r^t r \geq 4 : s &= v(r, 0) + \epsilon \sqrt{\frac{1}{4}r^t r - 1} \varepsilon_5
\end{aligned}$$

2. With signature (1,3)

$$\begin{aligned}
(a - v(r, w) + b\varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3) \cdot (a + v(r, w) + b\varepsilon_3 \cdot \varepsilon_2 \cdot \varepsilon_1 \cdot \varepsilon_0) &= 1 \\
&= a'' + v(r'', w'') + b'' \varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3
\end{aligned}$$

with :

$$\begin{aligned}
r'' &= \frac{1}{2}(-j(r)r + j(w)w) + ar - ar + bw - bw = 0 \\
w'' &= -\frac{1}{2}(-j(w)r - j(r)w) + aw - aw + br - br' \\
a'' &= a^2 - b^2 + \frac{1}{4}(-w^t w + r^t r) \\
b'' &= ab + ba - \frac{1}{4}(-w^t r - r^t w)
\end{aligned}$$

we get the same relations.

A.1.3 Homogeneous Space

The Clifford Algebras as well as the corresponding Spin groups, for any vector space E of the same dimension and bilinear form g of the same signature are isomorphic. The Clifford Algebra $Cl(3)$ is a subalgebra of $Cl(3, 1)$ and $Spin(3)$ a subgroup of $Spin(3, 1)$.

The Clifford algebras and Spin Group structures are built from the product of vectors. The structure $Cl(3)$ can be defined from a set of vectors only if their scalar product is always definite positive. So, in a given vector space (E, g) with Clifford Algebra isomorphic to $Cl(3, 1)$ the set isomorphic to $Cl(3)$ is not unique : there is one set for each choice of a vector $\varepsilon_0 \in E$ such that $g(\varepsilon_0, \varepsilon_0) = -1$.

The sets isomorphic to $Cl(3)$

In E let be E^\perp the orthogonal complement to $\varepsilon_0 : E^\perp = \{u \in E : g(\varepsilon_0, u) = 0\}$. This is a 3 dimensional vector space. The scalar product g^\perp induced on E^\perp by g is definite positive : in a basis of E^\perp the matrix of g^\perp has 3 positive eigen values, otherwise with ε_0 we would have another signature. The Clifford Algebra $Cl(E^\perp, g^\perp)$ generated by (E^\perp, g^\perp) is a subset of $Cl(E, g)$, Clifford isomorphic to $Cl(3)$. There is a morphism of Clifford algebras : $\pi : Cl(E, g) \rightarrow Cl(E^\perp, g^\perp)$ such that : $\forall u \in E : u = \pi(u) + u^0 \varepsilon_0, g(\pi(u), \pi(u)) > 0$ iff $u \neq 0$.

The Spin Group $Spin(E, g)$ isomorphic to $Spin(3, 1)$ is projected into a subgroup $Sp(E^\perp, g^\perp)$ isomorphic to $Spin(3)$.

$Spin(E^\perp, g^\perp)$ is characterized by the following theorem, which can be proven with any orthonormal basis of E containing ε_0 :

Theorem 110 *The subset of $Spin(3, 1)$ of the elements which commute with ε_0 is a subgroup of $Spin(3, 1)$. They leave ε_0 unchanged : $\mathbf{Ad}_{s_r, \varepsilon_0} = s_r \cdot \varepsilon_0 \cdot s_r^{-1} = \varepsilon_0 = [g]_0^i \varepsilon_i$. They read : $s_r = a + v(r, 0)$*

Proof. $s_r = a + v(r, w) + b\varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3$

$$s_r \cdot \varepsilon_0 = \varepsilon_0 \cdot s_r$$

In $Cl(3, 1)$:

$$s \cdot \varepsilon_0 = a\varepsilon_0 + v(r, w) \varepsilon_0 - b\varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3 = \varepsilon_0 \cdot s = a\varepsilon_0 + \varepsilon_0 v(r, w) + b\varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3$$

$$v(r, w) \varepsilon_0 =$$

$$\frac{1}{2} (w^1 \varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_0 + w^2 \varepsilon_0 \cdot \varepsilon_2 \cdot \varepsilon_0 + w^3 \varepsilon_0 \cdot \varepsilon_3 \cdot \varepsilon_0 + r^3 \varepsilon_2 \cdot \varepsilon_1 \cdot \varepsilon_0 + r^2 \varepsilon_1 \cdot \varepsilon_3 \cdot \varepsilon_0 + r^1 \varepsilon_3 \cdot \varepsilon_2 \cdot \varepsilon_0)$$

$$= \frac{1}{2} (w^1 \varepsilon_1 + w^2 \varepsilon_2 + w^3 \varepsilon_3 - r^3 \varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 + r^2 \varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_3 - r^1 \varepsilon_0 \cdot \varepsilon_2 \cdot \varepsilon_3)$$

$$\varepsilon_0 v(r, w)$$

$$= \frac{1}{2} (w^1 \varepsilon_0 \varepsilon_0 \cdot \varepsilon_1 + w^2 \varepsilon_0 \varepsilon_0 \cdot \varepsilon_2 + w^3 \varepsilon_0 \varepsilon_0 \cdot \varepsilon_3 + r^3 \varepsilon_0 \varepsilon_2 \cdot \varepsilon_1 + r^2 \varepsilon_0 \varepsilon_1 \cdot \varepsilon_3 + r^1 \varepsilon_0 \varepsilon_3 \cdot \varepsilon_2)$$

$$= \frac{1}{2} (-w^1 \varepsilon_1 - w^2 \varepsilon_2 - w^3 \varepsilon_3 - r^3 \varepsilon_0 \varepsilon_1 \cdot \varepsilon_2 + r^2 \varepsilon_0 \varepsilon_1 \cdot \varepsilon_3 - r^1 \varepsilon_0 \varepsilon_2 \cdot \varepsilon_3)$$

$$a\varepsilon_0 + \frac{1}{2} (w^1 \varepsilon_1 + w^2 \varepsilon_2 + w^3 \varepsilon_3 - r^3 \varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 + r^2 \varepsilon_0 \varepsilon_1 \cdot \varepsilon_3 - r^1 \varepsilon_0 \varepsilon_2 \cdot \varepsilon_3) - b\varepsilon_1 \cdot$$

$$\varepsilon_2 \cdot \varepsilon_3$$

$$= a\varepsilon_0 + \frac{1}{2} (-w^1 \varepsilon_1 - w^2 \varepsilon_2 - w^3 \varepsilon_3 - r^3 \varepsilon_0 \varepsilon_1 \cdot \varepsilon_2 + r^2 \varepsilon_0 \varepsilon_1 \cdot \varepsilon_3 - r^1 \varepsilon_0 \varepsilon_2 \cdot \varepsilon_3) +$$

$$b\varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3$$

$$\Rightarrow w = 0, b = 0$$

In $Cl(1, 3)$:

$$s \cdot \varepsilon_0 = a\varepsilon_0 - v(g) \varepsilon_0 - b\varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3 = \varepsilon_0 \cdot s = a\varepsilon_0 - \varepsilon_0 v(g) + b\varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3 \Rightarrow b = 0$$

$$v(g) \varepsilon_0 =$$

$$\frac{1}{2} (w^1 \varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_0 + w^2 \varepsilon_0 \cdot \varepsilon_2 \cdot \varepsilon_0 + w^3 \varepsilon_0 \cdot \varepsilon_3 \cdot \varepsilon_0 + r^3 \varepsilon_2 \cdot \varepsilon_1 \cdot \varepsilon_0 + r^2 \varepsilon_1 \cdot \varepsilon_3 \cdot \varepsilon_0 + r^1 \varepsilon_3 \cdot \varepsilon_2 \cdot \varepsilon_0)$$

$$= \frac{1}{2} (-w^4 \varepsilon_1 - w^2 \varepsilon_2 - w^3 \varepsilon_3 - r^3 \varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 + r^2 \varepsilon_0 \varepsilon_1 \cdot \varepsilon_3 - r^1 \varepsilon_0 \varepsilon_2 \cdot \varepsilon_3)$$

$$\varepsilon_0 v(g) = \frac{1}{2} (w^1 \varepsilon_0 \varepsilon_0 \cdot \varepsilon_1 + w^2 \varepsilon_0 \varepsilon_0 \cdot \varepsilon_2 + w^3 \varepsilon_0 \varepsilon_0 \cdot \varepsilon_3 + r^3 \varepsilon_0 \varepsilon_2 \cdot \varepsilon_1 + r^2 \varepsilon_0 \varepsilon_1 \cdot \varepsilon_3 + r^1 \varepsilon_0 \varepsilon_3 \cdot \varepsilon_2)$$

$$= \frac{1}{2} (w^4 \varepsilon_1 + w^2 \varepsilon_2 + w^3 \varepsilon_3 - r^3 \varepsilon_0 \varepsilon_1 \cdot \varepsilon_2 + r^2 \varepsilon_0 \varepsilon_1 \cdot \varepsilon_3 - r^1 \varepsilon_0 \varepsilon_2 \cdot \varepsilon_3)$$

$$\Rightarrow w = 0 \quad \blacksquare$$

So the elements such that $s = v(r, 0) + \epsilon \sqrt{\frac{1}{4} r^t r - 1} \varepsilon_5$ are excluded and we are left with

$$s = \epsilon \left(\sqrt{1 - \frac{1}{4} r^t r} + v(r, 0) \right) \quad (\text{A.8})$$

These elements constitute a subgroup, as it can easily be checked with the formula for the product. They are generated by vectors belonging to the subspace spanned by the vectors $(\varepsilon_i)_{i=1}^3$ so they belong to $Cl(E^\perp, g^\perp)$ and can be identified with $Spin(E^\perp, g^\perp)$, isomorphic to $Sp(3)$.

The scalars $\epsilon = \pm 1$ belong to the group. The group is not connected. The elements $s = \sqrt{1 - \frac{1}{4}r^t r} + v(r, 0)$ constitute the component of the identity.

Decomposition of the Lie algebra

The Lie algebra $T_1 Spin(3, 1)$ is the direct sum of the two vector vector subspaces :

$$X \in T_1 Spin(3, 1) :: X = v(r, w) = v(r, 0) + v(0, w)$$

The Lie bracket reads :

$$[v(r, w), v(r', w')] = v(j(r)r' - j(w)w', j(w)r' + j(r)w')$$

thus :

$$[v(r, 0), v(r', 0)] = v(j(r)r', 0)$$

$$[v(0, w), v(0, w')] = v(-j(w)w', 0)$$

The vector subspace $v(r, 0)$ can be identified with $T_1 Spin(3)$. It is generated by the vectors $(\varepsilon_1, \varepsilon_2, \varepsilon_3)$ if $\langle \varepsilon_0, \varepsilon_0 \rangle = -1$. And similarly in $Cl(1, 3)$.

For any element $v(r, w)$ of $T_1 Spin(3, 1)$ we have the identity :

$$v(r, w) \cdot \varepsilon_0 - \varepsilon_0 \cdot v(r, w) = w$$

$$\text{Thus } v(r, w) \in T_1 Spin(3) \Leftrightarrow v(r, w) \cdot \varepsilon_0 - \varepsilon_0 \cdot v(r, w) = 0 \Leftrightarrow w = 0$$

Homogeneous space

The quotient space $Spin(3, 1)/Spin(3)$ (called a homogeneous space) is not a group but a 3 dimensional manifold. It is characterized by the equivalence relation :

$$s = a + v(r, w) + b\varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3 \sim s' = a' + v(r', w') + b'\varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3$$

$$\Leftrightarrow \exists s_r \in Spin(3) : s' = s \cdot s_r$$

As any quotient space its elements are *subsets* of $Spin(3, 1)$.

Theorem 111 *In each class of the homogeneous space there are two elements, defined up to sign, which read : $s_w = \pm (a_w + v(0, w))$*

Proof. Each coset $[s]$ is in bijective correspondence with $Spin(3)$.

$$\text{So } [s] = \left\{ s' = s \cdot \left(\epsilon \sqrt{1 - \frac{1}{4}\rho^t \rho} + v(\rho, 0) \right), \rho^t \rho \leq 4 \right\}$$

In $Spin(3, 1)$:

$$s = a + v(r, w) + b\varepsilon_5$$

$$s' = a' + v(r', w') + b'\varepsilon_5$$

$$a' = a\epsilon \sqrt{1 - \frac{1}{4}\rho^t \rho} - \frac{1}{4}r^t \rho$$

$$b' = b\epsilon \sqrt{1 - \frac{1}{4}\rho^t \rho} - \frac{1}{4}w^t \rho$$

$$r' = \frac{1}{2}j(r)\rho + r\epsilon \sqrt{1 - \frac{1}{4}\rho^t \rho} + a\rho$$

$$\begin{aligned} w' &= \frac{1}{2}j(w)\rho + w\epsilon\sqrt{1 - \frac{1}{4}\rho^t\rho} + b\rho \\ a^2 - b^2 &= 1 + \frac{1}{4}(w^t w - r^t r) \\ ab &= -\frac{1}{4}r^t w \end{aligned}$$

We can always choose in the class an element s' such that : $r' = 0$. It requires

$$: \left(\frac{1}{2}j(r) + aI\right)\rho = -r\epsilon\sqrt{1 - \frac{1}{4}\rho^t\rho}$$

This linear equation in ρ has always a unique solution :

$$\begin{aligned} \rho &= -\epsilon\sqrt{1 - \frac{1}{4}\rho^t\rho} \frac{1}{a(4a^2 + r^t r)} \left((4a^2 + r^t r) I - 2aj(r) + j(r)j(r) \right) r \\ &= -\epsilon\sqrt{1 - \frac{1}{4}\rho^t\rho} \frac{1}{a(4a^2 + r^t r)} (4a^2 + r^t r) r \end{aligned}$$

$$= -\epsilon\sqrt{1 - \frac{1}{4}\rho^t\rho} \frac{1}{a} r$$

$$\begin{aligned} \rho^t\rho &= \left(1 - \frac{1}{4}\rho^t\rho\right) \frac{1}{a^2} (r^t r) \Rightarrow \\ (a^2 + \frac{1}{4}(r^t r)) \rho^t\rho &= (r^t r) \end{aligned}$$

$$\rho^t\rho = \frac{4(r^t r)}{4a^2 + (r^t r)} \leq 4$$

$$\sqrt{1 - \frac{1}{4}\rho^t\rho} = \sqrt{\frac{4a^2}{4a^2 + (r^t r)}} = \frac{2a}{\sqrt{4a^2 + (r^t r)}}$$

$$\rho = -\epsilon \frac{2}{\sqrt{4a^2 + (r^t r)}} r$$

$$a' = a\epsilon\sqrt{1 - \frac{1}{4}\rho^t\rho} - \frac{1}{4}r^t\rho = \frac{1}{2} \frac{\epsilon}{\sqrt{4a^2 + (r^t r)}} (4a^2 + r^t r) = \frac{1}{2}\epsilon\sqrt{4a^2 + r^t r}$$

$$w' = \frac{1}{2}j(w)\rho + w\epsilon\sqrt{1 - \frac{1}{4}\rho^t\rho} + b\rho = \epsilon \frac{2}{\sqrt{4a^2 + (r^t r)}} \left(\frac{1}{2}j(r)w + aw - br\right)$$

$$b' = b\epsilon\sqrt{1 - \frac{1}{4}\rho^t\rho} - \frac{1}{4}w^t\rho = \epsilon \frac{2}{\sqrt{4a^2 + (r^t r)}} \left(ab + \frac{1}{4}w^t r\right) = 0$$

$$s = (\epsilon\sqrt{4a^2 + r^t r}) \times$$

$$\left(\frac{1}{2} + v\left(0, \epsilon \frac{2}{\sqrt{4a^2 + (r^t r)}} \left(\frac{1}{2}j(r)w + aw - br\right)\right)\right) \cdot \left(\epsilon \frac{2}{\sqrt{4a^2 + (r^t r)}}\right) (a + v(r, 0))$$

So any element of $Spin(3, 1)$ can be written uniquely (up to sign) :

$$s = a + v(r, w) + b\varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3 = s_w \cdot s_r = (a_w + v(0, w_w)) \cdot (a_r + v(0, r_r))$$

and :

$$(a_w + v(0, w_w)) \cdot (a_r + v(0, r_r)) = a_w a_r + v(a_w r_r, a_r w_w) - \frac{1}{2}(w_w^t r_r) \varepsilon_5$$

In $Cl(1, 3)$ we have the same decomposition with the same components.

$$s = a + v(r, w) + b\varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3 =$$

$$r'' = a_w r_r = \frac{1}{2}\epsilon\sqrt{4a^2 + r^t r} \epsilon \frac{2}{\sqrt{4a^2 + (r^t r)}} r = r$$

$$w'' = \frac{1}{2}j(w_w) r_r + a_r w_w$$

$$= \frac{1}{2}j\left(\left(\epsilon\sqrt{4a^2 + r^t r}\right) \epsilon \frac{2}{\sqrt{4a^2 + (r^t r)}} \left(\frac{1}{2}j(r)w + aw - br\right)\right) \left(\epsilon \frac{2}{\sqrt{4a^2 + (r^t r)}}\right) r$$

$$+ \left(\epsilon \frac{2}{\sqrt{4a^2 + (r^t r)}}\right) a \left(\epsilon\sqrt{4a^2 + r^t r}\right) \epsilon \frac{2}{\sqrt{4a^2 + (r^t r)}} \left(\frac{1}{2}j(r)w + aw - br\right)$$

$$= 2j\left(\epsilon \frac{1}{\sqrt{4a^2 + (r^t r)}} \left(\frac{1}{2}j(r)w + aw - br\right)\right) r + a\epsilon \frac{4}{\sqrt{4a^2 + (r^t r)}} \left(\frac{1}{2}j(r)w + aw - br\right)$$

$$= \left(\epsilon \frac{2}{\sqrt{4a^2 + (r^t r)}}\right) \left(j\left(\left(\frac{1}{2}j(r)w + aw - br\right)\right) r + a2\left(\frac{1}{2}j(r)w + aw - br\right)\right)$$

$$= \left(\epsilon \frac{2}{\sqrt{4a^2 + (r^t r)}}\right) \left(\frac{1}{2}j(j(r)w)r - aj(w)r + aj(r)w + 2a^2w - 2abr\right)$$

$$\begin{aligned}
&= \left(\epsilon \frac{2}{4a^2 + (r^t r)} \right) \left(\frac{1}{2} (wr^t - rw^t) r + 2a^2 w + \frac{1}{2} (r^t w) r \right) \\
&= \left(\epsilon \frac{2}{4a^2 + (r^t r)} \right) \left(\frac{1}{2} w (r^t r) - \frac{1}{2} r (w^t r) + 2a^2 w + \frac{1}{2} (r^t w) r \right) \\
&= \left(\epsilon \frac{1}{4a^2 + (r^t r)} \right) \left((4a^2 + (r^t r)) w \right) = w \quad \blacksquare
\end{aligned}$$

Remark : the elements $\pm s_w$ are equivalent :

$$(a_w + v(0, w_w)) \sim -(a_w + v(0, w_w))$$

Take $s_r = -1 \in Spin(3)$: $-s_w = s_w \cdot s_r$

So $\pm s_w$ belong to the same class of equivalence. In the decomposition :
 $s = \epsilon s_w \cdot \epsilon s_r$, ϵs_w is a specific projection of s on the homogenous space.

A.1.4 Adjoint map

The translations on $Spin(3, 1)$ are :

$$L_g h = g \cdot h, R_g h = h \cdot g$$

and their derivatives :

$$L'_g h : T_h Spin(3, 1) \rightarrow T_{g \cdot h} Spin(3, 1) :: L'_g h(X_h) = g \cdot X_h$$

$$R'_g h : T_h Spin(3, 1) \rightarrow T_{h \cdot g} Spin(3, 1) :: R'_g h(X_h) = X_h \cdot g$$

Their inverse are, as in any Lie groups :

$$(L'_g h)^{-1} = L'_{g^{-1}}(g \cdot h); (R'_g h)^{-1} = R'_{g^{-1}}(h \cdot g)$$

$T_g Spin(3, 1) \subset Cl(3, 1)$ and there are two linear maps :

$$L'_{g^{-1}} g : T_g Spin(3, 1) \rightarrow T_1 Spin(3, 1) :: L'_{g^{-1}} g(Z_g) = g^{-1} \cdot Z_g$$

$$R'_{g^{-1}} g : T_g Spin(3, 1) \rightarrow T_1 Spin(3, 1) :: R'_{g^{-1}} g(Z_g) = Z_g \cdot g^{-1}$$

And the adjoint map:

$$Ad_g : T_1 Spin(3, 1) \rightarrow T_1 Spin(3, 1) :: Ad_g = L'_g g^{-1} \circ R'_{g^{-1}} 1 = R'_{g^{-1}} g \circ L'_g 1$$

$$Ad_g Z = L'_g g^{-1} \circ R'_{g^{-1}} 1(Z) = L'_g g^{-1}(Z \cdot g^{-1}) = g \cdot Z \cdot g^{-1}$$

$$Ad_g Z = (Ad_g Z)'_{x=1} = Ad_g Z$$

With

$$g = a + v(r, w) + b\epsilon_5$$

$$Z = v(x, y)$$

$$Ad_g X = (a + v(r, w) + b\epsilon_5) \cdot v(x, y) \cdot (a - v(r, w) + b\epsilon_5)$$

A straightforward computation gives :

$$Ad_g v(x, y) = v(X, Y)$$

with

$$X =$$

$$\left[(a+b)^2 - ab + \frac{1}{4} r^t r + (a+b)j(r) + \frac{1}{4} (j(w)j(r) - 2j(w)j(w) + 3j(r)j(r)) \right] x$$

$$+ \left[ab + \frac{1}{4} r^t r + (b-a)j(w) - \frac{1}{4} (3j(w)j(r) + 2j(r)j(w) - j(r)j(r)) \right] y$$

$$Y =$$

$$\left[\frac{1}{4} w^t w - ab + (a+b)j(w) + \frac{1}{4} (j(w)j(w) + 2j(w)j(r) + 3j(r)j(w)) \right] x$$

$$+ \left[(a-b)^2 - \frac{1}{4} w^t w - ab + (a-b)j(r) - \frac{1}{4} (3j(w)j(w) - j(r)j(w) - 2j(r)j(r)) \right] y$$

A.1.5 Derivatives

Let $g : M \rightarrow Spin(3, 1) :: g(m) = a(m) + v(r(m), w(m)) + b(m)\epsilon_5$

$$g'(m) : T_m M \rightarrow T_g Spin(3, 1) ::$$

$g'(m)u_m = a'(m)u_m + v(r'(m)u_m, w'(m)u_m) + b'(m)u_m\varepsilon_5$
 where $u_m \in T_mM$, $a'(m)u_m, b'(m)u_m \in \mathbb{R}$, $r'(m)u_m, w'(m)u_m \in \mathbb{R}^3$

Thus :

$g^{-1} \cdot (a'(m)u_m + v(r'(m)u_m, w'(m)u_m) + b'(m)u_m\varepsilon_5) \in T_1Spin(3, 1)$
 $(a'(m)u_m + v(r'(m)u_m, w'(m)u_m) + b'(m)u_m\varepsilon_5) \cdot g^{-1} \in T_1Spin(3, 1)$

$L'_{g^{-1}}g(g'(m)u_m) =$

$v(-\frac{1}{2}(j(r)r' - j(w)w') - a'r + ar' + b'w - bw'),$
 $-\frac{1}{2}(j(w)r' + j(r)w') - a'w + aw' - b'r + br')$

$R'_{g^{-1}}g(g'(m)u_m) =$

$v(-\frac{1}{2}(j(r)r' - j(w)w') - a'r + ar' + b'w - bw'),$
 $-\frac{1}{2}(j(w)r' + j(r)w') - a'w + aw' - b'r + br')$

A.1.6 Exponential on T_1Spin

The exponential on a Lie algebra is the flow of left invariant vector fields (Maths.22.2.6).

i) Left invariant vector fields on $Spin(3, 1)$

As $Spin(3, 1) \subset Cl(3, 1)$ which is a vector space, a vector field $X \in \mathfrak{X}(TSpin(3, 1))$

reads $X(\sigma) \in Cl(3, 1)$ with the relation :

$L'_g\sigma(X(\sigma)) = X(L_g\sigma) = g \cdot X(\sigma) = X(g \cdot \sigma)$

Thus the left invariant vector fields read :

$X(\sigma) = \sigma \cdot v(R, W)$ with $v(R, W) \in T_1Spin(3, 1)$

ii) The flow of $X = \sigma \cdot v(R, W) \in \mathfrak{X}(TSpin(3, 1))$ reads:

$\Phi_X(t, 1) = a(t) + v(r(t), w(t)) + b(t)\varepsilon_5 \in Spin(3, 1)$

$\Phi_X(t, 1) = \exp tX = \exp v(tR, tW)$

$\exp v(tR, tW) = a(t) + v(r(t), w(t)) + b(t)\varepsilon_5$

$\frac{d}{dt} \exp v(tR, tW) |_{t=\theta} = \exp v(\theta R, \theta W) \cdot v(R, W)$

$\frac{d}{dt} (a(t) + v(r(t), w(t)) + b(t)\varepsilon_5) |_{t=\theta} = (a(\theta) + v(r(\theta), w(\theta)) + b(\theta)\varepsilon_5) \cdot v(R, W)$

with :

$a^2 - b^2 = 1 + \frac{1}{4}(w^t w - r^t r)$

$ab = -\frac{1}{4}r^t w$

$1 = a(0) + v(r(0), w(0)) + b(0)\varepsilon_5$

The derivations give :

$\frac{\partial a}{\partial t} + v(\frac{\partial r}{\partial t}, \frac{\partial w}{\partial t}) + \frac{\partial b}{\partial t}\varepsilon_5 |_{t=\theta} = (a(\theta) + v(r(\theta), w(\theta)) + b(\theta)\varepsilon_5) \cdot v(R, W)$

$= a(\theta)v(R, W) + b(\theta)v(R, -W) + v(r(\theta), w(\theta)) \cdot v(R, W)$

$= v((a(\theta) + b(\theta))R, (a(\theta) - b(\theta))W)$

$+ \frac{1}{4}(W^t w - R^t r) + \frac{1}{2}v(-j(R)r + j(W)w, -j(W)r - j(R)w) - \frac{1}{4}(W^t r + R^t w)\varepsilon_5$

$\frac{\partial a}{\partial t} |_{t=\theta} = \frac{1}{4}(W^t w - R^t r)$

$\frac{\partial b}{\partial t} |_{t=\theta} = -\frac{1}{4}(W^t r + R^t w)$

$\frac{\partial r}{\partial t} |_{t=\theta} = (a(\theta) + b(\theta))R + \frac{1}{2}(-j(R)r + j(W)w)$

$\frac{\partial w}{\partial t} |_{t=\theta} = (a(\theta) - b(\theta))W + \frac{1}{2}(-j(W)r - j(R)w)$

$a \frac{\partial a}{\partial t} |_{t=\theta} - b \frac{\partial b}{\partial t} |_{t=\theta} = \frac{1}{4}(w^t \frac{\partial w}{\partial t} - r^t \frac{\partial r}{\partial t})$

$\frac{\partial a}{\partial t} b + a \frac{\partial b}{\partial t} = -\frac{1}{4}r^t \frac{\partial w}{\partial t} - \frac{1}{4}w^t \frac{\partial r}{\partial t}$

The last two equations give :

$$b(W + R)^t(w + r) = 0$$

$$b(W + R)^t(w - r) = 0$$

iii) We have the morphism :

$\Pi : Spin(3, 1) \rightarrow SO(3, 1) :: \Pi(\pm\sigma) = [h(\sigma)]$ such that :

$$\forall u \in \mathbb{R}^4 : \mathbf{Ad}_\sigma u = \sigma \cdot u \cdot \sigma^{-1} = [h(\sigma)] u = \Pi(\pm\sigma) u$$

Take a vector field $X(\sigma) = \sigma \cdot v(R, W) \in \mathfrak{X}(TSpin(3, 1))$ then (Maths.1460)

:

$$\Pi \circ \Phi_X = \Phi_{\Pi_* X} \circ \Pi$$

$$\Pi_* X([h(\sigma)]) = \Pi'(\sigma) X(\sigma)$$

$$\Pi'(1) X(1) = K(W) + J(R)$$

$$\Pi(\exp tv(R, W)) = \Phi_{\Pi'(1)v(R, W)}(t, \Pi(1)) = \exp t(K(W) + J(R)) = \exp tK(W) \exp tJ(R)$$

$$\Pi(\exp tv(R, 0)) = \exp tJ(R)$$

$$\Pi(\exp tv(0, W)) = \exp tK(W)$$

$$\Pi(\exp tv(R, W)) = \Pi(\exp tv(0, W)) \Pi(\exp tv(R, 0))$$

and because this is a morphism :

$$\exp tv(R, W) = \exp tv(0, W) \cdot \exp tv(R, 0)$$

iv) Coming back to the previous equations :

For $\exp v(0, tW)$:

$$\frac{\partial a}{\partial t} \Big|_{t=\theta} = \frac{1}{4} W^t w$$

$$\frac{\partial b}{\partial t} \Big|_{t=\theta} = -\frac{1}{4} (W^t r)$$

$$\frac{\partial r}{\partial t} \Big|_{t=\theta} = \frac{1}{2} j(W) w$$

$$\frac{\partial w}{\partial t} \Big|_{t=\theta} = (a(\theta) - b(\theta)) W - \frac{1}{2} j(W) r$$

$$bW^t(w + r) = 0$$

$$bW^t(w - r) = 0$$

if $b \neq 0$:

$$W^t w = -W^t r = W^t r = 0$$

$$b \frac{\partial b}{\partial t} \Big|_{t=\theta} = -\frac{1}{4} b (W^t r) = 0 \Rightarrow b^2 = Ct \Rightarrow b = Ct \Rightarrow W^t r = 0,$$

$$\Rightarrow W^t w = 0 \Rightarrow a = Ct$$

$$\Rightarrow r, w = Ct$$

$$\text{Thus } b = 0 \Rightarrow W^t r = 0$$

$$\frac{d^2 w}{dt^2} = \frac{1}{4} (W^t w) W - \frac{1}{4} j(W) j(W) w = \frac{1}{4} (W^t w) W - \frac{1}{4} (WW^t w - (W^t W) w) = \frac{1}{4} (W^t W) w$$

$$w(t) = w_1 \exp \frac{1}{2} t \sqrt{W^t W} + w_2 \exp \left(-\frac{1}{2} t \sqrt{W^t W} \right)$$

$$w(0) = 0 = w_1 + w_2$$

$$\frac{dw}{dt}(0) = W = w_1 - w_2$$

$$w(t) = \frac{1}{2} \left(\exp \frac{1}{2} t \sqrt{W^t W} - \exp \left(-\frac{1}{2} t \sqrt{W^t W} \right) \right) W = W \sinh \frac{1}{2} t \sqrt{W^t W}$$

$$\frac{\partial r}{\partial t} \Big|_{t=\theta} = \frac{1}{2} j(W) W \sinh \frac{1}{2} t \sqrt{W^t W} = 0$$

$$r(0) = R = 0 \Rightarrow r(t) = 0$$

$$w^t w = W^t W \sinh^2 \frac{1}{2} t \sqrt{W^t W}$$

$$a^2 - b^2 = a^2 = 1 + \frac{1}{4} \left(W^t W \sinh^2 \frac{1}{2} t \sqrt{W^t W} \right)$$

$$\exp v(0, tW) = \sqrt{1 + \frac{1}{4} \left(W^t W \sinh^2 \frac{1}{2} t \sqrt{W^t W} \right)} + \sinh \frac{1}{2} t \sqrt{W^t W} v(0, W)$$

For $\exp v(tR, 0)$

$$\begin{aligned}
\frac{\partial a}{\partial t}|_{t=\theta} &= -\frac{1}{4}R^t r \\
\frac{\partial b}{\partial t}|_{t=\theta} &= -\frac{1}{4}R^t w \\
\frac{\partial r}{\partial t}|_{t=\theta} &= (a(\theta) + b(\theta))R - \frac{1}{2}j(R)r \\
\frac{\partial w}{\partial t}|_{t=\theta} &= -\frac{1}{2}j(R)w \\
bR^t(w+r) &= 0 \\
bR^t(w-r) &= 0 \\
\Rightarrow b &= 0, R^t w = 0 \\
\frac{d^2 r}{dt^2} &= -\frac{1}{4}(R^t r)R - \frac{1}{2}j(R)(aR - \frac{1}{2}j(R)r) = -\frac{1}{4}(R^t r)R - \frac{1}{2}aj(R)R + \\
&\frac{1}{4}j(R)j(R)r \\
\frac{d^2 r}{dt^2} &= -\frac{1}{4}(R^t r)R + \frac{1}{4}((R^t r)R - R^t R r) \\
\frac{d^2 r}{dt^2} &= -\frac{1}{4}(R^t R)r \\
r(t) &= r_1 \exp it \frac{1}{2} \sqrt{R^t R} + r_2 \exp \left(-it \frac{1}{2} \sqrt{R^t R} \right) \\
r(0) = 0 &= r_1 + r_2 \\
\frac{dr}{dt}(0) &= R = r_1 - r_2 \\
r(t) &= R \sin t \frac{1}{2} \sqrt{R^t R} \\
r^t r &= R^r R \sin^2 t \frac{1}{2} \sqrt{R^t R} \\
a^2 - b^2 = a^2 &= 1 - \frac{1}{4}R^r R \sin^2 t \frac{1}{2} \sqrt{R^t R} \\
\exp tv(R, 0) &= \sqrt{1 - \frac{1}{4}R^r R \sin^2 t \frac{1}{2} \sqrt{R^t R}} + \sin t \frac{1}{2} \sqrt{R^t R} v(R, 0) \\
\exp tv(R, W) &= \left(\sqrt{1 + \frac{1}{4}(W^t W \sinh^2 \frac{1}{2} t \sqrt{W^t W})} + \sinh \frac{1}{2} t \sqrt{W^t W} v(0, W) \right) \cdot \\
&\left(\sqrt{1 - \frac{1}{4}R^r R \sin^2 t \frac{1}{2} \sqrt{R^t R}} + \sin t \frac{1}{2} \sqrt{R^t R} v(R, 0) \right)
\end{aligned}$$

A.2 COMPUTATION OF X

$$\begin{aligned}
(\sigma^{-1} \partial_\alpha \sigma + Ad_{\sigma^{-1}} G_\alpha) &= v(X_\alpha, Y_\alpha) \\
\sigma &= \sigma_w \cdot \sigma_r \\
\sigma^{-1} \partial_\alpha \sigma + Ad_{\sigma^{-1}} G_\alpha &= \sigma_r^{-1} \cdot \sigma_w^{-1} \cdot \partial_\alpha \sigma_w \cdot \sigma_r + \sigma_r^{-1} \cdot \sigma_w^{-1} \cdot \sigma_w \cdot \partial_\alpha \sigma_r + Ad_{\sigma_r^{-1}} Ad_{\sigma_w^{-1}} G_\alpha \\
&= \sigma_r^{-1} \cdot \left(\sigma_w^{-1} \cdot \partial_\alpha \sigma_w \cdot \sigma_r + \partial_\alpha \sigma_r + \left(Ad_{\sigma_w^{-1}} G_\alpha \right) \cdot \sigma_r \right) \\
&= \sigma_r^{-1} \cdot \left(\sigma_w^{-1} \cdot \partial_\alpha \sigma_w + \partial_\alpha \sigma_r \cdot \sigma_r^{-1} + \left(Ad_{\sigma_w^{-1}} G_\alpha \right) \right) \cdot \sigma_r \\
&= Ad_{\sigma_r^{-1}} \left(\sigma_w^{-1} \cdot \partial_\alpha \sigma_w + \partial_\alpha \sigma_r \cdot \sigma_r^{-1} + \left(Ad_{\sigma_w^{-1}} G_\alpha \right) \right) \\
&= Ad_{\sigma_r^{-1}} \left(\sigma_w^{-1} \cdot \partial_\alpha \sigma_w + \partial_\alpha \sigma_r \cdot \sigma_r^{-1} + \left(Ad_{\sigma_w^{-1}} G_\alpha \right) \right) \\
\sigma_w &= a_w + v(0, w) \\
\sigma_r &= a_r + v(r, 0) \\
a_w^2 &= 1 + \frac{1}{4} w^t w \\
a_r^2 &= 1 - \frac{1}{4} r^t r \\
\partial_\alpha a_w &= \frac{1}{4a_w} w^t \partial_\alpha w \\
\partial_\alpha a_r &= -\frac{1}{4a_r} r^t \partial_\alpha r \\
\sigma_w^{-1} \cdot \partial_\alpha \sigma_w &
\end{aligned}$$

$$\begin{aligned}
&= v \left(\frac{1}{2} j(w) \partial_\alpha w, -w \partial_\alpha a_w + a_w \partial_\alpha w \right) \\
&= v \left(\frac{1}{2} j(w) \partial_\alpha w, -\frac{1}{4a_w} w w^t \partial_\alpha w + a_w \partial_\alpha w \right) \\
&= v \left(\frac{1}{2} j(w) \partial_\alpha w, -\frac{1}{4a_w} (j(w) j(w) + w^t w) \partial_\alpha w + a_w \partial_\alpha w \right) \\
&= v \left(\frac{1}{2} j(w) \partial_\alpha w, \left(-\frac{1}{4a_w} j(w) j(w) - 4\frac{1}{4a_w} (a_w^2 - 1) + a_w \right) \partial_\alpha w \right) \\
&= v \left(\frac{1}{2} j(w) \partial_\alpha w, \left(\frac{1}{a_w} - \frac{1}{4a_w} j(w) j(w) \right) \partial_\alpha w \right) \\
&\quad \partial_\alpha \sigma_r \cdot \sigma_r^{-1} \\
&= v \left(-\frac{1}{2} j(r) \partial_\alpha r - r \partial_\alpha a_r + a_r \partial_\alpha r, 0 \right) \\
&= v \left(-\frac{1}{2} j(r) \partial_\alpha r + \frac{1}{4a_r} r r^t \partial_\alpha r + a_r \partial_\alpha r, 0 \right) \\
&= v \left(\left(-\frac{1}{2} j(r) + \frac{1}{4a_r} (j(r) j(r) + r^t r) + a_r \right) \partial_\alpha r, 0 \right) \\
&= v \left(\left(-\frac{1}{2} j(r) + \frac{1}{4a_r} j(r) j(r) + \frac{4(1-a_r^2)}{4a_r} + a_r \right) \partial_\alpha r, 0 \right) \\
&= v \left(\left(\frac{1}{a_r} - \frac{1}{2} j(r) + \frac{1}{4a_r} j(r) j(r) \right) \partial_\alpha r, 0 \right) \\
&\quad Ad_{\sigma_w^{-1}} G_\alpha = \sigma_w^{-1} \cdot G_\alpha \cdot \sigma_w \\
&= v \left([a_w^2 - \frac{1}{2} j(w) j(w)] G_{r\alpha} + a_w j(w) \widehat{G}_w, \right. \\
&\quad [a_w^2 - 1 - a_w j(w) + \frac{1}{4} j(w) j(w)] G_{r\alpha} + [1 - \frac{3}{4} j(w) j(w)] \widehat{G}_w \left. \right) \\
&\quad \sigma_w^{-1} \cdot \partial_\alpha \sigma_w + \partial_\alpha \sigma_r \cdot \sigma_r^{-1} + \left(Ad_{\sigma_w^{-1}} G_\alpha \right) \\
&= v \left(\frac{1}{2} j(w) \partial_\alpha w + \left(\frac{1}{a_r} - \frac{1}{2} j(r) + \frac{1}{4a_r} j(r) j(r) \right) \partial_\alpha r + [a_w^2 - \frac{1}{2} j(w) j(w)] G_{r\alpha} + \right. \\
&\quad \left. a_w j(w) G_{w\alpha}, \right. \\
&\quad \left. \left(\frac{1}{a_w} - \frac{1}{4a_w} j(w) j(w) \right) \partial_\alpha w + [a_w^2 - 1 - a_w j(w) + \frac{1}{4} j(w) j(w)] \widehat{G}_r + [1 - \frac{3}{4} j(w) j(w)] G_{w\alpha} \right) \\
&\quad Ad_{\sigma_r^{-1}} v(x, y) = v(X, Y) \\
&\quad \text{with} \\
&\quad X = [1 - a_r j(r) + \frac{3}{4} j(r) j(r)] x + [1 - a_r^2 + \frac{1}{4} j(r) j(r)] y \\
&\quad Y = [a_r^2 - a_r j(r) + \frac{1}{2} j(r) j(r)] y \\
&\quad x = \left(\frac{1}{a_r} - \frac{1}{2} j(r) + \frac{1}{4a_r} j(r) j(r) \right) \partial_\alpha r + \frac{1}{2} j(w) \partial_\alpha w + (a_w^2 - \frac{1}{2} j(w) j(w)) G_{r\alpha} + \\
&\quad a_w j(w) G_{w\alpha} \\
&\quad y = \frac{1}{a_w} \left(1 - \frac{1}{4} j(w) j(w) \right) \partial_\alpha w + [a_w^2 - 1 - a_w j(w) + \frac{1}{4} j(w) j(w)] G_{r\alpha} + [1 - \frac{3}{4} j(w) j(w)] G_{w\alpha}
\end{aligned}$$

$$X_\alpha = [A_r(r)] \partial_\alpha r + [A_w(r, w)] \partial_\alpha w + [D_r(r, w)] G_{r\alpha} + [D_w(r, w)] G_{w\alpha} \quad (\text{A.9})$$

$$[A_r(r)] = [B_r(r)] \left[\frac{1}{a_r} - \frac{1}{2} j(r) + \frac{1}{4a_r} j(r) j(r) \right] = \left[\frac{1}{a_r} + \frac{2-5a_r^2}{2} j(r) + \frac{5a_r^2+1}{4a_r} j(r) j(r) \right]$$

$$[A_w(r, w)] = [B_r(r)] [B_w(w)] + [C_r(r)] [C_w(w)]$$

$$[D_r(r, w)] = [B_r(r)] [B_{wr}(w)] + [C_r(r)] [C_{wr}(w)]$$

$$[D_w(r, w)] = [B_r(r)] [B_{ww}(w)] + [C_r(r)] [C_{ww}(w)]$$

with :

$$[B_r(r)] = [1 - a_r j(r) + \frac{3}{4} j(r) j(r)]$$

$$[B_w(w)] = \frac{1}{2} j(w)$$

$$[C_r(r)] = [1 - a_r^2 + \frac{1}{4} j(r) j(r)]$$

$$[C_w(w)] = \frac{1}{a_w} [1 - \frac{1}{4} j(w) j(w)]$$

$$\begin{aligned}
[B_{wr}(w)] &= [a_w^2 - \frac{1}{2}j(w)j(w)] \\
[B_{ww}(w)] &= a_w [j(w)] \\
[C_{wr}(w)] &= [a_w^2 - 1 - a_w j(w) + \frac{1}{4}j(w)j(w)] \\
[C_{ww}(w)] &= [1 - \frac{3}{4}j(w)j(w)] \\
a_r^2 &= 1 - \frac{1}{4}r^t r \\
a_w^2 &= 1 - \frac{1}{4}w^t w
\end{aligned}$$

A.3 COMPUTATION OF THE CURRENTS

$$\begin{aligned}
a = 1, 2, 3 : J_G &= \frac{C_I}{4C_G} \mu_i^{\frac{1}{2}} \langle \psi, \gamma C(\vec{\kappa}_a) \psi \rangle V \otimes \vec{\kappa}_a \in T_1 Spin(3, 1) \otimes TM \\
a = 4, 5, 6 : J_G &= -\frac{C_I}{4C_G} \mu_i^{\frac{1}{2}} \langle \psi, \gamma C(\vec{\kappa}_a) \psi \rangle V \otimes \vec{\kappa}_a \in T_1 Spin(3, 1) \otimes TM \\
\psi &= \gamma C(\sigma_w \cdot \sigma_r) \psi_0 \varrho(\mathfrak{X}) \\
\langle \psi, \gamma C(\vec{\kappa}_a) \psi \rangle &= \langle \gamma C(\sigma_w \cdot \sigma_r) \psi_0 \varrho(\mathfrak{X}), \gamma C(\vec{\kappa}_a) \gamma C(\sigma_w \cdot \sigma_r) \psi_0 \varrho(\mathfrak{X}) \rangle \\
&= \langle \psi_0, \gamma C(\sigma_r^{-1} \cdot \sigma_w^{-1}) \gamma C(\vec{\kappa}_a) \gamma C(\sigma_w \cdot \sigma_r) \psi_0 \rangle \\
&= \gamma C(\sigma_r^{-1} \cdot \sigma_w^{-1}) \gamma C(\vec{\kappa}_a) \gamma C(\sigma_w \cdot \sigma_r) \\
&= \gamma C(\sigma_r^{-1} \cdot \sigma_w^{-1} \cdot \vec{\kappa}_a \cdot \sigma_w \cdot \sigma_r) \\
&= \gamma C(\mathbf{Ad}_{\sigma_r^{-1}} \circ \mathbf{Ad}_{\sigma_w^{-1}} \vec{\kappa}_a) \\
a = 1, 2, 3 : \\
\mathbf{Ad}_{\sigma_w^{-1}} \vec{\kappa}_a &= v([a_w^2 - \frac{1}{2}j(w)j(w)] \varepsilon_a, [\frac{1}{4}w^t w - a_w j(w) + \frac{1}{4}j(w)j(w)] \varepsilon_a) \\
\mathbf{Ad}_{\sigma_r^{-1}} \circ \mathbf{Ad}_{\sigma_w^{-1}} \vec{\kappa}_a &= v(X, Y) \\
X &= ([1 - a_r j(r) + \frac{3}{4}j(r)j(r)] [a_w^2 - \frac{1}{2}j(w)j(w)] \\
&+ [1 - a_r^2 + \frac{1}{4}j(r)j(r)] [a_w^2 - 1 - a_w j(w) + \frac{1}{4}j(w)j(w)]) \varepsilon_a \\
Y &= ([a_r^2 - a_r j(r) + \frac{1}{2}j(r)j(r)] [a_w^2 - 1 - a_w j(w) + \frac{1}{4}j(w)j(w)]) \varepsilon_a \\
a = 4, 5, 6 : \\
\mathbf{Ad}_{\sigma_w^{-1}} \vec{\kappa}_a &= v(a_w j(w) \varepsilon_a, [a_w^2 - \frac{1}{4}w^t w - \frac{3}{4}j(w)j(w)] \varepsilon_a) \\
\mathbf{Ad}_{\sigma_r^{-1}} \circ \mathbf{Ad}_{\sigma_w^{-1}} \vec{\kappa}_a &= v(X, Y) \\
X &= ([1 - a_r j(r) + \frac{3}{4}j(r)j(r)] a_w j(w) + [1 - a_r^2 + \frac{1}{4}j(r)j(r)] [1 - \frac{3}{4}j(w)j(w)]) \varepsilon_a \\
Y &= ([a_r^2 - a_r j(r) + \frac{1}{2}j(r)j(r)] [1 + \frac{3}{4}j(w)j(w)]) \varepsilon_a
\end{aligned}$$

$$\begin{aligned}
\gamma C(v([X] \varepsilon_a, [Y] \varepsilon_a)) &= -i \frac{1}{2} \sum_{b=1}^3 \left([[Y] \varepsilon_a]^b \gamma_b \gamma_0 + [[X] \varepsilon_a]^b \tilde{\gamma}_b \right) \\
\langle \psi, \gamma C(\vec{\kappa}_a) \psi \rangle &= -i \frac{1}{2} \langle \psi_0, \left(\sum_{b=1}^3 \left([[Y] \varepsilon_a]^b \gamma_b \gamma_0 + [[X] \varepsilon_a]^b \tilde{\gamma}_b \right) \right) \psi_0 \rangle \\
\psi_0 &= \begin{bmatrix} \psi_R \\ \epsilon i \psi_L \end{bmatrix} \\
\langle \psi, \gamma C(\vec{\kappa}_a) \psi \rangle &= -i \frac{1}{2} Tr \{ \begin{bmatrix} \psi_R^* & -\epsilon i \psi_R^* \end{bmatrix} \gamma_0 \left(\sum_{b=1}^3 \left([[Y] \varepsilon_a]^b \gamma_b \gamma_0 + [[X] \varepsilon_a]^b \tilde{\gamma}_b \right) \right) \begin{bmatrix} \psi_R \\ \epsilon i \psi_R \end{bmatrix} \} \\
&= -i \frac{1}{2} Tr \begin{bmatrix} \psi_R^* & -\epsilon i \psi_R^* \end{bmatrix} \begin{bmatrix} \epsilon i \sum_{b=1}^3 \left(-[[Y] \varepsilon_a]^b - i [[X] \varepsilon_a]^b \right) \sigma_b \psi_R \\ \sum_{b=1}^3 \left(-[[Y] \varepsilon_a]^b + i [[X] \varepsilon_a]^b \right) \sigma_b \psi_R \end{bmatrix} \\
&= -i \frac{1}{2} Tr \left(\epsilon i \sum_{b=1}^3 \psi_R^* \left(-[[Y] \varepsilon_a]^b - i [[X] \varepsilon_a]^b \right) \sigma_b \psi_R - \epsilon i \sum_{b=1}^3 \psi_R^* \left(-[[Y] \varepsilon_a]^b + i [[X] \varepsilon_a]^b \right) \sigma_b \psi_R \right)
\end{aligned}$$

$$= \epsilon \frac{1}{2} \sum_{b=1}^3 \left(-[[Y] \varepsilon_a]^b - i [[X] \varepsilon_a]^b + [[Y] \varepsilon_a]^b - i [[X] \varepsilon_a]^b \right) Tr(\psi_R^* \sigma_b \psi_R)$$

$$= -i \epsilon \sum_{b=1}^3 [[X] \varepsilon_a]^b Tr(\psi_R^* \sigma_b \psi_R)$$

$$\text{With : } k = -\epsilon (Tr(\psi_R^* \sigma_a \psi_R))_{a=1}^3$$

$$\langle \psi, \gamma C(\vec{\kappa}_a) \psi \rangle = ik^t [X] \varepsilon_a$$

$a = 1, 2, 3 :$

$$J_G^a = \frac{C_I}{4C_G} \mu k^t \left\{ [1 - a_r j(r) + \frac{3}{4} j(r) j(r)] [a_w^2 - \frac{1}{2} j(w) j(w)] \right. \\ \left. + [1 - a_r^2 + \frac{1}{4} j(r) j(r)] [a_w^2 - 1 - a_w j(w) + \frac{1}{4} j(w) j(w)] \right\} \varepsilon_a V^\alpha$$

$$J_G^a = \frac{C_I}{4C_G} \mu k^t ([B_r(r)] [B_r(r)] + [C_r(r)] [C_{wr}(w)]) k V^\alpha$$

$a = 4, 5, 6 :$

$$J_G^a = -\frac{C_I}{4C_G} \mu k^t \left\{ [1 - a_r j(r) + \frac{3}{4} j(r) j(r)] a_w j(w) \right. \\ \left. + [1 - a_r^2 + \frac{1}{4} j(r) j(r)] [1 - \frac{3}{4} j(w) j(w)] \right\} \varepsilon_a V^\alpha$$

$$J_G^a = -\frac{C_I}{4C_G} \mu k^t ([B_r(r)] [B_{ww}(w)] + [C_r(r)] [C_{ww}(w)]) \varepsilon_a V^\alpha$$

A.4 FORMULAS

A.4.1 Relativist Geometry

Divergence of a vector field V :

$$\operatorname{div} V = \frac{1}{\det P'} \sum_{\alpha=0}^3 \frac{d}{d\xi^\alpha} (V^\alpha \det P') = \sum_{\alpha=0}^3 \frac{dV^\alpha}{d\xi^\alpha} + \frac{1}{\det P'} \sum_{\alpha=0}^3 V^\alpha \frac{d \det P'}{d\xi^\alpha}$$

For a motion on an integral curve of the vector field V : $\frac{1}{\det P'} \sum_{\alpha=0}^3 V^\alpha \frac{d \det P'}{d\xi^\alpha} = \frac{1}{\det P'} \frac{d \det P'}{d\tau}$

Between the proper time τ of a particle and the time t of an observer :

$$\frac{d\tau}{dt} = \sqrt{1 - \frac{\|\vec{v}\|^2}{c^2}} = \frac{1}{\gamma} = \frac{c}{u^0} = \frac{1}{2a_w^2 - 1}$$

Between the velocity u of a particle and the speed V as measured by an observer :

$$u = \frac{dp}{d\tau} = V \frac{c}{\sqrt{-\langle V, V \rangle}} = \frac{1}{\sqrt{1 - \frac{\|\vec{v}\|^2}{c^2}}} (\vec{v} + c\varepsilon_0(m))$$

$$= c \left((2a_w^2 - 1) \varepsilon_0 + \epsilon a_w \sum_{i=1}^3 w_i \varepsilon_i \right)$$

$$V = \frac{dp}{dt} = u \frac{\sqrt{-\langle V, V \rangle}}{c} = \vec{v} + c\varepsilon_0(m) = c \left(\varepsilon_0 + \epsilon \frac{a_w}{2a_w^2 - 1} \sum_{i=1}^3 w_i \varepsilon_i \right)$$

$$V(t) = \frac{dq}{dt} = c\varepsilon_0(q(t)) + \vec{v}$$

A.4.2 Operator j

Let $r \in \mathbb{C}^3, w \in \mathbb{C}^3$:

$$[j(r)]w = \begin{bmatrix} 0 & -r_3 & r_2 \\ r_3 & 0 & -r_1 \\ -r_2 & r_1 & 0 \end{bmatrix} \begin{bmatrix} w_1 \\ w_2 \\ w_3 \end{bmatrix} = \begin{bmatrix} r_2 w_3 - r_3 w_2 \\ -r_1 w_3 + r_3 w_1 \\ r_1 w_2 - r_2 w_1 \end{bmatrix}$$

$$[j(r)w]^a = \sum_{b,c=1}^3 \epsilon(a,b,c) r_b w_c$$

$$j(r)^t = -j(r) = j(-r)$$

$$j(x)y = -j(y)x$$

$$y^t j(x) = -x^t j(y)$$

$$j(x)y = 0 \Leftrightarrow \exists k \in R : y = kx$$

eigenvectors:

$$r = \sqrt{r^t r}$$

$$0 : \begin{bmatrix} r_1 \\ r_2 \\ r_3 \end{bmatrix}$$

$$ir : \begin{bmatrix} -(-r_1 r_2 + ir_3 r) \\ -(r_1^2 + r_3^2) \\ r_2 r_3 + ir_1 r \end{bmatrix}$$

$$-ir : \begin{bmatrix} -(r_1 r_2 + ir_3 r) \\ (r_1^2 + r_3^2) \\ -r_2 r_3 + r_1 ir \end{bmatrix}$$

Identities

$$\begin{aligned}
j(x)j(y) &= yx^t - (y^t x) I \\
j(x)j(x)j(x) &= j(x)(xx^t - (x^t x) I) = -(x^t x)j(x) \\
j(j(x)y) &= yx^t - xy^t = j(x)j(y) - j(y)j(x) \\
j(j(x)j(x)y) &= (y^t x)j(x) - (x^t x)j(y) \\
j(x)j(y)j(x) &= -(y^t x)j(x) \\
j(x)j(x)j(y) &= j(x)yx^t - (y^t x)j(x) = -j(y)xx^t - (y^t x)j(x) = -j(y)(j(x)j(x) + x^t x) - \\
&\quad (y^t x)j(x) \\
x^t j(r)j(s)y &= x^t(sr^t - r^t s I)y = (x^t s)(r^t y) - (x^t y)(r^t s) \\
\|j(x)y\|^2 &= (x^t x)(y^t y) - (x^t y)^2 \\
M \in L(3) : M^t j(Mx)M &= (\det M)j(x) \\
M \in O(3) : j(Mx)My &= Mj(x)y \Leftrightarrow Mx \times My = M(x \times y) \\
k > 0 : j(r)^{2k} &= (-r^t r)^{k-1}(rr^t - (r^t r)I) = (-r^t r)^{k-1}j(r)j(r) \\
k \geq 0 : J(r)^{2k+1} &= (-r^t r)^k j(r) \\
\exp[j(r)] &= I_3 + [j(r)] \frac{\sin \sqrt{r^t r}}{\sqrt{r^t r}} + [j(r)][j(r)] \frac{1 - \cos \sqrt{r^t r}}{r^t r}
\end{aligned}$$

Polynomials

The set of polynomials $aI + bj(z) + cj(z)j(z)$ where $z \in \mathbb{C}^3$ is fixed, $a, b, c \in \mathbb{C}$ is a commutative ring.

$$\begin{aligned}
&(a + bj(z) + cj(z)j(z))(a' + b'j(z) + c'j(z)j(z)) \\
&= aa' + (ab' + a'b - (r^t r)(c'b + b'c))j(z) + (ac' + b'b + a'c - (r^t r)c'c)j(z)j(z)
\end{aligned}$$

The inverse are :

$$\begin{aligned}
[aI + bj(z) + cj(z)j(z)]^{-1} &= \frac{1}{a}I - \frac{bc^2}{(a - cr^2)^2 + b^2 r^2} j(z) + \frac{(b^2 + c^2 r^2 - a)}{a((a - cr^2)^2 + b^2 r^2)} j(z)j(z) \\
\det(a + bj(r)j(r) + cj(r)j(k) + dj(k)j(r)) &= \\
cd^2(r^2 k^2 z - z^3) + ad^2 z^2 + c^2 d(r^2 k^2 z - z^3) + bcd(r^4 k^2 - r^2 z^2) + acd(-k^2 r^2 + 3z^2) + \\
2abdr^2 z - 2a^2 dz + 2abcr^2 z - 2a^2 cz - 2a^2 cz \\
+ ab^2 r^4 - 2a^2 br^2 + a^3 \\
\text{with } z = r^t k, r^2 = r^t r, k^2 = k^t k
\end{aligned}$$

A.4.3 Dirac's matrices

$$\begin{aligned}
\sigma_1 &= \begin{bmatrix} 0 & 1 \\ 1 & 0 \end{bmatrix}; \sigma_2 = \begin{bmatrix} 0 & -i \\ i & 0 \end{bmatrix}; \sigma_3 = \begin{bmatrix} 1 & 0 \\ 0 & -1 \end{bmatrix}; \sigma_0 = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} \\
\sigma_i \sigma_j + \sigma_j \sigma_i &= 2\delta_{ij} \sigma_0 \\
\sigma_i^2 &= \sigma_0 \\
j \neq k, l = 1, 2, 3 : \sigma_j \sigma_k &= \epsilon(j, k, l) i \sigma_l \\
\sigma_1 \sigma_2 \sigma_3 &= i \sigma_0 \\
\left(\sum_{a=1}^3 U_a \sigma_a \right) \left(\sum_{b=1}^3 V_b \sigma_b \right) &= \sum_{a < b=1}^3 (j(U)V)^a i \sigma_a + \sum_{a=1}^3 U_a V_a \sigma_0 \\
\sum_{a=1}^3 z_a \sigma_a &= \begin{bmatrix} z_3 & z_1 - iz_2 \\ z_1 + iz_2 & -z_3 \end{bmatrix} \\
\left(\sum_{a=1}^3 z_a \sigma_a \right)^{-1} &= \frac{1}{z_1^2 + z_2^2 + z_3^2} \sum_{a=1}^3 z_a \sigma_a
\end{aligned}$$

A.4.4 γ matrices

$$\gamma_0 = \begin{bmatrix} 0 & -i\sigma_0 \\ i\sigma_0 & 0 \end{bmatrix}; \gamma_1 = \begin{bmatrix} 0 & \sigma_1 \\ \sigma_1 & 0 \end{bmatrix}; \gamma_2 = \begin{bmatrix} 0 & \sigma_2 \\ \sigma_2 & 0 \end{bmatrix}; \gamma_3 = \begin{bmatrix} 0 & \sigma_3 \\ \sigma_3 & 0 \end{bmatrix};$$

$$\gamma_i \gamma_j + \gamma_j \gamma_i = 2\delta_{ij} I_4$$

$$\gamma_i = \gamma_i^* = \gamma_i^{-1}$$

$$\gamma_1 \gamma_2 \gamma_3 = i \begin{bmatrix} 0 & \sigma_0 \\ \sigma_0 & 0 \end{bmatrix}$$

$$\gamma_5 = \gamma_0 \gamma_1 \gamma_2 \gamma_3 = \begin{bmatrix} \sigma_0 & 0 \\ 0 & -\sigma_0 \end{bmatrix}$$

$$\gamma_5 \gamma_a = -\gamma_a \gamma_5$$

$$Cl(3, 1) : \gamma C(\varepsilon_j) = \gamma_j, j = 1, 2, 3; \gamma C(\varepsilon_0) = i\gamma_0; \gamma C(\varepsilon_5) = i\gamma_5$$

$$Cl(1, 3) : \gamma C'(\varepsilon_j) = i\gamma_j, j = 1, 2, 3; \gamma C'(\varepsilon_0) = \gamma_0; \gamma C'(\varepsilon_5) = \gamma_5$$

$$j = 1, 2, 3 : \tilde{\gamma}_j = \begin{bmatrix} \sigma_j & 0 \\ 0 & \sigma_j \end{bmatrix}$$

$$j \neq k, l = 1, 2, 3 : \gamma_j \gamma_k = -\gamma_k \gamma_j = i\epsilon(j, k, l) \tilde{\gamma}_l$$

$$a = 1, 2, 3 : \gamma C(\vec{\kappa}_a) = -\frac{1}{2} i \tilde{\gamma}_a$$

$$a = 4, 5, 6 : \gamma C(\vec{\kappa}_a) = \frac{1}{2} \begin{bmatrix} \sigma_a & 0 \\ 0 & -\sigma_a \end{bmatrix}$$

$$\gamma_0 \tilde{\gamma}_j = \tilde{\gamma}_j \gamma_0 = i \begin{bmatrix} 0 & -\sigma_j \\ \sigma_j & 0 \end{bmatrix}$$

$$\tilde{\gamma}_a \tilde{\gamma}_a = I_4$$

$$\tilde{\gamma}_a \tilde{\gamma}_b = i\epsilon(a, b, c) \tilde{\gamma}_c$$

$$\tilde{\gamma}_a \tilde{\gamma}_b + \tilde{\gamma}_b \tilde{\gamma}_a = 2\delta_{ab} I_4$$

$$\gamma_a \tilde{\gamma}_b = i\epsilon(a, b, c) \gamma_c$$

$$\gamma_a \tilde{\gamma}_a = \begin{bmatrix} 0 & \sigma_0 \\ \sigma_0 & 0 \end{bmatrix}$$

$$\gamma_0 \tilde{\gamma}_a = \begin{bmatrix} 0 & -i\sigma_a \\ i\sigma_a & 0 \end{bmatrix} = (\gamma_0 \tilde{\gamma}_a)^* = \begin{bmatrix} 0 & -i\sigma_a^* \\ i\sigma_a^* & 0 \end{bmatrix}$$

$$j = 1, 2, 3 : \gamma_j \gamma_0 = -\gamma_0 \gamma_j = i \begin{bmatrix} \sigma_j & 0 \\ 0 & -\sigma_j \end{bmatrix} = i\gamma_5 \tilde{\gamma}_j$$

$$\gamma_1 \gamma_2 = -\gamma_2 \gamma_1 = i \begin{bmatrix} \sigma_3 & 0 \\ 0 & \sigma_3 \end{bmatrix}; \gamma_2 \gamma_3 = -\gamma_3 \gamma_2 = i \begin{bmatrix} \sigma_1 & 0 \\ 0 & \sigma_1 \end{bmatrix};$$

$$\gamma_3 \gamma_1 = -\gamma_1 \gamma_3 = i \begin{bmatrix} \sigma_2 & 0 \\ 0 & \sigma_2 \end{bmatrix}$$

$$\gamma_1 \gamma_0 = -\gamma_0 \gamma_1 = i \begin{bmatrix} \sigma_1 & 0 \\ 0 & -\sigma_1 \end{bmatrix}; \gamma_2 \gamma_0 = -\gamma_0 \gamma_2 = i \begin{bmatrix} \sigma_2 & 0 \\ 0 & -\sigma_2 \end{bmatrix};$$

$$\gamma_3 \gamma_0 = -\gamma_0 \gamma_3 = i \begin{bmatrix} \sigma_3 & 0 \\ 0 & -\sigma_3 \end{bmatrix}$$

$$\gamma C(a + v(r, w) + b\varepsilon_5) = aI + ib\gamma_5 - i\frac{1}{2} \sum_{a=1}^3 (w_a \gamma_a \gamma_0 + r_a \tilde{\gamma}_a)$$

In $Cl(3, 1)$:

$$\gamma C(v(r, w)) = -i\frac{1}{2} \sum_{a=1}^3 (w_a \gamma_a \gamma_0 + r_a \tilde{\gamma}_a)$$

$$= \frac{1}{2} \begin{bmatrix} \sum_{a=1}^3 (w_a - ir_a) \sigma_a & 0 \\ 0 & -\sum_{a=1}^3 (w_a + ir_a) \sigma_a \end{bmatrix}$$

In $Cl(1, 3)$:

$$\gamma C'(v(r, w)) = -i\frac{1}{2} \sum_{a=1}^3 (w_a \gamma_a \gamma_0 - r_a \tilde{\gamma}_a)$$

A.4.5 Clifford algebra

$$\begin{aligned} \varepsilon_i \cdot \varepsilon_j + \varepsilon_j \cdot \varepsilon_i &= 2\eta_{ij} \\ \varepsilon_5 \cdot \varepsilon_5 &= -1 \end{aligned}$$

Lie Algebras

$$\begin{aligned} v(r, w) &= \frac{1}{2} (w^1 \varepsilon_0 \cdot \varepsilon_1 + w^2 \varepsilon_0 \cdot \varepsilon_2 + w^3 \varepsilon_0 \cdot \varepsilon_3 + r^3 \varepsilon_2 \cdot \varepsilon_1 + r^2 \varepsilon_1 \cdot \varepsilon_3 + r^1 \varepsilon_3 \cdot \varepsilon_2) \\ \vec{K}_1 &= v((1, 0, 0), (0, 0, 0)) = \frac{1}{2} \varepsilon_3 \cdot \varepsilon_2, \\ \vec{K}_2 &= v((0, 1, 0), (0, 0, 0)) = \frac{1}{2} \varepsilon_1 \cdot \varepsilon_3, \\ \vec{K}_3 &= v((0, 0, 1), (0, 0, 0)) = \frac{1}{2} \varepsilon_2 \cdot \varepsilon_1, \\ \vec{K}_4 &= v((0, 0, 0), (1, 0, 0)) = \frac{1}{2} \varepsilon_0 \cdot \varepsilon_1, \\ \vec{K}_5 &= v((0, 0, 0), (0, 1, 0)) = \frac{1}{2} \varepsilon_0 \cdot \varepsilon_2, \\ \vec{K}_6 &= v((0, 0, 0), (0, 0, 1)) = \frac{1}{2} \varepsilon_0 \cdot \varepsilon_3 \\ v(r, w) \cdot \varepsilon_0 &= \frac{1}{2} (w^1 \varepsilon_1 + w^2 \varepsilon_2 + w^3 \varepsilon_3 + r^3 \varepsilon_2 \cdot \varepsilon_1 \cdot \varepsilon_0 + r^2 \varepsilon_1 \cdot \varepsilon_3 \cdot \varepsilon_0 + r^1 \varepsilon_3 \cdot \varepsilon_2 \cdot \varepsilon_0) \\ \varepsilon_0 \cdot v(r, w) &= \frac{1}{2} (-w^1 \varepsilon_1 - w^2 \varepsilon_2 + w^3 \varepsilon_3 + r^3 \varepsilon_2 \cdot \varepsilon_1 \cdot \varepsilon_0 + r^2 \varepsilon_1 \cdot \varepsilon_3 \cdot \varepsilon_0 + r^1 \varepsilon_3 \cdot \varepsilon_2 \cdot \varepsilon_0) \\ v(r, w) \cdot \varepsilon_0 - \varepsilon_0 \cdot v(r, w) &= w \\ v(r, w) \cdot \varepsilon_0 + \varepsilon_0 \cdot v(r, w) &= 2v(r, 0) \cdot \varepsilon_0 \end{aligned}$$

In $Cl(3, 1)$:

$$\begin{aligned} v(r, w) \cdot \varepsilon_5 &= \varepsilon_5 \cdot v(r, w) = v(r, -w) \\ v(r', w') \cdot v(r, w) &= \frac{1}{4} (w^t w' - r^t r') + \frac{1}{2} v(-j(r) r' + j(w) w', -j(w) r' - j(r) w') - \frac{1}{4} (w^t r' + r^t w') \varepsilon_5 \\ [v(r, w), v(r', w')] &= v(j(r) r' - j(w) w', j(w) r' + j(r) w') \end{aligned}$$

$$\mathbf{Ad}_s v(x, y) = v(X, Y)$$

with

$$s = a + v(r, w) + b\varepsilon_5$$

$$X =$$

$$\left[(a+b)^2 - ab + \frac{1}{4} r^t r + (a+b)j(r) + \frac{1}{4} (j(w)j(r) - 2j(w)j(w) + 3j(r)j(r)) \right] x$$

$$+ \left[ab + \frac{1}{4} r^t r + (b-a)j(w) - \frac{1}{4} (3j(w)j(r) + 2j(r)j(w) - j(r)j(r)) \right] y$$

$$Y =$$

$$\left[\frac{1}{4} w^t w - ab + (a+b)j(w) + \frac{1}{4} (j(w)j(w) + 2j(w)j(r) + 3j(r)j(w)) \right] x$$

$$+ \left[(a-b)^2 - \frac{1}{4} w^t w - ab + (a-b)j(r) - \frac{1}{4} (3j(w)j(w) - j(r)j(w) - 2j(r)j(r)) \right] y$$

In $Cl(1, 3)$:

$$v(r, w) \cdot \varepsilon_5 = \varepsilon_5 \cdot v(r, w) = v(w, r)$$

$$v(r, w) \cdot v(r', w')$$

$$= \frac{1}{4} (w^t w' - r^t r') - \frac{1}{2} v(-j(r) r' + j(w) w', j(w) r' + j(r) w') - \frac{1}{4} (w^t r' + r^t w') \varepsilon_5$$

$$[v(r, w), v(r', w')] = -v(j(r) r' - j(w) w', j(w) r' + j(r) w')$$

Spin groups

$$s = a + v(r, w) + b\varepsilon_5$$

$$a^2 - b^2 = 1 + \frac{1}{4} (w^t w - r^t r)$$

$$ab = -\frac{1}{4} r^t w$$

$$\text{if } r = 0 \text{ then } a = \epsilon \sqrt{1 + \frac{1}{4} w^t w}; b = 0$$

if $w = 0$ then

$$r^t r \leq 4 : a = \epsilon \sqrt{1 - \frac{1}{4} r^t r}; b = 0$$

$$r^t r \geq 4 : b = \epsilon \sqrt{-1 + \frac{1}{4} r^t r}; a = 0$$

Product :

$$(a + v(r, w) + b\varepsilon_5)^{-1} = a - v(r, w) + b\varepsilon_5$$

$$s \cdot s' = a'' + v(r'', w'') + b'' \varepsilon_0 \cdot \varepsilon_1 \cdot \varepsilon_2 \cdot \varepsilon_3$$

with :

$$a'' = aa' - b'b + \frac{1}{4} (w^t w' - r^t r')$$

$$b'' = ab' + ba' - \frac{1}{4} (w^t r' + r^t w')$$

i) In $Cl(3, 1)$:

$$r'' = \frac{1}{2} (j(r) r' - j(w) w') + a' r + ar' - b' w - bw'$$

$$w'' = \frac{1}{2} (j(w) r' + j(r) w') + a' w + aw' + b' r + br'$$

$$(a + v(0, w)) \cdot (a' + v(0, w')) = aa' + \frac{1}{4} w^t w' + v(-\frac{1}{2} (j(w) w', a' w + aw'))$$

$$(a + v(r, 0)) \cdot (a' + v(r', 0)) = aa' - \frac{1}{4} r^t r' + v(\frac{1}{2} j(r) r' + (a' r + ar'), 0)$$

$$(a_w + v(0, w)) \cdot (a_r + v(r, 0)) = a_w a_r + v(a_w r, \frac{1}{2} j(w) r + a_r w) - \frac{1}{4} (w^t r) \varepsilon_5$$

ii) In $Cl(1, 3)$:

$$r'' = \frac{1}{2} (j(r) r' - j(w) w') + a' r + ar' + b' w + bw'$$

$$w'' = -\frac{1}{2} (j(w) r' + j(r) w') + a' w + aw' + b' r + br'$$