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A Unified Framework for Double Sweep Methods for the Helmholtz Equation

Nacime Bouziani*, Frédéric Nataf and Pierre-Henri Tournier †

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Abstract

We consider sweeping domain decomposition preconditioners to solve the Helmholtz equation in the case of stripwise domain decomposition with or without overlaps. We unify their derivation and convergence studies as Jacobi, Gauss-Seidel or Symmetric Gauss-Seidel for different numbering of the unknowns. This enables the theoretical comparisons of the double sweep methods in [23, 31] with that of [27, 30, 28]. It also makes possible the introduction of two new sweeping algorithms. We provide numerical test cases that assess the validity of the theoretical studies.

1 Introduction

Solving the Helmholtz equation numerically is a difficult task, especially when dealing with high-frequency regimes, heterogeneous media or reflecting boundary conditions. Over the last decades a lot of effort and progress has been made in developing efficient algorithms to solve the ill-conditioned linear system resulting from the Helmholtz operator's discretization. Domain decomposition methods (DDM) try to overcome these difficulties. They are hybrid methods that combine

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direct solvers in subdomains and iterative matching of the solutions across the subdomains. The original domain decomposition method introduced by Schwarz [26] only works for overlapping domain decomposition. P. L. Lions [20] introduced a new variant of this algorithm where the Dirichlet interface conditions are replaced by Robin interface conditions, his method can be applied to both overlapping and nonoverlapping subdomains. He showed convergence for the elliptic case for a non overlapping domain decomposition. The proof was extended by Després [9] to the Helmholtz equation and later on to the time-harmonic Maxwell equations [10]. More recently, sweeping-type domain decomposition methods have been made popular due to their capability to achieve nearly-linear asymptotic complexity. A sweeping algorithm was first proposed and analyzed in [23] for convection-diffusion operators. Sweeping approaches for Helmholtz problems have recently seen their interest renewed as a preconditioner to speed up the convergence of the solver: the double sweep preconditioners of Stolk for overlapping decomposition [27, 28] and of Vion and Geuzaine for non overlapping decomposition with high order interface conditions [30, 31], the PML-based sweep method of Stolk [27], and the polarized traces method of Zepeda-Núñez and Demanet [32]. There also exists sweeping-type methods that are not domain decomposition based methods, such as the sweeping PML preconditioner of Engquist and Ying [13, 12], the source transfer method [5], see [15] for a complete panorama and relations between these methods.

The highlights of the article are:

- New formulation of [27, 30] which allows for two new variants introduced in § 4.4
- Unified convergence analysis for the three above mentioned algorithms which enables a comparison with the algorithms proposed in [10], [23, 31], see Table 1.
- Theoretical and numerical comparisons that show the advantage of the double sweep method in [23, 31] over the double sweep method in [27, 30].

We first state the problem in § 2. Then we explain in § 3 how to substructure the problem in terms of interface unknowns and how to apply classical linear algebra preconditioners (Jacobi,Gauss-Seidel and Symmetric Gauss-Seidel) to two different unknown numberings. After these preparatory tools have been introduced, we present and analyze in § 4 the convergence of six sweeping algorithms. Numerical results are shown in § 5.

2 Statement of the problem and two classical algorithms

We consider the Helmholtz equation in a bounded domain $\Omega \subset \mathbb{R}^2$ with frequency ω , velocity c and wavenumber k defined by $k = \omega/c$:

$$(-k^2 - \Delta) u = f \text{ in } \Omega + \text{ appropriate boundary conditions on } \partial\Omega .$$
 (1)

We consider a layered decomposition of Ω into N slices $(\Omega_i)_{1 \le i \le N}$ with or without overlap, see Figure 1. The boundary $\partial \Omega_i \setminus \partial \Omega$ is written as the disjoint union of $\Gamma_{i,l}$ and $\Gamma_{i,r}$ where $\Gamma_{i,l}$ is on the left of Ω_i and $\Gamma_{i,r}$ is on its right $(\Omega_{1,l} = \emptyset$ and $\Omega_{N,r} = \emptyset$) (see Fig. 2). The outward normal from Ω_i on $\Gamma_{i,l}$ (resp. $\Gamma_{i,r}$) is denoted by $\vec{n}_{i,l}$ (resp. $\vec{n}_{i,r}$). The problem (1) can



Figure 1: Decomposition into vertical strips

be solved iteratively using a domain decomposition method where we solve locally on each subdomain Ω_i the equation (1) with appropriate boundary conditions on the physical boundaries and interfaces [9]. The method writes: Solve in parallel:

$$\begin{cases} \left(-k^{2}-\Delta\right)u_{i}^{n+1}=f \text{ in }\Omega_{i}, \ 1\leq i\leq N\\ \mathcal{B}_{i,l}\left(u_{i}^{n+1}\right)=\mathcal{B}_{i,l}\left(u_{i-1}^{n}\right) \text{ on }\Gamma_{i,l}, \ 2\leq i\leq N\\ \mathcal{B}_{i,r}\left(u_{i}^{n+1}\right)=\mathcal{B}_{i,r}\left(u_{i+1}^{n}\right) \text{ on }\Gamma_{i,r}, \ 1\leq i\leq N-1\\ + \text{ appropriate boundary conditions on }\partial\Omega\cap\partial\Omega_{i}, \end{cases}$$

$$(2)$$

where $\mathcal{B}_{i,l}$ and $\mathcal{B}_{i,r}$ are the interface conditions. Here *I* denotes the square root of -1 ($I^2 = -1$). For sake of simplicity, we consider here either zeroth-order ABC (ABC0):

$$\begin{cases} \mathcal{B}_{i,l} = \partial_{\vec{n}_{i,l}} + Ik \\ \mathcal{B}_{i,r} = \partial_{\vec{n}_{i,r}} + Ik , \end{cases}$$
(3)

exact ABC as interface conditions or subdomains coupling via Perfectly Matched Layer (PML) as in [27]. In practice, when used as truncation conditions on artificial boundaries, ABC0 yields high non physical reflection of the order of 10% of the incoming wave whereas by definition exact ABC leads to no reflection at all. ABC0 boundary conditions are easy and cheap to use whereas exact ABC, always defined in theory, are sometimes impossible to use in practice (e.g. variable coefficients problems). As a result, there is room for compromise and a great deal of literature has been devoted to introduce various high order interface conditions. The most notable techniques are based on partial differential operators (see [11] and [2]) or on PML (see [19] and [6]), see [22] as well for a gentle introduction to this question. When used as truncation conditions, the final accuracy of the computation does depend on the choice of the ABC.

But let us stress that here, ABCs are used as interface conditions in domain decomposition methods so that the final accuracy of the computed result is not impacted by the choice of the ABC, only the iteration counts to solution are impacted. High-order ABC tailored to domain decomposition improve the iteration counts with respect to ABC0, see e.g. [14, 1]. A remarkable super convergence result noticed in [17] for the two subdomain case and in [24] for a decomposition into N strips (see Fig. 1) is that the use of exact ABCs as interface conditions yields convergence in a number of iterations equal to the number of subdomains. Since the solution in a subdomain depends on the value of the right everywhere and that in algorithm (2) a subdomain receives data only from its neighbors it is not possible to achieve convergence in less than N iterations. When looking at the proof in [24], it appears that somehow the correct information flows from the extreme subdomain labelled 1 to the right and at the same time from subdomain labelled N to the left.

This motivated the search for algorithms which would sweep over the subdomains to reach convergence in one iteration consisting of a double sweep. In the sequel we will consider the double sweep algorithms introduced in [23, 31] and in [27, 30] since they converge in one double sweep if exact ABCs are used as interface conditions. Note that both algorithms were named double sweep algorithms which could be confusing. But as we shall see they are actually not the same and have different convergence rates when implemented with non exact ABCs. We start with the double sweep algorithm introduced in [23]. It consists in double sweeps over the subdomains:

left to right sweep:

$$\begin{cases} \left(-k^{2}-\Delta\right)u_{i}^{n+1/2} = f \text{ in }\Omega_{i}, \ 1 \leq i \leq N\\ \mathcal{B}_{i,l}\left(u_{i}^{n+1/2}\right) = \mathcal{B}_{i,l}\left(u_{i-1}^{n+1/2}\right) \text{ on }\Gamma_{i,l}, \ 2 \leq i \leq N\\ \mathcal{B}_{i,r}\left(u_{i}^{n+1/2}\right) = \mathcal{B}_{i,r}\left(u_{i+1}^{n}\right) \text{ on }\Gamma_{i,r}, \ 1 \leq i \leq N-1\\ + \text{ appropriate boundary conditions on }\partial\Omega \cap \partial\Omega_{i}, \end{cases}$$

$$\tag{4}$$

then, right to left sweep:

$$\begin{cases} \left(-k^{2}-\Delta\right)u_{i}^{n+1}=f \text{ in }\Omega_{i}, \ 1\leq i\leq N\\ \mathcal{B}_{i,l}\left(u_{i}^{n+1}\right)=\mathcal{B}_{i,l}\left(u_{i-1}^{n+1/2}\right) \text{ on }\Gamma_{i,l}, \ 2\leq i\leq N\\ \mathcal{B}_{i,r}\left(u_{i}^{n+1}\right)=\mathcal{B}_{i,r}\left(u_{i+1}^{n+1}\right) \text{ on }\Gamma_{i,r}, \ 1\leq i\leq N-1\\ + \text{ appropriate boundary conditions on }\partial\Omega\cap\partial\Omega_{i}. \end{cases}$$

$$(5)$$

It can be seen as a Symmetric Gauss-Seidel version of the Jacobi algorithm (2). This statement will be made more precise in section 4. Its study is made easier in its substructured formulation which is moreover needed to introduce the double sweep algorithm of [27, 30]. We devote the next section to substructuring.

3 Substructuring

In this section, we introduce the substructured problem related to algorithm (2). This will be the basis for the unified framework of DD sweeping methods. Substructuring consists in reformulating the iterative method considering only surface unknowns on the interfaces:

$$\begin{cases} h_{i,l}^{n} \coloneqq \mathcal{B}_{i,l}\left(u_{i}^{n}\right), \text{ on } \Gamma_{i,l} \text{ for } 2 \leq i \leq N \\ h_{i,r}^{n} \coloneqq \mathcal{B}_{i,r}\left(u_{i}^{n}\right), \text{ on } \Gamma_{i,r} \text{ for } 1 \leq i \leq N-1. \end{cases}$$

$$\tag{6}$$

Considering the global vector h^n containing the local unknowns $(h_{i,l}^n)_{2 \le i \le N}$ and $(h_{i,r}^n)_{1 \le i \le N-1}$, we can reformulate the additive Schwarz method (2) as a Jacobi algorithm on h^n :

$$h^{n+1} := \mathscr{T}(h^n) + G, \tag{7}$$



Figure 2: Local problem on the subdomain Ω_i

where the iteration operator \mathscr{T} can be written in the form of an operator valued matrix and G refers to the contribution of the right-hand side f, see [24]. The above equation is what is called a substructured formulation of the volumic algorithm (2). Taking the limit as n tends to infinity, we see that we look for a vector h such that,

$$(Id - \mathscr{T})(h) = G.$$
(8)

Equation (8) is what is called the substructured formulation of the domain decomposition problem. In order to define more precisely the operator \mathscr{T} , we introduce for each subdomain an operator S_i which takes three arguments, two surface functions h_l and h_r and a volume function f and maps them to the local solution v:

$$S_i(h_l, h_r, f) \coloneqq v \tag{9}$$

where $v: \Omega_i \mapsto \mathbb{C}$ satisfies:

$$\begin{cases} \left(-k^{2}-\Delta\right)v = f \text{ in }\Omega_{i} \\ \mathcal{B}_{i,l}(v) = h_{l} \text{ on }\Gamma_{i,l} \quad (2 \leq i \leq N) \\ \mathcal{B}_{i,r}(v) = h_{r} \text{ on }\Gamma_{i,r} \quad (1 \leq i \leq N-1) \\ + \text{ appropriate boundary conditions on }\partial\Omega \cap \partial\Omega_{i}, \end{cases}$$
(10)

for 1 < i < N. For i = 1, the definition of S_1 is similar except that it takes only the two arguments (h_r, f) since domain Ω_1 has no left interface and similarly operator S_N takes only the two arguments (h_l, f) since domain Ω_N has no right interface. As of now, for sake of simplicity and by abuse of notation, $S_1(h_l, h_r, f)$ (resp. $S_N(h_l, h_r, f)$) will refer to $S_1(h_r, f)$ (resp. $S_N(h_l, f)$).

Next, we introduce the surface right hand-side G(f) by

$$\begin{aligned}
G_{i,l}(f) &:= & \mathcal{B}_{i,l}(S_{i-1}(0,0,f)), & 2 \le i \le N \\
G_{i,r}(f) &:= & \mathcal{B}_{i,r}(S_{i+1}(0,0,f)), & 1 \le i \le N-1.
\end{aligned} \tag{11}$$

and the substructured operator \mathcal{T} by:

$$\mathcal{T}(h)_{i+1,l} := \mathcal{B}_{i+1,l}(S_i(h_{i,l}, h_{i,r}, 0)), \quad 1 \le i \le N - 1
\mathcal{T}(h)_{i-1,r} := \mathcal{B}_{i-1,r}(S_i(h_{i,l}, h_{i,r}, 0)), \quad 2 \le i \le N.$$
(12)

The operator $\mathcal T$ has thus the following possibly non zero entries:

$$\begin{aligned}
\mathscr{T}_{(i+1,l)(i,l)} &:= \mathcal{B}_{i+1,l}(S_i(\cdot,0,0)), \ 1 \le i \le N-1 \\
\mathscr{T}_{(i+1,l)(i,r)} &:= \mathcal{B}_{i+1,l}(S_i(0,\cdot,0)), \ 1 \le i \le N-1 \\
\mathscr{T}_{(i-1,r)(i,r)} &:= \mathcal{B}_{i-1,r}(S_i(0,\cdot,0)), \ 2 \le i \le N \\
\mathscr{T}_{(i-1,r)(i,l)} &:= \mathcal{B}_{i-1,r}(S_i(\cdot,0,0)), \ 2 \le i \le N .
\end{aligned}$$
(13)

3.1 Double sweep algorithm of [27, 30]

The rationale behind the algorithm in [27, 30] is that for exact ABC used as interface conditions, we have a decoupling of left and right interface unknowns since

$$\mathscr{T}_{(i+1,l)(i,r)}^{EABC} \equiv 0 \text{ and } \mathscr{T}_{(i-1,r)(i,l)}^{EABC} \equiv 0.$$
(14)

Then, the operator \mathscr{T}^{EABC} is nilpotent of order N-1. This is related to the convergence in N iterations of algorithm (2) with exact ABC as interface conditions. In practice, ABC and even PML truncation techniques are not perfect and the nilpotency effect is lost. In order to force it, a new operator is introduced:

$$\mathcal{T}_{SDS}(h)_{i+1,l} := \mathcal{B}_{i+1,l}(S_i(h_{i,l}, 0, 0)), \ 1 \le i \le N - 1$$

$$\mathcal{T}_{SDS}(h)_{i-1,r} := \mathcal{B}_{i-1,r}(S_i(0, h_{i,r}, 0)), \ 2 \le i \le N$$
(15)

which is by construction nilpotent of order N - 1 even with non exact ABCs as interface conditions. A fixed point method based on this reads:

$$(Id - \mathscr{T}_{SDS})(h^{n+1}) = (\mathscr{T} - \mathscr{T}_{SDS})(h^n) + G.$$
(16)

More efficiently, the operator $I - \mathcal{T}_{SDS}$ can then be used as a preconditioner in order to solve the substructured problem (8). For instance the left preconditioned system

$$(Id - \mathscr{T}_{SDS})^{-1} (Id - \mathscr{T}) (h) = (Id - \mathscr{T}_{SDS})^{-1} G$$

$$(17)$$

can be solved by a Krylov type method. A closer look at the operator $I - \mathcal{T}_{SDS}$ shows that inverting it can be made by two concurrent sweeps, hence the name double sweep.

In the sequel, algorithms other than (7) or (17) will be introduced as classical Jacobi, Gauss-Seidel, symmetric Gauss-Seidel applied to the substructured problem (8) with two different numberings of the interfaces. This will give a unified view to the methods considered in [8, 23, 27, 30, 31] and it will enable the introduction of two new algorithms in § 4.4.

3.2 Three subdomain case

For sake of clarity and to give a taste of the general case, we start with a three-domain decomposition of the whole plane \mathbb{R}^2 , see Figure 3.

3.2.1 Subdomain wise (SW) ordering

The vector h has four components which are interfaces functions living respectively on $\Gamma_{1,r}$, $\Gamma_{2,l}$, $\Gamma_{2,r}$ and $\Gamma_{3,l}$, see Figure 3:

$$h := (h_{1,r} h_{2,l} h_{2,r} h_{3,l})^T$$
.

This is the natural geometric ordering that we will refer to as the subdomain wise numbering (SW). Then for arbitrary interface conditions $\mathcal{B}_{i,l \text{ or } r}$, the sparsity pattern of the substructured problem is:

$$(I - \mathscr{T}_{SW})(h_{SW}) = \begin{pmatrix} I & X & X & 0 \\ X & I & 0 & 0 \\ 0 & 0 & I & X \\ 0 & X & X & I \end{pmatrix} \begin{pmatrix} h_{1,r} \\ h_{2,l} \\ h_{2,r} \\ h_{3,l} \end{pmatrix},$$
(18)

where X denotes a possibly non zero entry.

3.2.2 Left-Right (LR) ordering

Following [23], we consider now an ordering where left interfaces are numbered first and then the right interfaces in the reverse order. In our three subdomain case, we define;

$$h := (h_{2,l} h_{3,l} h_{2,r} h_{1,r})^T$$

For arbitrary interface conditions $\mathcal{B}_{i,l \text{ or } r}$, the sparsity pattern of the substructured operator is:

$$(I - \mathscr{T}_{LR})(h_{LR}) = \begin{pmatrix} I & 0 & 0 & X \\ X & I & X & 0 \\ \vdots & \vdots & \vdots & \vdots \\ 0 & X & I & 0 \\ X & 0 & X & I \end{pmatrix} \begin{pmatrix} h_{2,l} \\ h_{3,l} \\ \vdots \\ h_{2,r} \\ h_{1,r} \end{pmatrix}.$$
 (19)

3.2.3 Dependence of the preconditioners on ordering

Of course, systems (18) and (19) are strictly equivalent. But as a consequence of the different numberings, the approximate inverses obtained from Jacobi or Gauss-Seidel type mechanisms may not be the same. As an example, consider the Gauss-Seidel preconditioner for the SW numbering:

$$(I - \mathscr{T}_{SW,GS})(h_{SW}) := \begin{pmatrix} I & 0 & 0 & 0\\ \mathscr{T}_{(2,l)(1,r)} & I & 0 & 0\\ 0 & 0 & I & 0\\ 0 & \mathscr{T}_{(3,l)(2,l)} & \mathscr{T}_{(3,l)(2,r)} & I \end{pmatrix} \begin{pmatrix} h_{1,r} \\ h_{2,l} \\ h_{2,r} \\ h_{3,l} \end{pmatrix},$$
(20)

and the Gauss-Seidel preconditioner for the LR numbering:

$$(I - \mathscr{T}_{LR,GS})(h_{LR}) := \begin{pmatrix} I & 0 & 0 & 0 \\ \mathscr{T}_{(3,l)(2,l)} & I & 0 & 0 \\ 0 & \mathscr{T}_{(2,r)(3,l)} & I & 0 \\ \mathscr{T}_{(1,r)(2,l)} & 0 & \mathscr{T}_{(1,r)(2,r)} & I \end{pmatrix} \begin{pmatrix} h_{2,l} \\ h_{3,l} \\ 0 \\ \ddots \\ h_{2,r} \\ h_{1,r} \end{pmatrix}.$$
(21)

Even up to a reordering these two preconditioners are different since the entries of \mathscr{T} which are kept in these two approximations are not the same. In § 4, we will develop this analysis for other preconditioners with also an arbitrary number of subdomains.

Another important fact already mentioned above, see eq. (14), is that if the interface conditions are exact absorbing boundary conditions (EABC), four additional entries cancel in (18) or (19) so that \mathscr{T}_{SW} and \mathscr{T}_{LR} are actually nilpotent operators of order 2. Indeed the entry $\mathscr{T}_{(1,r)(2,l)}$ is the operator that maps $h_{2,l}$ to $\mathcal{B}_{1,r}(v_{2,l})$ where $v_{2,l}$ satisfies:

$$\mathcal{L}(v_{2,l}) = 0 \text{ in } \Omega_2,
\mathcal{B}_{2,r}(v_{2,l}) = 0 \text{ on } \Gamma_{2,r},
\mathcal{B}_{2,l}(v_{2,l}) = h_{2,l} \text{ on } \Gamma_{2,l}.$$
(22)

Since $\mathcal{B}_{2,r}$ is an EABC, $v_{2,l}$ can be seen as the restriction of a harmonic function defined on the half plane at the right of $\Gamma_{2,l}$ so that $\mathcal{B}_{1,r}(v_{2,l}) = 0$ since $\mathcal{B}_{1,r}$ is an EABC as well. That is, $\mathscr{T}_{(1,r)(2,l)} = 0$. In the same way, we can prove three other cancellations, namely: $\mathcal{T}_{(2,r)(3,l)} = 0$ and $\mathcal{T}_{(i+1,l)(i,r)} = 0$ for i = 1, 2. The only entries left are $\mathscr{T}_{(i,r)(i+1,r)}$ and $\mathscr{T}_{(i+2,l)(i+1,l)}$ for i = 1. Let us denote by \mathscr{T}_{SW}^{EABC} the operator \mathscr{T}_{SW} when the interface conditions are EABC. We have just proved that its sparsity pattern is then:

$$\mathscr{T}_{SW}^{EABC} = \begin{pmatrix} 0 & 0 & X & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & X & 0 & 0 \end{pmatrix}.$$
(23)

As for the left-right numbering, let us denote by \mathscr{T}_{LR}^{EABC} the operator \mathscr{T}_{LR} when the interface conditions are EABC so that with four entries cancelled, its sparsity pattern is then:

$$\mathscr{T}_{LR}^{EABC} = \begin{pmatrix} 0 & 0 & 0 & 0 \\ X & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & X & 0 \end{pmatrix} .$$
(24)

Another way to see that numbering impacts the preconditoners is to remark that even for EABCs, the Gauss-Seidel preconditioner for the left-right numbering is exact:

$$(I-\mathscr{T}_{LR,GS}^{EABC})=(I-\mathscr{T}_{LR}^{EABC})$$

which is not the case for the Gauss-Seidel preconditioner for the subdomain wise numbering

$$(I - \mathscr{T}_{SW,GS}^{EABC}) \neq (I - \mathscr{T}_{SW}^{EABC})$$

4 Analysis in the many subdomain case

We come back to the case of a stripwise decomposition into N subdomains as in Figures 1 or 2. As in the three subdomain case, we consider two numberings for the substructured system (8).



Figure 3: Decomposition of the plane into three subdomains

4.1 Subdomain wise (SW) numbering

The most natural numbering is to order the interface unknowns subdomain-wise (SW), i.e. one subdomain after the other. It yields the following substructured system:

4.2 Left-Right (LR) numbering

The following left-right (LR) numbering, where left interface unknowns are numbered first and then the right interface unknowns, leads to the following substructured system:

$$(I - \mathscr{T}_{LR})(H_{LR}) = \begin{bmatrix} I & 0 & 0 & -\mathscr{T}_{(2,l)(1,r)} \\ -\mathscr{T}_{(3,l)(2,l)} & \ddots & & & \ddots & \\ 0 & -\mathscr{T}_{(N,l)(N-1,l)} & I & -\mathscr{T}_{(N,l)(N-1,r)} & 0 \\ 0 & -\mathscr{T}_{(N-1,r)(N,l)} & I & 0 \\ 0 & -\mathscr{T}_{(N-1,r)(N,l)} & -\mathscr{T}_{(N-2,r)(N-1,r)} & \ddots & \\ & \ddots & & & \ddots & \\ -\mathscr{T}_{(1,r)(2,l)} & 0 & 0 & 0 & -\mathscr{T}_{(1,r)(2,r)} & I \end{bmatrix} \begin{bmatrix} h_{2,l} \\ \vdots \\ h_{N,l} \\ h_{N-1,r} \\ \vdots \\ h_{1,r} \end{bmatrix}$$
(26)

In order to highlight nilpotency, we introduce the restriction operator to the left (resp. right) interface unknowns R_l (resp. R_r) as well as four $(2N - 2) \times (2N - 2)$ submatrices of the operator \mathscr{T}_{LR} :

$$\mathcal{M}_{l} := R_{l}^{T} \mathcal{T}_{LR} R_{l} \text{ (top left)}, \ \mathcal{A}_{l} := R_{l}^{T} \mathcal{T}_{LR} R_{r} \text{ (top right)}$$
$$\mathcal{A}_{r} := R_{r}^{T} \mathcal{T}_{LR} R_{l} \text{ (bottom left)}, \ \mathcal{M}_{r} := R_{r}^{T} \mathcal{T}_{LR} R_{r} \text{ (bottom right)}$$

so that we have $I - \mathscr{T}_{LR} = I - \mathscr{M}_l - \mathscr{A}_l - \mathscr{M}_r - \mathscr{A}_r$. It is easy to check that we have the following cancellation relations:

$$\mathcal{M}_{r}^{N-1} = \mathcal{M}_{l}^{N-1} = 0; \quad \mathcal{M}_{l} \mathcal{M}_{r} = \mathcal{M}_{r} \mathcal{M}_{l} = 0; \quad \mathcal{A}_{l}^{2} = \mathcal{A}_{r}^{2} = 0$$

$$\mathcal{A}_{l} \mathcal{M}_{l} = \mathcal{A}_{r} \mathcal{M}_{r} = 0; \quad \mathcal{M}_{l} \mathcal{A}_{r} = \mathcal{M}_{r} \mathcal{A}_{l} = 0.$$
(27)

It is worth noticing that these relations come from the structure of the matrices and do not depend on the value of the entries. Let us introduce the following operators that will play a crucial role in the sequel:

$$C_r := (I - \mathcal{M}_r)^{-1} \mathcal{A}_r = (\sum_{i=0}^{N-2} \mathcal{M}_r^i) \mathcal{A}_r \text{ and } C_l := (I - \mathcal{M}_l)^{-1} \mathcal{A}_l = (\sum_{i=0}^{N-2} \mathcal{M}_l^i) \mathcal{A}_l.$$
(28)

Note that using cancellation relations (27), we have

$$C_r^2 = C_l^2 = 0. (29)$$

4.3 Jacobi, Gauss-Seidel and Symmetric Gauss-Seidel for the subdomain wise numbering

With the notations introduced above, the substructured operator (25) reads:

When considering the decomposition of matrix $I - \mathcal{T}_{SW}$ into its diagonal, lower and upper parts as $I - \mathcal{T}_{SW} = L + D + U$, its upper part U involves only entries of the matrices \mathcal{A}_r and \mathcal{M}_r and similarly, its lower part L involves only entries of the matrices \mathcal{A}_l and \mathcal{M}_l and the diagonal part D is simply the identity. *Expressed now in the LR numbering*, it is then easy to check that we have D = I, $L = -\mathcal{M}_l - \mathcal{A}_l$ and $U = -\mathcal{M}_r - \mathcal{A}_r$. Thus we introduce in a classical way several preconditioners whose formulas are given in the LR numbering:

• the Jacobi preconditioner D^{-1} :

$$M_{Jacobi}^{-1} := I \,,$$

• the left Gauss-Seidel (GS) preconditioner $(L + D)^{-1}$:

$$M_{GS}^{-1} := (I - \mathscr{M}_l - \mathscr{A}_l)^{-1},$$

• the Symmetric Gauss-Seidel (SGS) preconditioner $[(L+D)D^{-1}(D+U)]^{-1} = (D+U)^{-1}D(L+D)^{-1}$:

$$M_{SGS}^{-1} := (I - \mathscr{M}_r - \mathscr{A}_r)^{-1} (I - \mathscr{M}_l - \mathscr{A}_l)^{-1}.$$

The Jacobi preconditioner amounts to Desprès-Lions algorithm and the two other preconditioners had been introduced and studied in [23, 31]. In [23], the GS preconditioner was referred to as FDA and the SGS preconditioner as the Double Sweep (DS) algorithm.

When used as solvers, their convergence depends on the spectral radius of the error propagation operators defined as:

Definition 4.1. For a preconditioner M^{-1} of a linear operator A, we denote by R its error propagation operator as: $R := M^{-1}(M - A)$.

It was proved in [23] that the spectral radius of their error propagation operators denoted by $\rho(R)$ satisfy the following bounds: $\rho(R_{Jacobi}) \leq \rho^{1/N}$, $\rho(R_{GS}) \leq \rho^{2/N}$ and $\rho(R_{SGS}) \leq \rho$ where N is the number of subdomains and $\rho = \rho(C_r) \rho(C_l)$. The proof was written with the help of the formal language theory, see e.g. [21]. The SGS algorithm can be studied more classically since we have:

Proposition 4.2. The formula for R_{SGS} is:

$$R_{SGS} = (I + C_r) C_l \left(\mathscr{M}_r + \mathscr{A}_r \right)$$

As for the power of R_{SGS} , we have the following formula:

$$R_{SGS}^n = (I + C_r) C_l (C_r C_l)^{n-1} (\mathscr{M}_r + \mathscr{A}_r).$$

Proof. By definition, we have:

$$R_{SGS} = (Id - \mathscr{A}_r - \mathscr{M}_r)^{-1} (Id - \mathscr{M}_l - \mathscr{A}_l)^{-1} (\mathscr{M}_l + \mathscr{A}_l) (\mathscr{A}_r + \mathscr{M}_r) + \mathscr{M}_r (\mathcal{M}_r - \mathscr{M}_r)^{-1} (\mathscr{M}_l - \mathscr{M}_l)^{-1} ($$

Using the equalities: $\mathcal{M}_l \, \mathcal{A}_r = \mathcal{M}_l \, \mathcal{M}_r = 0$ and

$$Id - \mathscr{A}_r - \mathscr{M}_r = (Id - \mathscr{A}_r) (Id - \mathscr{M}_r)$$

we have:

$$R_{SGS} = \left((Id - \mathscr{M}_r)^{-1} + C_r \right) C_l \left(\mathscr{A}_r + \mathscr{M}_r \right).$$

Note that since $\mathcal{M}_r \mathcal{M}_l = \mathcal{M}_r \mathcal{A}_l = 0$, we have $\mathcal{M}_r C_l = 0$ and thus $(Id - \mathcal{M}_r)^{-1} C_l = C_l$. As for the powers of R_{SGS} , we first note that:

$$R_{SGS}^{n} = (I + C_r) C_l \left[(\mathscr{M}_r + \mathscr{A}_r) (I + C_r) C_l \right]^{n-1} (\mathscr{M}_r + \mathscr{A}_r).$$

Using cancellation relations (27) to simplify the middle term, we get:

$$(\mathscr{M}_r + \mathscr{A}_r)(I + C_r) C_l = \left[(\sum_{i=1}^{N-2} \mathscr{M}_r^i) \mathscr{A}_r + \mathscr{M}_r + \mathscr{A}_r \right] C_l = C_r C_l ,$$

where the last equality comes from $\mathcal{M}_r C_l = 0$.

4.4 Block Jacobi, Gauss-Seidel and Symmetric Gauss-Seidel for the left-right numbering

Considering system (26) as a block two by two matrix, the matrix $I - \mathscr{T}_{LR}$ can be decomposed into its block diagonal, lower and upper parts as L + D + U where $D = I - \mathscr{M}_r - \mathscr{M}_l$, $L = -\mathscr{A}_r$ and $U = -\mathscr{A}_l$. Thus we introduce in a classical way several preconditioners:

- the Block Jacobi preconditioner (BJ) D^{-1}
- the Block Gauss-Seidel (BGS) preconditioner $(L + D)^{-1}$
- the Block Symmetric Gauss-Seidel (BSGS) preconditioner $[(L+D)D^{-1}(D+U)]^{-1} = (D+U)^{-1}D(L+D)^{-1}$

Remark : The BJ preconditioner was introduced in [27, 28] for overlapping decompositions and in [30] for non overlapping decompositions. Note that a diagonal block solve consists in two independent forward substitutions on the left and right interface unknowns. So the method was called Double Sweep although it is different from the one previously introduced in [23] and named as well Double Sweep. We hope to have clarified this possible confusion.

We have thus three preconditioners with the following formulas:

$$\begin{array}{lll}
M_{BJ}^{-1} &:= & (I - \mathcal{M}_r - \mathcal{M}_l)^{-1}, \\
M_{BGS}^{-1} &:= & (I - \mathcal{M}_r - \mathcal{M}_l - \mathcal{A}_r)^{-1}, \\
M_{BSGS}^{-1} &:= & (I - \mathcal{M}_r - \mathcal{M}_l - \mathcal{A}_l)^{-1} (I - \mathcal{M}_r - \mathcal{M}_l) (I - \mathcal{M}_r - \mathcal{M}_l - \mathcal{A}_r)^{-1}.
\end{array}$$
(31)

To our knowledge, the last two algorithms have not been introduced before (except for the BGS algorithm mentioned briefly in [27] eq. (17-23), page 246). Their convergence rates nor that of BJ had not been studied before. They arise from our way to introduce the BJ algorithm which is different from the one developed in [27, 30] and that has been recalled above in § 3.1.

We have the following propositions.

Proposition 4.3. The error propagation operator R_{BJ} has the following expression:

$$R_{BJ} = C_r + C_l \,,$$

and for n even we have:

$$R_{BJ}^n = (C_r C_l)^{n/2} + (C_l C_r)^{n/2}.$$

Proof. Thanks to relations (27), we have $M_{BJ}^{-1} = Id + \sum_{i=1}^{N-2} \mathscr{M}_l^i + \sum_{i=1}^{N-2} \mathscr{M}_r^i$ and:

$$R_{BJ} = (Id + \sum_{i=1}^{N-2} \mathscr{M}_l^i + \sum_{i=1}^{N-2} \mathscr{M}_r^i)(\mathscr{A}_l + \mathscr{A}_r) = \sum_{i=0}^{N-2} \mathscr{M}_r^i \mathscr{A}_r + \sum_{i=0}^{N-2} \mathscr{M}_l^i \mathscr{A}_l = C_r + C_l \,.$$

Next using cancellation relation (29), the formula for R_{BJ}^n can easily be proved by induction.

Proposition 4.4. The error propagation operator R_{BGS} has the following expression:

$$R_{BGS} = (I + C_r) C_l ,$$

and for any integer n we have:

$$R_{BGS}^{n} = (I + C_r) C_l (C_r C_l)^{n-1}$$

Proof. Using the following formulas:

$$(I - \mathcal{M}_r - \mathcal{M}_l)^{-1} = (I - \mathcal{M}_r)^{-1} (I - \mathcal{M}_l)^{-1} = (I - \mathcal{M}_l)^{-1} (I - \mathcal{M}_r)^{-1} (I - \mathcal{M}_r - \mathcal{M}_l - \mathcal{A}_r)^{-1} = (I - \mathcal{M}_r)^{-1} (I + \mathcal{A}_r) (I - \mathcal{M}_l)^{-1} (I - \mathcal{M}_r - \mathcal{M}_l - \mathcal{A}_l)^{-1} = (I - \mathcal{M}_l)^{-1} (I + \mathcal{A}_l) (I - \mathcal{M}_r)^{-1},$$
(32)

in addition to (27), we have:

$$\begin{aligned} R_{BGS} &= M_{BGS}^{-1} \mathscr{A}_l = (I - \mathscr{M}_r)^{-1} (I + \mathscr{A}_r) (I - \mathscr{M}_l)^{-1} \mathscr{A}_l \\ &= (I - \mathscr{M}_r)^{-1} (I - \mathscr{M}_l)^{-1} \mathscr{A}_l + (I - \mathscr{M}_r)^{-1} \mathscr{A}_r (I - \mathscr{M}_l)^{-1} \mathscr{A}_l \\ &= (I - \mathscr{M}_l)^{-1} (I - \mathscr{M}_r)^{-1} \mathscr{A}_l + C_r C_l = C_l + C_r C_l \,. \end{aligned}$$

As for the *n*-th power of R_{BGS} , the formula can be proved by induction using the fact that:

$$(I + C_r) C_l (I + C_r) C_l = (I + C_r) C_l C_r C_l$$

Proposition 4.5. The error propagation operator R_{BSGS} has the following expression:

$$R_{BSGS} = (I + C_l) C_r C_l ,$$

and for any integer n we have:

$$R_{BSGS}^n = (I + C_l) \left(C_r \, C_l \right)^n$$

Proof. Using cancellation relations (27) and (32), we have:

$$\begin{aligned} R_{BSGS} &= M_{BGS}^{-1}L \, D^{-1} \, U = M_{BGS}^{-1} \mathscr{A}_r \, (I - \mathscr{M}_r)^{-1} \, (I - \mathscr{M}_l)^{-1} \, \mathscr{A}_l = M_{BGS}^{-1} \, \mathscr{A}_r \, C_l \\ &= (I - \mathscr{M}_l)^{-1} \, (I + \mathscr{A}_l) \, (I - \mathscr{M}_r)^{-1} (I - \mathscr{M}_r) (I - \mathscr{M}_l) (I - \mathscr{M}_r)^{-1} \, (I + \mathscr{A}_r) \, (I - \mathscr{M}_l)^{-1} \, \mathscr{A}_r \, C_l \\ &= (I - \mathscr{M}_l)^{-1} \, (I + \mathscr{A}_l - \mathscr{M}_l) (I - \mathscr{M}_r)^{-1} \, (I + \mathscr{A}_r) \, (I - \mathscr{M}_l)^{-1} \, \mathscr{A}_r \, C_l \\ &= (I + C_l) \, (I - \mathscr{M}_r)^{-1} \, (I + \mathscr{A}_r) \, \mathscr{A}_r \, C_l = (I + C_l) \, C_r \, C_l \, . \end{aligned}$$

As for the last formula of the proposition, it comes from the nullity of both C_r^2 and C_l^2 .

Remark 4.1. It is clear from the above results that the operators C_r and C_l (see eq. (28)) are a key measure of the efficiency of these algorithms. If the left interface conditions are exact absorbing conditions, operator $\mathscr{A}_r = 0$, see equation (22) and thus $C_r = 0$ as well. Similarly if right interface conditions are exact absorbing conditions, operator $C_l = 0$. More generally, the norms of C_r and C_l are proportional to that of $\mathscr{A}_r = 0$ and $\mathscr{A}_l = 0$. Thus as expected the more absorbing the interface conditions are, the better the convergence is. Another parameter is the number of subdomains since for given interface conditions, as the number of subdomains increases, the norm of C_r and C_l will grow. It echoes what was noticed in [27] at the end of § 4:

"With the Robin transmission conditions the iteration numbers grow roughly linearly in N_x , or as $N^{1/2}$ in 2-D."

In this respect, note that R_{BSGS} satisfies

$$(I+C_l)^{-1} R_{BSGS} (I+C_l) = C_r C_l$$

so that from all algorithms listed in Table 1 it has the most favorable amplification error operator iff the norm of $C_r C_l$ is smaller than one. In a recent article [16], these norms have been estimated in the context of the study of the convergence rate of the Jacobi preconditioner.

4.5 Theoretical comparison between the SGS and BJ preconditioners

We devote a paragraph to this comparison since these algorithms have been misleadingly coined the same name Double Sweep algorithm in the articles in which they had been introduced, in [23, 31] for SGS and in [27, 30] for BJ. From Propositions 4.2 and 4.3, we have for n even:

$$R_{B,I}^n = (C_r C_l)^{n/2} + (C_l C_r)^{n/2},$$

and for any \boldsymbol{n}

$$R_{SGS}^n = (I + C_r) C_l (C_r C_l)^{n-1} (\mathscr{M}_r + \mathscr{A}_r).$$

Let us denote by

the spectral radius of
$$C_r C_l$$
,

which is the same as that of $C_l C_r$. We have for any operator norm $\|\cdot\|$ on the matrices the following estimates:

$$\|R_{BJ}^{n}\| \le \|(C_{r} C_{l})^{n/2}\| + \|(C_{l} C_{r})^{n/2}\| \le \|(C_{r} C_{l})^{n/2}\| + \|C_{l}\|\|(C_{r} C_{l})^{n/2-1}\|\|C_{l}\|,$$
(34)

so that taking the *n*-th square root, we get the following estimate for the spectral radius of R_{BJ} :

ſ

$$\rho(R_{BJ}) \le \rho^{1/2} \,. \tag{35}$$

(33)

As for the SGS algorithm, we get:

$$\|R_{SGS}^{n}\| \le \|I + C_{r}\| \, \|C_{l}\| \, \|\mathcal{M}_{r} + \mathcal{A}_{r}\| \, \|(C_{r} \, C_{l})^{n-1}\| \, . \tag{36}$$

We thus get the following estimate for the spectral radius of R_{SGS} :

$$\rho(R_{SGS}) \le \rho \,. \tag{37}$$

Thus when $\rho < 1$, an advantage of SGS over BJ is the square root of ρ in the ratio of the convergence rates. This is coherent with the factor two in the iteration counts in favour of SGS observed in most tables of the numerical section 5 when PMLs are interface conditions. Nevertheless, note that at the expense of doubling the number of cores used in the application of the BJ preconditioner, this difference is nullified in terms of elapsed time. Indeed, during one iteration of the SGS algorithm, only one subdomain is active at a time in the order $1 \rightarrow 2 \rightarrow ... \rightarrow N$ and then in the reverse order $N \rightarrow N - 1 \rightarrow ... \rightarrow 1$. Whereas during one iteration of BJ, two subdomains are active at a time in the order $1 \rightarrow 2 \rightarrow ... \rightarrow N$ for one subdomain and concurrently in the reverse order $N \rightarrow N - 1 \rightarrow ... \rightarrow 1$ for the other one. Thus, the elapsed time of one iteration of SGS is twice as much as one iteration of BJ using two cores. This compensates for the higher iteration counts of BJ compared to SGS when the number of cores is not limited.

Ref.	Abbr.	Linear Algebra	Definition	Ampl. Error	Radius
Després (1991)	Jacobi	Jacobi-SW	Ι	$\mathscr{M}_r + \mathscr{M}_l + \mathscr{A}_r + \mathscr{A}_l$	$ ho^{1/N}$
Nataf-Nier (1997)	GS	Gauss-Seidel-SW	$(I - \mathscr{M}_l - \mathscr{A}_l)^{-1}$	$(I+C_l)(\mathscr{M}_r+\mathscr{A}_r)$	$\rho^{2/N}$
{ Nataf-Nier (1997) Vion-Geuzaine(2018)	SGS	Symm. Gauss-Seidel-SW	$(I - \mathscr{M}_r - \mathscr{A}_r)^{-1} \ imes (I - \mathscr{M}_l - \mathscr{A}_l)^{-1}$	$(I+C_r) C_l \left(\mathscr{M}_r + \mathscr{A}_r\right)$	ρ
{ Stolk (2013) Vion-Geuzaine (2014)	BJ	Block Jacobi-LR	$(I - \mathcal{M}_r - \mathcal{M}_l)^{-1}$	$C_r + C_l$	$ ho^{1/2}$
Here (2022)	BGS	Block GS-LR	$(I - \mathcal{M}_r - \mathcal{M}_l - \mathcal{A}_r)^{-1}$	$(I+C_r) C_l$	ρ
Here (2022)	BSGS	Symm. Block GS-LR	$ \begin{vmatrix} (I - \mathcal{M}_r - \mathcal{M}_l - \mathcal{A}_l)^{-1} \\ \times (I - \mathcal{M}_r - \mathcal{M}_l) \\ \times (I - \mathcal{M}_r - \mathcal{M}_l - \mathcal{A}_r)^{-1} \end{vmatrix} $	$(I+C_l) C_r C_l$	ρ

Table 1: Algorithms and their convergence properties. SW means subdomain wise numbering and LR means left right numbering

4.6 Volumic preconditioners

The substructured algorithms demand exact local solves and data structures adapted to substructured algorithms. In order to allow for the use of fast approximate solvers in subdomains and of volumic data structures, an extension to the volumic case

is defined below. For the SGS algorithm, the volumic formulation is simply given by equations (4)-(5) followed by the use of a partition of unity in order to iterate on functions which are uniquely defined in the overlaps. For this, we first introduce for $1 \le i \le N$ partition of unity functions $\chi_i : \Omega_i \to \mathbb{R}^+$ and for $v_i : \Omega_i \to \mathbb{C}$, $E_i(v_i)$ denotes the extension by zero of v_i to Ω so that for any function $v : \Omega \to \mathbb{C}$, we have:

$$v = \sum_{i=1}^{N} E_i(\chi_i v_{|\Omega_i}).$$

We introduce the operator MC (MakeCoherent) that maps a collection of local functions $(v_i)_{1 \le i \le N}$ to a global function defined as follows:

$$MC((v_i)_{1 \le i \le N}) := \sum_{i=1}^{N} E_i(\chi_i \, v_i) \,. \tag{38}$$

Let f be a source term and h be interface sources, it is then natural to introduce the following linear parallel reconstruction algorithm Vol:

$$Vol(h, f) := MC((S_i(h_{i,l}, h_{i,r}, f))_{1 \le i \le N}.$$
(39)

Let u be the solution to the original problem (1), we have:

$$u = \operatorname{Vol}((I - \mathscr{T})^{-1}G(f), f).$$
(40)

Since a substructured preconditioner M^{-1} may be seen as an approximate inverse to $(I - \mathscr{T})$, it is natural to define the related volumic preconditioner as:

Definition 4.6. Let M^{-1} be a substructured preconditioner for problem (8), the related volumic preconditioner to problem (1) M_{vol}^{-1} is defined as follows:

$$M_{vol}^{-1}(f) := Vol(M^{-1}G(f), f) = Vol(M^{-1}G(f), 0) + Vol(0, f).$$
(41)

Note that if the surface variant of the preconditioner is exact as it happens for some if the interface conditions are exact absorbing boundary conditions, this property will be inherited by its volumic counterpart.

As an example, we detail below after some rewriting the action of the volumic preconditioner associated to the surface variant of the BJ preconditioner.

Volumic BJ Preconditioner

Left to right sweep	Right to left sweep
Subdomain 1	Subdomain N
$w_1 \leftarrow S_1(0,0,f)$	$w_N \leftarrow S_N(0,0,f)$
Subdomain 2	Subdomain $N-1$
$G_{2,l}(f) \leftarrow \mathcal{B}_{2,l}(w_1)$	$G_{N-1,r}(f) \leftarrow \mathcal{B}_{N-1,r}(w_N)$
$h_{2,l} \leftarrow G_{2,l}(f)$	$h_{N-1,r} \leftarrow G_{N-1,r}(f)$
$v_{2,l} \leftarrow S_2(h_{2,l}, 0, 0)$	$v_{N-1,r} \leftarrow S_{N-1}(0, h_{N-1,r}, 0)$
$w_2 \leftarrow S_2(0,0,f)$	$w_{N-1} \leftarrow S_{N-1}(0,0,f)$
Subdomain 3	Subdomain $N-2$
$G_{3,l}(f) \leftarrow \mathcal{B}_{3,l}(w_2)$	$G_{N-2,r}(f) \leftarrow \mathcal{B}_{N-2,r}(w_{N-1})$
$h_{3,l} \leftarrow G_{3,l}(f) + \mathcal{B}_{3,l}(v_{2,l})$	$h_{N-2,r} \leftarrow G_{N-2,r}(f) + \mathcal{B}_{N-2,r}(v_{N-1,r})$
$v_{3,l} \leftarrow S_3(h_{3,l}, 0, 0)$	$v_{N-2,r} \leftarrow S_{N-2}(0, h_{N-2,r}, 0)$
$w_3 \leftarrow S_3(0,0,f)$	$w_{N-2} \leftarrow S_{N-2}(0,0,f)$
Subdomain 4	Subdomain $N-3$
$G_{4,l}(f) \leftarrow \mathcal{B}_{4,l}(w_3)$	$G_{N-3,r}(f) \leftarrow \mathcal{B}_{N-3,r}(w_{N-2})$
$h_{4,l} \leftarrow G_{4,l}(f) + \mathcal{B}_{4,l}(v_{3,l})$	$h_{N-3,r} \leftarrow G_{N-3,r}(f) + \mathcal{B}_{N-3,r}(v_{N-2,r})$
÷	÷
Subdomain N	Subdomain 1
$w_N \leftarrow S_N(0,0,f)$	$w_1 \leftarrow S_1(0,0,f)$
$v_{N,l} \leftarrow S_N(h_{N,l}, 0, 0)$	$v_{1,r} \leftarrow S_1(h_{1,r}, 0, 0)$

Then we perform local summations:

 $u_i := w_i + v_{i,l} + v_{i,r}, \quad 2 \le i \le N - 1$

and

 $u_1 := w_1 + v_{1,r}$ and $u_N := w_N + v_{N,l}$.

Finally the action of the preconditioner M_{BJvol}^{-1} reads:

$$M_{BJ,vol}^{-1}(f) := \sum_{i=1}^{N} E_i(\chi_i(u_i)).$$
(42)

5 Numerical results

In this section, we present numerical results for the volumic variants of the preconditioners of Table 1 used with the GM-RES algorithm [25] with two different relative tolerances, $TOL=10^{-6}$ and $TOL=10^{-3}$. The Helmholtz equation is discretized with P1 or P2 finite elements using the FreeFem++ domain specific language [18]. The following test cases are shown: homogeneous waveguide (§ 5.1 and 5.2), open cavity (§ 5.3), Marmousi (§ 5.4) and 3D Overthrust (§ 5.5).

We also compare the use of zeroth-order ABC with that of Perfectly Matched Layers (PML) as interface conditions for the various sweeping algorithms. It is worth noting that in order to avoid transmitting attenuated solutions between subdomains, the width of the overlap region needs to be taken larger that the width of the PML. In the following numerical results, the overlap δ is equal to 4 mesh elements and the width of the PML is 2. The PML is defined following [3]. As we can see, using PML interface conditions, iteration counts are slightly better for all cases and much better for the Marmousi test case. Also note that as the number of subdomains increases, the superiority of PML becomes more pronounced.

When comparing the various algorithms, we see that the iteration counts are qualitatively in agreement with the spectral radius estimates summarized in Table 1 but only when using PML as interface conditions. For the Jacobi method, the increase is linear w.r.t. to the number of subdomains. The sweeping methods have iteration counts that increase sublinearly with the number of subdomains. When using PML, the Block Jacobi (BJ) methods needs twice as many iterations as the other sweeping methods. We also see that the two new preconditioners BGS and BSGS offer no iteration count improvement over the SGS preconditioner although they are more expensive.

5.1 Homogeneous waveguide

First, we consider the homogeneous waveguide test case with a layered decomposition into N subdomains. More specifically, we consider a rectangular geometry ($\Omega = [0, N] \times [0, 1]$) made of a homogeneous medium. On the upper and lower sides of the waveguide, we impose homogeneous Dirichlet conditions (cf. black lines in Figure 4). In addition, we perform a multimode excitation on the left side and we impose an absorbing boundary condition on the right side. The global problem is written as

$$\begin{cases} \left(-k^{2} - \Delta\right) u = f \text{ in } \Omega \\ \left(\partial_{\vec{n}} + Ik\right) u = 0 \text{ on } \{x = N\} \times [0, 1] \\ \left(\partial_{\vec{n}} + Ik\right) u = u_{g} \text{ on } \{x = 0\} \times [0, 1] \\ u = 0 \text{ on } [0, N] \times \{y = 0, y = 1\} \end{cases}$$

$$\tag{43}$$

where $u_g = e^{-120(y-0.5)^2} sin(\pi y)$.



Figure 4: Homogeneous waveguide ($k = 20\pi$)

We considered two values for the wave number: k = 20 and $k = 20\pi$ and results are given in Tables 2 and 4 for the zeroth order interface conditions and in Tables 3 and 5 for PML interface conditions. The iteration counts of the various sweeping algorithms are in agreement with the spectral radius estimates of Table 1. In this simple case, the superiority of PML over zeroth order ABC is not so significant except for large number of subdomains.

5.2 Influence of the overlap

We have also tested the effect of the width of the overlap on the convergence in the case of the homogeneous waveguide with zeroth-order ABC. In this case, there is little if no effect of the overlap as shown in Table 6 where the overlap varies from 2 mesh sizes up to 16 mesh sizes.

N	Jacobi	BJ	BGS	BSGS	SGS
5	29 (18)	9 (5)	5 (3)	5 (3)	5 (3)
10	62 (39)	12 (7)	7 (4)	6 (4)	6 (4)
20	135 (81)	18 (10)	10 (6)	9 (6)	9 (6)
40	283 (163)	26 (12)	14 (8)	14 (7)	14 (8)
80	744 (329)	42 (20)	23 (12)	22 (12)	23 (12)

Table 2: Volumic preconditioner, homogeneous waveguide with zeroth-order ABC, k = 20, $\delta = 4h$, TOL= $10^{-6}(10^{-3})$, nppwl = 24, P1

N	Jacobi	BJ	BGS	BSGS	SGS
5	33 (21)	9 (5)	5 (2)	4 (2)	5 (3)
10	71 (40)	11 (6)	6 (3)	5 (3)	6 (3)
20	150 (80)	13 (7)	7 (4)	6 (3)	7 (4)
40	293 (143)	17 (9)	9 (5)	8 (5)	9 (5)
80	690 (276)	22 (14)	12 (7)	12 (7)	12 (7)

Table 3: Volumic preconditioner with PML interface conditions, homogeneous waveguide, k = 20, $\delta = 4h$, TOL= $10^{-6}(10^{-3})$, nppwl = 24, P1

N	Jacobi	BJ	BGS	BSGS	SGS
5	35	15	10	9	9
10	72	22	14	13	14
20	150	34	22	23	23
40	335	58	38	45	42

Table 4: Volumic preconditioner, homogeneous waveguide with zeroth-order ABC, $k = 20\pi$, $\delta = 4h$, TOL= 10^{-6} , nppwl = 24, P1

N	Jacobi	BJ	BGS	BSGS	SGS
5	38	10	6	5	6
10	78	11	7	5	6
20	162	17	9	7	9
40	304	23	12	10	12

Table 5: Volumic preconditioner with PML interface conditions, homogeneous waveguide, $k = 20\pi$, $\delta = 4h$, TOL= 10^{-6} , nppwl = 24, P1

δ	Jacobi	BJ	BGS	BSGS	SGS
2	136	19	11	10	10
4	135	18	10	9	9
8	137	17	9	8	9
16	149	21	11	11	11

Table 6: Volumic preconditioner, homogeneous waveguide with zeroth-order ABC, k = 20, δ varies, TOL= 10^{-6} , nppwl = 24, P1

5.3 Open cavity test

Same as before, the domain is rectangular with an homogeneous medium and its length increases with the number of subdomains. The open cavity test is challenging due to the homogeneous Dirichlet conditions imposed on three sides (cf. black lines in Figure 5). In addition, we perform an excitation on the left side. The Dirichlet conditions create rebounds leading to an increase in the number of reflections, this phenomenon is exacerbated for high-frequency regimes. The global problem can be written as

$$\begin{cases} \left(-k^{2}-\Delta\right) u = f \text{ in } \Omega\\ \left(\partial_{\vec{n}}+Ik\right) u = g \text{ on } \Gamma\\ u = 0 \text{ on } \partial\Omega \backslash \Gamma \end{cases}$$

$$\tag{44}$$

where $\Gamma := \{x = 0\} \times [0, 1]$ and $g = exp^{-ik(x\cos(\theta) + y\sin(\theta))}$, $\theta = \frac{\pi}{8}$. It corresponds to an incident plane wave propagating at an angle $\frac{\pi}{8}$ w.r.t. to the horizontal direction, see Fig. 5. This creates numerous reflections on the lateral boundaries of the open cavity. Here as well, there is roughly a factor two in the iteration counts in favour of the SGS, BGS and BSGS algorithms compared to BJ algorithm.

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Figure 5: Open cavity solution ($k = 20\pi$)

N	Jacobi	BJ	BGS	BSGS	SGS
5	44 (29)	12 (8)	7 (4)	6 (4)	7 (5)
10	98 (62)	16(11)	9 (6)	8 (5)	8 (6)
20	202 (142)	21 (14)	11 (8)	10 (8)	11 (8)
40	478 (255)	30 (22)	16 (12)	15 (11)	16 (12)

Table 7: Volumic preconditioner with zeroth-order ABC, open cavity, k = 20, $\delta = 4h$, TOL= $10^{-6}(10^{-3})$, nppwl = 24, P1

N	Jacobi	BJ	BGS	BSGS	SGS
5	49 (33)	12 (8)	6 (4)	6 (4)	7 (4)
10	106 (71)	14 (10)	7 (5)	7 (5)	7 (5)
20	216 (145)	16(11)	9 (6)	8 (5)	8 (6)
40	427 (260)	19 (12)	10 (7)	10 (6)	10 (7)

Table 8: Volumic preconditioner with PML interface conditions, open cavity, k = 20, $\delta = 4h$, TOL= $10^{-6}(10^{-3})$, nppwl = 24, P1

N	Jacobi	BJ	BGS	BSGS	SGS
5	67 (37)	23 (14)	14 (7)	13 (8)	14 (8)
10	137 (77)	32 (19)	19 (11)	19 (12)	20 (13)
20	298 (152)	50 (31)	29 (21)	31 (23)	32 (23)
40	789 (351)	77 (52)	48 (37)	53 (44)	53 (41)

Table 9: Volumic preconditioner with zeroth-order ABC, open cavity, $k = 20\pi$, $\delta = 4h$, TOL= $10^{-6}(10^{-3})$, nppwl = 24, P1

5.4 Marmousi

Here we consider the Marmousi benchmark [4]. The velocity profile of the model is depicted in Figure 6. The domain is of size 9.2 km \times 3 km. A Neumann boundary condition is imposed at the top boundary, and PMLs are used on the other three boundaries. The source is located at the top. The problem is discretized with P2 elements on a regular mesh with 8 points per wavelength for the reference wavelength corresponding to c = 2. Figure 7 shows the real part of the acoustic field at 100

N	Jacobi	BJ	BGS	BSGS	SGS
5	72 (42)	18 (11)	9 (6)	9 (5)	10 (6)
10	153 (84)	19 (11)	10 (6)	9 (5)	11 (6)
20	321 (162)	26 (16)	14 (9)	12 (7)	14 (9)
40	638 (313)	34 (21)	18 (11)	16 (10)	19 (11)

Table 10: Volumic preconditioner with PML interface conditions, open cavity, $k = 20\pi$, $\delta = 4h$, TOL= $10^{-6}(10^{-3})$, nppwl = 24, P1

Hz frequency. We present results for 25 Hz, 50 Hz and 100 Hz frequencies using the volumic preconditioner. The number of degrees of freedom grows from 1.1 million for f = 25 Hz to 4.5 million for f = 50 Hz and 17.8 million for f = 100Hz. Table 11 shows numerical results for zeroth order ABC interface conditions and Table 12 for PML. Interestingly we see that for ABC results are quite bad with high iteration counts and dependence on both the number of subdomains and the wave number. Whereas with PML interface conditions, both the iteration counts and their behaviour with respect to the number of subdomains and wavenumber is similar to the previous test cases on waveguides and open cavity. This very bad behaviour of zeroth order ABC at higher frequency and subdomain count can be explained by the variability of the coefficients.



Figure 6: Velocity model of the Marmousi test case



Figure 7: Real part of the solution for f = 100 Hz for the Marmousi test case

5.5 3D Overthrust

Here we consider the 3D Overthrust acoustic benchmark. The velocity profile of the model is depicted in Figure 6. The domain is of size 20 km \times 20 km \times 4.65 km. A Neumann boundary condition is imposed at the top boundary, and PMLs are used on the other five boundaries. The source is located at the top, at (2.5 km, 2.5 km). The problem is discretized with P1 elements on a regular mesh with 10 points per wavelength for the reference wavelength corresponding to c = 2. Figure 9 shows the real part of the acoustic field at 4 Hz frequency.

Tables 13 and 14 present results for 1 Hz, 2 Hz and 4 Hz frequencies using the volumic preconditioner with zeroth order ABC and PML interface conditions. The number of degrees of freedom grows from 0.33 million for f = 1 Hz to 16.2 million

	25 Hz							50 Hz		
N	Jacobi	BJ	BGS	BSGS	SGS	Jacobi	BJ	BGS	BSGS	SGS
3	22 (12)	11 (6)	10 (5)	7 (4)	7 (4)					
7	46 (25)	17 (9)	13 (7)	10 (6)	11 (6)	49 (25)	19 (9)	16 (8)	13 (8)	14 (8)
14	94 (51)	25 (14)	18 (10)	19 (12)	20(12)	98 (51)	28 (15)	22 (12)	21 (13)	21 (11)
28	185 (100)	41 (24)	27 (16)	30 (20)	30 (19)	195 (101)	49 (26)	38 (21)	60 (48)	47 (32)
56	382 (220)	98 (69)	65 (47)	101 (91)	70 (49)	426 (222)	123(81)	90 (61)	X (192)	143 (116)
112						1505 (690)	X (324)	X(X)	X(X)	X (X)

			100 Hz		
N	Jacobi	BJ	BGS	BSGS	SGS
56	476 (216)	147 (87)	120 (77)	X (X)	208 (171)
112	1701 (691)	X (X)	X (375)	X(X)	X (X)
224	X (X)	X (X)	X(X)	X(X)	X (X)

Table 11: Volumic preconditioner with zeroth-order ABC, Marmousi test case, $\delta = 4h$, TOL= $10^{-6}(10^{-3})$, nppwl = 8, P2. X means that the algorithm did not reach the convergence criterion in 2000 iterations for Jacobi, or in 400 iterations for the sweeping methods.

	25 Hz							50 Hz		
N	Jacobi	BJ	BGS	BSGS	SGS	Jacobi	BJ	BGS	BSGS	SGS
3	14 (8)	7 (4)	6 (3)	4 (2)	5 (3)					
7	33 (17)	10 (5)	8 (4)	5 (3)	6 (4)	34 (18)	10 (6)	9 (4)	6 (3)	6 (4)
14	64 (35)	11 (6)	8 (5)	6 (4)	7 (4)	69 (35)	12 (6)	10 (5)	7 (4)	8 (4)
28	126 (66)	13 (7)	9 (5)	7 (4)	8 (4)	133 (66)	14 (7)	11 (6)	7 (4)	8 (5)
56	247 (124)	18 (11)	12(7)	11 (6)	12 (6)	260 (125)	18 (9)	13 (7)	10 (5)	11 (5)
112						531 (240)	32 (17)	22 (12)	20 (11)	21 (11)

	100 Hz									
N	Jacobi	BJ	BGS	BSGS	SGS					
56	242 (103)	18 (9)	14 (6)	10 (5)	11 (5)					
112	483 (198)	29 (13)	20 (9)	15 (7)	18 (8)					
224	1069 (417)	66 (32)	42 (20)	62 (44)	38 (18)					

Table 12: Volumic preconditioner with PML interface conditions, Marmousi test case, $\delta = 4h$, TOL= $10^{-6}(10^{-3})$, nppwl = 8, P2

for f = 4 Hz. We observe similar trends as for the Marmousi test case, although the zeroth order ABC case is not so bad, as the frequency and number of wavelengths in one direction is not as high for this 3D test case. Moreover, we can see that the iteration count is much more dependent on the number of subdomains than on the frequency.



Figure 8: Velocity model of the 3D Overthrust benchmark



Figure 9: Real part of the solution for f = 4 Hz for the 3D Overthrust benchmark

6 Conclusion

We have introduced a unified framework for several sweeping algorithms which eases their derivation and theoretical studies, see Table 1. We have focused on two double sweep methods: SGS of [23, 31] and BJ of [27, 30]. We conclude both from their theoretical convergence rates and from our numerical results that the algorithm SGS has a better behavior than the BJ algorithm. The numerical experiments also illustrate the superiority of PML as interface conditions compared to zeroth order ABCs, especially in the variable coefficient case. Although these algorithms are more easily derived in their substructured form, they lend themselves to volumic variants that allow for inexact solves in the subdomains, see § 4.6. In this form, these methods should have then a behavior similar to the sweeping algorithm of [13].

	1 Hz							2 Hz		
N	Jacobi	BJ	BGS	BSGS	SGS	Jacobi	BJ	BGS	BSGS	SGS
3	14 (7)	7 (4)	6 (3)	4 (2)	4 (2)					
7	32 (17)	9 (5)	7 (4)	5 (3)	6 (3)	34(17)	10 (5)	8 (4)	5 (3)	6 (3)
14	94 (45)	18 (9)	10 (5)	9 (5)	9 (5)	68 (34)	15 (8)	10 (5)	8 (4)	9 (4)
28						163 (80)	30 (16)	16 (9)	17 (10)	14 (8)

			4 Hz		
Ν	Jacobi	BJ	BGS	BSGS	SGS
28	162 (80)	32 (17)	19 (10)	24 (15)	17 (9)
56	365 (187)	87 (48)	42 (22)	168 (133)	36 (20)

Table 13: Volumic preconditioner with zeroth-order ABC, Overthrust test case, $\delta = 4h$, TOL= $10^{-6}(10^{-3})$, nppwl = 10, P1

			1 Hz			2 Hz				
Ν	Jacobi	BJ	BGS	BSGS	SGS	Jacobi	BJ	BGS	BSGS	SGS
3	11 (7)	5 (3)	4 (2)	3 (2)	3 (2)					
7	28 (15)	7 (4)	5 (3)	4 (2)	5 (3)	30(16)	8 (4)	6 (3)	4 (2)	5 (3)
14	59 (31)	9 (5)	6 (4)	5 (3)	6 (3)	63 (31)	11 (6)	7 (4)	6 (3)	7 (3)
28						123 (61)	16 (8)	10 (5)	7 (4)	9 (5)

			4 Hz		
Ν	Jacobi	BJ	BGS	BSGS	SGS
28	121 (63)	18 (9)	11 (6)	10 (6)	11 (6)
56	224 (122)	28 (15)	18 (9)	19 (12)	16 (9)

Table 14: Volumic preconditioner with PML interface conditions, Overthrust test case, $\delta = 4h$, TOL= $10^{-6}(10^{-3})$, nppwl = 10, P1

Also, an intrinsic problem with double sweep methods is that due to the sequentiality of the algorithm, subdomains are idle most of the time. To be more precise, in SGS, only one process is active at a time whereas in BJ, two processes are active at a time (one for the left sweep and one for the right sweep performed concurrently). A solution to overcome this is to introduce a pipelining technique that can be applied to multiple right-hand sides problems to improve parallelism and achieve significant speed-ups, see [27, 30].

We have considered stripwise decompositions and it would be interesting to see if the present analysis can be extended to L-sweeps preconditioners (see e.g. [29, 7] and references therein) used for checkerboard type decompositions.

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