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STABILIZING CONTROL DESIGN FOR LIQUID PROPELLED ROCKET ENGINES

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ABSTRACT:

The recent development of reusable rocket launchers has led to new requirements for liquid propelled rocket engines as the reusable trajectory planning and landing of the engine stage leads to performing complex maneuvers. In this paper, a Port-Hamiltonian model is derived from the state space equations of a liquid propelled rocket engine. From this modeling, conditions are searched out to show the passivity of the system around a functioning point of the engine. This paper additionally proposes the use of contraction theory to build a suiting stabilizing controller for the system. The controller is illustrated with simulations.

1. INTRODUCTION

In the wake of reusable rocket engines, the requirements for maneuverability and knowledge on Liquid Propelled Rocket Engines (LPRE) have increased. Following the need of completing more complex maneuvers, the development of electrical actuators has left room for the design of new control laws. LPREs constitute complex non-linear systems that contain several unknown parameters, which are often approximated. As the modeling of LPREs becomes more precise [1] and evolves, the control law methods used require more flexibility. In a previous work [2], the stability of a transient model of a rocket engine has been tackled, however, the determination of a Lyapunov function for this type of system proves to be a difficult task. Moreover, such method for the determination of a Lyapunov function copes badly with the addition of new dynamics to the system. This can lead to the stability demonstration not to be viable anymore. A solution would be to develop an approach in which the dynamics added present guarantees for the system to remain stable, as well as an intrinsic Lyapunov function that can be added to the previous one. Following this observation, the choice has been made to introduce a Port-Hamiltonian [3] formulation of the non-linear systems constituting

an LPRE. Such a formulation allows for a closer analysis of the different elements and the simplified introduction of new dynamics. Port-Hamiltonian systems present an intrinsic energy function, that under certain conditions prevails as a Lyapunov function, guaranteeing the stability of the system around a chosen functioning point.

To perform complex maneuvers however, fixedpoint stability is not sufficient. In this paper, the determination of the appropriate thrust for the maneuver is not computed, and we only consider pre-set scenarios for the thrust. For such trajectory tracking, it is possible to define enough intermediary points to ensure the stability along the trajectory; however, such a method requires additional computation during the process. An alternative solution is to analyse the convergence of the trajectories themselves. Contraction theory [4] considers the flow of possible trajectories of a system and relies on proving the convergence of the flow. Recent progress in contracting control design [5] has led to the development of control methods using contraction theory. In [6], a contraction based controller is used for friction compensation in trajectory tracking, while in [7] the control of mechanical systems using contraction theory is depicted. In this paper, we use contraction theory on an affine input nonlinear system, for trajectory tracking objectives. A previous work [8] tackled the problem on a simplified model of an LPRE, while providing a good understanding of the contracting controller.

The paper is organized as follows. The model used is presented in section 2. The Port-Hamiltonian approach is then depicted in section 3 and the stabilizing control using contraction theory is described in section 4. Concluding remarks close the paper in section 5.

2. MODEL PRESENTATION

2.1. Principle of an LPRE

An LPRE is a complex system that provides a thrust to a body through a combustion. For a reusable engine, liquid propellers are used, in this case oxygen O2 and hydrogen H2 are chosen for the fuel and oxidizer species. Storage of the liquid species requires cryogenic temperatures, leading to low temperatures in the supply lines. The fuels are supplied using turbo-pumps situated between the liquid tanks and the combustion chambers (see figure 1). To power the turbo-pumps, a gas generator (GG) consumes a portion of the fuels into a secondary combustion chamber and the resulting hot gases are redirected into the turbine end of the turbo-pumps. Each pump is independent and the fraction of gas supplied to each pump is controlled using a valve (Z_{HG} on figure 1). Four additional valves (S_{VGH} , S_{VGO} , S_{VCH} , S_{VCO}) are placed to control the



Figure 1: LPRE model considered

mass flows in the different pipes (into the combustion chamber and into the gas generator). The modeling of the LPRE stands in the modeling of several physical phenomena, divided here in different domains:

- The mechanical domain is composed of the equation of the conservation of the angular moment of the turbo-pump axis of rotation.
- The **hydrodynamical** domain contains the evolution of the mass flows in the pipes, as well as the pressure increase in the cavities due to the hydraulic effect of the flows.
- The **chemical** domain represents the evolution of the pressures in the combustion chambers following the chemical reaction of the combustion.
- The **thermic** domain, containing the evolution of the heat exchanges between the different elements.

For simplification, several hypothesis are considered for the modeling of the LPRE:

The thermic phenomena are neglected in this approach, since the impact on the chamber temperature of the heat exchange with the cooling circuit is negligible. However, a static temperature increase is modeled by raising the temperature of the fluid flowing into the chamber.

- The different densities of the fluids are considered constant over time in the delivery circuit to the combustion chambers. The density is approximated in the hot gases lines using the perfect gas equation.
- The actuators dynamics are neglected, and the controller is set to provide a control surface to the system with no additional dynamics.
- The symmetry of the system is exploited, and by supposing that the Oxygen half of the system is well-regulated, it is possible to fix the mixture ratios in order to only consider the Hydrogen part of the system.
- The turbo-pumps' dynamics are computed using manufacturer data, under the form of polynomial functions. The functions' validity domain is supposed to be the study domain.

2.2. State equations of the model

In this paragraph, the state equations of the system are given following the modeling in [2]. The equations are expressed for the Hydrogen half of the model, with the complete equations given in 2.3.

2.2.1.Turbo-pumps

The dynamics of the turbo-pump rotational speed $\omega_{\rm H}$ is obtained using the conservation of the kinetic moment:

$$\dot{\omega}_H = \frac{1}{J_H} (T_{mH} - T_{rH})$$
 Eq.1

where J_H is the angular inertia of the turbine, T_{mH} and T_{rH} are the motor and resistive torques respectively. The modeling of the turbo-pump dynamics requires to determine the resisting torque T_{rH} applied by the fluid in the line and the motor torque T_{mH} generated by the turbine. The resistive torque is expressed following

$$T_{rH} = -\left|\frac{a_{cH}}{a_{H}}q_{H}^{2} + b_{cH}\omega_{H}q_{H} + c_{cH}\rho_{H}\omega_{H}^{2}\right|, \quad \text{Eq.2}$$

where a_{cH} , b_{cH} , c_{cH} are manufacturer parameters for the turbo-pump, !H is the rotational speed of the pump, ρ_H is the density of the fluid and q_H is the mass flow through the pump with $q_H = q_{ccH} + q_{ggH}(q_{ccH}, q_{ggH}$ are the mass-flows in the CC and GG respectively). T_{mH} is then expressed considering the evolution of the hot gases produced by the GG,

$$N_R = \frac{\omega_H Rad_T}{\sqrt{\gamma R\Theta}}$$
 Eq.3

$$W = q_{outH} \frac{\sqrt{\gamma RT}}{\gamma - 1} \left(1 - \pi_T^{\frac{1 - \gamma}{\gamma}} \right) Rad_T \qquad \text{Eq.4}$$

 $ST = (a_1 + N_R(A_2 + N_Ra_3) + \pi_T(a_4 + \pi_Ta_5) + a_6\pi_TN_R + a_7ln\pi_T + a_8lnN_R). Corr$ Eq.5

$$T_{mH} = ST.W$$
 Eq.6

$$q_{outH} = \frac{p_{tuH}A_{tuH}}{\gamma} \sqrt{\frac{R\Theta}{\gamma} (\frac{2}{\gamma+1})^{-\frac{\gamma+1}{\gamma-1}}} \quad \text{Eq.7}$$

Where q_{outH} is the mass flow through the turbine, $a_1, ..., a_8$ are specific coefficients for the turbo-pump and N_R is a reduced rotational speed.

The chamber pressure's evolution follows the energy conservation equation:

$$\frac{d}{dx}\left(\frac{pV}{\gamma-1}\right) = (h_{in}q_{in} - h_{out}q_{out}), \qquad \text{Eq.8}$$

where *h* is the enthalpy of the species consumed and created. The thermodynamic parameter γ is assumed constant during the evolution of the system. The enthalpy of the system is composed of the formation enthalpy h_f , the kinetic energy e_c and the loss due to the heating of the fluid $C_v L_v \Theta$,

$$h = h_f + e_c + C_v L_v \Theta \qquad \qquad \text{Eq.9}$$

where $C_v L_v$ is the vaporisation heat of the fluid. The equation Eq.8 can be written as follows,

$$\dot{p}_{cc} = (k_{1cc} + k_{3cc} - k_{4cc}\Theta_{cc})(q_{ccH} + q_{cc0}) - k_{2cc}p_{cc}$$

Eq.10

Where k_{1cc} to k_{4cc} are obtained by first order Taylor expansion of the equation Eq.8, Θ_{cc} is the temperature in the CC and q_{ccH} , q_{cc0} are the mass flows of the species entering the CC. With the use of the mixture ratio, the equation is

$$\dot{p}_{cc} = (k_{1cc} + k_{3cc} - k_{4cc}\Theta_{cc})(MR + 1)q_{ccH} - k_{2cc}p_{cc} \qquad \text{Eq.11}$$

The evolution of the gas generator pressure p_{gg} is modeled with Eq.8, using the output mass flows rather than the internal pressure.

$$\dot{p}_{gg} = (k_{1gg} + k_{3gg} - k_{4gg}\Theta_{gg})(q_{ggH} + q_{gg0}) - k_{2gg}(q_{tuH} + q_{tu0})$$
Eq.12

Where k_{1gg} to k_{4gg} are also obtained by first order Taylor expansion of Eq.8, Θ_{gg} is the temperature, q_{ggH}, q_{gg0} are the entering mass flows in and q_{tuH}, q_{tu0} are the output mass flows of the GG. The temperature Θ of the combustion chambers is approximated by a function f_{Θ} of the mixture ration, i. e. $\Theta = f_{\Theta}(MR)$.

Turbine pressures

0

The evolution of the turbine pressures follows equation Eq.8, where no chemical reaction occurs.

$$\dot{p}_{tuH} = k_{1tuH}\Theta_{gg}q_{tuH} - k_{2tuH}\sqrt{\Theta_{gg}}p_{tuH}$$
 Eq.13

Where k_{1tuH}, k_{2tuH} are obtained as the other coefficients.

Mass-flows

The dynamics of the mass flows are derived from the equation of momentum-conservation applied to the lines,

$$I \dot{q} = p_{in} - p_{out} - Res_{eq}q|q|.$$
 Eq.14

Where *I* is the inertia of the line, of the form $I = \frac{A}{L}$, *A* the area and *L* the length of the line, p_{in} , p_{out} are the upstream and downstream pressures, Res_{eq} is the equivalent resistance coefficient associated with pressure drops and *q* is the mass flow considered. Pressure drops are due to the valves S_{VGH} , S_{VCO} , S_{VCH} , S_{VCO} , injectors and friction in the lines.

Pressure drop due to the valves is explicited by

$$\Delta p = \frac{1}{2\rho A_v^2} q^2$$

$$Res_v = \frac{1}{2\rho A_v^2}$$
Eq.15

Where A_{ν} is the surface associated with the opening angle of the valve, ρ is the density of the fluid in the line. Pressure drops in the exhaust of the GG express as

$$\Delta_{ptuH-gg} = \frac{Khy_{tuH}R_{gg}\Theta_{gg}}{p_{gg}}q_{tuH}^{2},$$

$$\Delta_{ptuO-gg} = \frac{ZresR_{gg}\Theta_{gg}}{p_{gg}}q_{tuO}^{2},$$

Eq.16

Where Z_{res} , Khy_{tuH} are equivalent resistance coefficients for the valve and the line. Due to the burned gases, ρ is not supposed constant in the turbine lines (q_{tuH}). Using the perfect gas formula, its variation is expressed as a function of p_{gg} . Pressure drops associated with friction in the lines also follows equation Eq.15.

The pressure increase due to the pumps follows the equation

$$p_{spH} = p_{epH} + \frac{a_{pH}}{\rho_H} q_H^2 + b_{pH} \omega_H q_H + c_{pH} \rho_H \omega_H^2$$
. Eq.17

Where p_{spH} is the output pressure of the pump, p_{epH} is the output pressure of the tank. The other parameters are similar to Eq.2.

2.3. State equations of the global system:

$$\begin{split} \dot{p}_{cc} &= (k_{1cc} + k_{3cc} - k_{4cc}\Theta_{cc})(q_{ccH} + q_{ccO}) - k_{2cc}p_{cc} \\ \dot{p}_{gg} &= (k_{1gg} + k_{3gg} - k_{4gg}\Theta_{gg})(q_{ggH} + q_{ggO}) - k_{2gg}(q_{tuH} + q_{tuO}) \\ \dot{p}_{tuH} &= k_{1tuH}\Theta_{gg}q_{tuH} - k_{2tuH}\sqrt{\Theta_{gg}}p_{tuH} \end{split}$$

$$\begin{split} \dot{p}_{tu0} &= k_{1tu0} \Theta_{gg} q_{tu0} - k_{2tu0} \sqrt{\Theta_{gg} p_{tu0}} \\ \dot{q}_{ccH} &= \frac{A_{VCH}}{L_{CH}} \left(p_{spH} - p_{cc} - \left(\frac{1}{2\rho A_{VCH}^2} + Res_{ccH} \right) q_{ccH}^2 \right) \\ \dot{q}_{cc0} &= \frac{A_{VCO}}{L_{c0}} \left(p_{sp0} - p_{cc} - \left(\frac{1}{2\rho A_{VC0}^2} + Res_{cc0} \right) q_{cc0}^2 \right) \end{split}$$

q_{ggH}

$$=\frac{A_{VGH}}{L_{GH}}\left(p_{spH}-p_{gg}-\left(\frac{1}{2\rho A_{VGH}^{2}}+Res_{ggH}\right)q_{ggH}^{2}\right)$$

q_{gg0}

$$= \frac{A_{VGO}}{L_{GO}} \left(p_{spo} - p_{gg} - \left(\frac{1}{2\rho A_{VGO}^2} + Res_{ggO} \right) q_{ggO}^2 \right)$$

$$\dot{q}_{tuH} = \frac{1}{I_{tuH}} \left(p_{gg} - p_{tuH} - \frac{Khy_{tuH}R_{gg}\Theta_{gg}}{p_{gg}} q_{tuH}^2 \right)$$

$$\dot{q}_{tuO} = \frac{1}{I_{tuO}} \left(p_{gg} - p_{tuO} - \frac{ZresR_{gg}\Theta_{gg}}{p_{gg}} q_{tuO}^2 \right)$$

$$\dot{\omega}_H = \frac{1}{J_H} (T_{mH} - T_{rH})$$

$$\dot{\omega}_O = \frac{1}{I_O} (T_{mO} - T_{rO}) \qquad \text{Eq.18}$$

3. PORT-HAMILTONIAN FORMULATION OF AN LPRE

The LPRE's state's evolution requires complex and highly non-linear equations, that are not adequate for an easy design of a control law. A modeling solution proposed here is to make use of the Port-Hamiltonian systems framework, and reformulate the LPRE model established under this formulation.

3.1. Port-Hamiltonian Modeling

Port-Hamiltonian systems, described in [9], [10] and [3], used for the modeling of mechanical systems, have been extended to various physical domains as hydrodynamics in [11] and chemical processes in [12], [13]. Port-Hamiltonian modeling makes use of physical invariants of a system, such as the total energy, to highlight the energy transfers between the different elements composing a system. The complete model is broken down to the level a simple elements that each represent a simple equation. They are represented as independent subsystems, with two links: the effort *e* and the flow *f*. Both the effort and flow represent coupled physical quantities (e.g. pressure and flow). Elements can be classified into three distinct functions:

- Energy-storing, denoted by *e_s*, *f_s*, are elements that store energy (kinetic, potential, ...).
- Energy-dissipating, denoted by e_r, f_r dissipate the energy, using friction for example.
- Energy-routing, denoted by e_t , f_t transfers energy, with no storage nor loss. These elements are by construction perfect

transformers, therefore are seldom used in real systems.

Additionally, interactions with the exterior environment is denoted by e_p , f_p . The product $e^T f$ is the power resulting of the flow and effort. Reconstructing an equation consists in regrouping the efforts to build the equation for the flow.

An important property resides in the energy balance of such a system:

$$e_s^T f_s + e_r^T f_r + e_p^T f_p = 0 Eq.19$$



Figure 2: Physical domains of an LPRE

By definition, the dissipative elements yield $e_r^T f_r \le 0$, therefore

$$e_s^T f_s \le e_p^T f_p$$
 Eq.20

Equation Eq.20 is a classical equation for the passivity of a system where the energetic input is e_p and the output y is f_p . Let the function H be defined by:

$$\frac{dH}{dt} = e_s^T f_s = \frac{\partial H}{\partial x} \dot{x}$$
. Eq.21

Such function *H* is called the Hamiltonian of the system that represents the total energy of the system. For H(x) > 0 for all *x*, *H* is a candidate storage function for the passivity of the system. From these definitions, a system under PH form is expressed

$$\dot{x} = (J(x) - R(x))\frac{\partial H}{\partial x} + G(x)u$$
 Eq.22

Where the matrix J(x) expresses the energy storage terms, $J = -J^T$ and R(x) expresses the energy dissipation $R(x) \ge 0$.

3.2. LPRE formulation

In this section, a reformulation under Port-Hamiltonian form is proposed for the system in Eq.18. The different physical domains as expressed in **2.2**. can be connected using the different energy transfers between the domains. On figure 2 the energy transfers are represented between the different domains. While each domain can be represented on its own, the interconnection between the domains has to be studied to provide a complete PH system. In previous works [14], the PH framework has been extended to cover many physical domains. Moreover, the interconnection of several physical domains has been proposed in [15].

3.2.1.Expression of subsystems under PH formulation

In this subsection, the individual subsystems are expressed under Port-Hamiltonian form, referring to previous works in the literature. The hydraulic part of the system is composed of the evolution of the mass flows and pressures. The coupled flow and efforts (e, f) are $(q, I\dot{q})$ and $(p, k_p \dot{p})$, which both have the dimensions of a pressure/mass flow couple. From Eq.14, the different elements that drive the mass flow behavior are the resistive inertia of the pipe and the upstream and downstream pressures, respectively a resistive effort, the energetic input and the pressure effort (stoarge effort). This reasoning leads to the structure:

$$\begin{bmatrix} I\dot{q} \\ k_p\dot{p} \end{bmatrix} = \begin{bmatrix} -Rq & -1 \\ 1 & -R_p \end{bmatrix} \begin{bmatrix} q \\ p \end{bmatrix} + \begin{bmatrix} p_{in} \\ 0 \end{bmatrix}$$
Eq.23

where R_q and R_p are the resistive terms associated to the evolution of the mass flow and p_{in} is the input pressure of the pipe. This equation is applicable to the different mass flows of the LPRE, using the expressions of the input and output pressures of the flow. The equations of the pressures in the combustion chambers p_{cc} and p_{gg} allow for the chemical domain to be expressed under the same form provided the expression of R_p and k_p . In the case of the turbine pressure p_{tu} , no chemical domain is involved as there is no combustion.

The mechanical subsystem is expressed using the port variables for a torque and a rotational speed

$$[J\dot{\omega}] = [-R_{-}\omega][\omega] + [T_{mot}] \qquad \text{Eq.24}$$

3.2.2. Interconnection of subsystems

Interconnection of several Port-Hamiltonian systems has been described in [3] [16]. Two Dirac structures D_A and D_B are considered on product spaces respectively $F_1 \times F_2$ and $F_2 \times F_3$, where F_2 is the space of shared flow variables. The interconnection is realized by equating the flows and efforts on the shared space of variables.

The interconnection of the different hydraulic subsystems can be performed directly as the shared space of variables covers both the mass flows and pressures. However, interconnection between mechanical and hydraulic spaces requires to find a shared space of variables. Similarly to the work performed in [15], the method used consists in finding a function to express the torque and the pressure under a common equation leading to the creation of the shared space. In this paper, the pressure and torque have non-reversible equations. A proposition to solve the interconnection is to use an upper value of the torque, and considering the other terms as resistive terms. Let $\alpha_q = \frac{b_p \rho}{a_c}$. Let also $\bar{p} = b_p q \omega$ be an upper value for the pressure increase of the pump (see Eq.17), and $\underline{T} = a_c (q_{cc} + q_{gg})^2$ be a lower value of the resistive torque generated by the fluids on the pump (see Eq.2). By multiplying the equation of the resistive torque by α_q , the evolution equations reads:

$$\begin{aligned} \alpha_q J \dot{\omega} &= -b_p q^2 - \alpha_q (b_c q + c_c \omega) \omega + \alpha_q T_m \\ I \dot{q} &= b_p q \omega - \left(Z_{eq} q - \frac{a_p}{\rho} q - \frac{c_p \omega^2}{q} \right) q + p_{ep}. \end{aligned} \quad \text{Eq.25} \\ \text{This yields the PHS interconnection:} \end{aligned}$$

$$\begin{bmatrix} \alpha_q J \dot{\omega} \\ I \dot{q} \end{bmatrix} = \begin{bmatrix} -R_{\omega} & -b_p q \\ b_n q & -R_q \end{bmatrix} \begin{bmatrix} \omega \\ q \end{bmatrix} + \begin{bmatrix} \alpha_q T_m \\ p_{ep} \end{bmatrix} \quad \text{Eq.26}$$

where $R_q = \left(Z_{eq}q - \frac{a_p}{\rho}q - \frac{c_p\omega^2}{q}\right)$ and $R_\omega = \alpha_q(b_cq + c_c\omega)$ have to be positive to satisfy the upper and lower value hypothesis. Here two conditions are derived for the validity of the PH formulation of the system, which can be used as validation test for the stability of the system.

Similarly, by posing $\alpha_t = \alpha_q ST(\omega)w_{tu}$, where $w_{tu} = \frac{w}{p_{tu}}$ it is possible to obtain the interconnection for the turbine:

$$\begin{bmatrix} \alpha_q J \dot{\omega} \\ k_{ptu} \dot{p}_{tu} \end{bmatrix} = \begin{bmatrix} -R_{\omega} & \alpha_t \\ -\alpha_t & -R_{ptu} \end{bmatrix} \begin{bmatrix} \omega \\ p_{tu} \end{bmatrix}, \qquad \text{Eq.27}$$

Where R_{ptu} is computed to match the original equation Eq.13.

The complete Hamiltonian function expresses as :

$$H(x) = \frac{1}{2}\alpha_q J \omega^2 + \frac{1}{2}I_{cc}q_{cc}^2 + \frac{1}{2}I_{gg}q_{gg}^2 + \frac{1}{2}I_{tu}q_{tu}^2 + \frac{1}{2}k_{pcc}p_{cc}^2 + \frac{1}{2}k_{pgg}p_{gg}^2 \frac{1}{2}k_{ptu}p_{tu}^2, \quad \text{Eq.28}$$

The hydrogen subsystem under Port-Hamiltonian form is given in annex.

3.3. Derivation of passivity conditions

Denoting that the Hamiltonian function given in Eq.28 is positive, *H* is a candidate Lyapunov function, if the system verifies the conditions for the PHS structure. The construction of the interconnection focuses on the structure of the energy transfers, yielding $J = -J^T$. However, to justify the dissipativity of the system it is necessary to verify that the terms R_{ω} , R_q , R_{ptu} are positive and do not lead to the creation of energy. With the three conditions fulfilled, the system is passive with respect to the energetic input p_ep .

The condition for the passivity relies on the construction of the model. During the model construction, it is possible to optimize the passivity analysis when given a choice between different representations. The most adapted form for passivity analysis is the overestimation of a parameter, and using correction terms that consume energy.

4. STABILIZATION AND CONTRACTION THEORY

In the section 3., a method to prove stability around a functioning point of the system has been given. However, in the case of a maneuver to be performed, it is necessary to ensure stability along the trajectory and between the known functioning points. To address this problem, two solutions are available:

- Study the stability around each functioning point on the trajectory even by adding rallying points along the trajectory.
 - Study the stability around the trajectory

The use of contraction theory allows the latter option to be chosen. The idea is to generate a trajectory that corresponds to the maneuver to be performed, and stabilize the system around this reference trajectory. In this section, a method for the application of contraction theory is presented.

4.1. Contraction theory

Let a system of the form

$$\dot{x} = f(x, t) \qquad \qquad \mathsf{Eq.29}$$

We define a virtual displacement δx [17],

$$\delta \dot{x} = \frac{\partial f}{\partial x}(x,t)\delta x$$
 Eq.30

A trajectory in the flow field $\dot{x} = f(x, t)$ is a possible solution of the system Eq.29. The squared distance between two neighbouring trajectories is defined in [17] ad $\delta x^T \delta x$ and its rate of change as

$$\frac{d}{dt}(\delta x^T \delta x) = 2 \,\delta x^T \delta \dot{x} = 2 \delta x^T \frac{\partial f}{\partial x} \delta x \qquad \text{Eq.31}$$

Definition 1. A contraction region [17], is a region of the state if the Jacobian $\frac{\partial f}{\partial x}$ is uniformly negative definite in that region,

$$\exists \beta > 0, \forall x, \forall t \ge 0, \frac{1}{2} \left(\frac{\partial f}{\partial x} + \frac{\partial f^T}{\partial x} \right) \le -\beta I < 0 \quad \text{Eq.32}$$

Theorem 1. [17] Given the system equations $\dot{x} = f(x, t)$, any trajectory which starts in a ball of constant radius centered about a given trajectory and contained at all times in a contraction region, remains in that ball and converges exponentially to this trajectory.

This sufficient condition for contraction is seldom fulfilled with a system, therefore, we introduce a sufficient condition introduced in [18]

If there exists a constant matrix $P \in \mathbb{R}^{n \times n}$, $P^T = P$, such that:

$$P\frac{\partial f}{\partial x} + \frac{\partial f^T}{\partial x}P < -\beta I. \qquad \text{Eq.33}$$

Then the system defines a contraction on the

region. Moreover, for nonlinear systems, it is possible to decompose the system equation:

$$\dot{x} = f(x,t) = Ax + M\phi(y,t), y = Lx$$
 Eq.34

where the linear and nonlinear parts of the system are separated, y is a vector composed of the states used in the nonlinear part of the equation, ϕ is a C^1 mapping and A, L, M are matrices of appropriate dimensions.

The condition Eq.33 reads:

$$P\left(A + M\frac{\partial\phi}{\partial y}(y,t)L\right) + \left(A + M\frac{\partial\phi}{\partial y}(y,t)L\right)^{T}P < -\beta I$$

Eq.35

Remark 1. Such expression of the system pairs efficiently with a PH system, as the matrix *J* used for the energy storing is often taken as close as the identity as possible, resulting in a good candidate for the matrix *A* in Eq.34.

In the case of the system studied not being contracting, it is possible to define a control to force the contraction of the closed-loop system. In [5], such a design is proposed, using an LMI solving approach to answer the problem. The benefit of using an LMI approach is the reduction of computation time, which is appropriate when dealing with embarked systems.

Let a system under the form Eq.34. Assume that for all (y, t).

$$0 \le \frac{\partial \phi}{\partial y}(y,t) + \frac{\partial \phi^T}{\partial y}(y,t)$$
 Eq.36

Then, for a controlled system such that,

$$\dot{x} = f(x,t) = Ax + M\phi(y,t) + Bu, y = Lx \qquad \text{Eq.37}$$

 $u = Kx + N\phi(y, t)$ is a contracting controller if we can find three matrices *Z*, *W*, *N* such that:

$$K = ZW^{-1}$$

 $(AW + BZ) + (AW + BZ)^T \le 0, W > 0$
 $WL^T = -(M + BN).$ Eq.38

4.2. Design of a contracting controller for an LPRE

In this section, the method for designing a contracting controller is depicted for the application of the LPRE. The controller must regulate the mass flows that enter the different combustion chambers. However, rather than a energy input to the system, such interaction with the system is a friction-based controller, where the controller action performs as a friction coefficient. The term associated to the friction exercised by the controller is under the form:

$$\frac{1}{2\rho A_v^2}q^2.$$
 Eq.39

For simplification, the equivalent resistance of the

valve is used to represent the controller:

Denoting $u = Z_{eq}$ in Eq.37, the matrix *B* is time-varying;

$$B(x) = [0 \ 0 - q^2 \ 0 \dots 0]^T$$
, Eq.41

which leads to the LMI Eq.38 being state dependent. This is a contradiction with the hypothesis Eq.33, where we suppose that the matrix $P = W^{-1}$ is time-invariant. Note that a time-variant $\mathcal{P}(t)$ matrix consists in an other formulation of the problem [5].

The method proposed here consists in fixing the LMI parameter $P = P_0$, which locks the value $W = W_0 = P_0^{-1}$. The value P_0 is called the constant metric for the system. Additionally, let the introduction of the notion of virtual systems [19].

Let a system

$$\Sigma: \dot{x} = f(x) + g(x)u \qquad \qquad \text{Eq.42}$$

And introduce the associated virtual system Σ_v with associated virtual state x_v and virtual control u_v ,

$$\Sigma_{v}: \dot{x}_{v} = f(x_{v}) + g(x)u_{v}$$
 Eq.43

Here the state *x* is considered a constant value regarding the variations of virtual state x_v . Now suppose a contracting control $u_v(x, x_v, t)$ for the virtual system Eq.43, for all *x* with respect to x_v . Then the controller $u_v(x, x, t)$ is contracting for the real system Eq.42, (see [19], Theorem 2).

4.3. Affine controlled system

In the event of an affine controlled system of the form Eq.42, the decomposition looking forward to the LMI resolution Eq.37 yields:

$$\dot{x} = f(x,t) + g(x)u = Ax + M\phi(y,t) + B(x)u$$

$$y = Lx$$
Eq.44

Where x is considered constant regarding the variations of x_v . In particular, for $x = x_0$, we are able to solve the LMI:

$$\begin{split} K_0 &= Z_0 W_0^{-1} \\ (AW_0 + B(x_0) Z_0) + (AW_0 + B(x_0) Z_0)^T \leq 0, W_0 > 0 \\ W_0 L^T &= -(M + B(x_0) N_0). \end{split}$$
 Eq.45

 $P_0 = W_0^{-1}$ is a constant metric of contraction for the virtual system Eq.44 for $x = x_0$. It is possible to conserve such constant metric of contraction for different values of *x*. Let $\hat{B}(x)$ be the generalized left inverse of B(x), then posing $Z(x) = \hat{B}(x)B(x_0)Z_0$ and $N(x) = \hat{B}(x)B(x_0)N_0$ yields

$$K_{0} = Z_{0}W_{0}^{-1} (AW_{0} + B(x)\hat{B}(x)B(x_{0})N_{0}) + (AW_{0} + B(x)\hat{B}(x)B(x_{0})N_{0})^{T} \le 0, W_{0} > 0 W_{0}L^{T} = -(M + B(x)\hat{B}(x)B(x_{0})N_{0}).$$
Eq.46

which by replacing the terms gives back Eq.45. Therefore, then $Z(x) = \hat{B}(x)B(x_0)Z_0$ and $N(x) = \hat{B}(x)B(x_0)N_0$ are solutions of the LMI problem, with a constant $P_0 = W_0^{-1}$. To design this controller, no state-dependent LMI is computed. The LMI is computed once for the chosen state x_0 , for example a particular functioning point, and then the control action can be deduced from K_0, N_0 .

For the LPRE system, it can be shown that the control action $u = K_0 \frac{q_{gg0}^2}{q_{gg}^2} x + N_0 \frac{q_{gg0}^2}{q_{gg}^2} \phi(y,t)$ is a contractive controller in the sense of [5].

4.4. Results

In this subsection, application of contraction theory to the model described in section 2.3. and reformulated in 3.2. is presented. The model used is valid for the range 90% - 150% of the states' nominal value x_0 . The different results and graphs are given in normalized form. The normalized form is $x_{norm} = \frac{x}{x_0}$. This paper focuses on the stability of the engine and the provided controller. The scenario proposed are computed beforehand, and are designed to cover different situations, and highlight the limits of controllers.



Figure 3: Tracking performances of the states of an LPRE for a simple reference using a contracting controller



Figure 4: Evolution of the actuators of an LPRE for a simple reference using a contracting controller

To illustrate the tracking performance of the controller, a first non-realistic simple trajectory is applied to the system. The objective of this trajectory is to evaluate the tracking performance of the controller. On figure 3, the evolution of both mass flows incoming in the GG is represented, as well as the main combustion chamber pressure evolution. Figure 4 completes the illustration by depicting the evolution of the surface actuators during the same scenario.

The next illustration of the controller proposed makes use of a realistic reference for a maneuver. For a typical landing scenario, the returning module of the rocket handles three phases *[20]*. A first reentry burn is performed, for the safety of the returning module. Due to air friction, it is necessary to reduce speed in the first phase of the landing. The second phase of the scenario consists in atmospheric gliding using the fins of the craft, while the rocket engine is kept at a lower thrust value. Finally, the final burn drastically reduces speed for a safe landing. The challenge of such a scenario relies on the amplitude of the changes in the



Figure 5: Evolution of the states during a re-entry scenario using a contracting controller



Figure 6: Evolution of the actuators during a re-entry scenario using a contracting controller

trajectory of the states. Figure 5 and 6 represent the contracting controller action following such a reference. Noise has been added on the chamber temperature to represent combustion fluctuations. As illustrated on figure 5, the controller provides a good precision and speed for the scenario, and the actuators' evolution on figure 6 remains feasible. The controller's contracting property is illustrated on figure 7. In this scenario, the reference is reaching the validity limit of the model. Additionally, a noised signal has been added on the chamber temperature to illustrate uncertainty on the parameter Θ_{cc} . Two controllers are proposed to illustrate the behavior. First, the contracting controller design provides a stable trajectory. The second controller is a PID that is designed around the functioning point x_0 . The second controller fails to provide stability to the system and the simulation is unable to complete, due to divergence of states (see q_{cc} on figure 6). In this scenario, the contracting controller fulfills its role as it manages to maintain the system in the contraction region, and thus ensuring the stability of the LPRE.



Figure 7: Evolution of the states in an unstable scenario for a contracting controller and a PID controller



Figure 8: Evolution of the pressure for different initial conditions using a contracting controller

A property of contraction theory is to forget initial conditions of a system. On figure 8 are represented different initial conditions for a system regulated around a nominal trajectory $p_{cc} = 1$). All trajectories converge to the same reference trajectory, validating the notion of forgetting initial conditions.

5. CONCLUSION

This paper summarizes the work performed on the stability of an LPRE. Two approaches were considered, for fixed-point stability and state-transition stability.

Using Port-Hamiltonian systems framework, it is possible to reformulate a complex non linear model under a form optimized for passivity analysis. Using the reformulation, a Hamiltonian function of the LPRE has been deduced, additionally proving the passivity of the system. The framework also enables flexibility in the design, as the addition of elements is simplified. As an example, this paper provides the addition of actuator dynamics. Port-Hamiltonian framework opens for new control methods and the derived conditions for the passivity enables the user to verify the controller proposed.

Contraction theory is a powerful alternative to feedback controller, ensuring the stability of the system over a range of trajectories defined by the user. The analytical formulation of the controller is suitable for embedded systems, and provides good performances in terms of speed and precision. Moreover, the reference provided by the user is subject to optimization, leading to different solutions for the problem.

Future work includes the design of a passivity base controller.

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ANNEX: PORT-HAMILTONIAN FORMULATION OF THE LPRE

$$J(X) = \begin{cases} \alpha_{q}J\dot{\omega} \\ I_{cc}\dot{q}_{cc} \\ I_{gg}\dot{q}_{gg} \\ I_{tu}\kappa\dot{q}_{tu} \\ k_{pcc}\dot{p}_{cc} \\ k_{pgg}\dot{p}_{gg} \\ k_{ptu}\dot{p}_{tu} \end{cases} = [J(X) - R(X)] \begin{bmatrix} \omega \\ q_{cc} \\ q_{gg} \\ q_{tu} \\ p_{cc} \\ p_{gg} \\ p_{tu} \end{bmatrix} + \begin{bmatrix} 0 \\ 0 \\ 0 \\ 0 \\ 0 \\ 0 \end{bmatrix}$$
$$J(X) = \begin{bmatrix} 0 & -\lambda & -\lambda & 0 & 0 & 0 & \alpha_t \\ \lambda & 0 & 0 & 0 & -1 & 0 & 0 \\ \lambda & 0 & 0 & 0 & 0 & -1 & 0 \\ 0 & 0 & 0 & 0 & 0 & -1 & 0 \\ \lambda & 0 & 0 & 0 & 0 & -1 & 0 \\ 0 & 0 & 0 & 0 & 0 & \kappa_{tu} & -\kappa_{tu} \\ 0 & 1 & 0 & 0 & 0 & 0 & 0 \\ \alpha_t & 0 & 0 & \kappa_{tu} & 0 & 0 & 0 \end{bmatrix}$$
$$R(X) = diag[R_{\omega}R_{qcc}R_{qgg}R_{qtu}R_{pcc}R_{pgg}\beta_t]$$

Where $\lambda = b_p (q_{cc} + q_{gg})$. Expression of the resistive terms: $R_\omega = \alpha_q (b_c (q_{cc} + q_{gg}) - c_c \rho \omega)$ $R_{qcc} = Z_{eqcc} q_{cc} - \frac{a_p}{\rho} (q_{cc} + q_{gg}) - \frac{c_p \rho \omega^2}{q_{cc}}$ $R_{qgg} = Z_{eqgg} q_{gg} - \frac{a_p}{\rho} (q_{cc} + q_{gg}) - \frac{c_p \rho \omega^2}{q_{gg}}$ $R_{qtu} = \kappa_{tu} Z_{hg} q_{tu}$ $R_{pcc} = k_{2cc} k_{pcc}$ $R_{pgg} = 0$ $R_{ptu} = \beta_t$ Expression of the constant scaling terms:

$$\kappa_{tu} = k_{2gg}k_{gg}$$

$$k_{ptu} = \frac{\kappa_{tu}}{k_{1tu}}$$

$$k_{pcc} = \frac{1}{\frac{1}{k_{1cc} + k_{3cc} - \Theta_{cc}k_{4cc}}}$$

$$k_{pgg} = \frac{1}{\frac{1}{k_{1gg} + k_{3gg} - \Theta_{gg}k_{4gg}}}$$

Expression of the interconnection terms:

$$\alpha_q = \frac{b_p \rho}{a_c}$$
$$\alpha_t = \alpha_q ST(\omega) w_{tu}$$