

**Schwarz Methods: To Symmetrize
or Not to Symmetrize**

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SCHWARZ METHODS: TO SYMMETRIZE OR NOT TO SYMMETRIZE*

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Abstract. A preconditioning theory for Schwarz methods is presented. The theory establishes sufficient conditions for multiplicative and additive Schwarz algorithms to yield self-adjoint positive definite preconditioners. It allows for the analysis and use of non-variational and non-convergent linear methods as preconditioners for conjugate gradient methods, and it is applied to domain decomposition and multigrid. It is illustrated why symmetrizing may be a bad idea for linear methods. It is conjectured that enforcing minimal symmetry achieves the best results when combined with conjugate gradient acceleration. Also, it is shown that absence of symmetry in the linear preconditioner is advantageous when the linear method is accelerated by using the Bi-CGstab method. Numerical examples are presented for two test problems which illustrate the theory and conjectures.

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1. Introduction. Domain decomposition (DD) and multigrid (MG) methods have been studied extensively in recent years, both from a theoretical and numerical point of view. DD methods were first proposed in 1869 by H. A. Schwarz as a theoretical tool in the study of elliptic problems on non-rectangular domains [20]. More recently, DD methods have been reexamined for use as practical computational tools in the (parallel) solution of general elliptic equations on complex domains [8]. MG methods were discovered much more recently [9]. They have been extensively developed both theoretically and practically since the late seventies [5, 10], and they have proven to be extremely efficient for solving very broad classes of partial differential equations. Recent insights in the product nature of certain MG methods have led to a unified theory of MG and DD methods, collectively referred to as *Schwarz methods* [4, 25].

In this paper, we consider additive and multiplicative Schwarz methods and their acceleration with Krylov methods, for the numerical solution of self-adjoint positive definite (SPD) operator equations arising from the discretization of elliptic partial differential equations. The standard theory of conjugate gradient acceleration of linear methods requires that a certain operator associated with the linear method – the preconditioner – be symmetric and positive definite. Often, however, as in the case of Schwarz-based preconditioners, the preconditioner is known only implicitly, and symmetry and positive definiteness are not easily verified. Here, we try to construct natural sets of sufficient conditions that are easily verified and do not require the explicit formulation of the preconditioner. More precisely, we derive conditions for the constituent components of MG and DD algorithms (smoother, subdomain solver, transfer operators, etc.), that guarantee symmetry and positive definiteness of the preconditioning

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operator which is (explicitly or implicitly) defined by the resulting Schwarz method.

We examine the implications of these conditions for various formulations of the standard DD and MG algorithms. The theory we develop helps to explain the often observed behavior of a poor or even divergent MG or DD method which becomes an excellent preconditioner when accelerated by a conjugate gradient method. We also investigate the role of symmetry in linear methods and preconditioners. Both analysis and numerical evidence suggest that linear methods should not be symmetrized when used alone, and only minimally symmetrized when accelerated by conjugate gradients, in order to achieve the best possible convergence results. In fact, the best results are often obtained when a very nonsymmetric linear iteration is used in combination with a nonsymmetric system solver such as Bi-CGstab, even though the original problem is SPD.

The outline of the paper is as follows. We begin in §2 by reviewing basic linear methods for SPD linear operator equations, and examine Krylov acceleration strategies. In §3 and §4, we analyze multiplicative and additive Schwarz preconditioners. We develop a theory that establishes sufficient conditions for the multiplicative and additive algorithms to yield SPD preconditioners. This theory is used to establish sufficient conditions for multiplicative and additive DD and MG methods, and allows for analysis of non-variational and even non-convergent linear methods as preconditioners. A simple lemma, given in §5, illustrates why symmetrizing may be a bad idea for linear methods. In §6, results of numerical experiments obtained with finite-element-based DD and MG methods applied to some non-trivial test problems are reported.

2. Krylov acceleration of linear iterative methods. In this section, we review some background material on self-adjoint linear operators, linear methods, and conjugate gradient acceleration. A more thorough reviews can be found in [11, 16].

2.1. Background material, terminology and notation. Let \mathcal{H} be a real finite-dimensional Hilbert space equipped with the inner-product (\cdot, \cdot) inducing the norm $\|\cdot\| = (\cdot, \cdot)^{1/2}$. \mathcal{H} can be thought of as, for example, the Euclidean space \mathbb{R}^n , or as an appropriate finite element space.

The *adjoint* of a linear operator $A \in \mathbf{L}(\mathcal{H}, \mathcal{H})$ with respect to (\cdot, \cdot) is the unique operator A^T satisfying $(Au, v) = (u, A^T v)$, $\forall u, v \in \mathcal{H}$. An operator A is called *self-adjoint* or *symmetric* if $A = A^T$; a self-adjoint operator A is called *positive definite* or simply *positive*, if $(Au, u) > 0$, $\forall u \in \mathcal{H}$, $u \neq 0$. If A is self-adjoint positive definite (SPD) with respect to (\cdot, \cdot) , then the bilinear form (Au, v) defines another inner-product on \mathcal{H} , which we denote as $(\cdot, \cdot)_A$. It induces the norm $\|\cdot\|_A = (\cdot, \cdot)_A^{1/2}$.

The adjoint of an operator M with respect to $(\cdot, \cdot)_A$, the *A-adjoint*, is the unique operator M^* satisfying $(Mu, v)_A = (u, M^*v)_A$, $\forall u, v \in \mathcal{H}$. From this definition it follows that

$$(1) \quad M^* = A^{-1} M^T A .$$

An operator M is called *A-self-adjoint* if $M = M^*$, and *A-positive* if $(Mu, u)_A > 0$, $\forall u \in \mathcal{H}$, $u \neq 0$.

If $N \in \mathbf{L}(\mathcal{H}_1, \mathcal{H}_2)$, then the adjoint satisfies $N^T \in \mathbf{L}(\mathcal{H}_2, \mathcal{H}_1)$, and relates the inner-products in \mathcal{H}_1 and \mathcal{H}_2 as follows:

$$(Nu, v)_{\mathcal{H}_2} = (u, N^T v)_{\mathcal{H}_1}, \quad \forall u \in \mathcal{H}_1, \quad \forall v \in \mathcal{H}_2.$$

Since it is usually clear from the arguments which inner-product is involved, we shall drop the subscripts on inner-products (and norms) throughout the paper, except when necessary to avoid confusion.

We denote the spectrum of an operator M as $\sigma(M)$. The spectral theory for self-adjoint linear operators states that the eigenvalues of the self-adjoint operator M are real and lie in the closed interval $[\lambda_{\min}(M), \lambda_{\max}(M)]$ defined by the Raleigh quotients:

$$(2) \quad \lambda_{\min}(M) = \min_{u \neq 0} \frac{(Mu, u)}{(u, u)}, \quad \lambda_{\max}(M) = \max_{u \neq 0} \frac{(Mu, u)}{(u, u)}.$$

Similarly, if an operator M is A -self-adjoint, then its eigenvalues are real and lie in the interval defined by the Raleigh quotients generated by the A -inner-product. A well-known property is that if M is self-adjoint, then the spectral radius of M , denoted as $\rho(M)$, satisfies $\rho(M) = \|M\|$. This property can also be shown to hold in the A -norm for A -self-adjoint operators (or, more generally, for A -normal operators [1]).

LEMMA 2.1. *If A is SPD and M is A -self-adjoint, then $\rho(M) = \|M\|_A$.*

2.2. Linear methods. Given the equation $Au = f$, where $A \in \mathbf{L}(\mathcal{H}, \mathcal{H})$ is SPD, consider the *preconditioned* equation $BAu = Bf$, with $B \in \mathbf{L}(\mathcal{H}, \mathcal{H})$. The operator B , the *preconditioner*, is usually chosen so that the linear iteration:

$$(3) \quad u^{n+1} = u^n - BAu^n + Bf = (I - BA)u^n + Bf,$$

has some desired convergence properties. The convergence of (3) is determined by the properties of the so-called *error propagation operator*,

$$(4) \quad E = I - BA.$$

The spectral radius of the error propagator E is called the *convergence factor* for the linear method, whereas the norm is referred to as the *contraction number*. We recall two well-known lemmas; see for example [15] or [18].

LEMMA 2.2. *For arbitrary f and u^0 , the condition $\rho(E) < 1$ is necessary and sufficient for convergence of the linear method (3).*

LEMMA 2.3. *The condition $\|E\| < 1$, or the condition $\|E\|_A < 1$, is sufficient for convergence of the linear method (3).*

We now state a series of simple lemmas that we shall use repeatedly in the following sections. Their short proofs are added for the reader's convenience.

LEMMA 2.4. *If A is SPD, then BA is A -self-adjoint if and only if B is self-adjoint.*

Proof. Note that: $(ABAu, v) = (BAu, Av) = (Au, B^T Av)$. The lemma follows since $BA = B^T A$ if and only if $B = B^T$. \square

LEMMA 2.5. *If A is SPD, then E is A -self-adjoint if and only if B is self-adjoint.*

Proof. Note that: $(AEu, v) = (Au, v) - (ABAu, v) = (Au, v) - (Au, (BA)^*v) = (Au, (I - (BA)^*)v)$. Therefore, $E^* = E$ if and only if $BA = (BA)^*$. By Lemma 2.4, this holds if and only if B is self-adjoint. \square

LEMMA 2.6. *If A and B are SPD, then BA is A -SPD.*

Proof. By Lemma 2.4, BA is A -self-adjoint. Also, $(ABAu, u) = (BAu, Au) = (B^{1/2}Au, B^{1/2}Au) > 0$, $\forall u \neq 0$. Hence, BA is A -positive, and the result follows. \square

LEMMA 2.7. *If A is SPD and B is self-adjoint, then $\|E\|_A = \rho(E)$.*

Proof. By Lemma 2.5, E is A -self-adjoint. By Lemma 2.1 the result follows. \square

LEMMA 2.8. *If E^* is the A -adjoint of E , then $\|E\|_A^2 = \|EE^*\|_A$.*

Proof. The proof follows that of a familiar result for the Euclidean 2-norm [11]. \square

LEMMA 2.9. *If A and B are SPD, and E is A -non-negative, then $\|E\|_A < 1$.*

Proof. By Lemma 2.5, E is A -self-adjoint. As E is A -non-negative, it holds that $(Eu, u)_A \geq 0$, or $(BAu, u)_A \leq (u, u)_A$. By Lemma 2.6, BA is A -SPD, and we have that $0 < (BAu, u)_A \leq (u, u)_A$, $\forall u \neq 0$, which, by (2), implies that $0 < \lambda_i \leq 1$, $\forall \lambda_i \in \sigma(BA)$. Thus, $\rho(E) = 1 - \min_i \lambda_i < 1$. Finally, by Lemma 2.7, we have $\|E\|_A = \rho(E)$. \square

We will have use for the following two simple lemmas, appearing previously in [24].

LEMMA 2.10. *If A is SPD and B is self-adjoint, and E is such that:*

$$-C_1(u, u)_A \leq (Eu, u)_A \leq C_2(u, u)_A, \quad \forall u \in \mathcal{H},$$

for $C_1 \geq 0$ and $C_2 \geq 0$, then $\rho(E) = \|E\|_A \leq \max\{C_1, C_2\}$.

Proof. By Lemma 2.5, E is A -self-adjoint, and by (2) $\lambda_{\min}(E)$ and $\lambda_{\max}(E)$ are bounded by $-C_1$ and C_2 , respectively. The result then follows by Lemma 2.7. \square

LEMMA 2.11. *If A and B are SPD, then Lemma 2.10 holds for some $C_2 < 1$.*

Proof. By Lemma 2.6, BA is A -SPD, which implies that the eigenvalues of BA are real and positive. Hence, we must have that $\lambda_i(E) = 1 - \lambda_i(BA) < 1$, $\forall i$. Since C_2 in Lemma 2.10 bounds the largest positive eigenvalue of E , we have that $C_2 < 1$. \square

2.3. Krylov acceleration of SPD linear methods. The conjugate gradient method was developed by Hestenes and Stiefel [12] as a method for solving linear systems $Au = f$ in a space \mathcal{H} , with SPD operators A . In order to improve convergence, it is common to *precondition* the linear system by an SPD *preconditioning operator* $B \approx A^{-1}$, in which case the generalized or preconditioned conjugate gradient method results ([7]). Our goal in this section is to briefly review some relationships between the contraction number of a basic linear preconditioner and that of the resulting preconditioned conjugate gradient algorithm.

We start with the well-known conjugate gradient contraction bound ([11]):

$$(5) \quad \|e^{i+1}\|_A \leq 2 \left(1 - \frac{2}{1 + \sqrt{\kappa_A(BA)}} \right)^{i+1} \|e^0\|_A = 2 \delta_{\text{cg}}^{i+1} \|e^0\|_A.$$

The ratio of extreme eigenvalues of BA appearing in the derivation of the bound gives rise to the generalized condition number $\kappa_A(BA)$ appearing above. This ratio is often mistakenly called the (spectral) condition number $\kappa(BA)$; in fact, since BA is not self-adjoint, this ratio is not in general equal to the usual condition number (this point is

discussed in great detail in [1]). However, the ratio does yield a condition number in the A -norm. The following lemma is a special case of Corollary 4.2 in [1].

LEMMA 2.12. *If A and B are SPD, then*

$$(6) \quad \kappa_A(BA) = \|BA\|_A \|(BA)^{-1}\|_A = \frac{\lambda_{\max}(BA)}{\lambda_{\min}(BA)}.$$

Remark 2.1. Often a linear method requires a parameter α in order to be convergent, leading to an error propagator of the form $E = I - \alpha BA$. Equation (6) shows that the A -condition number does not depend on the particular choice of α . Hence, one can use the conjugate gradient method as an accelerator for the method without a parameter, avoiding the possibly costly estimation of a good α .

The following result gives a bound on the condition number of the operator BA in terms of the extreme eigenvalues of the error propagator $E = I - BA$; such bounds are often used in the analysis of linear preconditioners (cf. Proposition 5.1 in [24]). We give a short proof of this result for completeness.

LEMMA 2.13. *If A and B are SPD, and E is such that:*

$$(7) \quad -C_1(u, u)_A \leq (Eu, u)_A \leq C_2(u, u)_A, \quad \forall u \in \mathcal{H},$$

for $C_1 \geq 0$ and $C_2 \geq 0$, then the above must hold with $C_2 < 1$, and it follows that:

$$\kappa_A(BA) \leq \frac{1 + C_1}{1 - C_2}.$$

Proof. First, since A and B are SPD, by Lemma 2.11 we have that $C_2 < 1$. Since $(Eu, u)_A = (u, u)_A - (BAu, u)_A$, it is clear that

$$(1 - C_2)(u, u)_A \leq (BAu, u)_A \leq (1 + C_1)(u, u)_A, \quad \forall u \in \mathcal{H}.$$

By Lemma 2.6, BA is A -SPD. Its eigenvalues are real and positive, and lie in the interval defined by the Raleigh quotients generated by the A -inner-product. Hence, that interval is given by $[(1 - C_2), (1 + C_1)]$, and by Lemma 2.12 the result follows. \square

Remark 2.2. Even if a linear method is not convergent, it may still be a good preconditioner. If it is the case that $C_2 \ll 1$, and if $C_1 > 1$ does not become too large, then $\kappa_A(BA)$ will be small and the conjugate gradient method will converge rapidly, even though the linear method diverges. This implication of Lemma 2.13 was first noticed in [24].

If only a bound on the norm of the error propagator $E = I - BA$ is available, then the following result can be used to bound the condition number of BA . This result is used for example in [25].

COROLLARY 2.14. *If A and B are SPD, and $\|I - BA\|_A \leq \delta < 1$, then*

$$(8) \quad \kappa_A(BA) \leq \frac{1 + \delta}{1 - \delta}.$$

Proof. This follows immediately from Lemma 2.13 with $\delta = \max\{C_1, C_2\}$. \square

The next result connects the contraction number of the preconditioner to the contraction number of the preconditioned conjugate gradient method. It shows that the conjugate gradient method always accelerates a linear method (if the conditions of the lemma hold).

LEMMA 2.15. *If A and B are SPD, and $\|I - BA\|_A \leq \delta < 1$, then $\delta_{\text{cg}} < \delta$.*

Proof. An abbreviated proof appears in [25], a more detailed proof in [13]. \square

2.4. Krylov acceleration of nonsymmetric linear methods. The convergence theory of the conjugate gradient iteration requires that the preconditioned operator BA be A -self-adjoint (see [2] for more general conditions), which from Lemma 2.4 requires that B be self-adjoint. If a Schwarz method is employed which produces a nonsymmetric operator B , then although A is SPD, the theory of the previous section does not apply, and a nonsymmetric solver such as conjugate gradients on the normal equations [2], GMRES [19], CGS [21], or Bi-CGstab [23] must be used for the now non- A -SPD preconditioned system, $BAu = Bf$.

The conjugate gradient method for SPD problems has several nice properties (good convergence rate, efficient three-term recursion, and minimization of the A -norm of the error at each step), some of which must be given up in order to generalize the method to nonsymmetric problems. For example, while GMRES attempts to maintain a minimization property and a good convergence rate, the three-term recursion must be sacrificed. Conjugate gradients on the normal equations maintains a minimization property as well as the efficient three-term recursion, but sacrifices convergence speed (the effective condition number is the square of the original system). Methods such as CGS and Bi-CGstab sacrifice the minimization property, but maintain good convergence speed and the efficient three-term recursion. For these reasons, methods such as CGS and Bi-CGstab have become the methods of choice in many applications that give rise to nonsymmetric problems. Bi-CGstab has been shown to be more attractive than CGS in many situations due to the more regular convergence behavior [23].

In §6, we shall use the preconditioned Bi-CGstab algorithm to accelerate nonsymmetric Schwarz methods. In a sequence of numerical experiments, we shall compare the effectiveness of this approach with unaccelerated symmetric and nonsymmetric Schwarz methods, and with symmetric Schwarz methods accelerated with conjugate gradients.

3. Multiplicative Schwarz methods. We develop a preconditioning theory of product algorithms which establishes sufficient conditions for producing SPD preconditioners. This theory is used to establish sufficient SPD conditions for multiplicative DD and MG methods.

3.1. A product operator. Consider a product operator of the form:

$$(9) \quad E = I - BA = (I - \bar{B}_1 A)(I - B_0 A)(I - B_1 A),$$

where \bar{B}_1, B_0 and B_1 are linear operators on \mathcal{H} , and where A is, as before, an SPD operator on \mathcal{H} . We are interested in conditions for \bar{B}_1, B_0 and B_1 , which guarantee

that the implicitly defined operator B is self-adjoint and positive definite and, hence, can be accelerated by using the conjugate gradient method.

LEMMA 3.1. *Sufficient conditions for symmetry and positivity of operator B , implicitly defined by (9), are:*

1. $\bar{B}_1 = B_1^T$;
2. $B_0 = B_0^T$;
3. $\|I - B_1 A\|_A < 1$;
4. B_0 non-negative on \mathcal{H} .

Proof. By Lemma 2.5, in order to prove symmetry of B , it is sufficient to prove that E is A -self-adjoint. By using (1), we get

$$\begin{aligned} E^* &= A^{-1} E^T A \\ &= A^{-1} (I - AB_1^T)(I - AB_0^T)(I - A\bar{B}_1^T)A \\ &= (I - B_1^T A)(I - B_0^T A)(I - \bar{B}_1^T A) \\ &= (I - \bar{B}_1 A)(I - B_0 A)(I - B_1 A) = E, \end{aligned}$$

which follows from conditions 1 and 2.

Next, we prove that $(Bu, u) > 0$, $\forall u \in \mathcal{H}$, $u \neq 0$. Since A is non-singular, this is equivalent to proving that $(BAu, Au) > 0$. Using condition 1, we have that

$$\begin{aligned} (BAu, Au) &= ((I - E)u, Au) \\ &= (u, Au) - ((I - B_1^T A)(I - B_0 A)(I - B_1 A)u, Au) \\ &= (u, Au) - ((I - B_0 A)(I - B_1 A)u, A(I - B_1 A)u) \\ &= (u, Au) - ((I - B_1 A)u, A(I - B_1 A)u) + (B_0 w, w), \end{aligned}$$

where $w = A(I - B_1 A)u$. By conditions 2 and 4, we have that $(B_0 w, w) \geq 0$. Condition 3 implies that $((I - B_1 A)u, A(I - B_1 A)u) < (u, Au)$ for $u \neq 0$. Thus, the first two terms above are together positive, while the third is non-negative, so that B is positive. \square

COROLLARY 3.2. *If $B_1 = B_1^T$, then condition 3 in Lemma 3.1 is equivalent to $\rho(I - B_1 A) < 1$.*

Proof. This follows directly from Lemma 2.1 and Lemma 2.5. \square

3.2. Multiplicative domain decomposition. Given the finite-dimensional Hilbert space \mathcal{H} , consider J spaces \mathcal{H}_k , $k = 1, \dots, J$, together with linear operators $I_k \in \mathbf{L}(\mathcal{H}_k, \mathcal{H})$, $\text{null}(I_k) = \{0\}$, such that $I_k \mathcal{H}_k \subseteq \mathcal{H} = \sum_{k=1}^J I_k \mathcal{H}_k$. We also assume the existence of another space \mathcal{H}_0 , the associated operator I_0 such that $I_0 \mathcal{H}_0 \subseteq \mathcal{H}$, and some linear operators $I^k \in \mathbf{L}(\mathcal{H}, \mathcal{H}_k)$, $k = 0, \dots, J$. We shall assume that the inner-products (and hence also their induced norms) on \mathcal{H}_k are inherited from the “parent” space \mathcal{H} , and we shall denote them by (\cdot, \cdot) .

In a domain decomposition context, the spaces \mathcal{H}_k , $k = 1, \dots, J$ are typically associated with *local subdomains* of the original domain on which the partial differential equation is defined. The space \mathcal{H}_0 is then a space associated with some global coarse mesh. The operators I_k , $k = 1, \dots, J$ are usually inclusion operators, while I_0 is an interpolation or prolongation operator (as in a two-level MG method). The operators

$I^k, k = 1, \dots, J$ are usually orthogonal projection operators, while I^0 is a restriction operator (again, as in a two-level MG method).

The error propagator of a multiplicative DD method on the space \mathcal{H} employing the subspaces $I_k \mathcal{H}_k$ has the general form [8]:

$$(10) \quad E = I - BA = (I - I_J \bar{R}_J I^J A) \cdots (I - I_0 R_0 I^0 A) \cdots (I - I_J R_J I^J A) ,$$

where \bar{R}_k and $R_k, k = 1, \dots, J$, are linear operators on \mathcal{H}_k , and R_0 is a linear operator on \mathcal{H}_0 . Usually the operators \bar{R}_k and R_k are constructed so that $\bar{R}_k \approx A_k^{-1}$ and $R_k \approx A_k^{-1}$, where A_k is the operator defining the subdomain problem in \mathcal{H}_k . Similarly, R_0 is constructed so that $R_0 \approx A_0^{-1}$. Actually, quite often R_0 is a “direct solve”, i.e., $R_0 = A_0^{-1}$. The subdomain problem operator A_k is related to the restriction of A to \mathcal{H}_k . We say that A_k is *variationally* defined or satisfies the *Galerkin conditions* when

$$(11) \quad A_k = I^k A I_k, \quad I^k = c_k I_k^T, \quad c_k > 0.$$

In the case of finite element, finite volume, or finite difference discretization of an elliptic problem, conditions (11) can be shown to hold naturally for both the matrices and the abstract weak form operators, with $c_k = 1$, for all subdomains $k = 1, \dots, J$. For the coarse space \mathcal{H}_0 , often (11) must be imposed algebraically, perhaps with $c_0 \neq 1$.

Propagator (10) can be thought of as the product operator (9), by choosing

$$I - \bar{B}_1 A = \prod_{k=J}^1 (I - I_k \bar{R}_k I^k A) , \quad B_0 = I_0 R_0 I^0 , \quad I - B_1 A = \prod_{k=1}^J (I - I_k R_k I^k A) ,$$

where \bar{B}_1 and B_1 are known only implicitly. (Note that we take the convention that the first term in the product appears on the left.) This identification allows for the use of Lemma 3.1 to establish sufficient conditions on the subdomain operators \bar{R}_k, R_k and R_0 to guarantee that multiplicative domain decomposition yields an SPD operator B .

THEOREM 3.3. *Sufficient conditions for symmetry and positivity of the multiplicative domain decomposition operator B , implicitly defined by (10), are:*

1. $I^k = c_k I_k^T$, $c_k > 0$, $k = 0, \dots, J$;
2. $\bar{R}_k = R_k^T$, $k = 1, \dots, J$;
3. $R_0 = R_0^T$;
4. $\left\| \prod_{k=1}^J (I - I_k R_k I^k A) \right\|_A < 1$;
5. R_0 non-negative on \mathcal{H}_0 .

Proof. We show that the sufficient conditions of Lemma 3.1 are satisfied. First, we prove that $\bar{B}_1 = B_1^T$, which, by Lemma 2.5, is equivalent to proving that $(I - B_1 A)^* = (I - \bar{B}_1 A)$. By using (1), we have

$$\left(\prod_{k=1}^J (I - I_k R_k I^k A) \right)^* = A^{-1} \left(\prod_{k=1}^J (I - I_k R_k I^k A) \right)^T A = \prod_{k=J}^1 (I - (I^k)^T R_k^T (I_k)^T A) ,$$

which equals $(I - \bar{B}_1 A)$ under conditions 1 and 2 of the theorem. The symmetry of B_0 follows immediately from conditions 1 and 3; indeed,

$$B_0^T = (I_0 R_0 I^0)^T = (I^0)^T R_0^T (I_0)^T = (c_0 I_0) R_0 (c_0^{-1} I^0) = I_0 R_0 I^0 = B_0 .$$

By condition 4 of the theorem, condition 3 of Lemma 3.1 holds trivially. The theorem follows by realizing that condition 4 of Lemma 3.1 is also satisfied, since,

$$(B_0 u, u) = (I_0 R_0 I^0 u, u) = (R_0 I^0 u, I_0^T u) = c_0^{-1} (R_0 I^0 u, I^0 u) \geq 0, \quad \forall u \in \mathcal{H}.$$

□

Remark 3.3. Note that one sweep through the subdomains, followed by a coarse problem solve, followed by another sweep through the subdomains in reversed order, gives rise an error propagator of the form (10).

Remark 3.4. Note that no conditions are imposed on the nature of the operators A_k associated with each subdomain. In particular, the theorem *does not* require that the variational conditions are satisfied. While it is natural for condition (11) to hold between the fine space and the spaces associated with each subdomain, these conditions are often difficult to enforce for the coarse problem. Violation of variational conditions can occur, for example, when complex coefficient discontinuities do not lie along element boundaries on the coarse mesh (we present numerical results for such a problem in §6).

Condition 1 of the theorem (with $c_k = 1$) for $k = 1, \dots, J$ is usually satisfied trivially for domain decomposition methods. For $k = 0$, it may have to be imposed explicitly. Condition 2 of the theorem allows for several alternatives which give rise to an SPD preconditioner, namely: (1) use of exact subdomain solvers (if A_k is a symmetric operator); (2) use of identical symmetric subdomain solvers in the forward and backward sweeps; (3) use of the adjoint of the subdomain solver on the second sweep. Condition 3 is satisfied when the coarse problem is symmetric and the solve is an exact one, which is usually the case. If not, the coarse problem solve has to be symmetric. Condition 5 is satisfied for example when the coarse problem is SPD and the solve is exact.

Condition 4 in Theorem 3.3 is clearly a non-trivial one; it is essentially the assumption that the multiplicative DD method without a coarse space is convergent. Convergence theories for DD methods can be quite technical and depend on such things as the discretization, the subdomain number, shape, and size, and the regularity of the solution [4, 8, 25]. However, since variational conditions hold naturally between the fine space and each subdomain space for nearly any formulation of a DD method, very general convergence theorems can be derived, if one is not concerned about the actual rate of convergence. Using the Schwarz theory framework in any of [4, 8, 25], it can be shown that Condition 4 in Theorem 3.3 (convergence of multiplicative DD without a coarse space) hold if the variational conditions (11) holds, and if the subdomain solvers R_k are SPD. A proof of this result may be found for example in [13].

Remark 3.5. Note that the theorem does not require that the overall multiplicative DD method be convergent. In particular, the conditions on the coarse problem and coarse problem solver are very relaxed.

3.3. Multiplicative multigrid. Given are the Hilbert space \mathcal{H} , J spaces \mathcal{H}_k together with linear operators $I_k \in \mathbf{L}(\mathcal{H}_k, \mathcal{H})$, $\text{null}(I_k) = 0$, such that the spaces $I_k \mathcal{H}_k$ are nested and satisfy $I_1 \mathcal{H}_1 \subseteq I_2 \mathcal{H}_2 \subseteq \dots \subseteq I_{J-1} \mathcal{H}_{J-1} \subseteq \mathcal{H}_J \equiv \mathcal{H}$. As before we shall

denote the \mathcal{H}_k -inner-products by (\cdot, \cdot) , and assume that they are inherited from the parent inner-product on \mathcal{H} . We assume also the existence of operators $I^k \in \mathbf{L}(\mathcal{H}, \mathcal{H}_k)$.

In a multigrid context, the spaces \mathcal{H}_k are typically associated with a nested hierarchy of successively refined meshes, with \mathcal{H}_1 being the coarsest mesh, and \mathcal{H}_J being the fine mesh on which the PDE solution is desired. The linear operators I_k are prolongation operators, constructed from given interpolation or prolongation operators that operate between subspaces, i.e., $I_{k-1}^k \in \mathbf{L}(\mathcal{H}_{k-1}, \mathcal{H}_k)$. The operator I_k is then constructed (only as a theoretical tool) as a composite operator

$$(12) \quad I_k = I_{J-1}^J I_{J-2}^{J-1} \cdots I_{k+1}^{k+2} I_k^{k+1}, \quad k = 1, \dots, J-1.$$

The composite restriction operators I^k , $k = 1, \dots, J-1$, are constructed similarly from some given restriction operators $I_k^{k-1} \in \mathbf{L}(\mathcal{H}_k, \mathcal{H}_{k-1})$.

The coarse problem operators A_k are related to the restriction of A to \mathcal{H}_k . As in the case of DD methods, we say that A_k is *variationally* defined or satisfies the *Galerkin conditions* when conditions (11) hold. It is not difficult to see that conditions (11) are equivalent to the following recursively defined variational conditions:

$$(13) \quad A_k = I_{k+1}^k A_{k+1} I_k^{k+1}, \quad I_{k+1}^k = c_k (I_k^{k+1})^T, \quad c_k > 0,$$

when the composite operators I_k appearing in (11) are defined as in (12).

In a finite element setting, conditions (13) with $c_k = 1$ can be shown to hold in ideal situations, for both the stiffness matrices and the abstract weak form operators, for a nested sequence of successively refined finite element meshes. In the finite difference or finite volume method setting, conditions (13) must often be imposed algebraically, in a recursive fashion, typically with $c_k \neq 1$.

The error propagator of a multiplicative V-cycle MG method is defined implicitly:

$$(14) \quad E = I - BA = I - D_J A_J,$$

where $A_J = A$, and where operators D_k , $k = 2, \dots, J$ are defined recursively,

$$(15) \quad I - D_k A_k = (I - \bar{R}_k A_k)(I - I_{k-1}^k D_{k-1} I_k^{k-1} A_k)(I - R_k A_k), \quad k = 2, \dots, J,$$

$$(16) \quad D_1 = R_1.$$

Operators \bar{R}_k and R_k are linear operators on \mathcal{H}_k , usually called *smoothers*. The linear operators $A_k \in \mathbf{L}(\mathcal{H}_k, \mathcal{H}_k)$ define the coarse problems. They often satisfy the variational condition (11).

The error propagator (14) can be thought of as an operator of the form (9) with

$$\bar{B}_1 = \bar{R}_J, \quad B_0 = I_{J-1}^J D_{J-1} I_J^{J-1}, \quad B_1 = R_J.$$

Such an identification with the product method allows for the use of the result in Lemma 3.1. The following theorem establishes sufficient conditions for the subspace operators R_k , \bar{R}_k and A_k in order to generate an (implicitly defined) SPD operator B that can be accelerated with conjugate gradients.

THEOREM 3.4. *Sufficient conditions for symmetry and positivity of the multiplicative multigrid operator B , implicitly defined by (14), (15), and (16), are*

1. A_k is SPD on \mathcal{H}_k , $k = 1, \dots, J-1$;
2. $I_k^{k-1} = c_k(I_{k-1}^k)^T$, $c_k > 0$, $k = 2, \dots, J$;
3. $\bar{R}_k = R_k^T$, $k = 2, \dots, J$;
4. $R_1 = R_1^T$;
5. $\|I - R_J A\|_A < 1$, ;
6. $\|I - R_k A_k\|_{A_k} \leq 1$, $k = 2, \dots, J-1$;
7. R_1 non-negative on \mathcal{H}_1 .

Proof. Since $\bar{R}_J = R_J^T$, we have that $\bar{B}_1 = B_1^T$, which gives condition 1 of Lemma 3.1. Now, B_0 is symmetric if and only if

$$B_0 = I_{J-1}^J D_{J-1} I_J^{J-1} = (c_J^{-1} I_J^{J-1})^T D_{J-1}^T (c_J I_{J-1}^J)^T = B_0^T,$$

which holds under condition 2 and a symmetry requirement for D_{J-1} . We will prove that $D_{J-1} = D_{J-1}^T$ by induction. First, $D_1 = D_1^T$ since $R_1 = R_1^T$. By Lemma 2.5 and condition 1, D_k is symmetric if and only if $E_k = I - D_k A_k$ is A_k -self-adjoint. By using (1), we have that

$$\begin{aligned} E_k^* &= A_k^{-1} \left((I - \bar{R}_k A_k)(I - I_{k-1}^k D_{k-1} I_k^{k-1} A_k)(I - R_k A_k) \right)^T A_k \\ &= A_k^{-1} (I - A_k^T R_k^T)(I - A_k^T (I_k^{k-1})^T D_{k-1}^T (I_{k-1}^k)^T)(I - A_k^T \bar{R}_k^T) A_k \\ &= (I - R_k^T A_k) A_k^{-1} (I - A_k^T (I_k^{k-1})^T D_{k-1}^T (I_{k-1}^k)^T) A_k (I - \bar{R}_k^T A_k) \\ &= (I - \bar{R}_k A_k)(I - (c_k I_{k-1}^k) D_{k-1}^T (c_k^{-1} I_k^{k-1}) A_k)(I - R_k A_k), \end{aligned}$$

where we have used conditions 1, 2 and 3. Therefore, $E_k^* = E_k$, if $D_{k-1} = D_{k-1}^T$. Hence, the result follows by induction on k .

Condition 3 of Lemma 3.1 follows trivially by condition 5 of the theorem.

It remains to verify condition 4 of Lemma 3.1, namely that B_0 is non-negative. This is equivalent to showing that D_{J-1} is non-negative on \mathcal{H}_{J-1} . This will follow again from an induction argument. First, note that $D_1 = R_1$ is non-negative on \mathcal{H}_1 . Next, we prove that $(D_k v_k, v_k) \geq 0$, $\forall v_k \in \mathcal{H}_k$, or, equivalently, since A_k is non-singular, that $(D_k A_k v_k, A_k v_k) \geq 0$. So, for all $v_k \in \mathcal{H}_k$,

$$\begin{aligned} (D_k A_k v_k, A_k v_k) &= (A_k v_k, v_k) - (A_k E_k v_k, v_k) \\ &= (A_k v_k, v_k) - (A_k (I - \bar{R}_k A_k)(I - I_{k-1}^k D_{k-1} I_k^{k-1} A_k)(I - R_k A_k) v_k, v_k) \\ &= (A_k v_k, v_k) - (A_k (I - I_{k-1}^k D_{k-1} I_k^{k-1} A_k)(I - R_k A_k) v_k, (I - R_k A_k) v_k) \\ &= (A_k v_k, v_k) - (A_k (I - R_k A_k) v_k, (I - R_k A_k) v_k) \\ &\quad + (A_k I_{k-1}^k D_{k-1} I_k^{k-1} A_k (I - R_k A_k) v_k, (I - R_k A_k) v_k) \\ &= (v_k, v_k)_{A_k} - (S_k v_k, S_k v_k)_{A_k} + c_k^{-1} (D_{k-1} v_{k-1}, v_{k-1}) \end{aligned}$$

where $S_k = I - R_k A_k$ and $v_{k-1} = I_k^{k-1} A_k (I - R_k A_k) v_k \in \mathcal{H}_{k-1}$. By condition 6, the first two terms add up to a non-negative value. Hence, D_k is non-negative if D_{k-1} is non-negative. Condition 4 of Lemma 3.1 follows. \square

COROLLARY 3.5. *If the fine grid smoother is symmetric, i.e., $R_J = \bar{R}_J^T$, then condition 5 in Theorem 3.4 is equivalent to $\rho(I - R_J A) < 1$.*

Proof. This follows directly from Corollary 3.2. \square

Remark 3.6. The coarse grid operators A_k , $k = 1, \dots, J-1$, need only be SPD. They need not satisfy the Galerkin conditions (11).

Condition 1 of the theorem requires that all coarse grid operators be SPD. This is easily satisfied when they are constructed either by discretization or by explicitly using the Galerkin or variational condition. Condition 2 requires restriction and prolongation to be adjoints in the inherited inner-product, possibly multiplied by an arbitrary constant. Condition 3 of the theorem is satisfied when the number of pre-smoothing steps equals the number of post-smoothing steps, and in addition one of the following is imposed: (1) use of the same symmetric smoother for both pre- and post-smoothing; (2) use of the adjoint of the pre-smoothing operator as the post-smoother. Condition 4 requires a symmetric coarsest mesh solver. When the coarsest mesh problem is SPD, the symmetry of R_1 is satisfied when it corresponds to an exact solve (as is typical for MG methods). Condition 5 is a convergence requirement on the fine space smoother. Condition 6 requires the coarse grid smoothers to be non-divergent. The nonnegativity requirement for R_1 is a non-trivial one; however, since A_1 is SPD, it is immediately satisfied when the operator corresponds to an exact solve.

If variational conditions are satisfied on all levels, and convergent smoothers are used on all levels, then there is a simple proof which shows that in addition to defining an SPD operator B , the conditions of the theorem are sufficient to prove the convergence of the MG method itself. The result is as follows.

THEOREM 3.6. *If in addition to the conditions for Theorem 3.4, it holds that $A_k = I^k A I_k$, $I^k = c_k I_k^T$, $c_k \equiv 1$, and $R_1 = A_1^{-1}$, then the MG error propagator satisfies:*

$$\rho(E) \leq \|E\|_A < 1.$$

Proof. Under the conditions of the theorem, the MG error propagator can be written explicitly as the product ([3, 17]):

$$E = (I - I_J R_J^T I_J^T A) \cdots (I - I_1 R_1 I_1^T A) \cdots (I - I_J R_J I_J^T A) .$$

Since the coarse problem is solved exactly, and since variational conditions hold, the coarse product term is an A -orthogonal projector:

$$I - I_1 R_1 I_1^T A = I - I_1 (I_1 A I_1^T)^{-1} I_1^T A = (I - I_1 (I_1 A I_1^T)^{-1} I_1^T A)^2 = (I - I_1 R_1 I_1^T A)^2 .$$

Therefore, we may define $\bar{E} = (I - I_1 R_1 I_1^T A) \cdots (I - I_J R_J I_J^T A)$, and represent E as the product $E = \bar{E}^* \bar{E}$. Now, since A is SPD, we have that:

$$(AEv, v) = (A\bar{E}v, \bar{E}v) \geq 0 .$$

Hence, E is A -non-negative. Under the conditions of the theorem, Lemma 3.1 implies that the preconditioner is SPD, and so by Lemma 2.9 it holds that $\|E\|_A < 1$. \square

4. Additive Schwarz methods. We now present an analysis of additive Schwarz methods. We establish sufficient conditions for additive algorithms to yield SPD preconditioners. This theory is then employed to establish sufficient SPD conditions for additive DD and MG methods.

4.1. A sum operator. Consider a sum operator of the following form:

$$(17) \quad E = I - BA = I - \omega(B_0 + B_1)A, \quad \omega > 0,$$

where, as before, A is an SPD operator, and B_0 and B_1 are linear operators on \mathcal{H} .

LEMMA 4.1. *Sufficient conditions for symmetry and positivity of B , defined in (17), are*

1. B_1 is SPD in \mathcal{H} ;
2. B_0 is symmetric and non-negative on \mathcal{H} .

Proof. We have that $B = \omega(B_0 + B_1)$, which is symmetric by the symmetry of B_0 and B_1 . Positivity follows since $(B_0u, u) \geq 0$ and $(B_1u, u) > 0, \forall u \in \mathcal{H}, u \neq 0$. \square

Remark 4.7. The parameter ω is usually required to make the additive method a convergent one. Its estimation is often nontrivial, and can be very costly. As was noted in Remark 2.1, the parameter ω is not required when the linear additive method is used as a preconditioner in a conjugate gradients algorithm. This is exactly why additive multigrid and domain decomposition methods are used almost exclusively as preconditioners.

4.2. Additive domain decomposition. As in §3.2, we consider the Hilbert space \mathcal{H} , and J subspaces $I_k\mathcal{H}_k$ such that $I_k\mathcal{H}_k \subseteq \mathcal{H} = \sum_{k=1}^J I_k\mathcal{H}_k$. Again, we allow for the existence of a “coarse” subspace $I_0\mathcal{H}_0 \subseteq \mathcal{H}$.

The error propagator of an additive DD method on the space \mathcal{H} employing the subspaces $I_k\mathcal{H}_k$ has the general form (see [25]):

$$(18) \quad E = I - BA = I - \omega(I_0R_0I^0 + I_1R_1I^1 + \cdots + I_JR_JI^J)A.$$

The operators R_k are linear operators on \mathcal{H}_k , constructed in such a way that $R_k \approx A_k^{-1}$, where the A_k are the subdomain problem operators. Propagator (18) can be thought of as the sum method (17), by taking

$$B_0 = I_0R_0I^0, \quad B_1 = \sum_{k=1}^J I_kR_kI^k.$$

This identification allows for the use of Lemma 4.1 in order to establish conditions to guarantee that additive domain decomposition yields an SPD preconditioner. Before we state the main theorem, we need the following lemma, which characterizes the splitting of \mathcal{H} into the subspaces $I_k\mathcal{H}_k$ in terms of a positive *splitting constant* S_0 .

LEMMA 4.2. *Given any $v \in \mathcal{H}$, there exists a splitting $v = \sum_{k=1}^J I_kv_k$, $v_k \in \mathcal{H}_k$, and a constant $S_0 > 0$, such that*

$$(19) \quad \sum_{k=1}^J \|I_kv_k\|_A^2 \leq S_0 \|v\|_A^2.$$

Proof. Since $\sum_{k=1}^J I_k\mathcal{H}_k = \mathcal{H}$, we can construct subspaces $\mathcal{V}_k \subseteq \mathcal{H}_k$, such that

$$I_k\mathcal{V}_k \cap I_l\mathcal{V}_l = \{0\}, \text{ for } k \neq l \text{ and } \mathcal{H} = \sum_{k=1}^J I_k\mathcal{V}_k.$$

Any $v \in \mathcal{H}$, can be decomposed uniquely as $v = \sum_{k=1}^J I_k v_k$, $v_k \in \mathcal{V}_k$. Define the projectors $Q_k \in L(\mathcal{H}, \mathcal{V}_k)$ such that $Q_k v = I_k v_k$. Then,

$$\sum_{k=1}^J \|I_k v_k\|_A^2 = \sum_{k=1}^J \|Q_k v\|_A^2 \leq \sum_{k=1}^J \|Q_k\|_A^2 \|v\|_A^2.$$

Hence, the result follows with $S_0 = \sum_{k=1}^J \|Q_k\|_A^2$. \square

THEOREM 4.3. *Sufficient conditions for symmetry and positivity of the additive domain decomposition operator B , defined in (18), are*

1. $A_k = I^k A I_k$, $k = 1, \dots, J$;
2. $I^k = c_k I_k^T$, $c_k > 0$, $k = 0, \dots, J$;
3. R_k is SPD on \mathcal{H}_k , $k = 1, \dots, J$;
4. R_0 is symmetric and non-negative on \mathcal{H}_0 .

Proof. Symmetry of B_0 and B_1 follow trivially from the symmetry of R_k and R_0 , and from $I^k = c_k I_k^T$. That B_0 is non-negative on \mathcal{H} follows immediately from the non-negativity of R_0 on \mathcal{H}_0 .

Finally, we prove positivity of B_1 . By conditions 1 and 2, and the full rank nature of I_k , we have that A_k is SPD. Now, since R_k is also SPD, the product $R_k A_k$ is A_k -SPD. Hence, there exists an $\omega_0 > 0$ such that $0 < \omega_0 < \lambda_i(R_k A_k)$, $k = 1, \dots, J$. This is used together with (19) to bound the following sum,

$$\begin{aligned} \sum_{k=1}^J c_k^{-1} (R_k^{-1} v_k, v_k) &= \sum_{k=1}^J c_k^{-1} (A_k A_k^{-1} R_k^{-1} v_k, v_k) \\ &\leq \sum_{k=1}^J c_k^{-1} (A_k v_k, v_k) \max_{v_k \neq 0} \frac{(A_k A_k^{-1} R_k^{-1} v_k, v_k)}{(A_k v_k, v_k)} \leq \sum_{k=1}^J c_k^{-1} \omega_0^{-1} (A_k v_k, v_k) \\ &= \sum_{k=1}^J \omega_0^{-1} (A I_k v_k, I_k v_k) = \sum_{k=1}^J \omega_0^{-1} \|I_k v_k\|_A^2 \leq \left(\frac{S_0}{\omega_0} \right) \|v\|_A^2. \end{aligned}$$

We can now employ this result to establish positivity of B_1 .

$$\|v\|_A^2 = (Av, v) = \sum_{k=1}^J (Av, I_k v_k) = \sum_{k=1}^J (I_k^T Av, v_k) = \sum_{k=1}^J (R_k c_k^{1/2} I_k^T Av, R_k^{-1} c_k^{-1/2} v_k).$$

By using the Cauchy-Schwarz inequality in the R_k -norm, we have that

$$\begin{aligned} \|v\|_A^2 &\leq \left(\sum_{k=1}^J (R_k R_k^{-1} c_k^{-1/2} v_k, R_k^{-1} c_k^{-1/2} v_k) \right)^{1/2} \left(\sum_{k=1}^J (R_k c_k^{1/2} I_k^T Av, c_k^{1/2} I_k^T Av) \right)^{1/2} \\ &\leq \left(\frac{S_0}{\omega_0} \right)^{1/2} \|v\|_A \left(\sum_{k=1}^J (I_k R_k c_k I_k^T Av, Av) \right)^{1/2} \\ &= \left(\frac{S_0}{\omega_0} \right)^{1/2} \|v\|_A (B_1 Av, Av)^{1/2}. \end{aligned}$$

Finally, division by $\|v\|_A$ and squaring yields

$$(B_1 A v, A v) \geq \frac{\omega_0}{S_0} \|v\|_A^2 > 0, \quad \forall v \in \mathcal{H}, \quad v \neq 0.$$

□

Remark 4.8. Variational conditions are required for the subdomain operators A_k , $k \neq 0$. However, this is a very natural condition with domain decomposition methods. No variational conditions are needed for the coarse space operator A_0 .

As explained in the previous remark, condition 1 of the theorem is usually satisfied. Condition 2 is also naturally satisfied for $k = 1, \dots, J$, with $c_k = 1$, since the associated I_k and I^k are usually inclusion and orthogonal projection operators (which are natural adjoints). The fact that $I^0 = c_0 I_0^T$ needs to be satisfied explicitly. Condition 3 requires the use of SPD subdomain solvers. The condition will hold, for example, when the subdomain solve is exact. (Note that A_k is SPD by condition 1 and the full rank nature of I_k .) Finally, condition 4 is nontrivial, and needs to be checked explicitly. The condition holds when A_0 is SPD and the solve is exact.

4.3. Additive multigrid. As in §3.3, given are the Hilbert space \mathcal{H} , and $J - 1$ nested subspaces $I_k \mathcal{H}_k$ such that $I_1 \mathcal{H}_1 \subseteq I_2 \mathcal{H}_2 \subseteq \dots \subseteq I_{J-1} \mathcal{H}_{J-1} \subseteq \mathcal{H}_J \equiv \mathcal{H}$. The operators I_k , I^k , I_{k-1}^k and I_k^{k-1} are the usual linear operators between the different spaces, as in the previous sections.

The error propagator of an additive MG method is defined explicitly:

$$(20) \quad E = I - BA = I - \omega(I_1 R_1 I^1 + I_2 R_2 I^2 + \dots + I_{J-1} R_{J-1} I^{J-1} + R_J)A.$$

This can be thought of as the sum method analyzed earlier, by taking

$$B_0 = \sum_{k=1}^{J-1} I_k R_k I^k, \quad B_1 = R_J.$$

This identification allows for the use of Lemma 4.1 to establish sufficient conditions to guarantee that additive MG yields an SPD preconditioner.

THEOREM 4.4. *Sufficient conditions for symmetry and positivity of the additive multigrid operator B , defined in (20), are:*

1. $I^k = c_k I_k^T$, $c_k > 0$, $k = 1, \dots, J - 1$;
2. R_J is SPD in \mathcal{H} ;
3. R_k is symmetric non-negative in \mathcal{H}_k , $k = 1, \dots, J - 1$.

Proof. Symmetry of B_0 and B_1 is obvious. B_1 is positive by condition 2. Non-negativity of B_0 follows from

$$(B_0 u, u) = \sum_{k=1}^{J-1} (I_k R_k (c_k I_k)^T u, u) = \sum_{k=1}^{J-1} c_k (R_k I_k^T u, I_k^T u) \geq 0, \quad \forall u \in \mathcal{H}, u \neq 0.$$

□

Remark 4.9. Variational conditions for the subspace operators A_k are not required.

Condition 1 of the theorem has to be imposed explicitly. Conditions 2 and 3 require the smoothers to be symmetric. The positivity of R_J is satisfied when the fine grid smoother is convergent, although this is not a necessary condition. The non-negativity of R_k , $k < J$, has to be checked explicitly. When the coarse problem operators A_k are SPD, this condition is satisfied, for example, when the smoothers are non-divergent.

5. To symmetrize or not to symmetrize. The following lemma illustrates why symmetrizing is a bad idea for linear methods. It exposes the convergence rate penalty incurred by symmetrization of a linear method.

LEMMA 5.1. *For any $E \in \mathbf{L}(\mathcal{H}, \mathcal{H})$, it holds that:*

$$\rho(E) \leq \|EE\|_A \leq \|E\|_A^2 = \|EE^*\|_A = \rho(E E^*).$$

Proof. The first and second inequalities hold for any norm. The first equality follows from Lemma 2.8, and the second follows from Lemma 2.1. \square

Note that this is an inequality not only for the spectral radii, which is only an asymptotical measure of convergence, but also for the A -norms of the nonsymmetric and symmetrized error propagators. The lemma illustrates that one may actually see the differing convergence rates early in the iteration as well.

Based on this lemma, and Corollary 2.14, we conjecture that when symmetrization of a linear method is required for its use as a preconditioner, the best results will be obtained by enforcing only a minimal amount of symmetry.

6. Numerical results. We present numerical results obtained by using multiplicative and additive finite-element-based DD and MG methods applied to two test problems, and we illustrate the theory of the preceding sections.

6.1. Example 1. Violation of variational conditions can occur in DD and MG methods when, for example, complex coefficient discontinuities do not lie along element boundaries on coarse meshes. An example of this occurs with the following test problem. The Poisson-Boltzmann equation describes the electrostatic potential of a biomolecule lying in an ionic solvent (see, e.g., [6] for an overview). This nonlinear elliptic equation for the dimensionless electrostatic potential $u(\mathbf{r})$ has the form:

$$-\nabla \cdot (\epsilon(\mathbf{r}) \nabla u(\mathbf{r})) + \bar{\kappa}^2 \sinh(u(\mathbf{r})) = \left(\frac{4\pi e_c^2}{k_B T} \right) \sum_{i=1}^{N_m} z_i \delta(\mathbf{r} - \mathbf{r}_i), \quad \mathbf{r} \in \mathbb{R}^3, \quad u(\infty) = 0.$$

The coefficients appearing in the equation are discontinuous by several orders of magnitude. The placement and magnitude of atomic charges are represented by source terms involving delta-functions. Analytical techniques are used to obtain boundary conditions on a finite domain boundary.

We will compare several MG and DD methods for a two-dimensional, linearized Poisson-Boltzmann problem, modeling a molecule with three point charges. The surface of the molecule is such that the discontinuities do not align with the coarsest mesh or with the subdomain boundaries. Beginning with the coarse mesh shown on the left

TABLE 1
Normalized operation counts per iteration, Example 1.

Method	UNACCEL	CG	Bi-CGstab
multiplicative MG	1.0	1.3	2.6
additive MG	1.1	1.4	2.7
multiplicative DD	3.5	3.8	7.5
additive DD	3.1	3.4	6.7

in Figure 1, we uniformly refine the initial mesh of 10 elements (9 nodes) five times, leading to a fine mesh of 2560 elements (1329 nodes). Piecewise linear finite elements, combined with one-point Gaussian quadrature, are used to discretize the problem. The three coarsest meshes used to formulate the MG methods are given in Figure 1. For the DD methods, the subdomains, corresponding to the initial coarse triangulation, are given a small overlap of one fine mesh triangle. The DD methods also employ a coarse space constructed from the initial triangulation. Figure 2 shows three overlapping subdomains overlaying the initial coarse mesh.

Computed results are presented in Tables 2 to 5. Given for each experiment is the number of iterations required to satisfy the error criterion (reduction of the A -norm of the error by 10^{-10}). We report results for the unaccelerated, CG-accelerated, and Bi-CGstab-accelerated methods. Since the cost of one iteration differs for each method, Table 1 gives the operation counts per iteration, normalized by the cost of a single multigrid iteration. For the MG operation counts, two smoothing iterations (one pre- and one post-smoothing) are used. The DD operation counts are for methods employing two sweeps through the subdomains, each approximate subdomain solve consisting of four sweeps of a Gauss-Seidel iteration.

Table 1 shows that multiplicative MG is slightly less costly than additive MG, since it is formulated in the usual recursive fashion, requiring fewer prolongations and restrictions. On the other hand, multiplicative DD is somewhat more costly than additive DD, due to the need to update boundary information after the solution of each subdomain problem. Table 1 should not be used to compare MG and DD methods for efficiency. Similar experiments [14] with more carefully optimized DD and MG methods show DD to be often competitive with MG for difficult elliptic equations such as those with discontinuous coefficients, although there may be some debate as to which approach is more effective on parallel computers [22].

Multiplicative multigrid. The results for multiplicative V-cycle MG are presented in Table 2. Each row corresponds to a different smoothing strategy, and is annotated by (ν_1, ν_2) , with ν_1 : pre-smoothing strategy, and ν_2 : post-smoothing strategy. An “f” indicates the use of a single forward Gauss-Seidel sweep, while a “b” denotes the use of the adjoint of the latter, i.e., a backward Gauss-Seidel sweep. $(\nu_1, \nu_2) = (ff, fb)$, for example, corresponds to two forward Gauss-Seidel pre-smoothing steps, and a symmetric (forward/backward) post-smoothing step. Two series of results are given. For the first set, we explicitly imposed the Galerkin conditions when constructing the coarse op-

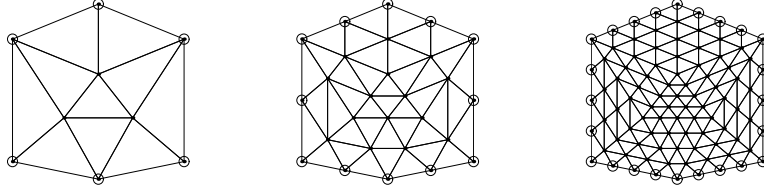


FIG. 1. *Example 1: Nested finite element meshes for MG.*

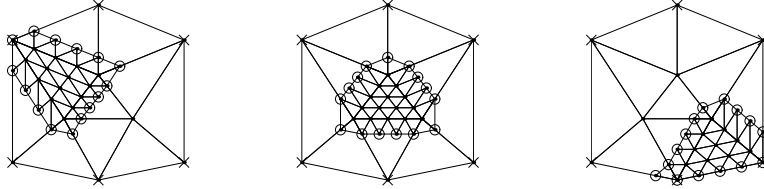


FIG. 2. *Example 1: Overlapping subdomains for DD.*

erators. In this case, the multigrid algorithm is guaranteed to converge by Theorem 3.6. In the second series of tests, corresponding to the numbers in parentheses, the coarse mesh operators are constructed using standard finite element discretization. In that case, Galerkin conditions are not satisfied everywhere due to coefficient discontinuities appearing within coarse elements; hence, the MG method may diverge (DIV).

The unaccelerated MG results clearly illustrate the symmetry penalty discussed in §5. The nonsymmetric methods are always superior to the symmetric ones (the cases (f,b), (ff,bb), and (fb,fb)). Note that minimal symmetry (ff,bb) leads to a better convergence than maximal symmetry (fb,fb). The correctness of Lemma 5.1 is illustrated by noting that two iterations of the (f,0) strategy are actually faster than one iteration of the (f,b) strategy; also, compare the (ff,0) strategy to the (ff,bb) one. CG-acceleration leads to a guaranteed reduction in iteration count for the symmetric preconditioners (see Lemma 2.15). We observe that the unaccelerated method need not be convergent for CG to be effective (recall Remarks 2.1 and 4.7, and the (f,b) result). CG appears to accelerate also some non-symmetric linear methods. Yet, it seems difficult to predict failure or success beforehand in such cases. The most robust method appears to be the Bi-CGstab method. The number of iterations with this method depends only marginally on the symmetric or nonsymmetric nature of the linear method. Note the tendency to favor the nonsymmetric V-cycle strategies. Overall, the fastest method proves to be the Bi-CGstab-acceleration of a (very nonsymmetric) V(1,0)-cycle.

Multiplicative domain decomposition. Some numerical results for multiplicative DD with different subdomain solvers, and different subdomain sweeps are given in Table 3. In the column “forw”, the iteration counts reported were obtained with a single sweep though the subdomains on each multiplicative DD iteration. The other columns correspond to a symmetric forward/backward sweep or to two forward sweeps. Four different subdomain solvers are used: an *exact* solve, a *symmetric* method consisting of two symmetric Gauss-Seidel iterations, a *nonsymmetric* method consisting of four Gauss-Seidel iterations, and, finally, a method using four forward Gauss-Seidel

iterations in the forward subdomain sweep and using their *adjoint*, i.e., four backward Gauss-Seidel iterations, in the backward subdomain sweep. The latter leads to an symmetric iteration; see Remarks 3.3 and 3.4. Note that the cost of the three inexact subdomain solvers is identical.

Although apparently not as sensitive to operator symmetries as MG, the same conclusions can be drawn for DD as for MG. In particular, the symmetry penalty is seen for the pure DD results. Lemma 5.1 is confirmed since two iterations in the column “forw” are always more efficient than one iteration of the corresponding symmetrized method in column “forw/back”. The CG results indicate that using minimal symmetry (the “adjointed” column) is a more effective approach than the fully symmetric one (the “symmetric” column). Again, the most robust acceleration is the Bi-CGstab one.

Additive multigrid. Results obtained with an additive multigrid method are reported in Table 4. The number and nature of the smoothing strategy is given in the first column of the table.

In the case of an unaccelerated additive method, the selection of a good damping parameter is crucial for convergence of the method. We did not search extensively for an optimal parameter; a selection of $\omega = 0.45$ seemed to provide good results in the case when the coarse problem is variationally defined. No ω -value leading to satisfactory convergence was found in the case when the coarse problem is obtained by discretization. In the case of CG acceleration the observed convergence behavior was completely independent of the choice of ω ; see Remarks 3.3 and 3.4. The symmetric methods ($\nu = fb, ffb, fbf$) are accelerated very well. Some of the nonsymmetric methods are accelerated too, especially when the number of smoothing steps is sufficiently large. In the case of Bi-CGstab-acceleration, there appeared to be a dependence of convergence on ω (only with use of non-variational coarse problem). In that case we took $\omega = 1$. The overall best method appears to be the Bi-CGstab acceleration of the nonsymmetric multigrid method with a single forward Gauss-Seidel sweep on each grid-level.

Additive domain decomposition. The results for additive DD are given in Table 5. The subdomain solver is either an exact solver, a symmetric solver based on two symmetric (forward/backward) Gauss-Seidel sweeps, or a nonsymmetric solver based on four forward Gauss-Seidel iterations.

No value of ω was found that led to satisfactory convergence of the unaccelerated method. CG-acceleration performs well when the linear method is symmetric; it performs less well for the nonsymmetric method. Again, the best overall method is the Bi-CGstab-acceleration of the nonsymmetric additive solver.

6.2. Example 2. The second test problem is the Laplace equation on a semi-adapted L-shaped domain, with Dirichlet boundary conditions chosen in such a way that the equation has the following solution (in polar coordinates):

$$u(r, \theta) = \sqrt{r} \sin(\theta/2) .$$

Note that the one-point Gaussian quadrature rule which we employ to construct the stiffness matrix entries is an exact integrator here. Hence, the variational conditions (11)

TABLE 2

Example 1: Multiplicative MG with variational (discretized) coarse problem

ν_1	ν_2	UNACCEL	CG	Bi-CGstab
f	0	65 (DIV)	$\gg 100$ ($\gg 100$)	14 (16)
f	b	55 (DIV)	16 (18)	10 (15)
f	f	40 (31)	30 ($\gg 100$)	9 (9)
ff	0	39 (48)	$\gg 100$ ($\gg 100$)	8 (10)
fb	0	53 (DIV)	$\gg 100$ ($\gg 100$)	10 (11)
0	ff	39 (29)	29 ($\gg 100$)	8 (9)
0	fb	53 (DIV)	17 (99)	10 (12)
fb	fb	34 (27)	12 (13)	8 (8)
ff	bb	28 (18)	11 (11)	7 (7)
ff	ff	24 (15)	12 (12)	6 (6)
fff	f	24 (15)	17 (27)	6 (6)
fff	0	25 (17)	$\gg 100$ ($\gg 100$)	7 (6)

TABLE 3

Example 1: Multiplicative DD with variational (discretized) coarse problem

Accel.	subdomain solve	forw	forw/back	forw/forw
UNACCEL	exact	40 (42)	38 (39)	20 (21)
	symmetric	279 (282)	146 (149)	140 (141)
	adjointed	–	110 (112)	102 (103)
	nonsymmetric	189 (191)	102 (104)	95 (96)
CG	exact	$\gg 500$ ($\gg 500$)	13 (13)	20 (20)
	symmetric	140 (56)	24 (24)	29 (27)
	adjointed	– –	21 (21)	25 (26)
	nonsymmetric	135 (83)	22 (23)	28 (28)
Bi-CGstab	exact	9 (9)	9 (9)	6 (6)
	symmetric	23 (23)	17 (16)	16 (16)
	adjointed	– –	14 (14)	14 (13)
	nonsymmetric	19 (20)	13 (13)	13 (13)

TABLE 4

Example 1: Additive MG with variational (discretized) coarse problem

ν	UNACCEL	CG	Bi-CGstab
f	175 ($\gg 1000$)	$\gg 100$ ($\gg 100$)	23 (52)
ff	110 ($\gg 1000$)	119 (168)	19 (43)
fb	146 ($\gg 1000$)	34 (54)	23 (49)
fff	95 ($\gg 1000$)	28 (67)	17 (37)
ffbb	100 ($\gg 1000$)	27 (47)	17 (34)
fbfb	95 ($\gg 1000$)	28 (48)	20 (43)

TABLE 5
Example 1: Additive DD with variational (discretized) coarse problem

subdomain solve	UNACCEL	CG	Bi-CGstab
exact	$\gg 1000$ ($\gg 1000$)	34 (34)	25 (27)
symmetric	$\gg 1000$ ($\gg 1000$)	57 (57)	50 (49)
nonsymmetric	$\gg 1000$ ($\gg 1000$)	69 (65)	38 (41)

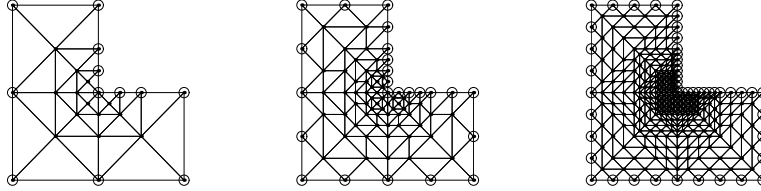


FIG. 3. *Example 2: Nested finite element meshes for MG.*

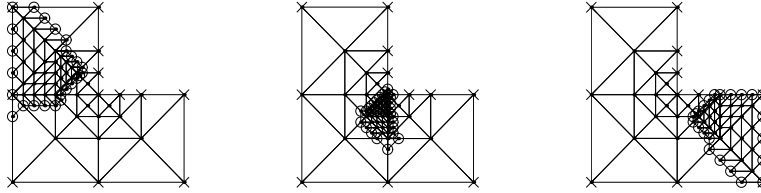


FIG. 4. *Example 2: Overlapping subdomains for DD.*

hold automatically between the fine space and all subdomain and coarse spaces for both the MG and the DD methods.

Figure 3 shows a nested sequence of uniform mesh refinements used to formulate the MG methods. Figure 4 shows several overlapping subdomains constructed from a piece of the fine mesh of 9216 elements (4705 nodes) overlaying the initial coarse mesh of 36 elements (25 nodes).

Multiplicative Methods. The results for multiplicative MG are given in Table 6, whereas the results for multiplicative DD are given in Table 7. The results are similar to those for Example 1; in particular, imposing minimal symmetry is the most effective CG-accelerated approach to the problem. Employing the least symmetric linear method alone is the most effective linear method, and the same nonsymmetric linear method yields the most effective Bi-CGstab-accelerated approach.

Additive Methods. As for Example 1, in the case of the unaccelerated additive methods the selection of the damping parameter was crucial for convergence of the methods. We did not search extensively for an optimal parameter; a selection of $\omega = 0.45$ seemed to provide acceptable results for DD. Note that improved convergence behavior might be obtained by allowing different ω values for each subdomain solver (this will not be further investigated here). No satisfactory value for ω was found for additive MG. In the case of CG acceleration, the observed convergence behavior was

completely independent of the choice of ω . However, again in the case of the additive methods with discretized (non-variational) coarse problems accelerated by Bi-CGstab, there was convergence rate dependence on ω . For uniform comparisons we took $\omega = 1$ in those cases. The results for additive MG are given in Table 8, whereas the results for additive DD are given in Table 9. The effect of the symmetry of the linear method's error propagator on its convergence, and on the convergence behavior of CG and Bi-CGstab, was as for Example 1.

7. Concluding remarks. In this paper, we have developed a preconditioning theory for additive and multiplicative Schwarz methods. We established sufficient conditions which guarantee that abstract multiplicative and additive algorithms yield SPD preconditioners. We then analyzed four specific methods: MG and DD methods, in both their additive and multiplicative forms. In all four cases, we used the general theory to establish sufficient conditions that guarantee the resulting preconditioner is SPD. As discussed in Remarks 3.4, 3.6, 4.8, and 4.9, the sufficient conditions for the theory, in the case of all four methods, are easily satisfied for non-variational, and *even non-convergent* methods. The analysis shows that by simply taking some care in the way a Schwarz method is formulated, one can guarantee that the method is convergent when accelerated with the conjugate gradient method. These results hold for finite difference or finite element-based methods, even if variational conditions are violated.

We also investigated the role of symmetry in linear methods and preconditioners. A certain penalty lemma (Lemma 5.1) was stated and proved, illustrating why symmetrizing is actually a bad idea for linear methods. It was conjectured that enforcing minimal symmetry in a linear preconditioner achieves the best results when combined with the conjugate gradient method, and our numerical examples illustrate this behavior almost uniformly. A sequence of experiments with two non-trivial test problems showed that the most efficient approach may be to abandon symmetry in the preconditioner altogether, and to employ a nonsymmetric solver such as Bi-CGstab. While acceleration with CG was strongly dependent on the symmetric nature of the preconditioner, Bi-CGstab always converged rapidly. In addition, Bi-CGstab appeared to benefit from the behavior predicted by Lemma 5.1, namely that a nonsymmetric linear preconditioner should have better convergence properties than its symmetrized form.

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TABLE 6
Example 2: Multiplicative MG

ν_1	ν_2	UNACCEL	CG	Bi-CGstab
f	b	23	12	7
f	f	19	22	6
ff	0	21	$\gg 100$	7
fb	0	25	$\gg 100$	8
0	ff	20	42	7
0	fb	23	17	8
fb	fb	16	9	6
ff	bb	15	9	5
ff	ff	14	9	5
fff	f	14	12	5
fff	0	16	36	5
f	0	33	$\gg 100$	11

TABLE 7
Example 2: Multiplicative DD

Accel.	subdomain solve	forw	forw/back	forw/forw
UNACCEL	exact	73	60	37
	symmetric	402	205	207
	adjointed	–	153	146
	nonsymmetric	267	144	134
CG	exact	116	17	17
	symmetric	164	37	38
	adjointed	–	32	33
	nonsymmetric	121	31	32
Bi-CGstab	exact	11	11	7
	symmetric	37	25	26
	adjointed	–	22	23
	nonsymmetric	27	21	21

TABLE 8
Example 2: Additive MG

ν	UNACCEL	CG	Bi-CGstab
f	91	$\gg 1000$	21
ff	62	31	16
fb	74	29	18
fff	126	25	14
ffbb	136	27	15
fbfb	98	27	15

TABLE 9
Example 2: Additive DD

subdomain solve	UNACCEL	CG	Bi-CGstab
exact	$\gg 1000$	42	29
symmetric	$\gg 1000$	86	56
nonsymmetric	$\gg 1000$	82	49

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