# DOMAIN DECOMPOSITION WITH NONMATCHING GRIDS FOR EXTERIOR TRANSMISSION PROBLEMS VIA FEM AND DTN MAPPING *1) 

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Dedicated to the 70th birthday of Professor Lin Qun


#### Abstract

In this paper, we are concerned with a non-overlapping domain decomposition method (DDM) for exterior transmission problems in the plane. Based on the natural boundary integral operator, we combine the DDM with a Dirichlet-to-Neumann (DtN) mapping and provide the numerical analysis with nonmatching grids. The weak continuity of the approximation solutions on the interface is imposed by a dual basis multiplier. We show that this multiplier space can generate optimal error estimate and obtain the corresponding rate of convergence. Finally, several numerical examples confirm the theoretical results.


Mathematics subject classification: 65N30, 65N55.
Key words: Domain decomposition, Nature boundary element, Nonmatching grids, Weak continuity, D-N alternating, Dual basis, Projection operator, Error estimate.

## 1. Introduction

Domain decomposition method (DDM) with nonmatching grids is a kind of nonconforming finite element methods. In the past few years, there is a fast growing interest in this field (see [1], [2], [5], [7] ). This kind of DDM allows different discretizations in different nonoverlapping subdomains by some Lagrange multiplier. This nonconforming element method also allows for local refinement in only certain subregions of the computational domain. Hence, it is suitable for parallel computing (see [6]).

The key point to deal with the nonmatching grids is how to choose the matching condition so that the resulting approximation problem possesses the optimal error estimate. The approximate solutions must satisfy some weak continuity such as the integration matching condition, whereas the pointwise matching.

In this paper, we propose a new class of multiplier space for the exterior unbounded problems with annular interfaces, which is based on the idea of dual basis multiplier (refer to [7]). We impose weak continuity conditions in the sense that the jump of the DDM solution across the interface is required to be orthogonal to a space of test functions. Due to the character of annular interface that there is no intersections between any of two subregions, it is easier for us to construct efficient and practical multiplier. The basis functions of the multiplier spaces

[^0]are generated by a set of simple functions with local compact supports. The resulting discrete system is still symmetric and positive definite. It will be shown such construction guarantee the optimal energy error estimate for the approximation solutions and the discrete formulation is easy to be solved.

The outline of the paper is as follows. In Section 2, we present the coupled variational formulation for the exterior transmission problem by the finite element method and the natural boundary element method (FEM-NBEM). Then we make a finite element discretization with nonmatching grids for this coupled system in Section 3, and the construction of the multiplier spaces is also introduced. It will be shown that the nonconforming approximation posseses the optimal energy error estimate. In Section 4, we give a D-N alternating method to solve the discrete system and show that this D-N algorithm is convergent and independent of the finite element meshes. Finally, in Section 5, we illustrate these theoretical results by using some numerical examples.

## 2. FEM-NBEM Coupling

As a model problem, we consider a second order elliptic equation in divergence form coupled with the Laplace equation in the exterior unbounded region. Let $\Omega_{0}$ be a bounded domain of $\mathbb{R}^{2}$ with a Lipschitz-continuous boundary $\Gamma_{0} . \Omega_{1}$ is the annular region bounded by $\Gamma_{0}$ and another smooth closed curve $\Gamma_{1}$ that is strictly contained in $\mathbb{R}^{2} \backslash \bar{\Omega}_{0}$ (see Figure 1 ). We denote by $\Omega_{c}$ the complement of $\bar{\Omega}_{0} \bigcup \bar{\Omega}_{1}$. Assume that $g \in H^{1 / 2}\left(\Gamma_{0}\right)$ and $f \in L^{2}\left(\Omega_{1}\right)$, then the exterior transmission problem reads as: find $u$ such that

$$
\begin{array}{r}
u_{1}=g, \quad \text { on } \Gamma_{0} \quad-\operatorname{div}\left(A \nabla u_{1}\right)=f \quad \text { in } \Omega_{1} \\
u_{1}=u_{c} \quad \text { and } \quad\left(A \nabla u_{1}\right) \cdot \boldsymbol{n}=\frac{\partial u_{c}}{\partial \boldsymbol{n}} \quad \text { on } \Gamma_{1} \\
-\Delta u_{c}=0, \quad \text { in } \Omega_{c} \quad u_{c}(x)=O(1) \quad \text { as }|x| \rightarrow \infty \tag{2.1c}
\end{array}
$$

where $\boldsymbol{n}=\left(n_{1}, n_{2}\right)^{T}$ denotes the unit outward normal to $\Gamma_{1}$ and $A$ is uniformly symmetric pos-


Figure 1: The domain of transmission problem
itive definite matrix with Lipschitz-continuous coefficients, that is to say, there exists constants
$\alpha_{1}$ and $\alpha_{2}$ such that

$$
\begin{equation*}
\alpha_{1}\|\boldsymbol{\eta}\|^{2} \leq(A \boldsymbol{\eta}) \cdot \boldsymbol{\eta} \leq \alpha_{2}\|\boldsymbol{\eta}\|^{2} \quad \forall \boldsymbol{\eta} \in \mathbb{R}^{2} \tag{2.2}
\end{equation*}
$$

Since $\Gamma_{1}$ is not a circle generally, we draw a auxiliary circle $\Gamma_{2}$ with radius $R$, centered at the origin, such that its interior region contains $\bar{\Omega}_{0} \bigcup \bar{\Omega}_{1}$ properly. The auxiliary boundary divides the exterior region of $\Gamma_{1}$ into two nonoverlapping subdomains: one bounded annular domain denoted by $\Omega_{2}$, another unbounded subdomain denoted by $\Omega_{3}$. Set $\Omega:=\Omega_{1} \bigcup \Gamma_{1} \bigcup \Omega_{2}$, then $\Omega_{3}=\mathbb{R}^{2} \backslash \bar{\Omega}$, and define $u_{i}=\left.u\right|_{\Omega_{i}}, i=1,3$. For the picture see Figure 2.

Define

$$
\begin{equation*}
H_{\Gamma_{0}}^{1}(\Omega):=\left\{v \in H^{1}(\Omega):\left.v\right|_{\Gamma_{0}}=g\right\} \quad \text { and } \quad H_{0}^{1}(\Omega):=\left\{v \in H^{1}(\Omega):\left.v\right|_{\Gamma_{0}}=0\right\} \tag{2.3}
\end{equation*}
$$

Then we rewrite our exterior transmission (1.1a-1.1c) as follows: Find $u \in H_{\Gamma_{0}}^{1}(\Omega)$ such that

$$
\begin{array}{r}
u_{1}=g, \text { on } \Gamma_{0}-\operatorname{div}\left(A \nabla u_{1}\right)=f \quad \text { in } \Omega_{1} \\
u_{1}=u_{2} \text { and }\left(A \nabla u_{1}\right) \cdot \boldsymbol{n}=\frac{\partial u_{2}}{\partial \boldsymbol{n}} \quad \text { on } \Gamma_{1} \\
-\Delta u_{2}=0, \quad \text { in } \Omega_{2} \\
u_{2}=u_{3}, \quad \frac{\partial u_{2}}{\partial \boldsymbol{n}}=\frac{\partial u_{3}}{\partial \boldsymbol{n}} \quad \text { on } \Gamma_{2} \\
-\Delta u_{3}=0, \quad \text { in } \Omega_{3}, \quad u_{3}(x)=O(1) \quad \text { as }|x| \rightarrow \infty \tag{2.4e}
\end{array}
$$

Here we use standard notations for Sobolev spaces and their norms and semi-norms. $(\cdot, \cdot)_{\Omega_{i}}$,


Figure 2: $\Omega=\Omega_{1} \bigcup \Gamma_{1} \bigcup \Omega_{2}$ and $\Gamma$ is an auxiliary circle
$\langle\cdot, \cdot\rangle_{\Gamma_{i}}$ denote the $L^{2}$ inner product in $\Omega_{i}$ and on $\Gamma_{i}$, respectively.
Applying the natural boundary reduction principle([10],[11]) in the exterior region $\Omega_{3}$, we obtain the Poisson integral formula

$$
\begin{equation*}
u(r, \theta)=\frac{r^{2}-R^{2}}{2 \pi} \int_{0}^{2 \pi} \frac{\lambda\left(\theta^{\prime}\right)}{R^{2}+r^{2}-2 \operatorname{Rr} \cos \left(\theta-\theta^{\prime}\right)} d \theta^{\prime}, \quad r>R \tag{2.5}
\end{equation*}
$$

and the natural integral equation

$$
\begin{equation*}
\frac{\partial u(\theta)}{\partial \boldsymbol{n}}=-\frac{1}{4 \pi R} \int_{0}^{2 \pi} \frac{\lambda\left(\theta^{\prime}\right)}{\sin ^{2} \frac{\theta-\theta^{\prime}}{2}} d \theta^{\prime}=-\frac{1}{4 \pi \sin ^{2} \frac{\theta}{2}} * \lambda(\theta) \tag{2.6}
\end{equation*}
$$

Here $*$ denotes the convolution with respect to $\theta$. Define $\mathcal{K}: H^{1 / 2}(\Gamma) \rightarrow H^{-1 / 2}(\Gamma)$ as the natural integral operator, then (2.6) can be written

$$
\begin{equation*}
\frac{\partial u(\theta)}{\partial \boldsymbol{n}} \equiv-\mathcal{K} \lambda(\theta) \tag{2.7}
\end{equation*}
$$

Next, we multiply the divergence partial differential equation in (1.1) by any test function $v \in H_{0}^{1}(\Omega)$ and apply the Green formula to yield

$$
\begin{equation*}
\int_{\Omega_{1}}\left(A \nabla u_{1}\right) \cdot \nabla v d x-\int_{\Gamma_{1}}\left(A \nabla u_{1}\right) \cdot \boldsymbol{n} v d s=\int_{\Omega_{1}} f v d x \tag{2.8}
\end{equation*}
$$

According to the interface condition (2.4b), we have

$$
\begin{equation*}
\int_{\Omega_{1}}\left(A \nabla u_{1}\right) \cdot \nabla v d x-\int_{\Gamma_{1}} \frac{\partial u_{2}}{\partial \boldsymbol{n}} v d s=\int_{\Omega_{1}} f v d x \tag{2.9}
\end{equation*}
$$

In the same way, on $\Omega_{2}$, we get

$$
\begin{equation*}
\int_{\Omega_{2}} \nabla u_{2} \cdot \nabla v d x+\int_{\Gamma_{1}} \frac{\partial u_{2}}{\partial \boldsymbol{n}} v d s-\int_{\Gamma_{2}} \frac{\partial u_{2}}{\partial \boldsymbol{n}} v d s=0 \tag{2.10}
\end{equation*}
$$

which, due to the natural integral equation (2.7), becomes

$$
\begin{equation*}
\int_{\Omega_{2}} \nabla u_{2} \cdot \nabla v d x+\int_{\Gamma_{1}} \frac{\partial u_{2}}{\partial \boldsymbol{n}} v d s+\int_{\Gamma_{2}} v \mathcal{K} u_{2} d s=0 \tag{2.11}
\end{equation*}
$$

Adding (2.9) and (2.11), we obtain the coupled FEM-NBEM variational problem of (2.4):

$$
\left\{\begin{array}{l}
\text { find } u \in H_{\Gamma_{0}}^{1}(\Omega) \text { such that }  \tag{2.12}\\
a(u, v)=f(v), \quad \forall v \in H_{0}^{1}
\end{array}\right.
$$

where $a(u, v)$ is the bilinear form

$$
\begin{equation*}
a(u, v):=(A \nabla u, \nabla v)_{\Omega_{1}}+(\nabla u, \nabla v)_{\Omega_{2}}+\langle\mathcal{K} u, v\rangle_{\Gamma_{2}} \tag{2.13}
\end{equation*}
$$

and $f(v):=(f, v)_{\Omega_{1}}$ is the linear functional.
Lemma 2.1. The natural integral operator $\mathcal{K}: H^{\frac{1}{2}}\left(\Gamma_{2}\right) \rightarrow H^{-\frac{1}{2}}\left(\Gamma_{2}\right)$ is just the DirichletNeumann operator (Steklov-Poincaré operator) for the exterior domain $\Omega_{3}$. It is symmetric and semi-positive definite with respect to the inner product $\langle\cdot, \cdot\rangle_{\Gamma_{2}}$, (see [10], [11]), i.e. there is a positive constant $c$ such that

$$
\begin{equation*}
\langle\mathcal{K} v, v\rangle_{\Gamma_{2}} \geq c\|v\|_{\frac{1}{2}, \Gamma_{2}}^{2}, \quad \forall v \in H^{1 / 2}\left(\Gamma_{2}\right) / P_{0} \tag{2.14}
\end{equation*}
$$

where $P_{0}$ denotes the set of all constants.
which together with the strongly elliptic condition (2.2) of $A$, yields the following coercivity lemma.

Lemma 2.2. Suppose that the matrix valued function $A$ satisfies the condition (2.2). Then, for any function $v \in H_{0}^{1}(\Omega)$ there exists a positive constant $C$ such that

$$
\begin{equation*}
a(v, v) \geq C\left(|v|_{1, \Omega_{1}}^{2}+|v|_{1, \Omega_{2}}^{2}+\|v\|_{\frac{1}{2}, \Gamma_{2}}^{2}\right) \tag{2.15}
\end{equation*}
$$

Therefore, the coercivity and the continuity of $a(u, v)$ and the boundedness of $f(v)$ give the uniqueness solvability of the variational problem (2.12) according to the Lax-Milgram Lemma.

## 3. Finite Element Discretization with Non-matching Grids

In this section, we make a finite element discretization for the subdomains and introduce the non-matching grids method (See [8]) and the construction of basis functions of the Lagrange multiplier space. The main motivation to do this is that we can couple different discretizations in different subdomains in this way. It seems very reasonable especially for the case of singularities of the solution.

Families of finite element triangulations $\mathcal{T}_{h_{i}}, i=1,2$, are associated with $\Omega_{1}$ and $\Omega_{2}$ (e.g. some regular quasi-uniform triangles and curved triangles at the interfaces). We denotes by $h_{i}$ the maximum diameter of the elements of $\mathcal{T}_{h_{i}}$. But in most real calculation, the curved triangles nearby the interfaces are approximated by the straight triangles which has the same nodes with the curved triangles. This simplified method generates only small error. Let $V_{h_{i}}\left(\Omega_{i}\right) \subset$ $H^{1}\left(\Omega_{i}\right), i=1,2$, be the finite element spaces on $\Omega_{i}$ with respect to $\mathcal{T}_{h_{i}}, i=1,2$. Next, we discretize the auxiliary circle $\Gamma_{2}$. Given $n \in \mathbb{N}$, we let $0=t_{0} \leq t_{1} \leq \cdots \leq t_{n}=2 \pi$ be a uniform partition of $[0,2 \pi]$ with $h_{3}=t_{i+1}-t_{i}=\frac{2 \pi}{n}, j=0,1, \cdots, n-1$, which generates a division $\mathcal{T}_{h_{3}}$ on $\Gamma_{2}$. We denote this boundary element space by $V_{h_{3}}\left(\Gamma_{2}\right)$.

The division in $\Omega_{i}, i=1,2$ leads to a division on the interface $\Gamma_{1}$ and $\Gamma_{2}$, so we set

$$
\begin{equation*}
V_{h_{i}}\left(\Gamma_{1}\right)=\left\{\left.v\right|_{\Gamma_{1}}: v \in V_{h_{i}}\left(\Omega_{i}\right), i=1,2\right\} \quad \text { and } \quad V_{h_{2}}\left(\Gamma_{2}\right)=\left\{\left.v\right|_{\Gamma_{2}}: v \in V_{h_{2}}\left(\Omega_{2}\right)\right\} \tag{3.1}
\end{equation*}
$$

We note that the meshes need not match at the interface between any two subdomains. Thus in order to discretize the space $H_{\Gamma_{0}}^{1}(\Omega)$, we have to introduce a Lagrange multipliers space $M_{h}$ used to impose a weak continuity constraint across the interface.

We set the product spaces $Q_{h}$

$$
\begin{equation*}
Q_{h}:=V_{h_{1}}\left(\Omega_{1}\right) \times V_{h_{2}}\left(\Omega_{2}\right) \times V_{h}\left(\Gamma_{2}\right) \tag{3.2}
\end{equation*}
$$

Define

$$
\begin{gather*}
V_{h}=\left\{v_{h}=\left(v_{h_{1}}, v_{h_{2}}, v_{h_{3}}\right) \in Q_{h}: \int_{\Gamma_{1} \cup \Gamma_{2}}\left[v_{h}\right] \cdot \mu d s=0, \forall \mu \in M_{h}\left(\Gamma_{i}\right), i=1,2\right\},  \tag{3.3}\\
V_{h}^{0}=\left\{v_{h} \in V_{h}:\left.v_{h_{1}}\right|_{\Gamma_{0}}=0\right\} \tag{3.4}
\end{gather*}
$$

where [.] denotes the jump of the function $v_{h}$ across the interfaces.
Then we obtain the discrete problem of (2.12) with respect to this kind of non-matching grids discretization.

$$
\left\{\begin{array}{l}
\text { find } u_{h} \in V_{h} \text { such that }  \tag{3.5}\\
a\left(u_{h}, v_{h}\right)=f\left(v_{h}\right), \quad \forall v_{h} \in V_{h}^{0}
\end{array}\right.
$$



Figure 3: Dual basis functions

Remark 3.1. The setting of Lagrange multipliers space would guarantee uniform ellipticity of this discrete problem(see [8]). Then it can be shown that the coupled discrete problem (3.5) has unique solution $\left(u_{h_{1}}, \varphi_{h_{2}}\right) \in V_{h}$.

As we have seen, the construction of Lagrange multiplier space is of great importance for the unique solvability. It is proven in [7] the dual basis mortar method leads to a stable and optimally convergent approximation. Here we apply the same dual basis approach and introduce a new non-matching grids method for unbounded domain problems (see [8]).

Here and below we only discuss the interface $\Gamma_{1}$. A similar definition is to the interface $\Gamma_{2}$. To avoid confusion for the subscript, let us denote by $\Gamma=\Gamma_{i}, i=1,2$. Let $N$ be the number of nodes on $\Gamma$ and $\left\{a_{i}\right\}_{i=1}^{N}$ be the set of nodal points in $\Gamma$. For the nodal basis $\left\{\Phi_{i}\right\}_{i=1}^{N}$, we define by the dual basis $\left\{\Psi_{i}(\theta)\right\}_{i=1}^{N}$ (see Figure 3 for piecewise linear basis funcitons)

$$
\begin{equation*}
\left\langle\Phi_{i}(\theta), \Psi_{i}(\theta)\right\rangle_{\Gamma}=\delta_{i, j}\left\langle\Phi_{i}(\theta), 1\right\rangle_{\Gamma}, \quad 1 \leq i, j \leq N \tag{3.6}
\end{equation*}
$$

where $\delta_{i j}$ is the Kronecker symbol.
Before we begin the analysis of error estimate, we will introduce two important projection operator. Since each interface has two sides, we denote by $\Gamma_{12}$ and $\Gamma_{21}$. Define the projection operator $\Pi_{h}$ in such way : it maps the space $V_{h}\left(\Gamma_{12}\right)$ into $V_{h}\left(\Gamma_{21}\right)$ or maps $V_{h}\left(\Gamma_{21}\right)$ into $V_{h}\left(\Gamma_{12}\right)$. We can see that the choice of side is rather arbitrary. In our case, we choose the fine mesh side as the beginning such as $\Gamma_{12}$. That is to say, Given $v \in L^{2}(\Gamma)$, the values of $\Pi_{h} v \in V_{h}\left(\Gamma_{21}\right)$ can be determined by

$$
\begin{equation*}
\int_{\Gamma}\left(v-\Pi_{h} v\right) \mu d s=0, \quad \forall \mu \in M_{h}(\Gamma) . \tag{3.7}
\end{equation*}
$$

Since $V_{h_{i}}(\Gamma) \subset H^{\frac{1}{2}}(\Gamma)$, the multiplier space $M_{h}(\Gamma)$ generated by the dual basis may be embedded in the dual space of $H^{\frac{1}{2}}(\Gamma)$ with respect to the $L^{2}$-inner product. Therefore, $M_{h}(\Gamma) \subset H^{-\frac{1}{2}}(\Gamma)$. This operator was used in ([2],[7],[8]) and plays a central role in the error analysis of the nonmatching grid finite discretization.

Then define by $P_{h}: L^{2}(\Gamma) \rightarrow M_{h}(\Gamma)$ the usual orthogonal projection operator. We recall its approximation properties in the following lemma. We can verify it in the standard manner and do not include the proof here. See the proof in [7].

Lemma 3.1. For any real number $s, 0 \leq s \leq 1$, there exist constants such that the following estimates hold for any function $v$ in $H^{s}(\Gamma)$ :

$$
\begin{gather*}
\left\|v-P_{h} v\right\|_{0, \Gamma} \leq c h^{s}\|v\|_{s, \Gamma}  \tag{3.8}\\
\left\|v-P_{h} v\right\|_{\left(H^{\frac{1}{2}}(\Gamma)\right)^{\prime}} \leq c h^{s+\frac{1}{2}}\|v\|_{s, \Gamma} . \tag{3.9}
\end{gather*}
$$

Here the dual norm is defined by

$$
\begin{equation*}
\|f\|_{X^{\prime}}:=\sup _{v \in X} \frac{\langle f, v\rangle}{\|v\|_{X}} \tag{3.10}
\end{equation*}
$$

where $X^{\prime}$ is the dual space of the Hilbert space $X$. The definition of operator $P_{h}$ yields to:

$$
\begin{equation*}
\int_{\Gamma}\left(v-P_{h} v\right) \mu d s=0, \quad \forall \mu \in M_{h}(\Gamma) \tag{3.11}
\end{equation*}
$$

where $P_{h} v \in M_{h}(\Gamma)$. Then

$$
\begin{equation*}
\int_{\Gamma}\left(\Pi_{h} v-P_{h} v\right) \mu d s=0 \quad \forall \mu \in M_{h}(\Gamma) \tag{3.12}
\end{equation*}
$$

which means $P_{h} v$ is also the projection of $\Pi_{h} v$ into the multiplier space $M_{h}(\Gamma)$.
The next lemma shows the stability property of the projection operator $\Pi_{h}$ in $L^{2}(\Gamma)$ and $H^{1}(\Gamma)$.

Lemma 3.2. There exist a constant $c>0$ such that for $\forall v \in L^{2}(\Gamma)$

$$
\begin{equation*}
\left\|\Pi_{h} v\right\|_{0, \Gamma} \leq c\|v\|_{0, \Gamma} \tag{3.13}
\end{equation*}
$$

Let $v \in H^{1}(\Gamma)$, then for uniform meshes we have

$$
\begin{equation*}
\left|\Pi_{h} v\right|_{1, \Gamma} \leq c^{\prime}|v|_{1, \Gamma} \tag{3.14}
\end{equation*}
$$

Proof. For any $v \in L^{2}(\Gamma), \Pi_{h} v$ can be written as

$$
\begin{equation*}
\Pi_{h} v=\sum_{i=1}^{N} \alpha_{i} \Phi_{i}(\theta) \tag{3.15}
\end{equation*}
$$

Substitute $\Pi_{h} v$ for (3.15) in (3.7). Due to the global orthogonality relation (3.6) between the nodal basis $\left\{\Phi_{i}\right\}$ and its dual basis $\left\{\Psi_{i}\right\}$, the values of $\alpha_{i}$ can be direct calculated by the formula

$$
\begin{equation*}
\alpha_{i}=\Pi_{h} v\left(a_{i}\right)=\frac{\left.\int_{\Gamma} v\right|_{\Omega_{1}} \Psi_{i} d s}{\int_{\Gamma} \Phi_{i} d s} \tag{3.16}
\end{equation*}
$$

Then some primary inequalities lead to

$$
\begin{align*}
\left\|\Pi_{h} v\right\|_{0, \Gamma}^{2} & \leq \sum_{i=1}^{n} \int_{\Gamma}\left(\Pi_{h} v\left(a_{i}\right)\right)^{2} \Phi_{i}^{2} d s \\
& \leq \sum_{i=1}^{n} \frac{\int_{\Gamma} \Phi_{i}^{2} d s \int_{\gamma} \Psi_{i}^{2} d s}{\left(\int_{\Gamma} \Phi_{i} d s\right)^{2}}\|v\|_{0, \gamma}^{2} \leq c\|v\|_{0, \Gamma}^{2} \tag{3.17}
\end{align*}
$$

in which $\gamma$ denotes $\overline{\operatorname{supp} \Phi_{i}}$ and $c$ is a constant. Thus, (3.13) is obtained.
Let $0=a_{0} \leq a_{1} \leq \cdots \leq a_{N}=2 \pi$ be a corresponding uniform partition on $\Gamma$, which generates a division $\mathcal{T}_{h}$ for the interface $\Gamma$. Let $\hat{T}$ be the 1-dimensional reference element. Let $\left\{\hat{\phi}_{i}\right\}$ be a nodal basis for $\hat{T}$ and let $\left\{\hat{\psi}_{i}\right\}$ be the dual basis with respect to the inner product for $L^{2}(\hat{T})$.

In order to derive the stability in $H^{1}(\Gamma)$, we introduce the piecewise affine mapping

$$
\begin{equation*}
F(\hat{\theta})=\beta \hat{\theta}+\theta_{0} \tag{3.18}
\end{equation*}
$$

which maps $\hat{T}$ one-to-one and onto some $T \in T_{i}=\left[a_{i-1}, a_{i}\right]$. Let $\left\{\hat{\phi}_{i}\right\}$ be the nodal basis for $\hat{T}$ and let $\left\{\hat{\psi}_{i}\right\}$ be the dual basis with respect to the inner product for $L^{2}(\hat{T})$. Then we have (see [3]), for any $v \in H^{1}(T)$,

$$
\begin{gather*}
\|v\|_{1, T} \leq h^{-\frac{1}{2}}\|\hat{v}\|_{1, \hat{T}}  \tag{3.19}\\
\|v\|_{L^{1}(T)} \leq h \left\lvert\,\|\hat{v}\|_{L^{1}(\hat{T})} \leq h\|\hat{v}\|_{1, \hat{T}} \leq\left(h^{\frac{1}{2}}|v|_{0, T}+h^{\frac{3}{2}}|v|_{1, T}\right)\right. \tag{3.20}
\end{gather*}
$$

here, $h:=\max _{T \in \mathcal{T}_{h}} \operatorname{diam}(T)$.
For any $v \in H^{1}(\Gamma)$, we can prove the local estimat

$$
\begin{align*}
\left\|\Pi_{h} v\right\|_{1, T} & \leq \sum_{i=1}^{k}\left|\Pi_{h} v\left(a_{i}\right)\right|\left\|\phi_{i}\right\|_{1, T} \leq C h^{-\frac{1}{2}} \max _{1 \leq i \leq n}\left\|\hat{\phi}_{i}\right\|_{1, \hat{T}} \sum_{i=1}^{k}\left|\Pi_{h} v\left(a_{i}\right)\right| \\
& \leq h^{-\frac{1}{2}} \sum_{i=1}^{k}\left|\frac{\int_{T} \psi_{i} v d s}{\int_{T} \phi_{i} d s}\right| \leq C h^{-\frac{1}{2}} \sum_{i=1}^{k}\left\|\psi_{i}\right\|_{L^{\infty}(T)}\|v\|_{L^{1}(T)} \\
& \leq C h^{-\frac{1}{2}} h^{-1}\left(h^{\frac{1}{2}}|v|_{0, T}+h^{\frac{3}{2}}|v|_{1, T}\right) \\
& \leq C\left(h^{-1}|v|_{0, T}+|v|_{1, T}\right) \tag{3.21}
\end{align*}
$$

where $\phi_{i}$ and $\psi_{i}$ are the nodal basis and dual basis for $T_{i}$. The global estimate (3.14) is obtained by summing over all local contributions.

Then, by an interpolation argument, the following estimate holds for any function $v$ in $H^{\frac{1}{2}}(\Gamma)$ :

$$
\begin{equation*}
\left\|\Pi_{h} v\right\|_{\frac{1}{2}, \Gamma} \leq C\|v\|_{\frac{1}{2}, \Gamma} \tag{3.22}
\end{equation*}
$$

Define the norm

$$
\begin{equation*}
\left\|v_{h}\right\|=\left(\left\|v_{h_{1}}\right\|_{1, \Omega_{1}}^{2}+\left\|v_{h_{2}}\right\|_{1, \Omega_{2}}^{2}+\left\|v_{h_{3}}\right\|_{\frac{1}{2}, \Gamma}^{2}\right)^{1 / 2} \tag{3.23}
\end{equation*}
$$

For our nonconforming situation we use the well-known second Strang's lemma. Let $u=$ $\left(u_{1}, u_{2}, \lambda\right)$ and $u_{h}=\left(u_{h_{1}}, u_{h_{2}}, \lambda_{h_{3}}\right)$ be the solutions of (2.4) and (3.5), respectively. The error can be formulated as follows:

$$
\begin{equation*}
\left\|u-u_{h}\right\| \leq \inf _{\forall v_{h} \in V_{h}^{0}}\left\|u-v_{h}\right\|+\sup _{\forall 0 \neq v_{h} \in V_{h}^{0}} \frac{\int_{\Gamma_{1} \cup \Gamma_{2}} \frac{\partial u}{\partial \boldsymbol{n}}\left[v_{h}\right] d s}{\left\|v_{h}\right\|} \tag{3.24}
\end{equation*}
$$

where $v_{h}=\left(v_{h_{1}}, v_{h_{2}}, v_{h_{3}}\right)$. We note that the first term of the right hand of (3.24) is the approximation error, while the second term is the consistency error. The best approximation error can be estimated by using interpolation inequalities for conforming finite elements and stability properties of the projection $\Pi_{h}$; For estimation of the consistency error, we use the fact
the jump of the solution is orthogonal to the multiplier space $M_{h}$. We summarize the results in the following lemmas.

Lemma 3.3. Assume that the solution $u$ of problem 2.4 is satisfy, for any real number $s$, $\frac{1}{2} \leq \varepsilon_{i} \leq 1, i=1,2,\left.u\right|_{\Omega_{1}} \in H^{1+\varepsilon_{1}}\left(\Omega_{1}\right),\left.u\right|_{\Omega_{2}} \in H^{1+\varepsilon_{2}}\left(\Omega_{2}\right)$ and $\left.u\right|_{\Gamma_{2}} \in H^{\frac{3}{2}}\left(\Gamma_{2}\right)$. Then there exists a function $v_{h} \in V_{h}^{0}$ such that

$$
\begin{equation*}
\left\|u-v_{h}\right\| \leq C\left(h_{1}^{\varepsilon_{1}}\|u\|_{1+\varepsilon_{1}, \Omega_{1}}+h_{2}^{\varepsilon_{2}}\|u\|_{1+\varepsilon_{2}, \Omega_{2}}+h_{3}\|u\|_{\frac{3}{2}, \Gamma_{2}}\right) \tag{3.25}
\end{equation*}
$$

Proof. First, we estimate the error bound on $\Gamma_{1}$. Let $\pi_{h_{i}}, i=1,2$ are the Lagrange interpolation operators in $\Omega_{i}, i=1,2$, respectively. Then we define $v_{h}$ by

$$
\begin{equation*}
v_{h_{1}}=\pi_{h_{1}} u_{1}, \quad v_{h_{2}}=\pi_{h_{2}} u_{2}+\Pi_{h}\left[\pi_{h_{1}}\left(\left.u_{1}\right|_{\Gamma_{1}}\right)-\pi_{h_{2}}\left(\left.u_{2}\right|_{\Gamma_{1}}\right)\right] \tag{3.26}
\end{equation*}
$$

In this way, recalling the definition of projection $\Pi_{h}$, we have

$$
\begin{align*}
\left\langle v_{h_{1}}-v_{h_{2}}, \mu\right\rangle_{\Gamma_{1}} & =\left\langle\left\{\pi_{h_{1}}\left(\left.u_{1}\right|_{\Gamma_{1}}\right)-\pi_{h_{2}}\left(u_{2} \mid \Gamma_{1}\right)\right\}-\Pi_{h}\left\{\pi_{h_{1}}\left(\left.u_{1}\right|_{\Gamma_{1}}\right)-\pi_{h_{2}}\left(\left.u_{2}\right|_{\Gamma_{1}}\right)\right\}, \mu\right\rangle_{\Gamma_{1}} \\
& =0 \tag{3.27}
\end{align*}
$$

For the interface $\Gamma_{2}$, we also define

$$
\begin{equation*}
v_{h_{2}}=\pi_{h_{2}} u_{2}, \quad v_{h_{3}}=\pi_{h_{3}} \lambda+\Pi_{h}\left[\pi_{h_{2}}\left(\left.u_{2}\right|_{\Gamma_{2}}\right)-\pi_{h_{3}} \lambda\right], \tag{3.28}
\end{equation*}
$$

where $\pi_{h_{3}}$ is the usual Lagrange interpolation operators on $\Gamma_{2}$.
Then the trace theorem and the stability properties of $\Pi_{h}$ lead to

$$
\begin{align*}
\inf _{\forall v_{h} \in V_{h}^{0}}\left\|u-v_{h}\right\| & \leq \inf _{\forall v_{h} \in V_{h}^{0}}\left(\left\|u_{1}-v_{h_{1}}\right\|_{1, \Omega_{1}}+\left\|u_{2}-v_{h_{2}}\right\|_{1, \Omega_{2}}+\left\|\lambda-v_{h_{3}}\right\|_{\frac{1}{2}, \Gamma_{2}}\right) \\
& \leq\left\|u-\pi_{h_{1}} u\right\|_{1, \Omega_{1}}+\left\|u-\pi_{h_{2}} u\right\|_{1, \Omega_{2}}+\left\|u-\pi_{h_{3}} u\right\|_{\frac{1}{2}, \Gamma_{2}} \\
& +\left\|\Pi_{h}\left(\pi_{h_{1}} u_{1}-\pi_{h_{2}} u_{2}\right)\right\|_{\frac{1}{2}, \Gamma_{1}}+\left\|\Pi_{h}\left(\pi_{h_{2}} u_{2}-\pi_{h_{3}} \lambda\right)\right\|_{\frac{1}{2}, \Gamma_{2}} \\
& \leq c h_{1}^{\varepsilon_{1}}\left\|u_{1}\right\|_{1+\varepsilon_{1}, \Omega_{1}}+c h_{2}^{\varepsilon_{2}}\left\|u_{2}\right\|_{1+\varepsilon_{2}, \Omega_{2}}+c h_{3}\|\lambda\|_{\frac{3}{2}, \Gamma_{2}} \tag{3.29}
\end{align*}
$$

which complete the proof.

Lemma 3.4. Assume that the solution $u$ of problem 2.4 satisfies the same regularity conditions as the lemma 3.3. then there exists a function $v_{h} \in V_{h}^{0}$ such that

$$
\begin{equation*}
\sup _{\forall 0 \neq v_{h} \in V_{h}^{0}} \frac{\int_{\Gamma_{1} \cup \Gamma_{2}} \frac{\partial u}{\partial \boldsymbol{n}}\left[v_{h}\right] d s}{\left\|v_{h}\right\|} \leq C\left(h_{1}^{\varepsilon_{1}}\|u\|_{1+\varepsilon_{1}, \Omega_{1}}+h_{3}\|u\|_{\frac{3}{2}, \Gamma_{2}}\right) \tag{3.30}
\end{equation*}
$$

Proof. First, we fix our attention to the error on $\Gamma_{1}$. From (3.24) and by using the definition
of the projection operators $P_{h}$ and $\Pi_{h}$, we have

$$
\begin{align*}
\left|\int_{\Gamma_{1}} \frac{\partial u}{\partial \boldsymbol{n}}\left[v_{h}\right] d s\right| & =\left|\int_{\Gamma_{1}} \frac{\partial u}{\partial \boldsymbol{n}}\left(v_{h_{1}}-\Pi_{h} v_{h_{1}}\right) d s\right| \\
& =\left|\int_{\Gamma_{1}}\left(\frac{\partial u}{\partial \boldsymbol{n}}-P_{h} \frac{\partial u}{\partial \boldsymbol{n}}\right)\left(v_{h_{1}}-\Pi_{h} v_{h_{1}}\right) d s\right| \\
& \leq\left.\left\|\frac{\partial u}{\partial \boldsymbol{n}}-P_{h} \frac{\partial u}{\partial \boldsymbol{n}}\right\|\right|_{-\frac{1}{2}, \Gamma_{1}}\left\|v_{h_{1}}-\Pi_{h} v_{h_{1}}\right\|_{\frac{1}{2}, \Gamma_{1}} \\
& \leq\left\|\frac{\partial u}{\partial \boldsymbol{n}}-P_{h} \frac{\partial u}{\partial \boldsymbol{n}}\right\|_{-\frac{1}{2}, \Gamma_{1}}\left(\left\|v_{h_{1}}\right\|_{\frac{1}{2}, \Gamma_{1}}+\left\|v_{h_{2}}\right\|_{\frac{1}{2}, \Gamma_{1}}\right) \tag{3.31}
\end{align*}
$$

Applying the lemma 3.1 and the trace theorem for $v_{h_{i}}$ we deduce that

$$
\begin{align*}
\left|\int_{\Gamma_{1}} \frac{\partial u}{\partial \boldsymbol{n}}\left[v_{h}\right] d s\right| & \leq C h_{1}^{\varepsilon_{1}}\left\|\frac{\partial u}{\partial \boldsymbol{n}}\right\| \|_{\frac{1}{2}+\varepsilon_{1}}\left(\left\|v_{h_{1}}\right\|_{1, \Omega_{1}}+\left\|v_{h_{2}}\right\|_{1, \Omega_{2}}\right) \\
& \leq C h_{1}^{\varepsilon_{1}}\|u\|_{1+\varepsilon_{1}, \Omega_{1}}\left(\left\|v_{h_{1}}\right\|_{1, \Omega_{1}}+\left\|v_{h_{2}}\right\|_{1, \Omega_{2}}\right) \tag{3.32}
\end{align*}
$$

The error bound on $\Gamma_{2}$ can be found in [8]:

$$
\begin{equation*}
\left|\int_{\Gamma_{2}} \frac{\partial u}{\partial \boldsymbol{n}}\left[v_{h}\right] d s\right| \leq C h_{3}\|u\|_{\frac{3}{2}, \Gamma_{2}}\left(\left\|v_{h_{2}}\right\|_{1, \Omega_{2}}+\left\|v_{h_{3}}\right\|_{\frac{1}{2}, \Gamma_{2}}\right) \tag{3.33}
\end{equation*}
$$

Combining (3.32) and (3.33), we obtain (3.30).
The following theorem conclude the discrete error analysis. We can see that the weak constraint across the interfaces we have defined, for the circle interface, is sufficient to guarantee the optimal error estimate.

Theorem 3.1. Assume that the solution $u$ of problem 2.4 have the same regularity as the lemma 3.4 and lemma 3.3, then the exact solution $u$ of (2.4) and the approximate solution $u_{h}$ of (3.5) satisfy

$$
\begin{equation*}
\left\|u-u_{h}\right\| \leq C\left(h_{1}^{\varepsilon_{1}}\|u\|_{1+\varepsilon_{1}, \Omega_{1}}+h_{2}^{\varepsilon_{2}}\|u\|_{1+\varepsilon_{2}, \Omega_{2}}+h_{3}\|u\|_{\frac{3}{2}, \Gamma_{2}}\right) \tag{3.34}
\end{equation*}
$$

where $C>0$ is a constant independent of the mesh parameters $h_{i}, i=1,3$.
Remark 3.2. In order to obtain the optimal error estimation in $V_{h}^{0}$, we should balance the finite element grids in $\Omega_{i}, i=1,3$ such that the fine mesh size $h_{i}, i=1,3$ satisfy $h_{1}^{\varepsilon_{1}} \approx h_{2}^{\varepsilon_{2}} \approx h_{3}$. The main results given in above theorem can be extended to multi-sub-domains case for the exterior problems.

## 4. D-N Alternating Method

The exterior transmission problem (2.4) can be solved by a D-N alternating scheme as follows:

1. Choose initial value $\lambda^{0} \in H^{\frac{1}{2}}\left(\Gamma_{2}\right), \lambda_{1}^{0} \in H^{\frac{1}{2}}\left(\Gamma_{1}\right)$ and set $n:=0$.
2. Solve the Dirichlet problem on the exterior subdomain $\Omega_{3}$ :

$$
\left\{\begin{align*}
-\Delta u_{3}^{n} & =0 & & \text { in } \Omega_{3},  \tag{4.1}\\
u_{3}^{n} & =\Pi_{h} \lambda^{n} & & \text { on } \Gamma_{2}, \\
u_{3}^{n} & =O(1) & & \text { as }|x| \rightarrow \infty .
\end{align*}\right.
$$

3. Solve the mixed boundary value problem in the annular subdomain $\Omega_{2}$ :

$$
\left\{\begin{align*}
-\Delta u_{2}^{n} & =0 & & \text { in } \Omega_{2}  \tag{4.2}\\
\frac{\partial u_{2}^{n}}{\partial \boldsymbol{n}} & =\frac{\partial u_{3}^{n}}{\partial \boldsymbol{n}} & & \text { on } \Gamma_{2} \\
u_{2}^{n} & =\Pi_{h} \lambda_{1}^{n} & & \text { on } \Gamma_{1} .
\end{align*}\right.
$$

4. Solve nonhomogeneous boundary value problem in the annular subdomain $\Omega_{1}$ :

$$
\left\{\begin{align*}
-\operatorname{div}\left(A \nabla u_{1}^{n}\right) & =f & & \text { in } \Omega_{1},  \tag{4.3}\\
\left(A \nabla u_{1}^{n}\right) \cdot \boldsymbol{n} & =\frac{\partial u_{2}^{n}}{\partial \vec{n}} & & \text { on } \Gamma_{1}, \\
u_{1}^{n} & =g & & \text { on } \Gamma_{0} .
\end{align*}\right.
$$

5. Let $\theta_{n}$ be the $n$-th linear relaxation factor selected in computation. Set

$$
\begin{equation*}
\lambda^{n+1}=\theta_{n} u_{2}^{n}+\left(1-\theta_{n}\right) \lambda^{n}, \quad \text { on } \Gamma_{2}, \tag{4.4}
\end{equation*}
$$

and

$$
\begin{equation*}
\lambda_{1}^{n+1}=\theta_{n} u_{1}^{n}+\left(1-\theta_{n}\right) \lambda_{1}^{n}, \quad \text { on } \Gamma_{1} . \tag{4.5}
\end{equation*}
$$

6. Set $n:=n+1$, then goto the second step.

Note that only the approximation of the normal derivative of $u_{3}^{n}$ on the interface $\Gamma_{2}$ is required for solving the mixed boundary value problem in the annular subdomain $\Omega_{2}$. Thus, in practical computation, it is not necessary to actually solve the Dirichlet problem (4.1). Applying the natural integral equation (2.7) and the projection operator $\Pi_{h}$ defined in previous section, we can directly compute the value of $\frac{\partial u_{3}}{\partial \boldsymbol{n}}$ on $\Gamma_{2}$ :

$$
\begin{equation*}
\frac{\partial u_{3}^{n}}{\partial \boldsymbol{n}}=-\mathcal{K}\left(\Pi_{h} u_{2}^{n}\right) . \tag{4.6}
\end{equation*}
$$

And then through the second and third equation of (4.3), we can solve the PDE system of (4.3) in subdomain $\Omega_{1}$.

For the analysis of the convergence of our D-N alternating algorithm, we divide it into two parts. The first part is discussed between the unbounded subdomain $\Omega_{3}$ and the annular subdomain $\Omega_{2}$ (see [8]).

Define

$$
\begin{gathered}
R_{1}: H^{1 / 2}\left(\Gamma_{1}\right) \rightarrow H^{1}\left(\Omega_{1}\right), \\
\phi \mapsto R_{1} \phi,
\end{gathered}
$$

It follows that for any $\phi \in H^{1 / 2}\left(\Gamma_{1}\right)$, if $w=R_{1} \phi$, then $w \in H^{1}\left(\Omega_{1}\right)$ and satisfies

$$
\begin{cases}-\operatorname{div}(A \nabla w)=0, & \text { in } \Omega_{1}  \tag{4.7}\\ w=\phi, & \text { on } \Gamma_{1} \\ w=0, & \text { on } \Gamma_{0}\end{cases}
$$

$$
\begin{gathered}
R_{2}: H^{1 / 2}\left(\Gamma_{1}\right) \rightarrow H^{1}\left(\Omega_{2}\right), \\
\phi \mapsto R_{2} \phi,
\end{gathered}
$$

for any $\phi \in H^{1 / 2}\left(\Gamma_{1}\right)$, if $w=R_{2} \phi$, then $w \in H^{1}\left(\Omega_{2}\right)$ and satisfies

$$
\begin{cases}-\Delta w=0, & \text { in } \Omega_{1}  \tag{4.8}\\ w=\phi, & \text { on } \Gamma_{1} \\ w=0, & \text { on } \Gamma_{2}\end{cases}
$$

Assume that $u_{1}, u_{2}$ satisfy

$$
\begin{gather*}
\begin{cases}-\operatorname{div}\left(A \nabla u_{1}\right)=f, & \text { in } \Omega_{1}, \\
u_{1}=\lambda, & \text { on } \Gamma_{1} \\
u_{1}=g, & \text { on } \Gamma_{0}\end{cases}  \tag{4.9}\\
\begin{cases}-\Delta u_{2}=0, & \text { in } \Omega_{2}, \\
u_{2}=\Pi_{h} \lambda, & \text { on } \Gamma_{1}, \\
u_{2}=g_{2}, & \text { on } \Gamma_{2}\end{cases} \tag{4.10}
\end{gather*}
$$

and $Q_{1}, Q_{2}$ satisfy

$$
\begin{gather*}
\begin{cases}-\operatorname{div}\left(A \nabla Q_{1}\right)=0, & \text { in } \Omega_{1}, \\
Q_{1}=0, & \text { on } \Gamma_{1} \\
Q_{1}=g, & \text { on } \Gamma_{0}\end{cases}  \tag{4.11}\\
\begin{cases}-\Delta Q_{2}=0, & \text { in } \Omega_{2}, \\
Q_{2}=0, & \text { on } \Gamma_{1} \\
Q_{2}=g_{2}, & \text { on } \Gamma_{2}\end{cases} \tag{4.12}
\end{gather*}
$$

Then it's easy to see

$$
\begin{equation*}
u_{1}=R_{1} \lambda+Q_{1}, \quad u_{2}=R_{2} \Pi_{h} \lambda+Q_{2} \tag{4.13}
\end{equation*}
$$

On the interface $\Gamma_{1}, \lambda$ should satisfies

$$
\begin{equation*}
A \nabla u_{1}(\lambda) \cdot \boldsymbol{n}=\nabla u_{2}(\lambda) \cdot \boldsymbol{n} \tag{4.14}
\end{equation*}
$$

Set

$$
\begin{gather*}
S_{1}=-\left(\left(n_{x} a_{11}+n_{x} a_{12}\right) \frac{\partial}{\partial x}\left(R_{1} \cdot\right)+\left(n_{y} a_{21}+n_{y} a_{22}\right) \frac{\partial}{\partial y}\left(R_{1} \cdot\right)\right)  \tag{4.15}\\
S_{2}=\frac{\partial}{\partial \boldsymbol{n}}\left(R_{2} \Pi_{h} \cdot\right) \tag{4.16}
\end{gather*}
$$

where the coefficients $a_{11}, a_{12}, a_{21}, a_{22}$ are the elements of the matrix function $A$. Let $S=$ $S_{1}+S_{2}$. Then we obtain the interface equation

$$
\begin{equation*}
S \lambda=\chi \tag{4.17}
\end{equation*}
$$

here $S$ is just the Steklov-Poicaré operator on the interface $\Gamma_{1}$ and $\chi$ is independent of $\lambda_{1}$ and can be solved beforehand in the subdomains.
Theorem 4.1. The $D-N$ alternating method is equivalent to the preconditioned Richardson iterative method

$$
\begin{equation*}
S_{1}\left(\lambda^{n+1}-\lambda^{n}\right)=\theta_{n}\left(\chi-S \lambda^{n}\right) \tag{4.18}
\end{equation*}
$$

Proof. We consider the error $e_{k}^{n}=u-u_{k}^{n}, k=1,2$ and $\mu^{n}=\Pi_{h}\left(\lambda-\lambda^{n}\right)$, where $\lambda=\left.u\right|_{\Gamma_{1}}$. Then the error terms $e_{1}^{n}$ and $e_{2}^{n}$ satisfy the following equations, respectively,

$$
\begin{align*}
& \left\{\begin{aligned}
-\Delta e_{2}^{n}=0, & \text { in } \Omega_{2}, \\
e_{2}^{n}=\mu^{n}, & \text { on } \Gamma_{1}, \\
e_{2}^{n}=0, & \text { on } \Gamma_{2},
\end{aligned}\right.  \tag{4.19a}\\
& \left\{\begin{array}{rll}
-\operatorname{div}\left(A e_{1}^{n}\right)=0, & \text { in } \Omega_{1} \\
e_{1}^{n}=0, & \text { on } \Gamma_{0} \\
\frac{\partial e_{1}^{n}}{\partial \boldsymbol{n}} & =-\mathcal{K}\left(\mu^{n}\right), & \text { on } \Gamma_{1}
\end{array}\right. \tag{4.19b}
\end{align*}
$$

and

$$
\begin{equation*}
\mu^{n+1}=\theta_{n} \Pi_{h}\left(\left.e_{1}^{n}\right|_{\Gamma_{1}}\right)+\left(1-\theta_{n}\right) \mu^{n} \tag{4.20}
\end{equation*}
$$

Therefore,

$$
\begin{equation*}
e_{1}^{n}=R_{1}\left(\left.e_{1}^{n}\right|_{\Gamma_{1}}\right), \quad e_{2}^{n}=R_{2}\left(\left.e_{2}^{n}\right|_{\Gamma_{1}}\right)=R_{2} \mu^{n} \tag{4.21}
\end{equation*}
$$

Then we have

$$
\begin{equation*}
S_{1}\left(\left.e_{1}^{n}\right|_{\Gamma_{1}}\right)=-\frac{\partial}{\partial \boldsymbol{n}}\left[R_{1}\left(\left.e_{1}^{n}\right|_{\Gamma_{1}}\right)\right]=\mathcal{K}\left(\mu^{n}\right)=-\frac{\partial}{\partial \boldsymbol{n}}\left[\Pi_{h}\left(\lambda-\lambda^{n}\right)\right]=-S_{2}\left(\lambda-\lambda^{n}\right) \tag{4.22}
\end{equation*}
$$

Since

$$
\begin{equation*}
\lambda^{n+1}-\lambda^{n}=\theta_{n}\left(\left.u_{1}^{n}\right|_{\Gamma_{1}}-\lambda^{n}\right) \tag{4.23}
\end{equation*}
$$

finally we derive

$$
\begin{align*}
S_{1}\left(\lambda^{n+1}-\lambda^{n}\right) & =S_{1}\left[\theta_{n}\left(\left.u_{1}^{n}\right|_{\Gamma_{1}}-\lambda^{n}\right)\right]=\theta_{n}\left[S_{1}\left(\left.u_{1}^{n}\right|_{\Gamma_{1}}-\lambda\right)+S_{1}\left(\lambda-\lambda^{n}\right)\right] \\
& =\theta_{n}\left(S_{1}+S_{2}\right)\left(\lambda-\lambda^{n}\right)=\theta_{n} S\left(\lambda-\lambda^{n}\right)=\theta_{n}\left(\chi-S \lambda^{n}\right) \tag{4.24}
\end{align*}
$$

The proof is completed.
Define by $R_{h}^{i}: H^{1 / 2}\left(\Gamma_{1}\right) \rightarrow V_{h_{i}}\left(\Omega_{i}\right)$ the discrete harmonic extension operators, i.e. for any $\lambda_{h} \in H^{1 / 2}\left(\Gamma_{1}\right), R_{h}^{i} \lambda_{h} \in V_{h_{i}}\left(\Omega_{i}\right)$ satisfies

$$
\left\{\begin{array}{ll}
a_{i}\left(R_{h}^{i} \lambda_{h}, v_{h}\right)=0, & \forall v_{h} \in V_{h_{i}}\left(\Omega_{i}\right),  \tag{4.25}\\
R_{h}^{i} \lambda_{h}=\lambda_{h}, & \text { on } \Gamma_{1}, \\
R_{h}^{i} \lambda_{h}=0, & \text { on } \partial \Omega_{i} \backslash \Gamma_{1},
\end{array} \quad i=1,2\right.
$$

where $a_{i}(u, v), i=1,2$, are the bilinear forms corresponding to the harmonic problems in $\Omega_{i}$.
The discrete scheme of the D-N alternating methods is as follows:

1. Choose an initial value $\lambda_{h}^{0} \in V_{h_{2}}\left(\Gamma_{2}\right), \lambda_{1 h}^{0} \in V_{h_{1}}\left(\Gamma_{1}\right)$ and set $n:=0$.
2. Apply the natural integral equation on $\Gamma_{2}$ and compute:

$$
\begin{equation*}
\frac{\partial u_{h_{3}}^{n}}{\partial \boldsymbol{n}}=-\mathcal{K} \Pi_{h} \lambda_{h}^{n}, \quad \text { on } \Gamma_{2} . \tag{4.26}
\end{equation*}
$$

3. Solve the discrete mixed boundary value problem in $\Omega_{2}$ :

$$
\left\{\begin{array}{l}
\text { find } u_{h_{2}} \in V_{h_{2}}\left(\Omega_{2}\right) \text { such that } \forall v_{h_{2}} \in V_{h_{2}}\left(\Omega_{2}\right)  \tag{4.27}\\
\int_{\Omega_{2}} \nabla u_{h_{2}}^{n} \cdot \nabla v_{h_{2}} d x=\int_{\Gamma_{2}} \frac{\partial u_{h_{3}}^{n}}{\partial \boldsymbol{n}} \Pi_{h} v_{h_{2}} d s+\int_{\Omega_{2}} \nabla R_{h}^{2} \Pi_{h} \lambda_{1 h}^{n} \cdot \nabla v_{h_{2}} d s \\
u_{h_{2}}^{n}=\Pi_{h} \lambda_{1 h}^{n} \quad \text { on } \Gamma_{1} .
\end{array}\right.
$$

4. Solve the discrete mixed boundary value problem in $\Omega_{1}$ :

$$
\left\{\begin{array}{l}
\text { find } u_{h_{1}} \in V_{h_{1}}\left(\Omega_{1}\right) \text { such that } \forall v_{h_{1}} \in V_{h_{1}}\left(\Omega_{1}\right),  \tag{4.28}\\
\int_{\Omega_{1}} A \nabla u_{h_{1}}^{n} \cdot \nabla v_{h_{1}} d x=\int_{\Omega_{1}} f v_{h_{1}} d x+\int_{\Gamma_{1}} \Pi_{h} v_{h_{1}} \frac{\partial}{\partial \boldsymbol{n}} u_{h_{2}}^{n} d s
\end{array}\right.
$$

5. Set $\lambda_{h}^{n+1}=\theta_{n} u_{h_{1}}^{n}+\left(1-\theta_{n}\right) \lambda_{h}^{n}$ on $\Gamma_{2}$, and $\lambda_{1 h}^{n+1}=\theta_{n} u_{h_{1}}^{n}+\left(1-\theta_{n}\right) \lambda_{1 h}^{n}$ on $\Gamma_{1}$,

6 . Let $n:=n+1$, then goto the second step.
Theorem 4.2. The discrete $D-N$ alternating method is equivalent to the associated preconditioned Richardson iterative method

$$
\begin{equation*}
S_{h}^{1}\left(\lambda_{h}^{n+1}-\lambda_{h}^{n}\right)=\theta_{n}\left(\chi-S_{h} \lambda^{n}\right) \tag{4.29}
\end{equation*}
$$

and converges if $0<\theta_{n}<1$, where $\theta_{n}$ is the $n$-th relaxation factor in the computations.
The proof is similar to [8].

## 5. Numerical Experiments

We now provide some numerical examples to illustrate the theoretical results in precious sections. We use the D-N alternating method to solve the discrete problem (3.5).
Example 1. The solution domain and the triangular mesh are illustrated in Figure 4. In this example we consider $A$ equals the identity matrix $I$. The exact solution of the problem (2.4) is given by

$$
\begin{equation*}
u_{1}(x, y)=\frac{1}{x^{2}+y^{2}}, \quad u_{2}(x, y)=\frac{x}{x^{2}+y^{2}} \tag{5.1}
\end{equation*}
$$

with

$$
\begin{equation*}
f(x, y)=-\frac{4}{\left(x^{2}+y^{2}\right)^{2}} \tag{5.2}
\end{equation*}
$$

and the boundary value $g$ on $\Gamma_{0}$ is computed from the exact solution $\left.u_{1}\right|_{\Gamma_{0}}$.
The discrete system (3.5) is solved by the D-N alternating method proposed in Section 4. In our computation we compute the natural boundary element matrix $\mathbf{K}$, and then solve the


Figure 4: The uniform triangular division for Example 1


Figure 5: The uniform triangular division for Example 2
corresponding linear systems by the conjugate gradient method. In Table 1, the discretization errors are given in the $L^{2}$-norm and energy norm as well as $L^{\infty}$-norm for the two subdomains, where NEM1 and NEM2 are the number of elements in $\Omega_{i}, i=1,2$, respectively. iters indicates the number of iterations required by the D-N alternating method. Here we adopt the notation

$$
\begin{equation*}
\left\|u-u_{h_{i}}\right\|_{\infty, \Omega_{i}}:=\max _{j \in\{1, \cdots, N N M i\}}\left|u\left(x_{j}\right)-u_{h_{j}}\left(x_{j}\right)\right|, \quad i=1,2 \tag{5.3}
\end{equation*}
$$

We observe that the energy error is of order $h$ and the error in the $L^{2}$-norm is of order $h^{2}$. In Figure 6, the errors in the energy norm and the $L^{2}$-norm are shown versus the number of elements, which also indicate that the errors are asymptotic optimal. The exact solutions and numerical solutions are visualized in Figure 9 and Figure 8.

Table 1: Comparison of errors for CG solvers.

| NEM1 | NEM2 | $\left\\|u_{1}-u_{h_{1}}\right\\|_{0, \Omega_{1}}$ | Ratio | $\left\\|u_{1}-u_{h_{1}}\right\\|_{1, \Omega_{1}}$ | Ratio | $\left\\|u_{1}-u_{h_{1}}\right\\|_{\infty, \Omega_{1}}$ |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 64 | 16 | $1.3957 \times 10^{-1}$ | - | $8.9967 \times 10^{-1}$ | - | $4.0603 \times 10^{-2}$ |
| 256 | 64 | $3.7894 \times 10^{-2}$ | 3.683 | $4.5687 \times 10^{-1}$ | 1.969 | $8.9620 \times 10^{-3}$ |
| 1024 | 256 | $9.7081 \times 10^{-3}$ | 3.903 | $2.2980 \times 10^{-1}$ | 1.988 | $2.0970 \times 10^{-3}$ |
| 4096 | 1024 | $2.4428 \times 10^{-3}$ | 3.974 | $1.1510 \times 10^{-1}$ | 1.997 | $5.0595 \times 10^{-4}$ |
| 16384 | 4096 | $6.1171 \times 10^{-4}$ | 3.993 | $5.7574 \times 10^{-2}$ | 1.999 | $1.2403 \times 10^{-4}$ |
| 64512 | 16384 | $1.5902 \times 10^{-4}$ | 3.847 | $2.9207 \times 10^{-2}$ | 1.971 | $3.1644 \times 10^{-5}$ |
| 258043 | 64512 | $3.9781 \times 10^{-5}$ | 3.997 | $1.4604 \times 10^{-2}$ | 2.000 | $7.9193 \times 10^{-6}$ |


| $\left\\|u_{2}-u_{h_{2}}\right\\|_{0, \Omega_{2}}$ | Ratio | $\left\\|u_{2}-u_{h_{2}}\right\\|_{1, \Omega_{2}}$ | Ratio | $\left\\|u_{2}-u_{h_{2}}\right\\|_{\infty, \Omega_{2}}$ | iters |
| :---: | :---: | :---: | :---: | :---: | :---: |
| $1.2019 \times 10^{-1}$ | - | $4.4410 \times 10^{-1}$ | - | $6.1547 \times 10^{-2}$ | 11 |
| $3.0114 \times 10^{-2}$ | 3.783 | $2.1131 \times 10^{-1}$ | 2.007 | $1.6227 \times 10^{-2}$ | 11 |
| $7.4141 \times 10^{-3}$ | 3.860 | $1.0449 \times 10^{-1}$ | 1.996 | $3.3571 \times 10^{-3}$ | 11 |
| $1.8516 \times 10^{-3}$ | 3.916 | $5.2155 \times 10^{-2}$ | 1.998 | $7.2441 \times 10^{-4}$ | 12 |
| $4.6399 \times 10^{-4}$ | 4.004 | $2.6070 \times 10^{-2}$ | 1.999 | $1.9160 \times 10^{-4}$ | 12 |
| $1.2697 \times 10^{-4}$ | 3.836 | $1.3035 \times 10^{-2}$ | 2.000 | $4.8141 \times 10^{-5}$ | 13 |
| $6.4486 \times 10^{-5}$ | 3.290 | $6.5416 \times 10^{-3}$ | 1.984 | $1.2032 \times 10^{-5}$ | 13 |

Example 2. Let $\Omega_{0}$ be a square with center at ( 0,0 ) and side lengths given by 1 (see Figure $5)$. The exact solution is given by

$$
\begin{equation*}
u_{1}(x, y)=\sin (\pi x) \sin (\pi y) \quad u_{2}(x, y)=\frac{x^{2}-y^{2}}{\left(x^{2}+y^{2}\right)^{2}}+\frac{x}{x^{2}+y^{2}} \tag{5.4}
\end{equation*}
$$

Here we choose the matrix valued function as

$$
A=\left(\begin{array}{ll}
\epsilon & 0  \tag{5.5}\\
0 & 1
\end{array}\right)
$$

with

$$
\begin{equation*}
f(x, y)=(\epsilon+1) \pi^{2} \sin (\pi x) \sin (\pi y) \tag{5.6}
\end{equation*}
$$

In Table 2, order $h$ for the energy norm and the order $h^{2}$ for the $L^{2}$-norm can be observed. The discretization errors are plotted in Figure 10.


Figure 6: Discretization errors in $L^{2}$-norm and $H^{1}$-norm versus number of elements for Example 1.


Figure 7: Finite element solutions in $\Omega_{1}$ for example 1.


Figure 8: Finite element solutions in $\Omega_{2}$ for example 1.


Figure 9: Finite element errors in $\Omega_{1}$ for example 1.

Table 2: Comparison of errors for CG solvers.

| NEM1 | NEM2 | $\left\\|u_{1}-u_{h_{1}}\right\\|_{0, \Omega_{1}}$ | Ratio | $\left\\|u_{1}-u_{h_{1}}\right\\|_{1, \Omega_{1}}$ | Ratio | $\left\\|u_{1}-u_{h_{1}}\right\\|_{\infty, \Omega_{1}}$ |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 64 | 16 | $6.8473 \times 10^{-1}$ | - | 4.18518 | - | $2.7447 \times 10^{-1}$ |
| 256 | 64 | $2.7053 \times 10^{-1}$ | 2.531 | 2.52416 | 1.658 | $1.6151 \times 10^{-1}$ |
| 1024 | 256 | $7.6572 \times 10^{-2}$ | 3.533 | 1.32712 | 1.902 | $3.0370 \times 10^{-2}$ |
| 4096 | 1024 | $1.9716 \times 10^{-2}$ | 3.884 | 0.67223 | 1.974 | $8.6012 \times 10^{-3}$ |
| 16384 | 4096 | $4.9566 \times 10^{-3}$ | 3.978 | 0.33723 | 1.993 | $2.2386 \times 10^{-3}$ |
| 64512 | 16384 | $1.2375 \times 10^{-3}$ | 4.005 | 0.16964 | 1.988 | $5.6412 \times 10^{-4}$ |
| 258043 | 64512 | $2.9445 \times 10^{-4}$ | 4.203 | 0.08484 | 2.000 | $1.4442 \times 10^{-4}$ |


| $\left\\|u_{2}-u_{h_{2}}\right\\|_{0, \Omega_{2}}$ | Ratio | $\left\\|u_{2}-u_{h_{2}}\right\\|_{1, \Omega_{2}}$ | Ratio | $\left\\|u_{2}-u_{h_{2}}\right\\|_{\infty, \Omega_{2}}$ | iters |
| :---: | :---: | :---: | :---: | :---: | :---: |
| $1.8424 \times 10^{-1}$ | - | 0.67003 | - | $6.1547 \times 10^{-2}$ | 11 |
| $5.1630 \times 10^{-2}$ | 3.568 | 0.34369 | 1.950 | $4.0774 \times 10^{-2}$ | 8 |
| $1.2182 \times 10^{-2}$ | 4.238 | 0.16975 | 2.025 | $8.9008 \times 10^{-3}$ | 9 |
| $3.0108 \times 10^{-3}$ | 4.046 | 0.08475 | 2.003 | $1.8041 \times 10^{-3}$ | 7 |
| $7.5245 \times 10^{-4}$ | 4.001 | 0.04238 | 2.000 | $4.2226 \times 10^{-4}$ | 7 |
| $1.9506 \times 10^{-4}$ | 3.858 | 0.02119 | 2.000 | $1.2427 \times 10^{-4}$ | 7 |
| $7.4485 \times 10^{-5}$ | 2.620 | 0.01064 | 1.992 | $5.5621 \times 10^{-5}$ | 7 |



Figure 10: Discretization errors in $L^{2}$-norm and $H^{1}$-norm versus number of elements for Example 2.

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