

## ANALYSIS AND COMPARISON OF GEOMETRIC AND ALGEBRAIC MULTIGRID FOR CONVECTION-DIFFUSION EQUATIONS\*

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**Abstract.** The discrete convection-diffusion equations obtained from streamline diffusion finite element discretization are solved on both uniform meshes and adaptive meshes. Estimates of error reduction rates for both geometric multigrid (GMG) and algebraic multigrid (AMG) are established on uniform rectangular meshes for a model problem. Our analysis shows that GMG with line Gauss–Seidel smoothing and bilinear interpolation converges if  $h \gg \epsilon^{2/3}$ , and AMG with the same smoother converges more rapidly than GMG if the interpolation constant  $\beta$  in the approximation assumption of AMG satisfies  $\beta \ll (\frac{h}{\sqrt{\epsilon}})^\alpha$ , where  $\alpha = \begin{cases} 1, & h < \sqrt{\epsilon}, \\ 2, & h \geq \sqrt{\epsilon}. \end{cases}$  On unstructured triangular meshes, the performance of GMG and AMG, both as solvers and as preconditioners for GMRES, are evaluated. Numerical results show that GMRES with AMG preconditioning is a robust and reliable solver on both type of meshes.

**Key words.** convection-diffusion equations, multigrid, adaptive mesh refinement

**AMS subject classifications.** Primary, 65N30, 65N50, 65N55; Secondary, 65F10

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**1. Introduction.** The purpose of this paper is to evaluate different solution strategies for the linear systems obtained from discretization of the convection-diffusion equation

$$(1.1) \quad \begin{cases} -\epsilon \Delta u + b \cdot \nabla u = f & \text{on } \Omega, \\ u = g & \text{on } \partial\Omega, \end{cases}$$

where  $b$  and  $f$  are sufficiently smooth and the domain  $\Omega$  is convex with Lipschitz boundary  $\partial\Omega$ . We are interested in the convection-dominated case, i.e.,  $|b| \gg \epsilon$ . In this setting, the solution typically has steep gradients in some parts of the domain. These may take the form of boundary layers caused by Dirichlet conditions on the outflow boundaries or internal layers caused by discontinuities in the inflow boundaries.

Let  $\mathfrak{S}_h$  be a given quasi-uniform mesh of triangles on  $\Omega$  and let  $V_h$  be the linear finite element space on  $\mathfrak{S}_h$ . It is well known that the standard Galerkin finite element discretization on uniform grids produces inaccurate oscillatory solutions to convection-diffusion problems. Here, we will discretize (1.1) on  $V_h$  using the streamline diffusion finite element method (SDFEM) [7], a variant of the standard Galerkin method, where extra diffusion in the streamline direction is introduced. The SDFEM formulation entails finding  $u_h \in V_h$  such that

$$(1.2) \quad B_{sd}(u_h, v_h) = f_{sd}(v_h) \quad \text{for all } v_h \in V_h,$$

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where

$$B_{sd}(u_h, v_h) = \epsilon(\nabla u_h, \nabla v_h) + (b \cdot \nabla u_h, v_h) + \sum_{T \in \mathfrak{S}_h} \delta_T (b \cdot \nabla u_h, b \cdot \nabla v_h)_T,$$

$$f_{sd}(v_h) = (f_h, v_h) + \sum_{T \in \mathfrak{S}_h} (f, \delta_T b \cdot \nabla v_h),$$

and  $\delta_T$  is a stabilization parameter. With a carefully chosen value of  $\delta_T$ , the streamline diffusion finite element discretization is able to eliminate most oscillations and produce accurate solutions in the regions where no layers are present [7]. As shown in [6], a good choice of  $\delta_T$  is

$$(1.3) \quad \delta_T = \begin{cases} \frac{1}{2\|b\|_T} \left(1 - \frac{1}{P_{e_T}}\right) & \text{if } P_{e_T} > 1, \\ 0 & \text{if } P_{e_T} < 1, \end{cases}$$

where

$$P_{e_T} = \frac{\|b\|_T h_T}{2\epsilon} \quad \text{for } T \in \mathfrak{S}_h \text{ with diameter } h_T$$

is the mesh Peclet number.

This strategy does not produce accurate solutions in regions containing layers that are not resolved by the grid. Accuracy can be achieved at reasonable cost in such regions by adaptive mesh refinement, in which an a posteriori error estimation strategy is used to identify regions where errors are large, and a marking strategy is used to select elements to be refined. An adaptive solution strategy is to start with a coarse mesh, compute the discrete system, solve the linear system, and then locally refine the mesh using the information provided by an a posteriori error estimation. This adaptive solution process can be applied repeatedly until the a posteriori error estimator is less than a prescribed tolerance or the maximal number of mesh refinement steps is reached.

In this paper, we use the local error estimator proposed by Kay and Silvester [8], with the maximum marking strategy [11], and regular mesh refinement strategy for the adaptive refinement process. We are concerned with the costs of performing this computation with emphasis on the costs of solving the discrete systems that arise at each step of the mesh refinement. Let  $A_h u_h = f_h$  denote the discrete system of equations derived from (1.2). Since adaptive mesh refinement produces a sequence of nested meshes, multigrid methods are natural candidates for solving this linear system. Here, we would like to explore the effectiveness of the geometric multigrid (GMG) [5] and algebraic multigrid (AMG) [17] methods. Since the linear system is nonsymmetric, Krylov subspace linear solvers such as the generalized minimal residual method (GMRES) [18] are also good choices. In order to accelerate the convergence of the Krylov subspace linear solvers, good preconditioners are needed. We would also like to investigate the performance of GMRES preconditioned by GMG and AMG.

This paper is organized as follows: Details of the error estimator and refinement strategy are given in section 2 where numerical studies that demonstrate the efficiency of the error indicator are presented. In section 3, multigrid algorithms and some estimates of two-level multigrid convergence are analyzed and numerical results that support our theoretical analysis are given. A comparison of the solvers is shown in section 4. This comparison examines both standalone versions of multilevel multigrid as well as versions in which multigrid is used as a preconditioner for GMRES. We also

compare these with unpreconditioned GMRES, and with a preconditioned GMRES algorithm in which the smoother used for multigrid is used as a preconditioner. These experiments are performed using the following two benchmark problems.

*Problem 1: Constant flow with characteristic and downstream layers.* Equation (1.1) is given with the coefficient  $b = (\sin(\phi), \cos(\phi))$ , where  $\phi$  measures the angle of the flow direction from the  $y$ -axis, and the right-hand side  $f = 0$  on the domain  $\Omega = [-1, 1] \times [-1, 1]$ . The Dirichlet boundary condition is set as  $g = 1$  on the segments  $y = -1 \cap x > 0$  and  $x = 1$ , and  $g = 0$  elsewhere.

*Problem 2: Flow with closed characteristics:* Here, the coefficient vector  $(b_1, b_2)$  is  $(2y(1 - x^2), 2x(1 - y^2))$  and the right-hand side  $f = 0$  on the domain  $\Omega = [-1, 1] \times [-1, 1]$ . The Dirichlet boundary condition is  $g = 1$  on the segments  $y = 1$  and  $g = 0$  elsewhere.

These examples are chosen because they exhibit many of the important features of flow problems so that their effect on discretization and multigrid solvers can be identified. In particular, Problem 1 has an exponential boundary layer at the outflow boundary and a characteristic internal layer induced by a discontinuity on the inflow boundary. Both of these can be treated using adaptive mesh refinement. Moreover, by varying the angle  $\phi$ , we can assess the effect of flow direction on performance of solvers. Problem 2 also has a (characteristic) boundary layer caused by discontinuous boundary conditions, and its recirculating flow requires that care be taken in choosing smoothers for multigrid computations.

Finally, in section 5, we draw our conclusions.

**2. Adaptive mesh refinement by a posteriori error estimation.** One common technique to increase the accuracy of the finite element solution is mesh refinement. Unlike uniform mesh refinement, the adaptive mesh refinement process refines meshes only in the regions where errors between the weak solution of the partial differential equation and the corresponding finite element solution are large. In general, the adaptive mesh refinement process consists of loops of the following form:

$$\underbrace{\text{Solve}}_1 \rightarrow \underbrace{\text{Compute error indicator}}_2 \rightarrow \underbrace{\text{Refine mesh}}_3$$

In step 3, elements in which the value of the a posteriori error estimator is large are marked for refinement according to some element selection algorithm. On a given triangulation  $\mathfrak{S}_h$ , for any element  $T \in \mathfrak{S}_h$ , let  $\eta_T$  be the a posteriori error indicator of element  $T$ . A heuristic marking strategy is the *maximum marking strategy* where an element  $T^*$  is marked for refinement if

$$(2.1) \quad \eta_{T^*} > \theta \max_{T \in \mathfrak{S}_h} \eta_T,$$

with a prescribed threshold  $0 < \theta < 1$ . For other marking strategies, we refer to [11]. Once decisions on where to refine are made, commonly used mesh refinement schemes are regular refinement, where a triangle is divided into four triangles equally, and longest side bisection [15], [16], where a triangle is divided into two triangles by adding a new node to the midpoint of the longest edge. The maximum marking strategy and regular refinement scheme are used in our experiments.

A reliable computable a posteriori error estimator in step 2 is the key for the adaptive mesh refinement process to succeed. For the convection-diffusion equation discretized by SDFEM, the first a posteriori error estimation where the error is estimated by computing the residual was proposed by Verfürth [21], and the first a posteriori error estimation where the error is estimated by solving a local problem was

developed by Kay and Silvester [8]. In this study, we use the Kay and Silvester a posteriori error estimator, which we have found to be a more effective choice in [23]. Hereafter, we call this indicator the KS-indicator. First, let us introduce the following abbreviations. Let  $\|\cdot\|_{0,\Omega}$  and  $\|\cdot\|_{0,T}$  denote the  $L^2$  norm on domain  $\Omega$  and element  $T$ , respectively. Let  $\mathcal{E}(T)$  be the set of edges of  $T$  and  $\omega_T = \cup_{T' \cap T \in \mathcal{E}(T)} T'$ , i.e.,  $\omega_T$  is the set of triangles  $T'$  that share a common edge with  $T$ . Let  $P_T^0$  be the  $L^2$  projection onto the space of polynomials of degree 0 on element  $T$ . The interior residual  $R_T$  of element  $T$  and the interelement flux jump  $R_E$  of edge  $E$  are defined as follows:

$$\begin{aligned} R_T &= (f - b \cdot \nabla u_h)|_T, \\ R_T^0 &= \mathcal{P}_T^0(R_T), \\ R_E &= \begin{cases} [\frac{\partial u_h}{\partial n_E}]_E & \text{if } E \in \Omega, \\ 0 & \text{if } E \in \partial\Omega, \end{cases} \end{aligned}$$

where  $[\cdot]_E$  is the jump across edge  $E$ .<sup>1</sup> Let  $\Phi$  be the element affine mapping from the physical domain to the computational domain and let  $\chi_i$  be the nodal basis function of node  $i$ . The approximation space is denoted as  $Q_T = Q_T \oplus B_T$ , where

$$Q_T = \text{span}\{\psi_E \circ \Phi^{-1} \mid \psi_E = 4\chi_i\chi_j, \ i, j \text{ are the endpoints of } E \text{ and } E \in \partial T \cap (\Omega \cup \Gamma_N)\}$$

is the space spanned by quadratic edge bubble functions and

$$B_T = \text{span}\left\{ \psi_T \circ \Phi^{-1} \mid \psi_T = 27 \prod_{i=1}^3 \chi_i \right\}$$

is the space spanned by cubic interior bubble functions. For details, see [8]. On each element  $T$ , the error estimator is then given by  $\eta_{h,T} = \|\nabla e_T\|_{0,T}$ , where  $e_T \in Q_T$  satisfies

$$(2.2) \quad \epsilon(\nabla e_T, \nabla v)_T = (R_T^0, v)_T - \frac{1}{2}\epsilon \sum_{E \in \partial T} (R_E, v)_E.$$

Let  $e_h = u - u_h$ . The a posteriori error estimation is specified as follows:

(global upper bound)

$$(2.3) \quad \|\nabla(e_h)\|_{0,\Omega} \leq C \left( \sum_{T \in \mathfrak{S}_h} \eta_{h,T}^2 + \sum_{T \in \mathfrak{S}_h} \left(\frac{h}{\epsilon}\right)^2 \|R_T - R_T^0\|_{0,T}^2 \right)^{1/2}$$

(local lower bound)

$$(2.4) \quad \eta_{h,T} \leq c \left( \|e_h\|_{0,\omega_T} + \sum_{T \subset \omega_T} \frac{h_T}{\epsilon} \|b \cdot \nabla e_h\|_{0,T} + \sum_{T \subset \omega_T} \frac{h_T}{\epsilon} \|R_T - R_T^0\|_{0,T} \right),$$

where constants  $C$  and  $c$  are independent of the diffusion parameter  $\epsilon$  and mesh size  $h$ .

<sup>1</sup>This definition of  $R_E$  for  $E \in \partial\Omega$  is for Dirichlet boundary conditions. We don't consider Neumann conditions in this study; in case  $E \in \partial\Omega$  and a Neumann condition holds, the flux jump  $R_E$  would be set to  $-2(\frac{\partial u_h}{\partial n_E})$  [8].

The sharpness of the error indicator can be revealed by computing the local effectivity index,

$$E_T = \max_{T \in \mathfrak{S}_h} \frac{\eta_{h,T}}{|u - u_h|_{1,T}},$$

and the global effectivity index,

$$E_\Omega = \frac{(\sum_{T \in \mathfrak{S}_h} \eta_{h,T}^2)^{1/2}}{|u - u_h|_{1,\Omega}},$$

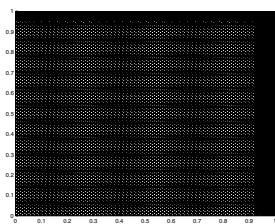
where  $|\cdot|_{1,\Omega}$  and  $|\cdot|_{1,T}$  represent the  $H^1$  seminorms on domain  $\Omega$  and element  $T$ , respectively. If both indices are close to 1, the error indicator is reliable and efficient. The global upper bound provides an estimate of the error over the whole domain, and the local lower bound locates where the error is large. It has been shown that  $E_\Omega$  grows in proportion to  $\sqrt{Pe}$  and  $E_T$  grows in proportion to  $Pe$  for the KS-indicator [8]. The following numerical results support this conclusion.

Consider here equation (1.1) on the domain  $\Omega = [0, 1] \times [0, 1]$  with  $b = (\beta_1, \beta_2) = (\sin 75^\circ, \cos 75^\circ)$  and the Dirichlet boundary condition is obtained from the analytic solution

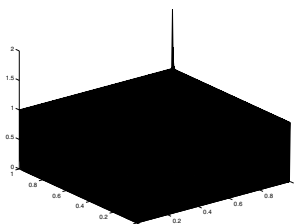
$$(2.5) \quad u(x, y) = \frac{e^{\beta_1 x/\varepsilon} - 1}{e^{\beta_1/\varepsilon} - 1} + \frac{e^{\beta_2 y/\varepsilon} - 1}{e^{\beta_2/\varepsilon} - 1}.$$

Both global effectivity index  $E_\Omega$  and local effectivity index  $E_T$  are computed. In order to see how the effectivity indices change in terms of the diffusion parameter  $\varepsilon$  and mesh size  $h$ , the problem is solved over uniform meshes with mesh size  $h = \frac{1}{8}, \frac{1}{16}, \frac{1}{32}$ , and  $\frac{1}{64}$  for  $\varepsilon = \frac{1}{64}, \frac{1}{256}, \frac{1}{1024}$ , and  $\frac{1}{4096}$ . For accurately approximating the true error  $u - u_h$ , a very fine mesh  $\mathfrak{S}_f$ , as shown in Figure 2.1, consisting of 11272 nodes and 21379 elements is generated by first solving the above test problem with  $\varepsilon = 10^{-3}$  on a  $64 \times 64$  mesh followed by three mesh refinement steps where a large threshold  $\theta = 0.75$  in the maximum marking strategy (2.1) is used to prevent unnecessary over refinement. An approximation of the exact error is then computed on  $\mathfrak{S}_f$  by

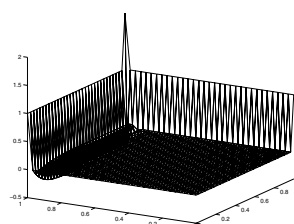
$$|u - u_h|_{1,\Omega^*} = \left( \sum_{\tau \in \mathfrak{S}_f \cup \Omega^*} |u - u_h|_{1,\tau}^2 \right)^{1/2} \quad \text{for } \Omega^* = \Omega \text{ or } T \in \mathfrak{S}_h,$$



(a) Mesh  $\mathfrak{S}_f$ .



(b) Solution  $u$  on  $\mathfrak{S}_f$ .



(c) SDFEM solution  $u_h$  on  $\mathfrak{S}_h$ , where  $h = \frac{1}{32}$ .

FIG. 2.1.

TABLE 2.1  
Effectivity indices of the KS-indicator.

$\epsilon$	$8 \times 8$	$16 \times 16$	$32 \times 32$	$64 \times 64$
$\frac{1}{64}$	1.156	0.951	0.979	1.022
$\frac{1}{256}$	2.044	1.457	1.105	0.929
$\frac{1}{1024}$	4.016	2.772	1.952	1.422
$\frac{1}{4096}$	8.470	5.842	4.075	2.867

$\epsilon$	$8 \times 8$	$16 \times 16$	$32 \times 32$	$64 \times 64$
$\frac{1}{64}$	2.504	1.714	1.687	1.557
$\frac{1}{256}$	9.242	4.637	2.536	1.750
$\frac{1}{1024}$	36.95	18.48	9.239	4.629
$\frac{1}{4096}$	147.8	73.90	36.95	18.48

(a) Global index  $E_\Omega$ .

(b) Local index  $E_T$ .

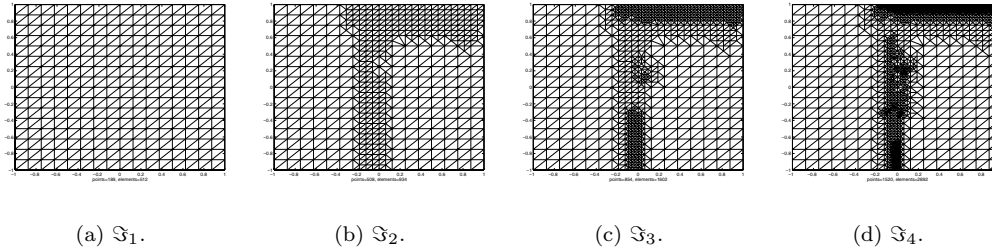


FIG. 2.2. Adaptively refined mesh with threshold value  $\theta = 0.01$ .

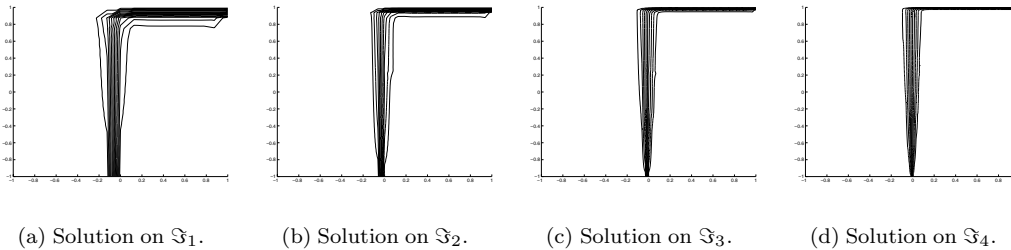


FIG. 2.3. Contour plots of the solutions of Problem 1 on adaptively refined meshes starting with a uniform  $16 \times 16$  mesh.

where the discrete true solution  $u$  is evaluated directly from (2.5) on  $\mathfrak{S}_f$  and the SDFEM solution  $u_h$  is prolonged from  $\mathfrak{S}_h$  onto  $\mathfrak{S}_f$  by standard bilinear interpolation. In Table 2.1, for  $h \gg \epsilon$ ,  $E_\Omega$  and  $E_T$  decrease by factors approximately of size  $\sqrt{2}$  and 2, respectively, as the mesh size is halved. Also,  $E_\Omega$  and  $E_T$  double and quadruple, respectively, when  $\epsilon$  is reduced by a factor of four. Thus,  $E_T = O(P_e)$  and  $E_\Omega = O(\sqrt{P_e})$ , as shown in [8].

Although there is deterioration in reliability and efficiency for these error indicators, the regions with large errors can still be located when a good threshold value  $\theta$  is chosen. Next, we give a representative picture of how the adaptive solution process improves the solution quality. We use the KS-indicator for mesh refinement in solving Problem 1 with the diffusion parameter  $\epsilon = 10^{-3}$  and  $\phi = 0$ . From the local lower bound (2.4), we suspect that the value of the indicator  $\eta_{h,T}$  in the outflow boundary layer regions is  $O(\frac{h}{\epsilon})$  larger than its value in the internal layer regions. Therefore, for the case considered here, a threshold value  $\theta = 0.01 \approx \frac{\epsilon}{h}$  is chosen so that both the boundary and internal layer regions can be refined by the maximum marking strategy (2.1). A sequence of adaptively refined meshes is shown in Figure 2.2 in which ele-



where the block tridiagonal matrices

$$\begin{aligned}
 (3.2) \quad D &= h \times \text{tridiag} \left[ \frac{1}{6} - \frac{1}{3} \frac{\epsilon}{h}, \frac{2}{3} + \frac{8}{3} \frac{\epsilon}{h}, \frac{1}{6} - \frac{1}{3} \frac{\epsilon}{h} \right], \\
 L &= h \times \text{tridiag} \left[ \frac{1}{6} + \frac{1}{3} \frac{\epsilon}{h}, \frac{2}{3} + \frac{1}{3} \frac{\epsilon}{h}, \frac{1}{6} + \frac{1}{3} \frac{\epsilon}{h} \right], \\
 U &= \epsilon \times \text{tridiag} \left[ \frac{1}{3}, \frac{1}{3}, \frac{1}{3} \right].
 \end{aligned}$$

In the following, we describe each component of the multigrid methods used here. First, we use the same smoother in both GMG and AMG. On uniform rectangular meshes, one step of a horizontal line Gauss–Seidel (HLGS) smoother is used. The HLGS method uses the block lower triangle of  $A_k$  as a splitting operator. (Thus, the HLGS sweep can be viewed as proceeding along horizontal lines, one line at a time, in the vertical direction.) Second, in GMG, (1) the coarse grids are obtained directly from the mesh refinement process; (2) the grid transfer is via linear interpolation for prolongation, and the restriction operator  $I_k^{k-1}$  is then taken to be the transpose of the prolongation operator  $I_{k-1}^k$ ; and (3) the matrix  $A_k$  is obtained directly from SDFEM discretization in the refinement process. For AMG, we use the method developed by Ruge and Stüben [17] (AMG): (1) a coarse grid at level  $k - 1$  is generated by coarsening the fine grid along the so-called strong connection direction of matrix graph  $\mathcal{G}$  of  $A_k = (a_{i,j})$  where a directive edge  $\vec{e}_{i,j} \in \mathcal{G}$  is defined to be strongly connected from node  $i$  to node  $j$  if  $\mu \leq \frac{-a_{i,j}}{\max_{m \neq i} (-a_{i,m})}$  for a given parameter  $0 < \mu < 1$ ; (2) the interpolation operator  $I_{k-1}^k$  is defined dynamically by a sophisticated algebraic formula during the AMG coarsening process [17]; and (3) the matrix  $A_{k-1}$  for a coarse grid is computed by  $I_k^{k-1} A_k I_{k-1}^k$ . The above process is repeated until all coarse grids are generated.

Next, we show the convergence behavior of 2-level V-cycle multigrid for Problem 1. Let  $E_k^s$  be the error reduction operator of the HLGS smoother and let  $E_k^c = I - I_{k-1}^k A_{k-1}^{-1} I_k^{k-1} A_k$  be the coarse grid correction operator. The error reduction operators of multigrid methods can then be written as  $E_k^s E^c E_k^s$ . The notation  $E_k^{mg}$  and  $E_k^{amg}$  denotes the error reduction operators of geometric multigrid and the algebraic multigrid, respectively. The notation  $\|\cdot\|$  denotes the usual discrete  $L^2$  norm for any vector and matrix. The notation  $\|\cdot\|_0$  and  $\|\cdot\|_1$  represents the continuous  $L^2$  and  $H^1$  norms. Variants of these norms are specified using matrix notation in (3.17) below.

In order to show convergence, the following auxiliary lemmas are needed.

LEMMA 3.1. *Given two symmetric matrices  $B_1$  and  $B_2$ , assume that  $B_1, B_2 \geq 0$ ,  $B_1$  is irreducible, and  $B_2$  is positive definite. The following properties hold.*

1. *There exists a positive eigenvector  $x^+$  of  $B_2^{-1} B_1$  such that*

$$(3.3) \quad B_2^{-1} B_1 x^+ = \rho(B_2^{-1} B_1) x^+.$$

2. *If  $\alpha I - B_2^{-1} B_1$  is nonsingular and  $(\alpha I - B_2^{-1} B_1)^{-1} \geq 0$ , then  $\rho(B_2^{-1} B_1) < \alpha$ .*

*Proof.* The existence of a positive eigenvector satisfying (3.3) is essentially a generalization of the well-known Perron–Frobenius theorem [20, Theorem 2.7]. Using (3.3), one can prove the second result using a standard argument of the Perron–Frobenius theory; see Theorem 3.16 in [20].  $\square$

LEMMA 3.2. *Let  $L$  and  $D$  be the matrices defined in (3.2). For any  $\delta \geq (1 + \frac{2\epsilon}{h}) \frac{h}{\epsilon}$ , the matrix  $\delta(D - L) - D$  is an M-matrix.*



*Proof.* Let us choose  $\delta = \frac{h\gamma}{\epsilon}$  for some  $\gamma > 0$ . From (3.2),  $D - L = \frac{\epsilon}{3} \times \text{tridiag}[-2, 7, -2]$ . Therefore, we have

$$\begin{aligned} \delta(D - L) - D &= h \times \text{tridiag} \left[ \frac{-2\gamma}{3}, \frac{7\gamma}{3}, \frac{-2\gamma}{3} \right] - h \times \text{tridiag} \left[ \frac{1}{6} - \frac{\epsilon}{3h}, \frac{2}{3} + \frac{8\epsilon}{h}, \frac{1}{6} - \frac{\epsilon}{3h} \right] \\ &= h \times \text{tridiag} \left[ -\left( \frac{2\gamma}{3} + \frac{1}{6} - \frac{\epsilon}{3h} \right), \frac{7\gamma}{3} - \frac{2}{3} - \frac{8\epsilon}{3h}, -\left( \frac{2\gamma}{3} + \frac{1}{6} - \frac{\epsilon}{3h} \right) \right]. \end{aligned}$$

Since

$$\frac{7\gamma}{3} - \frac{2}{3} - \frac{8\epsilon}{3h} - 2 \left( \frac{2\gamma}{3} + \frac{1}{6} - \frac{\epsilon}{3h} \right) = \gamma - 1 - \frac{2\epsilon}{h},$$

clearly, for  $\gamma \geq 1 + \frac{2\epsilon}{h}$ , the matrix  $\delta(D - L) - D$  is irreducible and weakly diagonal dominant. This implies that the matrix  $\delta(D - L) - D$  and  $(\delta(D - L) - D)^{-1}$  are positive definite. Moreover, since the off-diagonal entries of  $(\delta(D - L) - D)$  are all negative, the matrix  $\delta(D - L) - D$  is an M-matrix for  $\delta \geq (1 + \frac{2\epsilon}{h})\frac{h}{\epsilon}$ .  $\square$

LEMMA 3.3. For  $h \gg \epsilon$ , the error reduction matrix  $E_k^s$  of the HLGS iterative method for the matrix obtained from SDFEM discretization of Problem 1 satisfies the following inequality:

$$(3.4) \quad \|E_k^s\| \leq 3 \frac{\epsilon}{h^2} \min \left\{ \frac{3h}{\sqrt{\epsilon}}, 1 \right\}.$$

*Proof.* With  $A_k$  from (3.1),  $E_k^s$  is the product of the inverse of the block lower triangular part of  $A_k$  and the strict block upper triangular part of  $A_k$ . By direct computation, one has  $E_k^s = G_1 G_2$ , where

$$G_1 = \begin{bmatrix} 0 & I & & & \\ 0 & D^{-1}L & & \ddots & \\ \vdots & \vdots & & \ddots & I \\ 0 & (D^{-1}L)^{n-2} & \dots & D^{-1}L & I \\ 0 & (D^{-1}L)^{n-1} & \dots & (D^{-1}L)^2 & D^{-1}L \end{bmatrix}, G_2 = \begin{bmatrix} 0 & & & & \\ & D^{-1}U & & & \\ & & \ddots & & \\ & & & D^{-1}U & \end{bmatrix}.$$

From the Gerschgorin circle theorem, the following inequalities hold for  $h \gg \epsilon$ :

$$\begin{aligned} \|U\| &= \rho(U) \leq \epsilon, \\ \|D^{-1}\| &= \rho(D^{-1}) = \frac{1}{\lambda_{\min}(D)} < \frac{1}{\frac{h}{3} + \frac{10\epsilon}{3}} \leq \frac{3}{h}. \end{aligned}$$

Therefore, we have

$$(3.5) \quad \|G_2\| \leq 3 \frac{\epsilon}{h}.$$

Next, we estimate  $\|G_1\|$ . First, let us estimate  $\|D^{-1}L\|$ . Considering  $D^{-1}L = I - D^{-1}(D - L)$ , we have

$$(3.6) \quad \alpha I - D^{-1}L = D^{-1}(D - L) - (1 - \alpha)I = (1 - \alpha) \left\{ D^{-1} \left[ \frac{1}{1 - \alpha} (D - L) - D \right] \right\}.$$

Let us choose  $\alpha$  satisfying  $\frac{1}{1-\alpha} = \delta = (1 + \frac{2\epsilon}{h})\frac{h}{\epsilon}$ . Lemma 3.2 implies that the matrix  $\frac{1}{1-\alpha}(D - L) - D$  is an M-matrix. Consequently,  $(\frac{1}{1-\alpha}(D - L) - D)^{-1} \geq 0$ . Then using (3.6) and the fact that  $D \geq 0$ , it follows that the matrix  $\alpha I - D^{-1}L > 0$ . Since  $D$  is also positive definite, by Lemma 3.1 we can conclude that

$$(3.7) \quad \rho(D^{-1}L) < \alpha = 1 - \frac{1}{\delta} = 1 - \frac{\epsilon}{h} \left( \frac{1}{1 + 2\frac{\epsilon}{h}} \right) < 1 - \frac{\epsilon}{3h}.$$

Let  $x = (x_1, x_2, \dots, x_N) \in V_h$ , where  $x_i \in R^N$  and  $\sum_{i=1}^N \|x_i\|^2 = 1$ . We have that  $\|G_1x\| < \|G_1y\|$  for  $y = (\|x_1\| x^+, \|x_2\| x^+, \dots, \|x_N\| x^+)$ , where  $x^+$  is the Perron-Frobenius eigenvector of  $D^{-1}L$ . Therefore, the eigenvector corresponding to the maximum eigenvalue has the following form:

$$y = (0, \beta_1x^+, \beta_2x^+, \dots, \beta_{N-1}x^+), \text{ where } \sum_{i=1}^{N-1} \beta_i^2 = 1.$$

By direct computation,

$$\begin{aligned} \|G_1y\| &\leq \left\{ \sum_{i=1}^N \left\| \sum_{k=1}^i \beta_k (D^{-1}L)^{i-k} x^+ \right\|^2 \right\}^{1/2} \\ &= \left\{ \sum_{i=1}^N \left( \sum_{k=1}^i \beta_k l^{i-k} \right)^2 \right\}^{1/2}, \end{aligned}$$

where  $l = \rho(D^{-1}L)$ . Since  $l < 1$  and  $\sum_{i=1}^{N-1} \beta_i^2 = 1$ , the inequalities

$$(3.8) \quad \|G_1\| \leq \left\{ \sum_{i=1}^{N-1} \left( \sum_{k=1}^i \beta_k \right)^2 \right\}^{1/2} \leq \left( \sum_{i=1}^{N-1} i \right)^{1/2} \leq N$$

and

$$(3.9) \quad \begin{aligned} \|G_1\| &\leq \left\{ \sum_{i=1}^{N-1} \left[ \sum_{k=1}^i \beta_k^2 \right] \left[ \sum_{k=1}^i (l^{i-k})^2 \right] \right\}^{1/2} \leq \left\{ \sum_{i=1}^{N-1} \sum_{k=1}^i (l^{i-k})^2 \right\}^{1/2} \\ &\leq \left\{ \frac{1}{1-l^2} \left( \sum_{i=1}^{N-1} 1 - l^{2i} \right) \right\}^{1/2} \leq \left( \frac{N}{1-l} \right)^{1/2} \end{aligned}$$

hold. Recall that  $l < 1 - \frac{\epsilon}{3h}$  from (3.7). By combining (3.8) and (3.9), we have

$$(3.10) \quad \|G_1\| \leq \frac{1}{h} \min \left\{ \frac{3h}{\sqrt{\epsilon}}, 1 \right\}.$$

Therefore, from (3.5) and (3.10), inequality (3.4) holds.  $\square$

LEMMA 3.4 (smoothing property). *Let  $E_k^s$  be the error reduction operator of the HLGs iterative method on  $V_k$ . The following inequality holds:*

$$(3.11) \quad \|A_k(E_k^s)\| \leq \epsilon \left( 1 + 3 \frac{\epsilon}{h_k^2} \min \left\{ \frac{3h_k}{\sqrt{\epsilon}}, 1 \right\} \right).$$

*Proof.* By directly multiplying  $A_k$  and  $E_k^s$ , we have

$$A_k E_k^s = \begin{bmatrix} 0 & U - U(D^{-1}L)D^{-1}U & -UD^{-1}U & & & \\ 0 & -U(D^{-1}L)^2D^{-1}U & \ddots & & \ddots & \\ \vdots & \vdots & \ddots & -U(D^{-1}L)D^{-1}U & -UD^{-1}U & \\ 0 & -U(D^{-1}L)^{n-1}D^{-1}U & \dots & -U(D^{-1}L)^2D^{-1}U & -U(D^{-1}L)D^{-1}U & \\ 0 & 0 & \dots & 0 & 0 & \end{bmatrix}$$

$$= \text{diag}[U](T_1 - T_2G_2),$$

where

$$T_1 = \begin{bmatrix} 0 & I & & & \\ 0 & 0 & \ddots & & \\ \vdots & \vdots & \ddots & I & \\ 0 & 0 & \dots & 0 & I \\ 0 & 0 & \dots & 0 & 0 \end{bmatrix} \quad \text{and} \quad T_2 = \begin{bmatrix} 0 & D^{-1}L & I & & \\ 0 & (D^{-1}L)^2 & \ddots & \ddots & \\ \vdots & \vdots & \ddots & D^{-1}L & I \\ 0 & (D^{-1}L)^{n-1} & \dots & (D^{-1}L)^2 & D^{-1}L \\ 0 & 0 & \dots & 0 & 0 \end{bmatrix}.$$

Using the same argument as in the proof of Lemma 3.3, one can show

$$\|T_2G_2\| \leq 3\frac{\epsilon}{h_k^2} \min\left\{\frac{3h_k}{\sqrt{\epsilon}}, 1\right\}.$$

Therefore,

$$\begin{aligned} \|A_k E_k^s\| &\leq \epsilon \left(1 + 3\frac{\epsilon}{h_k^2} \min\left\{\frac{3h_k}{\sqrt{\epsilon}}, 1\right\}\right) \\ &= \epsilon \left(1 + 3\frac{\epsilon}{h_k^2} \min\left\{\frac{3h_k}{\sqrt{\epsilon}}, 1\right\}\right). \quad \square \end{aligned}$$

Using the inequalities in Lemmas 3.3 and 3.4, our multigrid convergence results can be shown. In the following, estimates of the error reduction operators of GMG and AMG are given in Theorems 3.5 and 3.7, respectively.

**THEOREM 3.5.** *Let  $E_k^{mg}$  be the error reduction operator of the 2-level V-cycle geometric multigrid in Algorithm 1 with the HLGS smoother and bilinear interpolation prolongation on a uniform rectangular mesh  $\mathfrak{S}_k$  for Problem 1. The following estimate for  $E_k^{mg}$  holds:*

$$(3.12) \quad \|E_k^{mg}\| \leq c\frac{\epsilon}{h^{3/2}} \text{ for some constant } c.$$

*Proof.* First, in general, the error reduction operator of GMG can be written as

$$(3.13) \quad E_k^{mg} = E_k^s E^c E_k^s = E_k^s (I - I_{k-1}^k (I - E_{k-1}^{mg}) A_{k-1}^{-1} I_k^{k-1} A_k) E_k^s.$$

For 2-level GMG, since  $E_{k-1}^{mg} = 0$ , we have

$$(3.14) \quad \begin{aligned} \|E_k^{mg}\| &= \|E_k^s (I - I_{k-1}^k A_{k-1}^{-1} I_k^{k-1} A_k) E_k^s\| \\ &\leq \|E_k^s\| \|(I - I_{k-1}^k A_{k-1}^{-1} I_k^{k-1} A_k) E_k^s\|. \end{aligned}$$

Let  $P_k : H^1 \mapsto V_k$  be the projection operator defined by  $B_{sd}(P_k u, v) = B_{sd}(u, v)$  for all  $v \in V_k$  for all  $k$ . We have  $I_k^{k-1} A_k = A_{k-1} P_{k-1}$  on  $V_k$ . Therefore, let  $e^s = E_k^s(e)$  for any  $e \in V_k \subset H^1$ ,

$$\begin{aligned} \|(I - I_{k-1}^k A_{k-1}^{-1} I_k^{k-1} A_k) E_k^s(e)\| &= \|(I - I_{k-1}^k P_{k-1}) e^s\| \\ &\leq ch_k^{-2} \|(I - P_{k-1}) e^s\|_0 \leq ch_k^{-2} \frac{1}{\sqrt{\epsilon}} \|(I - P_{k-1}) e^s\|_1 \\ &\leq ch_k^{-2} \sqrt{\frac{h_k}{\epsilon}} |e^s|_1 \text{ by the a priori error estimation,} \\ &\leq ch_k^{-2} \sqrt{\frac{h_k}{\epsilon^2}} \|A_k E_k^s(e)\|_0 \text{ by the regularity estimate,} \\ &\leq c\sqrt{h_k} \left(1 + 3\frac{\epsilon}{h_k^2} \min\left\{\frac{3h_k}{\sqrt{\epsilon}}, 1\right\}\right) \|e\| \text{ by (3.11),} \end{aligned}$$

where  $c$  is a constant independent of  $h$  and  $\epsilon$ . By combining the above estimate and (3.4), inequality (3.14) implies

$$\|E_k^{mg}\| \leq c\frac{\epsilon}{h^{3/2}} \min\left\{\frac{h}{\sqrt{\epsilon}}, 1\right\} \left(1 + \frac{\epsilon}{h^2} \min\left\{\frac{h}{\sqrt{\epsilon}}, 1\right\}\right) \leq c\frac{\epsilon}{h^{3/2}}. \quad \square$$

*Remark 3.6.* From inequality (3.12), it is clear that the geometric multigrid converges when  $h_k \gg \epsilon^{2/3}$ . The above 2-level analysis can be generalized by mathematical induction and estimation of (3.13). In [23], we have found that the multigrid V-cycle converges when  $h \gg \sqrt{\epsilon}$ .

Next, we analyze AMG convergence. The framework to prove the convergence of algebraic multigrid is based on the *smoothing assumption*,

(3.15)  $\exists \alpha > 0$  such that  $\|E_k^s e_h\|_1^2 \leq \|e_h\|_1^2 - \alpha \|e_h\|_2^2$  for any  $e_h \in V_k$ ,

and the *approximation assumption*,

(3.16)  $\min_{e_H \in V_{k-1}} \|e_h - I_H^h e_H\|_0^2 \leq \beta \|e_h\|_1^2$  with  $\beta > 0$  independent with  $e_h \in V_k$ ,

where

(3.17)  $\|v\|_0^2 = \langle Dv, v \rangle, \|v\|_1^2 = \langle A_k v, v \rangle, \text{ and } \|v\|_2^2 = \langle D^{-1} A_k v, A_k v \rangle$

for any  $v \in V_k$ . For problems where the coefficient matrix is an M-matrix, Ruge and Stüben have proved that 2-level AMG converges when both the smoothing assumption and the approximation assumption hold for M-matrices. It has also been shown that the smoothing assumption holds for the usual point Gauss-Seidel relaxation, and the approximation assumption holds for some simple coarse grid selection algorithms

with algebraic interpolation formula [17]. The classical algebraic multigrid coarsening algorithm is designed to satisfy the approximation assumption based on the property of algebraically smooth error, whereby the smooth errors vary slowly along the strong connection directions.

For the discrete convection-diffusion Problem 1, the matrix obtained from SDFEM discretization is nonsymmetric and not an M-matrix. As a result, the usual framework of AMG analysis cannot be applied directly. In Lemma 3.2 of [14], Reusken has shown that there is a constant  $c$  independent of  $\epsilon$  and  $h$  such that  $\|E^c\| \leq c$  for a two-grid method with a coarse grid from semicoarsening and a matrix-dependent prolongation. In the following, we take this as an assumption and estimate the error reduction rate of AMG in terms of the parameter  $\beta$  in the approximation assumption (3.16) by using (3.4) and (3.11).

**THEOREM 3.7.** *Assume the approximation assumption (3.16) holds. Then the following inequality holds:*

$$(3.18) \quad \|E_k^{amg}\| \leq \sqrt{3}c \frac{\epsilon^2}{h^3} \min \left\{ \frac{h}{\sqrt{\epsilon}}, 1 \right\} \frac{\beta}{\sqrt{h}},$$

where  $c$  is some constant independent of  $\epsilon$  and  $h$ .

*Proof.* For any  $e \in V_k$ , we have

$$(3.19) \quad \|E_k^{amg}(e)\| = \|E_k^s E^c E_k^s(e)\| \leq \frac{\epsilon}{h^2} \min \left\{ \frac{h}{\sqrt{\epsilon}}, 1 \right\} \|E^c E_k^s(e)\|.$$

Let  $\bar{e} = E_k^s(e)$ . Since  $\text{Ker}(E^c) = \text{Rang}(I_{k-1}^k)$ , we have

$$(3.20) \quad \begin{aligned} \|E^c(\bar{e})\| &\leq c \min_{e_H \in V_{k-1}} \|\bar{e} - I_{k-1}^k e_H\| \leq c \sqrt{\frac{3}{h}} \min_{e_H \in V_{k-1}} \|\bar{e} - I_{k-1}^k e_H\|_0 \\ &\leq c \sqrt{\frac{3}{h}} \beta \|\bar{e}\|_1 \text{ by approximation assumption (3.16),} \\ &\leq \sqrt{\frac{3}{h}} \beta (\|A_k E_k^s(e)\| \|E_k^s(e)\|)^{1/2} \\ &\leq c \sqrt{\frac{3}{h}} \beta \left( \epsilon \left[ 1 + \frac{\epsilon}{h^2} \min \left\{ \frac{h}{\sqrt{\epsilon}}, 1 \right\} \right] \left[ \frac{\epsilon}{h^2} \min \left\{ \frac{h}{\sqrt{\epsilon}}, 1 \right\} \right] \right)^{1/2} \|e\| \\ &\quad \text{by (3.4) and (3.11)} \\ &\leq c \sqrt{\frac{3}{h}} \frac{\epsilon}{h} \beta \|e\|. \end{aligned}$$

By combining (3.19) and (3.20), inequality (3.18) holds.  $\square$

*Remark 3.8.* Inequality (3.18) can be rewritten as

$$\|E_k^{amg}\| \leq c \frac{\epsilon}{h^{3/2}} \min \left\{ \frac{h}{\sqrt{\epsilon}}, 1 \right\} \frac{\beta \epsilon}{h^2} = c \frac{\epsilon}{h^{3/2}} \left( \frac{\sqrt{\epsilon}}{h} \right)^\alpha \beta,$$

where  $\alpha = \begin{cases} 1, & h < \sqrt{\epsilon}, \\ 2, & h \geq \sqrt{\epsilon}. \end{cases}$  Comparison with the estimate of  $E_k^{mg}$  in (3.12) suggests that algebraic multigrid will converge more rapidly than geometric multigrid if the parameter  $\beta$  in the approximation assumption is less than  $(\frac{h}{\sqrt{\epsilon}})^\alpha$ . Although we have not

TABLE 3.1  
Comparison of GMG and AMG for Problem 1 on uniform rectangular meshes.

$\epsilon$	$10^{-1}$	$10^{-2}$	$10^{-3}$	$10^{-4}$	$\epsilon$	$10^{-1}$	$10^{-2}$	$10^{-3}$	$10^{-4}$
$h = \frac{1}{8}$	5	4	2	2	$h = \frac{1}{8}$	4	3	2	2
$h = \frac{1}{16}$	5	4	3	2	$h = \frac{1}{16}$	4	3	3	2
$h = \frac{1}{32}$	5	5	4	2	$h = \frac{1}{32}$	5	4	3	2

(a) GMG iterative steps.

(b) AMG iterative steps.

yet estimated  $\beta$  for the interpolation operator from AMG coarsening, AMG converges more rapidly than GMG, as shown in Table 3.1. Moreover, in Table 3.1, the facts that both GMG and AMG converge more rapidly as  $\epsilon$  becomes smaller and converge more slowly as the mesh size  $h$  decreases are also consistent with our theoretical analysis. In these numerical studies, the stopping tolerance is set to be

$$(3.21) \quad \|r_m\| \leq 10^{-6} \|r_0\|,$$

where  $r_0$  is the initial residual and  $r_m$  is the residual at the  $m$ th iteration.

*Remark 3.9.* Since the row-sum  $\sum_j(a_{i,j})$  and column-sum  $\sum_i(a_{i,j})$  are equal for the matrix  $A_k$  arising from Problem 1, the following equality holds:

$$(3.22) \quad \|e\|_1^2 = (A_k e, e) = \sum_{i,j} a_{i,j} e_i e_j = \frac{1}{2} \left( \sum_{i,j} (-a_{i,j})(e_i - e_j)^2 \right) + \sum_i s_i e_i^2,$$

where  $s_i = \sum_j(a_{i,j})$ . Recall that the algebraically smooth error  $e_s$  is characterized by  $E_k^s(e_s) \approx e_s$ . Clearly, inequalities (3.4) and (3.11) imply

$$\begin{aligned} \langle A_k e_s, e_s \rangle &\approx \langle A_k E_k^s e_s, E_k^s e_s \rangle \leq \frac{c}{h} \|A_k E_k^s\| \|E_k^s\| \|e_s\|_0^2 \\ &\leq c \frac{\epsilon^2}{h^3} \min \left\{ \frac{3h}{\sqrt{\epsilon}}, 1 \right\} \left( 1 + \frac{3\epsilon}{h^2} \min \left\{ \frac{3h}{\sqrt{\epsilon}}, 1 \right\} \right) \|e_s\|_0^2 \\ &\leq c \frac{\epsilon^2}{h^3} \|e_s\|_0^2. \end{aligned}$$

For  $h \gg \epsilon^{2/3}$ ,  $\|e_s\|_1^2 \ll \|e_s\|_0^2$  follows from the above estimate. Moreover, by (3.22), this is equivalent to

$$\begin{aligned} &\frac{1}{2} \sum_i \sum_{j \neq i-1, i+1} (-a_{i,j})(e_i - e_j)^2 + \sum_i s_i e_i^2 \\ &\ll \sum_i a_{i,i} e_i^2 + a_{i,i-1}(e_i - e_{i-1})^2 + a_{i,i+1}(e_i - e_{i+1})^2 \\ &\Rightarrow \frac{1}{2} \sum_i \sum_{j \neq i-1, i+1} \left( \frac{-a_{i,j}}{a_{i,i}} \right) (e_i - e_j)^2 \ll \sum_i e_i^2 + \frac{1}{4}(e_i - e_{i-1})^2 + \frac{1}{4}(e_i - e_{i+1})^2. \end{aligned}$$

It can be seen from (3.1) that the weights  $\frac{a_{i,j}}{a_{i,i}} = O(1)$  for downstream indices  $j$  and  $O(\frac{\epsilon}{h})$  for upstream. Therefore, we can conclude that the smooth error tends to vary more slowly in downstream directions than in upstream directions. Since the strong

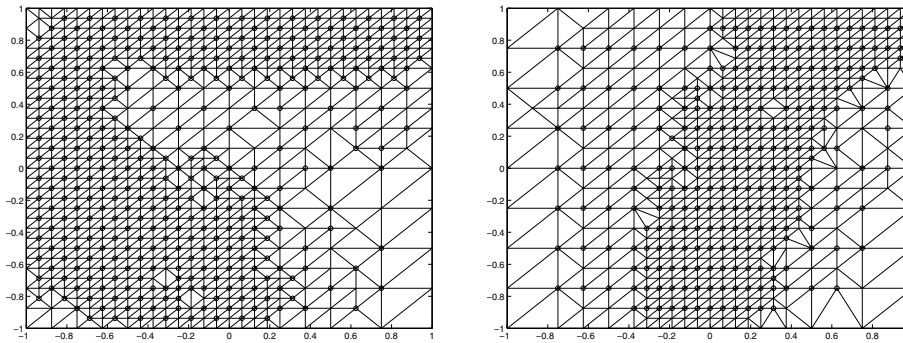
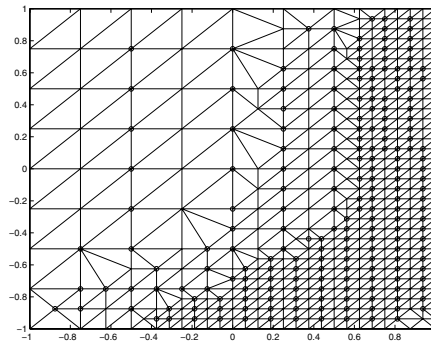
(a)  $\phi = -45^\circ$  on an adaptive mesh.(b)  $\phi = 15^\circ$  on an adaptive mesh.(c)  $\phi = 75^\circ$  on an adaptive mesh.

FIG. 3.1. Coarse grid points selected by AMG coarsening.

connectivity of the matrix  $A_k$  is also along the downstream direction, we expect the AMG coarsening to select a large number of downstream nodes, and the resulting strategy tends to resemble one of semicoarsening.

In the following, we show the AMG coarse grids for the Problem 1 with  $\epsilon = 10^{-3}$  and  $\phi = -45^\circ$ ,  $15^\circ$ , and  $75^\circ$ , where the connection parameter is  $\mu = 0.25$ . Each fine grid is generated by two adaptive mesh refinement steps from an initial  $8 \times 8$  mesh and the threshold value  $\theta$  in (2.1) is set to 0.01. In Figure 3.1, the fine grid is plotted and the coarse grid points, selected by AMG coarsening, are marked by  $\circ$ . Evidently, the AMG coarsening process is sensitive to the flow direction mentioned in Remark 3.9 and generates meshes by means of semicoarsening. The numerical results in Table 3.1, and Table 4.3 and 4.4 of the next section, shows that AMG converges more quickly than GMG for Problem 1. This leads us to conjecture that the algebraically smooth errors are well in the range of the AMG interpolation operator, i.e., the parameter  $\beta$  in the approximation assumption is small. Analysis of the AMG interpolation and approximation assumption will be considered in a future study.

For Problem 2, multigrid convergence is much harder to prove. In fact, the presence of a stagnation point leads to poor performance of multigrid methods when the equation is discretized by a first order upwind scheme [10], [4]. However, many

numerical studies have shown that GMG- and AMG-accelerated Krylov space methods, such as GMRES and BiCGSTAB, achieve nearly mesh-independent convergence [10], [19]. Recently, Ramage has demonstrated that GMRES with GMG preconditioning achieves mesh-independent convergence when SDFEM with an optimal stabilization parameter is employed for discretization on uniform rectangular mesh [13]. In the next section, GMG and AMG are used as a preconditioner for GMRES in some numerical tests.

**4. Numerical results.** In this section, the performances of different linear solvers, including GMG, AMG, and preconditioned GMRES, for the discrete convection-diffusion equation are compared on both adaptively refined meshes and uniform meshes. We use both Problem 1 with  $\phi = 0$  and Problem 2. In both problems, (1.1) is discretized by SDFEM with the stabilization parameter  $\delta_T$  defined in (1.3) on both a uniform  $32 \times 32$  mesh and an adaptively refined mesh for  $\epsilon = 10^{-2}, 10^{-3}$ , and  $10^{-4}$ . Here, a  $32 \times 32$  mesh is generated from 3 uniform refinements starting with a  $4 \times 4$  initial mesh, and the adaptively refined mesh is generated by refining an initial  $8 \times 8$  mesh 4 times, where full multigrid is used with those coarse meshes. The threshold value  $\theta$  in the maximum marking strategy is chosen such that elements in the regions containing large errors can be refined for both problems.

For Problem 1,  $\theta = 0.1, 0.01$ , and  $0.001$  for  $\epsilon = 10^{-2}, 10^{-3}$ , and  $10^{-4}$ , respectively. The adaptive meshes are shown in Figure 4.1. For Problem 2,  $\theta = 0.1$  for all  $\epsilon$  and the adaptive meshes are shown in Figure 4.2.

Since there are no natural “lines” in unstructured meshes or in the coarse meshes of AMG, only point-versions of the Gauss–Seidel method are used for smoothing. Grid points are numbered in lexicographical order where the  $y$ -coordinate is the primary

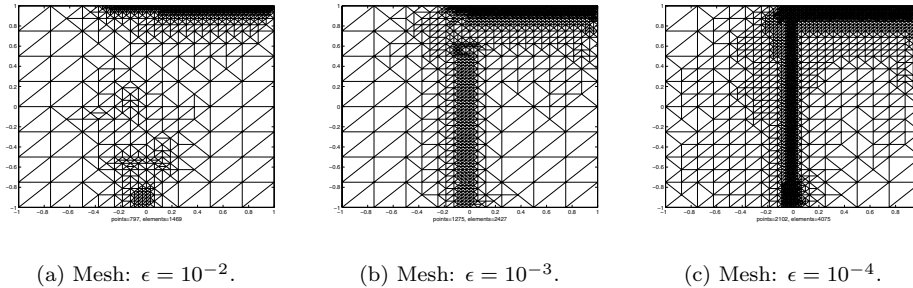


FIG. 4.1. Problem 1: Adaptively refined meshes after 4 refinements starting from an  $8 \times 8$  grid for various  $\epsilon$ .

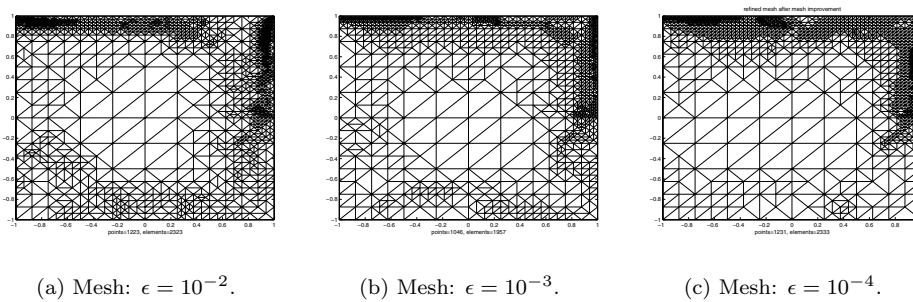


FIG. 4.2. Problem 2: Adaptively refined meshes for various  $\epsilon$ .



key and the  $x$ -coordinate is the secondary key. The point Gauss–Seidel method associated with this node ordering is called the forward horizontal Gauss–Seidel (forward-HGS) method. Naturally, one can obtain another lexicographical order where the  $x$ -coordinate is the primary key and the  $y$ -coordinate is the secondary key. The point Gauss–Seidel method associated with this node ordering is then called the forward vertical Gauss–Seidel (forward-VGS) method. The backward-HGS and backward-VGS methods are obtained from the forward-HGS and forward-VGS methods, respectively, by simply reversing the node ordering. In Problem 1, one step of forward-HGS is used in both presmoothing and postsmoothing. In Problem 2, the smoother consists of four point Gauss–Seidel sweeps (forward-HGS, forward-VGS, backward-HGS, and backward-VGS). Hereafter, this smoother is abbreviated as ADGS (namely, alternating direction Gauss–Seidel).

Both problems are solved by GMG, AMG, GMRES, and right-preconditioned GMRES with zero initial guess, and the stopping tolerance (3.21) is used here for these iterative solvers. The preconditioned GMRES methods with HGS, ADGS, GMG, and AMG preconditioners are denoted as GMRES-HGS, GMRES-ADGS, GMRES-GMG, and GMRES-AMG, respectively. In the following tables, increases in level numbers correspond to finer meshes, where level 1 represents the coarsest mesh in multigrid solver. The notation “-” represents that the number of iterations is greater than 400.

First, Tables 4.1 and 4.2 show some details about the meshes used by GMG and AMG. Table 4.1 shows the number of coarse grid points obtained starting from a uniform  $32 \times 32$  fine grid of elements. Listed under “GMG” is simply the number of points in the hierarchical set of uniform coarse grids. Listed under “AMG” is the number of grid points derived using the AMG coarsening strategy. Table 4.2 shows what happens when adaptive grid refinement is used. That is, GMG uses the hierarchical set of grids obtained from adaptive refinement. In contrast, AMG starts with the fine grid produced by grid refinement, but then it generates *its own* set of coarse grids using its coarsening strategy. Table 4.1 shows that when AMG starts with a uniform fine mesh, it generates more coarse grid points than are found in the hierarchical coarse grids. On the other hand, Table 4.2 shows that if AMG starts with a fine mesh obtained from adaptive refinement, the coarse grids it generates often contain fewer points than are found in the adaptively refined coarse grids.

Tables 4.3–4.6 examine the iteration counts for both multigrid strategies as a function of the levels at which the solver is used. Tables 4.3 and 4.4 show that both solvers display “textbook” performance on Problem 1 for  $\epsilon = 10^{-2}$ . As  $\epsilon$  is reduced, however, performance of GMG decreases and becomes mesh-dependent, especially on uniform grids. In contrast, the performance of AMG is robust and shows weaker

TABLE 4.1  
Number of coarse grid points from uniform refinement and AMG coarsening.

	GMG	AMG				GMG	AMG		
$\log_{10} \frac{1}{\epsilon}$	2, 3, 4	2	3	4	$\log_{10} \frac{1}{\epsilon}$	2, 3, 4	2	3	4
Level = 4	1089	1089	1089	1089	Level = 4	1089	1089	1089	1089
Level = 3	289	480	479	479	Level = 3	289	499	507	513
Level = 2	81	307	331	231	Level = 2	81	294	283	289
Level = 1	25	157	108	108	Level = 1	25	174	164	169

(a) Problem 1.

(b) Problem 2.

TABLE 4.2

Number of coarse grid points from adaptive refinement and AMG coarsening.

$\log_{10} \frac{1}{\epsilon}$	GMG			AMG		
	2	3	4	2	3	4
Level = 5	797	1275	2102	797	1275	2102
Level = 4	410	649	1047	359	586	997
Level = 3	215	320	528	193	328	547
Level = 2	122	176	239	112	179	308
Level = 1	81	81	81	55	101	153

(a) Problem 1.

$\log_{10} \frac{1}{\epsilon}$	GMG			AMG		
	2	3	4	2	3	4
Level = 5	1233	946	922	1233	946	922
Level = 4	625	632	823	597	496	445
Level = 3	305	373	387	338	291	258
Level = 2	150	196	199	205	168	165
Level = 1	81	81	81	121	95	97

(b) Problem 2.

TABLE 4.3

Problem 1: Iteration counts for GMG and AMG on uniform meshes.

Level	GMG	AMG
3	13	7
2	13	6
1	12	6

(a)  $\epsilon = 10^{-2}$ .

Level	GMG	AMG
3	27	8
2	26	7
1	16	6

(b)  $\epsilon = 10^{-3}$ .

Level	GMG	AMG
3	51	11
2	35	8
1	17	6

(c)  $\epsilon = 10^{-4}$ .

TABLE 4.4

Problem 1: Iteration counts for GMG and AMG on adaptive meshes.

Level	GMG	AMG
4	9	6
3	8	6
2	7	5
1	7	5

(a)  $\epsilon = 10^{-2}$ .

Level	GMG	AMG
4	22	6
3	24	6
2	18	6
1	17	6

(b)  $\epsilon = 10^{-3}$ .

Level	GMG	AMG
4	59	11
3	57	9
2	47	7
1	34	6

(c)  $\epsilon = 10^{-4}$ .

dependence on  $\epsilon$  and mesh size. Problem 2, with a recirculating flow, is more difficult for both solvers. On a uniform grid, GMG diverges for  $\epsilon = 10^{-4}$ , but AMG manages to converge, as shown in Table 4.5. On adaptive meshes, as shown in Table 4.6, textbook-like performance is achieved for both methods. However, AMG is considerably slower than GMG, especially for small  $\epsilon = 10^{-4}$  on the finest mesh. Our speculation

TABLE 4.5  
 Problem 2: Iteration counts for GMG and AMG on uniform meshes.

Level	GMG	AMG	Level	GMG	AMG	Level	GMG	AMG
3	25	9	3	179	11	3	-	52
2	25	8	2	135	21	2	-	43
1	20	7	1	43	11	1	57	13

(a)  $\epsilon = 10^{-2}$ .(b)  $\epsilon = 10^{-3}$ .(c)  $\epsilon = 10^{-4}$ .

TABLE 4.6  
 Problem 2: Iteration counts for GMG and AMG on adaptive meshes.

Level	GMG	AMG	Level	GMG	AMG	Level	GMG	AMG
4	9	12	4	27	26	4	30	39
3	9	11	3	21	26	3	32	38
2	7	10	2	24	21	2	40	44
1	6	7	1	20	17	1	30	36

(a)  $\epsilon = 10^{-2}$ .(b)  $\epsilon = 10^{-3}$ .(c)  $\epsilon = 10^{-4}$ .

TABLE 4.7  
 Problem 1: Iteration counts for various GMRES methods on finest grids.

$\epsilon$	$10^{-2}$	$10^{-3}$	$10^{-4}$	$\epsilon$	$10^{-2}$	$10^{-3}$	$10^{-4}$
GMG	13	27	51	GMG	9	22	59
AMG	7	8	11	AMG	6	6	11
GMRES	56	76	95	GMRES	83	146	-
GMRES-HGS	24	26	36	GMRES-HGS	19	25	48
GMRES-GMG	11	16	22	GMRES-GMG	8	13	27
GMRES-AMG	6	7	9	GMRES-AMG	5	5	9

(a) Iterative steps on uniform mesh.

(b) Iterative steps on adaptive mesh.

is that this is caused by inadequacy of the grid transfer operator used in AMG and the fact that the number of coarse grid points generated by AMG coarsening is only about half the number of coarse grid points used in GMG, as shown in Table 4.2(b). AMG interpolation can be improved by the recently developed AMGe methods [3]. As we will show below, performance can also be remedied using a Krylov space acceleration.

Next, we compare the performance of preconditioned GMRES methods on both the finest uniform mesh and the finest adaptive mesh. Tables 4.7 and 4.8 examine the iteration counts for various versions of preconditioned GMRES. The Krylov space acceleration of the convergence of GMG and AMG is observed in most cases. The numerical results indicate that the AMG-preconditioned GMRES requires the fewest iterations to converge, compared to GMG, AMG, and other versions of GMRES. Particularly in Problem 2, despite the fact that AMG converges more slowly than GMG, GMRES-AMG outperforms GMRES-GMG. This observation suggests that AMG can be a good preconditioner for Krylov space solvers in solving complex flow problems. Moreover, in Table 4.9, iterations of GMRES-AMG on each level of the adaptive meshes are shown for  $\epsilon = 10^{-2} \dots 10^{-4}$ . It is clear that the convergence of GMRES-AMG depends only weakly on the mesh size and  $\epsilon$  on adaptive meshes.

*Remark 4.1.* Notice a difference between performance of GMG for Problem 1

TABLE 4.8  
 Problem 2: Iteration counts for various GMRES methods on finest grids.

$\epsilon$	$10^{-2}$	$10^{-3}$	$10^{-4}$	$\epsilon$	$10^{-2}$	$10^{-3}$	$10^{-4}$
GMG	25	179	-	GMG	9	27	30
AMG	9	11	52	AMG	11	26	39
GMRES	392	-	-	GMRES	199	-	-
GMRES-ADGS	33	54	73	GMRES-ADGS	30	37	35
GMRES-GMG	11	29	42	GMRES-GMG	7	12	12
GMRES-AMG	6	8	15	GMRES-AMG	6	8	9

(a) Iterative steps on uniform mesh.

(b) Iterative steps on adaptive mesh.

TABLE 4.9  
 Iteration counts of GMRES-AMG for various  $\epsilon$  on adaptive meshes.

Level	$\epsilon = 10^{-2}$	$10^{-3}$	$10^{-4}$	Level	$\epsilon = 10^{-2}$	$10^{-3}$	$10^{-4}$
4	5	5	9	4	6	8	9
3	5	5	8	3	6	7	9
2	5	5	6	2	5	7	9
1	4	5	6	1	4	6	7

(a) Iterative steps in Problem 1.

(b) Iterative steps in Problem 2.

here and the bounds shown in section 3 which suggest convergence is essentially independent of  $\epsilon$ . The latter result holds only in special settings and depends on having a very accurate smoother. Similar  $\epsilon$ -independent results for this problem can be found in [9], [12], [14]. In contrast, the better performance of AMG is due to matrix-dependent grid transfer operators and the superior choice of coarse mesh identified by AMG coarsening, for which the smoother is better able to mimic the flow characteristics.

**5. Conclusion.** In this work, we have explored solution strategies for the convection-diffusion equation using adaptive gridding techniques and a multigrid algorithm. In section 3, the convergence analysis for both GMG with bilinear interpolation and line Gauss-Seidel smoother and AMG with the same smoother not only shows that GMG converges for Problem 1 when  $h \gg \epsilon^{2/3}$ , but also suggests that AMG converges faster than GMG on uniform rectangular meshes. Although the AMG interpolation and approximation assumption remain to be analyzed, our numerical results support this observation on both triangular and rectangular uniform meshes. In addition, numerical studies in section 4 show that the convergence rates of both GMG and AMG deteriorates as the diffusion parameter  $\epsilon$  decreases. To overcome this drawback, we have found that Krylov space acceleration significantly improves the convergence of both GMG and AMG, especially when  $\epsilon$  is small. Moreover, GMRES-AMG converges more rapidly than GMRES-GMG in all of our test cases on both uniform and adaptive meshes.

With adaptive mesh refinement, GMG seems to be a natural choice because the coarse grids, correction operators, and interpolations are ready to be used in GMG. However, GMG convergence is slower than AMG for some examples of both tested problems even though more coarse grid points are employed in GMG. On the other hand, the most important strength of AMG is that it can be treated as a black-box solver. There is no need to compute coarse grids and interpolations explicitly in AMG. The facts that AMG is applicable to a wider range of applications, AMG

converges for all of our test cases, and GMRES-AMG outperforms all other solvers in our numerical studies make AMG very attractive for solving the SDFEM-discretized convection-diffusion problems.

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