Abstract

Accuracy is critical if we are to trust the simulation’s predictions. In settings such as fluid-structure interaction it is all the more important to obtain reliable results to understand, for example, the impact of pathologies on blood flows in the cardiovascular system. In [1] we have proposed a framework for high order (in space and time) fluid structure interaction in 2D using an efficient high order ALE map construction which is described in [2]. In the first part of the paper, we propose a new high order ALE construction which allows for any type — not only vertical as in [2] — of not too large displacements that applies to 2D as well as 3D. This construction relies on a first order approximation and a correction step that recovers the high order accuracy. In the second part of the paper we present an update of our fluid-structure interaction framework.

Keywords: High order methods, Arbitrary Lagrangian Eulerian transformation, Fluid-Structure Interaction

1. Introduction

Over the last few years, we have been working on building a mathematical and computational framework for high order fluid-structure interaction, see [1–6], in 2D and 3D, on simplices and hypercubes, with a wide range of applications and in particular, bio-mechanics (e.g. blood flows in arteries). In this paper we present the progress made since our last publications [1, 2] as well as a brief overview of the framework we have built so far.

The paper is organized in the following way: first, we introduce some notations and present a brief overview of the status of [6] in section 1.2 (which will be described in another publication). In section 2 we state our latest advances in one of our central ingredients to achieve high order fluid-structure interaction, namely the Arbitrary Lagrangian-Eulerian framework. Next, the fluid and structure models are presented, associated with their respective discretisation and solution methods. These methods are...
validated based on the benchmark [7]. In the last section, we propose a fluid-structure interaction solver, combination of the tools presented throughout the paper. The solver is validated by the benchmark [7] and some 3D preliminary results are presented.

1.1. Notations

Given an elementary simplex domain $K^* \subset \mathbb{R}^d$, $d = 1, 2, 3$, and a positive integer $N$, let us denote by $P^N(K^*)$ the space of polynomials of total degree less or equal than $N$, defined in $K^*$. We fix a reference element, $\hat{K}$, and consider a transformation $\varphi_{K^*} : \hat{K} \rightarrow K^*$, called the geometrical transformation.

We consider now two domains, $\Omega^*$ and $\Omega \subset \mathbb{R}^d$, $d = 1, 2, 3$, which we later refer as the reference and the computational domains, respectively. We further assume that the reference domain has a straight edge/face mesh associated with it, $T^*$, i.e., $\varphi_{K^*} \in P^1(\hat{K}), \forall K^* \in T^*$. Furthermore we admit that the mesh $T^*$ covers exactly the domain $\Omega^*$, i.e., $\Omega^* = \bigcup_{K^* \in T^*} \overline{K^*}$.

We denote $P^N_{c,h}(\Omega^*)$, $P^N_{c,h}(\Omega^*)$ the spaces of piecewise scalar, respectively vectorial, polynomial of total degree $N$, continuous functions in $\Omega^*$, and $P^N_{td,h}(\Omega^*)$ the space of piecewise polynomial of total degree $N$, totally discontinuous functions in $\Omega^*$.

Finally let us denote by $\eta : \partial \Omega^* \rightarrow \partial \Omega$ a displacement function. Through $\eta$, we classify three subsets of the boundary: (i) $\Gamma^*_M$, the portion of the boundary that moves according to the displacement $\eta$, (ii) $\Gamma^*_F$, the portion of the boundary that stays fixed (i.e., $\eta(s) = s, \forall s \in \Gamma^*_F$) and (iii) $\Gamma^*_N$, the part of the boundary on which we do not prescribe a displacement. The image of each subset, $\Gamma^*_M$, $\Gamma^*_F$ and $\Gamma^*_N$ by $\eta$ is denoted by $\Gamma_M$, $\Gamma_F$ and $\Gamma_N$, respectively. These three sets do not overlap and they verify $\partial \Omega^* = \overline{\Gamma_M} \cup \overline{\Gamma_F} \cup \overline{\Gamma_N}$. Denote $T^{*b} = \{ K^* \in T^* : \partial K^* \cap \Gamma^*_M \neq \emptyset \}$ the set of elements $K^*$ sharing a face with the boundary of $\Omega^*$.

1.2. Computational framework for Galerkin methods

Our computational framework builds upon Feel++ [6] which allows for arbitrary order cG and dG Galerkin methods (finite element, spectral elements, ...) in 1D, 2D and 3D on simplices and hypercubes. The computational domain can also be high order, that is to say, the geometrical transformation of each element $K$ of the mesh is a polynomial of degree greater than one. These high order meshes can be generated by Gmsh [8] — up to order five in 2D and order four in 3D. — High order approximations come at a cost both in terms of implementation and computational points of view. The former is addressed by a very generic framework based on modern C++ programming (meta-programming, expression templates, ...) and a language mimicking the mathematical language. The latter is addressed by a careful implementation and optimisation. One of the optimisations that allows to have a huge gain in computational effort is to straighten all the high order elements except for the boundary faces of the computational mesh. This is achieved by moving all the nodes associated to the high order transformation to the position...
these nodes would have if a first order geometrical transformation were applied. This procedure can be formalized in the following operator

$$\eta_{K}^{\text{straightening}}(\varphi^{N}_{K}(x^{*})) = (\varphi^{1}_{K}(x^{*}) - \varphi^{N}_{K}(x^{*})) - (\varphi^{1}_{K\cap\Gamma}(x^{*}) - \varphi^{N}_{K\cap\Gamma}(x^{*}))$$

(1)

where \(x^{*}\) is any point in \(K^{*}\) and \(\varphi^{1}_{K}(x^{*})\) and \(\varphi^{N}_{K}(x^{*})\) its images by the geometrical transformation of order one and order \(N\), respectively. On one hand, the first two terms ensure that for all \(K\) not intersecting \(\Gamma\), the order one and \(N\) transformations produce the same image. On the other hand, the last two terms are 0 unless the image of \(x^{*}\) in on \(\Gamma\) and, in this case, we don’t move the high order image of \(x^{*}\). This allows to have straight internal elements and elements touching the boundary to remain high order. When applying numerical integration, specific quadratures are considered when dealing with internal elements or elements sharing a face with the boundary. The performances, thanks to this transformation, are similar to the ones obtained with first order meshes. However, it needs to be used with care as it can generate folded meshes.

2. High order ALE

We now turn to our high order Arbitrary Lagrangian-Eulerian (ALE) framework. A fundamental piece in performing simulations in the ALE framework is the transformation that maps the reference configuration onto the computational domain, at each timestep. This is called ALE map. In Pena and Prud’homme [1], the authors propose a high order ALE map that allows for an accurate description of the boundary of the computational domain, while inducing a straight edges in the interior elements of the computational domain’s mesh. However, this construction has the disadvantage of relying upon the Gordon-Hall transformations, see Gordon-Hall [9, 10], which makes it (implementation-wise) intricate to extend to three dimensional domains.

To overcome this difficulty, we replace the stage based on Gordon-Hall transformations, by the solution of a local differential problem in each element in contact with the curved boundary. We review here the construction from [1, 2]. The first step is to perform a modified harmonic extension (according to Masud and Kanchi [11]) of the displacement \(\eta\) to the interior of the reference domain using piecewise linear polynomial functions. The corresponding ALE transformation, \(A_{1}\), satisfies a discrete element-weighted Laplace equation.

The second step is a correction performed in each element that touches the curved boundary in order to build a high order approximation. In each element \(K^{*} \in T^{*}b\) we look for \(A^{N}_{K^{*}} \in [P^{N}(K^{*})]^{d}\) such that

$$\begin{cases}
\int_{K^{*}} (1 + \tau) \nabla A^{N}_{K^{*}} : \nabla z \, dx = 0, \quad \forall z \in [P^{N}(K^{*})]^{d} \\
A^{N}_{K^{*}}(x^{*}) = \eta(x^{*}) + x^{*} - A^{1}(x^{*}), \quad \forall x^{*} \in \partial K^{*} \cap \Gamma^{*}_{M} \\
A^{N}_{K^{*}} = 0, \quad \text{elsewhere on } \partial K^{*}.
\end{cases}$$
where \( \Gamma^* \) is the portion of boundary in the reference domain that is curved in the computational domain. The final ALE map, \( A^N \) is obtained by adding to \( A^1 \) the correction \( A^N_{K^*} \) on each element of \( T^*_{b} \)

\[
A^N(x^*) = A^1(x^*) + \sum_{K^* \in T^*_{b}} A^N_{K^*}(x^*) + x^*
\]

**Proposition 1** (Properties of \( A^N \)). Under the previous assumptions and by construction, \( A^N \in P_{c,h}^N(T^*_{b}) \) — enjoys optimal approximation properties i.e. the boundary approximation is \( O(h^{N+1}) \) in the \( L_2 \) norm — and \( A^N \in P_{c,h}^1(T^* \setminus T^*_{b}) \).

**2.1. The Harmonic extension and Winslow smoother**

In both papers [1, 2], the piecewise linear map created in the first step is calculated by performing a harmonic extension (or modified harmonic extension) of the boundary data. However, if the displacement is too large, these operators can induce meshes that are not valid due to, for instance, mesh folding. A way to circumvent this problem, that steams from the structure of the proposed ALE map construction, is to replace the harmonic extension by a more suitable and flexible operator that avoids these issues or improves the mesh quality. An example of such an operator is the Winslow smoother [12]. From a continuous point of view, the Winslow smoother enforces that the inverse of the ALE map is harmonic, not the map itself. This accounts for solving a quasi-linear system of PDE’s, which can be done using fixed point iterations. In Figure 1 we show the effect of the modified harmonic extension and the Winslow smoother for the same testcase.

![Figure 1: Comparison of first order meshes generated by the harmonic extension (left) and Winslow smoother (right) operators respectively](image-url)
2.2. Numerical verification of Proposition 1

We present now some numerical experiments to verify Proposition 1. In order to conduct these experiments we use (i) Gmsh [8] to build the initial high order meshes in 2D and 3D and visualize the meshes computed by our methods and (ii) Feel++ [3, 6] which provides the framework for arbitrary order Galerkin methods to solve the partial differential equations and handle the computational meshes. Note that the mesh $\mathcal{T}^*$ used to solve equation (1) is built automatically from the initial high order mesh through the straightening process.

We consider the reference domain depicted in Figure 2(a) defined by

$$\Omega^{*\cdot\text{cy}} = \left\{ (x^*, y^*, z^*) \in \mathbb{R}^3 : x^* \in [0, 5], y^*^2 + z^*^2 \leq 0.5^2 \right\} \quad (2)$$

and the associated displacement of its boundary $\eta^y(x^*) = 0.2 \exp\left(\frac{x^*}{5}\right) \sin\left(\frac{\pi x^*}{2.5}\right) n^*$. Figure 2(b) displays the computational domains colored by the corresponding ALE map.

Finally, Figure 3(a) displays the convergence rate of the quantity $\|A^N(x^*) - (x^* + \eta(x^*))\|_{[L^2(\Gamma_M^*)]^d}$ which confirms the result of Proposition 1. Figure 3(b) represents the convergence rate with respect to geometric order $N$. 

Figure 2: Reference (left) and computational (right) meshes of order 4 displayed using Gmsh colored by the displacement 2-norm
3. Models and their discretizations methods

We now turn to the structure and fluid models, their associated discretization and solution methods. In each case, we display some of the results of our strategy applied to the benchmark [7]. Note that our fluid solver has also been benchmarked for bi-fluid simulations [13] following the proposal [14].

Two benchmarks presented in [7] are performed to verify our code. The first one corresponds to the oscillation of an elastic beam only subjected to gravity. We monitor the coordinates of the tip of the beam. The second is the simulation of a flow in rectangular domain with a rigid obstacle: a flag clamped to a cylinder. We monitor the lift and drag.

The algebraic systems arising from the discretization proposed in the next sections are solved using a Newton or quasi-Newton algorithm with a cubic line search method. At each step, the linear solver applies the GMRES method with a LU preconditioner. The preconditioner is typically built only once throughout the nonlinear iterations unless the nonlinearity is stiff and the preconditioner needs to be recalculated. In the quasi-newton instance, the Jacobian can be rebuilt once in a while during the nonlinear iterations or just once, which is often preferred when simulating time-dependent problems. The underlying framework for the linear and nonlinear solvers is PETSc [15].

3.1. Structure

3.1.1. Models and discretizations

We first introduce the deformation gradient tensor, which allows to measure the solid deformation from the displacement $\eta_s$, $F_s = I + \nabla \eta_s$.

Other useful tensors are the right Cauchy-Green tensor $C_s$ and the Green-Lagrange tensor $E_s$ which can be expressed by $C_s = F_s^T F_s$, $E_s = \frac{1}{2} (C_s - I)$ where $E_s$ has the property to be divided in two terms, $\epsilon_s$ (linear) and $\gamma_s$ (quadratic)

$$ E_s = \frac{1}{2} \left( \nabla \eta_s + (\nabla \eta_s)^T \right) + \frac{1}{2} \left( \nabla \eta_s \right)^T \nabla \eta_s. \tag{3} $$
The simplest model of this type is the **elastic linear model**, which is valid for small displacements and deformations. It reads as
\[ \rho_s \frac{\partial^2 \eta_s}{\partial t^2} - \nabla \cdot (\Sigma_s) = f_s, \quad \Sigma_s = \lambda_s (\text{tr} \epsilon_s) I + 2\mu_s \epsilon_s \]
with \( \lambda_s \) and \( \mu_s \) being the Lamé coefficients.

The next model is the **hyper-elastic and compressible model** which is valid for large deformations. The generic balance equation of hyper-elasticity in Lagrangian description is given by
\[ \rho_s \frac{\partial^2 \eta_s}{\partial t^2} - \nabla \cdot (F_s \Sigma_s) = f_s, \quad \Sigma_s = \lambda_s (\text{tr} E_s) I + 2\mu_s E_s. \]

where \( \Sigma_s \) represents the second Piola-Kirchhoff stress tensor.

To take into account the material incompressibility, we use the **Hyper-elastic and incompressible model**. The pressure \( p_s \) acts as a lagrange multiplier to enforce incompressibility. The model reads as
\[ \rho_s \frac{\partial^2 \eta_s}{\partial t^2} - \nabla \cdot (F_s \Sigma_s) = f_s, \quad \text{det} F_s = 1, \quad \Sigma_s = -p_s (\text{det} F_s) C_s^{-1} + \lambda_s (\text{tr} E_s) I + 2\mu_s E_s. \]

Finally we also make use of a **1D reduced model** for thin structures like shells, also known as generalized strings, see [5, 16].

For the numerical results presented in the next section we considered a hyper-elastic model for the structure. We discretized the equations using \([P]^d\) elements and a continuous approximation in space while we used the Newmark method to get an order 2 discretization in time.

**3.1.2. Benchmark**

The results of the CSM3 benchmark from [7] are displayed in Table 1 and the \( x, y \) coordinates of the tip of the beam on the Figure 4. The results are in accordance with the reference values REF.
\[ \begin{array}{|c|c|c|c|c|} \hline N_{\text{element}} & N_{\text{dof}} & x \times 10^{-3} & y \times 10^{-3} \\ \hline \text{REF} & -14.305 \pm 14.305 \ [1.0995] & -63.607 \pm 65.160 \ [1.0995] \\ \hline 4199 & 17536(P_2) & -14.585 \pm 14.59 \ [1.0953] & -63.981 \pm 65.521 \ [1.093] \\ 4199 & 38900(P_3) & -14.589 \pm 14.594 \ [1.0953] & -63.998 \pm 65.522 \ [1.093] \\ 1043 & 17422(P_4) & -14.591 \pm 14.596 \ [1.0953] & -64.009 \pm 65.521 \ [1.093] \\ 4199 & 68662(P_4) & -14.59 \pm 14.595 \ [1.0953] & -64.003 \pm 65.522 \ [1.093] \\ \hline 4199 & 17536(P_2) & -14.636 \pm 14.64 \ [1.0969] & -63.937 \pm 65.761 \ [1.0945] \\ 4199 & 38900(P_3) & -14.642 \pm 14.646 \ [1.0969] & -63.949 \pm 65.771 \ [1.0945] \\ 1043 & 17422(P_4) & -14.645 \pm 14.649 \ [1.0961] & -63.955 \pm 65.778 \ [1.0945] \\ 4199 & 68662(P_4) & -14.627 \pm 14.629 \ [1.0947] & -63.916 \pm 65.739 \ [1.0947] \\ \hline 4199 & 17536(P_2) & -14.645 \pm 14.645 \ [1.0966] & -64.083 \pm 65.623 \ [1.0951] \\ 4199 & 38900(P_3) & -14.649 \pm 14.65 \ [1.0966] & -64.092 \pm 65.637 \ [1.0951] \\ 1043 & 17422(P_4) & -14.652 \pm 14.653 \ [1.0966] & -64.099 \pm 65.645 \ [1.0943] \\ \hline \end{array} \]

Table 1: Results for CSM3 with $\Delta t = 0.02, 0.01, 0.005$. Line REF displays the reference values for [7].

3.2. Fluid

3.2.1. Models and discretization

We mainly use a Newtonian fluid model which neglects shear-thinning and viscoelastic effects. The governing equations are the classical Navier-Stokes equations which read as

\[
\rho_f \frac{\partial u_f}{\partial t} + \rho_f (u_f \cdot \nabla) u_f - \nabla \cdot \sigma_f = f_f \\
\nabla \cdot u_f = 0
\]

where $u_f$ is the fluid velocity and $\rho_f$ its density and

\[
\sigma_f = -p_f I + \tau_f, \quad \tau_f = 2\mu_f D_f, \quad D_f = \frac{1}{2} \left( \nabla u_f + (\nabla u_f)^T \right)
\]
where \( p_f \) is the pressure and \( \mu_f \) the fluid viscosity.

The previous equations are discretized using the standard Taylor-Hood element \([\mathbb{P}^N]^d \times \mathbb{P}^{N-1}\) in space and the BDF of order 2 or 3 discretizations in time. In 2D the geometry is discretized using order 1 to 5 geometric transformations while in 3D from order 1 to 4, see [8].

### 3.2.2. Benchmark

The results of the CFD3 benchmark from [7] are displayed in Table 3.2.2 for various geometrical \( (N_{geo}) \) and velocity/pressure \( N/N-1 \) approximations. The results in Table 3.2.2 for a BDF2 scheme and time step \( 5 \times 10^{-3} \) are in accordance with the REF results. However, we made some extra computations using a smaller time step \( \Delta t = 0.002 \) and BDF2/BDF3 schemes and we observe a significant shift in the Drag and Lift coefficients. This shift is already present with BDF3 and \( \Delta t = 0.005 \). Several computations were made using various meshes and discretization for the velocity, pressure and geometry. BDF2 and BDF3 for \( \Delta t = 0.002 \) are very much in accordance with the reference results except perhaps for the mean of the Lift which tells us that we are probably resolving properly the time and spatial scales while \( \Delta t = 0.005 \) using BDF2 or BDF3 is not sufficient.

<table>
<thead>
<tr>
<th>( N_{geo} )</th>
<th>( N_{element} )</th>
<th>( N_{dof} ) ( (N, N-1) )</th>
<th>( N_{bdf} )</th>
<th>Drag ( \times 10^{-3} )</th>
<th>Lift ( \times 10^{-3} )</th>
</tr>
</thead>
<tbody>
<tr>
<td>( P_1 )</td>
<td>8042</td>
<td>37514(( P_2/)P_1)</td>
<td>2</td>
<td>437.47 ± 5.3750 [4.3457]</td>
<td>−9.7865 ± 437.54 [4.3457]</td>
</tr>
<tr>
<td>( P_2 )</td>
<td>2334</td>
<td>26706(( P_3/)P_2)</td>
<td>2</td>
<td>439.27 ± 5.1620 [4.3457]</td>
<td>−8.887 ± 429.06 [4.3457]</td>
</tr>
<tr>
<td>( P_2 )</td>
<td>7970</td>
<td>89790(( P_3/)P_2)</td>
<td>2</td>
<td>439.56 ± 5.2335 [4.3457]</td>
<td>−11.719 ± 425.81 [4.3457]</td>
</tr>
<tr>
<td>( P_3 )</td>
<td>3509</td>
<td>39843(( P_3/)P_2)</td>
<td>2</td>
<td>438.24 ± 5.5375 [4.3945]</td>
<td>−11.024 ± 433.90 [4.3945]</td>
</tr>
<tr>
<td>( P_3 )</td>
<td>8042</td>
<td>90582(( P_3/)P_2)</td>
<td>2</td>
<td>439.25 ± 5.6130 [4.3945]</td>
<td>−10.988 ± 437.70 [4.3945]</td>
</tr>
<tr>
<td>( P_2 )</td>
<td>2334</td>
<td>26706(( P_3/)P_2)</td>
<td>2</td>
<td>439.49 ± 5.5985 [4.3945]</td>
<td>−10.534 ± 441.02 [4.3945]</td>
</tr>
<tr>
<td>( P_2 )</td>
<td>7970</td>
<td>89790(( P_3/)P_2)</td>
<td>2</td>
<td>439.71 ± 5.6410 [4.3945]</td>
<td>−11.375 ± 438.37 [4.3945]</td>
</tr>
<tr>
<td>( P_3 )</td>
<td>3499</td>
<td>73440(( P_4/)P_3)</td>
<td>3</td>
<td>439.93 ± 5.8072 [4.492]</td>
<td>−14.511 ± 440.96 [4.3945]</td>
</tr>
<tr>
<td>( P_4 )</td>
<td>2314</td>
<td>78168(( P_5/)P_4)</td>
<td>2</td>
<td>439.66 ± 5.6412 [4.3945]</td>
<td>−11.329 ± 438.93 [4.3945]</td>
</tr>
<tr>
<td>( P_3 )</td>
<td>2340</td>
<td>49389(( P_4/)P_3)</td>
<td>2</td>
<td>440.03 ± 5.7321 [4.3945]</td>
<td>−13.25 ± 439.64 [4.3945]</td>
</tr>
<tr>
<td>( P_3 )</td>
<td>2334</td>
<td>49266(( P_4/)P_3)</td>
<td>3</td>
<td>440.06 ± 5.7773 [4.3945]</td>
<td>−14.092 ± 440.07 [4.3945]</td>
</tr>
</tbody>
</table>

Table 2: Results for CFD3 with \( \Delta t = 0.01 \times 0.005 \). Line REF displays the reference values for [7]

### 4. Fluid structure interaction

In the fluid-structure interaction context, we chose to write the fluid dynamics equations in the Arbitrary Lagrangian Eulerian (ALE) framework. It allows to take into account the deformation of
the fluid domain. We need also to introduce into our model the domain’s velocity of deformation $w_f$, see e.g. [1, 2], the fluid equation set now reads in the moving domain $\Omega_t$ over the time interval $I$:

$$
\rho_f \frac{\partial \mathbf{u}_f}{\partial t} \bigg|_{x^*} - \operatorname{div}_x(2\mu_f \mathbf{D}_x(u_f)) + \rho_f((\mathbf{u}_f - w_f) \cdot \nabla_x) \mathbf{u}_f + \nabla_x p_f = \mathbf{f}, \quad \text{in } \Omega_t \times I \quad (10)
$$

$$
\operatorname{div}_x(u_f) = 0, \quad \text{in } \Omega_t \times I \quad (11)
$$

where all differential operators are defined w.r.t. the Eulerian coordinate system, except the ALE time derivative.

The fluid and the structure are coupled through a partitioned method with an implicit or semi-implicit scheme, see [1, 5] for more details.

4.1. 2D Benchmark

This final benchmark is a mix of the two previous tests. We remove the gravity and the flag part (not the cylinder) is now allowed to move. The results are in accordance with the reference values even though we used a really coarse mesh except for the drag. A complete study using high order approximation like for the fluid shall be available at the time of Acomen’11.

<table>
<thead>
<tr>
<th></th>
<th>$x \times 10^{-3}$</th>
<th>$y \times 10^{-3}$</th>
<th>Drag</th>
<th>Lift</th>
</tr>
</thead>
<tbody>
<tr>
<td>REF</td>
<td>$-2.69 \pm 2.53$</td>
<td>$1.48 \pm 34.38$</td>
<td>$457.3 \pm 22.66$</td>
<td>$2.22 \pm 149.78$</td>
</tr>
<tr>
<td></td>
<td>$-2.78 \pm 2.69$</td>
<td>$1.55 \pm 33.87$</td>
<td>$459.8 \pm 31.06$</td>
<td>$-2.19 \pm 174.45$</td>
</tr>
</tbody>
</table>

Table 3: Results for FSI3 with $\Delta t = 0.01$. There are 1226 elements in $\Omega_f$ and 26211 dofs associated ($P_4/P_3$), 260 elements in $\Omega_s$ and 2612 dofs associated ($P_3$). The geometry is first order and the BDF scheme for the fluid is order 2. REF values are found in [7]

4.2. 3D Benchmark

Finally we present a blood flow application in large arteries that has been proposed in [16]. The geometry is a straight pipe and a pressure pulse of $1.3332 \times 10^4 g/(\text{cm} s^2)$ has been imposed at the inlet boundary during 0.003s. The thin elastic vessel (0.1cm) is clamped at the inlet and outlet. Figure 6 shows the pressure wave propagation for different time steps. We have used a ($P_2/P_1$) space for the fluid and $P_1$ for the structure. The geometry for the fluid and structure is order one. The time scheme for the fluid is also order 1.

5. Conclusion

We have now a complete high order fluid-structure interaction framework in 2D and 3D. However much remains to be in various areas. Indeed we need to make a thorough study of our Navier-Stokes in moving domains framework in terms of approximations — the Arbitrary Lagrangian Eulerian in
Figure 5: Results for FSI3 with $\Delta t = 0.01$.

Figure 6: Pressure wave in a straight pipe. We show the fluid pressure and the fluid displacement of the pipe is magnified 15 times.
particular — as well the underlying linear algebra solvers updating the results of our previous paper [2].

We also have to compare various fluid-structure strategies as well as the wide range of space time and geometry approximations at our disposal to provide a complete overview.

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