

A parallel iterative approach for the Stokes–Darcy coupling

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Stokes–Darcy coupling

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Let us start describing the coupled problem. We denote by Ω_S **the fluid region**, and by Ω_D **the ground region**. Moreover, $\Gamma := \overline{\Omega_S} \cap \overline{\Omega_D}$ will be **the interface** between Ω_S and Ω_D .

Stokes–Darcy coupling (cont.)

The **Stokes system** is given by

$$(1) \quad \begin{cases} -\nu \Delta \mathbf{u} + \nabla p & = \mathbf{f} & \text{in } \Omega_S \\ \operatorname{div} \mathbf{u} & = 0 & \text{in } \Omega_S, \end{cases}$$

where \mathbf{u} is the velocity field, p is the pressure, $\nu > 0$ is the kinematic viscosity and \mathbf{f} is a given force field.

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The **Darcy equation** is given by

$$(2) \quad -\operatorname{div} \left(\frac{K}{N} \nabla \varphi \right) = 0 \text{ in } \Omega_D,$$

where φ is the piezometric head, K is the hydraulic conductivity tensor and $N > 0$ is the volumetric porosity.

Stokes–Darcy coupling (cont.)

For simplicity, as **boundary conditions** let us assume that:

$$(3) \quad \begin{aligned} \mathbf{u} &= \mathbf{0} && \text{on } \partial\Omega_S \setminus \Gamma \\ \varphi &= 0 && \text{on } \partial\Omega_D \setminus \Gamma. \end{aligned}$$

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The **interface conditions** [matching of the normal velocity and of the normal stress] are:

$$(4) \quad \begin{aligned} \mathbf{u} \cdot \mathbf{n} &= -\frac{K}{N} \nabla \varphi \cdot \mathbf{n} && \text{on } \Gamma \\ \mathbf{T}(\mathbf{u}, p) \cdot \mathbf{n} &= -g\varphi \mathbf{n} && \text{on } \Gamma \end{aligned}$$

where g is the gravity acceleration, and the fluid stress tensor is given by $T_{ij}(\mathbf{u}, p) := \nu(D_i u_j + D_j u_i) - p\delta_{ij}$. Here \mathbf{n} denotes the unit normal vector on Γ , pointing from Ω_S into Ω_D .

Well-posedness and iterative solution algorithms

It can be proved that the coupled problem (1)–(4) **has a unique solution** [Discacciati and Quarteroni, ENUMATH 2001; Layton, Schieweck and Yotov, SINUM 2003].

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Let us present some of the **iterative algorithms** that have been proposed for the solution of (1)–(4).

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- pose on Γ

$$\varphi^{m+1} := \theta \varphi^{m+1/2} + (1 - \theta) \varphi^m,$$

where $\theta > 0$ is an acceleration parameter.

Dirichlet/Neumann algorithm (cont.)

[The name Dirichlet/Neumann is somehow arbitrary: indeed, we are solving two Neumann problems. However, for the velocity field $(\mathbf{u}|_{\Omega_S}, -\nabla\varphi|_{\Omega_D})$ the step in Ω_S is a Neumann step, whereas the step in Ω_D is a Dirichlet step (for the normal component...)].

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However:

- for a finite element approximation, the convergence is **independent** of the mesh parameter h , but **depends heavily** on the viscosity ν and the conductivity K

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Here $\gamma_S > 0$ and $\gamma_D > 0$ are suitable acceleration parameters, and τ is a unit tangent vector on Γ .

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Moreover:

- it can be interpreted as an **alternating direction** algorithm (which is useful for tuning the parameters γ_S and γ_D)
- for a finite element approximation, the convergence is **independent** of the mesh parameter h , and, for suitable choices of γ_S and γ_D , in the numerical computations it looks also **independent** of the viscosity ν and the conductivity K

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Recalling that the Neumann/Neumann algorithm for the Laplace operator is indeed given by a couple of Dirichlet solvers, followed by a couple of (homogeneous) Neumann solvers, which play the role of correctors, we propose a **double parallel Robin/Robin algorithm**.

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- solve in parallel the Stokes problem and the Darcy problem with

$$(5) \quad \begin{aligned} \mathbf{n} \cdot \mathbf{T}(\mathbf{u}^{m+1}, p^{m+1}) \cdot \mathbf{n} - \gamma_1 \mathbf{u}^{m+1} \cdot \mathbf{n} &= \eta^m \\ &= -g\varphi^{m+1} + \gamma_1 \frac{K}{N} \nabla \varphi^{m+1} \cdot \mathbf{n} \quad \text{on } \Gamma \\ \boldsymbol{\tau} \cdot \mathbf{T}(\mathbf{u}^{m+1}, p^{m+1}) \cdot \mathbf{n} &= 0 \quad \text{on } \Gamma \end{aligned}$$

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$$(6) \quad \sigma^{m+1} := \mathbf{u}^{m+1} \cdot \mathbf{n} + \frac{K}{N} \nabla \varphi^{m+1} \cdot \mathbf{n}$$

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- solve in parallel the homogeneous ($\mathbf{f} = 0$) Stokes problem and the Darcy problem with

$$(7) \quad \begin{aligned} \mathbf{n} \cdot \mathbf{T}(\boldsymbol{\omega}^{m+1}, \pi^{m+1}) \cdot \mathbf{n} + \gamma_2 \boldsymbol{\omega}^{m+1} \cdot \mathbf{n} &= \gamma_2 \sigma^{m+1} \\ &= g \chi^{m+1} - \gamma_2 \frac{K}{N} \nabla \chi^{m+1} \cdot \mathbf{n} \quad \text{on } \Gamma \\ \boldsymbol{\tau} \cdot \mathbf{T}(\boldsymbol{\omega}^{m+1}, \pi^{m+1}) \cdot \mathbf{n} &= 0 \quad \text{on } \Gamma \end{aligned}$$

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$$(8) \quad \eta^{m+1} := \eta^m - \theta (\mathbf{n} \cdot \mathbf{T}(\boldsymbol{\omega}^{m+1}, \pi^{m+1}) \cdot \mathbf{n} + g\chi^{m+1}),$$

where $\gamma_1 > 0$, $\gamma_2 > 0$ and $\theta > 0$ are acceleration parameters.

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therefore **normal stresses of the corrections are matching**.

On the other hand, from (7) **normal velocities of the corrections are jumping**

$$\boldsymbol{\omega}^\infty \cdot \mathbf{n} + \frac{K}{N} \nabla \chi^\infty \cdot \mathbf{n} = \frac{2g}{\gamma_2} \chi^\infty.$$

Consistency of the Robin/Robin algorithm 2 (cont.)

But this jump gives an additional **positive** term in the energy of the problem, namely, one obtains

$$\int_{\Omega_S} 2\nu \sum_{ij} |D_i \omega_j^\infty|^2 + \int_{\Omega_D} \frac{g}{N} K \nabla \chi^\infty \cdot \nabla \chi^\infty + \int_{\Gamma} \frac{2g^2}{\gamma_2} |\chi^\infty|^2 = 0.$$

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Hence, $\omega^\infty = 0$ and $\chi^\infty = 0$, and consequently $\sigma^\infty = 0$.

Consistency of the Robin/Robin algorithm 2 (cont.)

Finally:

- from (6), σ^∞ is the jump of the normal velocities, therefore **they are matching**:

$$\mathbf{u}^\infty \cdot \mathbf{n} = -\frac{K}{N} \nabla \varphi^\infty \cdot \mathbf{n}$$

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We have found the right solution!

Convergence of the Robin/Robin algorithm 2

The parallel Robin/Robin algorithm can be rewritten as a **preconditioned Richardson scheme**:

$$\eta^{m+1} = \eta^m + \theta(K_S + K_D)[\psi - (H_S + H_D)\eta^m],$$

for a suitable right hand side ψ .

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The operators K_S , K_D , H_S and H_D are the main building blocks of the algorithm, and are defined as follows.

Convergence of the Robin/Robin algorithm 2 (cont.)

- $K_S \sigma := \mathbf{n} \cdot \mathbf{T}(\boldsymbol{\omega}, \pi) \cdot \mathbf{n}$, where $\boldsymbol{\omega}$ is a solution with the Robin datum

$$\begin{aligned}\mathbf{n} \cdot \mathbf{T}(\boldsymbol{\omega}, \pi) \cdot \mathbf{n} + \gamma_2 \boldsymbol{\omega} \cdot \mathbf{n} &= \gamma_2 \sigma \quad \text{on } \Gamma \\ \boldsymbol{\tau} \cdot \mathbf{T}(\boldsymbol{\omega}, \pi) \cdot \mathbf{n} &= 0 \quad \text{on } \Gamma.\end{aligned}$$

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- $H_S \eta := \mathbf{u} \cdot \mathbf{n}$, where \mathbf{u} is a solution (for $\mathbf{f} = 0$) with the Robin datum

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- $H_D \eta := \frac{K}{N} \nabla \varphi \cdot \mathbf{n}$, where φ is a solution with the Robin datum

$$-g\varphi + \gamma_1 \frac{K}{N} \nabla \varphi \cdot \mathbf{n} = \eta \quad \text{on } \Gamma.$$

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Therefore, both the operator $(H_S + H_D)$ and the preconditioner $(K_S + K_D)^{-1}$ are symmetric, continuous and positive definite: **convergence is achieved!** (for a suitable choice of the parameter θ)

Comments

- It is well-known that the Dirichlet-to-Neumann operator is symmetric, continuous and positive definite from the energy trace space $H^{1/2}(\Gamma)$ into its dual; here we have seen that:

the **Robin-to-Neumann operator** is symmetric, continuous and **positive definite** in $L^2(\Gamma)$,
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whereas:
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- the above results also hold for the finite element numerical approximation, **uniformly** with respect to the mesh parameter h