Application of the Fictitious Domain method to flow problems with complex geometries

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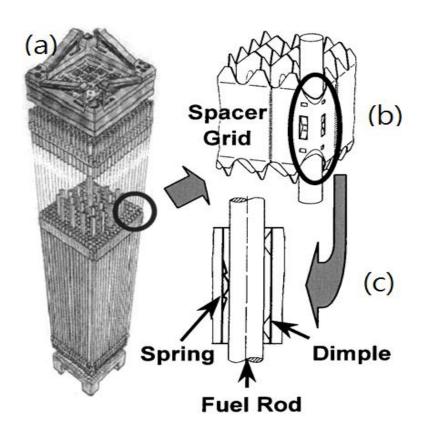


OUTLINE

- 1. Introduction
- 2. Case 1: Poisson problem
- 3. Case 2: Stokes flow problem
- 4. Future work



Introduction

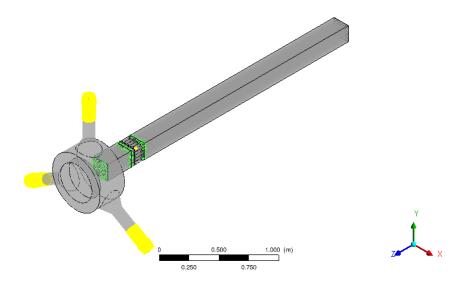


 Flow problems with complex geometries are widespread

PWR fuel assembly and spacer grid spring (Kim et al. 2001)



CFD simulation for one smaller sized assembly,
 68 million cells (Wei 2013,)

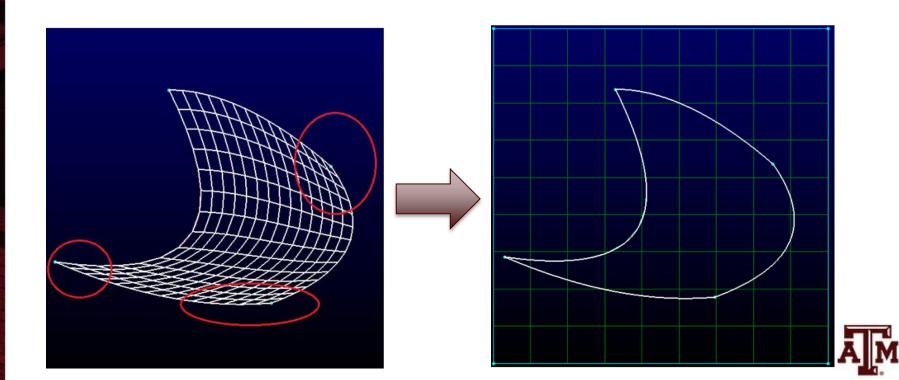


- Body-fitted mesh is not suitable
- Our choice: Fictitious Domain method (FD method)



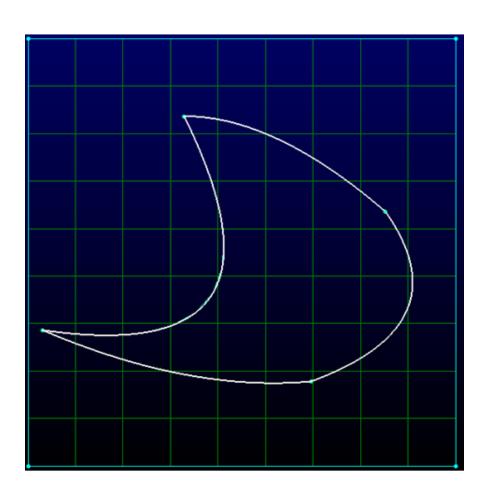
Basic concept regarding the FD method:

Whenever a problem needs to be solved on a domain with an irregular boundary, it may be useful to embed it into a larger domain of a simpler shape (Quarteroni and Valli).



Critical point regarding the FD method implementation:

How to include the influence from the "immersed" BC





Strategies regarding this immersed boundary issue:

1. Penalty function method (Ramière et al. 2005, Zhou and Saito 2014, Saito and Zhou 2014)

$$\frac{1}{\varepsilon} \int_{\Omega \setminus \omega} (v \cdot \tilde{u}) dx$$

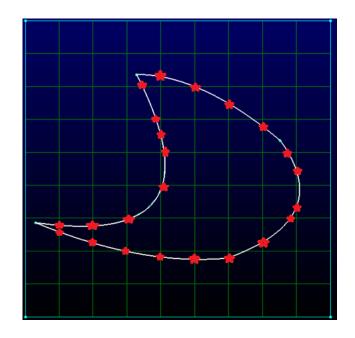
 \Rightarrow limitation exists $(e \sim C(\sqrt{\varepsilon} + h))$



Strategies regarding this immersed boundary issue:

2. Lagrange multiplier method (Glowinski et al. 1994, Glowinski et al. 1995, Glowinski et al. 1998)

$$\int_{\gamma} \lambda \cdot (u - g) dx$$

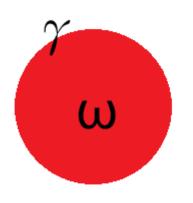


⇒ Powerful, but may be not suitable for problems with complex 3d geometries



Case 1: Poisson problem

Governing equation



$$-\Delta u = f \qquad in \ \omega,$$

$$u = g \qquad on \ \gamma,$$

where
$$f \in H^{-1}(\omega)$$
, $g \in H^{\frac{1}{2}}(\gamma)$, which is Dirichlet BC



Weak formulation

$$a_{\omega}(v,u) = \langle v, f \rangle \ \forall v \in H_0^1(\omega),$$

where

$$a_{\omega}(v,u) = \int_{\omega} (\nabla v \cdot \nabla u) dx \qquad \forall u, v \in H_0^1(\omega),$$
$$\langle v, f \rangle = \int_{\omega} (v \cdot f) dx.$$

Here, V_g is the solution space and its Definition is:

$$V_g = \{ v \mid v \in H^1(\omega), v = g \text{ on } \gamma \}.$$

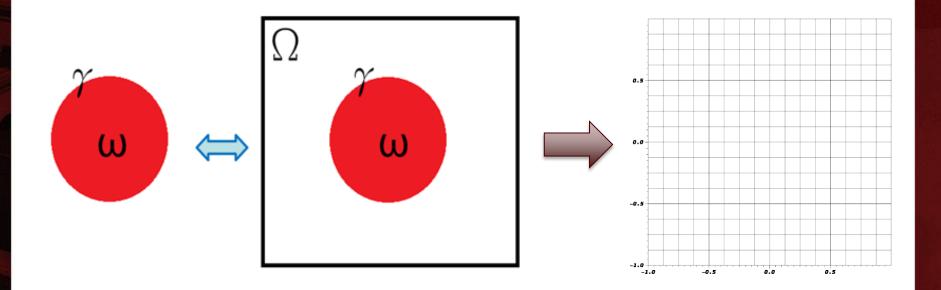
Lagrangian functional:

$$L(u,\lambda) = \frac{1}{2} \int_{\omega} (\nabla u \cdot \nabla u) dx - \int_{\omega} (f \cdot u) dx + \int_{\gamma} \lambda \cdot (u - g) dx \qquad in \ \omega$$



FD method implementation

1. Embed the original domain





FD method implementation

2. Define Lagrangian functional for the fictitious domain Ω :

$$L(\tilde{u},\lambda) = \frac{1}{2} \int_{\Omega} (\nabla \tilde{u} \cdot \nabla \tilde{u}) dx - \int_{\Omega} (\tilde{f} \cdot \tilde{u}) dx + \int_{\gamma} \lambda \cdot (\tilde{u} - \tilde{g}) dx \quad in \Omega.$$

Solution \tilde{u} is valid in the whole domain Ω and the Lagrange multiplier λ is valid on the immersed boundary γ . In addition, $\tilde{f}|_{\Omega} = f$ and $\tilde{g}|_{\gamma} = g$.

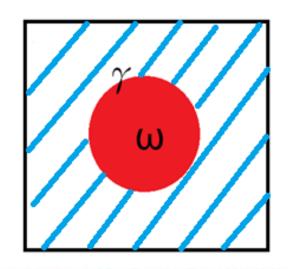


Corresponding weak formulation for Ω :

$$\begin{split} &\int_{\Omega} (\nabla v \cdot \nabla \tilde{u}) dx - \int_{\Omega} (v \cdot \tilde{f}) dx + \int_{\gamma} (v \cdot \lambda) d\gamma = 0, \\ &\int_{\gamma} \mu \cdot (\tilde{u} - \tilde{g}) d\gamma = 0. \end{split}$$

Our strategies

1. Approach 1: penalty lives everywhere



e.g.
$$\int_{\gamma} (v \cdot \lambda) d\gamma \Rightarrow \int_{\Omega \setminus \omega} (v \cdot \lambda) dx$$

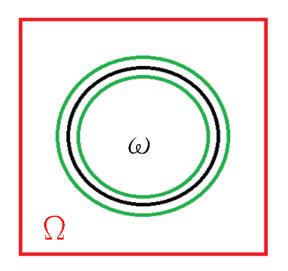
$$\Rightarrow \int_{\Omega} (w(x) \cdot v \cdot \lambda) dx$$

where
$$w(x) = \begin{cases} 1, & x \in \Omega \backslash \omega \\ 0, & x \in \omega \end{cases}$$



Our strategies

2. Approach 2:



$$\overline{\omega}$$
: $d(x) < r_a - \theta$,

$$\bar{\gamma}$$
: $r_a - \theta \le d(x) \le r_a + \theta$,

$$\Omega \setminus \overline{\omega} \setminus \overline{\gamma}$$
: $r_a + \theta \le d(x)$.

$$\theta = c \times h$$

e.g.

$$\int_{\gamma} (v \cdot \lambda) d\gamma \Rightarrow \int_{\Omega} (w(x) \cdot \boldsymbol{k}(\boldsymbol{x}) \cdot v \cdot \lambda) dx$$

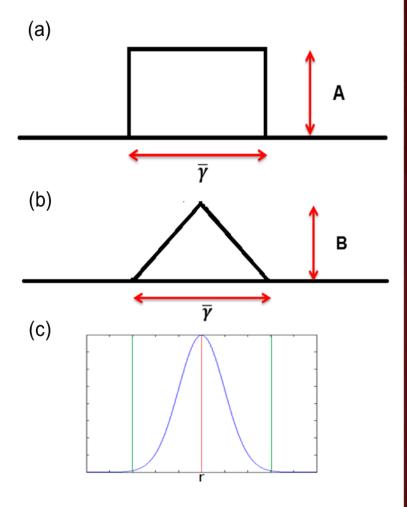
w(x) is equal to one in $\bar{\gamma}$ and equal to zero in the rest areas.



Our strategies

2. Approach 2:about function *k(x)*

$$\int_{\gamma} s(x) dx \approx \int_{\overline{\gamma}} k(x) \cdot s(x) dx$$



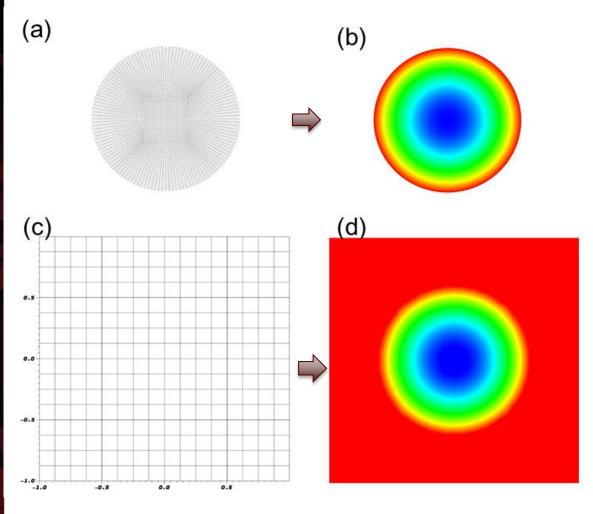
Three k(x) functions



Numerical details

- 1. FEM, u_h , $\lambda_h \in Q_1(K)$
- 2. Gaussian Quadrature for computing the integration
- 3. Direct solver UMFPACK for solving the resulting linear system
- 4. L-2 error norm for the error analysis



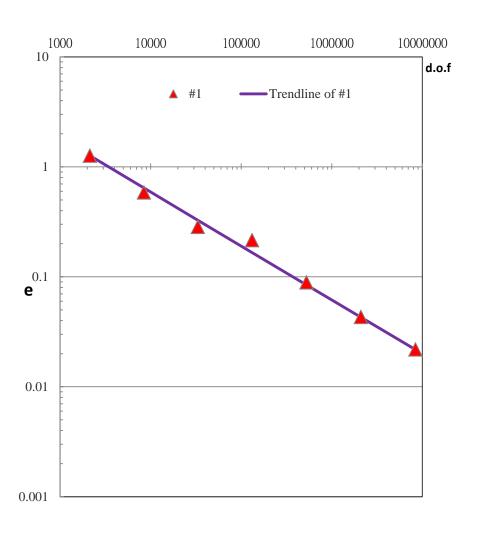


- Solution profile
- In quality, these two results are similar

- (a) and (b) are direct result, computed from step-6
- (c) and (d) are results from the FD method



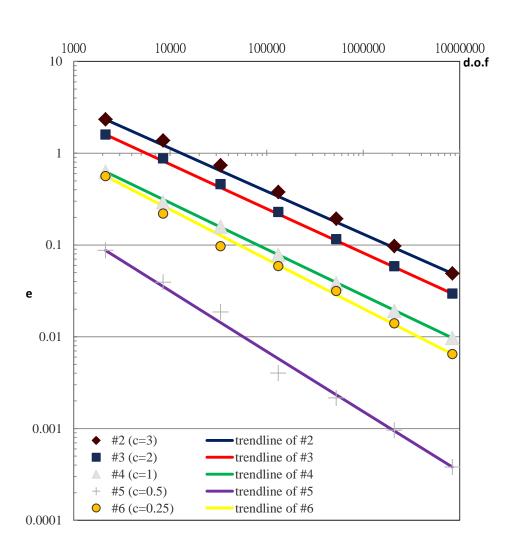
Error analysis: Approach 1



- 1. Oscillation appears in coarse mesh
- 2. $e \sim 0(h^{0.98})$, comparable to the penalty function method (Ramière et al. 2005)



Error analysis: Approach 2, with constant function



1. When θ is small (i.e., c is small), the result is not stable $\begin{pmatrix} \theta = c \times h, \\ c = 0.25, 0.5, 1.0, 2.0 \ and \ 3.0 \end{pmatrix}$

2. $e \sim 0(h^1)$, for c=1 $e \sim 0(h^{1.3})$, for c=0.5

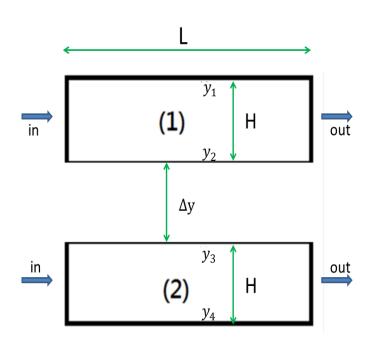


Summary 1

1. Both of our two approaches can get reasonable results



Case 2: Stokes flow problem



Two parallel, fully developed, Poiseuille flow with very low velocity

Governing equation

$$-\operatorname{div}[2\eta \cdot \varepsilon(u)] + \nabla p = 0 \quad in \, \omega,$$
$$-\operatorname{div} u = 0 \quad in \, \omega,$$

where

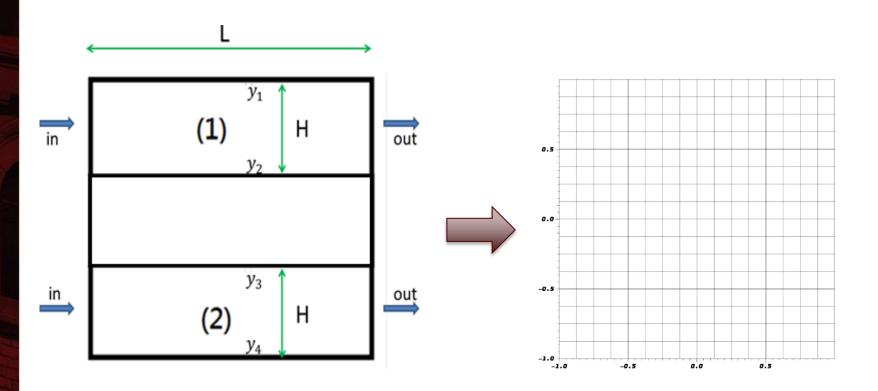
$$\varepsilon(u) = \frac{1}{2} [(\nabla u) + (\nabla u)^T],$$

$$u = 0.0 \text{ at } \Gamma.$$



FD method implementation

1. Embed the original domain





FD method implementation

 Define Lagrangian functional for the fictitious domain Ω:

$$L(\tilde{u}, \tilde{p}, \lambda) = 2\eta \frac{1}{2} \int_{\Omega} |\nabla \tilde{\varepsilon}|^{2} dx - \int_{\Omega} p \nabla \tilde{u} dx$$
$$- \int_{\Omega} (\tilde{F} \cdot \tilde{u}) dx + \int_{\gamma} \lambda \cdot (\tilde{u} - \tilde{g}) dx \text{ in } \Omega,$$

The solutions \tilde{u} and \tilde{p} are valid in the whole domain Ω . In addition, $\tilde{F}\big|_{\omega} = F = 0.0$ and $\tilde{g}\big|_{\gamma} = g = 0.0$.

For simplicity, we set $\tilde{F}\big|_{\Omega\setminus\omega}=0.0$.



Corresponding weak formulation for Ω :

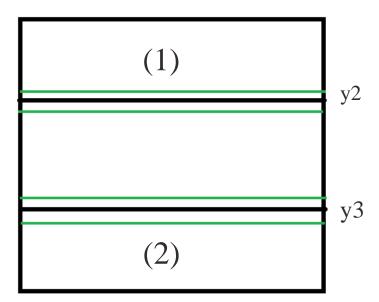
$$2\nu(\varepsilon(v), \varepsilon(\tilde{u}))_{\Omega} - (\nabla v, \tilde{p})_{\Omega} = (v, \tilde{F})_{\Omega} - (v, \lambda)_{\gamma},$$
$$-(q, \nabla \tilde{u})_{\Omega} = 0,$$
$$(\mu, \tilde{u})_{\gamma} = (\mu, g)_{\gamma}.$$



Our strategy for the boundary related term

Approach 2, boundary region $\bar{\gamma}$ is used here

$$\bar{\gamma}$$
: $y_{ib} - hf \le y_{qp} \le y_{ib} + hf$.



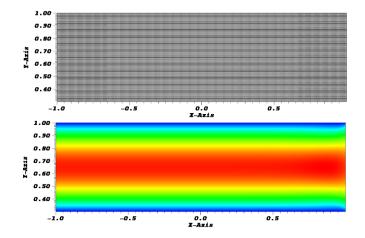


Numerical details

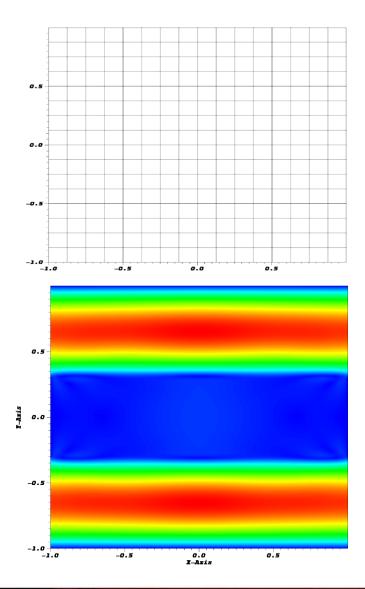
- 1. FEM, u_h , $\lambda_h \in Q_2(K)$, $p_h \in Q_1(K)$
- 2. Gaussian Quadrature for computing the integration
- 3. The resulting linear system is solved iteratively (modified from Glowinski et al. 1995, a variation of the Uzawa algorithm).
- 4. L-2 error norm for the error analysis



Direct simulation's velocity profile(step-22)

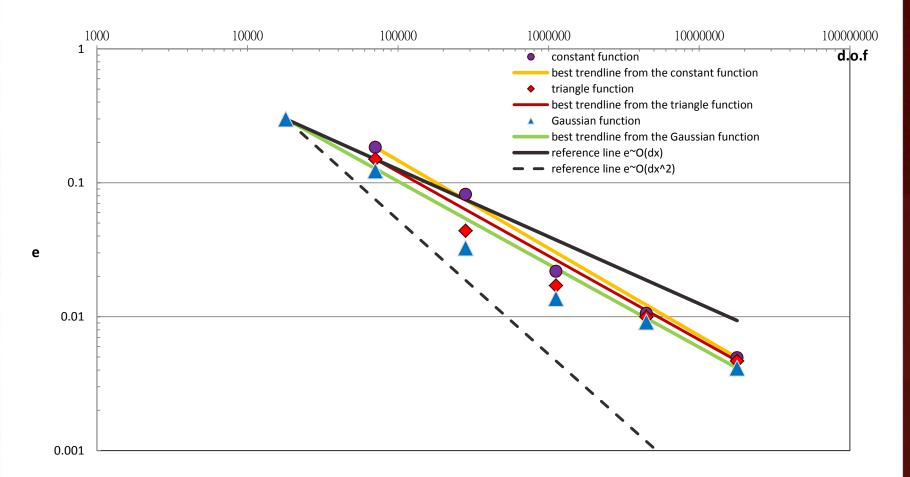


FD method's velocity profile





Error analysis



 Among the three k(x) functions, Gaussian function may be the best, whose error can be expressed as:

 $e \sim h^{1.236}$

Summary 2

- 1. The modified iterative algorithm used here can be parallelized
- 2. What we have from the Case 2 can be the basis for us to solve the Navier-Stokes problem



Future work

1. Parallelize the code for the Stokes flow problem

Solve the Navier-Stokes problem (Operator splitting method)



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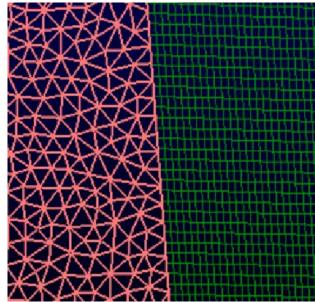


Questions?

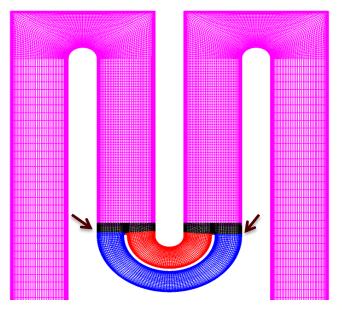


Supplement - introduction

 Several techniques are invented, such as multiblock or overset mesh, but they all cause extra burdens to the user or computers



Interface between two blocks with different meshes (Multi-block)



 Interpolation area between two blocks (Overset mesh)



Weak formulation

Find
$$u \in H^1(\omega), p \in L^2(\omega)$$

$$2\eta \left(\varepsilon(v), \varepsilon(u)\right)_{\omega} - (\nabla v, p)_{\omega} = \langle v, 0 \rangle_{\omega} \quad \forall v \in H^1_0(\omega),$$

$$-(q, \nabla u)_{\omega} = \langle v, 0 \rangle_{\omega} \quad \forall q \in L^2_0(\omega),$$

where

$$\left(\varepsilon(v), \varepsilon(u) \right)_{\omega} = 2\eta \int_{\omega} \varepsilon(v) \cdot \varepsilon(u) \, dx,$$

$$(\nabla v, p)_{\omega} = \int_{\omega} \nabla v \cdot p \, dx, \quad \langle v, 0 \rangle_{\omega} = \int_{\omega} v \cdot 0 \, dx,$$

$$(q, \nabla u)_{\omega} = \int_{\omega} q \cdot \nabla u \, dx, \quad \langle v, G \rangle_{\omega} = \int_{\omega} q \cdot 0 \, dx,$$



Lagrangian functional

$$L(u, p, \lambda) = 2\eta \frac{1}{2} \int_{\omega} |\nabla \varepsilon|^2 dx - \int_{\omega} p \nabla u dx$$

$$-\int_{\omega} (F \cdot u) dx + \int_{\gamma} \lambda \cdot (u - g) dx \text{ in } \omega.$$



Supplement Case 2 Approach

In Glowinski et al. 1995

Concerning the multiplier λ , its interpretation is very simple since it is equal to the jump of

$$v\frac{\partial u}{\partial n} - np$$



Supplement- iterative algorithm 1

Iterative a	lgorithm in	Clowinski	et al 1994
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Iterative algorithm used for solving our designed problem

Initial guess part

Guess λ⁰

2. Find the initial velocity U⁰ and the initial pressure P⁰ by solving the Stokes 2. solve problem

$$\nu \int_{\Omega} \nabla U^{0} \cdot \nabla v \, dx - \int_{\Omega} P^{0} \nabla \cdot v \, dx$$

$$= \int_{\Omega} f \cdot \mathbf{v} \, dx + \int_{\gamma} \lambda^0 \cdot \mathbf{v} \, d\gamma,$$

and

$$\int_{\Omega} q \nabla \cdot U^0 \, \mathrm{d} \mathbf{x} = 0.$$

In addition, impose:

$$U^0 = g_1$$
 on Γ ,

where Γ is the boundary condition for the fictitious domain Ω .

$$\begin{split} &2v\big(\varepsilon(v),\varepsilon(\tilde{u}^0)\big)_{\Omega}-(\nabla v,\tilde{p}^0)_{\Omega}\\ &=\big(v,\tilde{F}\big)_{\Omega}-(w(x)\cdot k(x)\cdot v,\lambda^0)_{\Omega}, \end{split}$$

and

$$-(q,\nabla \tilde{u}^0)_{\Omega}=0.$$

In addition, impose

$$U^0 = g_1$$
 on Γ ,

where

$$\tilde{F}\big|_{\omega}$$
=0.0, $\tilde{F}\big|_{\Omega\setminus\omega}$ =0.0.

$$w(x) = \{ \begin{matrix} 0, & in \ \Omega \setminus \overline{\gamma} \\ in \ \overline{\gamma} \end{matrix}.$$

There are three setting regarding the weight function k(x), when we use Approach 2

3. Compute g⁰:

$$\int_{\gamma} g^{0} \cdot \mu \, d\gamma = \int_{\gamma} (U^{0} - g_{2}) \cdot \mu \, d\gamma$$

3. Compute g^0 by solving

$$\begin{aligned} (w(x) \cdot k(x) \cdot \mu, g^0)_{\Omega} \\ + ([1 - w(x)] \cdot \mu, g^0)_{\Omega} \\ = (w(x) \cdot k(x) \cdot \mu, (\tilde{u}^0 - g))_{\Omega} \\ + ([1 - w(x)] \cdot \mu, 0)_{\Omega} \end{aligned}$$

g = 0.0 in this designed problem (i.e., no-slip boundary condition



Supplement- iterative algorithm 2

4. Set $W^0 = g^0$

Iteration loop starts from step-5

5. Find $\overline{\mathbb{U}}^n$ and \overline{P}^n which satisfy

$$\nu \int_{\Omega} \nabla \overline{U}^n \cdot \nabla v \, dx - \int_{\Omega} \overline{P}^n \nabla \cdot v \, dx = \int_{\gamma} W^n \cdot v \, d\gamma$$

and

$$\int_{\Omega} q \nabla \cdot U^0 \, \mathrm{d} \mathbf{x} = 0.$$

5. Solve

$$2\nu \left(\varepsilon(v), \varepsilon(\overline{U}^n)\right)_{\Omega} - (\overline{v}v, \overline{P}^n)_{\Omega}$$
$$= (w(x) \cdot k(x) \cdot v, W^n)_{\Omega},$$

and

$$-(q,\nabla \overline{U}^n)_{\Omega}=0.$$

6. Compute ρ_n by using:

$$\rho_n = \frac{\int_{\gamma} |g^n|^2 d\gamma}{\int_{\gamma} \overline{U}^n \cdot W^n d\gamma}$$

6. compute

$$o_n = \frac{\int_{\Omega} w(x) \cdot k(x) \cdot |g^n|^2 dx}{\int_{\Omega} w(x) \cdot k(x) \cdot \overline{u}^n \cdot W^n dx}$$

7. Get new λ, U and P:

$$\lambda^{n+1} = \lambda^n - \rho_n W^n$$

$$U^{n+1} = U^n - \rho_n \overline{U}^n$$

$$P^{n+1} = P^n - \rho_n \bar{P}^n$$

Get new λ , \tilde{u} and \tilde{p} :

$$\lambda^{n+1} = \lambda^n - \rho_n W^n$$

$$\tilde{u}^{n+1} = \tilde{u}^n - \rho_n \overline{U}^n$$

$$\tilde{p}^{n+1} = \tilde{p}^n - \rho_n \bar{P}^n$$



Supplement- iterative algorithm 3

8. Renew g^n by

$$\int_{\gamma} g^{n+1} \cdot \mu \, d\gamma$$

$$= \int_{\gamma} g^{n} \cdot \mu \, d\gamma - \rho^{n} \int_{\gamma} \overline{U}^{n} \cdot \mu \, d\gamma$$

8. Renew g^n

$$(w(x) \cdot k(x) \cdot \mu, g^{n+1})_{\Omega}$$

$$+([1 - w(x)] \cdot \mu, g^{n+1})_{\Omega}$$

$$= (w(x) \cdot k(x) \cdot \mu, g^{n})_{\Omega}$$

$$-\rho^{n}(w(x) \cdot k(x) \cdot \mu, \overline{U}^{n})$$

$$+([1 - w(x)] \cdot \mu, g^{n+1})_{\Omega}$$

9. If it reaches the stop criteria:

$$\frac{\int_{\gamma} |g^{n+1}|^2 d\gamma}{\int_{\gamma} |g^0|^2 d\gamma} \le \varepsilon$$

solutions are λ^{n+1} , U^{n+1} and P^{n+1} .

9. Compute the iteration error:

$$\frac{\int_{\Omega} w(x) \cdot k(x) \cdot |g^{n+1}|^2 dx}{\int_{\Omega} w(x) \cdot k(x) \cdot |g^0|^2 dx}$$

If it is equal or smaller than the stop criterion ε , then, the solutions are λ^{n+1} , $\tilde u^{n+1}$ and $\tilde p^{n+1}$

10. If it hasn't reached the stop criteria, compute:

$$r_n = \frac{\int_{\gamma} |g^{n+1}|^2 d\gamma}{\int_{\gamma} |g^n|^2 d\gamma},$$

and get

$$W^{n+1} = g^{n+1} + r_n W^n.$$

Restart to compute the step-5 by using this new W^{n+1}

10. If it hasn't reached the stop criteria, compute:

$$r_n = \frac{\int_{\Omega} w(x) \cdot k(x) \cdot |g^{n+1}|^2 dx}{\int_{\Omega} w(x) \cdot k(x) \cdot |g^n|^2 dx},$$

and get

$$W^{n+1} = g^{n+1} + r_n W^n.$$

Restart to compute the step-5 by using this new W^{n+1}

