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Marie Pomarede, Elisabeth Longatte-Lacazedieu, Jean-François Sigrist. Benchmark of numerical codes for coupled CSD/CFD computations on an elementary vortex induced vibration problem. ASME 2009 Pressure Vessels and Piping Conference, Jul 2009, Prague, Czech Republic. pp.537-546, <10.1115/PVP2009-77028>. <hal-04642661>

HAL Id: hal-04642661

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Submitted on 15 Feb 2025

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PVP2009-77028

**BENCHMARK OF NUMERICAL CODES FOR COUPLED CSD/CFD COMPUTATIONS ON AN
ELEMENTARY VORTEX INDUCED VIBRATION PROBLEM**

Marie POMAREDE

Service Technique et Scientifique
DCNS Propulsion
44620 LA MONTAGNE, France

Elisabeth LONGATTE

Département Mécanique des Fluides,
Energies et Environnement
EDF R&D
78401 CHATOU Cedex, France

Jean-François SIGRIST

Service Technique et Scientifique
DCNS Propulsion
44620 LA MONTAGNE, France
jean-francois.sigrist@dcnsgroup.com

ABSTRACT

Numerical simulation of vortex-induced-vibrations (VIV) of an elastically supported rigid circular cylinder in a fluid cross-flow has been thoroughly studied over the past years, both from the experimental and numerical points of view, because of its theoretical and practical interest in the understanding of flow-induced vibrations problems. In this context, the present paper aims at exposing a numerical study based on a coupled fluid-structure simulation, compared with previously published studies [34], [36]. The computational procedure relies on a partitioned method ensuring the coupling between fluid and structure solvers. The fluid solver involves a moving mesh formulation for simulation of the interface motion. Energy exchanges between both systems are ensured through convenient coupling schemes. The present study is devoted to a low Reynolds number configuration ($Re = 100$). Cylinder motion magnitude, hydrodynamic forces, oscillation frequency and fluid vortex shedding modes are investigated with the intention to observe the “lock-in” phenomenon. These numerical simulations are proposed for code validation purposes prior to industrial applications to tube bundle configurations [4].

INTRODUCTION

Studies on Vortex Induced Vibrations (VIV) encountered a large interest over the past decades because this is a phenomenon appearing in many fields: riser tubes on off-shore oil rigs [28], [29], steam exchangers [23] or bridge cables are VIV-prone. It may cause severe damages when vibrations are repetitive and/or large-amplitude phenomena. VIV are still a research challenge because of the lack of comprehension of certain observable facts, and because of long calculation times and many other requirements of numerical simulations. Fundamental studies on this subject can be found in [5], [8] and [35]. Experimental studies are proposed by many authors: [12], [13] for instance. Numerical works like [15], [27], [32], [36] and [38] give solidly-based results for the conduction of VIV numerical simulations.

In the present work, we are concerned with the numerical simulation of the transverse oscillations of an elastically

mounted rigid cylinder at low Reynolds number ($Re = 100$) without damping. Relative studies to these characteristics which constitute the main reference calculations for the present work are [34], [36]. We are leading numerical simulations using a fluid-structure coupling, performed thanks to a CFD code: *Code_Saturne*, which is free software developed by EDF R&D, based on a collocated finite volume approach. Momentum equations are solved considering an explicit mass flux. Velocity and pressure coupling is insured by a SIMPLEC prediction/correction method with outer iterations [3]. Within the framework of this work, a resolution of structure equations is directly performed in *Code_Saturne*. Both fluid and solid resolutions are numerically coupled thanks to the modeling of a fluid-structure interface. The main purpose of this paper is to compare our numerical results to those obtained by previous authors with other codes or techniques [14], [20], [31], [24] for simulating the lock-in. Under specific conditions, the vortex shedding frequency can be tuned to the natural frequency of the oscillating cylinder. This phenomenon is an important case in the framework of code validation because of its specificity in the fluid-structure interaction field. The main result of the present work is that “lock-in”¹ (or “synchronization”) features and the influence of the non-dimensional mass on the amplitude and frequency response are simulated with a good accuracy.

The paper is organized as follows: main VIV features for rigid circular cylinder are recalled in the first section, then basic principles and choices of numerical approaches using fluid-structure coupling are presented. Finally, in the third part

¹ Definition of “lock-in” as a simple matching of the vortex shedding frequency and the cylinder natural frequency refers to a rather crude description of the process. A thorough discussion on the “true” definition of lock-in is developed in [13] or [21].

numerical results are discussed by comparison with [34] and [36] regarding cylinder's displacement amplitude, frequency, hydrodynamic forces, vortex shedding mode, and influence of non-dimensional mass on these variables. The capability of the tool to reproduce the lock-in phenomenon and its performances are evaluated, giving encouraging results for further applications to simulations of problems involving fluid structure interactions and flow induced vibrations.

1. VORTEX-INDUCED VIBRATIONS OF AN ELASTICALLY MOUNTED RIGID CYLINDER

Figure 1 shows flow past a cylinder in a plan (x, y) view with uniform cross flow. Flow upstream velocity is denoted U_∞ . For three-dimensional configurations, it is necessary to consider the z -direction, in which a lot of complex phenomena interact with two-dimensional ones, especially when the Reynolds number increases.

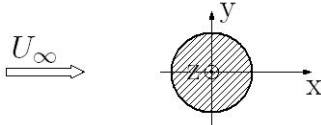


Fig. 1. Elastically mounted rigid cylinder in cross-flow

Main variables and their definition are depicted in table 1:

Variable	Definition
ρ	Fluid density (kg.m^{-3})
μ	Fluid dynamic viscosity ($\text{kg.m}^{-1}.\text{s}^{-1}$)
D	Cylinder diameter (m)
L	Cylinder length in the z -direction (m)
m	Structure mass (kg)
k	Structure stiffness
c	Structure damping

Tab. 1. Main variables of the system

From a physical point of view, it is interesting to recall dimensionless numbers commonly used to describe the flow and its interactions with a structure: the Reynolds number R_e which represents the ratio between inertia and viscous forces, the Strouhal number S_t which characterizes the reduced frequency of the system, describing the vortex shedding frequency (f_s), and drag and lift coefficients (resp. C_D and C_L), respectively calculated from the evaluation of in-line and transverse forces (F_x and F_y) on the cylinder induced by the flow.

Considering the cylinder oscillating in the transverse direction because of vortex shedding phenomenon assuming the cylinder oscillates around its mean position in the presence of flow, the cylinder motion equation is:

$$m\ddot{y} + c\dot{y} + ky = F_y \quad (1)$$

where F_y the transverse force induced by the flow is deduced from the Navier-Stokes equations. The variable y designates

the cylinder displacement due to vortex shedding. A dotted letter represents a time derivation. This dimensional equation can be re-written thanks to dimensionless parameters, as follow:

$$\frac{d^2 Y}{dt^{*2}} + \frac{4\pi\zeta}{U^*} \frac{dY}{dt^*} + \frac{4\pi^2}{U^{*2}} Y = \frac{2C_L}{\pi m^*} \quad (2)$$

Variable	Description
R_e	$\frac{\rho U_\infty D}{\mu}$
S_t	$\frac{f_s D}{U_\infty}$
C_D	$\frac{F_x}{0.5\rho U_\infty^2 DL}$
C_L	$\frac{F_y}{0.5\rho U_\infty^2 DL}$

Tab. 2. Dimensionless numbers for a fluid-structure system

Where dimensionless ratios (reduced variables) are gathered in table 3:

Symbol	Dimensionless ratio	Description
t^*	$t^* = \frac{tU_\infty}{D}$	Dimensionless time
Y	$Y = \frac{y}{D}$	Dimensionless displacement
ζ	$\zeta = \frac{c}{\sqrt{k(m + m_d)}}$	Damping ratio
m^*	$m^* = \frac{m}{\rho\pi L(D^2/4)}$	Mass ratio
U^*	$U^* = \frac{U_\infty}{f_N D}$	Reduced velocity
f^*	$f^* = \frac{f}{f_N}$	Frequency ratio

Tab. 3. Dimensionless ratios

f_N is the proper frequency of the cylinder in quiescent fluid [13], m and m_d are respectively the cylinder mass and the added mass (equal to the fluid displaced mass in the case of a simple cylinder). Maximum dimensionless amplitude A^* is deduced from an average of dimensionless displacement maxima.

Equation (2) shows that the amplitude of cylinder's vibration *a priori* depends on mass and damping ratios. The problem can be described using only one coupled mass-damping parameter ($m^* - \zeta$) if $(m^* + C_a)\zeta > 0.006$, where the added-mass coefficient C_a represents the ratio m/m_d [21]. Amplitude response is function of the mass-damping parameter:

- For high $m^* - \zeta$: amplitude response drawn versus U^* is a two-branch response, respectively called “initial excitation branch” and “lower branch”.
- For low $m^* - \zeta$: there are three distinct branches: initial and lower branches are still present, and there is another one, the so-called “upper branch”.

Details on this cylinder response (see figure 2) show a hysteretic phenomenon (“H” on figure 2) between initial and lower branch in both cases: depending on if U^* is increasing or decreasing, the obtained response is not the same. Moreover, in case of low $m^* - \zeta$, during the transition from upper to lower branch, an intermittent switching (“I”) in amplitude values is observed. Maximal amplitude is governed by the coupled mass-damping parameter for a large range of mass and damping ratios [21], while when $m^* - \zeta$ remains constant, the mass ratio variation does mainly control the width of U^* range where the system is in “synchronization” regime. Therefore, the higher the mass ratio m^* , the more significant is the tendency to develop a “lock-in” zone [36]. There is an m^* -dependency of the U^* -zone where the “lock-in” is observed: results of [36] for two different mass ratio at low $m^* - \zeta$ and low Reynolds number are presented on figure 3.

Many studies are made presenting variables as function of the reduced velocity. A more pertinent parameter is proposed in [36], which provides a unified scaling for system behavior, and which is defined when k^* or ζ falls to zero.

More precisely, induced vibrations are observed in [36] even for extreme conditions, where mass, stiffness and damping ratios tend to zero. In case of low damping (under 0.01), it is possible to completely define the system thanks to a single dimensionless parameter taking into account both mass m^* and stiffness κ^* ratios. This parameter is called “effective stiffness” (κ^*).

Considering in fact a negligible damping and a supposed purely sinusoidal cylinder’s response (and hydrodynamic forces too), the governing equation of motion can be re-written in the following form:

$$m^* \frac{d^2 Y}{dt^2} + k^* Y = C_L(t^*) \quad (3)$$

And since

$$\begin{cases} Y(t^*) = A^* \sin(\omega^* t^*) \\ C_L(t^*) = C_L \sin(\omega^* t^*) \end{cases},$$

$$\underbrace{(-m^* \omega^{*2} + k^*)}_{\kappa^*} A^* \sin(\omega^* t^*) = C_L \sin(\omega^* t^*) \quad (4)$$

Thus, effective stiffness actually combines m^* and k^* and for any choice (k^*, m^*) keeping κ^* constant, a single response is

consistent with the governing equation of motion (for further details, see [36]).

Therefore, this combined parameter is used in [36] to describe a zero- or low damping system at a given value of Reynolds number.

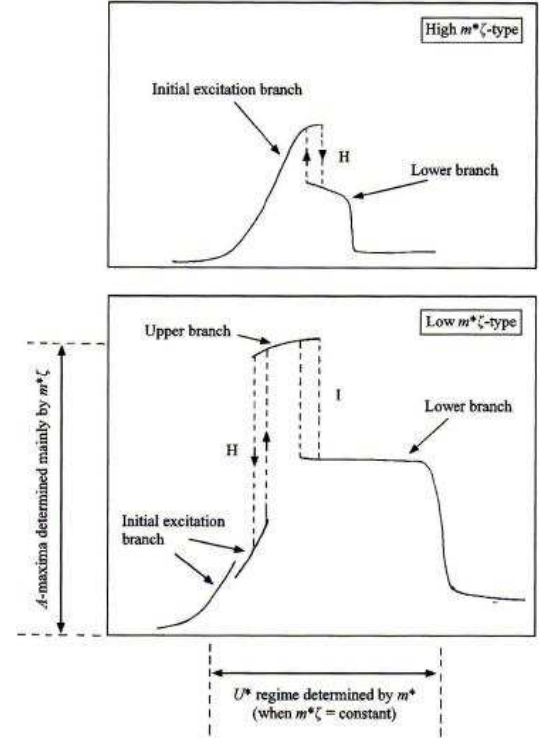


Fig. 2. Scheme for both high and low $m^* - \zeta$ types of amplitude response [21]

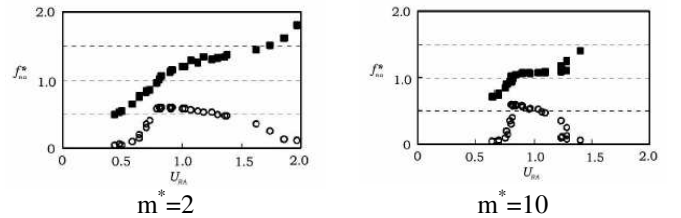


Fig. 3. Influence of mass ratio m^* on “lock-in” zone [36]

Finally, it is interesting to briefly describe the wake behavior behind the cylinder. Oscillation modes clearly depend on the response regime (initial, upper or lower branch) [21]. On the initial branch, vortex shedding occurs alternately from one side of the cylinder to the opposite: it is the so-called “2S” mode, for 2 single vortices shed per cycle (figure 4a). The regime in the lower branch is characterized by a shedding of two pairs or vortices during a cycle (“2P-mode”, see figure 4b). Concerning upper branch, conclusions are uncertain according to [21], while it is suggested in [32] that vortex shedding mode on upper branch corresponds to an intermittent switching between initial and lower branch, with a possible observation of a mixed regime 2S-2P. In [30], thanks to a

numerical simulation in good agreement with experiments of the literature, it is noticed that when the “lock-in” phenomenon is possible, a second vortex is shed from one side or the other, in addition to the 2 single vortices. This mode is called “P+S” (figure 4c).

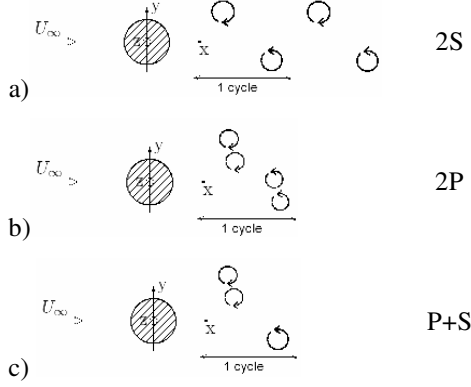


Fig. 4. Vortex shedding modes depend on the regime [41]

The Reynolds number has its influence too on the cylinder response, especially at low $m^* - \zeta$. Simulations from [38] evidence that for a fixed value of reduced velocity, changing vortex shedding modes are observed. Table 4 gathers results obtained in [38] for a cylinder free to oscillate in both x and y directions, at $U^* = 4.92$.

Reynolds number	Vortex shedding mode	
$R_e = 70$		2S
$R_e = 83$	Hysteresis	Increasing R_e : 2S
		Decreasing R_e : C(2S)
$R_e > 83 \rightarrow R_e \sim 300$		C(2S)
$R_e = 325$	Hysteresis	Increasing R_e : 2S
		Decreasing R_e : P+S
$R_e > 325 \rightarrow R_e \sim 500$		P+S

Tab. 4. Vortex shedding modes depending on the Reynolds number [38]

C(2S) stands for “2S Coalescent”: at the end of the wake, vortices merge.

Purpose of the present paper is to investigate the propensity of coupled codes to reproduce these features, namely “lock-in” behavior already observed in [36], and the effect of mass ratio on amplitude and frequency response. In the next section we present principles of fluid-structure coupling procedure, with a description of various coupling approaches, then time and space discretization choices. The fluid moving mesh formulation allowing the management of mobile meshes is also introduced. Then, in the last part, results obtained thanks to these numerical investigations are discussed. The question of using implicit or explicit coupling scheme is asked, as well as the existing limits of the use of

such a numerical partitioned procedure. But present calculations give quite encouraging results for relatively “standard” simulation conditions.

2. FLUID STRUCTURE COUPLING PROCEDURE FOR STUDYING VORTEX INDUCED VIBRATIONS

2.1. Fluid and solid models

Here we simply recall governing fluid and structure equations in order to grasp their numerical resolution (see also [23], [25] and [37]). For incompressible viscous flow, conservation equations are:

$$\begin{cases} \nabla \cdot \mathbf{u} = 0 \\ \rho_f \frac{d\mathbf{u}}{dt} = \nabla \cdot \overline{\overline{\boldsymbol{\sigma}_f}} + \mathbf{F}_f \end{cases} \quad (5)$$

where \mathbf{u} is the fluid velocity, \mathbf{F} represents volume forces and $\overline{\overline{\boldsymbol{\sigma}}}$ the stress tensor. Suffix f indicates fluid variables. The structure dynamic behavior in the transverse direction without damping is governed by:

$$m\ddot{y} + ky = F_y \quad (6)$$

Boundary conditions are the following $\begin{cases} u(t_0, x, y) = u_0(x, y) \\ y(t_0) = y_0 \end{cases}$

for the time, and the fluid velocity is imposed on Γ_v part of the fluid domain boundary:

$$\overline{\overline{\mathbf{u}}} = U_\infty \overline{\overline{\mathbf{e}_x}} \text{ on } \Gamma_v \quad \left\{ \begin{array}{l} u = \dot{y} \\ \overline{\overline{\boldsymbol{\sigma}_f}} \cdot \mathbf{n}_f = \overline{\overline{\boldsymbol{\sigma}_s}} \cdot \mathbf{n}_s \end{array} \right.$$

Fig. 5. Spatial boundary conditions on the studied domain

Finally, two continuity conditions are to be satisfied at the interface:

$$\begin{cases} \mathbf{u} = \dot{\mathbf{y}} \\ \overline{\overline{\boldsymbol{\sigma}_f}} \cdot \mathbf{n}_f = \overline{\overline{\boldsymbol{\sigma}_s}} \cdot \mathbf{n}_s \end{cases} \quad (7)$$

Fluid equations added to the pressure law $\overline{\overline{\boldsymbol{\sigma}_f}} = -p \overline{\overline{\mathbf{1}}} + 2\mu \overline{\overline{\mathbf{D}}}$ yield to Navier-Stokes equations:

$$\begin{cases} \nabla \cdot \mathbf{u} = 0 \\ \frac{d\mathbf{u}}{dt} + \nabla \cdot (\mathbf{u} \otimes \mathbf{u}) = -\frac{1}{\rho_f} \nabla p + \nu \Delta \mathbf{u} + \frac{\mathbf{F}_f}{\rho_f} \end{cases} \quad (8)$$

where ν is the kinematics viscosity. Here, we consider a rigid cylinder: consequently, we do not use stress-displacement laws.

Exchanges through the interface between fluid and structure domains have to be precise, that is why the space discretization is a real challenge for the information conservation. Moreover, diffusion or overestimations all the way through the time must be avoided as far as possible, so the time coupling represents another problem, still under investigations [2], [23] and [25].

2.2. Time matching of fluid/structure coupling

The main difficulties lie in the coupling scheme, which yields to a time-lag between fluid and structure equation resolutions. Moreover equation formulations are fundamentally different: fluid formulation is Eulerian while the structure one is Lagrangian. Interaction between both fluid and structure domains can be summed up on figure 5:

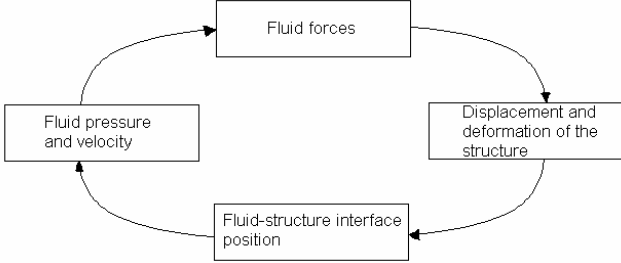


Fig. 6. Fluid and structure domains interaction

A partitioned procedure has been chosen for the present study, because it is strategically convenient to use existing solver for computation of both fluid and structure problems. A monolithic approach [6], [17] would imply an entire new code development. A partitioned scheme consists of the successive resolution of fluid and solid equations. Figure 6 simply describes the time-lag previously mentioned between both resolutions.

Knowing structure displacement at the $(n+1)^{\text{th}}$ time step, calculations can be performed to obtain a fluid state on date t^{n+1} and fluid forces are thus deduced. Now, these forces being known at t^{n+1} , structure displacement at t^{n+1} is calculated thanks to the n^{th} state and forces at t^{n+1} . Thus, because of this time-lag, energy exchanges are not exactly conserved through the interface; that is why a good displacement predictor choice is crucial.

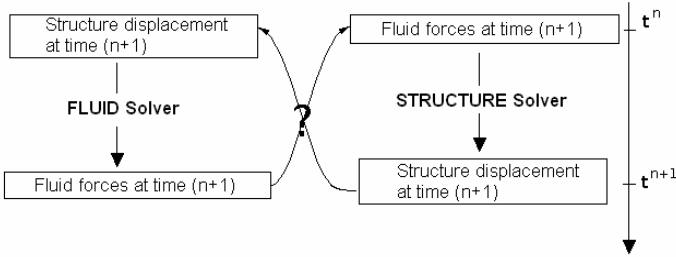


Fig. 7. Time-lag implied by partitioned approach (according to [18])

Among the partitioned schemes, we have to distinguish the weakly coupled ones from the strongly coupled. A scheme is said to be strongly coupled when (7) is exactly satisfied at each time-step, while a weakly coupled scheme does not enforce exactly (7). In the present work, we are interested in strongly coupled schemes, as far as possible.

Within the framework of our investigations, two partitioned coupled schemes are considered: explicit asynchronous [10], [33] and semi-implicit [1], [11], [16]. The first one consists in considering an intermediary time step $t^{n+1/2}$. Effort prediction is simply identified to calculated fluid force:

$$\mathbf{F}_s^{n+1,P} = \mathbf{F}_f \quad (9)$$

where the superscript P indicates a predicted value. Interface position X_f is given at time $t^{n+1/2}$ thanks to the variables calculated at t^n :

$$x_f^{n+1/2} = u_s^n + \frac{\Delta t}{2} \dot{u}_s^n \quad (10)$$

Force balance is respected with this first order scheme, as well as geometry. It is also possible to use a second order prediction for fluid forces of the form :

$$\mathbf{F}_s^{n+1,P} = \frac{1}{2} \mathbf{F}_f^n + \frac{1}{2} \mathbf{F}_f^{n+1}$$

The second scheme, the so-called ‘‘semi-implicit’’ one, is constructed in order to improve explicit predictions: fluid sub-cycles are performed at each time step, so that the position prediction should converge. The first part of the algorithm consists of an initialisation phase, and then predictions on the structure displacement are carried out thanks to the values of the previous sub-cycle [19], see Figure 7.

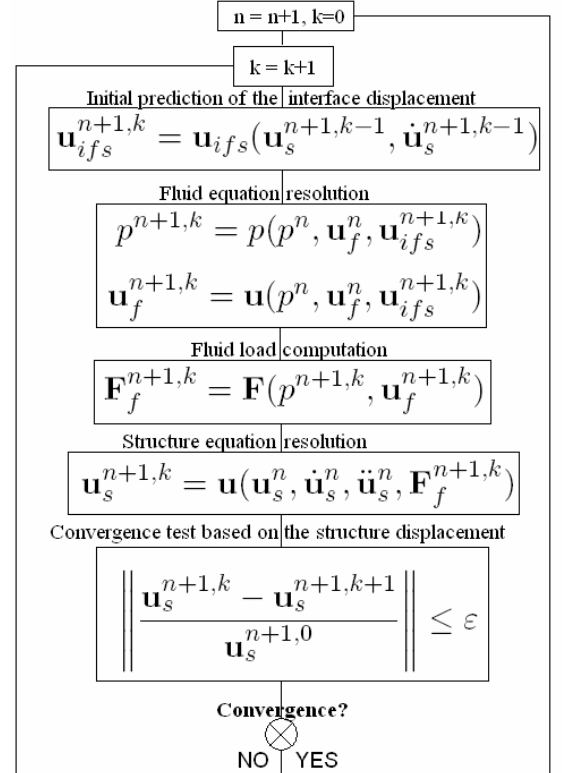


Fig. 8. Sub-cycling coupling scheme [19]

2.3. Data projection, spatial discretization of the interface

Sometimes geometries of fluid and structure models do not coincide at the interface and it is necessary to use specific techniques to keep the information conserved through the interface. However, in the present work, the cylinder is modelled as a single point mass associated with a stiffness and information on forces and displacements is interpolated at this point.

The present chosen interfacial conservation method consists of convenient projections ensuring data transfers through the interface. Fluid forces are estimated at the centre of each cell face located on the interface. They are averaged and condensed at the mass point. Conversely the structure displacement is imposed at each node of the fluid mesh located on the moving boundary. In the present case we are dealing with a projection method combined with a condensation technique. This procedure is also possible with a beam modelling for the structure.

From a general point of view, the main difficulty of data projection techniques is limited to the possible incompatibility between fluid and structure interface modelling. It is necessary to distinguish two cases: 1) cases where interfaces are jointed, when the physical interface exactly corresponds to fluid and structure interfaces at each time step; 2) cases where they are disconnected or incompatible as explained in [11], [18] and [26]. In the presence of incompatibilities several techniques are possible but master-slave techniques are suitable as far as fluid/structure interface problems are considered. An example is illustrated in [18].

2.4. Fluid moving mesh formulation

The fluid domain discretization presents difficulties because structure displacement, and hence interface, have to be followed up. As previously said, fluid resolution process is based on an Eulerian formulation, where a boundary displacement cannot be followed, while the structure resolution is based on a Lagrangian formulation, which demands big calculation times since the system encounters large-scale displacements. To get round this problem, we use the so-called ‘‘Arbitrary Lagrange Euler’’ formulation (ALE), which consists in giving a partial Eulerian and Lagrangian description to the system [9]. Thus, a new calculation domain is created, following the system physical boundaries, even for high displacement problems.

In [39], [40] ALE implementation in fluid equations is studied for fluid-structure interaction problems. We now present the analytic formulation changing fluid equations.

Let Ω_m be a material domain where fluid particles X are followed in their movements (Lagrangian description), Ω_s being a spatial fixed domain with fixed spatial positions x and Ω_{ale} a given domain whose arbitrary motion is different

from the Ω_m 's one (figure 8). In Ω_{ale} , the particular derivative is:

$$\frac{d\Phi}{dt} = \partial_t \hat{\Phi}(\xi, t)|_{\xi} = \partial_t \tilde{\Phi}(x, t)_x + \partial_{x_i} \tilde{\Phi}(x, t)_i \times \tilde{w}_i(x_i, t) \quad (11)$$

\tilde{w} represents the arbitrary referential velocity in Ω_s .

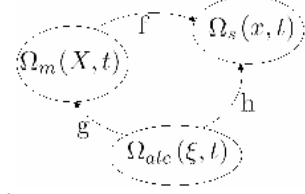


Fig. 9. Coordinates change between the three domains

Practically, it is interesting to choose Ω_{ale} in such a way that a part of its boundary is connected with the mobile interface.

Thus, ξ and w respectively correspond to the mesh position and velocity: from now on, Navier-Stokes equations are solved in the referential Ω_{ale} . So if the domain Ω_s is mobile, Eulerian formulation of Navier-Stokes equations are modified with the particular derivative definition in Ω_{ale} :

$$\begin{cases} \partial_t \tilde{p}|_x + \nabla_x \cdot (\tilde{p}\tilde{u}) = 0 \\ \tilde{\rho}[\partial_t \tilde{u}_\xi + (\tilde{u} - \tilde{w})\nabla_x \tilde{u}] = -\nabla_x \tilde{p} + \nabla_x \cdot (\tilde{\mu}\nabla_x \tilde{u}) \end{cases} \quad (12)$$

After variable changes and simplifications (see [7] and [35]), for incompressible flow, the new system becomes:

$$\begin{cases} \nabla_\xi \cdot u = 0 \\ \rho[\partial_t u_\xi + (u - w)\nabla_\xi u] = -\nabla_\xi p + \nabla_\xi \cdot (\mu\nabla_\xi u) \end{cases} \quad (13)$$

It is necessary to add the geometric conservation law (GCL) [22] to this equation system. It ensures numerical conservation of physical conserved quantities: this law states that the variation of an elementary volume during a time Δt must be equal to the volume covered by its boundary faces during the same time. This law is commonly used as a necessary condition to provide mesh stability.

The integral form of the GCL is (14):

$$\frac{\partial}{\partial t} \int_v dV|_\xi = \int_{\delta v} \tilde{w} \cdot ndS \quad (14)$$

In order to work with regular cells, the mesh velocity choice can either be based on the node displacement or on the velocity. In the case where the mesh displacement (or the mesh velocity) governing equation is linear, a first order approximation gives:

$$w_i^n = \frac{\xi_i^{n+1} - \xi_i^n}{\Delta t}$$

where ξ_i^{n+1} and ξ_i^n respectively the displacement at $(n+1)^{th}$ and n^{th} time step and w_i^n the mesh velocity at $t = n\Delta t$.

Furthermore, the mesh actualisation must be adapted to various cell forms and must avoid excessive modifications of the mesh topology. Moreover, as it is called at each time step, the algorithm has to spare calculation time and to be as robust as possible. The grid updating is performed as follows: considering a domain with moving boundaries, each node of

the mesh is linearly linked to two nodes of the boundary (see fig. 9), thanks to the relation between positions \mathbf{X} :

$$\mathbf{X}_{\text{node}} = \alpha \mathbf{X}_{\text{boundary-1}} + (1 - \alpha) \mathbf{X}_{\text{boundary-2}} \quad (15)$$

where α represents the parametric coordinates.

Thanks to this method, the boundary conditions are preserved.

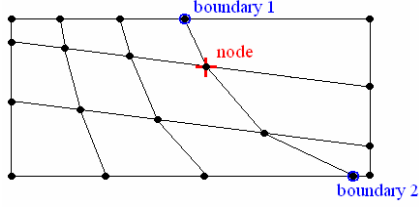


Fig. 10. Mesh connection example

Finally, a general algorithm for a finite volume Navier-Stokes resolution with ALE method is presented. We first have to pay attention to the fact that physical magnitudes are calculated at cell centres, but the mesh deformation requires the information at nodes. The result is that an interpolation from mesh deformation to determine node displacements, respecting the GCL, must be performed.

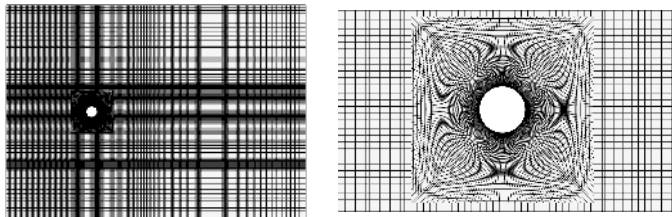
Thus, the general ALE algorithm is given by:

- *Time loop*
 - ALE displacement boundary condition defined at the fluid-structure interface
 - Transformation of these boundary conditions into velocity boundary conditions, GCL enforced
 - Mesh deformation velocity calculation
 - Mesh deformation thanks to velocity data and GCL
 - Navier-Stokes equation resolution

3. NUMERICAL RESULTS AND DISCUSSION

By default, the time coupling scheme is the explicit asynchronous one and an unsteady method is adopted, as we are interested in the setting up of an established regime. Pressure calculation is performed thanks to a second order scheme.

In order to avoid 3D effects, investigations are performed at low Reynolds number ($R_e = 100$). On Figure 10 we show the fluid mesh for present calculations.



a) Global fluid mesh b) Zoom around the cylinder

Fig. 11. Finite volume mesh

3.1. Vortex shedding in the wake of the fixed rigid cylinder in cross flow

A time-step convergence is set up, so that the Courant-Friedrichs-Levy condition should be satisfied. Our aim is to be able to observe the vortex shedding phenomenon, so a flow perturbation is imposed at the beginning of the calculation with the fixed cylinder. Then, when the steady state is reached, calculations with mobile mesh are launched. Final results are presented in table 5, in comparison with [14], [31], [34].

	$C_{D,\text{mean}}$	$C'_{L,\text{max}}$	$C'_{L,\text{rms}}$	S_t
Present simulation	1.389	0.328	0.224	0.166
[34]	1.374	0.327	0.227	0.169
[14]	1.386	0.323	-	0.165
[31]	-	-	0.227	0.164
Error present/[34] (%)	1.1	0.3	1.3	1.8
Error Present/[14] (%)	0.2	1.5	-	0.6
Error Present/[31] (%)	-	-	1.3	1.2

Tab. 5. Hydrodynamic forces and Strouhal number for a fixed cylinder

Here, $C_{D,\text{mean}}$, $C'_{L,\text{max}}$ and $C'_{L,\text{rms}}$ respectively represent a mean of the drag performed at each time step, the mean on local maxima of the fluctuating lift signal, and the RMS (root mean square) of the fluctuating lift signal. Results are reasonably encouraging for the rest of the investigations.

3.2. Vortex induced vibrations at mass ratio $m^*=3.3$

The cylinder is allowed to vibrate in the transverse direction (y direction), while the flow velocity is uniform in the x-direction. Mechanical characteristics are chosen to be the same as those of [34] and [36]. Thus, mass ratio is equal to 3.3, and the range of U^* (or κ^* , notation chosen to indicate κ^*_{eff} in [36]) is from $U^* = 3.00$ to $U^* = 11.32$ (or κ^* ranging from 1.47 to 17.64). Various dimensionless variables are plotted against effective stiffness κ^* : the maximal amplitude A^* , reduced frequency f^* (where $f^* = fd/U_\infty$ with f the actual vibration frequency of the cylinder), and both hydrodynamic coefficients ($C_{D,\text{mean}}$ and $C_{L,\text{max}}$).

These results are respectively gathered from figure 11a) to 11d).

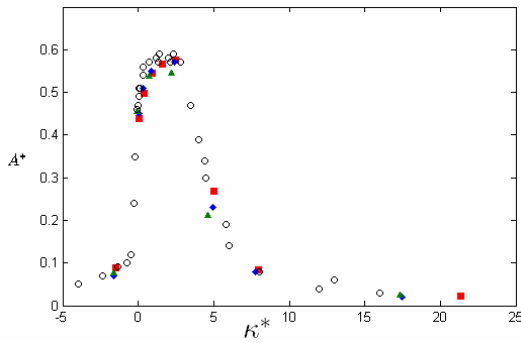
Good qualitative and quantitative reproduction of the “lock-in” phenomenon by both EDF and DCNS codes are noted. It is encouraging for the use of coupled schemes in fluid-structure interaction, even with an explicit partitioned scheme. In a next part, the advantages of an implicit scheme will be discussed.

A frequency analysis on maximal non-dimensional amplitude reveals that the Khalak et al. [21] dimensionless frequency $f^*_{\kappa\&w}$ (with $f^*_{\kappa\&w} = f/f_N$) according as κ^* (or

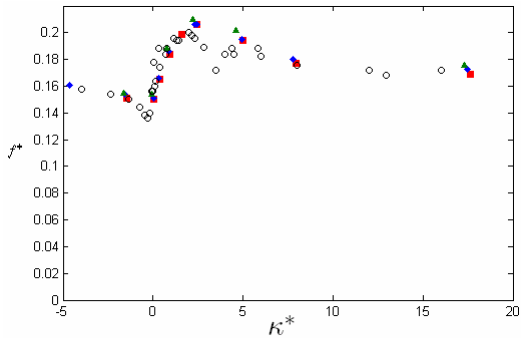
U^*) be high or low is not the same. For low κ^* , $f_{\kappa\&w}^*$ is close to 2, and for high κ^* , $f_{\kappa\&w}^*$ is almost equal to 0.5.

For κ^* around 2.44 ($U^* = 4.91$), $f_{\kappa\&w}^*$ is very close to unity: it confirms a “resonance” phenomenon, where the cylinder frequency is close to its proper frequency in water, all the more so since for each variable (A^* , $C_{D,mean}$ and $C_{L,max}$), the biggest value is obtained for $\kappa^* \cong 2.44$. On the other hand, when $U^* \cong 4.40$, we observe a beating phenomenon (several frequency existence is revealed by the FFT of signals), also observed in [21], [34] and [36]. This aspect is considered in [21] as a vortex shedding mode changing.

a)



b)



c)

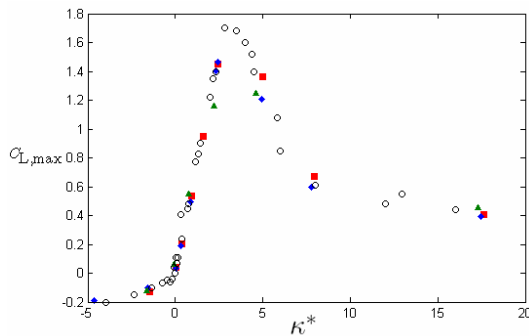
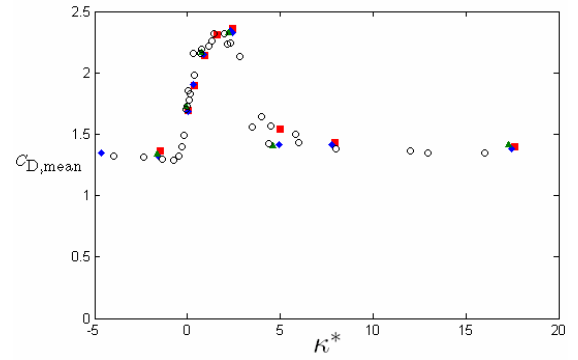


Fig. 12. Comparison between present simulations and [34], reference computation is [36], $R_e = 100$

- a) Dimensionless amplitude,
- b) Dimensionless frequency,
- c) Maximal lift coefficient, all variables are plotted versus κ^* .

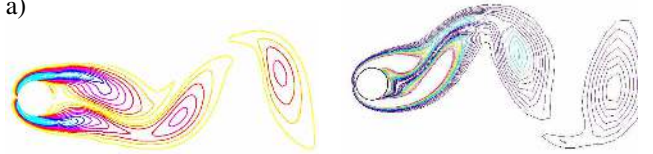
d)



- Present simulation, ◆ Placzek et al. (2007),
- ▲ Jus (2007), ○ Reference (Shiels et al. 2001)

Fig. 11. Comparison between present simulations and [34], reference computation is [36], $R_e = 100$ (continued)
d) Mean value of the drag coefficient, all variables are plotted versus κ^* .

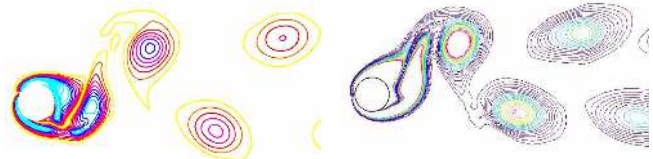
a)



Present: $\kappa^* = -1.47$

[34]: $\kappa^* = 0.05$

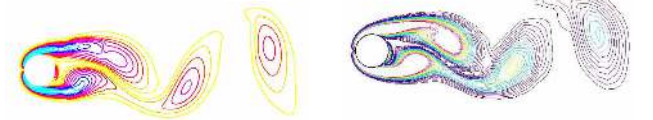
b)



Present: $\kappa^* = 2.44$

[34]: $\kappa^* = 2.32$

c)



Present: $\kappa^* = 17.64$

[34]: $\kappa^* = 17.5$

Fig. 13. Different vortex shedding modes observed for 3 values of κ^* (low position of the cylinder, and dimensionless vorticity $\omega^* = \omega d/U_\infty$ is between 0.2 and 2)

At $R_e = 100$, it is difficult to give a precise conclusion, but observing vorticity contours on figure 12, it is possible to distinguish two different emission modes, compared with [34]. Thus, when κ^* is low or high (limits of the “lock-in” zone), the shedding mode is close to the Strouhal’s one, (observed for a fixed cylinder) i.e. Von Kármán vortex street, while if κ^* has a midway value, i.e. $\kappa^* = 2.44$ (“lock-in”), vortices are not so stretched and the wake is clearly separated in two parts on both sides of the centerline. In spite of the fact that

this study is led at low Reynolds number, the change in vortex shedding modes usually observed in the literature for higher Reynolds number is here perceptible.

Figure 13 is a description of the link between the frequency behavior and the “synchronization”: straight lines represent the Strouhal frequency versus reduced velocity, for $Re = 100$ and Re situated in $[10^3, 10^4]$, respectively calculated from the present Strouhal number and from empirical formula proposed in [31]. The reduced frequency is $f^*_{K&W}$, plotted against reduced velocity. Data points that are situated close to Strouhal lines signify that for this range of reduced velocity, the Strouhal frequency (i.e. vortex shedding frequency) is governing the motion. On the other hand, data points situated just above $f^* = 1$ depict the range of reduced velocity where the “lock-in” phenomenon is present. For higher Reynolds numbers (reference experimental data come from [21]), we can clearly distinguish three response branches.

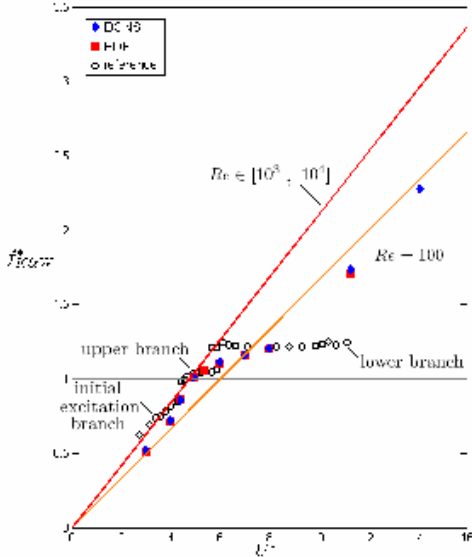


Fig.14. Representation of the frequency lock-in phenomenon

3.3. Mass ratio influence consideration

Now, we are interested in the propensity of the system to have a larger reduced velocity range where amplitude is the highest for a low mass ratio m^* , than for a high mass ratio. At the same time, reduced frequency f^* is most clearly “locked” close to unity for high mass ratio. Figure 14 gathers results obtained for $m^* = 1$, $m^* = 3.3$ and $m^* = 10$. The code reproduces well this mass-ratio dependence.

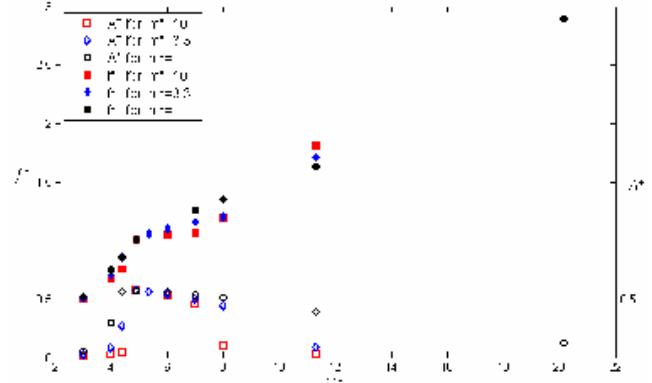


Fig.15. Dimensionless frequency and maximal amplitude for various mass ratios

It is worth noting that for the case $m^* = 1$, we have to choose the semi-implicit coupled scheme [7]: without sub-cycles, the explicit one is diverging. Comparing both synchronous explicit and semi-implicit schemes for $\kappa^* = 2.44$ ($U^* = 4.91$), we obtain that, as expected, the implicit scheme is more precise and an error between both amplitude response sets up. But this error remains low, and thanks to the good comparison with literature data, an explicit scheme is sufficient as long as convergence is assured. On the other hand, as soon as we want to work with a very low mass-ratio, neither of explicit or semi-implicit schemes is sufficient, even if a large number of sub-cycles is used. The mass ratio $m^* = 0.8$ seems to be a limit of the use of such a coupled algorithm unless an additional stabilization procedure is involved.

CONCLUSION

Numerical simulations of the VIV phenomenon and particularly the frequency “lock-in” of a cylinder in a uniform cross flow, without damping and at low Reynolds number has been performed and discussed in the present study. Simulations have been performed thanks to a CFD code, using a partitioned coupled scheme for simulation of coupling with the structure motion. An ALE formulation has been chosen to resolve fluid equations with a moving mesh. The aim of these investigations is to validate the computational process for this “lock-in” phenomenon, and for the mass ratio influence. Quantitative agreement has been highlighted, respectively for amplitude, frequency, hydrodynamic forces, and qualitative agreement has been observed for vortex shedding patterns. Moreover, mass ratio influence is quite well reproduced by the code. However, in the extreme case of very low mass ratio, a better coupled scheme should be considered.

Nevertheless, these results give encouraging hopes for the use of such a coupling method for simulation of fluid-structure interactions.

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ACKNOWLEDGEMENT

The authors wish to thank Vincent MELOT from DCNS Propulsion and François JUSSEMAND from EDF R&D Chatou for their helpful contribution in the development of coupling procedures within the CFX code and *Code_Saturne*.