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ANALYSIS AND OPTIMIZATION OF ROBIN-ROBIN PARTITIONED PROCEDURES IN FLUID-STRUCTURE INTERACTION PROBLEMS *

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Abstract. In the solution of Fluid-Structure Interaction problems, partitioned procedures are modular algorithms that involve separate fluid and structure solvers, that interact, in an iterative framework, through the exchange of suitable transmission conditions at the FS interface. In this work we study, using Fourier analysis, the convergence of partitioned algorithms based on Robin transmission conditions. We derive, for different models of the fluid and the structure, a frequency dependent reduction factor at each iteration of the partitioned algorithm, which is minimized by choosing optimal values of the coefficients in the Robin transmission conditions. Two-dimensional numerical results are also reported, which highlight the effectiveness of the optimization procedure.

Key words. Fluid-structure interaction, Robin conditions, Fourier analysis, optimized Schwartz methods

AMS subject classifications. 65B99, 65M60,

1. Introduction. In the framework of Domain-Decomposition methods for the solution of differential problems, one of the most popular non-overlapping method is the one based on successive exchanges of interface Robin data and called *Lions' method* (see [18] and, e.g., [7, 16]). Recently, this strategy has been applied to the Fluid-Structure Interaction (FSI) problem, which describes any physical phenomenon where a fluid and a structure interact by exchanging normal stresses through an interface (see [3]).

The solution of the FSI problem is problematic, since the fluid and the structure subproblems are coupled through the *geometry problem* (that is the determination of the unknown interface position) and through the *interface continuity conditions*, namely the continuity of the velocity and of the normal stresses at the FS interface.

Whatever treatment for the geometry problem is considered (e.g. fixed point, Newton, explicit extrapolation in time), a sequence of linearized FSI problems, implicitly coupled through the interface conditions, has to be solved (see Sect. 2.1). Implicit coupling of the interface conditions is required in many applications featuring a large added mass effect of the fluid on the structure (see [6]) and it allows to achieve at the numerical level perfect energy balance between fluid and structure.

One possible strategy for the solution of such problems is to consider *modular algorithms* (also referred to as *partitioned procedures*), that involve separate fluid and structure solvers in an iterative framework. In particular, they interact through the exchange of suitable transmission conditions at the FS interface, and guarantee, at convergence, the satisfaction of the continuity conditions. Modular algorithms can be reinterpreted, in the framework of the Domain-Decomposition method, as preconditioned Richardson iterations over a suitable interface equation (see [2, 8]).

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This allows one to introduce more performing Krylov methods for the solution of the FSI problem. In particular we mention the *Dirichlet-Neumann/GMRES* and the *Robin-Neumann/GMRES* schemes, which lead to different modular algorithms (see [2, 4]).

The introduction of *Robin-Robin* (RR) partitioned procedures in the framework of FSI problems, as generalization of the classic *Dirichlet-Neumann* (DN) scheme, has been motivated to overcome the limitations of the latter algorithm. In particular, the performances of the DN scheme when the *added mass effect* is high (that is when the fluid and structure densities are similar) are very poor, and a (sometimes big) relaxation is needed to reach convergence (see [6, 12, 23]). On the contrary, RR schemes highlighted better convergence properties in the presence of a high added-mass. In particular, in [2, 3] the *Robin-Neumann* (RN) scheme has been shown to converge without relaxation, in the test cases studied, and to feature a big saving in computational time with respect to DN scheme.

This behaviour has been mostly evidenced by numerical tests and only few theoretical results of the convergence properties of partitioned procedures for the FSI problems are available so far. In particular, at the best of the authors' knowledge, the only convergence analysis have been proposed in [6] for the DN scheme and in [3] for the RN scheme. We mention also the analysis in [4] and in [2] for DN-GMRES and RN-GMRES schemes. In all the cases, the analysis has been performed on a simplified problem where the fluid is described by a 2D potential flow and the structure by a 1D reduced model (the *independent rings* in [6] and the *generalized string* in the other works).

The first goal of this work is to extend the convergence analysis of RR schemes (and then of DN) to more general classes of subproblems. In particular, we consider the generalized Stokes equations to describe the discretized-in-time fluid problem and a 2D linear elastic incompressible structure in the half plane. In particular, the proposed analysis are based on the application of the Fourier transform (see, e.g. [1, 13, 14]) and on the determination of a reduction factor.

Secondly, we focus on parameters in Robin transmission conditions. Obviously, the convergence velocity of RR schemes heavily depends on the choice of these parameters. A proposal for these parameters has been given in [3], based on heuristic considerations. This choice has revealed to be satisfactory in order to overcome the limitations of the DN scheme, but a rigorous optimization procedure is still missing.

The second goal of this work is to optimize the values of parameters in the Robin interface conditions, starting from the new convergence analysis proposed in this work and following the literature on Optimization Schwartz Methods (see [1, 13, 14] for instance).

The outline of the paper is as follows. In Sect. 2 we introduce the FSI problem under investigation, we describe the different treatments of the geometry problem and we introduce the RR partitioned procedure. In Sect. 3 we perform the new convergence analysis, extending the results obtained in [3, 6] to more general classes of FSI problems. In Sect. 4 we propose an optimization procedure for the values of the parameters in Robin transmission conditions, starting from the convergence analysis. Finally, in Sect. 5 we present 2D numerical results, which confirm the effectiveness of the proposed optimal values with respect to the heuristic ones.

2. Problem setting. Let us consider a computational domain $\Omega^t \subset \mathbb{R}^d$ ($d=2, 3$, being the space dimension). This domain is divided into a sub-domain Ω_s^t occupied by an elastic structure and its complement Ω_f^t occupied by the fluid. The fluid-structure

interface Σ^t is the common boundary between Ω_s^t and Ω_f^t (see Fig. 2.1), while with Γ_i^t and $\Gamma_{i,s}^t$ we denote the fluid and structure artificial sections. Furthermore, \mathbf{n} is the outward normal on $\partial\Omega_f^t$. The initial configuration Ω^0 at $t = 0$ is considered as the reference one.

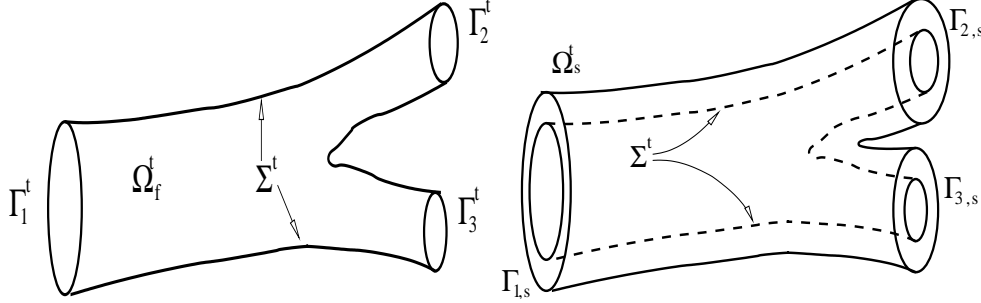


FIG. 2.1. Example of computational fluid domain Ω_f^t (left) and solid domain Ω_s^t (right).

We adopt a purely Lagrangian approach to describe the structure kinematics. We denote the reference (initial) configuration by $\Omega_s := \Omega_s^0$. Hereafter, $\bar{\eta}$ denotes the displacement of the solid medium with respect to Ω_s . For any function \bar{g} defined in the reference solid configuration, we denote by g its counterpart in the current configuration. The solid is assumed to be a linear elastic material, characterized by the Cauchy stress tensor

$$\mathbf{T}_s(\boldsymbol{\eta}) = \lambda_1(\nabla\boldsymbol{\eta} + (\nabla\boldsymbol{\eta})^T) + \lambda_2(\nabla \cdot \boldsymbol{\eta})\mathbf{I},$$

where

$$\lambda_1 = \frac{E}{1+\nu}, \quad \lambda_2 = \frac{E\nu}{(1+\nu)(1-2\nu)} + \lambda_1 \quad (2.1)$$

are the Lamé constants, E is the Young modulus, ν the Poisson modulus and \mathbf{I} is the identity tensor.

On the other hand, the fluid problem is stated in an *Arbitrary Lagrangian-Eulerian* (ALE) framework (see, e.g., [9, 17]). The ALE mapping is defined by an appropriate lifting of the structure displacement. A classical choice is to consider a harmonic extension operator in the reference domain. In order to write the fluid problem in ALE coordinates, we recall the definition of ALE time derivative of the velocity \mathbf{u} :

$$\frac{D^A \mathbf{u}}{Dt} = \frac{\partial \mathbf{u}}{\partial t} + \mathbf{w} \cdot \nabla \mathbf{u},$$

where $\partial \mathbf{u} / \partial t$ is the Eulerian derivative and \mathbf{w} is the velocity of the points of the fluid domain defined by the ALE map. Moreover, the fluid is assumed to be homogeneous, Newtonian and incompressible, with Cauchy stress tensor given by

$$\mathbf{T}_f(\mathbf{u}, p) = -p\mathbf{I} + \mu(\nabla \mathbf{u} + (\nabla \mathbf{u})^T),$$

where p is the pressure and μ the dynamic viscosity.

Then, the full FSI problem in strong form reads:

1. *Fluid-structure problem.* Find the fluid velocity \mathbf{u} , pressure p and the structure displacement $\bar{\boldsymbol{\eta}}$ such that

$$\left\{ \begin{array}{ll} \rho_f \frac{D^A \mathbf{u}}{Dt} + \rho_f ((\mathbf{u} - \mathbf{w}) \cdot \nabla) \mathbf{u} - \nabla \cdot \mathbf{T}_f = \mathbf{f}_f & \text{in } \Omega_f^t \times (0, T), \\ \nabla \cdot \mathbf{u} = 0 & \text{in } \Omega_f^t \times (0, T), \\ \rho_s \frac{\partial^2 \bar{\boldsymbol{\eta}}}{\partial t^2} - \bar{\nabla} \cdot \bar{\mathbf{T}}_s = \bar{\mathbf{f}}_s & \text{in } \Omega_s^0 \times (0, T), \\ \mathbf{u} = \frac{\partial \boldsymbol{\eta}}{\partial t} & \text{on } \Sigma^t \times (0, T), \\ \mathbf{T}_f \mathbf{n} - \mathbf{T}_s \mathbf{n} = \mathbf{0} & \text{on } \Sigma^t \times (0, T), \end{array} \right. \quad (2.2)$$

2. *Geometry problem.* Given the interface structure displacement $\boldsymbol{\eta}|_{\Sigma^t}$, find a map $\mathcal{A} : \Omega_f^0 \rightarrow \Omega_f^t$ e.g. through an harmonic extension Ext of the boundary displacement

$$\mathcal{A}^t(\mathbf{x}_0) = \mathbf{x}_0 + Ext(\bar{\boldsymbol{\eta}}|_{\Sigma^0}),$$

such that $\Omega_f^t = \mathcal{A}^t(\Omega_f^0)$.

Here, ρ_f and ρ_s are the fluid and structure density and \mathbf{f}_f and $\bar{\mathbf{f}}_s$ the forcing terms. System (2.2) has to be endowed with suitable Dirichlet or Neumann boundary conditions on $\partial\Omega_f^t \setminus \Sigma^t$ and $\partial\Omega_s^0 \setminus \Sigma^0$, and initial conditions in Ω^0 . Two transmission conditions are enforced at the interface: the *continuity of fluid and structure velocities* (2.2)₄ and the *continuity of stresses* (2.2)₅. The fluid and the structure are also coupled by the geometry problem, leading to a non-linear system of partial differential equations.

Remark 1. In the case of an incompressible structure, we have $\nu = 0.5$. In this case, the Cauchy stress tensor is given by

$$\mathbf{T}_s(\boldsymbol{\eta}) = \lambda(\nabla \boldsymbol{\eta} + (\nabla \boldsymbol{\eta})^T) - \chi \mathbf{I},$$

where χ is the structure pressure, and the structure equation (2.2)₃ becomes

$$\left\{ \begin{array}{ll} \rho_s \frac{\partial^2 \bar{\boldsymbol{\eta}}}{\partial t^2} - \nabla \cdot \bar{\mathbf{T}}_s = \bar{\mathbf{f}}_s & \text{in } \Omega_s^0 \times (0, T), \\ \nabla \cdot \bar{\boldsymbol{\eta}} = 0 & \text{in } \Omega_s^0 \times (0, T). \end{array} \right.$$

2.1. Time discretization and FS interface treatment. Let us introduce the backward Euler time discretization for the fluid and the first order BDF scheme for the structure. The investigation of this work is independent on the time discretization, so that other schemes can be considered as well.

The main source of non-linearity comes from the fact that the interface (and hence the fluid domain) is unknown (*geometrical non-linearity*). This can be treated numerically in several ways. We focus here either on *implicit* treatments, where the FSI problem is solved at each time step by Picard, Newton or quasi-Newton iterations (see, e.g. [11, 19]), or on *explicit* treatments, where the interface position is extrapolated from previous time steps. Similar considerations apply also to the non-linearity coming from the Navier-Stokes convective term.

Whatever strategy is adopted, a sequence of linearized FSI problems implicitly coupled through the interface conditions (2.2)_{4,5} has to be solved. In particular, let us denote by Ω_f^* , \mathbf{u}^* and \mathbf{w}^* the known fluid domain, fluid velocity and fluid domain velocity, where $*$ denotes either an extrapolated value (explicit algorithms) or

the value of the previous subiteration (implicit algorithm). Moreover, let g^m be the approximation of a function g at time $t^m := m\Delta t$, where Δt is the time discretization parameter. Then, by applying the backward Euler scheme also for the continuity condition (2.2)₄, we obtain the following linear system

Given Ω_f^* , \mathbf{u}^* , \mathbf{w}^* , \mathbf{u}^n , $\boldsymbol{\eta}^n$ and $\boldsymbol{\eta}^{n-1}$, find the fluid velocity \mathbf{u}^{n+1} , pressure p^{n+1} and the structure displacement $\bar{\boldsymbol{\eta}}^{n+1}$ such that

$$\left\{ \begin{array}{ll} \frac{\rho_f}{\Delta t} \mathbf{u}^{n+1} + \rho_f ((\mathbf{u}^* - \mathbf{w}^*) \cdot \nabla) \mathbf{u}^{n+1} - \nabla \cdot \mathbf{T}_f^{n+1} = \mathbf{f}_f^{n+1} + \frac{\rho_f}{\Delta t} \mathbf{u}^n & \text{in } \Omega_f^*, \\ \nabla \cdot \mathbf{u}^{n+1} = 0 & \text{in } \Omega_f^*, \\ \frac{\rho_s}{\Delta t^2} \bar{\boldsymbol{\eta}}^{n+1} - \nabla \cdot \bar{\mathbf{T}}_s^{n+1} = \bar{\mathbf{f}}_s^{n+1} + \frac{\rho_s}{\Delta t^2} (2\bar{\boldsymbol{\eta}}^n - \bar{\boldsymbol{\eta}}^{n-1}) & \text{in } \Omega_s^0, \\ \mathbf{u} = \frac{\boldsymbol{\eta}^{n+1} - \boldsymbol{\eta}^n}{\Delta t} & \text{on } \Sigma^*, \\ \mathbf{T}_f^{n+1} \mathbf{n} = \mathbf{T}_s^{n+1} \mathbf{n} & \text{on } \Sigma^*, \end{array} \right. \quad (2.3)$$

with suitable boundary conditions on the artificial sections.

For the sake of simplicity, here and in the sequel of the paper we drop the index $n+1$ referring to the time step.

2.2. Partitioned procedures. System (2.3) is still coupled through the interface conditions (2.3)₄ and (2.3)₅. In this work we consider schemes which guarantee strong enforcement of such conditions, thus achieving a perfect energy balance. In particular, we focus on the so called *modular algorithms* that involve separate fluid and structure solvers interacting through the exchange of suitable transmission conditions at the FS interface Σ^* . At convergence, they guarantee the continuity of the velocity and of the normal stress at Σ^* . In particular, an algorithm that uses an explicit treatment of the FS interface position and a strong coupling of the interface conditions is called *semi-implicit* (see [5, 10, 20]).

The most classical modular algorithm is the Dirichlet-Neumann (DN) scheme, which consists in solving iteratively the fluid problem with the structure velocity as Dirichlet boundary condition at the FS interface, and the structure problem with the fluid normal stress as Neumann boundary condition at Σ^* (see, e.g., [19, 21]). However, it has been shown in [6, 12, 23] that in the presence of a large added mass effect, this procedure needs a strong relaxation and features a very slow convergence.

A new class of iterative procedures based on Robin transmission conditions, which generalizes the DN approach, has been introduced in [3]. In particular, the Robin-Robin schemes are based on the following transmission conditions:

$$\begin{aligned} \alpha_f \mathbf{u} - \mathbf{T}_f \mathbf{n} &= \alpha_f \frac{\boldsymbol{\eta} - \boldsymbol{\eta}^n}{\Delta t} - \mathbf{T}_s \mathbf{n}, & \text{on } \Sigma^*, \\ \frac{\alpha_s}{\Delta t} \boldsymbol{\eta} + \mathbf{T}_s \mathbf{n} &= \frac{\alpha_s}{\Delta t} \boldsymbol{\eta}^n + \alpha_s \mathbf{u} + \mathbf{T}_f \mathbf{n}, & \text{on } \Sigma^*, \end{aligned} \quad (2.4)$$

obtained by a linear combination of (2.3)₄ and (2.3)₅, with coefficients α_f, α_s positive. We observe that with the choice $\alpha_f \rightarrow \infty$ and $\alpha_s = 0$ we recover the DN scheme.

In this work we consider a fixed point algorithm consisting of subsequent iterations of fluid subproblems with interface condition (2.4)₁ and structure subproblems with interface condition (2.4)₂:

Robin-Robin partitioned algorithm: Given Ω_f^* , \mathbf{u}^* , \mathbf{w}^* , \mathbf{u}^n , $\boldsymbol{\eta}^n$ and $\boldsymbol{\eta}^{n-1}$ and the solution at previous iteration $\boldsymbol{\eta}_{m-1}$, find at each iteration m the fluid velocity

\mathbf{u}_m , the fluid pressure p_m and the structure displacement $\bar{\boldsymbol{\eta}}_m$ until convergence, such that

Fluid problem:

$$\begin{aligned} \frac{\rho_f}{\Delta t} \mathbf{u}_m + \rho_f ((\mathbf{u}^* - \mathbf{w}^*) \cdot \nabla) \mathbf{u}_m - \nabla \cdot \mathbf{T}_{f,m} &= \mathbf{f}_f + \frac{\rho_f}{\Delta t} \mathbf{u}^n && \text{in } \Omega_f^*, \\ \nabla \cdot \mathbf{u}_m &= 0 && \text{in } \Omega_f^*, \\ \alpha_f \mathbf{u}_m - \mathbf{T}_{f,m} \mathbf{n} &= \alpha_f \frac{\boldsymbol{\eta}^{m-1} - \boldsymbol{\eta}^n}{\Delta t} - \mathbf{T}_{s,m-1} \mathbf{n}, && \text{on } \Sigma^*, \end{aligned}$$

Structure problem

$$\begin{aligned} \frac{\rho_s}{\Delta t^2} \bar{\boldsymbol{\eta}}_m - \nabla \cdot \bar{\mathbf{T}}_{s,m} &= \bar{\mathbf{f}}_s + \frac{\rho_s}{\Delta t^2} (2\bar{\boldsymbol{\eta}}^n - \bar{\boldsymbol{\eta}}^{n-1}) && \text{in } \Omega_s^0, \\ \frac{\alpha_s}{\Delta t} \bar{\boldsymbol{\eta}}_m + \mathbf{T}_{s,m} \mathbf{n} &= \frac{\alpha_s}{\Delta t} \boldsymbol{\eta}^n + \alpha_s \mathbf{u}_m + \mathbf{T}_{f,m} \mathbf{n}, && \text{on } \Sigma^*. \end{aligned}$$

This strategy can also be interpreted as a preconditioned Richardson algorithm applied to a suitable interface condition (see [2]).

3. Convergence analysis of simplified problems. In this Section we analyze the convergence of the RR scheme applied to three different reduced FSI problems, which are simple enough to be analyzed theoretically. On the other hand, they feature a behavior similar to the more complex system (2.3). Therefore, we expect that the theoretical results obtained from the analysis of these models give insightful information also for system (2.3). This analysis is performed by following the optimization strategy proposed in the Optimized Schwarz Methods literature (see [1, 13, 14] for instance). In these works, a convergence analysis of two and three dimensional coupled problems with Robin transmission conditions, based on the application of the Fourier transform, has been derived, for the Laplace, Helmholtz and Maxwell equations.

In order to apply this strategy to the FSI problem, we introduce suitable simplifying assumptions and reduced models. In particular, in all the convergence analysis, we consider a fixed FSI domain. The fluid domain is the half plane $\Omega_f = \{\mathbf{x} = (x, y) \in \mathbb{R}^2 : x < 0\}$, the FS interface the line $\Sigma = \{(x, y) \in \mathbb{R}^2 : x = 0\}$, the structure domain either coincides with the interface Σ (*generalized string model*) or is the thick region $\Omega_s = [0, H_s] \times \mathbb{R}$ in the case of linear incompressible elasticity. In all the cases considered, we allow the FS interface to move only in the normal direction.

We will base our convergence analysis on a Fourier transform in the direction tangential to the FS interface (corresponding to the y variable in the case at hand), which is defined, for $w(x, y) \in L^2(\mathbb{R}^2)$, as

$$\mathcal{F} : w(x, y) \mapsto \hat{w}(x, k) = \int_{\mathbb{R}} e^{-iky} w(x, y) dy,$$

where k is the frequency variable. We will then be able to quantify the error, in the frequency space, between the pressure at the m -th iteration, $\hat{p}_m(x, k)$, and the exact pressure $\hat{p}(x, k)$. This allows us to define, on the FS interface, the *reduction factor* at iteration m for each frequency as

$$\rho_m(k) := \frac{|\hat{p}_m(0, k) - \hat{p}(0, k)|}{|\hat{p}_{m-1}(0, k) - \hat{p}(0, k)|}.$$

The RR algorithm converges if, at each iteration m , we have $\rho_m(k) < 1$ for all the relevant frequencies of the problem, namely for $0 \leq k \leq k_{max}$, where k_{max} is the maximal frequency supported by the numerical grid, and is of order π/h (h being the mesh parameter).

3.1. Potential flow - Generalized string model (P/GS). The first simplified model is obtained by considering a potential flow described by the Darcy equations for the fluid and the *generalized string model* for the structure (see [22])

$$\begin{aligned}
\rho_f \delta_t \mathbf{u} + \nabla p &= \mathbf{0} && \text{in } \Omega_f, \\
\nabla \cdot \mathbf{u} &= 0 && \text{in } \Omega_f, \\
\mathbf{u} \cdot \mathbf{n} &= \delta_t \eta && \text{on } \Sigma, \\
p &= \rho_s H_s \delta_{tt} \eta + \beta H_s \eta - G H_s \partial_{yy} \eta, && \text{on } \Sigma,
\end{aligned} \tag{3.1}$$

where $G = K\lambda_1/2$, with K the Timoshenko correction factor, and we have set $\delta_t w := \frac{w - w^n}{\Delta t}$, $\delta_{tt} w := \frac{\delta_t w - \delta_t w^n}{\Delta t}$, where w could be a scalar or a vector function and as usual we have omitted the time index $n+1$. The reaction term β , arising in the derivation of the generalized string model in the 3D case, is introduced also in the 2D case to take into account for transversal membrane effects. Problem (3.1) has to be completed with initial conditions and with boundary conditions here reducing to the assumption of boundedness for $x \rightarrow -\infty$ and $|y| \rightarrow \infty$. We point out that the structure displacement is in the x direction, which is the normal direction to the FS interface. The structure equation (3.1)₄ enforces the continuity of the normal stress at the interface along the normal direction. We observe that the analysis of this symplified problem has been already performed in [3]. For completeness, we report here the analysis via Fourier transform, which is functional for the further developments.

By combining linearly (3.1)₃ and (3.1)₄ with coefficients $(\alpha_f, -1)$ and $(\alpha_s, 1)$, respectively, we obtain two Robin boundary conditions. Observe that in this problem the viscous terms have been neglected so the fluid Cauchy stress tensor reduces to the only pressure. By setting $u_x = \mathbf{u} \cdot \mathbf{n}$, with \mathbf{n} the normal unit vector to the FS interface, and using coefficients $(\alpha_f, -1)$, we obtain

$$\alpha_f u_x - p = \alpha_f \delta_t \eta - \rho_s H_s \delta_{tt} \eta - \beta H_s \eta + G H_s \partial_{yy} \eta,$$

that is

$$\alpha_f \Delta t \delta_t u_x - p = \alpha_f \delta_t \eta - \rho_s H_s \delta_{tt} \eta - \beta H_s \eta + G H_s \partial_{yy} \eta - \alpha_f u_x^n.$$

Then, the Robin transmission condition for the fluid problem can be rearranged as

$$\alpha_f \Delta t \delta_t u_x - p = \left(\frac{\alpha_f}{\Delta t} - \frac{\rho_s H_s}{\Delta t^2} - \beta H_s \right) \eta + G H_s \partial_{yy} \eta + F_1(u_x^n, \eta^n, \eta^{n-1}), \tag{3.2}$$

where $F_1(u_x^n, \eta^n, \eta^{n-1})$ accounts for terms at previous time steps.

By using coefficients $(\alpha_s, 1)$, we have

$$\rho_s H_s \delta_{tt} \eta + \beta H_s \eta - G H_s \partial_{yy} \eta + \alpha_s \delta_t \eta = p + \alpha_s u_x,$$

and the Robin condition for the structure problem can be rearranged as

$$\left(\frac{\rho_s H_s}{\Delta t^2} + \beta H_s + \frac{\alpha_s}{\Delta t} \right) \eta - G H_s \partial_{yy} \eta = \alpha_s \Delta t \delta_t u_x + p + F_2(u^n, \eta^n, \eta^{n-1}), \tag{3.3}$$

where again $F_2(u^n, \eta^n, \eta^{n-1})$ accounts for terms at previous time steps.

Let m be an iteration index. Then, the corresponding iterative Robin-Robin algorithm reads:

Given $\mathbf{u}^0, p^0, \eta^0$, solve for $m \geq 0$

1. Fluid problem

$$\rho_f \delta_t \mathbf{u}_{m+1} + \nabla p_{m+1} = \mathbf{0} \quad \text{in } \Omega_f,$$

$$\nabla \cdot \mathbf{u}_{m+1} = 0 \quad \text{in } \Omega_f,$$

$$\begin{aligned} \alpha_f \Delta t \delta_t u_{x,m+1} - p_{m+1} = & \left(\frac{\alpha_f}{\Delta t} - \frac{\rho_s H_s}{\Delta t^2} - \beta H_s \right) \eta_m + \\ & + G H_s \partial_{yy} \eta_m + F_1(u_x^n, \eta^n, \eta^{n-1}) \quad \text{on } \Sigma; \end{aligned} \quad (3.4)$$

2. Structure problem

$$\left(\frac{\rho_s H_s}{\Delta t^2} + \beta H_s + \frac{\alpha_s}{\Delta t} \right) \eta_{m+1} - G H_s \partial_{yy} \eta_{m+1} = \quad (3.5)$$

$$= \alpha_s \Delta t \delta_t u_{x,m+1} + p_{m+1} + F_2(u_x^n, \eta^n, \eta^{n-1}) \quad \text{on } \Sigma.$$

We have the following

Proposition 1. *The reduction factor of iterations (3.4)-(3.5) is given by*

$$\rho^m(k) = \rho^{P/GS}(k) = \left| \frac{\frac{\rho_s H_s}{\Delta t} + \beta H_s \Delta t + G H_s \Delta t k^2 - \alpha_f}{\frac{\rho_s H_s}{\Delta t} + \beta H_s \Delta t + G H_s \Delta t k^2 + \alpha_s} \cdot \frac{\rho_f - \alpha_s \Delta t k}{\rho_f + \alpha_f \Delta t k} \right|, \quad (3.6)$$

which is independent of the iteration m .

Moreover, for all Fourier modes $k \in [0, k_{max}]$, there exist $\alpha_f^{max} = \alpha_f^{max}(k)$ and $\alpha_s^{max} = \alpha_s^{max}(k)$ such that, given $\mathbf{u}^0, p^0, \eta^0$, iterations (3.4)-(3.5) converge if $\alpha_f \in [0, \alpha_f^{max})$ and $\alpha_s \in [0, \alpha_s^{max})$ and if at least one of the two values is different from 0.

Proof. Thanks to the relation

$$\partial_x p = -\rho_f \delta_t u_x, \quad \text{on } \Sigma, \quad (3.7)$$

obtained by restricting the first equation of the fluid problem (3.4)₁ on the FS interface, it is possible to rewrite the Robin interface conditions (3.4)₃ and (3.5) in terms of the sole pressure, obtaining, on Σ ,

$$\begin{aligned} \frac{\alpha_f \Delta t}{\rho_f} \partial_x p^{m+1} + p^{m+1} = & \left(\frac{\rho_s H_s}{\Delta t^2} + \beta H_s - \frac{\alpha_f}{\Delta t} \right) \eta^m - G H_s \partial_{yy} \eta^m - F_1(u_x^n, \eta^n, \eta^{n-1}), \\ \left(\frac{\rho_s H_s}{\Delta t^2} + \beta H_s + \frac{\alpha_s}{\Delta t} \right) \eta^m - G H_s \partial_{yy} \eta^m = & p^m - \frac{\alpha_s \Delta t}{\rho_f} \partial_x p^m + F_2(u_x^n, \eta^n, \eta^{n-1}). \end{aligned} \quad (3.8)$$

We point out that for the structure problem we wrote the m -th iteration instead of the $m+1$ -th.

Since the problems involved are linear, we analyze without loss of generality the convergence to the zero solution when the forcing terms vanish, namely for $F_1(u_x^n, \eta^n, \eta^{n-1}) = F_2(u_x^n, \eta^n, \eta^{n-1}) = 0$.

The divergence free condition on \mathbf{u} allows us to rewrite the fluid problem in the unknown pressure

$$\Delta p = 0 \quad \text{in } \Omega_f. \quad (3.9)$$

Applying the Fourier transform in the y direction, we obtain the following ordinary differential equation for the pressure

$$-\partial_{xx} \widehat{p}^{m+1} + k^2 \widehat{p}^{m+1} = 0 \quad \text{in } (-\infty, 0),$$

whose solution is given by $\widehat{p}^{m+1}(x, k) = A(k)e^{kx} + B(k)e^{-kx}$. The boundedness assumption on the solution entails $B(k) = 0$, thus

$$\widehat{p}(x, k) = A(k)e^{kx}, \quad (3.10)$$

for a suitable $A(k)$.

Now, the interface conditions (3.8) become

$$\begin{aligned} \widehat{p}^{m+1} + \frac{\alpha_f \Delta t}{\rho_f} \partial_x \widehat{p}^{m+1} &= \left(\frac{\rho_s H_s}{\Delta t^2} + \beta H_s + G H_s k^2 - \frac{\alpha_f}{\Delta t} \right) \widehat{\eta}^m, & \text{for } x = 0, \\ \left(\frac{\rho_s H_s}{\Delta t^2} + \beta H_s + G H_s k^2 + \frac{\alpha_s}{\Delta t} \right) \widehat{\eta}^m &= \widehat{p}^m - \frac{\alpha_s \Delta t}{\rho_f} \partial_x \widehat{p}^m, & \text{for } x = 0. \end{aligned}$$

By noticing that $\partial_x \widehat{p}^{m+1} = A^{m+1}(k) k e^{kx}$, we have

$$\begin{aligned} \left(\frac{\rho_f + \alpha_f \Delta t k}{\rho_f} \right) A^{m+1}(k) &= \left(\frac{\rho_s H_s}{\Delta t^2} + \beta H_s + G H_s k^2 - \frac{\alpha_f}{\Delta t} \right) \widehat{\eta}^m, & \text{for } x = 0, \\ \left(\frac{\rho_s H_s}{\Delta t^2} + \beta H_s + G H_s k^2 + \frac{\alpha_s}{\Delta t} \right) \widehat{\eta}^m &= \frac{\rho_f - \alpha_s \Delta t k}{\rho_f} A^m(k), & \text{for } x = 0. \end{aligned}$$

We thus have $|A^{m+1}(k)| = \rho(k)|A^m(k)|$, with reduction factor ρ given by (3.6).

In order to guarantee convergence, we have to find suitable values of α_f and α_s such that $\rho^{P/GS} < 1$. To do this, if we consider the two factors d_1 and d_2 such that $\rho^{P/GS} = |d_1 d_2|$ in (3.6), we obtain sufficient conditions by imposing that $|d_1| < 1$ and $|d_2| < 1$, separately. For the first term, $d_1 = \frac{\rho_s H_s + \beta H_s \Delta t + G H_s \Delta t k^2 - \alpha_f}{\rho_f + \alpha_f \Delta t k}$, it is sufficient to have $\alpha_f \in [0, \alpha_f^{max})$, with

$$\alpha_f^{max}(k) = 2 \left(\frac{\rho_s H_s}{\Delta t} + \beta H_s \Delta t + G H_s \Delta t k^2 \right). \quad (3.11)$$

For the second term, $d_2 = \frac{\rho_f - \alpha_s \Delta t k}{\rho_f + \alpha_f \Delta t k}$, it is sufficient to have $\alpha_s \in [0, \alpha_s^{max})$, where

$$\alpha_s^{max}(k) = \frac{2\rho_f}{\Delta t k}.$$

A less sharp upper bound, but independent of k , is given by

$$\alpha_s^{max} = \frac{2\rho_f}{\Delta t k_{max}}. \quad (3.12)$$

Remark 2. We observe that function $A(k)$ is in general complex. However, in the particular case where the initial condition for the x component of the fluid velocity is symmetric with respect to the x -axis, then its Fourier transform is real and therefore A is real for each time (see (3.7)) and u_x remains symmetric at each time.

Remark 3. For the DN scheme, that is for $\alpha_f \rightarrow \infty$ and $\alpha_s = 0$, we obtain

$$\rho^{P/GS}(k) = \left| \frac{\rho_f}{k(\rho_s H_s + \beta H_s \Delta t^2 + G H_s \Delta t^2 k^2)} \right|.$$

Taking $k = 0$, we have that no convergence of the DN scheme is achieved, even with relaxation ($\rho^{P/GS}(0) \rightarrow \infty$). This is due to the unboundedness of the fluid domain. If

we consider a bounded domain $\mathbb{R} \times [-L, 0]$, we have that $k \in (k_{min}, k_{max})$, with k_{min} of the order of $\pi/L > 0$ (see [6]). In this case, the reduction factor of the DN scheme is always bounded, but it highlights a high dependence of the convergence performances on the ratio ρ_f/ρ_s , as well on k_{min} . In particular, if ρ_f/ρ_s is not small enough, for k sufficiently small we have $\rho^{P/GS}(k) > 1$, so that a relaxation is needed to reach convergence. On the contrary, for a general RR scheme the sensitivity of $\rho^{P/GS}$ on the ratio ρ_f/ρ_s is minimal and, as stated by Proposition 1, there exist always suitable values of α_f and α_s such that for each k convergence is guaranteed without relaxation.

3.2. Stokes - Generalized string model (S/GS). Let us now consider the coupling between the unsteady Stokes problem for the fluid and the generalized string model for the structure. The fluid and structure domains are the same of the previous problem and again we consider only normal displacement of the FS interface.

In this case, the fluid stress tensor is given by $\mathbf{T}_f = p\mathbf{I} + \mu(\nabla\mathbf{u} + \nabla^T\mathbf{u})$, since the viscous term is now considered and the normal component of the normal stress at the interface Σ is $\mathbf{n}^T \cdot \mathbf{T}_f \mathbf{n} = -p + \mu\partial_x u_x$. Therefore, the Robin-Robin sequential algorithm for the generalized Stokes-generalized string model coupling is given by

1. Fluid problem

$$\begin{aligned} \rho_f \delta_t \mathbf{u}^{m+1} - \mu \Delta \mathbf{u}^{m+1} + \nabla p^{m+1} &= \mathbf{0} && \text{in } \Omega_f \\ \nabla \cdot \mathbf{u}^{m+1} &= 0 && \text{in } \Omega_f \\ \alpha_f u_x^{m+1} - p^{m+1} + \mu \partial_x u_x^{m+1} &= \\ &= \left(\frac{\alpha_f}{\Delta t} - \frac{\rho_s H_s}{\Delta t^2} - \beta H_s \right) \eta^m + GH_s \partial_{yy} \eta^m + F_1(\eta^n, \eta^{n-1}) && \text{on } \Sigma \end{aligned} \quad (3.13)$$

2. Structure problem

$$\begin{aligned} \left(\frac{\rho_s H_s}{\Delta t^2} + \beta H_s + \frac{\alpha_s}{\Delta t} \right) \eta^{m+1} - GH_s \partial_{yy} \eta^m &= \\ &= \alpha_s u_x^{m+1} + p^{m+1} - \mu \partial_x u_x^{m+1} + F_2(\eta^n, \eta^{n-1}) && \text{on } \Sigma, \end{aligned} \quad (3.14)$$

where we have set $u_y|_{\Sigma} = 0$ and, again, F_1 and F_2 account for terms at previous time steps. We have the following

Proposition 2. *The reduction factor of iterations (3.13)-(3.14) is given by*

$$\rho^{S/GS}(k) = \left| \frac{\frac{\rho_s H_s}{\Delta t} + \beta H_s \Delta t + GH_s \Delta t k^2 - \alpha_f}{\frac{\rho_s H_s}{\Delta t} + \beta H_s \Delta t + GH_s \Delta t k^2 + \alpha_s} \cdot \frac{\alpha_s k - \mu \gamma_f (k + \gamma_f)}{\alpha_f k + \mu \gamma_f (k + \gamma_f)} \right|, \quad (3.15)$$

where

$$\gamma_f(k) = \sqrt{\rho_f / (\mu \Delta t) + k^2}, \quad (3.16)$$

which is independent of the iteration m .

Moreover, for all Fourier modes $k \in [0, k_{max}]$ there exist $\alpha_f^{max} = \alpha_f^{max}(k)$ and $\alpha_s^{max} = \alpha_s^{max}(k)$ such that, given $\mathbf{u}^0, p^0, \eta^0$, iterations (3.13)-(3.14) converge if $\alpha_f \in [0, \alpha_f^{max})$ and $\alpha_s \in [0, \alpha_s^{max})$ and if at least one of the two values is different from 0.

Proof. In order to write the fluid Robin interface condition (3.13)₃ in terms of the sole pressure, we need again to write the interface velocity \mathbf{u} as a function of p . However, in this case we have

$$\partial_x p = -\rho_f \delta_t u_x + \mu \Delta u_x \quad \text{on } \Sigma,$$

so that the expression of u_x at the interface in terms of p is not straightforward anymore. However, by applying the divergence operator to the fluid problem (3.13)₁ we obtain again problem (3.9) for the pressure solely. We notice that the solution of this problem after applying the Fourier transform in the y direction is again of the form (3.10), for a suitable $A(k)$, which in this case depends also on Δu_x .

Now, the fluid problem in the x direction after the application of the Fourier transform in the y direction, reads

$$\frac{\rho_f}{\mu \Delta t} \widehat{u}_x - \partial_{xx} \widehat{u}_x + k^2 \widehat{u}_x + \frac{A}{\mu} k e^{kx} = \frac{\rho_f}{\mu \Delta t} \widehat{u}_x^n, \quad \text{in } \Omega_f, \quad (3.17)$$

having noticed that $\partial_x \widehat{p} = kA e^{kx}$. The homogeneous solutions of this equation are $\widehat{u}_{x,1}^o(x, k) = B(k) e^{\gamma_f x}$ and $\widehat{u}_{x,2}^o(x, k) = C(k) e^{-\gamma_f x}$, for suitable B and C and where γ_f is given by (3.16). From the boundedness assumption, it follows $C \equiv 0$. As before, in the convergence analysis the terms at the previous time steps are discarded, so that we drop them from now on. As particular solution let us consider $\widehat{u}_x^p(x, k) = -\frac{kA}{2\mu\gamma_f(k-\gamma_f)} e^{\gamma_f x} + \frac{kA}{\mu(k+\gamma_f)(k-\gamma_f)} e^{kx}$. Adding this solution to the homogeneous ones, we obtain

$$\widehat{u}_x(x, k) = \left(B - \frac{kA}{2\mu\gamma_f(k-\gamma_f)} \right) e^{\gamma_f x} + \frac{kA}{\mu(k+\gamma_f)(k-\gamma_f)} e^{kx}.$$

The equation for u_y reads

$$\frac{\rho_f}{\mu \Delta t} \widehat{u}_y - \partial_{xx} \widehat{u}_y + k^2 \widehat{u}_y - ik \frac{A}{\mu} e^{kx} = \frac{\rho_f}{\mu \Delta t} \widehat{u}_y^n, \quad \text{in } \Omega_f,$$

having noticed that $\partial_y \widehat{p} = -ikA e^{kx}$. Then, as before, it is possible to show that the solution is given by

$$\widehat{u}_y(x, k) = \left(D + \frac{ikA}{2\mu\gamma_f(k-\gamma_f)} \right) e^{\gamma_f x} - \frac{ikA}{\mu(k+\gamma_f)(k-\gamma_f)} e^{kx}, \quad (3.18)$$

for a suitable $D(k)$.

From the incompressibility constraint, we have $\partial_x \widehat{u}_x - ik \widehat{u}_y = 0$, obtaining

$$\left(\gamma_f B - ikD + \frac{kA}{2\mu\gamma_f} \right) e^{\gamma_f x} = 0$$

and then

$$\gamma_f B - ikD + \frac{kA}{2\mu\gamma_f} = 0. \quad (3.19)$$

From the interface condition for the velocity u_y , namely $u_y = 0$ for $x = 0$, we obtain

$$\widehat{u}_y(0, k) = D + \frac{ikA}{2\mu\gamma_f(k-\gamma_f)} - \frac{ikA}{\mu(k+\gamma_f)(k-\gamma_f)} = 0,$$

and then $D = -\frac{ikA}{2\mu\gamma_f(k+\gamma_f)}$, which inserted in (3.19) leads to

$$\gamma_f B + \frac{kA}{2\mu(k+\gamma_f)} = 0. \quad (3.20)$$

Then, the normal stress at the FS interface in the x direction is

$$(\hat{p} - \mu\partial_x\hat{u}_x)|_{x=0} = A - \mu\left(\gamma_f B - \frac{kA}{2\mu(k-\gamma_f)} + \frac{k^2 A}{\mu(k+\gamma_f)(k-\gamma_f)}\right) = A, \quad (3.21)$$

where the last equality is obtained thanks to (3.20).

Let us finally compute the value of the velocity in the x direction at the FS interface:

$$\hat{u}_x|_{x=0} = B - \frac{kA}{2\mu\gamma_f(k-\gamma_f)} + \frac{kA}{\mu(k+\gamma_f)(k-\gamma_f)} = -\frac{kA}{\mu\gamma_f(k+\gamma_f)}, \quad (3.22)$$

where the last equality is obtained again thanks to (3.20).

Therefore, owing to (3.21) and (3.22), the interface Robin conditions (3.13)₃ and (3.14) read

$$\begin{aligned} -\frac{\alpha_f k A^{m+1}}{\mu\gamma_f(k+\gamma_f)} - A^{m+1} &= \left(-\frac{\rho_s H_s}{\Delta t^2} - \beta H_s - G H_s k^2 + \frac{\alpha_f}{\Delta t}\right) \hat{\eta}^m, \quad \text{for } x = 0, \\ \left(\frac{\rho_s H_s}{\Delta t^2} + \beta H_s + G H_s k^2 + \frac{\alpha_s}{\Delta t}\right) \hat{\eta}^m &= -\frac{\alpha_s k A^m}{\mu\gamma_f(\gamma_f+k)} + A^m, \quad \text{for } x = 0. \end{aligned}$$

We thus have $|A^{m+1}(k)| = \rho(k)|A^m(k)|$, where the reduction factor ρ is given by (3.15).

In order to guarantee convergence, we have to find suitable values of α_f and α_s such that $\rho^{S/GS} < 1$. Again, to do this, we obtain sufficient conditions by imposing that the two factors in (3.15) are separately less than 1. The first term is equal to the previous analysis so that the upper bound α_f^{max} is given again by (3.11). The second term is $d_2 = \frac{\alpha_s k - \mu\gamma_f(k+\gamma_f)}{\alpha_f k + \mu\gamma_f(k+\gamma_f)}$, so that it is sufficient to take as upper bound

$$\alpha_s^{max}(k) = \frac{2\mu\gamma_f(k+\gamma_f)}{k}. \quad (3.23)$$

A less sharp upper bound, but independent of k , is given by

$$\alpha_s^{max} = \frac{2\mu\gamma_f^*(k^* + \gamma_f^*)}{k^*}, \quad (3.24)$$

where $\gamma_f^* = \gamma_f(k^*)$ and $k^* = \sqrt{\frac{(\sqrt{5}-1)\rho_f}{2\mu\Delta t}}$ is such that α_s^{max} in (3.23) is minimal.

Remark 4. We point out that functions $A(k)$, $B(k)$ and $D(k)$ are in general complex. In particular, if the initial condition for the x component of the fluid velocity is symmetric with respect to the x -axis, A and B are real. In this case (3.17) has real coefficients so that u_x is symmetric for each time. Conversely, if the initial condition for the y component of the fluid velocity is emi-symmetric with respect to the x -axis, then equation (3.18) has imaginary coefficients, so that D has to be imaginary and u_y is emi-symmetric for each time. This is not surprising, since in the y direction we have two waves propagating in the two opposite directions.

Remark 5. For the DN scheme we obtain

$$\rho^{S/GS} = \left| \frac{\mu\gamma_f(k + \gamma_f)\Delta t}{k(\rho_s H_s + \beta H_s \Delta t^2 + G H_s \Delta t^2 k^2)} \right| = \left| \frac{\rho_f + k^2 \mu \Delta t + \sqrt{\rho_f \mu k^2 \Delta t + \mu^2 k^4 \Delta t^2}}{k(\rho_s H_s + \beta H_s \Delta t^2 + G H_s \Delta t^2 k^2)} \right|$$

which highlights again the high dependence of the convergence performances on the ratio ρ_f/ρ_s . Also in this case, for a general RR scheme the sensitivity of $\rho^{S/GS}$ on this ratio is minimal.

3.3. Stokes - Linear incompressible elasticity. Let us consider a thick structure, whose thickness is H_s . Again, the structure domain is unbounded in the y direction, so that $\Omega_s = [0, H_s] \times \mathbb{R}$. We assume that the structure is incompressible. This is an assumption verified in many applications, for example in haemodynamics. Moreover, as in the previous problems, we set $u_y|_\Sigma = 0$, that is we allow the FS interface to move only in the normal direction. This assumption is not realistic in many practical applications, but, in the case of haemodynamics, which inspired the present work, the transversal displacement is typically much smaller than the normal one. Moreover, as in the generalized string model, to emulate a ‘‘cylindrically-shaped’’ structure and take into account for transversal membrane effects, we add a reaction term with coefficient β .

Therefore, the (discretized in time) structure problem reduces to a generalized Stokes problem and the RR algorithm reads:

1. Fluid problem

$$\begin{aligned} \rho_f \delta_t \mathbf{u}^{m+1} - \mu \Delta \mathbf{u}^{m+1} + \nabla p^{m+1} &= \mathbf{0}, & \text{in } \Omega_f, \\ \nabla \cdot \mathbf{u}^{m+1} &= 0, & \text{in } \Omega_f, \\ \alpha_f u_x^{m+1} - p^{m+1} + \mu \partial_x u_x^{m+1} &= \frac{\alpha_f}{\Delta t} \eta_x^m - \chi^m + \lambda \partial_x \eta_x^m + F_1(\eta_x^n, \eta_x^{n-1}), & \text{on } \Sigma. \end{aligned} \quad (3.25)$$

2. Structure problem

$$\begin{aligned} \rho_s \delta_{tt} \boldsymbol{\eta}^m - \lambda \Delta \boldsymbol{\eta}^m + \beta \boldsymbol{\eta}^m + \nabla \chi^m &= \mathbf{0}, & \text{in } \Omega_s, \\ \nabla \cdot \boldsymbol{\eta}^m &= 0, & \text{in } \Omega_s, \\ \frac{\alpha_s}{\Delta t} \eta_x^m + \chi^m - \lambda \partial_x \eta_x^m &= \alpha_s u_x^{m+1} + p^{m+1} - \mu \partial_x u_x^{m+1} + F_2(\eta_x^n, \eta_x^{n-1}), & \text{on } \Sigma, \end{aligned} \quad (3.26)$$

where χ is the pressure for the structure problem and we have set $u_y|_\Sigma = \eta_y|_\Sigma = 0$. We have the following

Proposition 3. *The reduction factor of iterations (3.25)-(3.26) is given by*

$$\rho^{S/S}(k) = \left| \frac{2\alpha_f \delta - \varepsilon \Delta t - \Delta t}{2\alpha_s \delta + \varepsilon \Delta t + \Delta t} \cdot \frac{\alpha_s k - \mu\gamma_f(k + \gamma_f)}{\alpha_f k + \mu\gamma_f(k + \gamma_f)} \right| \quad (3.27)$$

where γ_f is given by (3.16), and

$$\begin{aligned} \varepsilon(k) &= -1 + \frac{2k\gamma_s^2}{k^2 - \gamma_s^2} H_s - \frac{2k^2\gamma_s^4}{(k^2 - \gamma_s^2)^2} H_s^2 + O(H_s^3), \\ \delta(k) &= \frac{(\varepsilon(k)+1)}{2\lambda\gamma_s} \left(\frac{1}{\gamma_s H_s} - \frac{\gamma_s H_s}{6} \right) + O(H_s^2) = \frac{k}{\lambda(k^2 - \gamma_s^2)} \left(1 - \frac{k\gamma_s^2 H_s}{k^2 - \gamma_s^2} \right) + O(H_s^2), \\ \gamma_s(k) &= \sqrt{\rho_s / (\lambda \Delta t^2) + \beta / \lambda + k^2}. \end{aligned} \quad (3.28)$$

Moreover, for all Fourier modes $k \in [0, k_{max}]$ there exist $\alpha_f^{max} = \alpha_f^{max}(k)$ and $\alpha_s^{max} = \alpha_s^{max}(k)$ such that, given $\mathbf{u}^0, p^0, \eta^0$, iterations (3.25)-(3.26) converge if $\alpha_f \in [0, \alpha_f^{max})$ and $\alpha_s \in [0, \alpha_s^{max})$ and if at least one of the two values is different from 0.

Proof. Let us apply the divergence operator to the structure equation (3.26)₁, obtaining again a problem for the pressure solely:

$$\Delta \chi = 0, \quad \text{in } \Omega_s,$$

with suitable boundary conditions depending on the value of $\boldsymbol{\eta}$. Again, we apply the Fourier transform in the y direction, so that the solution can be written as

$$\widehat{\chi}(x, k) = E(k) e^{kx} + F(k) e^{-kx}, \quad x > 0, \quad (3.29)$$

for suitable functions E and F . In this case it is not possible anymore to exploit a boundedness assumption at infinity, so that both terms in (3.29) have to be considered.

Applying the Fourier transform along the y direction to the first component of the structure problem (3.26)₁, we obtain

$$\frac{\rho_s}{\lambda \Delta t^2} \widehat{\eta}_x - \partial_{xx} \widehat{\eta}_x + k^2 \widehat{\eta}_x + \frac{\beta}{\lambda} \widehat{\eta}_x + \frac{kE}{\lambda} e^{kx} - \frac{kF}{\lambda} e^{-kx} = 0, \quad x > 0,$$

having noticed that $\partial_x \widehat{\chi} = kE e^{kx} - kF e^{-kx}$. The solution of this problem is given by

$$\begin{aligned} \widehat{\eta}_x(x, k) &= \frac{kE(k)}{\lambda(k^2 - \gamma_s^2)} e^{kx} - \frac{kF(k)}{\lambda(k^2 - \gamma_s^2)} e^{-kx} + \left(L(k) - \frac{kE(k)}{2\lambda\gamma_s(k - \gamma_s)} - \frac{kF(k)}{2\lambda\gamma_s(k + \gamma_s)} \right) e^{\gamma_s x} + \\ &+ \left(M(k) + \frac{kE(k)}{2\lambda\gamma_s(k + \gamma_s)} + \frac{kF(k)}{2\lambda\gamma_s(k - \gamma_s)} \right) e^{-\gamma_s x}, \end{aligned} \quad (3.30)$$

where γ_s is given by (3.28)₃ and $L(k)$ and $M(k)$ have to be properly determined.

The equation for $\widehat{\eta}_y$ reads

$$\frac{\rho_s}{\lambda \Delta t^2} \widehat{\eta}_y - \partial_{xx} \widehat{\eta}_y + k^2 \widehat{\eta}_y + \frac{\beta}{\lambda} \widehat{\eta}_y - \frac{ikE}{\lambda} e^{kx} - \frac{ikF}{\lambda} e^{-kx} = 0, \quad x > 0,$$

having noticed that $\partial_y \widehat{\chi} = -ikE e^{kx} - ikF e^{-kx}$. The solution of this problem is given by

$$\begin{aligned} \widehat{\eta}_y(x, k) &= \frac{-ikE(k)}{\lambda(k^2 - \gamma_s^2)} e^{kx} - \frac{ikF(k)}{\lambda(k^2 - \gamma_s^2)} e^{-kx} + \left(N(k) + \frac{ikE(k)}{2\lambda\gamma_s(k - \gamma_s)} - \frac{ikF(k)}{2\lambda\gamma_s(k + \gamma_s)} \right) e^{\gamma_s x} + \\ &+ \left(P(k) - \frac{ikE(k)}{2\lambda\gamma_s(k + \gamma_s)} + \frac{ikF(k)}{2\lambda\gamma_s(k - \gamma_s)} \right) e^{-\gamma_s x}, \end{aligned} \quad (3.31)$$

where $N(k)$ and $P(k)$ are suitable functions of k .

Let us impose the incompressibility constraint $\partial_x \widehat{\eta}_x - ik \widehat{\eta}_y = 0$. We obtain

$$\left(\gamma_s L - ikN + \frac{kE}{2\lambda\gamma_s} - \frac{kF}{2\lambda\gamma_s} \right) e^{\gamma_s x} + \left(-\gamma_s M - ikP - \frac{kE}{2\lambda\gamma_s} + \frac{kF}{2\lambda\gamma_s} \right) e^{-\gamma_s x} = 0.$$

This equation has to be satisfied for all $x > 0$, so that it leads to the following conditions

$$\begin{aligned} \gamma_s(L - M) - ik(N + P) &= 0 \\ \gamma_s L - ikN + \frac{k(E-F)}{2\lambda\gamma_s} &= 0. \end{aligned} \quad (3.32)$$

Let us now impose that the velocity in the y direction at the FS interface is zero. From (3.31), we obtain

$$\hat{\eta}_y|_{x=0} = N + P = 0,$$

so that from (3.32)₁ we have $L = M$.

We observe that in the expressions of $\hat{\eta}_x$ and $\hat{\eta}_y$ in (3.30) and (3.31) we do not rely on any boundedness assumption, since the structure domain is bounded in the x direction. However, we can prescribe that the normal stress at the outlet $x = H_s$ is zero. In particular, we have

$$\begin{aligned} (\lambda\partial_x\hat{\eta}_x - \hat{\chi})|_{x=H_s} &= 0 \\ \frac{\lambda}{2}(\partial_x\hat{\eta}_y - ik\hat{\eta}_x)|_{x=H_s} &= 0. \end{aligned}$$

The first condition allows to give a relation among E, F and L , namely

$$\begin{aligned} L &= \frac{2\gamma_s^2 E e^{kH_s} + 2\gamma_s^2 F e^{-kH_s} - (kE(k + \gamma_s) + kF(k - \gamma_s))e^{\gamma_s H_s}}{2\lambda\gamma_s(k^2 - \gamma_s^2)(e^{-\gamma_s H_s} - e^{\gamma_s H_s})} + \\ &- \frac{(kE(k - \gamma_s) + kF(k + \gamma_s))e^{-\gamma_s H_s}}{2\lambda\gamma_s(k^2 - \gamma_s^2)(e^{-\gamma_s H_s} - e^{\gamma_s H_s})} = \frac{E + F}{2\lambda\gamma_s} \left(\frac{1}{\gamma_s H_s} - \frac{\gamma_s H_s}{6} \right) + O(H_s^2). \end{aligned} \quad (3.33)$$

The second condition allows to write N as function of E and F , namely

$$\begin{aligned} N &= -\frac{ikL}{\gamma_s} + \frac{1}{\gamma_s(e^{\gamma_s H_s} + e^{-\gamma_s H_s})} \left[\frac{2ik^2 E}{\lambda(k^2 - \gamma_s^2)} e^{kH_s} - \frac{2ik^2 F}{\lambda(k^2 - \gamma_s^2)} e^{-kH_s} + \right. \\ &- \frac{ik}{2\lambda\gamma_s(k^2 - \gamma_s^2)} \left((k + \gamma_s)^2 E + (k - \gamma_s)^2 F \right) e^{\gamma_s H_s} + \\ &\left. + \frac{ik}{2\lambda\gamma_s(k^2 - \gamma_s^2)} \left((k - \gamma_s)^2 E + (k + \gamma_s)^2 F \right) e^{-\gamma_s H_s} \right] = \\ &= -\frac{ik}{2\lambda\gamma_s^2} (E + F) \left(\frac{1}{\gamma_s H_s} - \frac{\gamma_s H_s}{6} + \frac{k - \gamma_s}{k + \gamma_s} k H_s \right) + O(H_s^2). \end{aligned}$$

This expression, together with (3.32)₂ allows to derive a relation between E and F , namely

$$F(k) = \varepsilon(k) E(k),$$

with ε given by (3.28)₁. From (3.33), we have $L(k) = \delta(k) E(k)$, with δ given by (3.28)₂.

Let us compute the values of the displacement and of the normal stress at the FS interface. From (3.30), we obtain

$$\widehat{\eta}_x|_{x=0} = 2L = 2\delta E,$$

and from (3.29) and (3.30)

$$(\lambda\partial_x\widehat{\eta}_x - \widehat{\chi})|_{x=0} = -E - F = -(1 + \varepsilon)E.$$

The fluid problem is identical to the one analyzed in the previous section (Stokes/generalized string). We have therefore

$$\begin{aligned} p - \mu\partial_x\widehat{u}_x &= A & \text{for } x = 0, \\ \widehat{u}_x &= -\frac{kA}{\mu\gamma_f(k+\gamma_f)} & \text{for } x = 0. \end{aligned}$$

Then the Robin interface conditions (3.25)₃ and (3.26)₃ read

$$\begin{aligned} -\frac{\alpha_f k}{\mu\gamma_f(k+\gamma_f)}A^{m+1} - A^{m+1} &= \left(\frac{2\alpha_f\delta}{\Delta t} - \varepsilon - 1\right)E^m & \text{for } x = 0, \\ \left(\frac{2\alpha_s\delta}{\Delta t} + \varepsilon + 1\right)E^m &= -\frac{\alpha_s k}{\mu\gamma_f(k+\gamma_f)}A^m + A^m & \text{for } x = 0, \end{aligned}$$

and the reduction factor is given by (3.27).

In order to guarantee convergence, we have to find suitable values of α_f and α_s such that $\rho^{S/GS} < 1$. Again, to do this, we obtain sufficient conditions by imposing that the two factors in (3.27) are separately less than 1. The second term is equal to the previous analysis so that the upper bound α_s^{max} is given again by (3.23). The first term is $d_1 = \frac{2\alpha_f\delta - \varepsilon\Delta t - \Delta t}{2\alpha_s\delta + \varepsilon\Delta t + \Delta t}$, so that it is sufficient to take as upper bound

$$\alpha_f^{max}(k) = \frac{(\varepsilon + 1)\Delta t}{\delta}.$$

Remark 6. For the DN scheme we obtain

$$\rho^{S/S} = \left| \frac{2\delta\mu\gamma_f(k+\gamma_f)}{k\Delta t(\varepsilon+1)} \right| = \left| \frac{\mu\gamma_f(k+\gamma_f)}{\lambda\gamma_s^2 H_s k\Delta t} \right| = \left| \frac{a_1\rho_f + a_2\sqrt{\rho_f} + a_3}{a_4\rho_s + a_5} \right|,$$

for suitable $a_j = a_j(k, \mu, \lambda, \beta, H_s, \Delta t)$, $j = 1, \dots, 5$, and again the DN scheme highlights a high dependence of the convergence performances on the ratio ρ_f/ρ_s . In particular, we point out that $\lim_{\Delta t \rightarrow 0} \rho^{S/S} = \rho_f/(k\rho_s H_s)$. Also in this case, for a general RR scheme the sensitivity of $\rho^{S/S}$ on this ratio is minimal.

4. Optimization of α_f and α_s . With the aim of constructing RR schemes with good convergence properties, the problem of determining suitable values for parameters α_f and α_s has to be addressed. Indeed, the convergence performances of RR schemes heavily depend on this choice (see, in a different context, [15]). By maximizing the convergence rate of the corresponding algorithm, it is possible to determine either optimal or optimized values of the coefficients α_f and α_s in the Robin interface conditions. The classical approach in the Optimized Schwarz Methods literature consists in looking for parameters that, if possible, annihilate identically the reduction factor, ensuring convergence for the scheme in just two iterations for a two-domain decomposition. Such parameters are referred to as *optimal*. In particular, in a two-domain decomposition framework, we recall that the optimal interface conditions

are of Robin type, where the Dirchlet-to-Neumann operator of one domain is applied as a boundary condition for the other domain. When such approach is not viable, the idea is to look for parameters that minimize the reduction factor.

In this section we look for optimized values for α_f and α_s either k -independent $\alpha_f = \alpha_{f,M}$, $\alpha_s = \alpha_{s,M}$, which lead to standard Robin boundary conditions (with standard mass matrix at the interface), or k -dependent values of the form $\alpha_f = \alpha_{f,M} + \alpha_{f,K}k^2$, $\alpha_s = \alpha_{s,M} + \alpha_{s,K}k^2$, which correspond to generalized Robin boundary conditions on Σ

$$\alpha_{f,M}\mathbf{u} - \alpha_{f,K}\Delta^\Sigma\mathbf{u} + \mathbf{T}_f\mathbf{n} = \alpha_{f,M}\left(\frac{\boldsymbol{\eta} - \boldsymbol{\eta}^n}{\Delta t}\right) - \alpha_{f,K}\Delta^\Sigma\left(\frac{\boldsymbol{\eta} - \boldsymbol{\eta}^n}{\Delta t}\right) + \mathbf{T}_s\mathbf{n}, \quad (4.1)$$

$$\frac{\alpha_{s,M}}{\Delta t}\boldsymbol{\eta} - \alpha_{s,K}\Delta^\Sigma\boldsymbol{\eta} + \mathbf{T}_s\mathbf{n} = \frac{\alpha_{s,M}}{\Delta t}\boldsymbol{\eta}^n - \alpha_{s,K}\Delta^\Sigma\boldsymbol{\eta}^n + \alpha_{s,M}\mathbf{u} - \alpha_{s,K}\Delta^\Sigma\mathbf{u}^n + \mathbf{T}_f\mathbf{n}, \quad (4.2)$$

where Δ^Σ is the Laplace-Beltrami operator over the interface Σ .

Before presenting the optimization procedures, we recall in the next subsection the choices proposed in [3] for a heuristic determination of the parameters.

4.1. Determination of α_f and α_s by heuristic methods. In [20] a membrane model has been proposed to describe in an approximated way the behaviour of thin structures. It is possible to “insert” the membrane equation into the fluid one as a boundary term (we highlight that in this case the structure domain coincides with the FS interface). This leads to a fluid problem with a suitable Robin boundary condition at the FS interface, which accounts for the structure (see [20]). In [2,3], this formulation has been used to obtain a plausible value of α_f in RR schemes, namely

$$\alpha_f^{heur} = \frac{\rho_s H_s}{\Delta t} + \beta H_s \Delta t, \quad (4.3)$$

where H_s is the structure thickness and

$$\beta = \frac{E}{1 - \nu^2}(4\rho_1^2 - 2(1 - \nu)\rho_2), \quad (4.4)$$

with ρ_1 and ρ_2 the mean and Gaussian curvatures of the FS interface, respectively. We point out that α_f^{heur} is a function of the position on the interface.

RN schemes with α_f given by (4.3) exhibit very good performances for a wide range of added mass and are by far more efficient than DN strategies (see [2,3]). In particular, the theoretical analysis presented in these works, have shown that RN schemes with α_f^{heur} are less sensitive than DN schemes to the added-mass effect. Moreover, numerical experiments with high added-mass effect, highlighted that the RN scheme converges without any relaxation and faster than the DN scheme. However, in [2] RN has been shown to be sensitive to the choice of α_f , leading to a deterioration of the performances when the curvature of the FS interface could not be computed accurately. On the contrary, the RN-GMRES scheme proposed in [2] is more robust with respect to the choice of α_f .

For what concerns the parameter α_s , in [3] it has been proved that the operator describing the added mass effect on the structure is not algebraic and its approximation by an algebraic relationship is not evident. In particular a generalized Robin condition (that involves differential operators) is obtained with $\alpha_s = (\rho_f/\Delta t^2)\mathcal{M}(\cdot)$, where \mathcal{M} is the added-mass operator, that is the Neumann-to-Dirichlet map (see [3,6]). In

order to obtain a “classical” Robin condition, in [3] it has been proposed to approximate the operator \mathcal{M} by $\theta\mu_{max}I$, where μ_{max} is the maximum eigenvalue of the added-mass operator, I is the identity operator and θ is a coefficient suitably chosen, getting

$$\alpha_s^{heur} = \theta \frac{\rho_f \mu_{max}}{\Delta t}. \quad (4.5)$$

The drawback of this choice is that the scaling factor θ has to be tuned “by hand” to obtain good convergence properties. The numerical experiments in [3] showed that the tuned value θ seems to be very robust and practically independent of ρ_f , Δt and of some geometrical parameters defining the physical domain (and then μ_{max}). However, the performance of the RR scheme with the choices (4.3) and (4.5) is only slightly better than RN with (4.3).

4.2. Optimization procedures. Our goal is to minimize the reduction factor ρ in the different cases considered, by choosing the Robin parameters α_f , α_s in the class of frequency dependent functions $\mathcal{A} \equiv \{\alpha(k) = \alpha_M + \alpha_K k^2, \alpha_M, \alpha_K \geq 0\}$. Functions in this class will lead to generalized Robin boundary conditions having a Mass and a Stiffness matrix on the boundary (see (4.1) and (4.2)).

In all the problems considered the reduction factor $\rho(k)$ can be split as the product of two factors: $\rho(k) = d_1(k) \cdot d_2(k)$ where

$$d_1(k) = \left| \frac{\frac{\rho_s H_s}{\Delta t} + \beta H_s \Delta t + G H_s \Delta t k^2 - \alpha_f}{\frac{\rho_s H_s}{\Delta t} + \beta H_s \Delta t + G H_s \Delta t k^2 + \alpha_s} \right| \quad (\text{Generalized string model})$$

$$d_1(k) = \left| \frac{2\alpha_f \delta - (\varepsilon + 1) \Delta t}{2\alpha_s \delta + (\varepsilon + 1) \Delta t} \right| \quad (\text{Linear incompressible elasticity})$$

and

$$d_2(k) = \left| \frac{\rho_f - \alpha_s \Delta t k}{\rho_f + \alpha_f \Delta t k} \right|, \quad (\text{Potential flow})$$

$$d_2(k) = \left| \frac{\alpha_s k - \mu \gamma_f (k + \gamma_f)}{\alpha_f k + \mu \gamma_f (k + \gamma_f)} \right|, \quad (\text{Stokes})$$

Since the global optimization of ρ for all frequencies is too difficult, we try to minimize separately the two factors $d_1 = d_1(k, \alpha_f, \alpha_s)$ and $d_2 = d_2(k, \alpha_f, \alpha_s)$. The first can be associated to the structure problem and will lead to an optimal selection of $\alpha_f \in \mathcal{A}$ for all possible frequencies in $[0, k_{max}]$ and all possible functions $\alpha_s \in \mathcal{A}$. Similarly, the second factor can be associated to the fluid problem and leads to the optimal selection of $\alpha_s \in \mathcal{A}$ uniformly with respect to k and $\alpha_f \in \mathcal{A}$.

We set therefore the two optimization problems

$$\alpha_f^{opt}(k) = \operatorname{argmin}_{\alpha_f(k) \in \mathcal{A}} \max_{k \in [0, k_{max}]} \sup_{\alpha_s \in \mathcal{A}} d_1(k, \alpha_f(k), \alpha_s(k)),$$

$$\alpha_s^{opt}(k) = \operatorname{argmin}_{\alpha_s(k) \in \mathcal{A}} \max_{k \in [0, k_{max}]} \sup_{\alpha_f \in \mathcal{A}} d_2(k, \alpha_f(k), \alpha_s(k)).$$

Of course, by this procedure we will not get the global optimum but, hopefully, improve the heuristic choice of the Robin coefficients.

4.2.1. Optimization of α_f . We start with the optimization of the factor d_1 .

Generalized string model. In this case we obtain $d_1 \equiv 0$ for

$$\alpha_f^{GS} = \frac{\rho_s H_s}{\Delta t} + \beta H_s \Delta t + G H_s \Delta t k^2. \quad (4.6)$$

We point out that $\alpha_f^{GS} < \alpha_f^{max}$, with α_f^{max} given in (3.11). The previous expression represents the symbol, in the Fourier space, of the Dirichlet-to-Neumann operator in the structure domain. We notice that $\alpha_f^{GS} = \alpha_f^{heur} + G H_s \Delta t k^2$, and then the convergence analysis in Sect. 3.1 provides a correction of the optimal value of α_f with respect to the heuristic choice (4.3). The dependence of this correction on k introduces in the Robin interface condition a differential term. In particular, we obtain condition (4.1) with

$$\alpha_{f,M}^{GS} = \alpha_f^{heur} = \frac{\rho_s H_s}{\Delta t} + \beta H_s \Delta t, \quad \alpha_{f,K}^{GS} = G H_s \Delta t.$$

Therefore, the correction given by the convergence analysis of the potential flow/generalized string problem requires to build a “stiffness” interface matrix.

Linear incompressible elasticity. We have $d_1 \equiv 0$ for $\alpha_f(k) = \frac{(1+\varepsilon)\Delta t}{2\delta}$, which, however, is not in the class \mathcal{A} . By considering an approximation of such expression for H_s small, that is by using the expansions of ε and δ in powers of H_s , given in (3.28), we obtain the optimized value

$$\alpha_f^{LIE} = \Delta t \lambda \gamma_s^2 H_s = \frac{\rho_s H_s}{\Delta t} + \beta H_s \Delta t + \lambda H_s \Delta t k^2,$$

which leads to $d_1 = O(H_s^3)$. Observe that we obtain an expression very similar to α_f^{GS} .

The previous result shows that the approximated reduction factor obtained by considering H_s small is the same as the one that is obtained when a membrane model is considered for the structure.

4.2.2. Optimization of α_s . We consider now the optimization of the factor d_2 .

Potential flow. In this case the optimization problem reads

$$\alpha_s^{opt}(k) = \operatorname{argmin}_{\alpha_s(k) \in \mathcal{A}} \max_{k \in [0, k_{max}]} \sup_{\alpha_f \in \mathcal{A}} \left| \frac{\rho_f - \alpha_s \Delta t k}{\rho_f + \alpha_f \Delta t k} \right| = \operatorname{argmin}_{\alpha_s(k) \in \mathcal{A}} \max_{k \in [0, k_{max}]} \left| 1 - \frac{\alpha_s \Delta t k}{\rho_f} \right|$$

It is clear that for a generic choice $\alpha_s(k) = \alpha_{s,M} + \alpha_{s,K} k^2$ with $\alpha_{s,K} > 0$ the factor $d_2(k)$ is unbounded for $k \rightarrow \infty$ (that is for $h \rightarrow 0$). Therefore we restrict the optimization problem only to frequency-independent functions, namely $\alpha_s(k) = \alpha_{s,M}$. In this case, we have

$$\max_{k \in [0, k_{max}]} \left| 1 - \frac{\alpha_s \Delta t k}{\rho_f} \right| = \begin{cases} 1 & \text{if } \alpha_{s,M} \leq \frac{2\rho_f}{\Delta t k_{max}}, \\ > 1 & \text{otherwise.} \end{cases}$$

Therefore, no choice of $\alpha_{s,M}$ will lead to a factor d_2 strictly smaller than 1, and we can choose any $\alpha_{s,M} \in [0, \frac{2\rho_f}{\Delta t k_{max}}]$ to have exactly $d_2 = 1$. However, in the expressions of d_1 , α_s is at the denominator, so that in view of the optimization of the factor d_1

we have interest in choosing $\alpha_{s,M}$ as large as possible, leading to the *optimized* Robin coefficient

$$\alpha_s^P = \frac{2\rho_f}{\Delta t k_{max}}. \quad (4.7)$$

Notice that this choice coincides with the upper bound of the stability interval given in (3.12)

Stokes flow. We proceed in a similar way as for the Potential flow. In this case we have

$$\begin{aligned} \alpha_s^{opt}(k) &= \operatorname{argmin}_{\alpha_s(k) \in \mathcal{A}} \max_{k \in [0, k_{max}]} \sup_{\alpha_f \in \mathcal{A}} \left| \frac{\alpha_s k - \mu\gamma_f(k + \gamma_f)}{\alpha_f k + \mu\gamma_f(k + \gamma_f)} \right| = \\ &= \operatorname{argmin}_{\alpha_s(k) \in \mathcal{A}} \max_{k \in [0, k_{max}]} \left| \frac{\alpha_s k}{\mu\gamma_f(k + \gamma_f)} - 1 \right| \end{aligned}$$

and again we restrict the optimization only to constant values of α_s or otherwise the factor d_2 will be unbounded for $k \rightarrow \infty$.

Let us denote $f(k) = \frac{k}{\mu\gamma_f(k + \gamma_f)} = \frac{k}{\mu\sqrt{a+k^2}(k + \sqrt{a+k^2})}$, with $a = \frac{\rho_f}{\mu\Delta t}$. The function f has a global maximum in $[0, \infty)$ at $k^* = \sqrt{a\frac{\sqrt{5}-1}{2}}$. It is a straightforward calculation to show that

$$\max_{k \in [0, k_{max}]} |\alpha_s f(k) - 1| = \begin{cases} 1 & \text{if } \alpha_s f(k^*) \leq 2 \\ > 1 & \text{otherwise} \end{cases}$$

Proceeding as for the potential flow we take the largest value for which $d_2 = 1$ leading to the *optimized coefficient for the Stokes flow*

$$\alpha_s^S = \frac{2}{f(k^*)} = \frac{2}{\Delta t k^*} \sqrt{\rho_f + \mu\Delta t (k^*)^2} \left(\sqrt{\mu\Delta t k^*} + \sqrt{\rho_f + \mu\Delta t (k^*)^2} \right). \quad (4.8)$$

4.3. Alternative optimization procedure. Another possible way to proceed to derive optimized parameters for the RR algorithm consists in taking a different split of the reduction factor, namely $\rho(k, \alpha_f, \alpha_s) = \tilde{d}_1(k, \alpha_f) \cdot \tilde{d}_2(k, \alpha_s)$. In such a case, we can look for parameters α_f and α_s that optimize separately the two factors, namely:

$$\alpha_f^{opt}(k) = \operatorname{argmin}_{\alpha_f \in \mathcal{A}} \max_{k \in [0, k_{max}]} \tilde{d}_1(k, \alpha_f(k)), \quad \alpha_s^{opt}(k) = \operatorname{argmin}_{\alpha_s \in \mathcal{A}} \max_{k \in [0, k_{max}]} \tilde{d}_2(k, \alpha_s(k)).$$

This optimization procedure will lead to the same values for α_f as in the previous section since these are the values that zero the factor d_1 . However, this procedure will lead to different values for α_s which will depend, this time, both on the fluid and the structure problem chosen.

4.3.1. Optimization of α_s . Potential flow / Generalized string. For the Potential Flow / Generalized string model, we have

$$\tilde{d}_2(k, \alpha_s) = \left| \frac{\rho_f - \alpha_s \Delta t k}{\frac{\rho_s H_s}{\Delta t} + \beta H_s \Delta t + G H_s \Delta t k^2 + \alpha_s} \right|.$$

First, we observe that if $\alpha_s = \alpha_{s,M} + \alpha_{s,K} k^2$ with $\alpha_{s,K} > 0$, the factor \tilde{d}_2 is unbounded for $k \rightarrow \infty$, so we restrict the optimization to constant values. Let us denote

$$\tilde{f}(k, \alpha_s) = \frac{\rho_f - \alpha_s \Delta t k}{\frac{\rho_s H_s}{\Delta t} + \beta H_s \Delta t + G H_s \Delta t k^2 + \alpha_s}.$$

The function $\tilde{f}(k, \alpha_s)$ is such that $\tilde{f}(0, \alpha_s) = \frac{\rho_f}{\frac{\rho_s H_s}{\Delta t} + \beta H_s \Delta t + \alpha_s} > 0$, $\frac{\partial \tilde{f}}{\partial k}(0, \alpha_s) < 0$ and $\lim_{k \rightarrow \infty} \tilde{f}(k, \alpha_s) = 0$, and has exactly one global minimum in $k^*(\alpha_s) \in [0, \infty)$, with $\tilde{f}(k^*(\alpha_s), \alpha_s) < 0$. We have, therefore that the optimal value α_s^{opt} satisfies (see Fig. 4.1)

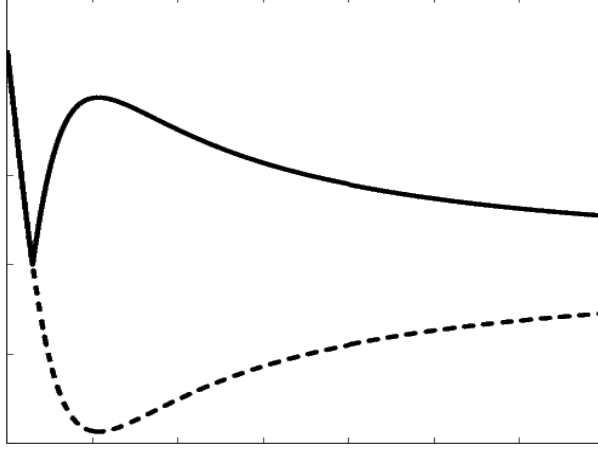


FIG. 4.1. \tilde{f} as a function of k .

$$\min_{\alpha_s} \max_k |\tilde{f}(k, \alpha_s)| = \min_{\alpha_s} \max\{\tilde{f}(0, \alpha_s), -\tilde{f}(k^*(\alpha_s), \alpha_s)\}.$$

We observe that $\tilde{f}(0, \alpha_s)$ is a decreasing function of α_s and $-\tilde{f}(k^*(\alpha_s), \alpha_s)$ is an increasing function of α_s . Moreover, for $\alpha_s = 0$ we have that $\tilde{f}(k, 0) > 0$ is a monotone decreasing function in k , so that, thanks to the continuity of \tilde{f} , we have $\tilde{f}(0, 0^+) > -\tilde{f}(k^*(0^+), 0^+)$. Therefore, the minimum is achieved for

$$\alpha_s^{P,2} \text{ such that } \tilde{f}(0, \alpha_s^{P,2}) = -\tilde{f}(k^*(\alpha_s^{P,2}), \alpha_s^{P,2}). \quad (4.9)$$

This non-linear problem can be solved numerically given all the parameters of the fluid and structure model.

Stokes / Generalized String In this case we have

$$\tilde{d}_2(k, \alpha_s) = \left| \frac{\mu \gamma_f (k + \gamma_f) - \alpha_s \Delta t k}{\frac{\rho_s H_s}{\Delta t} + \beta H_s \Delta t + G H_s \Delta t k^2 + \alpha_s} \right|.$$

Again, we observe that if $\alpha_s = \alpha_{s,M} + \alpha_{s,K} k^2$ with $\alpha_{s,K} > 0$, the factor \tilde{d}_2 is unbounded for $k \rightarrow \infty$, so we restrict the optimization to constant values. Let us denote

$$\tilde{f}(k, \alpha_s) = \frac{\mu \gamma_f (k + \gamma_f) - \alpha_s \Delta t k}{\frac{\rho_s H_s}{\Delta t} + \beta H_s \Delta t + G H_s \Delta t k^2 + \alpha_s}.$$

The function $\tilde{f}(k, \alpha_s)$ is such that $\tilde{f}(0, \alpha_s) = \frac{\rho_f}{\rho_s H_s + \beta H_s \Delta t^2 + \alpha_s \Delta t} > 0$, $\frac{\partial \tilde{f}}{\partial k}(0, \alpha_s) < 0$ and $\lim_{k \rightarrow \infty} \tilde{f}(k, \alpha_s) = \tilde{F} := \frac{\mu}{G H_s \Delta t} > 0$, and has exactly one local minimum in $k^*(\alpha_s) \in [0, \infty)$, with $\tilde{f}(k^*(\alpha_s), \alpha_s) < 0$. Let us assume that k_{max} is big enough to approximate $\tilde{f}(k_{max}, \alpha_s)$ with \tilde{F} . We have, therefore that the optimal value α_s^{opt} satisfies

$$\min_{\alpha_s} \max_k |\tilde{f}(k, \alpha_s)| = \min_{\alpha_s} \max\{\tilde{f}(0, \alpha_s), -f(k^*(\alpha_s), \alpha_s), \tilde{F}\}.$$

Again, $\tilde{f}(0, \alpha_s)$ is a decreasing function of α_s and $-f(k^*(\alpha_s), \alpha_s)$ is an increasing function of α_s , with $\tilde{f}(0, 0^+) > -f(k^*(0^+), 0^+)$. Moreover, under the hypothesis on the data $4\mu^2 \rho_s \beta > \rho_f^2 G^2$, we have that $\tilde{f}(0, 0) > \tilde{F}$. Therefore, under this hypothesis, minimum is achieved for

$$\alpha_s^{opt} = \begin{cases} \alpha_{s,1} & \text{if } \tilde{f}(0, \alpha_{s,1}) > \tilde{F}, \\ \alpha_{s,2} & \text{if } \tilde{F} > \tilde{f}(0, \alpha_{s,1}), \end{cases}$$

where

$$\alpha_{s,1} \text{ such that } \tilde{f}(0, \alpha_{s,1}) = -f(k^*(\alpha_{s,1}), \alpha_{s,1}),$$

and

$$\alpha_{s,2} \text{ such that } \tilde{f}(0, \alpha_{s,2}) = \tilde{F}.$$

Again, this non-linear problem can be solved numerically given all the parameters of the fluid and structure model.

5. Numerical Results. In this section we present some numerical results with the aim of testing the optimized parameters found in the previous analysis. We focus on two-dimensional numerical simulations, even if the application of the proposed parameters to three-dimensional problems is under investigation. This is in agreement with the fact that the convergence analysis has been performed in two-dimensional domains. However, the application of the optimal coefficients found in this analysis to three-dimensional cases is under investigation.

For the structure, we consider the following linear elasticity equation

$$\rho_s \partial_{tt} \boldsymbol{\eta} - \lambda_1 \nabla \cdot (\nabla \boldsymbol{\eta} + (\nabla \boldsymbol{\eta})^t) - \lambda_2 \nabla \cdot ((\nabla \cdot \boldsymbol{\eta}) \mathbf{I}) + \beta \boldsymbol{\eta} = \mathbf{0},$$

where the Lamé constants are given by (2.1) and a reaction term β , arising in the derivation of the generalized string model in the 3D case, is introduced also in 2D to reproduce the behaviour of a cylindrical structure. For its expression we use the particularization of (4.4) to the rectangular domain, that is $\beta = \frac{E}{(1-\nu^2)R^2}$, where R is half the height of the rectangle.

For the numerical solution we use a 2D Finite Element Code written in Matlab at MOX - Dipartimento di Matematica - Politecnico di Milano and at the Department of Information Technology and Mathematical Methods - Università degli Studi di Bergamo. Moreover, we consider \mathbb{P}_1 -bubble/ \mathbb{P}_1 elements for the fluid and \mathbb{P}_1 elements for the structure. We use the residual normalized to the initial one as stopping criterion (see [3]), with a tolerance equal to 10^{-4} .

In all the cases, we prescribe an impulse pressure jump between the inlet and the outlets

$$\Delta P(t) = \begin{cases} 10^4 \text{ dyne/cm}^2 & t \leq 0.005 \text{ s}, \\ 0 & t > 0.005 \text{ s}. \end{cases}$$

Finally, we set $\mu = 0.035 \text{ poise}$, $\rho_f = 1 \text{ g/cm}^3$ and the following set of reference values: $\Delta t = 10^{-3} \text{ s}$, $\rho_s = 1.1 \text{ g/cm}^3$, $E = 1.3 \cdot 10^6 \text{ dyne/cm}^2$ and the thickness of the structure $H_s = 0.1 \text{ cm}$.

5.1. Results in a rectangle domain. The numerical simulations of this section are performed in a rectangular domain both for the fluid and for the two structures, whose size is $6 \times 1 \text{ cm}$ and $6 \times H_s \text{ cm}$, respectively (see Fig. 5.1). We set $\nu = 0.3$, $R =$

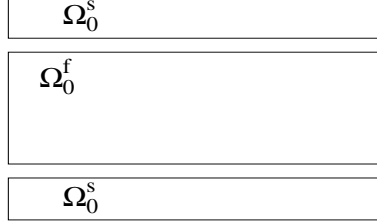


FIG. 5.1. Computational fluid and structure domains.

0.5 and the space discretization parameter is $h = H_s/2$, both for the fluid and the structure domains.

The aim of this test is to compare the performance of the proposed parameters given in (4.6), (4.7) and (4.8) with respect to the heuristic choice. In particular, for the reference values, we have the following values: $\alpha_{f,M}^{GS} = 681.43$, $\alpha_{f,K}^{GS} = 41.12$, $\alpha_s^P = 31.83$ and $\alpha_s^S = 38.95$.

In Table 5.1 we show the average number of iterations to reach convergence having considered as final time $T = 0.008 \text{ s}$. In brackets, we show the computational time, normalized to that obtained with the heuristic choice of α_f and with $\alpha_s = 0$.

$\alpha_f =$ $\alpha_s =$	α_f^{heur} 0	α_f^{GS} 0	α_f^{GS} α_s^P	α_f^{GS} α_s^S
Ref.	5.87 (1.00)	5.37 (0.91)	4.62 (0.79)	4.50 (0.77)
$\rho_s = 0.1$	6.00 (1.00)	5.62 (0.94)	4.87 (0.81)	4.87 (0.81)
$\rho_s = 0.01$	6.00 (1.00)	5.75 (0.96)	5.00 (0.83)	4.87 (0.81)
$H_s = 0.05$	6.75 (1.00)	5.87 (0.87)	5.00 (0.74)	5.00 (0.74)
$H_s = 0.0125$	10.75 (1.00)	7.87 (0.73)	7.62 (0.71)	5.75 (0.53)
$\Delta t = 5 \cdot 10^{-4}$	7.00 (1.00)	5.69 (0.81)	4.94 (0.71)	4.87 (0.70)
$\Delta t = 2.5 \cdot 10^{-4}$	6.94 (1.00)	5.56 (0.80)	4.81 (0.69)	4.41 (0.64)
$E = 6.5 \cdot 10^5$	6.50 (1.00)	5.62 (0.86)	4.87 (0.75)	4.87 (0.75)
$E = 1.3 \cdot 10^5$	7.50 (1.00)	5.50 (0.73)	4.37 (0.58)	5.00 (0.67)

TABLE 5.1

Average number of iterations and relative CPU times (in brackets) for the rectangular domain for different values of some parameters. Ref. means reference values.

These results show that the Robin-Robin schemes are very robust with respect to the added-mass effect. Indeed also for structure density 100 times lighter than the fluid one, convergence is reached without any relaxation and with a small number of iterations. However, a worsening in the performance is observed when the structure thickness H_s , the time step Δt and the Young modulus E decrease. Moreover, these results show that there is a systematic improvement in the performance of the Robin-Neumann scheme when adopting the optimized value α_f^{GS} instead of the heuristic

one, with a saving in the computational time up to 27%. A further improvement is obtained by considering the Robin-Robin scheme with coefficient α_s^P or α_s^S . In these cases the saving in computational time is up to 47%. We point out that the proposed choice is robust, since numerical simulation highlights that the number of iterations is independent of the value of the space discretization parameter h .

In the second set of simulations, we want to compare the two different optimization strategies proposed for the parameter α_s in Sect. 4.2 and 4.3, respectively. We limit the analysis to the Potential flow case, that is we compare the performances of the RR scheme with α_s given by (4.7) and (4.9).

In Table 5.2 we show the average number of iterations to reach convergence having considered 16 time steps, by changing the Young modulus and the time step. These

$\alpha_s =$	α_s^P	$\alpha_s^{P,2}$
Ref.	4.62	10.87
$H_s = 0.0125$	7.73	X
$\Delta t = 6.4 \cdot 10^{-5}$	6.00	5.00
$\Delta t = 3.2 \cdot 10^{-5}$	10.94	9.00
$E = 1.3 \cdot 10^4$	4.62	6.37
$E = 1.3 \cdot 10^3$	26.75	15.00

TABLE 5.2

Average number of iterations to perform 16 time steps for the rectangular domain. Ref. means reference values. X means no convergence.

results clearly show that the optimized parameter $\alpha_s^{P,2}$ shows a better performance when the structure stiffness term is small in comparison to the other terms. However, as highlighted by the performances obtained with the reference values and by the non-convergence with $H_s = 0.0125$, this choice seems less robust than α_s^P .

5.2. An application to a 2D bifurcation geometry. In this section we investigate the performance of Robin-Robin algorithm with optimized parameters considering a 2D geometry which is an idealization of a realistic domain, namely the human carotid. We use the reference values described in the previous sections and we set $\nu = 0.49$ in order to model the quasi-incompressibility of arterial vessels.

In Figure 5.2 the pressure in the deformed fluid domain is shown, at three time instants. In Table 5.3 we show the average number of iterations and the normalized

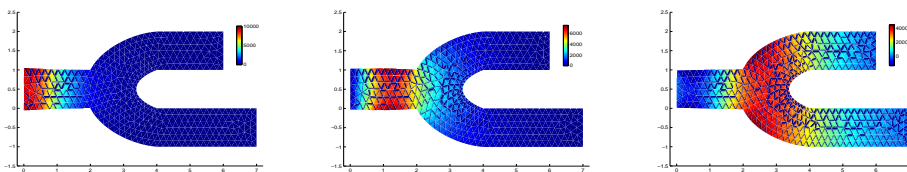


FIG. 5.2. Pressure in the deformed fluid domain obtained with $\alpha_f = \alpha_f^{GS}$ and $\alpha_s = \alpha_s^P$; $t = 0.004$ (left), $t = 0.008$ s (middle), $t = 0.012$ s (right).

computational times (in brackets) for two values of the structure thickness, having considered as final time $T = 0.012$ s.

These results show that the improvement in the performances of RR scheme in using the optimized parameters α_f^{GS} , α_s^P and α_s^S is even slightly higher than for the rectangular case.

$\alpha_f =$ $\alpha_s =$	α_f^{heur} 0	α_f^{GS} 0	α_f^{GS} α_s^P	α_f^{GS} α_s^S
Ref.	13.08 (1.00)	9.92 (0.76)	7.00 (0.54)	6.75 (0.52)
$H_s = 0.05$	23.25 (1.00)	14.08 (0.61)	13.50 (0.58)	12.67 (0.54)

TABLE 5.3

Average number of iterations and relative CPU times (in brackets) for the carotid domain for different values of the structure thickness H_s . Ref. means reference values.

6. Conclusions. In this work, we have proposed new convergence analysis of the Robin-Robin partitioned procedure for the Fluid-Structure Interaction problem. In particular, we consider 2D generalized Stokes problem both for the fluid and for structure. These analysis improve and generalize the ones presented in [3, 6] which referred to a potential flow model for the fluid and a 1D reduced model for the structure. The analysis have been performed applying the Fourier transform in one direction, and highlight the dependence of the Dirichlet-Neumann scheme on the ratio between the densities of the fluid and of the structure, and the robustness of a general RR scheme with respect to this ratio.

Moreover, in the spirit of the Optimized Schwarz Methods, the reduction factor of the analysis have been minimized by determining optimal values of the parameters in the Robin transmission conditions.

The 2D numerical results show that the proposed optimized parameters improve considerably the speed of convergence of RR schemes.

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