

A NITSCHE-BASED FORMULATION FOR FLUID-STRUCTURE INTERACTIONS WITH CONTACT

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Abstract. We derive a Nitsche-based formulation for fluid-structure interaction (FSI) problems with contact. The approach is based on the work of Chouly and Hild (*SIAM J. Numer. Anal.* **51** (2013) 1295–1307) for contact problems in solid mechanics. We present two numerical approaches, both of them formulating the FSI interface and the contact conditions simultaneously in equation form on a joint interface-contact surface $\Gamma(t)$. The first approach uses a relaxation of the contact conditions to allow for a small mesh-dependent gap between solid and wall. The second alternative introduces an artificial fluid below the contact surface. The resulting systems of equations can be included in a consistent fashion within a monolithic variational formulation, which prevents the so-called “chattering” phenomenon. To deal with the topology changes in the fluid domain at the time of impact, we use a fully Eulerian approach for the FSI problem. We compare the effect of slip and no-slip interface conditions and study the performance of the method by means of numerical examples.

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1. INTRODUCTION

Contact problems have to be considered in many physical processes in engineering, medicine and nature. To name only a few consider for example the contact of balls and races in roller bearings, While an extensive amount of literature exists for the numerical simulation of contact in a purely solid mechanics context (see *e.g.* Wohlmuth [65] for an overview), *i.e.* disregarding the gas or liquid that mostly lies between contacting structures, much less works can be found considering full fluid-structure interaction with contact. The flow between contacting surfaces might however be of great importance for the contact dynamics. In the example of heart valves, the pulsating blood flow is even the driving force that enables opening and closure [4, 49]. In the case of ball bearings, fluid forces in the lubricant between ball and bearing may have a significant influence on the performance and wear of the bearing [11, 50].

Contact between different structures is typically formulated by means of variational inequalities [65]. In the context of full fluid-structure interaction (FSI), first results and algorithms can be found using either an

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artificial penalty force [63] or Lagrange multipliers [4, 27, 54] to obtain a well-posed and computationally feasible variational formulation.

However, these approaches have certain drawbacks: The use of a penalty force prevents real contact. The force is typically an artificial force and involves the choice of penalty parameters. If they are chosen too small, the structures might overlap in a numerical simulation. If they are chosen too large, the contact dynamics might be significantly perturbed [33]. In the case of Lagrange multipliers, additional variables are introduced on the contacting surfaces and an *inf-sup* condition is needed to ensure the well-posedness of the system. To tackle the variational inequality numerically, classical methods are based on *active-set strategies* including additional loops, which means in particular that the system of equations has to be solved several times in each time step. There are several possibilities to overcome these issues. Instead of an *active-set strategy*, semi-smooth Newton methods can be used to solve the system of equations [45]. In order to avoid additional variables, dual mortar methods, for example based on biorthogonal Lagrange multipliers, have been introduced [46].

Another possibility is to use Nitsche's method [57] to incorporate the contact conditions. This approach is based on an equivalent re-formulation of the contact conditions in equality form, which goes back to Alart and Curnier [1] and has been used in many works for the iterative solution of variational inequalities since then (see e.g. [45, 46]). In the case of the contact of an elastic body with a wall, the mechanical contact conditions on the contact surface $\Gamma_C(t)$ read

$$d_n \leq 0, \quad \sigma_{s,n}(d) \leq 0, \quad \sigma_{s,n}(d)d_n = 0,$$

where d_n denotes the solid displacement in normal direction and $\sigma_{s,n}(d)$ is the normal stress component. It can be shown that these conditions are equivalent to the equality

$$\sigma_{s,n}(d) = -\frac{1}{\gamma} \max\{0, d_n - \gamma\sigma_{s,n}(d)\} \quad \text{on } \Gamma_C(t), \quad (1.1)$$

for arbitrary $\gamma > 0$ (see e.g. [1]). This equality can be incorporated weakly in the variational formulation using Nitsche's method. While this technique has originally been used as a solution technique for the iterative solution of variational inequalities, it has only recently been considered and analysed as a discretisation technique by Chouly and co-workers [20–22], including a complete analysis of the related discretisation errors.

The Nitsche approach has the advantage that it is fully consistent. Furthermore, no additional variables have to be introduced and no additional loop within each time-step is needed. Numerical convergence has been proven in a series of papers for friction-free and frictional contact [19, 22, 24]. Following these results, Burman and co-workers used the re-formulation (1.1) to derive a Galerkin Least Squares formulation in equality form for the obstacle problem [17] and a Galerkin Least Squares and a Lagrange multiplier formulation for membrane contact [18].

All these studies disregard however the fluid that usually lies between the contacting structures. The aim of the present paper is to derive a Nitsche-based formulation for FSI with contact. To this purpose, we will introduce Lagrange multipliers on the FSI interface which can then be eliminated following the arguments reported by Burman and Hansbo [16].

The mathematical modeling of contact in an FSI context brings along a further issue: It is unclear, whether the incompressible Navier–Stokes equations with standard boundary conditions are an appropriate model in the fluid part, when it comes to contact. Theoretical studies show, that for a smooth, rigid solid body, no contact with an exterior wall can happen, when no-slip conditions are used on the interface and the outer boundary of the fluid domain, see Hillairet [43] and Hesla [42] in 2 space dimensions and Hillairet *et al.* [36, 44] in 3 space dimensions. This changes, when slip- or Navier-slip conditions are used on both the interface and the wall [36] or when the boundary of the solid is non-smooth [35, 64]. Gerard-Varet and Hillairet [35] found in a model example that it comes to contact for a solid with a $C^{1,\alpha}$ -parametrised boundary for $\alpha < 1/2$, while no contact happens for $\alpha \geq 1/2$. In the context of fluid-structure interactions, the regularity of the solid boundary depends on the solid displacement d , for which such regularity can usually not be guaranteed.

For a full FSI problem with a thin-walled structure, a no-collision result has been shown by Grandmont and Hillairet [38] in the no-slip case. For an overview on further results regarding existence of fluid-structure

interaction problems, we refer to Grandmont *et al.* [39]. Recently, Muha and Čanić [56] showed the well-posedness of a fluid-structure interaction system with slip-conditions.

Motivated by these theoretical results, we will study both no-slip and slip conditions on the FSI interface $\Gamma(t)$ in this work. It will turn out that the latter transits naturally into a “no-friction” condition when it comes to contact, while the prior leads to frictional contact. In this work, we will therefore consider friction-free contact, when a slip-condition is used on $\Gamma(t)$ and the specific frictional contact condition that follows from the transition of the interface conditions, when a no-slip condition is used on $\Gamma(t)$. For recent works on the incorporation of different friction laws (in particular Coulomb and Tresca friction), we refer to Chouly and co-workers [19, 24]. Moreover, we will study only contact of a deformable elastic structure with a fixed and straight wall for simplicity. Efficient algorithms to treat contact between more complex structures can be found, for example, in Puso [60], Yang *et al.* [66] and Chouly *et al.* [23, 55].

Concerning the governing equations, we focus on linear model equations for the fluid and solid part, *i.e.* the incompressible Stokes equations in the fluid and linear elasticity in the solid sub-domain. These simplifications must be seen as a first step towards the derivation of Nitsche-based contact formulations for complex FSI-contact problems. We consider, however, the case of a moving interface, which is a major challenge from the numerical point of view and leads to a non-linear FSI system, already in absence of contact.

The change of topology in the fluid domain causes additional numerical difficulties. Standard numerical approaches as the Arbitrary Lagrangian Eulerian method are not able to deal with topology changes, as the map from the reference domain to the Eulerian domain degenerates necessarily in this situation. The dynamics shortly before the impact can only be handled robustly when a Eulerian description of the fluid equations is used.

In the last years, several numerical approaches have been developed that are able to deal with topology changes. The methods can be split into two categories, according to the coordinate systems that are used for the solid system: Fully Eulerian approaches, where also the structure equations are formulated in Eulerian coordinates [26, 29, 33, 41, 61]; and Euler–Lagrangian techniques, where Lagrangian coordinates are used for the solid equations [2, 8, 13, 27, 37, 51, 53, 58, 67], such as in the *Immersed Boundary* or *Immersed Finite Element* methods.

Regarding the *Euler–Lagrangian* techniques, one can further distinguish between methods using Lagrange multipliers for the coupling of fluid and structure [37, 51] and methods based on Nitsche-techniques [2, 13, 40, 49, 53]. For the latter, a theoretical stability and convergence analysis has been derived [13]. The reader is referred to Boilevin-Kayl *et al.* [9] for a comparative study on the accuracy of some of these approaches.

The FSI approach we use here is based on the monolithic *Fully Eulerian* approach [26, 29, 33, 41, 61]. As the complete system of equations is formulated in Eulerian coordinates, the incorporation of contact conditions is straight-forward by means of variational principles. While some of the early works in this context suffered from stability and accuracy issues, see *e.g.* Dunne [28], accurate and robust discretisation and stabilisation techniques have been developed recently [32–34, 41]. We remark, however, that the algorithms we derive to incorporate contact can be combined in a straight-forward way with different FSI coupling techniques, *e.g.* Fictitious Domain or Immersed Boundary methods.

Concerning discretisation, we allow both for unfitted and fitted finite element approaches. For the unfitted case, so-called “ghost penalty” stabilisations can be used to guarantee the coercivity of the system [12, 13]. In order to simplify the presentation, we concentrate on fitted discretisations in this work and will comment on the unfitted case in a remark. In the numerical examples at the end of this paper, we will use the fitted *locally modified finite element method* [32].

The remainder of this paper is organised as follows: In Section 2, we first introduce the equations and the contact model. Before tackling the full contact problem, we derive a variational formulation for a model problem of a virtual obstacle within the fluid domain in Section 3, where we already have to deal with inequalities, but without any topology changes in the fluid domain. Then, we investigate contact with an exterior wall in Section 4. We show a stability result in Section 5 and present detailed numerical studies in Section 6, investigating the

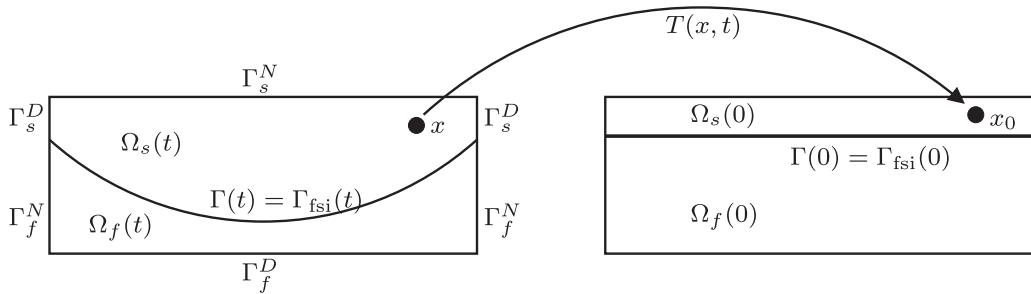


FIGURE 1. Illustration of an FSI problem without contact (left sketch). The domain affiliation in the current state can be determined by mapping back to the initial configuration, which is shown on the right.

influence of contact parameters, interface conditions and different contact formulations as well as convergence under mesh refinement. We conclude in Section 7.

2. MODEL

We begin by presenting the models for the fluid part, the solid part and the fluid-structure interaction on one hand in Section 2.1 and on the other hand the contact model in Section 2.2. For both models, Nitsche-based variational formulations are introduced. Different possibilities to combine the two models will then be presented in Sections 3 and 4.

2.1. Fluid-structure interaction without contact

We consider a fluid-structure interaction problem that is given on an overall domain $\Omega \subset \mathbb{R}^2$ which is split into a (variable) sub-domain $\Omega_f(t)$ occupied by a viscous fluid, a sub-domain $\Omega_s(t)$ occupied by an elastic solid and a lower-dimensional interface $\Gamma(t)$ separating them, such that

$$\Omega = \Omega_f(t) \cup \Gamma(t) \cup \Omega_s(t).$$

The boundary of the fluid domain is partitioned as follows $\partial\Omega_f(t) = \Gamma_{\text{fsi}}(t) \cup \Gamma_f^D \cup \Gamma_f^N$, where $\Gamma_{\text{fsi}}(t)$ stands for the fluid-solid interface. As regards the solid boundary, we assume that $\partial\Omega_s(t) = \Gamma(t) \cup \Gamma_s^D \cup \Gamma_s^N$, where the boundary part $\Gamma(t) = \Gamma_{\text{fsi}}(t) \cup \Gamma_C(t)$ is decomposed into the terms of $\Gamma_{\text{fsi}}(t)$ and the contact zone $\Gamma_C(t)$ (see Fig. 1 for a configuration without contact and the left sketch of Fig. 2 for a configuration with contact). The restriction to two dimensions is made only to simplify the presentation. The models and the methods derived in this paper can be generalised conceptually in a straight-forward way to three space dimensions.

In this work we will use a Eulerian description for the complete FSI problem. As already mentioned in the introduction this is not necessary for the contact algorithms derived below, but one convenient way to deal with (possible) topology changes in the fluid domain $\Omega_f(t)$. In an Eulerian description, the solid sub-domain and the interface are implicitly defined by the (unknown) solid displacement d

$$\Omega_s(t) = \{x \in \Omega \mid T(x, t) \in \Omega_s(0)\}, \quad \Gamma(t) = \{x \in \Omega \mid T(x, t) \in \Gamma(0)\}, \quad (2.1)$$

where $T : \Omega(t) \rightarrow \Omega$ is a bijective map, that is given by $T(x, t) = x - d(x, t)$ in the solid domain $\Omega_s(t)$ and by an arbitrary (smooth) extension in $\Omega_f(t) = \Omega \setminus (\Omega_s(t) \cup \Gamma(t))$. For the details, we refer to the textbook of Richter [62] or to Frei [30].

In the variable fluid sub-domain $\Omega_f(t)$, we consider the linear incompressible Stokes equations

$$\partial_t u - \operatorname{div} \sigma_f(u, p) = f_f, \quad \operatorname{div} u = 0,$$

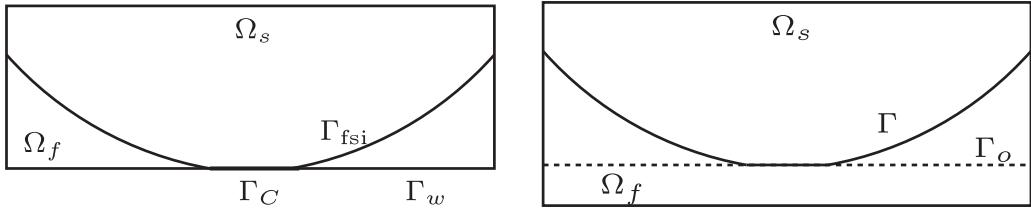


FIGURE 2. *Left:* body in contact with the wall Γ_w . *Right:* virtual obstacle line Γ_o within the fluid domain Ω_f .

where the Cauchy stress tensor σ_f is defined by

$$\sigma_f(u, p) = \nu_f (\nabla u^T + \nabla u) - pI,$$

u denotes the fluid velocity, p stands for pressure and $\nu_f > 0$ is a constant viscosity. In the solid-subdomain $\Omega_s(t)$, we assume a linear elastic material

$$\partial_t \dot{d} - \operatorname{div} \sigma_s(d) = f_s, \quad \partial_t d = \dot{d},$$

where the Cauchy stress tensor σ_s is given by

$$\sigma_s(d) = 2\mu_s E(d) + \lambda_s \operatorname{str}(E(d))I, \quad E(d) = \frac{1}{2} (\nabla d + \nabla d^T),$$

$\dot{d} = \partial_t d$ denotes the solid velocity and $\lambda_s, \mu_s > 0$ are positive constants.

For the coupling across the fluid-solid interface $\Gamma_{\text{fsi}}(t)$, the continuity of velocities and normal stresses

$$\dot{d} = u, \quad \sigma_f(u, p)n = \sigma_s(d)n \quad \text{on } \Gamma_{\text{fsi}}(t) \tag{2.2}$$

is typically considered for viscous fluids, where $n = n_s$ denotes the outer normal vector of the solid domain $\Omega_s(t)$. We recall that, since in this section there is no contact in the solid (*i.e.* $\Gamma_C(t) = \emptyset$), we have $\Gamma(t) = \Gamma_{\text{fsi}}(t)$. When it comes to contact, it is however questionable, whether the no-slip condition in (2.2) is still a reasonable approximation of the underlying physics, see the discussion on the no-collision paradox in the introduction. Therefore, we will study slip-conditions in this work as well (see Sect. 4.3), where the continuity across $\Gamma_{\text{fsi}}(t)$ is only imposed for the normal velocity

$$u \cdot n = \dot{d} \cdot n, \quad \tau^T \sigma_f n = 0, \quad \sigma_f n = \sigma_s n \quad \text{on } \Gamma_{\text{fsi}}(t). \tag{2.3}$$

In order to close the system of equations, we define exterior boundary conditions for the fluid and solid

$$\begin{aligned} u &= 0 \quad \text{on } \Gamma_f^D, \quad \sigma_f(u, p)n = 0 \quad \text{on } \Gamma_f^N \\ d &= 0 \quad \text{on } \Gamma_s^D, \quad \sigma_s(d)n = 0 \quad \text{on } \Gamma_s^N \end{aligned}$$

and the initial conditions

$$u(x, 0) = u^0(x) \text{ in } \Omega_f(0), \quad d(x, 0) = d^0(x), \quad \dot{d}(x, 0) = \dot{d}^0(x) \text{ in } \Omega_s(0).$$

We introduce the finite element spaces \mathcal{V}_h , \mathcal{Q}_h and \mathcal{W}_h on a quasi-uniform family of triangulations $(\mathcal{T}_h)_{h>0}$ and use Nitsche's method to combine the equations and interface conditions into a monolithic variational formulation (see *e.g.* [13, 40]). The monolithic system of equations reads in the no-slip case: *Find* $u(t) \in \mathcal{V}_h$, $p(t) \in \mathcal{Q}_h$, $d(t) \in \mathcal{W}_h$, *such that* $\dot{d} = \partial_t d$ and

$$\mathcal{A}_{\text{fsi, no-slip}}(u, p, d, \dot{d})(v, q, w) = (f_f, v)_{\Omega_f(t)} + (f_s, w)_{\Omega_s(t)} \quad \forall v, q, w \in \mathcal{V}_h \times \mathcal{Q}_h \times \mathcal{W}_h, \tag{2.4}$$

where

$$\begin{aligned} \mathcal{A}_{\text{fsi, no-slip}}(u, p, d, \dot{d})(v, q, w) := & (\partial_t u, v)_{\Omega_f(t)} + a_f(u, p; v, q) + (\partial_t \dot{d}, w)_{\Omega_s(t)} \\ & + (\sigma_s(d), \nabla w)_{\Omega_s(t)} - (T_f(u, p, \dot{d}), w - v)_{\Gamma(t)} - (\dot{d} - u, \sigma_f(v, -q)n)_{\Gamma(t)} \end{aligned} \quad (2.5)$$

and

$$a_f(u, p; v, q) := (\sigma_f(u, p), \nabla v)_{\Omega_f(t)} + (\operatorname{div} u, q)_{\Omega_f(t)} + S(p, q).$$

The numerical fluid traction on the interface is defined by $T_f(u, p, \dot{d}) := \sigma_f(u, p)n - \gamma_{\text{fsi}}(\dot{d} - u)$. The Nitsche parameter is chosen as $\gamma_{\text{fsi}} := \gamma_{\text{fsi}}^0 \nu_f h^{-1}$. The term $S(p, q)$ stands for a pressure stabilisation term that is non-zero in case that the discrete fluid spaces do not fulfil a discrete *inf-sup condition*. Note that at the FSI interface, we have used the following relation for the interface terms arising from integration by parts

$$(\sigma_s n, w)_{\Gamma(t)} - (\sigma_f n, v)_{\Gamma(t)} = (\sigma_f n, w - v)_{\Gamma(t)} + ([\![\sigma n]\!], w)_{\Gamma(t)}, \quad (2.6)$$

where we have dropped the dependencies of σ_f and σ_s for better readability. In the absence of contact, the jump of the stresses defined by

$$[\![\sigma n]\!] := \sigma_s n - \sigma_f n, \quad (2.7)$$

vanishes everywhere on $\Gamma(t)$. Furthermore, we have added the term

$$- (\dot{d} - u, \sigma_f(v, -q)n)_{\Gamma(t)}$$

as in Burman and Fernandez [13] for stability reasons.

Using a slip-condition at the FSI interface, the variational formulation reads: *Find $u \in \mathcal{V}_h, p \in \mathcal{Q}_h, d \in \mathcal{W}_h$, such that $\dot{d} = \partial_t d$ and*

$$\mathcal{A}_{\text{fsi, slip}}(u, p, d, \dot{d})(v, q, w) = (f_f, v)_{\Omega_f(t)} + (f_s, w)_{\Omega_s(t)} \quad \forall v, q, w \in \mathcal{V}_h \times \mathcal{Q}_h \times \mathcal{W}_h,$$

where

$$\begin{aligned} \mathcal{A}_{\text{fsi, slip}}(u, p, d, \dot{d})(v, q, w) := & (\partial_t u, v)_{\Omega_f(t)} + a_f(u, p; v, q) + (\partial_t \dot{d}, w)_{\Omega_s(t)} + (\sigma_s(d), \nabla w)_{\Omega_s(t)} \\ & - (T_f(u, p, \dot{d}) \cdot n, (w - v) \cdot n)_{\Gamma(t)} - ((\dot{d} - u) \cdot n, n^T \sigma_f(v, -q)n)_{\Gamma(t)}. \end{aligned} \quad (2.8)$$

2.2. Contact model without fluid

We assume that the solid is at a positive distance to the boundary at initial time and that contact can only happen with the lower wall (see Fig. 2, left sketch, where the situation at contact is shown)

$$\Gamma_w = \{(x_1, x_2) \in \partial\Omega \mid x_2 = 0\}.$$

We denote the outer normal vector of the fluid domain $\Omega_f(t)$ at Γ_w by $n_w = -e_2$. Moreover, let $g_0(x_2) > 0$ be the function describing the initial distance of a point $(x_1, x_2) \in \Gamma(t)$ to the wall Γ_w .

When contact with Γ_w occurs on a part $\Gamma_C(t) \subset \Gamma(t)$, suitable contact conditions are (see e.g. [1, 45])

$$d \cdot n_w \leq g_0, \quad \sigma_{s, n_w} := n_w^T \sigma_s n_w \leq 0, \quad (d \cdot n_w - g_0) \sigma_{s, n_w} = 0 \quad \text{on } \Gamma(t). \quad (2.9)$$

The first inequality in (2.9) ensures that the solid can not pass through Γ_w , the second inequality describes that the normal stress is zero (in the absence of contact) or negative (during contact) and the third condition is a complementarity condition that guarantees that at least one of the inequalities is “active”.

For arbitrary $\gamma_C > 0$ the first line in (2.9) is equivalent to (see *e.g.* [1])

$$\sigma_{s,n_w}(d) = -\gamma_C[P_\gamma(\sigma_{s,n_w}(d), d)]_+, \quad (2.10)$$

where $[f]_+ := \max\{f, 0\}$ and

$$P_\gamma(\lambda, d) := d \cdot n_w - g_0 - \frac{1}{\gamma_C} \lambda. \quad (2.11)$$

It remains to specify a contact condition for the tangential stresses. The simplest possibility is to consider frictionless contact

$$\tau_w^T \sigma_s n_w = 0 \quad \text{on } \Gamma_C(t). \quad (2.12)$$

Choosing $\gamma_C = \gamma_C^0 \mu_s h^{-1}$, the variational formulation reads: *Find* $d(t) \in \mathcal{W}_h$ *such that* $\dot{d} = \partial_t d$ *and*

$$(\partial_t \dot{d}, w)_{\Omega_s(t)} + (\sigma_s(d), \nabla w)_{\Omega_s(t)} + \gamma_C ([P_\gamma(\sigma_{s,n_w}(d), d)]_+, w \cdot n_w)_{\Gamma(t)} = (f_s, w)_{\Omega_s(t)} \quad \forall w \in \mathcal{W}_h. \quad (2.13)$$

We will discuss the tangential contact conditions in the context of fluid-structure interaction with contact below.

We close this section by mentioning that Chouly *et al.* [22] proposed a more general contact formulation that makes use of the consistency of the term

$$(\gamma_C[P_\gamma(\sigma_{s,n_w}(d), d)]_+ + \sigma_{s,n_w}(d), \sigma_{s,n_w}(w))_{\Gamma(t)}.$$

For $\theta \in [-1, 1]$, the contact term can be generalised to

$$\begin{aligned} & \gamma_C([P_\gamma(\sigma_{s,n_w}(d), d)]_+, w \cdot n_w)_{\Gamma(t)} - \theta (\gamma_C[P_\gamma(\sigma_{s,n_w}(d), d)]_+ + \sigma_{s,n_w}(d), \sigma_{s,n_w}(w))_{\Gamma(t)} \\ &= (1 - \theta) \gamma_C ([P_\gamma(\sigma_{s,n_w}(d), d)]_+, w \cdot n_w)_{\Gamma(t)} + \theta (\gamma_C[P_\gamma(\sigma_{s,n_w}(d), d)]_+, P_\gamma(\sigma_{s,n_w}(w), w))_{\Gamma(t)} \\ & \quad - \theta (\sigma_{s,n_w}(d), \sigma_{s,n_w}(w))_{\Gamma(t)}. \end{aligned} \quad (2.14)$$

For $\theta = 0$, we recover the formulation (2.13). Besides that, the case $\theta = 1$ is of particular interest, as it yields a symmetric formulation, for which a stability result has been shown [21].

3. VIRTUAL OBSTACLE WITHIN THE FLUID DOMAIN

Before considering full fluid-structure interaction with contact, we present our approach for a simplified model problem, where the movement of the solid is constraint by a virtual obstacle within the fluid domain, which is invisible to the fluid. This setting allows us to consider the numerical treatment of the interface conditions without accounting for the issues related to topology changes within the fluid. In Section 4, we will then show how this numerical setting can be extended to model contact of the solid with an exterior wall.

The setting of the model problem is shown on the right sketch of Figure 2. The fluid domain contains a horizontal obstacle line Γ_o which is invisible to the fluid, but an obstacle to the solid. One may consider for example a perfectly rigid thin membrane, that is laterally fixed and perfectly permeable for the fluid, but not for the solid, or a magnetic field below the obstacle that prevents the solid from crossing the line. We assume for simplicity that the obstacle line is parallel to the fluid boundary and denote its distance by $\alpha > 0$. The initial distance of the FSI interface $\Gamma(0)$ to the obstacle line is then given by

$$g_\alpha(x_2) := g_0(x_2) - \alpha.$$

We assume that $g_\alpha \geq 0$.

When considering fluid-structure interactions, the body is pre-loaded before it reaches the obstacle by means of the balance of normal forces

$$\sigma_s n = \sigma_f n \quad \text{on } \Gamma(t). \quad (3.1)$$

If the interface $\Gamma(t)$ reaches the obstacle Γ_o , the additional constraint $d \cdot n_w \leq g_\alpha$ has to be fulfilled. This gives rise to an additional surface force, that acts in the direction $-n_w = e_2$ (normal to Γ_o)

$$\llbracket \sigma n \rrbracket - \lambda n_w = 0 \quad \text{on} \quad \Gamma(t). \quad (3.2)$$

Note that we distinguish between the normal and tangential vectors of the wall n_w, τ_w and the normal and tangential vector of the FSI interface n, τ . The vectors are equal on $\Gamma_C(t)$, where the body is in contact with the wall, but different in the absence of contact.

The variable λ defined through (3.2) can be seen as a Lagrange multiplier, which is zero in absence of contact with the obstacle line (due to (3.1)) and can become negative during contact. This is described by the complementarity conditions

$$d \cdot n_w \leq g_\alpha, \quad \lambda \leq 0, \quad (d \cdot n_w - g_\alpha)\lambda = 0. \quad (3.3)$$

Splitting into normal and tangential contributions, (3.2) reads

$$n_w^T \llbracket \sigma n \rrbracket - \lambda = 0, \quad \tau_w^T \llbracket \sigma n \rrbracket = 0 \quad \text{on} \quad \Gamma(t). \quad (3.4)$$

Using the same trick as in Section 2.2, (3.3) is equivalent to

$$\lambda = -\gamma_C [P_\gamma(\lambda, d)]_+ \quad \text{on} \quad \Gamma(t), \quad (3.5)$$

with P_γ defined in (2.11).

Eliminating the Lagrange multiplier by using (3.4), this reads

$$\lambda = -\gamma_C [P_\gamma(\llbracket \sigma_n(u, p, d) \rrbracket, d)]_+ \quad \text{on} \quad \Gamma(t), \quad (3.6)$$

where we have used the abbreviations

$$\sigma_{s,n} := n_w^T \sigma_s n, \quad \sigma_{f,n} := n_w^T \sigma_f n, \quad \llbracket \sigma_n \rrbracket := n_w^T \llbracket \sigma n \rrbracket.$$

The natural formulation in the discrete setting is to consider the numerical stresses T_f in (2.7) and in (3.2). Let us derive these conditions first for the no-slip case. Adding the additional surface force to the FSI-Nitsche formulation (2.4), the discrete variational formulation reads

$$\mathcal{A}_{\text{FSI, no-slip}}(u, p, d, \dot{d})(v, q, w) - (\lambda n_w, w)_{\Gamma(t)} = (f_f, v)_{\Omega_f(t)} + (f_s, w)_{\Omega_s(t)} \quad \forall v, q, w \in \mathcal{V}_h \times \mathcal{Q}_h \times \mathcal{W}_h. \quad (3.7)$$

Due to the additional Nitsche interface terms in (3.2), this formulation includes the interface condition

$$\llbracket \tilde{\sigma}_n(u, p, d) \rrbracket - \lambda = 0 \quad \text{on} \quad \Gamma(t), \quad (3.8)$$

where the numerical stress jump in the FSI-Nitsche formulation across the interface is given by

$$\llbracket \tilde{\sigma}_n(u, p, d) \rrbracket := \sigma_{s,n}(d) - T_f(u, p, \dot{d}) \cdot n_w. \quad (3.9)$$

Eliminating the Lagrange multiplier by means of (3.8), the identity (3.5) reads

$$\lambda = -\gamma_C [P_\gamma(\llbracket \tilde{\sigma}_n(u, p, d) \rrbracket, d)]_+ \quad \text{on} \quad \Gamma(t). \quad (3.10)$$

For the definition of the numerical stresses in the slightly more complicated case of slip-interface conditions, we refer to Section 4.3. We will in the following analyse both contact formulations (3.6) and (3.10) and in particular their effect on the weakly imposed interface conditions. In order to avoid too much repetition, we use a general formulation that includes the Lagrange multiplier $\lambda = \lambda(u, p, d)$, keeping in mind that $\lambda(u, p, d)$ will be either chosen as the jump of normal stresses $\llbracket \sigma_n(u, p, d) \rrbracket$ or the jump of normal discrete stresses $\llbracket \tilde{\sigma}_n(u, p, d) \rrbracket$.

Using (3.5), the variational formulation reads:

Variational Formulation 1. Find $u(t) \in \mathcal{V}_h, p(t) \in \mathcal{Q}_h, d(t) \in \mathcal{W}_h$ such that $\dot{d} = \partial_t d$ and

$$\begin{aligned} \mathcal{A}_{\text{fsi},*}(u, p, d, \dot{d})(v, q, w) + \gamma_C ([P_\gamma(\lambda, d)]_+, w \cdot n_w)_{\Gamma(t)} \\ = (f_f, v)_{\Omega_f(t)} + (f_s, w)_{\Omega_s(t)} \quad \forall v, q, w \in \mathcal{V}_h \times \mathcal{Q}_h \times \mathcal{W}_h, \end{aligned} \quad (3.11)$$

where the bilinear form is one of the forms $\mathcal{A}_{\text{fsi, no-slip}}$ or $\mathcal{A}_{\text{fsi, slip}}$ defined in (2.5) and (2.8), respectively, and the contact parameter is chosen $\gamma_C = \gamma_C^0 \mu_s h^{-1}$, as in [20].

3.1. Weakly imposed interface conditions

Let us now analyse which interface conditions on $\Gamma(t)$ are implicitly included in Variational Formulation 1. For simplicity, we analyse the formulation with the bilinear form $\mathcal{A}_{\text{fsi, no-slip}}$ corresponding to no-slip conditions. Therefore, we integrate by parts in (3.11) and consider only the interface terms by formally neglecting the bulk and inter-element terms. For better readability, we drop all the dependencies of σ_f and σ_s . Collecting all terms with the fluid test function v , we obtain

$$-\sigma_f n + \sigma_f n - \gamma_{\text{fsi}}(\dot{d} - u) = -\gamma_{\text{fsi}}(\dot{d} - u) = 0 \quad \text{on } \Gamma(t), \quad (3.12)$$

i.e. the kinematic condition $\dot{d} = u$. Next, we collect the interface terms for the solid part w and split into a normal part ($w \cdot n_w$) and a tangential part ($w \cdot \tau_w$). We recall that since the boundary Γ_w is flat, the extension to $\Gamma(t)$ of its tangential and normal vectors are trivial. For the tangential part, we obtain as usual for Nitsche-based FSI

$$\tau_w^T \sigma_s n - \tau_w^T \sigma_f n + \gamma_{\text{fsi}}(\dot{d} - u) \cdot \tau_w = 0 \quad \text{on } \Gamma(t). \quad (3.13)$$

For the normal part, we have

$$\sigma_{s,n} - \sigma_{f,n} + \gamma_{\text{fsi}}(\dot{d} - u) \cdot n_w + \gamma_C [P_\gamma(d, \lambda)]_+ = 0 \quad \text{on } \Gamma(t).$$

Let us first consider the case that the contact force is not active. We obtain, as in the standard FSI-Nitsche formulation

$$[\![\sigma_n]\!] + \gamma_{\text{fsi}}(\dot{d} - u) \cdot n_w = 0.$$

If the contact force is active, we get

$$0 = [\![\sigma_n]\!] + \gamma_{\text{fsi}}(\dot{d} - u) \cdot n_w + \gamma_C (d \cdot n_w - g_\alpha) - \lambda = \begin{cases} \gamma_C (d \cdot n_w - g_\alpha) + \gamma_{\text{fsi}}(\dot{d} - u) \cdot n_w, & \lambda = [\![\sigma_n]\!] \\ \gamma_C (d \cdot n_w - g_\alpha), & \lambda = [\![\tilde{\sigma}_n]\!] \end{cases}. \quad (3.14)$$

In the first case, this is a combination of the “active” contact condition $d \cdot n_w = g$ and the continuity of velocities. As the continuity of velocities is imposed from the fluid side (3.12), this is not an issue for the model problem considered here. It will however lead to problems, when we consider contact of the solid with the lower wall Γ_w in Section 4. There, the second formulation, based on the discrete stresses (3.8) will be needed. As can be seen in (3.14), the pure contact condition $d \cdot n_w = g$ is valid from the solid side during contact.

4. FLUID-STRUCTURE INTERACTION WITH CONTACT

In this section, we consider contact of the solid body with the lower fluid boundary Γ_w . The formulation of a macroscopically relevant computational formulation without introducing the full physical model of all the interacting scales and elements, which is too costly in many applications, is an open problem to our knowledge and rarely discussed in literature.

When the solid enters into contact with a rigid wall, the action of the wall on the solid is expressed as a constraint, giving rise to a variational inequality. In principle there is no longer a fluid between the contacting

solids (“dry contact”). However, the presence of the fluid can not be discarded in this zone, since otherwise contact could be released in the interior of the contact zone through deformation of the elastic solid, thus creating a void, without neither fluid nor solid. A solution to this problem is to design a model where a fluid pressure remains in the contact zone, either on the boundary or in the bulk domain and prevents unphysical lift off creating vacuum. Our objective in the present work is to consider finite element formulations for contact between an elastic solid and a rigid wall assuming the existence of such a pressure model.

For the purpose of discussion, and without claiming that these models are optimal, we consider two simple models of seepage that introduce a fluid pressure in the contact zone. The first is a rather widespread approach which consists in relaxing the contact condition by introducing a small gap (see *e.g.* [4, 49]), so that the contact in the computational model takes place at an ϵ -distance from the wall (Sect. 4.1). Thus allowing for a thin fluid layer to remain. Here ϵ can be made to go to zero with reduced mesh-size. The second model consists in allowing the bulk fluid to penetrate into the solid wall in the form of a Navier–Stokes porous medium model where the permeability is driven to zero with reduced mesh-size (Sect. 4.2).

Our main objective is then to design a computational method that allows for contact in a consistent fashion, exploiting the presence of the modelled fluid contact pressure to get a smooth transition from contact to no contact without the appearance of non-physical inclusions or oscillations (“chattering”). To this end we study a nonlinear Nitsche-type coupling combining in a unified way the FSI coupling and the contact conditions.

4.1. Relaxed contact formulation

The idea of the relaxed formulation is to impose contact using a virtual barrier Γ_ϵ for the solid at a very small distance ϵ to Γ_w . This may be used for instance when the numerical discretisation of all the scales of the macroscopic model is unfeasible. We consider two examples.

If the solid has some fine scale structure on an unresolved scale ϵ , for example due to surface roughness, it is reasonable to consider that contact already takes place when the distance between the two structures is $O(\epsilon)$. Since by the modelling assumption the elastic solid can not penetrate into the rigid solid, the total forces from the fluid captured in the ϵ -layer and of the rigid solid must be sufficient to prevent penetration. This gives rise to equation (3.8), where λ represents the additional force coming from the rigid solid below Γ_ϵ that is necessary to prevent penetration.

On the other hand, as noted in the introduction, the continuous formulation does not allow for contact, when no-slip conditions are used on at least one of the interface and/or the boundary, due to the Navier–Stokes contact paradox. To reproduce this phenomenon in numerical simulations, an extremely high mesh resolution in the remaining small fluid layer between solid and rigid wall is typically required, in order to approximate the fluid tensor σ_f , which might become singular for $h \rightarrow 0$. In many situations, this might exceed the computational limits and moreover, it is *a priori* not clear how fine the mesh needs to be. In principle the condition $\sigma_{s,n} = \sigma_{f,n}$ is valid in the continuous formulation in this situation. As the singular behaviour of $\sigma_{f,n}$ might however not be accurately captured on coarser grids, it makes sense to re-inforce the fluid force by the contact force $[P_\gamma]_+$, if the line Γ_ϵ is passed, which indicates a failure of the force $\sigma_{f,n}$ to prevent contact.

In both these situations our idea to impose a virtual barrier Γ_ϵ for the solid at a very small distance ϵ to Γ_w permits numerical simulations already on reasonably coarse grids. In the numerical examples we will place the obstacle line at a distance $\alpha = \epsilon(h)$ to Γ_w , see Figure 3 on the left. We assume that $\epsilon(h) \rightarrow 0$, as the mesh size h tends to zero. The variational formulation is given by (3.11), where $\alpha = \epsilon(h)$, *i.e.*

$$g_\alpha(x_2) = g_{\epsilon(h)}(x_2) = g_0(x_2) - \epsilon(h).$$

Besides its simplicity, the main advantages of this contact formulation are:

- The numerical difficulties related to a topology change of the fluid domain are simplified.
- No-slip conditions can be used on both $\Gamma(t)$ and the lower wall Γ_w .

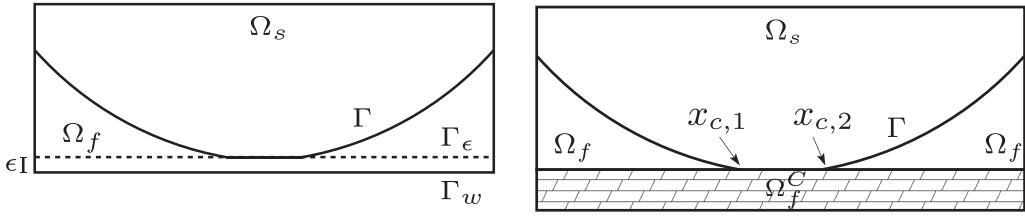


FIGURE 3. Illustration of the two approaches to include contact. *Left:* relaxed contact formulation: Body in contact with an obstacle line Γ_ϵ close to Γ_w . *Right:* introduction of an artificial fluid domain Ω_f^C below Γ_w .

4.2. Contact formulation using an artificial fluid

As a second possibility, we add an artificial fluid domain Ω_f^C below Γ_w , see the right sketch of Figure 3, and allow for a small, asymptotically for $h \rightarrow 0$ vanishing penetration into Ω_f^C . For the extension, we use a Stokes equation in combination with a penalty for the velocity u that drives the penetration to zero with $h \rightarrow 0$. This can be seen as a porous medium model with asymptotically vanishing permeability.

The variational formulation reads:

Variational Formulation 2. Find $u \in \mathcal{V}_h, p \in \mathcal{Q}_h, d \in \mathcal{W}_h$ such that $\dot{d} = \partial_t d$ and

$$\begin{aligned} \mathcal{A}_{FSI,*}^C(u, p, d, \dot{d})(v, q, w) &+ \gamma_C ([P_\gamma(\lambda, d)]_+, w \cdot n_w)_{\Gamma(t)} \\ &= (f_f, v)_{\Omega_f(t)} + (f_s, w)_{\Omega_s(t)} \quad \forall v, q, w \in \mathcal{V}_h \times \mathcal{Q}_h \times \mathcal{W}_h, \end{aligned} \quad (4.1)$$

where in the no-slip case

$$\begin{aligned} \mathcal{A}_{fsi,no-slip}^C(u, p, d, \dot{d})(v, q, w) &:= (\partial_t u, v)_{\Omega_f(t) \cup \Omega_f^C} + a_f^C(u, p; v, q) + \gamma_a(u, v)_{\Omega_f^C} + (\partial_t \dot{d}, w)_{\Omega_s(t)} \\ &\quad + (\sigma_s(d), \nabla w)_{\Omega_s(t)} - (T_f(u, p, \dot{d}), w - v)_{\Gamma(t)} - (\dot{d} - u, \sigma_f(v, -q)n)_{\Gamma(t)} \end{aligned} \quad (4.2)$$

and for slip interface conditions

$$\begin{aligned} \mathcal{A}_{fsi,slip}^C(u, p, d, \dot{d})(v, q, w) &:= (\partial_t u, v)_{\Omega_f(t) \cup \Omega_f^C} + a_f^C(u, p; v, q) + \gamma_a(u, v)_{\Omega_f^C} + (\partial_t \dot{d}, w)_{\Omega_s(t)} + (\sigma_s(d), \nabla w)_{\Omega_s(t)} \\ &\quad - (T_f(u, p, \dot{d}) \cdot n, (w - v) \cdot n)_{\Gamma(t)} - ((\dot{d} - u) \cdot n, n^T \sigma_f(v, -q)n)_{\Gamma(t)}. \end{aligned} \quad (4.3)$$

Note the presence of the penalty term $\gamma_a(u, v)_{\Omega_f^C}$ within the artificial fluid, where $\gamma_a := \gamma_a^0 h^\beta$ for $\beta > 0$. The fluid bilinear form is defined in $\Omega_f(t) \cup \Omega_f^C$

$$a_f^C(u, p; v, q) := (\sigma_f(u, p), \nabla v)_{\Omega_f(t) \cup \Omega_f^C} + (\operatorname{div} u, q)_{\Omega_f(t) \cup \Omega_f^C} + S(p, q).$$

Remark 4.1 (Porous medium analogy). The penalisation used in Variational Formulation 2 corresponds to the so-called penalty approach that is sometimes used for the coupling of free flow and flow through porous medium [25, 48]. There, the Stokes equations in the fluid part Ω_f and the Darcy equations in the porous medium Ω_p (which corresponds to Ω_f^C) are formulated simultaneously in the whole domain $\Omega = \Omega_f \cup \Omega_p$ in the spirit of the volume penalty approach

$$\partial_t u - \mu \Delta u + \nabla p + \frac{\mu}{K} u \chi_{\Omega_p} = 0, \quad \nabla \cdot u = 0 \quad \text{in } \Omega,$$

where K denotes the permeability of the porous medium and χ_{Ω_p} is the characteristic function of the domain Ω_p . In this sense the artificial fluid in our approach can be seen as a porous medium with asymptotically vanishing permeability $K = \mathcal{O}(h^\beta)$. A mathematical justification of this penalisation has been given by Angot [3]. Below we will choose the parameter $\beta = 2$, which is optimal in terms of the conditioning of the system. Following the argumentation of Angot a larger choice of β is also reasonable, with an optimal convergence rate reached for $\beta = 4$.

4.2.1. Discussion of the weakly imposed interface conditions: The no-slip case

We have already derived the weakly imposed interface conditions for the no-slip case in Section 3.1. Here, however, we have to consider that the fluid below the contact line is artificial, and we should in particular make sure that there is no feedback from the artificial fluid to the solid. In other words, we want that the artificial fluid acts as a *slave* to the solid during contact. Owing to (3.14), this naturally motivates the choice of λ in terms of the numerical stresses.

First, we note that the continuity of velocities (3.12) is imposed from the fluid side, such that no feedback to the solid is included from this equation. Considering the contact condition for the normal contact (3.14), we obtain

$$\gamma_C(d \cdot n_w - g_0) + \gamma_{\text{fsi}}(\dot{d} - u) \cdot n_w = 0 \quad (4.4)$$

when choosing $\lambda = [\![\sigma_n]\!]$. Instead of the condition $d \cdot n_w = g_0$, this induces an influence from the artificial velocity u from Ω_f^C onto the solid displacement. Moreover, if u is driven to zero in Ω_f^C , $\dot{d} \cdot n_w$ goes to zero as well, which might prevent the body from releasing from contact. On the other hand, using the jump of fluxes $\lambda = [\![\tilde{\sigma}_n]\!]$, we obtain the “pure” contact condition

$$d \cdot n_w = g_0 \quad (4.5)$$

as desired.

The weakly imposed tangential contact condition (3.13) reads

$$\tau_w^T \sigma_s n + \gamma_{\text{fsi}} \dot{d} \cdot \tau_w = 0, \quad (4.6)$$

when considering that the fluid velocity u is driven to zero asymptotically. As $\gamma_{\text{fsi}} \rightarrow \infty$ for $h \rightarrow 0$, this means (asymptotically) that the solid is not allowed to slide along the line Γ_w . While this might seem restrictive at first sight, this condition is in fact in some sense inherited from the no-slip condition at $\Gamma(t)$ before contact. This is due to the continuity of velocities on $\Gamma_{\text{fsi}}(t)$ and the fact that the velocity is driven to zero in Ω_f^C (and hence on Γ_w). Moreover, the no-slip condition on the fluid part of Γ_w implies that the solid can not slide on the end points $x_{C,1}$ and $x_{C,2}$ of the contact interval (see Fig. 3, right sketch). Altogether, this shows in agreement with a number of theoretical works (*e.g.* [36]) that the no-slip interface conditions are not an appropriate model for the case that it comes to contact with an exterior wall.

Remark 4.2 (Relation of the two contact formulations). The relaxed contact formulation derived in Section 4.1 can also be seen as an extension of the fluid forces σ_f to a region below the contact line (here Γ_ϵ), namely by using the Stokes equations in the extended domain. We will see in the numerical examples below that the two approaches yield similar results. In this way the use of the relaxed contact formulation might be justified also in cases, where real contact with the wall is expected. Moreover, it would be enough to use the extension in the artificial fluid approach only in a small layer of size $\mathcal{O}(\epsilon)$.

Remark 4.3 (Lagrange multiplier formulation). A further possibility would be to keep the Lagrange multiplier λ in the variational formulation (3.7) and (3.8) and to discretise additionally the Lagrange multiplier space, as in [37, 51] for the pure FSI case. Due to the difficulties concerning the discrete *inf-sup stability* and the additional computational effort, we will, however, not consider this alternative in the remainder of this work.

4.3. Slip conditions

Motivated by the above considerations, we study slip conditions in this section. The arguments that follow can be applied to the formulations proposed in Sections 3 and 4.

Observe that in the continuous case the discussion of (3.1) and (3.2) remains valid in the case of contact with slip conditions (noting that the tangential stresses vanish on both sides of the fluid-solid interfaces). However, at the discrete level, care has to be taken to use only the normal component of the numerical stress. Indeed, in this case the relation (3.2) becomes

$$(n^T \llbracket \sigma n \rrbracket) n - \lambda n_w = 0 \quad \text{on } \Gamma(t), \quad (4.7)$$

which yields $\lambda = (n^T \llbracket \sigma n \rrbracket)(n \cdot n_w) =: \llbracket \sigma_{n,\text{slip}} \rrbracket$.

At the discrete level, using the numerical stress this expression translates to

$$\lambda = (n^T \llbracket \tilde{\sigma} n \rrbracket)(n \cdot n_w) =: \llbracket \tilde{\sigma}_{n,\text{slip}} \rrbracket, \quad (4.8)$$

where the jump operator is given by (3.9). The resulting discrete formulations are given by Variational Formulation 1 or Variational Formulation 2, respectively, with the respective choice of λ .

4.3.1. Weakly imposed interface conditions: The slip case

Let us consider again which interface conditions are implicitly included in the variational formulation, when using the numerical stresses (4.8) for λ . Considering the interface terms with fluid test function v yields as usual the interface conditions $\tau^T \sigma_f n = 0$ and $(\dot{d} - u) \cdot n = 0$. Let us therefore concentrate on the terms with the solid test function w

$$(\sigma_s n, w)_{\Gamma(t)} - (T_f \cdot n, w \cdot n)_{\Gamma(t)} + \gamma_C([\tilde{P}_\gamma]_+, w \cdot n_w)_{\Gamma(t)} = 0.$$

In the case without contact, the last term vanishes and hence we retrieve the standard consistency of Nitsche's method for fluid-structure interaction with slip conditions. On the other hand, by developing the solid test function w in the local basis of $\Gamma(t)$ we have

$$(\tau^T \sigma_s n, w \cdot \tau)_{\Gamma(t)} + ((n^T \llbracket \tilde{\sigma} n \rrbracket), w \cdot n)_{\Gamma(t)} + \gamma_C([P_\gamma(d, \lambda)]_+, w \cdot n_w)_{\Gamma(t)} = 0.$$

Now, we use the identity $n = (n \cdot n_w)n_w + (n \cdot \tau_w)\tau_w$ to get

$$\begin{aligned} & (\tau^T \sigma_s n, w \cdot \tau)_{\Gamma(t)} + ((n^T \llbracket \tilde{\sigma} n \rrbracket), (w \cdot \tau_w)(n \cdot \tau_w))_{\Gamma(t)} + ((n^T \llbracket \tilde{\sigma} n \rrbracket), (w \cdot n_w)(n \cdot n_w))_{\Gamma(t)} \\ & \quad + \gamma_C([P_\gamma(d, \lambda)]_+, w \cdot n_w)_{\Gamma(t)} = 0. \end{aligned}$$

Hence, in the case of contact by using the definition of P_γ

$$(\tau^T \sigma_s n, w \cdot \tau)_{\Gamma(t)} + ((n^T \llbracket \tilde{\sigma} n \rrbracket), (w \cdot \tau_w)(n \cdot \tau_w))_{\Gamma(t)} + \gamma_C(d \cdot n_w - g_0, w \cdot n_w)_{\Gamma(t)} = 0. \quad (4.9)$$

When it comes to contact, we have in the asymptotic limit $\tau_w = \tau$ and $n \cdot \tau_w = 0$, so that (4.9) reduces to the "no-friction" condition $\tau^T \sigma_s n = 0$ and the non-penetration condition $d \cdot n_w = g_0$. This is the desired contact condition, as in the no-slip case, see (4.5).

We observe that in (4.9) both the tangential and the normal components (with respect to τ_w and n_w) have asymptotically vanishing perturbations. It is possible to eliminate the perturbation in the normal component, from the solid stress term $(\tau^T \sigma_s n, w \cdot \tau)_{\Gamma(t)}$, by adding the corresponding term to the definition of λ , *i.e.*

$$\lambda = \llbracket \tilde{\sigma}_{n,\text{slip}} \rrbracket + (\tau^T \sigma_s n)(\tau \cdot n_w). \quad (4.10)$$

In this case the relation (4.9) takes the form

$$(\tau^T \sigma_s n, w \cdot \tau_w(\tau \cdot \tau_w))_{\Gamma(t)} + ((n^T \llbracket \tilde{\sigma} n \rrbracket), (w \cdot \tau_w)(n \cdot \tau_w))_{\Gamma(t)} + \gamma_C(d \cdot n_w - g_0, w \cdot n_w)_{\Gamma(t)} = 0.$$

Testing with $w = n_w$ we see that here the non-penetration condition $d \cdot n_w = g_0$ is imposed unperturbed for all $h > 0$. By moving the perturbation to λ as in (4.10) it is instead the form P_γ that is perturbed, which implies a (weakly consistent) perturbation of the contact zone.

5. STABILITY

In this work, we will use fitted finite elements, *i.e.* we assume that both the interface $\Gamma(t)$ and the wall Γ_w are resolved by mesh lines. The approaches presented, however, can be combined with unfitted finite elements as well, with the only difference that some more stabilisation terms have to be added to the variational formulation. In order to simplify the presentation, we will concentrate on the fitted case first and discuss the extension to unfitted finite elements in a remark afterwards.

We will use equal-order finite elements in combination with a pressure stabilisation term $S_p(p, q)$ for the fluid equations. For the stability analysis, the only requirement on S_p is that it leads to a well-posed discrete fluid problem. Possibilities include the Brezzi–Pitkäranta stabilisation [10], local projections (LPS) [5], the pressure-stabilised Petrov–Galerkin approach (PSPG) [47] or the Continuous Interior Penalty method (CIP) [14].

In order to present the stability analysis in a general setting, we introduce some further notation splitting the contact force variable λ into a fluid part λ_f and a solid part λ_s , such that $\lambda = \lambda_s - \lambda_f$. For the no-slip case, we define $\lambda_s(w) = \sigma_{s,n}(w)$ and

$$\lambda_f(v, q, w) = \begin{cases} \sigma_{f,n}(v, q), & \text{if } \lambda = [\![\sigma_n]\!], \\ T_f(v, q, w) \cdot n, & \text{if } \lambda = [\![\tilde{\sigma}_n]\!], \end{cases} \quad (5.1)$$

see (3.4) and (3.8). For the slip case, we define $\lambda_s(w) = \sigma_{s,n}(w)(n \cdot n_w)$ and

$$\lambda_f(v, q, w) = \begin{cases} \sigma_{f,n}(v, q)(n \cdot n_w), & \text{if } \lambda = [\![\sigma_{n,\text{slip}}]\!], \\ (T_f(v, q, w) \cdot n)(n \cdot n_w), & \text{if } \lambda = [\![\tilde{\sigma}_{n,\text{slip}}]\!], \end{cases} \quad (5.2)$$

see (4.7) and (4.8).

5.1. Generalised contact formulation

Before we conduct the stability analysis, let us introduce a generalised contact formulation, following the ideas of Chouly *et al.* [22]. We have already briefly discussed their ideas for the case of a pure solid with contact in (2.14). The generalisation of (2.14) to the FSI-contact system (Variational Formulation 2) would be to add the terms

$$-(\gamma_C[P_\gamma(\lambda, d)]_+ + \lambda, \lambda(\partial_t v, \partial_t q, w))_{\Gamma(t)}. \quad (5.3)$$

The time derivatives on the test functions v and q are motivated by the stability analysis below, where we have to test the variational form with $v = u$, $q = p$ and $w = \partial_t d$, in order to show stability (see also Burman and Fernández [13]).

On the other hand, the term (5.3) is not usable within a time-stepping scheme due to the time derivatives on the test functions. A remedy is to shift the time derivatives to the first integrand (ignoring the boundary terms), *i.e.* adding the consistent terms

$$-(\gamma_C[P_\gamma(\lambda, d)]_+ + \lambda, \lambda_s(w))_{\Gamma(t)} - (\partial_t(\gamma_C[P_\gamma(\lambda, d)]_+ + \lambda), \lambda_f(v, q, w))_{\Gamma(t)}.$$

This yields the variational formulation:

Variational Formulation 3. Find $u(t) \in \mathcal{V}_h, p(t) \in \mathcal{Q}_h, d(t) \in \mathcal{W}_h$, such that $\dot{d} = \partial_t d$ and

$$\begin{aligned} \mathcal{A}(u, p, d)(v, q, w) &:= \mathcal{A}_{*,\text{FSI}}^*(u, p, d)(v, q, w) + \gamma_C([P_\gamma(\lambda, d)]_+, w \cdot n_w)_{\Gamma(t)} - \theta(\gamma_C[P_\gamma(\lambda, d)]_+ + \lambda, \lambda_s(w))_{\Gamma(t)} \\ &\quad - \theta(\partial_t(\gamma_C[P_\gamma(\lambda, d)]_+ + \lambda), \lambda_f(v, q, w))_{\Gamma(t)} \\ &= (f_f, v)_{\Omega_f(t)} + (f_s, w)_{\Omega_s(t)} \quad \forall v, q, w \in \mathcal{V}_h \times \mathcal{Q}_h \times \mathcal{W}_h, \end{aligned} \quad (5.4)$$

where $\mathcal{A}_{*,\text{FSI}}^*$ is one of the bilinear forms $\mathcal{A}_{\text{no-slip},\text{FSI}}$, $\mathcal{A}_{\text{slip},\text{FSI}}$, $\mathcal{A}_{\text{no-slip},\text{FSI}}^C$ or $\mathcal{A}_{\text{slip},\text{FSI}}^C$.

5.2. A stability result

In this section, we will investigate stability of the discrete formulation for different values of θ . In particular, we will show a stability result for the symmetric formulation ($\theta = 1$) and stability up to a specific term in the general case (including $\theta = 0$). These results correspond to the results that have been obtained by Chouly *et al.* [22] for the pure solid case. For the stability analysis, we will assume infinitesimal displacements, *i.e.* the sub-domains Ω_f and Ω_s as well as the interface Γ are fixed. We introduce the notation $\tilde{\Omega}_f = \Omega_f \cup \Omega_f^C$ for the combined fluid and artificial fluid domain and set $\Omega_f^C = \emptyset$ for the relaxed approach.

Theorem 5.1. *Let $u, p, d \in V_h$ and $\dot{d} = \partial_t d$. We have the following stability result for the form \mathcal{A} defined in (5.4), where $\theta \in [-1, 1]$ and γ_C^0 sufficiently large*

$$\begin{aligned} \|u(T)\|_{\tilde{\Omega}_f}^2 + \|\dot{d}(T)\|_{\Omega_s}^2 + \|d(T)\|_{H^1(\Omega_s)}^2 + \int_0^T \left(\nu_f \|\nabla u\|_{\tilde{\Omega}_f}^2 + S_p(p, p) + \gamma_a \|u\|_{\Omega_f^C}^2 \right) dt \\ + \int_0^T \gamma_{fsi} \|(\dot{d} - u) \cdot n\|_\Gamma^2 dt + \theta \left\| \gamma_C^{-1/2} \lambda(T) + \gamma_C^{1/2} [P_\gamma(\lambda, d)]_+(T) \right\|_\Gamma^2 \\ \leq C \left(\int_0^T \mathcal{A}(u, p, d; u, p, \dot{d}) - (1 - \theta) \gamma_C \left([P_\gamma(\lambda, d)]_+, \dot{d} \cdot n_w \right)_\Gamma dt + \|u_0\|_{\tilde{\Omega}_f}^2 \right. \\ \left. + \|\dot{d}_0\|_{\Omega_s}^2 + \|d_0\|_{H^1(\Omega_s)}^2 + \theta \left\| \gamma_C^{-1/2} \lambda_0 + \gamma_C^{1/2} [P_\gamma(\lambda_0, d_0)]_+ \right\|_\Gamma^2 \right), \end{aligned}$$

where we have used the abbreviation $\lambda_0 := \lambda(u_0, p_0, d_0)$. For the no-slip case, the term $\|(\dot{d} - u) \cdot n\|_\Gamma$ can be replaced by $\|(\dot{d} - u)\|_\Gamma$.

Remark 5.2 (Contact terms). The second line gives us control over the satisfaction of the FSI-contact condition for $\theta > 0$. In contrast to the work by Chouly *et al.* [22] for a pure solid problem, we obtain here discrete stability for $\theta = 1$, for the following positive discrete energy

$$E(T) := \|u(T)\|_{\tilde{\Omega}_f}^2 + \|\dot{d}(T)\|_{\Omega_s}^2 + \|d(T)\|_{H^1(\Omega_s)}^2 + \left\| \gamma_C^{-1/2} \lambda(T) + \gamma_C^{1/2} [P_\gamma(\lambda, d)]_+(T) \right\|_\Gamma^2.$$

For $\theta \neq 1$ on the other hand, the contact term $(1 - \theta) \gamma_C \left([P_\gamma(\lambda, d)]_+, \dot{d} \cdot n_w \right)_\Gamma$ appears on the right-hand side. For $\theta < 0$ we have additionally a negative contribution on the left-hand side. Both issues are directly inherited from the pure solid mechanical case, see [21]. The last term on the right-hand side vanishes, if we assume that the contact conditions are fulfilled at initial time, for example if the solid is not in contact with at $t = 0$.

Proof. We test (5.4) with $w = \dot{d} = \partial_t d, v = u$ and $q = p$ and integrate in time. We start by deriving a lower bound for $\mathcal{A}_{*,FSI}^*$. For the fluid part, we use the techniques from Burman and Fernandez [13], to show coercivity of the Stokes part including the coupling terms. For the no-slip case, the authors have shown that

$$\begin{aligned} a_f^C(u, p; u, p) + \gamma_a(u, u)_{\Omega_f^C} - (\sigma_f(u, p)n, \dot{d} - u)_\Gamma - (\dot{d} - u, \sigma_f(u, -p)n)_\Gamma + \gamma_{fsi} \|\dot{d} - u\|_\Gamma^2 \\ \geq c \left(\nu_f \|\nabla u\|_{\tilde{\Omega}_f}^2 + \gamma_{fsi} \|\dot{d} - u\|_\Gamma^2 + S_p(p, p) + \gamma_a \|u\|_{\Omega_f^C}^2 \right). \end{aligned}$$

Analogously, one can show in the slip-case that

$$\begin{aligned} a_f^C(u, p; u, p) + \gamma_a(u, u)_{\Omega_f^C} - (n^T \sigma_f(u, p)n, (\dot{d} - u) \cdot n)_\Gamma - ((\dot{d} - u) \cdot n, n^T \sigma_f(u, -p)n)_\Gamma + \gamma_{fsi} \|(\dot{d} - u) \cdot n\|_\Gamma^2 \\ \geq c \left(\nu_f \|\nabla u\|_{\tilde{\Omega}_f}^2 + \gamma_{fsi} \|(\dot{d} - u) \cdot n\|_\Gamma^2 + S_p(p, p) + \gamma_a \|u\|_{\Omega_f^C}^2 \right). \end{aligned}$$

Using the symmetry of σ_s , integration in time and a Korn's inequality, we obtain for the solid part

$$\begin{aligned} \int_0^T (\sigma_s(d), \nabla \dot{d})_{\Omega_s} dt &= \frac{1}{2} \int_0^T \partial_t (\sigma_s(d), \nabla d)_{\Omega_s} dt = \frac{1}{2} ((\sigma_s(d(T)), \nabla d(T))_{\Omega_s} - (\sigma_s(d(0)), \nabla d(0))_{\Omega_s}) \\ &\geq c_1 \|\nabla d(T)\|_{\Omega_s}^2 - c_2 \|\nabla d_0\|_{\Omega_s}^2. \end{aligned}$$

Moreover, we have

$$\int_0^T (\partial_t u, u)_{\tilde{\Omega}_f} + (\partial_t \dot{d}, \dot{d})_{\Omega_s} dt = \frac{1}{2} \left(\|u(T)\|_{\tilde{\Omega}_f}^2 + \|\dot{d}(T)\|_{\Omega_s}^2 - \|u_0\|_{\tilde{\Omega}_f}^2 - \|\dot{d}_0\|_{\Omega_s}^2 \right).$$

Together, we have shown that

$$\begin{aligned} \|u(T)\|_{\tilde{\Omega}_f}^2 + \|\dot{d}(T)\|_{\Omega_s}^2 + \|d(T)\|_{H^1(\Omega_s)}^2 + \int_0^T \nu_f \|\nabla u\|_{\tilde{\Omega}_f}^2 + S_p(p, p) + \gamma_a \|u\|_{\Omega_f^C}^2 + \gamma_{\text{fsi}} \|\dot{d} - u\|_{\Gamma}^2 dt \\ \leq C \left(\int_0^T \mathcal{A}_{*,FSI}^*(u, p, d; u, p, \dot{d}) dt + \|u_0\|_{\tilde{\Omega}_f}^2 + \|\dot{d}_0\|_{\Omega_s}^2 + \|d_0\|_{H^1(\Omega_s)}^2 \right). \end{aligned}$$

Let us now estimate the contact terms. We split the principal contact term into

$$\int_0^T \gamma_C ([P_\gamma(\lambda, d)]_+, \dot{d} \cdot n_w)_\Gamma dt = \int_0^T \theta \gamma_C ([P_\gamma(\lambda, d)]_+, \dot{d} \cdot n_w)_\Gamma + (1 - \theta) \gamma_C ([P_\gamma(\lambda, d)]_+, \dot{d} \cdot n_w)_\Gamma dt.$$

We have to estimate the terms

$$\begin{aligned} \theta \int_0^T \underbrace{\gamma_C ([P_\gamma(\lambda, d)]_+, \dot{d} \cdot n_w)_\Gamma}_{I_1} - \underbrace{(\gamma_C [P_\gamma(\lambda, d)]_+ + \lambda, \lambda_s(\dot{d}))_\Gamma}_{I_2} \\ - \underbrace{(\partial_t (\gamma_C [P_\gamma(\lambda, d)]_+ + \lambda), \lambda_f(u, p, \dot{d}))_\Gamma}_{I_3} dt. \end{aligned} \tag{5.5}$$

From the definition of P_γ we can write $d \cdot n_w = P_\gamma(\lambda, d) + \gamma_C^{-1} \lambda$. Hence, since the lower wall is assumed to be time independent, we have

$$\begin{aligned} \int_0^T I_1 dt &= \gamma_C \int_0^T ([P_\gamma(\lambda, d)]_+, \partial_t (P_\gamma(\lambda, d) + \gamma_C^{-1} \lambda))_\Gamma dt \\ &= \gamma_C \int_0^T ([P_\gamma(\lambda, d)]_+, \partial_t ([P_\gamma(\lambda, d)]_+ + \gamma_C^{-1} \lambda))_\Gamma dt. \end{aligned}$$

In the second line, we have used that [22]

$$\frac{1}{2} \partial_t [\phi]_+^2 = [\phi]_+ \partial_t [\phi]_+ = [\phi]_+ H(\phi) \partial_t [\phi]_+ = [\phi]_+ \partial_t \phi$$

where H is the Heaviside function. We insert $\pm\gamma_C^{-1}\lambda$ and integrate by parts

$$\begin{aligned} \int_0^T I_1 dt &= \gamma_C \int_0^T ([P_\gamma(\lambda, d)]_+ + \gamma_C^{-1}\lambda, \partial_t([P_\gamma(\lambda, d)]_+ + \gamma_C^{-1}\lambda))_\Gamma dt \\ &\quad - \int_0^T \underbrace{(\lambda, \partial_t([P_\gamma(\lambda, d)]_+ + \gamma_C^{-1}\lambda))_\Gamma}_{I_4} dt \\ &= \underbrace{\frac{\gamma_C}{2} \| [P_\gamma(\lambda, d)(T)]_+ + \gamma_C^{-1}\lambda(T) \|_{0,\Gamma}^2}_{D_T} - \underbrace{\frac{\gamma_C}{2} \| [P_\gamma(\lambda_0, d_0)]_+ + \gamma_C^{-1}\lambda_0 \|_{0,\Gamma}^2}_{D_0} - \int_0^T I_4 dt. \end{aligned}$$

Now the idea is to split the contribution from the term I_4 into fluid and solid stresses and to apply integrating by parts in time (only) in the solid stress contribution. By definition, we have

$$-\int_0^T I_4 dt = -\int_0^T (\lambda_s(d), \partial_t([P_\gamma(\lambda, d)]_+ + \gamma_C^{-1}\lambda))_\Gamma + (\lambda_f(u, p, \dot{d}), \partial_t([P_\gamma(\lambda, d)]_+ + \gamma_C^{-1}\lambda))_\Gamma dt$$

and, by integrating by parts in the first term of the right-hand side, we have

$$\begin{aligned} -\int_0^T I_4 dt &= \underbrace{\int_0^T (\lambda_s(\dot{d}), [P_\gamma(\lambda, d)]_+ + \gamma_C^{-1}\lambda)_\Gamma dt}_{I_5} - \underbrace{(\lambda_s(d(T)), [P_\gamma(\lambda, d)(T)]_+ + \gamma_C^{-1}\lambda(T))_\Gamma}_{I_6} \\ &\quad + \underbrace{(\lambda_s(d_0), [P_\gamma(\lambda_0, d_0)]_+ + \gamma_C^{-1}\lambda_0)_\Gamma}_{I_7} + \underbrace{\int_0^T (\lambda_f(u, p, \dot{d}), \partial_t([P_\gamma(\lambda, d)]_+ + \gamma_C^{-1}\lambda))_\Gamma dt}_{I_8}. \end{aligned}$$

The terms I_5 and I_8 cancel with the terms I_2 and I_3 in (5.5). The term I_6 is treated in a standard fashion using Young's inequality, an inverse inequality and the dissipation provided by D_T and the elastic energy $\frac{1}{2}a^s(d(T), d(T))$ for γ_C^0 sufficiently large. For $\theta > 0$, we estimate

$$\begin{aligned} I_6 &\geq -\gamma_C^{-1} \|\lambda_s(d(T))\|_{0,\Gamma}^2 - \frac{\gamma_C}{4} \| [P_\gamma(\lambda(T), d(T))]_+ + \gamma_C^{-1}\lambda(T) \|_{0,\Gamma}^2 \\ &\geq -\frac{1}{4} \|d(T)\|_{H^1(\Omega_s)}^2 - \frac{1}{2} D_T. \end{aligned}$$

In the same way, we get the upper bound

$$I_6 \leq \frac{1}{4} \|d(T)\|_{H^1(\Omega_s)}^2 + \frac{1}{2} D_T,$$

which is needed in the case $\theta < 0$. For I_7 we obtain analogously

$$-\frac{1}{4} \|d_0\|_{H^1(\Omega_s)}^2 - \frac{1}{2} D_0 \leq I_7 \leq \frac{1}{4} \|d_0\|_{H^1(\Omega_s)}^2 + \frac{1}{2} D_0,$$

which completes the proof. \square

Remark 5.3 (Unfitted finite elements). When using unfitted finite elements [13, 15], additional stabilisation terms S_u and S_d are needed, if the interface $\Gamma(t)$ is not resolved by mesh lines. Their purpose is to extend the coercivity of the fluid system from Ω_f (resp. Ω_s) to the extended domains Ω_h^f (resp. Ω_h^s) that consists of all element $T \in \mathcal{T}_h$, with a non-empty intersection with the respective sub-domain ($T \cap \Omega_i \neq \emptyset$). Suitable “ghost penalty” operators have been defined in Burman [12]. The same stability result as in Theorem 5.1 can then be shown with an analogous argumentation.

Remark 5.4 (Newton convergence). While the symmetric formulation ($\theta = 1$) seems beneficial from the theoretical point of view, the additional terms in (5.4) can cause severe difficulties for the non-linear solver. The reason is that the additional contact terms are not only highly non-linear, but also non-smooth, especially due to the time derivative acting on the maximum operator $\partial_t[\cdot]_+$. In our numerical tests, we were not able to obtain numerical convergence for different versions of generalised Newton methods [52]. The investigation of the case $\theta \neq 1$ by means of numerical tests and in particular the construction of a robust non-linear solver are subject to future research.

6. NUMERICAL RESULTS

In this section, we show some numerical results to analyse and to compare the different contact formulations. As mentioned in Remark 5.4, we were not able to obtain results for the generalised contact formulation with $\theta \neq 0$, due to divergence of the generalised Newton-type methods we have tried. Therefore, we only show results for $\theta = 0$, where Newton convergence was not an issue, at least when the time step δt was chosen reasonably small. Following the standard approach for contact in solid mechanics, we could in this case simply ignore the non-differentiability of the maximum operator when computing the Newton derivatives, as the term P_γ inside the bracket $[\cdot]_+$ is in practice typically never exactly zero. For all other values of P_γ the derivatives are well-defined. In the computations made for this paper, the Newton algorithm needed 1–2 iterations per time step to reduce the initial residual by a factor of 10^{-7} , if the contact force was not getting active during the iteration, and 1–5 iterations per time step in and around the interval of contact. This makes the method highly competitive in terms of computational costs compared to approaches using Lagrange multipliers and/or active-sets.

We first give some details in Section 6.1 on the fitted, equal-order finite element discretisation and the stabilisations we use. Then, in Section 6.2, we study the problem of a virtual obstacle within the fluid domain introduced in Section 3. The purpose of this example is to isolate the effect of the contact terms from issues related to discretisation during contact and the topology change in the fluid domain $\Omega_f(t)$. Then, we study in Section 6.3 a model problem with contact with the boundary of the fluid domain, where we compare among other aspects the two contact formulations introduced in Section 4, the different possibilities to choose the fluxes λ and the effect of slip and no-slip boundary and interface conditions.

6.1. Details on discretisation and stabilisation

For the numerical results in this paper, we will use a monolithic *Fully Eulerian* approach on a global mesh T_h covering $\Omega(t)$. In order to resolve the interface $\Gamma(t)$ within the discretisation, we use the *locally modified finite element method* introduced by Frei and Richter [32]. The idea of this approach is to use a fixed coarse triangulation T_{2h} of the overall domain $\Omega = \tilde{\Omega}_f(t) \cup \Gamma(t) \cup \Omega_s(t)$ that is independent of the position of the interface $\Gamma(t)$. Then, in each time step, this coarse grid is refined once by splitting each so-called “patch” element in either eight triangles or four quadrilaterals to resolve the interface in at least a linear approximation, see Figure 4 for an illustration.

The finite element space V_h is then defined as a combination of piece-wise linear and piece-wise bi-linear finite elements on the patches. It can be guaranteed that a maximum angle condition is fulfilled in each of the sub-cells, leading to optimal-order interpolation and error estimates [32].

For temporal discretisation, we split the time interval I into m equidistant-distant time intervals $I_j = (t_{j-1}, t_{j-1} + k]$ and use a time-stepping scheme that is based on a modified discontinuous Galerkin time discretisation of lowest order (dG(0)), see Frei and Richter [34]. The displacement-velocity relation $\partial_t d = \dot{d}$ is included by means of the L^2 -projection

$$(\partial_t d, z)_{\Omega_s(t)} - (\dot{d}, z)_{\Omega_s(t)} = 0 \quad \forall z \in V_h^s,$$

where V_h^s denotes the (modified) finite element space that is spanned by the degrees of freedom of the elements in the solid part Ω_h^s .

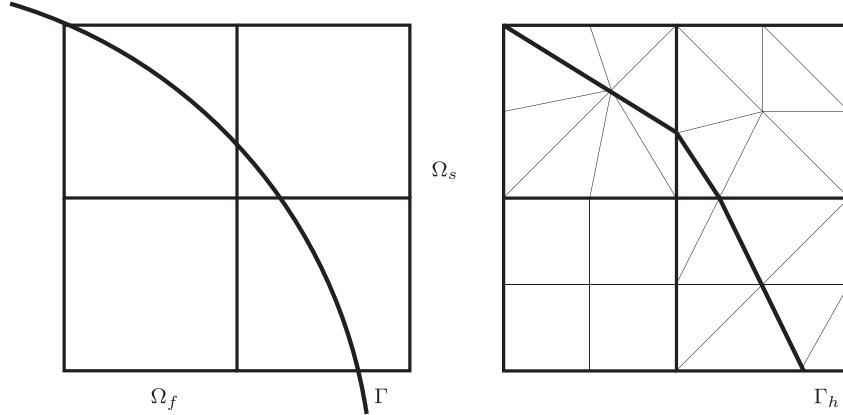


FIGURE 4. *Left:* fixed triangulation \mathcal{T}_{2h} of the domain Ω . *Right:* subdivision of the patches $P \in \mathcal{T}_{2h}$ such that the interface $\Gamma(t)$ is resolved in a linear approximation by the discrete interface Γ_h .

The domain affiliation of a point $x \in \Omega(t)$ is determined by means of the *Initial Point Set/Backward Characteristics method* [26, 29], that uses the displacement $d(t)$ in the solid domain and an extension to $\Omega_f(t)$ in order to trace back points to their initial position in $\Omega(0)$, following the definition (2.1).

For pressure stabilisation, we use an anisotropic variant of the Continuous Interior Penalty method, see Frei [30, 31]. In addition, we add the temporal pressure stabilisation term

$$S_{pt}(p, q) = \gamma_{pt} h(p^m - p^{m-1}, q)_{\Gamma(t)}$$

in each time interval I_m . This additional stabilisation is needed, as the mesh $\mathcal{T}_h(t_m)$, and hence the finite element spaces, change from time-step to time-step. The solution u^{m-1} from the previous time-step t_{m-1} is therefore not discrete divergence-free with respect to the new mesh $\mathcal{T}_h(t_m)$, which gives rise to pressure oscillations, see for example Besier and Wollner [7].

All the following results have been obtained using the finite element library *Gascoigne 3d* [6].

6.2. Virtual obstacle within the fluid domain

We begin by investigating the simplified problem introduced in Section 3. The initial fluid and solid domains are defined as

$$\Omega_f(0) = (0, 1) \times (0, 0.5), \quad \Omega_s(0) = (0, 1) \times (0.5, 0.6)$$

and a lower-dimensional obstacle $\Gamma_w = (0, 1) \times 0.25$ within the fluid domain. We consider a moving interface $\Gamma(t)$, which is resolved using the locally modified finite element method. The sub-domains $\Omega_s(t)$ and $\Omega_f(t)$ and the interface $\Gamma(t)$ depend on the solid displacement $d(t)$, see (2.1).

The constraint for the solid displacement is given by

$$d \cdot n_w \leq 0.25 \quad (= g_\alpha). \quad (6.1)$$

We use the elasticity parameters $\lambda_s = \mu_s = 2 \times 10^6$ and the fluid viscosity $\nu_f = 1$. The structure is pulled towards the bottom by fluid forces due to a prescribed pressure mean value at the left and right boundary of the fluid domain

$$\int_{\Gamma_{f,\text{left}}} p \, ds = \int_{\Gamma_{f,\text{left}}} \bar{P} \, ds, \quad \int_{\Gamma_{f,\text{right}}} p \, ds = \int_{\Gamma_{f,\text{right}}} \bar{P} \, ds$$

where $\bar{P} := 1.3 \times 10^5$. We consider the Variational Formulation 1 with no-slip conditions and $\lambda = [\![\tilde{\sigma}_n]\!]$, see (3.8), on a Cartesian mesh that consists of 5120 elements and with a small time step $\delta t = 10^{-5}$. The Nitsche constant at the FSI interface is chosen $\gamma_{\text{fsi}}^0 = 10^3$ and temporal pressure stabilisation with $\gamma_{pt} = 10^{-2}$ is used.

To analyse the results, we define the “minimal distance”

$$d_{\min} := \min_{x \in \Gamma(t)} x_2 - 0.25$$

of the interface $\Gamma(t)$ to Γ_w . To be precise the term “minimal distance” is only correct before contact, as d_{\min} gets negative in case of an overlap. Moreover, we define the following functionals in order to analyse the pressure p and the contact force at the interface $\Gamma(t)$

$$J_p := \left| \int_{\Gamma(t)} p \, ds \right|, \quad J_{P_\gamma} := \gamma_C \int_{\Gamma(t)} [P_\gamma(\lambda, d)]_+ \, ds.$$

In Figure 5, we plot these three functionals over time for two contact periods and the contact parameters $\gamma_C^0 = 10, 10^2, 10^3$. In the top left plot, we observe that the solid is pulled down until it reaches Γ_w at $t \approx 1.5 \times 10^{-3}$. After a short *contact* period, it is released again due to its elastic properties before it reaches the obstacle for a second time at $t \approx 5 \times 10^{-3}$.

The contact condition $d \cdot n_w \leq 0.25$ is only significantly violated for the smallest contact parameter, where d_{\min} reaches a minimum value of around -8×10^{-4} , see the zoom-in around the contact interval on the top left. This value is more than an order of magnitude smaller than the mesh size in vertical direction $h \approx 1.4 \times 10^{-2}$. For the larger values of γ_C^0 , the minimal value of d_{\min} is even much closer to zero. On the other hand, we observe that even for the largest value $\gamma_C^0 = 10^3$, the contact condition is slightly relaxed, allowing for very small overlaps of solid and contact line.

In the second row of Figure 5, we observe that the pressure shows a peak at the beginning ($t \approx 1.5 \times 10^{-3}$ and $t \approx 5 \times 10^{-3}$) of the contact periods, followed by some small oscillations. The peak is caused by the fluid dynamics and will be discussed below. The oscillations get smoother for larger values of γ_C^0 and are barely visible for $\gamma_C^0 = 10^3$.

Similarly, the contact force J_{P_γ} shows oscillations for $\gamma_C^0 = 10$ and a much smoother behaviour for $\gamma_C^0 \geq 10^2$. Note that this does not contradict the stability result in Theorem 5.1, where we have assumed that γ_C^0 is large enough. The relatively large value for γ_C^0 that is needed here is due to the anisotropic cells that appear in some of the time-steps, when using the *locally modified finite element method*. In the absence of extreme anisotropies a value of $\gamma_C^0 \approx 1$ seems to be enough to obtain stable numerical results. The optimal choice of the contact parameter γ_C in the context of anisotropic cells is subject to future research.

If γ_C^0 is chosen large enough, the contact force is roughly of the same size for different γ_C^0 . This is in agreement with the observations of Chouly *et al.* for the case of a pure solid problem [23], who showed that the consistency of the method makes the choice of the contact parameter much less sensitive compared to a pure penalty method.

6.2.1. Investigation of the pressure peak

The pressure peak at the beginning of the contact interval can be explained as follows. As the fluid does not “see” the obstacle before reaching it, the solid is pulled down towards it without reducing its velocity. At the moment when the obstacle is reached, its vertical velocity $d \cdot n_w$ has to decrease to zero in an instant. Due to the continuity of velocities, the same happens for the fluid velocity $u \cdot n_w$ at the interface, and due to the incompressibility constraint the velocity has to change globally in the fluid domain $\Omega_f(t)$. The pressure can be seen as a Lagrange multiplier and more specifically as sensitivity of a (jumping) energy functional with respect to the incompressibility constraint, which explains the peak.

To substantiate this explanation numerically, we add an artificial penalty for the velocity on the sub-domain Ω_f^0 below the contact line

$$S_a(u, v) := \gamma_a^0 h^{-2} (u, v)_{\Omega_f^0}.$$

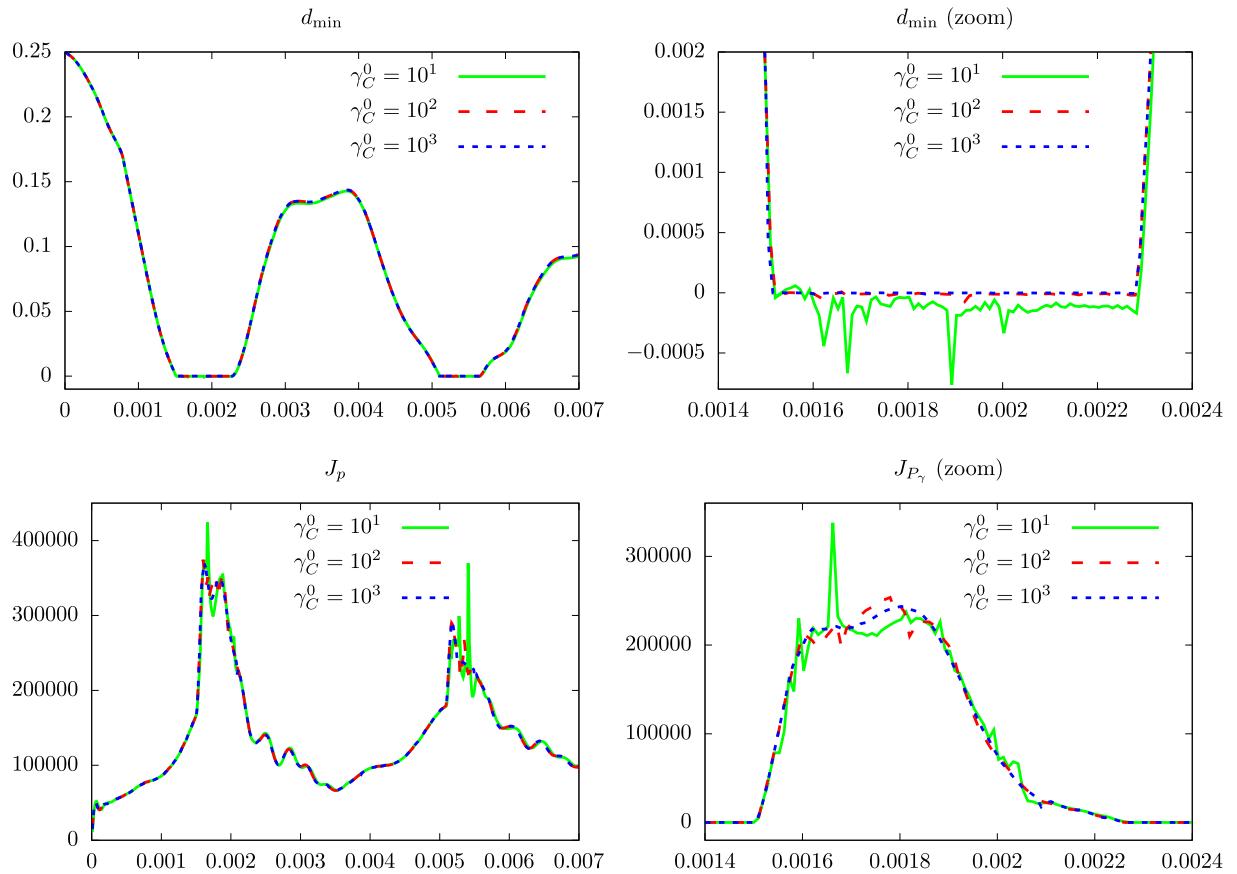


FIGURE 5. Top row: “Minimal distance” d_{\min} to Γ_w over time with two contact periods. Left: total time interval. Right: zoom-in at the first contact interval. Bottom left: pressure functional J_p over time. Bottom right: functional J_{P_γ} measuring the contact force around the contact interval over time.

For $\gamma_a := \gamma_a^0 h^{-2} \rightarrow \infty$, the fluid velocity is driven to zero below the obstacle. As this is already the case before contact, no abrupt changes in the fluid velocity are expected at the moment of the impact. Note that the problem with $\gamma_a^0 > 0$ is purely artificial, as the pressure mean values are still applied on the whole fluid boundary, including the boundary of Ω_f^0 .

In Figure 6 we compare the minimal distance and pressure functionals for computations without penalty ($\gamma_a^0 = 0$) to results for $\gamma_a^0 = 10$. First, we note that the vertical displacement is significantly influenced by the penalty, which has to be expected as the fluid dynamics are altered. Contact happens later at $t \approx 0.0022$ with the artificial penalty. Moreover, we observe indeed that the initial pressure peak at the time of impact is significantly reduced for $\gamma_a^0 = 10$.

6.3. Contact problem

Next, we study a problem, where it comes to real contact with the wall $\Gamma_w = \{(x, y) \in \Omega, y = 0.25\}$. At time $t = 0$, we define

$$\Omega_f(0) := (0, 1) \times (0.25, 0.5), \quad \Gamma(0) := (0, 1) \times 0.5, \quad \Omega_s(0) := (0, 1) \times (0.5, 0.6).$$

Below Γ_w , we define a fixed artificial fluid domain $\Omega_f^C := (0, 1) \times (0, 0.25)$.

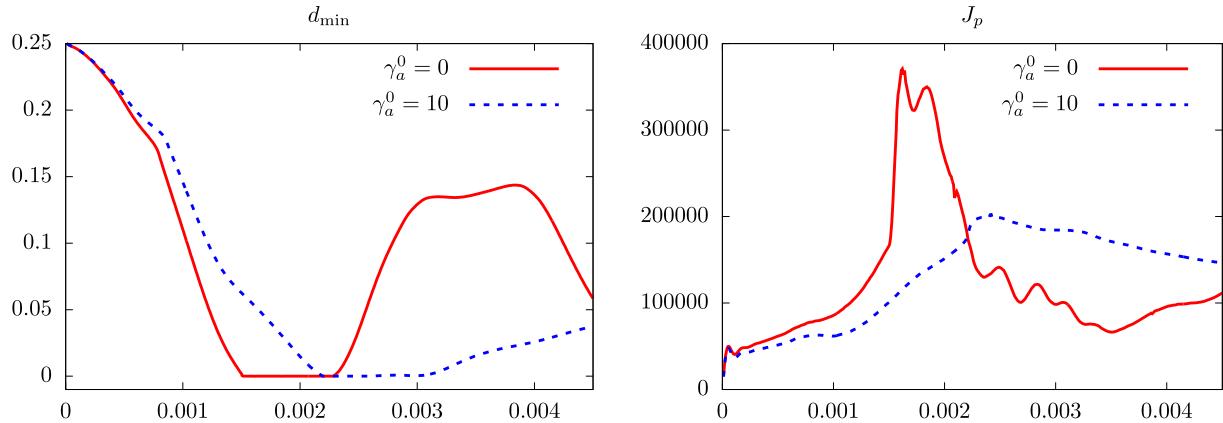


FIGURE 6. Minimal distance d_{\min} to the contact line Γ_w (left) and mean pressure over the interface J_p (right), plotted over time with an artificial penalty for the velocity below the virtual obstacle.

We apply again a pressure mean value \bar{P} on the lateral boundaries $\Gamma_{f,\text{left}}$ and $\Gamma_{f,\text{right}}$ of the fluid domain $\Omega_f(t)$. As the size of Γ_{left} and Γ_{right} is smaller and the viscous fluid forces acting against the closure of the channel are stronger than in the previous example, we have to set a larger pressure force $\bar{P} = 3 \times 10^5$ in order to obtain contact. On the other hand, the contact was never released again in our numerical experiments, when we used this constant boundary force for all times. Therefore, we decrease \bar{P} linearly from $t = 10^{-3}$ on until it reaches zero at $t = 1.2 \times 10^{-3}$. In order to avoid the issues related to no-slip conditions and contact, we use slip-interface conditions first, *i.e.* the Variational Formulation 2 with $\mathcal{A}_{\text{slip,FSI}}^C$. Unless stated differently, λ is chosen as the jump of numerical stresses $[\tilde{\sigma}_{n,\text{slip}}]$. Moreover, we use again a Cartesian mesh that consists of 5120 elements, a time step $\delta t = 10^{-5}$ and temporal pressure stabilisation with $\gamma_{pt} = 10^{-2}$. Unless explicitly stated, the Nitsche parameters are chosen as $\gamma_{\text{fsi}}^0 = \gamma_C^0 = 10^3$ and the penalty in the artificial fluid as $\gamma_a^0 = 10^2$.

The results on a coarser mesh are illustrated in Figure 7 at four time instants. Contact happens after the pressure on the lateral boundaries is released, as the solid continues moving downwards for some time. During contact, there is a very small overlap of the solid with the artificial fluid Ω_f^C . As the overlap is of order 10^{-5} , it can barely be seen in the bottom left picture. Notice however the triangular cells in Ω_f^C that are used only, when a patch is cut by $\Gamma(t)$.

6.3.1. Comparison of the two contact formulations

First, we compare the two contact strategies derived in Section 4, *i.e.* the relaxed contact formulation introduced in Section 4.1 with a small gap of size $\epsilon(h) = h/10$ between the solid and Γ_w and the strategy using an artificial fluid derived in Section 4.2. A comparison of the results for d_{\min} , the pressure norm $\|p\|_{L^2(\Omega_{f,\text{mid}}(t))}$, where

$$\Omega_{f,\text{mid}}(t) := \{x \in \Omega_f(t), 0.4 \leq x_1 \leq 0.6\}$$

denotes the central part of the fluid domain, and the contact force J_{P_γ} are shown in Figure 8 on two different meshes with 5120 and 20480 elements, respectively.

First, we observe from the plots in the top row that the interface stays at a distance to Γ_w of about $\epsilon \approx h/10 \approx 1.4 \times 10^{-3}$ on the coarser and $\epsilon \approx h/10 \approx 7 \times 10^{-4}$ on the finer mesh for the relaxed formulation. The much smaller overlap with Ω_f^C in the artificial fluid formulation is not visible, not even in the zoom-in on the right.

While the curves for d_{\min} look similar in the global picture (left), the zoom-in shows significant differences already before the impact. The contact happens earlier for the artificial fluid formulation: on the coarser mesh

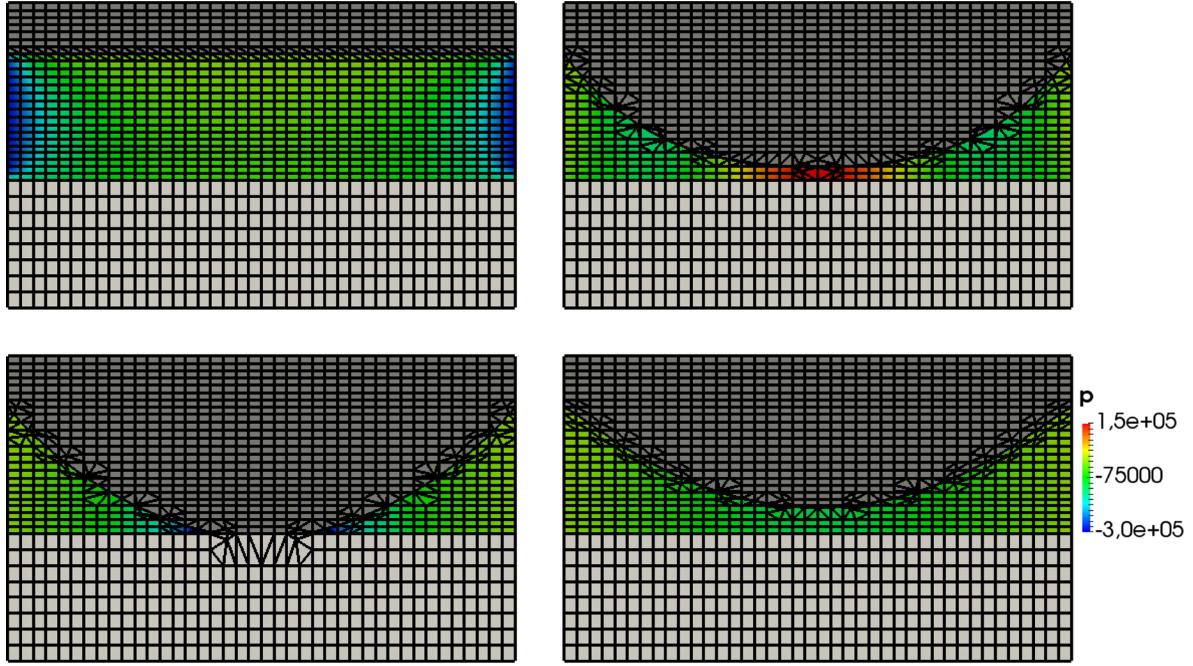


FIGURE 7. Illustration of the contact problem at four time instances on a coarse mesh: $t = 0$ (top left), $t = 1.2 \times 10^{-3}$ (top right), $t = 2 \times 10^{-3}$ (bottom left) and $t = 2.5 \times 10^{-3}$ (bottom right). The grey part corresponds to the structure $\Omega_s(t)$, the white part is the artificial fluid Ω_f^C . In the fluid domain $\Omega_f(t)$, values of the pressure p are visualised.

at time $t_{C,a} = 1.87 \times 10^{-3}$ compared to $t_{C,r} = 2.02 \times 10^{-3}$ for the relaxed formulation. This deviation is already much smaller on the finer mesh, where $t_{C,r} - t_{C,a} = 2 \times 10^{-5}$. The reason for this deviation is that in the artificial fluid formulation the wall Γ_w is only asymptotically for $\gamma_a \rightarrow \infty$ impermeable for the fluid. Therefore, in practice, the fluid forces acting against the contact, in particular the pressure p , are smaller for this formulation. As $\gamma_a = \gamma_a^0 h^{-2} \rightarrow \infty$ for $h \rightarrow 0$ the difference is significantly reduced on the finer mesh.

To substantiate this explanation, we plot the pressure norm $\|p\|_{L^2(\Omega_{f,mid}(t))}$ over the central part of the fluid domain $\Omega_{f,mid}(t)$ and the time period before the impact in the bottom left figure. The maximum value of the norm on the coarser mesh at time $t = 1.1 \times 10^{-3}$ is approximately 20.563 for the relaxed formulation and about 19.065 for the artificial fluid version. On the finer mesh, the pressure values are much closer. After that time the functional values decrease because the domain $\Omega_{f,mid}(t)$ gets smaller.

In the next paragraph, we will study the performance of both contact formulations under mesh refinement.

6.3.2. Convergence under mesh refinement

We solve the same problem on three different meshes with 1.280, 5.120 and 20.480 mesh elements, where the finer meshes are obtained from the coarsest one by global mesh refinement. The plots of the functionals d_{\min} and J_{P_γ} as well as the functionals

$$J_{\text{contact}} := \frac{\left\| \gamma_C^{1/2} [P_\gamma(\lambda, d)]_+ + \gamma_C^{-1/2} \lambda \right\|_{\Gamma(t)}}{\|\lambda\|_\Gamma}, \quad J_{\text{vel,fsi}} := \frac{\|(\dot{d} - u) \cdot n\|_{\Gamma_{\text{fsi}}(t)}}{\|u \cdot n\|_{\Gamma_{\text{fsi}}} + \|\dot{d} \cdot n\|_{\Gamma_{\text{fsi}}}}$$

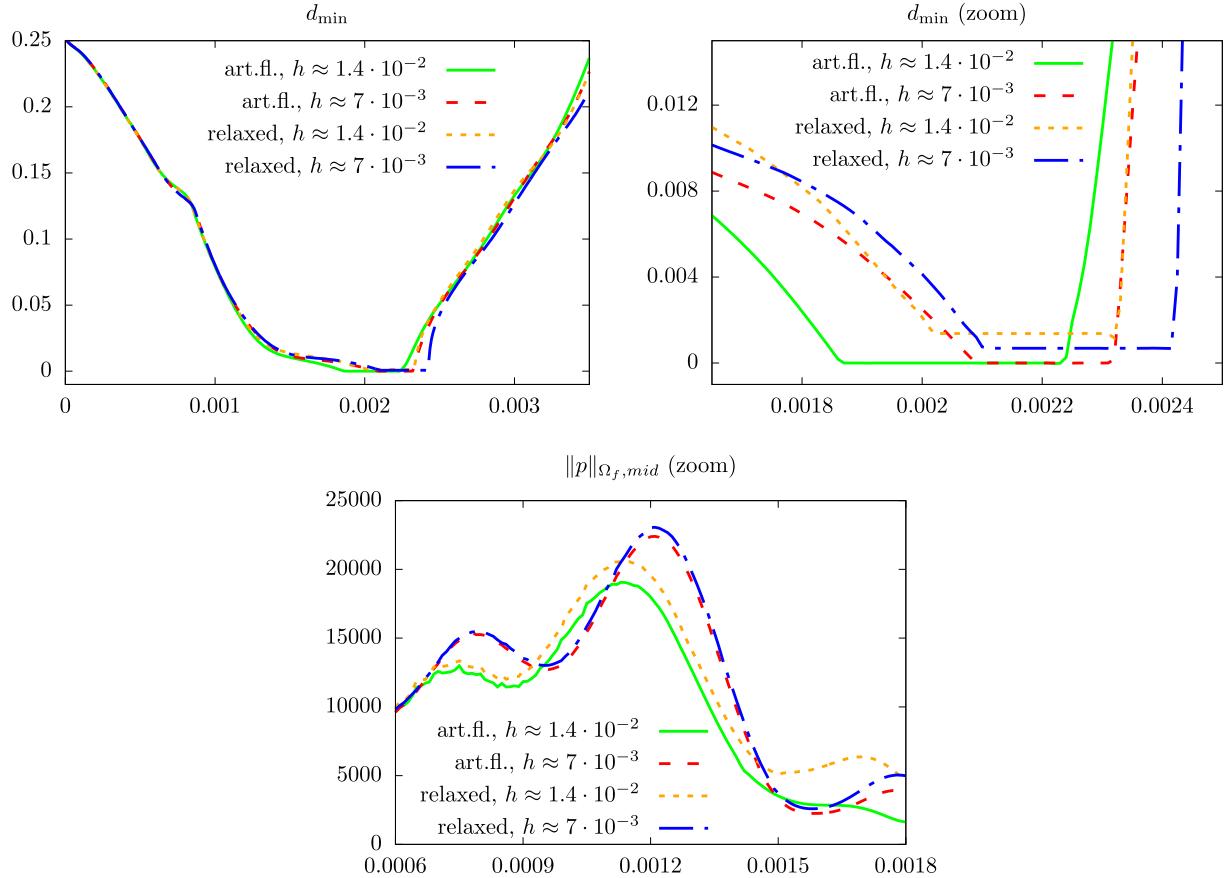


FIGURE 8. Comparison of the relaxed and the artificial fluid contact formulation. *Top*: minimal distance d_{\min} to Γ_w . *Right*: zoom-in around the contact interval. *Bottom*: pressure norm $\|p\|_{L^2(\Omega_f,mid)}$ over the central part of the fluid domain before contact over time.

measuring the fulfilment of the contact condition and the continuity of velocities on the part $\Gamma_{\text{fsi}}(t)$ of $\Gamma(t)$ that is not in contact with Γ_w

$$\Gamma_{\text{fsi}}(t) := \{x \in \Gamma(t) \mid P_\gamma(\lambda, d)(x) \leq 0\}$$

are shown in Figure 9 for the artificial fluid formulation and in Figure 10 for the relaxed contact formulation over time. The quantities $\|\cdot\|_*$ that are used to scale the functionals are temporal averages of the respective norms over the interval $I = [0, 0.004]$, computed on the finest grid.

First, we observe for both formulations in the plots on the top left that the contact happens later, the finer the discretisation is, as the fluid forces which act against the closure of the fluid channel are better resolved on the finer meshes (see also Fig. 8 and the related discussion above).

The curves for the contact force P_γ on the top right of both figures show significant differences between the two formulations. While the functional values seem to converge for the artificial fluid formulation (if we neglect the time shift), the contact force gets larger under mesh refinement for the relaxed formulation. The larger values for the relaxed formulation are due to the presence of the fluid forces $\sigma_{f,n}$ during the whole contact interval, that are not penalised in this formulation. As a Lagrange multiplier for the incompressibility constraint, the continuous pressure p gets singular when it comes to contact. The discrete pressure p_h gets larger and larger under mesh refinement in our computations.

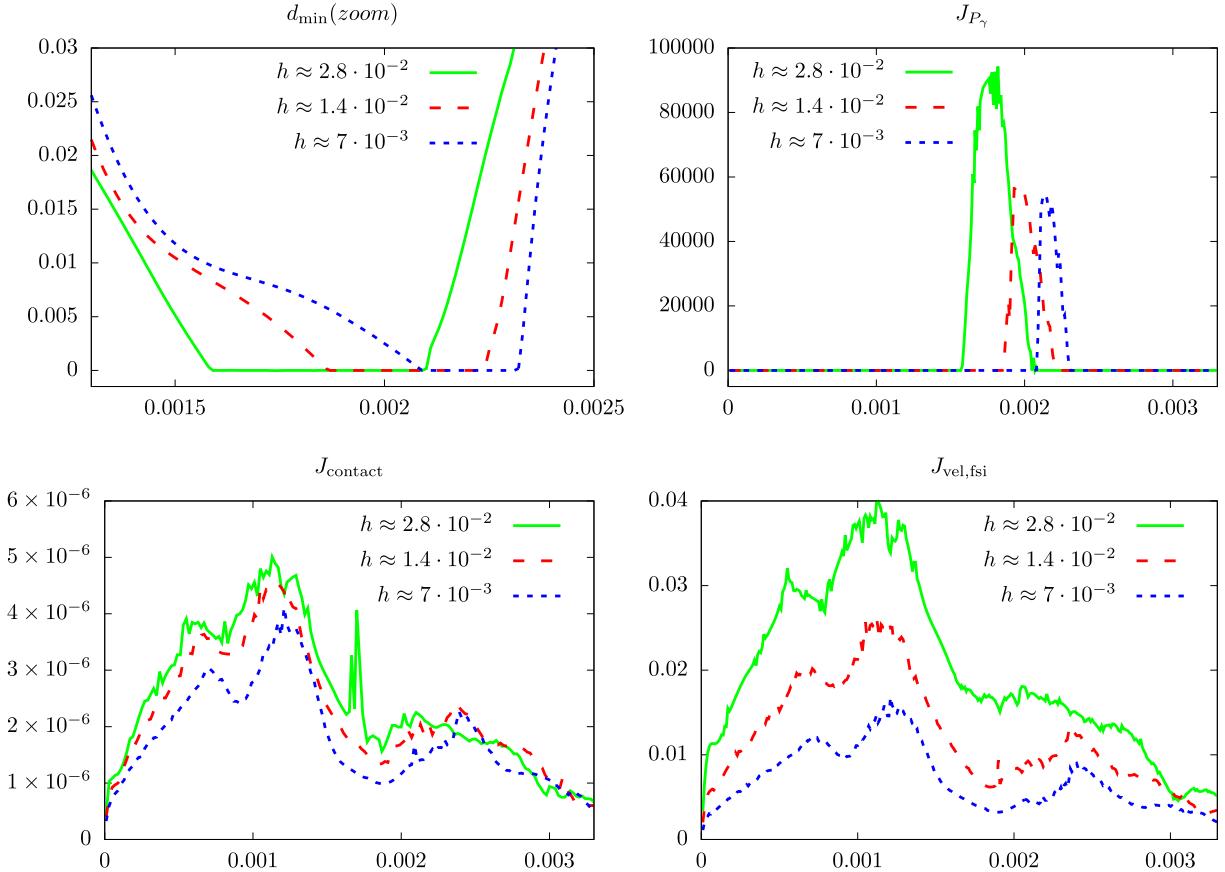


FIGURE 9. Convergence studies under mesh refinement for the *artificial fluid formulation* by means of the following functionals over time: *Top left*: minimal distance $d_{\min}(t)$ to Γ_w , *top right*: contact force J_{P_γ} , *Bottom left*: fulfilment of the contact condition J_{contact} . *Bottom right*: continuity of velocities $J_{\text{vel,fsi}}$.

On the other hand, the functional J_{contact} , that measures the difference between $-\gamma_C^{1/2}[P_\gamma]_+$ and $\gamma_C^{-1/2}\lambda$ decreases under mesh refinement for both formulations. Besides the differences in the contact force J_{P_γ} , the functional values on each of the mesh levels are actually very similar. The reason must be that the fluid forces $\sigma_{f,n}$ enter in both λ and $[P_\gamma]_+$. We conclude that the increase in the functional J_{P_γ} seems not to be an issue for the contact dynamics.

Both J_{contact} and the functional $J_{\text{vel,fsi}}$ are controlled by the stability estimate in Theorem 5.1 for $\theta = 1$. Although the parameter $\theta = 0$ is used here, we observe that both functionals decrease with mesh refinement before and during contact. While the convergence for the contact functional is quite slow, the values of the velocity functional indicate a convergence order $\mathcal{O}(h^\alpha)$ with $0.5 \leq \alpha \leq 1$ for both formulations. Note that in contrast to the term J_{contact} , $J_{\text{vel,fsi}}$ is controlled in the stability estimate in Theorem 5.1 even with the pre-factor $(\gamma_{\text{fsi}}^0 \mu_f)^{1/2} h^{-1/2}$.

6.3.3. Flux formulations

Next, we compare the different choices for λ . We show results exemplarily for the artificial fluid formulation with slip interface conditions. We will compare results using the jump of stresses $\lambda = [\![\sigma_{n,\text{slip}}]\!]$ (4.7), the jump

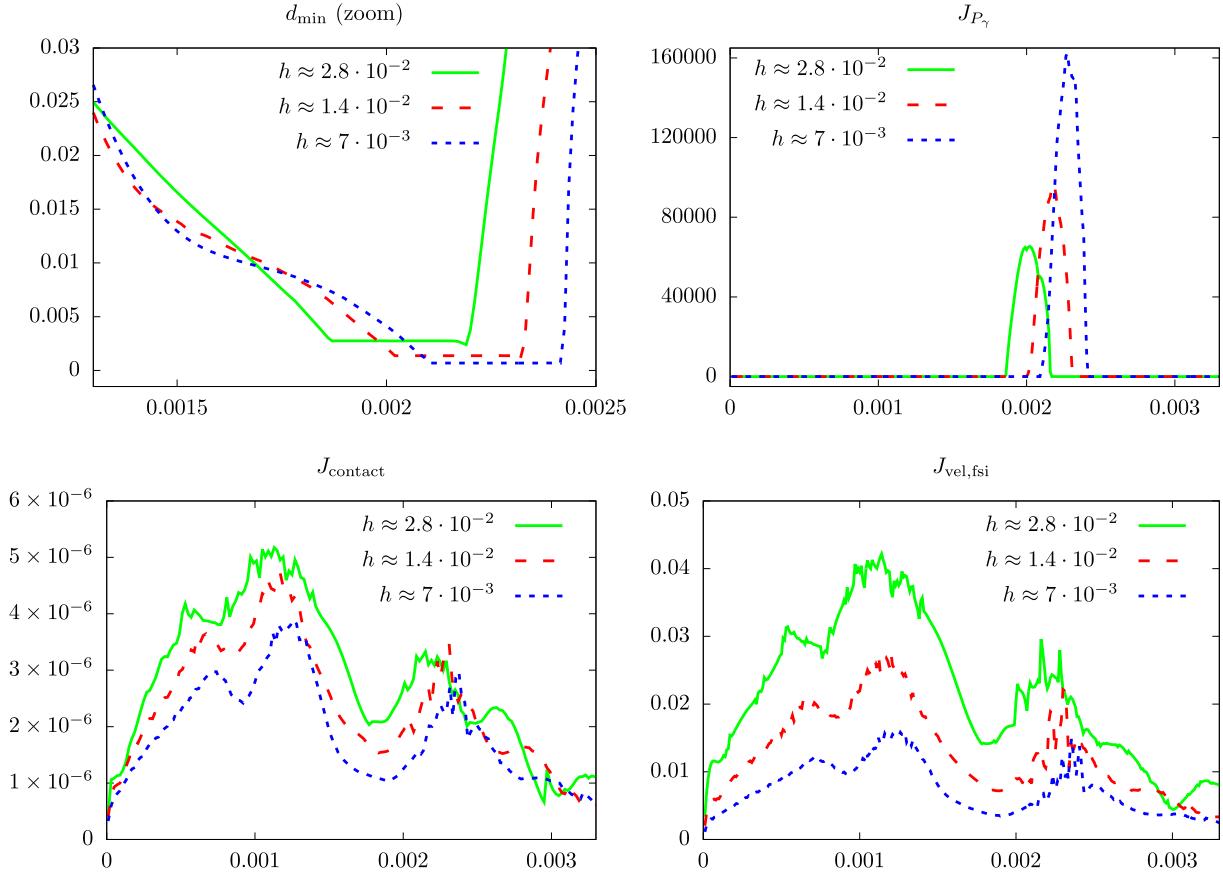


FIGURE 10. Convergence studies under mesh refinement for the *relaxed contact formulation* by means of the following functionals over time: *Top left*: minimal distance d_{\min} of $\Gamma(t)$ to Γ_w , *top right*: contact force J_{P_γ} , *Bottom left*: fulfilment of the contact condition J_{contact} . *Bottom right*: continuity of velocities $J_{\text{vel,fsi}}$.

of the numerical fluxes $\lambda = [\![\tilde{\sigma}_{n,\text{slip}}]\!]$ (4.8) and the extended fluxes $\lambda = [\![\tilde{\sigma}_{n,\text{slip}}]\!] + \tau^T \sigma_s n (\tau \cdot n_w)$ (see (4.10)). As the results for the latter two choices are nearly identical in this example, we show here only plots for the jump of stresses and the (non-extended) numerical fluxes. We use the artificial fluid formulation (Variational Formulation 2) and the previously used mesh with 5120 elements.

In Figure 11, we show the minimal distance d_{\min} to Γ_w , the contact force J_{P_γ} and the integral over the velocity difference across the contact part $\Gamma_C(t)$ of the interface over time

$$J_{\text{vel},C} := \int_{\Gamma_C(t)} (\dot{d} - u) \cdot n \, ds, \quad \Gamma_C(t) := \{x \in \Gamma(t) \mid P_\gamma(\lambda, d)(x) \leq 0\}.$$

The fluid velocity u is here artificial as it comes from Ω_f^C . When choosing $\lambda = [\![\tilde{\sigma}_{n,\text{slip}}]\!]$, we ensure that there is no feedback from this artificial velocity to the solid, see (4.5). For the jump of stresses $\lambda = [\![\sigma_{n,\text{slip}}]\!]$, we obtain a mixture of the solid contact condition and the continuity of normal velocities and a feedback might result. This follows analogously to the no-slip case, see (3.14).

In the left sketch of Figure 11, we see that the minimal distance in the stress-based formulation shows oscillations during the whole contact interval, especially in the second half. The interface jumps back and forth

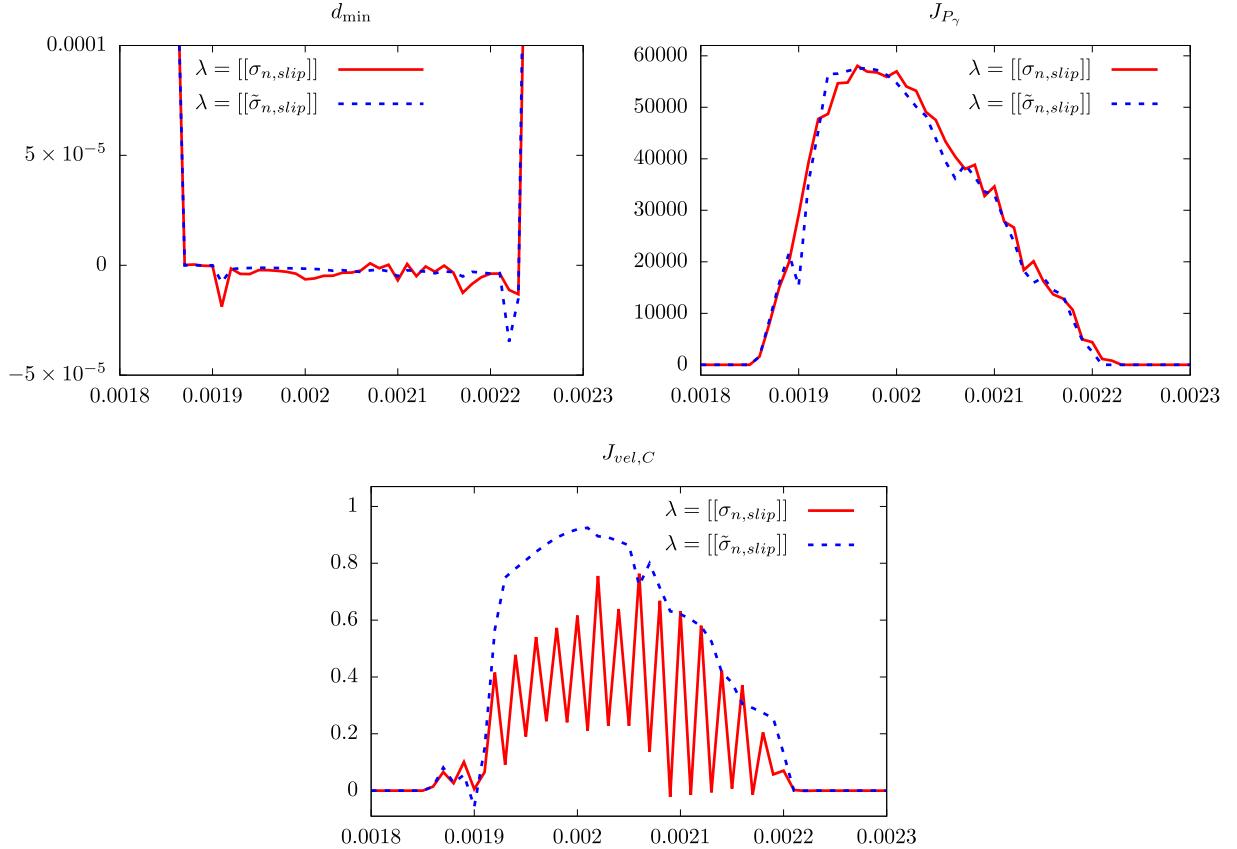


FIGURE 11. Comparison of the different possibilities to choose the fluxes λ . Minimal distance d_{min} to Γ_w (top left), contact force J_{P_γ} (top right) and velocity difference $J_{vel,C}$ (bottom) integrated over the contact part $\Gamma_C(t)$ of $\Gamma(t)$ over time for computations with $\lambda = [[\sigma_{n,slip}]]$ and $\lambda = [[\tilde{\sigma}_{n,slip}]]$ for the artificial fluid formulation.

over the contact line many times. The curve corresponding to the formulation using discrete fluxes is much smoother. Similarly, the contact force J_{P_γ} looks smoother, when the flux formulation is used. The reason for this behaviour is the mixture of the interface conditions during contact. On the bottom of Figure 11, we see that the velocity difference $J_{vel,C}$ shows wild oscillations for the stress formulation, while it looks much smoother when using $[[\tilde{\sigma}_{n,slip}]]$. As the artificial velocity in Ω_f^C has no physical meaning, it is not a drawback that the absolute values of $J_{vel,C}$ are larger. Due to the feedback of this velocity to the contact conditions, the oscillations appear in the displacement as well.

On the other hand, we should mention that the oscillations are relatively small. Especially those in d_{min} are almost by a factor 10^3 smaller than the mesh size $h \approx 1.4 \times 10^{-2}$ in vertical direction in this example and are therefore still acceptable.

6.3.4. Influence of the contact parameter γ_C^0

Next, we study the effect of different contact parameters γ_C^0 for the artificial fluid formulation and $\lambda = [[\tilde{\sigma}_{n,slip}]]$ on the mesh with 5120 elements. In Figure 12, we show the “minimal distance” d_{min} (top) and the contact force J_{P_γ} over time for different contact parameters γ_C^0 . The results are similar to the corresponding results for the virtual obstacle problem in Figure 5. For the smallest contact parameter $\gamma_C^0 = 10$, the contact condition is

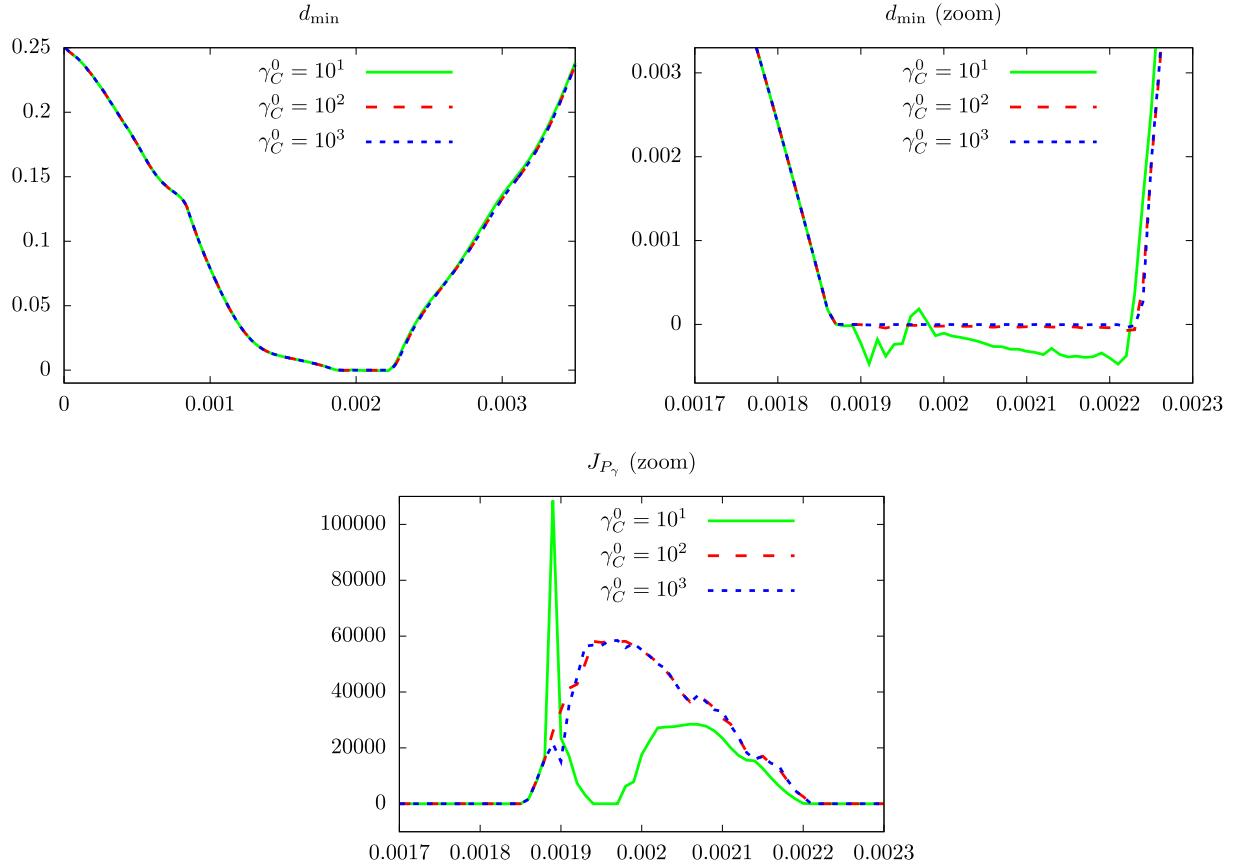


FIGURE 12. Parameter studies for the contact parameter γ_C^0 : Minimal distance of $\Gamma(t)$ to the wall Γ_w (*Top left*: total time interval, *top right*: zoom around the contact interval) and contact force J_{P_γ} (*bottom*) over time.

violated throughout the contact interval ($d_{\min} < 0$). The maximum overlap into the artificial fluid domain is again approximately by a factor 30 smaller than the mesh size $h \approx 1.4 \times 10^{-2}$. This violation gets smaller, the larger the contact parameter is chosen. The instabilities for the smallest parameter are still much better visible in the contact force J_{P_γ} . At time $t = 1.89 \times 10^{-3}$ the functional shows a huge peak, as the contact condition $d \cdot n_w \leq g_0$ is severely violated and it vanishes from $t = 1.94 \times 10^{-3}$ to $t = 1.97 \times 10^{-3}$, when the contact is in fact shortly released.

For the larger values $\gamma_C^0 \geq 10^2$, the curves are relatively smooth and very similar. Altogether, this shows again that the assumption “ γ_C^0 sufficiently large” in Theorem 5.1 is necessary in order to ensure stability.

6.3.5. Slip vs no-slip conditions

Next, we compare the effect of slip- and no-slip boundary and interface conditions in Figure 13. Due to the difficulties associated with the artificial fluid formulation and no-slip interface and boundary conditions (see the discussion at the end of Sect. 4.2), we use the relaxed contact formulation on the mesh with 5120 elements in this paragraph.

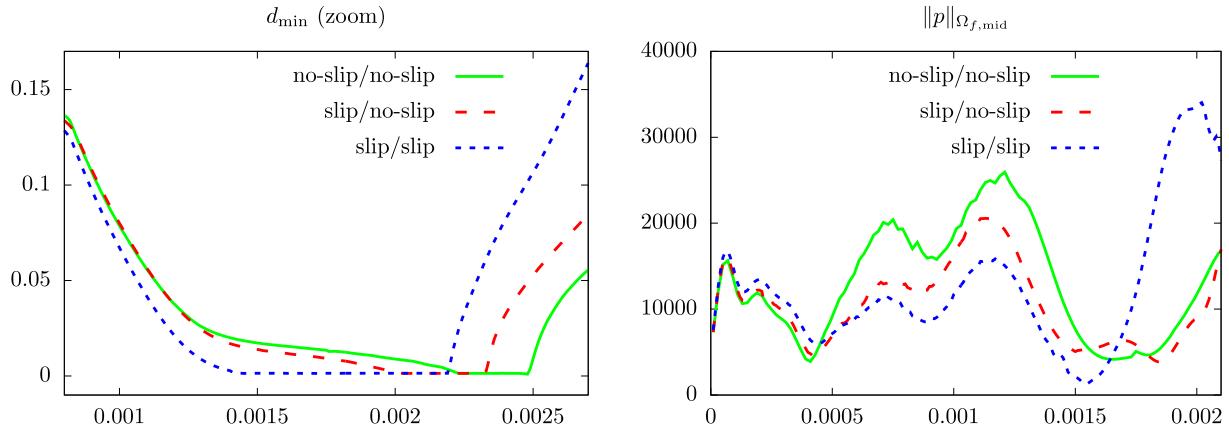


FIGURE 13. Comparison of slip- and no-slip interface/boundary conditions by means of the minimal distance d_{\min} of the interface $\Gamma(t)$ to Γ_w around the contact interval (*left*) and the L^2 -norm of the pressure over a region $\Omega_{f,\text{mid}}(t)$ around the contact surface before contact (*right*) over time. Due to the larger pressure before contact, the impact happens later when using no-slip conditions.

We show results for

- Slip conditions on the interface $\Gamma(t)$ and the lower wall Γ_w
- A slip condition on $\Gamma(t)$ and a no-slip condition on Γ_w
- No-slip conditions on $\Gamma(t)$ and Γ_w .

Note that the second option is possible, as for the relaxed contact formulation $\Gamma_w \cap \Gamma(t) = \emptyset$.

We observe that the contact condition (or more precisely the relaxed condition $d \cdot n_w \leq g_\epsilon$) is earlier active, when using slip-conditions: at $t_C = 1.42 \times 10^{-3}$ for slip/slip conditions compared to $t_C = 2.02 \times 10^{-3}$ for slip interface and no-slip boundary conditions and at $t_C = 2.23 \times 10^{-3}$ for no-slip conditions on interface and boundary. The reason is that the fluid forces, and in particular the pressure, that act against the contact are larger for no-slip conditions, as the fluid can not ‘slip’ out of the contact zone easily. This can be seen in the pressure plot on the right. The pressure is considerably larger from $t \approx 5 \times 10^{-4}$ for the no-slip conditions until contact is reached for the slip/slip case at $t_C = 1.42 \times 10^{-3}$.

As we are allowing for a small gap between the solid and the ground, these results do not contradict the theoretical results by Gerard-Varet *et al.* [36] discussed in Section 4.3, who showed that (in their configuration with a rigid body) contact can not happen, when no-slip conditions are used on the interface and/or the boundary. As discussed in Section 4.1, the basic assumption of the relaxed formulation is that a small or infinitesimal fluid layer remains during contact. On the contrary, the results confirm that contact is more likely to happen for slip-conditions, which is in agreement with the theoretical results.

6.3.6. Comparison with an explicit ad hoc approach

The probably simplest possibility to combine the FSI model introduced in Section 2.1 and the contact approach described in Section 2.2 is to split Γ explicitly in each time-step into a fluid-structure interface $\Gamma_{\text{fsi}}(t_{m-1})$ and a contact surface $\Gamma_C(t_{m-1})$ based on the displacement $d(t_{m-1})$ of the previous time-step and to use the interface condition (3.1) on $\Gamma_{\text{fsi}}(t_{m-1})$ and the contact condition (2.10) on $\Gamma_C(t_{m-1})$. A strategy of this type has been used by Hecht and Pironneau [41]. The system of equations reads in the slip case: *Find*

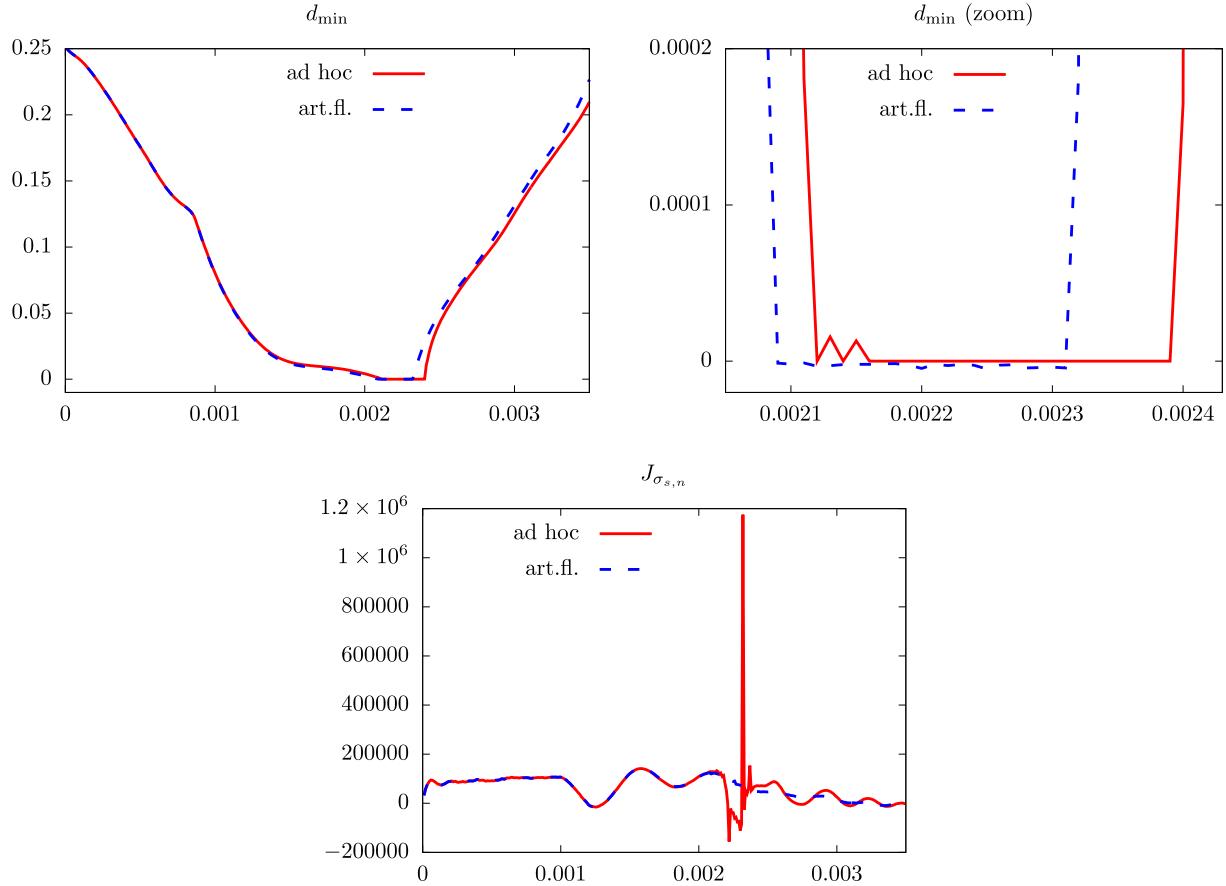


FIGURE 14. Comparison of an explicit *ad hoc* approach to include the contact and interface conditions with the approach using an artificial fluid described in Section 4.2. We show the minimal distance d_{\min} of the interface $\Gamma(t)$ to Γ_w on the top left and a zoom-in on the top right. Chattering, *i.e.* an alteration between contact and no contact is visible for the *ad hoc* approach at the beginning of the contact interval. The integral over the normal solid stresses $J_{\sigma_{s,n}}$ shown in the graph on the bottom shows large instabilities for the *ad hoc* approach.

$u \in \mathcal{V}, p \in \mathcal{Q}, d \in \mathcal{W}$ such that $\dot{d} = \partial_t d$ and

$$\begin{aligned}
 & (\partial_t u, v)_{\Omega_f(t)} + (\sigma_f(u, p), \nabla v)_{\Omega_f(t)} + (\operatorname{div} u, q)_{\Omega_f(t)} + (\partial_t \dot{d}, w)_{\Omega_s(t)} + (\sigma_s(d), \nabla w)_{\Omega_s(t)} \\
 & - (n^T \sigma_f(u, p) n, (w - v) \cdot n)_{\Gamma_{\text{fsi}}(t_{m-1})} - ((\dot{d} - u) \cdot n, n^T \sigma_f(v, -q) n)_{\Gamma_{\text{fsi}}(t_{m-1})} \\
 & + \gamma_{\text{fsi}} ((\dot{d} - u) \cdot n, (w - v) \cdot n)_{\Gamma_{\text{fsi}}(t_{m-1})} + \gamma_C (P_\gamma(\sigma_{s,n}(d), d), w \cdot n_w)_{\Gamma_C(t_{m-1})} \\
 & = (f_f, v)_{\Omega_f(t)} + (f_s, w)_{\Omega_s(t)} \quad \forall v, q, w \in \mathcal{V} \times \mathcal{Q} \times \mathcal{W}.
 \end{aligned} \tag{6.2}$$

We use the same numerical parameters as for the contact formulations presented in this work.

To compare this approach with the artificial fluid formulation we show the minimal distance to the ground d_{\min} and the integral over the normal solid stresses over $\Gamma(t) = \Gamma_{\text{fsi}}(t) \cup \Gamma_C(t)$

$$J_{\sigma_{s,n}} = \int_{\Gamma(t)} \sigma_{s,n} \, ds$$

on the finer mesh with 20 480 elements in Figure 14. While the curves for d_{\min} over the total time interval shown on the top left look similar, a zoom-in on the right shows again that the presence of the artificial fluid leads to an earlier time of impact. Moreover, we observe chattering for the ad-hoc approach at the beginning of the contact interval, *i.e.* contact is released twice again before the solid stays in contact with Γ_w . The interface jumps back to the fluid domain, with a (relatively small) minimal distance of approximately 1.6×10^{-5} .

In fact, the functional d_{\min} is not a good indicator to investigate stability for the *ad hoc* approach, as it is zero, as soon as one point of the interface lies on Γ_w . Note that this is different for the approaches presented in this work, where the interface can go beyond Γ_w (or Γ_ϵ for the relaxed approach). In the actual computation, the interface oscillates considerably in each time-step and contact is released and renewed frequently in different points. The functional $J_{\sigma_{s,n}}$ on the bottom left of Figure 14 serves to get a better impression of the instabilities during contact. It oscillates throughout the contact interval including a huge peak at $t = 2.32 \times 10^{-3}$. Moreover, we see that the elastic dynamics after the contact are also significantly influenced by these instabilities. Compared to the artificial fluid approach the oscillations in the displacement are significantly larger after contact.

We have also tried to iterate for the splitting into $\Gamma_C(t)$ and $\Gamma_{\text{fsi}}(t)$ within each time-step of the ad-hoc approach, which can be seen as an active-set strategy. This did however not cure the problem, as cycling between different active sets is not prevented.

7. CONCLUSIONS

We have presented two consistent formulations for fluid-structure interactions with contact, both including a continuous switch between the FSI interface and the contact condition depending on the contact force P_γ . In contrast to certain penalty approaches, the contact force is physically motivated and included in a consistent way in the variational formulations. Our numerical results indicate that the two proposed formulations have better stability properties than the usual *ad hoc* approaches and no chattering was observed in our computations.

Moreover, we have derived analytically a stability result for a generalised formulation including a parameter $\theta \in [-1, 1]$. As in the pure solid case [20], this result implies stability for $\theta = 1$ and stability up to a term including the contact force for $\theta \neq 1$. In our computations, we have however not observed any stability issues for the choice $\theta = 0$ either.

The contact formulations were derived here for the simplified configuration of contact with a fixed and straight wall and using linear models for the fluid and solid sub-problems. The algorithms can be applied to more complex contact configurations by using approaches from the literature to compute the projection and the distances between different surfaces [59, 65, 66]. In particular, the extension to the incompressible Navier–Stokes equations in the fluid and to non-linear elasticity in the solid can be addressed by combining the proposed approach with the arguments recently reported in Mlika *et al.* [55]. Moreover, Coulumb or Tresca friction can also be incorporated by following Chouly *et al.* [19, 24].

In our numerical examples, we have studied a two-dimensional model problem on a relatively simple and smooth geometry. In future, we plan to apply the methods on more complex geometries, including a direct comparison with experiments, in order to further validate the numerical approach.

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