# About the use of FreeFEM ++ to solve a diffusion model by domain decomposition methods. 

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#### Abstract

Domain decomposition methods have been particularly developed in recent years with the arrival of parallel computers, and currently represent a very active research direction in numerical analysis. These methods are an essential choice for solving problems on complex geometries, are very effective for large-scale systems, and are also used in the study of coupled models such as the Stokes-Darcy problem. In this work, we present our fundings on domain decomposition methods carried. A series of numerical tests using the FreeFem software were performed and demonstrated the robustness of these methods. The second part concerns a coupled Stokes-Darcy problem. The mathematical formulation as well as the finite element discretization of this problem are presented.


## 1 Introduction

The first domain decomposition method was developed in the late 19th century by mathematician H.A. Schwarz. The goal was to develop an analytical tool that could extend the Dirichlet principle to more complex domains. The most famous domain in this context is given by the union of a rectangle and a disk, as illustrated in the figure. The Schwarz method requires an overlap between the subdomains. The disadvantage of


Fig. 1: Domaine original de Schwarz
this type of partitioning is the complexity of its numerical implementation, especially in the case of 3D problems. Moreover, it is difficult to define the overlapping regions for highly complex geometries. Additionally, the convergence in this case is very slow. Non-overlapping methods are the most commonly used nowadays and are the ones we will adopt in the following. They allow reducing the global problem to a problem on the interface.

The rest of the work is outlined as follows: The first section presents the mathematical foundation of the non-overlapping domain decomposition method applied to an
elliptic problem on a domain $\Omega$ subdivided into two subdomains $\Omega_{1}$ and $\Omega_{2}$ separated by an interface $\Gamma$. The variational formulation is given for both the global problem and the subdomain problems. The finite element discretization is also presented, and the $L U$ factorization is used to obtain a condensed system at the interface called the Schur complement. The algebraic version of the Schur complement, called the SteklovPoincaré operator, is also applied to obtain the same condensed system at the interface. Two iterative algorithms, Dirichlet-Neumann and Neumann-Neumann, are presented for solving the system at the interface. It is shown that these algorithms are preconditioned Richardson methods.

In the second section, some numerical results are presented concerning the convergence and error of domain decomposition methods.

The last section is devoted to the mathematical study of the coupling of a fluid flow governed by the Stokes equation and a flow in porous media governed by the Darcy law. The mathematical model for both the Stokes equation and the Darcy model, as well as the interface conditions between the two zones, are presented. A variational formulation is given, as well as the finite element discretization. Finally, a reduction of the problem to a condensed problem at the interface is presented, which is the Schur complement for our coupled Stokes-Darcy problem.

## 2 Non-overlapping domain decomposition for the Poisson equation

The objective of this section is to study the mathematical foundations of non-overlapping domain decomposition methods. We mainly consider the elliptic boundary problem of Poisson.

Let $\Omega$ be a domain in $\mathbb{R}^{d}$ (where $d=2,3$ ) with a continuous and Lipschitz boundary $\partial \Omega$. Suppose that $\Omega$ is divided into two subdomains $\Omega_{1}$ and $\Omega_{2}$ with an interface $\Gamma$, such that $\bar{\Omega}=\overline{\Omega_{1} \cup \Omega_{2}}, \Omega_{1} \cap \Omega_{2}=\varnothing$, and $\Gamma=\partial \Omega_{1} \cap \partial \Omega_{2}$. Figure 2 illustrates a rectangular domain $\Omega$ divided into two subdomains.


Fig. 2: Décomposition en deux sous-domaines

For a heterogeneous medium, we consider the following elliptic problem:

$$
\begin{cases}-\nabla(\mu \nabla u)=f, & \text { dans } \Omega,  \tag{1}\\ u=0, & \text { sur } \partial \Omega\end{cases}
$$

such that $\mu \in \mathbb{L}^{\infty}(\Omega), \inf _{\Omega} \mu>0$ et $f \in \mathbb{L}^{2}(\Omega)$.
The function $\mu$ is given by :

$$
\mu(x)=\left\{\begin{array}{l}
\mu_{1}(x), \text { dans } \Omega_{1}, \\
\mu_{2}(x), \text { dans } \Omega_{2},
\end{array}\right.
$$

with $\mu_{i} \in \mathbb{L}^{\infty}\left(\Omega_{i}\right) \mathrm{i}=1,2$.
To simplify, the theoretical results will be presented for the case $\mu=1$, which corresponds to the following Poisson problem:

$$
\begin{cases}-\Delta u=f, & \text { dans } \Omega,  \tag{2}\\ u=0, & \text { sur } \partial \Omega\end{cases}
$$

The results obtained for the Poisson problem (2) remain valid for the heterogeneous problem (1) and vice versa.

We consider the following local problems:

$$
\begin{align*}
& \begin{cases}-\Delta u_{1}=f_{1}, & \text { dans } \Omega_{1}, \\
u_{1}=0, & \text { sur } \partial \Omega_{1} \backslash \Gamma\end{cases}  \tag{3}\\
& \begin{cases}-\Delta u_{2}=f_{2}, & \text { dans } \Omega_{2}, \\
u_{2}=0, & \text { sur } \partial \Omega_{2} \backslash \Gamma\end{cases} \tag{4}
\end{align*}
$$

où $f_{1}=f_{\mid \Omega_{1}}$ et $f_{2}=f_{\mid \Omega_{2}}$
with the following boundary conditions:

- Condition of admissibility in the interface

$$
\begin{equation*}
u_{1}=u_{2} \quad \text { sur } \quad \Gamma . \tag{5}
\end{equation*}
$$

- Equilibrium condition

$$
\begin{equation*}
\frac{\partial u_{1}}{\partial n_{1}}=-\frac{\partial u_{2}}{\partial n_{2}} \quad \text { sur } \quad \Gamma . \tag{6}
\end{equation*}
$$

with $u_{i}=u_{\Omega_{i}}, i=1,2$ is the restriction of u on $\Omega_{i}$ et $n_{i}$ is the outward normal of the domain $\Omega_{i}$. The domain decomposition methods consist of reformulating the problem under study into an equivalent problem whose unknowns are functions defined on the interface of the subdomains. The Schur complement methods (primal and dual) will be studied for the problem (2)-(6). The Steklov-Poincaré operator, which is an algebraic version of the Schur complement, will also be presented.

## 3 Weak formulation

In this section, we provide the weak form of the global problem and local problems. We then show that these problems are well-posed, have unique solutions, and that solving the global problem is equivalent to solving the two local problems with the two interface conditions. For the variational formulation of the global problem, we consider the following Hilbert spaces: $H^{1}(\Omega):=\left\{\mathrm{v} \in \mathbb{L}^{2}(\Omega) / D_{j} \mathrm{v} \in \mathbb{L}^{2}(\Omega), j=0, \ldots d\right\}$ et $H_{0}^{1}(\Omega):=\left\{v \in H^{1}(\Omega) /\left.v\right|_{\partial \Omega}=0\right\}$. The space $H^{1}(\Omega)$ has the following norm $\|v\|_{H^{1}}:=\left(\|v\|_{\mathbb{L}^{2}}^{2}+\|\nabla v\|_{\mathbb{L}^{2}}^{2}\right)^{\frac{1}{2}}$ with $\|v\|_{\mathbb{L}^{2}}:=(v, v)^{\frac{1}{2}}$ et $(u, v)=\int_{\Omega} u v d x$.

Let us denote: $V:=H_{0}^{1}(\Omega)$. Let $u$ be a fairly regular solution of (2) and $v \in V$. Multiplying the first equation of (2) by $v$ and integrating over $\Omega$ we obtain:

$$
\text { int }_{\Omega}-(\Delta u) v d x=\int_{\Omega} f v d x
$$

By the formula of Green we will have then,

$$
\text { int }_{\Omega} \nabla u . \nabla v d x-\int_{\partial \Omega} \partial_{n} u v d \sigma=\int_{\Omega} f v d x .
$$

Now $v=0$ on the edge, the term $i n t_{\partial \Omega} \partial_{n} u v d \sigma$ cancels out, so we get the following weak form:

$$
\begin{equation*}
\text { trouver } u \in V \quad a(u, v)=(f, v) \quad v \in V \tag{7}
\end{equation*}
$$

with $(f, v)=\int_{\Omega} f v d x$ et $a(u, v):=\int_{\Omega} \nabla u \cdot \nabla v$
Theorem 3.1. The problem (7) admits a unique solution.
The proof can be done through the Lax-Milligram theorem and the Poincaré inequality (For more details, we can refer to [5]).

### 3.1 Variational formulation of local problems : Trace operator and extension operator

The trace space of $H^{1}(\Omega)$ on the edge $\partial \Omega$ is denoted $H^{1 / 2}(\partial \Omega)$. For a non empty open $\Sigma \subset \partial \Omega$ the trace space will be denoted $H^{1 / 2}(\Sigma)$. The trace operator

$$
\begin{gathered}
\gamma_{0}: H^{1}(\Omega) \longrightarrow H^{1 / 2}(\partial \Omega) \\
v \longmapsto \gamma_{0}(v)=v_{\left.\right|_{\text {partial } \Omega}}
\end{gathered}
$$

is surjective and continuous and we have the following trace inequality: There exists a constant $C_{\Omega}>0$ such that

$$
\left\|v_{\mid \partial \Omega}\right\|_{H^{1 / 2}(\partial \Omega)} \leq C_{\Omega}\|v\|_{H^{1}(\Omega)} \quad \forall v \in H^{1}(\Omega)
$$

Let us consider, $\Lambda:=\left\{\eta \in H^{1 / 2}(\Gamma) / \eta=v_{\left.\right|_{\Gamma}}\right.$ pour $\left.\quad v \in V\right\}$. The space $\Lambda$ has the norm $\|\eta\|_{\Lambda}:=\inf _{v \in V_{i}, v_{\mid \Gamma}=\eta}\|v\|_{\mathbb{H}^{1}\left(\Omega_{i}\right)}$. Let the follosing Hilbert spaces $V_{i}:=\left\{v_{i} \in\right.$ $\left.H^{1}\left(\Omega_{i}\right) /\left.v_{i}\right|_{\partial \Omega \cap \partial \Omega_{i}}=0\right\} \mathrm{i}=1,2$. The trace operator

$$
\gamma_{i}: V_{i} \longrightarrow \Lambda
$$

is continuous with respect to the norm parallel. $\|_{\Lambda}$, hence the trace inequality follows:

$$
\exists C_{i}^{*} \quad \text { telle que: }\left\|\left.v_{i}\right|_{\Gamma}\right\|_{\Lambda} \leq C_{i}^{*}\left\|v_{i}\right\|_{\mathbb{H}^{1}\left(\Omega_{i}\right)} \quad v_{i} \in V_{i}
$$

We note by $\Re_{i}(i=1,2)$ the following extension operator:

$$
\begin{aligned}
\Re_{i}: \Lambda & \longrightarrow V_{i} \\
\eta & \longmapsto \Re_{i} \eta
\end{aligned}
$$

with $\left(\Re_{i} \eta\right)_{\left.\right|_{\Gamma}}=\eta$. This operator is continuous and is not unique.
Equivalence between the global problem and the local problems Let us consider: $a_{i}\left(u_{i}, v_{i}\right):=\left(\nabla u_{i}, \nabla v_{i}\right)_{\Omega_{i}} \mathrm{i}=1,2$, and $V_{i}^{0}:=\left\{v_{i} \in V_{i} /\left.v_{i}\right|_{\Gamma}=0\right\}=H_{0}^{1}\left(\Omega_{i}\right) \mathrm{i}=1,2$. The following lemma shows the equivalence between the weak form of the global problem and the local problems.
Lemma 3.2. The fish problem (2) is equivalent to:
Find $u_{1}=u_{\Omega_{1}}$ et $u_{2}=u_{\Omega_{2}}$, such that :

$$
(F V P) \begin{cases}a_{1}\left(u_{1}, v_{1}\right)=\left(f_{1}, v_{1}\right), & v_{1} \in V_{1}^{0}, \\ u_{1}=u_{2}, & \text { sur } \Gamma \\ a_{2}\left(u_{2}, v_{2}\right)=\left(f_{2}, v_{2}\right), & v_{2} \in V_{1}^{0} \\ a_{2}\left(u_{2}, \Re_{2} \eta\right)=\left(f_{2}, \Re_{2} \eta\right)_{\Omega_{2}}+\left(f_{1}, \Re_{1} \eta\right)_{\Omega_{1}}-a_{1}\left(u_{1}, \Re_{1} \eta\right), & \forall \eta \in \Lambda\end{cases}
$$

Proof. Let $u$ be the solution of (7). We have $u_{i}=u \mid \Omega_{i} \in V_{i}$ and the first three equations of (FVP) are satisfied. Moreover, for each $\eta \in \Lambda$, we can define the extension operator $\Re$ by:

$$
\begin{gathered}
\Re: \Lambda \longrightarrow V \\
\eta \longmapsto \Re \eta
\end{gathered}
$$

with

$$
\Re \eta:=\left\{\begin{array}{ll}
\Re_{1} \eta, \text { sur } & \Omega_{1} \\
\Re_{2} \eta, & \text { sur }
\end{array} \Omega_{2}\right.
$$

Thus we have $a(u, \Re \eta)=(f, \Re \eta)$, so $\sum i=1^{2} a_{i}\left(u_{i}, \Re_{i} \eta\right)=\sum_{i=1}^{2}\left(f_{i}, \Re_{i} \eta\right) \Omega$, and therefore the last equation is satisfied.
Conversely, if $u_{1}$ and $u_{2}$ are solutions of (FVP), let:

$$
u=\left\{\begin{array}{l}
u_{1}, \text { sur } \Omega_{1}, \\
u_{2}, \text { sur } \Omega_{2}
\end{array}\right.
$$

It is clear that $u \in V$. For each $v \in V$, we have $\eta:=v \mid \Gamma \in \Lambda$, and by definition of $\Re \eta$, we have $\left(v \mid \Omega_{i}-\Re{ }_{i} \eta\right) \in V_{i}^{0}$. Using equations 1,3 , and 4 of (FVP), we have:

$$
\begin{aligned}
a(u, v) & =\sum_{i=1}^{2}\left[a_{i}\left(u_{i},\left.v\right|_{\Omega_{i}}-\Re_{i} \eta\right)+a_{i}\left(u_{i}, \Re_{i} \eta\right)\right] \\
& =\sum_{i=1}^{2}\left[\left(f,\left.v\right|_{\Omega_{i}}-\Re_{i} \eta\right)_{\Omega_{i}}+\left(f, \Re_{i} \eta\right)_{\Omega_{i}}\right] \\
& =(f, v) \quad \forall v \in V .
\end{aligned}
$$

Therefore, $u$ is a solution of (7).

## 4 Discretization by Finite Element Method

We will give a finite element discretization of the (FVP) problem. For this, we consider a regular triangulation $\mathcal{T} h$ of the domain $\bar{\Omega}: \bar{\Omega}=\bigcup K \in \mathcal{T}_{h} K$, where each $K \in \mathcal{T}_{h}$ is a non-empty interior triangle with $\dot{K}_{1} \cap \grave{K}_{2}=\emptyset$ for all $K_{1}, K_{2} \in \mathcal{T} h$, and if $e=K_{1} \cap K_{2} \neq$ $\emptyset$, then $e$ is a face, edge, or vertex shared by $K_{1}$ and $K_{2}$. We consider the space

$$
X_{h}(\Omega):=v_{h} \in C^{0}(\bar{\Omega}) / v_{h} \mid K \in \mathbb{P}_{1}(K) \quad \forall K \in \mathcal{T}_{h},
$$

where $\mathbb{P} 1(K)$ is the space of polynomials on the element $K$ of degree less than or equal to 1 , and the space

$$
V_{h}=v_{h} \in X_{h}(\Omega) / \quad v_{h} \mid \partial \Omega=0=X_{h}(\Omega) \cap \mathbb{H}_{0}^{1}(\Omega),
$$

where $\mathbb{H}_{0}^{1}(\Omega)$ is the Sobolev space of functions in $\mathbb{H}^{1}(\Omega)$ with zero trace on the boundary $\partial \Omega$.

We denote $F(v)=(f, v)$, and therefore the problem (7) becomes:

$$
\begin{equation*}
\text { trouver } \quad u_{h} \in V_{h}: a\left(u_{h}, v_{h}\right)=F(v) \tag{8}
\end{equation*}
$$

Let $\varphi_{j}{ }_{j=1}^{N}$ be a basis for the space $V_{h}$. Suppose that the interface $\Gamma=\bar{\Omega}_{1} \cap \bar{\Omega}_{2}$ is a union of edges or faces of the triangulation $\mathcal{T} h$. We introduce the following partition of the nodes in the domain: let $x_{j}^{1}, 1 \leq j \leq N_{1}$ be the nodes in the subdomain $\Omega_{1}$, let $x_{j}^{2}, 1 \leq j \leq N_{2}$ be the nodes in the subdomain $\Omega_{2}$, and let $x_{j}^{\Gamma}, 1 \leq j \leq N \Gamma$ be the nodes on the interface $\Gamma$. We also partition the basis functions by denoting $\varphi_{j}^{i}$ the functions associated with nodes $x_{j}^{i} \quad\left(i=1,2 \quad j=1, \ldots N_{i}\right)$ and $\varphi_{j}^{\Gamma}$ the functions associated with nodes $x_{j}^{\Gamma}$ on the interface.

We take the $\varphi_{j}^{\alpha}(\alpha=1,2$ or $\Gamma)$ as test functions in the problem (8), and thus we have the following problem:

Find $u_{h} \in V_{h}$ such that:

$$
\left\{\begin{array}{l}
a\left(u_{h}, \varphi_{j}^{1}\right)=F\left(\varphi_{j}^{1}\right), \forall j=1 \ldots, N_{1}  \tag{9}\\
a\left(u_{h}, \varphi_{j}^{2}\right)=F\left(\varphi_{j}^{2}\right), \forall j=1 \ldots, N_{2} \\
a\left(u_{h}, \varphi_{j}^{\Gamma}\right)=F\left(\varphi_{j}^{\Gamma}\right), \forall j=1 \ldots, N_{\Gamma}
\end{array}\right.
$$

Let $a_{i}(.,$.$) and F_{i}($.$) be the restrictions of the form a(.,$.$) and F($.$) to the subdomain$ $\Omega_{i}, \mathrm{i}=1,2$. We consider the space $V_{i, h}$ of polynomial functions of degree less than or equal to 1 , which vanish on the boundary $\partial \Omega_{i}$ except for the interface $\Gamma$, and the space $V_{i, h}^{0}:=v_{h} \in V_{i, h} / v_{h} \mid \Gamma=0$. Let $u_{h}^{i}=u_{h} \mid \Omega_{i} \in V_{i, h}$, then the problem (9) can be written in the following multi-domain formulation:

$$
\begin{cases}a_{1}\left(u_{h}^{1}, \varphi_{k}^{1}\right)=F_{1}\left(\varphi_{k}^{1}\right), & \forall k=1 \ldots, N_{1},  \tag{10}\\ a_{2}\left(u_{h}^{2}, \varphi_{k}^{2}\right)=F_{2}\left(\varphi_{k}^{2}\right), & \forall k=1 \ldots, N_{2}, \\ a_{1}\left(u_{h}^{1},\left.\varphi_{k}^{\Gamma}\right|_{\Omega_{1}}\right)+a_{2}\left(u_{h}^{2},\left.\varphi_{k}^{\Gamma}\right|_{\Omega_{2}}\right)=F_{1}\left(\left.\varphi_{k}^{\Gamma}\right|_{\Omega_{1}}\right)+F_{2}\left(\left.\varphi_{k}^{\Gamma}\right|_{\Omega_{2}}\right), & \forall k=1 \ldots, N_{\Gamma},\end{cases}
$$

We decompose the function $u_{h}$ on the basis $\left\{\operatorname{varph} i_{j}\right\}$ of the space $V_{h}$, we have :

$$
\begin{equation*}
u_{h}(x)=\sum_{j=1}^{N_{1}}\left(u_{j}^{1} \varphi_{j}^{1}(x)\right)+\sum_{j=1}^{N_{2}}\left(u_{j}^{2} \varphi_{j}^{2}(x)\right)+\sum_{j=1}^{N_{\Gamma}}\left(u_{j}^{\Gamma} \varphi_{j}^{\Gamma}(x)\right) . \tag{11}
\end{equation*}
$$

where $u_{j}^{\alpha}:=u_{h}\left(x_{j}^{\alpha}\right)\left(j=1, \ldots, N_{\alpha}, \alpha=1,2, \Gamma\right)$ are the coefficients of the linear combination representing $u_{h}$ in the $\left\{\varphi_{j}\right\}$ basis. In the same way, we also obtain:

$$
\begin{equation*}
u_{h}^{i}(x)=\sum_{j=1}^{N_{i}}\left(u_{j}^{i} \varphi_{j}^{i}(x)\right)+\sum_{j=1}^{N_{\Gamma}}\left(u_{j}^{\Gamma} \varphi_{\left.j\right|_{\Omega_{i}}}^{\Gamma}(x)\right) . \tag{12}
\end{equation*}
$$

where $\varphi_{\left.j\right|_{\Omega_{i}}}^{\Gamma}$ is the restriction of $\varphi_{j}^{\Gamma}$ on $\Omega_{i}$.
By substituting (12) in the first two equations of (10), we find

$$
\begin{equation*}
\sum_{j=1}^{N_{1}} u_{j}^{1} a_{1}\left(\varphi_{j}^{1}(x), \varphi_{k}^{1}(x)\right)+\sum_{j=1}^{N_{\Gamma}} u_{j}^{\Gamma} a_{1}\left(\varphi_{j \mid \Omega_{1}}^{\Gamma}(x), \varphi_{k}^{1}(x)\right)=F\left(\varphi_{k}^{1}(x)\right) \tag{13}
\end{equation*}
$$

and

$$
\begin{equation*}
\sum_{j=1}^{N_{2}} u_{j}^{1} a_{2}\left(\varphi_{j}^{2}(x), \varphi_{k}^{2}(x)\right)+\sum_{j=1}^{N_{\Gamma}} u_{j}^{\Gamma} a_{2}\left(\varphi_{j \mid \Omega_{2}}^{\Gamma}, \varphi_{k}^{2}(x)\right)=F\left(\varphi_{k}^{2}(x)\right) \tag{14}
\end{equation*}
$$

Let define the following matrix

$$
\begin{gathered}
\left(A_{i i}\right)_{k j}=a_{i}\left(\varphi_{j}^{i}, \varphi_{k}^{i}\right), \quad i=1,2 \\
\left(A_{i \Gamma}\right)_{k j}=a_{i}\left(\varphi_{j \mid \Omega_{i}}^{\Gamma}, \varphi_{k}^{i}\right), \\
\left(A_{\Gamma i}\right)_{k j}=a_{i}\left(\varphi_{j}^{i}, \varphi_{k \mid \Omega_{i}}^{\Gamma}\right),
\end{gathered} \quad i=1,22
$$

$$
\begin{gathered}
\left(A_{\Gamma \Gamma}^{i}\right)_{k j}=a_{i}\left(\varphi_{j \mid \Omega_{i}}^{\Gamma}, \varphi_{k \mid \Omega_{i}}^{\Gamma}\right), \quad i=1,2 \\
\left(b_{i}\right)_{k}=F_{i}\left(\varphi_{k}^{i}\right), \quad i=1,2 \\
\left(b_{i}^{\Gamma}\right)_{k}=F_{i}\left(\varphi_{\left.k\right|_{\Omega_{i}}}^{\Gamma}\right), \quad i=1,2
\end{gathered}
$$

The system (10) can therefore be written as :

$$
\left\{\begin{array}{l}
A_{11} u_{1}+A_{1 \Gamma} u_{\Gamma}=b_{1},  \tag{15}\\
A_{22} u_{2}+A_{2 \Gamma} u_{\Gamma}=b_{2}, \\
A_{\Gamma 1} u_{1}+A_{\Gamma 2} u_{2}+\left(A_{\Gamma \Gamma}^{1}+A_{\Gamma \Gamma}^{2}\right) u_{\Gamma}=b_{1}^{\Gamma}+b_{2}^{\Gamma},
\end{array}\right.
$$

où $u_{1}=\left(u_{j}^{1}\right)_{j=1 \ldots N_{1}}, u_{2}=\left(u_{j}^{2}\right)_{j=1 \ldots N_{2}}, u_{\Gamma}=\left(u_{j}^{\Gamma}\right)_{j=1 \ldots N_{\Gamma}}, b_{1}=\left(b_{j}\right)_{j=1 \ldots N_{1}}, b_{2}=$ $\left(b_{j}\right)_{j=1 \ldots N_{2}}, b_{1}^{\Gamma}=\left(b_{1 j}^{\Gamma}\right)_{j=1 \ldots N_{1}}, b_{2}^{\Gamma}=\left(b_{2 j}^{\Gamma}\right)_{j=1 \ldots N_{2}}$ Then we have the system in the following matrix form:

$$
\left(\begin{array}{ccc}
A_{11} & 0 & A_{1 \Gamma}  \tag{16}\\
0 & A_{22} & A_{2 \Gamma} \\
A_{\Gamma 1} & A_{\Gamma 2} & A_{\Gamma \Gamma}
\end{array}\right)\left(\begin{array}{l}
u_{1} \\
u_{2} \\
u_{\Gamma}
\end{array}\right)=\left(\begin{array}{c}
b_{1} \\
b_{2} \\
b_{\Gamma}
\end{array}\right)
$$

with $A_{\Gamma \Gamma}=A_{\Gamma \Gamma}^{1}+A_{\Gamma \Gamma}^{2} \quad$ et $\quad b_{\Gamma}=b_{\Gamma}^{1}+b_{\Gamma}^{2}$.

## 5 Method of primal Schur's complement

The system (16) can be written in the following form:

$$
\left\{\begin{array}{l}
A_{i i} u_{i}+A_{i \Gamma} u_{\Gamma}=b_{i}, i=1,2  \tag{17}\\
A_{\Gamma 1} u_{1}+A_{\Gamma 2} u_{2}+\left(A_{\Gamma \Gamma}^{1}+A_{\Gamma \Gamma}^{2}\right) u_{\Gamma}=b_{\Gamma},
\end{array}\right.
$$

It is assumed that the blocks $A_{i i}$ admit a factorization $L U$, the first two equations of the system (17) then becomes :

$$
\begin{equation*}
u_{i}=A_{i i}^{-1}\left(b_{i}-A_{i \Gamma} u_{\Gamma}\right) \quad i=1,2 \tag{18}
\end{equation*}
$$

We replace $u_{i}, i=1,2$ in the third equation, so we have the condensed system at the interface

$$
\begin{equation*}
S_{\Gamma \Gamma} u_{\Gamma}=C_{\Gamma} \tag{19}
\end{equation*}
$$

with $S_{\Gamma \Gamma}=A_{\Gamma \Gamma}-A_{\Gamma 1} A_{11}^{-1} A_{1 \Gamma}-A_{\Gamma 2} A_{22}^{-1} A_{2 \Gamma}$ Schur's complement to the unknowns on $\Gamma$, and $C_{\Gamma}=\left(b_{1}^{(1)}-A_{\Gamma 1} A_{11}^{-1} b_{1}\right)+\left(b_{\Gamma}^{(2)}-A_{\Gamma 2} A_{22}^{-1} b_{2}\right)=C_{\Gamma}^{(1)}+C_{\Gamma}^{(2)}$.

The solution of the system (19) can be done by an iterative method. The corresponding $S_{\Gamma}$ matrix has a better conditioning compared to the starting matrix $A$, and thus allows a better convergence speed (see [1] page 98).

## 6 Steklov-Poincaré operator


V. A. Steklov $1864-1926$

H. Poincaré $1854-1912$

Let $\eta:=u_{\left.\right|_{\Gamma}}$ be the restriction of the solution $u$ of (2) on $\Gamma$. Consider the following Dirichlet problem:

$$
\begin{cases}-\Delta\left(w_{i}\right)=f, & \text { dans } \Omega_{i},  \tag{20}\\ w_{i}=0, & \text { sur }\left.\partial \Omega_{i}\right|_{\Gamma}, \quad i=1,2 \\ w_{i}=\eta, & \text { sur } \Gamma\end{cases}
$$

The solution $w_{i}$ of (20) can be written as

$$
\begin{equation*}
w_{i}=u_{i}^{0}+u_{i}^{*} . \tag{21}
\end{equation*}
$$

The solution $w_{i}$ of (20) can be written in the following form where $u_{i}^{0}$ and $u_{i}^{*}$ are the solutions of the following Dirichlet problems:

$$
\begin{cases}-\Delta\left(u_{i}^{0}\right)=0, & \text { dans } \Omega_{i},  \tag{22}\\ u_{i}^{0}=0, & \text { sur }\left.\partial \Omega_{i}\right|_{\Gamma}, \quad i=1,2 \\ u_{i}^{0}=\eta, & \text { sur } \Gamma\end{cases}
$$

and

$$
\begin{cases}-\Delta\left(u_{i}^{*}\right)=f, \text { dans } \Omega_{i},  \tag{23}\\ u_{i}^{0}=0, & \text { sur }\left.\partial \Omega_{i}\right|_{\Gamma}, \quad i=1,2 \\ u_{i}^{0}=0, & \text { sur } \Gamma\end{cases}
$$

The solution $u_{i}^{0}$ of (22) is called the harmonic extension of $\eta$ to $\Omega_{i}$ and will be denoted $H_{i} \eta$, while the solution $u_{i}^{*}$ of (23) will be denoted $G_{i} f$ and will be called the resolvent operator. We have then

$$
\begin{equation*}
w_{i}=H_{i} \eta+G_{i} f \tag{24}
\end{equation*}
$$

Remark 6.1. $G_{i} f$ depends only on $f$, while $H_{i} \eta$ depends only on the value of the solution $u$ on $\Gamma$.

The solution $w_{i}$ of (24) is equal to $u_{i}$ if and only if we have the condition $f r a c \partial w_{1} \partial n=$ $\frac{\partial w_{2}}{\partial n}$ on $\Gamma$, where $n$ is the normal on $\Gamma\left(n=n_{1}=-n_{2}\right)$. And since we have $\frac{\partial u_{1}}{\partial n}=$ $\frac{\partial u_{2}}{\partial n}$, then we will have $\left(\frac{\partial H_{1} \eta+\partial G_{1} f}{\partial n}\right)=\left(\frac{\partial H_{2} \eta+\partial G_{2} f}{\partial n}\right)$ therefore

$$
\begin{equation*}
\left(\frac{\partial H_{1}-\partial H_{2}}{\partial n}\right) \eta=\left(\frac{\partial G_{2} f-\partial G_{1} f}{\partial n}\right) \tag{25}
\end{equation*}
$$

Let $\chi$ be the function defined as follows:

$$
\begin{aligned}
\chi & =\frac{\partial G_{2} f}{\partial n}-\frac{\partial G_{1} f}{\partial n} \\
& =-\frac{\partial G_{2} f}{\partial n_{2}}-\frac{\partial G_{1} f}{\partial n_{1}} \\
& =-\sum_{i=1}^{2}\left(\frac{\partial G_{i} f}{\partial n_{i}}\right)
\end{aligned}
$$

This function depends only on f and the normal on $\Gamma$. Soit $\mathbb{S}$ l'opérateur de SteklovPoincaré définit par :

$$
\begin{aligned}
\mathbb{S} \eta & =\frac{\partial H_{1} \eta}{\partial n_{1}}-\frac{\partial H_{2} \eta}{\partial n_{2}} \\
& =\sum_{i=1}^{2}\left(\frac{\partial H_{i} \eta}{\partial n_{i}}\right) \\
& =\mathbb{S}_{1}+\mathbb{S}_{2}
\end{aligned}
$$

The equation (25) then becomes the Steklov-Poincaré equation on the following $\Gamma$ interface:

$$
\begin{equation*}
\mathbb{S} \eta=\chi \quad \text { sur } \quad \Gamma . \tag{26}
\end{equation*}
$$

## 7 Iterative methods for domain decomposition: Dirichlet-Neumann algorithm

The Dirichelt-Neumann method consists in solving a Dirichlet problem in the first domain, with a Dirichlet data $\lambda^{k}$ on $\Gamma$, and then solving a mixed Dirichlet-Neumann problem on the second domain where we use the value of the flux of the previous solution on $\Gamma$ as a Neumann condition, and a homogeneous condition on the remainder of the domain edge. The algorithm for the heterogeneous problem (1) is given as follows: • $\lambda^{0}$ given

- for each iteration $k \succeq 0$
- solve

$$
\begin{cases}-\nabla\left(\mu_{1} \nabla u_{1}^{k+1}\right)=f, \text { dans } \Omega_{1},  \tag{27}\\ u_{1}^{k+1}=0, & \text { sur }\left.\partial \Omega_{1}\right|_{\Gamma}, \\ u_{1}^{k+1}=\lambda^{k}, & \text { sur } \Gamma\end{cases}
$$

- solve :

$$
\begin{cases}-\nabla\left(\mu_{2} \nabla u_{2}^{k+1}\right)=f, & \text { dans } \Omega_{2},  \tag{28}\\ u_{2}^{k+1}=0, & \text { sur }\left.\partial \Omega_{2}\right|_{\Gamma}, \\ \mu_{2} \frac{\partial u_{2}^{k+1}}{\partial n}=\mu_{1} \frac{\partial u_{1}^{k+1}}{\partial n}, & \text { sur } \Gamma\end{cases}
$$

- update $\lambda^{k+1}$ :

$$
\begin{equation*}
\lambda^{k+1}=\theta u_{2}^{k+1}+(1-\theta) \lambda^{k} . \tag{29}
\end{equation*}
$$

où $\theta$ is a relaxation parameter to accelerate the convergence.

- Stop criterion:
$\left\|\lambda^{k+1}-\lambda^{k}\right\| \leq \epsilon$, avec $\epsilon$ une tolérance donnée.
The following theorem shows that the Dirichlet-Neumann method is nothing else than the preconditioned Richardson method.

Theorem 7.1. The Dirichlet-Neumann method is equivalent to the preconditioned Richardson method with preconditioner $S_{2}$ applied to the Steklov-Poincaré interface equation $S \lambda=\chi$ i.e. we have :

$$
\lambda^{k+1}=\lambda^{k}+\theta S_{2}^{-1}\left(\chi-S \lambda^{k}\right)
$$

### 7.1 Neumann-Neumann Algorithm

For Neumann-Neumann we first solve a Dirichlet problem in each subdomain with Dirichlet data $\lambda^{k}$ on $\Gamma$, then two Neumann problems where we use the difference of the fluxes of the previous solutions as Neumann data on $\Gamma$. We have the following algorithm to solve the heterogeneous problem (1): • $\lambda^{0}$ donnée sur $\Gamma$.

- for each $k \succeq 0$,
- Pour $\mathrm{i}=1,2$ solve :

$$
\begin{cases}-\nabla\left(\mu_{i} \nabla u_{i}^{k+1}\right)=f, & \text { dans } \Omega_{i},  \tag{30}\\ u_{i}^{k+1}=0, & \text { sur }\left.\partial \Omega_{i}\right|_{\Gamma}, \\ u_{i}^{k+1}=\lambda^{k}, & \text { sur } \Gamma\end{cases}
$$

- For $\mathrm{i}=1,2$ solve :

$$
\begin{cases}-\nabla\left(\mu_{i} \nabla \psi_{i}^{k+1}\right)=0, & \text { dans } \Omega_{i},  \tag{31}\\ \psi_{i}^{k+1}=0, & \text { sur }\left.\partial \Omega_{i}\right|_{\Gamma}, \\ \mu_{i} \frac{\partial \psi_{i}^{k+1}}{\partial n}=\mu_{1} \frac{\partial u_{1}^{k+1}}{\partial n}-\mu_{2} \frac{\partial u_{2}^{k+1}}{\partial n}, & \text { sur } \Gamma\end{cases}
$$

- mettre à jour $\lambda^{k+1}$ :

$$
\begin{equation*}
\lambda^{k+1}=\lambda^{k}-\theta\left(\sigma_{1} \psi_{1 \mid \Gamma}^{k+1}-\sigma_{2} \psi_{2 \mid \Gamma}^{k+1}\right) . \tag{32}
\end{equation*}
$$

- Stop criterion :
$\left\|\lambda^{k+1}-\lambda^{k}\right\| \leq \epsilon$, avec $\epsilon$ une tolérance donnée.

The parameter $\theta$ is a relaxation parameter to speed up the convergence. The positive weights $\sigma_{1}$ and $\sigma_{2}$ allow to weight the contributions of the flows following the $\Omega_{1}$ and $\Omega_{2}$ domains to attenuate the effects of the jumps of the diffusion functions $\mu_{1}$ and $\mu_{2}$ for example. In general we will take $\sigma_{1}+\sigma_{2}=1$ and in the case of constant diffusion coefficients by subdomain, we take $\sigma_{i}=\frac{\mu_{i}}{\mu_{1}+\mu_{2}}$.
Theorem 7.2. The Neumann-Neumann method is equivalent to the preconditioned Richardson method with preconditioner $N:=\left(\sigma_{1} S_{1}^{-1}+\sigma_{2} S_{2}^{-1}\right)^{-1}$ applied to the SteklovPoincaré interface equation $S \lambda=\chi$ c'est à dire qu'on $a$ :

$$
\lambda^{k+1}=\lambda^{k}+\theta\left(\sigma_{1} S_{1}^{-1}+\sigma_{2} S_{2}^{-1}\right)\left(\chi-S \lambda^{k}\right)
$$

## 8 Numerical results: Diffusion problem

FreeFem++ is a free software originally developed by Frédéric Hecht, researcher at the Jacques-Louis Lions Laboratory of the Pierre et Marie Curie University in Paris. It allows to numerically solve Partial Differential Equations (PDE) by the Finite Element Method (FEM). The FreeFEM + + language allows to quickly specify ( $2 D, 3 D$ ) PDEs, to manipulate several $(2 D, 3 D)$ meshes and to write scripts in $\mathrm{C}++$ to define numerical algorithms for problems: nonlinear, coupled...

FreeFEM + + allows to create meshes of the studied domains. Figures (3) and (4) show examples of meshes (uniform and non-uniform) created by FreeFEM.


Fig. 3: Maillage uniforme sous FreeFEM


Fig. 4: Maillage non-uniforme sous FreeFEM

We consider the following domain $\Omega=[0,1] \times[0,1]$, we fix the number of nodes on each interface of this domain to 5 nodes, then we generate a uniform mesh on the
whole domain. The problem to solve on all $\Omega$ is the following:

$$
\begin{cases}-\Delta u=1, & \text { dans } \Omega  \tag{33}\\ u=0, & \text { sur } \partial \Omega\end{cases}
$$

The graph (5) representing the solution on the whole domain $\Omega$.


Fig. 5: Solution sur tout le domaine $\Omega$.

We decompose the domain $\Omega$ in two subdomains $\Omega_{1}$ and $\Omega_{2}$, we solve the problem (33) on each subdomain using one of the Dirichlet-Neumann or Neumann-Neumann algorithms, the figure (??) represents the approximate solution on $\Omega_{1}$, while the figure (??) represents that on $\Omega_{2}$.


Fig. 6: Solution sur $\Omega_{1}$


Fig. 7: Solution sur $\Omega_{2}$

Let $u_{1}$ be the solution on the first domain $\Omega_{1}$, and $u_{2}$ that on the second domain $\Omega_{2}$, we characterize the stopping criterion as follows:

- We set a tolerance $\epsilon$.
- We initialize the error by $0(\mathrm{ER} 0=0)$.
- At each iteration, the following relative error is calculated:

$$
\begin{equation*}
E R 1=\frac{\left\|u_{1}-u_{2}\right\|_{\mathbb{L}^{2}}^{2}}{\left\|u_{1}\right\|_{\mathbb{L}^{2}}^{2}} \tag{34}
\end{equation*}
$$

- Case 1:

Si $|E R 1-E R 0|>\epsilon$, we continue, and we reset the calculated error : $E R 0=E R 1$.

- Case 2:

Si $|E R 1-E R 0| \leq \epsilon$, we stop the loop.
For the Poisson equation, both Dirichlet-Neumann and Neumann-Neumann algorithms converge in one iteration.

Let $\Omega=[0,1] \times[0,1]$ and

$$
\begin{cases}-\Delta u=\sin (\pi x) \sin (\pi y), & \text { dans } \Omega,  \tag{35}\\ u=0, & \text { sur } \partial \Omega .\end{cases}
$$

The analytical solution of the problem (??) is of the form :

$$
\begin{equation*}
u(x, y)=\frac{1}{2 \pi^{2}} \sin (\pi x) \sin (\pi y) \tag{36}
\end{equation*}
$$

We are going to make a comparison between the values taken on the $\Gamma$ interface between the two domains $\Omega_{1}$ and $\Omega_{2}$ by this analytical solution, by the solution without domain decomposition and the solution with domain decomposition. We fix y $=0.5$ then we vary
x and we recover the values of $u(x, 0.5)$, of the solution without domain decomposition and those with domain decomposition.

## Error calculation:

The following graph illustrates the errors $\left\|u_{A D D M}-u_{\text {exact }}\right\|_{\mathbb{L}^{2}}$, and $\left\|u_{\text {SDDM }}-u_{\text {exact }}\right\|_{\mathbb{L}^{2}}$, where $u_{A D D M}$ represents the approximate solution obtained with domain decomposition, $u_{S D D M}$ the approximate solution recovered without domain decomposition, and $u_{\text {exact }}$ the exact solution given in (??). We always use the criterion (34) as a stopping criterion.


Fig. 8: Error comparison on the interface with the NN algorithm.

The green curve represents the computed error between the analytical solution and the approximated solution without domain decomposition, while the brown curve represents the error between the analytical solution and the approximated solution with domain decomposition. The two curves are not identical, because of the errors between the approximated solution without domain decomposition and the one with domain decomposition on the ( $y=0.5$ ) interface.

### 8.1 Influence of the jumps of the coefficients for a heterogeneous problem

We are interested in the problem:

$$
\left\{\begin{array}{l}
-\nabla(\mu \nabla u)=1, \text { dans } \Omega=[0,1] \times[0,1],  \tag{37}\\
u=0, \\
\text { sur } \partial \Omega
\end{array}\right.
$$

with

$$
\mu(x)= \begin{cases}\mu_{1}, \text { dans } & \Omega_{1},  \tag{38}\\ \mu_{2}, \text { dans } & \Omega_{2},\end{cases}
$$

For each value of ( $\mu_{1}, \mu_{2}$ ), we refine the mesh (we vary the value of n ), and we recover the number of iterations needed for the convergence using the stopping criterion (34).

Remark 8.1. The error computed above is done on the interface ( $y=0.5$ ), while the solution on the interface is recovered in a different way according to the choice of the algorithm.

The table (1) presents the results for the two algorithms Dirichlet-Neuamnn and Neumann-Neumann.

| $\left(\mu_{1}, \mu_{2}\right)$ | $(1,5)$ |  |  |  | $\left(1,10^{-4}\right)$ |  |  |  | $\left(10^{-4}, 10^{-5}\right)$ |  |  |  |
| :--- | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| n | 10 | 20 | 40 | 80 | 10 | 20 | 40 | 80 | 10 | 20 | 40 | 80 |
| Nbre d'itérations DN | 1 | 1 | 1 | 1 | 12 | 12 | 12 | 12 | 14 | 14 | 14 | 14 |
| Nbre d'itérations NN | 1 | 1 | 1 | 1 | 22 | 22 | 22 | 22 | 23 | 23 | 23 | 23 |

Table 1: Results for the Dirichlet-Neumann algorithm and for the Neumann-Neumann algorithm.

Remark 8.2. In the case of two domains $\Omega_{1}$ and $\Omega_{2}$ of the same size, we see that the number of iterations increases with the jump size of the coefficients.

Even though the Dirichlet-Neumann(DN) algorithm converges in a smaller number of iterations than the Neumann-Neumann(NN) algorithm, the error calculation $\mid u_{S D D M}-u_{A D D M} \|_{\text {mathbbL }^{2}}$ with the stopping criterion (34), shows that the NeumannNeumann algorithm gives a good approximation than Dirichlet-Neumann. The table (??) presents the results obtained for different meshes.

| Taille du maillage (n) | 10 | 20 | 40 | 80 |
| :--- | :---: | :---: | :---: | :---: |
| ERREUR-DN | 0,0439327 | 0,0251683 | 0,0130351 | 0,00657722 |
| ERREUR-NN | 0,000623169 | 0,000234582 | $8,43998 \mathrm{E}-005$ | $2,99788 \mathrm{E}-005$ |

Table 2: Error calculation for different meshes.

It is clear that the error of both algorithms improves while refining the mesh.

## 9 Conclusion

Domain decomposition methods are numerical methods well adapted to parallel solving of partial differential equations. They are not an alternative but an essential choice for solving large, coupled or geometrically complex problems. These methods are more efficient than an iterative method applied to the global problem, because they combine direct methods and an iterative method to solve an interface problem, better conditioned than the global problem.

The study made on this work has treated the domain decomposition methods without overlap. The mathematical foundations of domain decomposition methods applied to the Poisson equation were studied, then the study was extended to a heterogeneous problem. The two algorithms Dirichlet-Neumann and Neumann-Neumann were presented to solve these systems, and to illustrate the importance of these methods as preconditioning methods. The discretization of the problem has been done by finite elements, and the numerical implementation has been done on the FreeFem software.

Numerical results concerning the convergence speed and the error estimation are presented in the numerical part.

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