## PRECONDITIONING FOR THE STEADY-STATE NAVIER–STOKES EQUATIONS WITH LOW VISCOSITY\*

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**Abstract.** We introduce a preconditioner for the linearized Navier–Stokes equations that is effective when either the discretization mesh size or the viscosity approaches zero. For constant coefficient problems with periodic boundary conditions, we show that the preconditioning yields a system with a single eigenvalue equal to 1, so that performance is independent of both viscosity and mesh size. For other boundary conditions, we demonstrate empirically that convergence depends only mildly on these parameters and we give a partial analysis of this phenomenon. We also show that some expensive subsidiary computations required by the new method can be replaced by inexpensive approximate versions of these tasks based on iteration, with virtually no degradation of performance.

Key words. Navier-Stokes, Oseen, preconditioning, iterative methods

AMS subject classifications. Primary, 65F10, 65N20; Secondary, 15A06

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**1. Introduction.** This paper introduces a methodology for preconditioning the discrete steady-state incompressible Navier–Stokes equations

(1) 
$$-\nu \Delta \mathbf{u} + (\mathbf{u} \cdot \operatorname{grad}) \mathbf{u} + \operatorname{grad} p = \mathbf{f} \\ -\operatorname{div} \mathbf{u} = 0 \quad \text{in } \Omega,$$

subject to suitable boundary conditions on  $\partial\Omega$ , where  $\Omega$  is an open bounded domain in  $\mathbb{R}^2$  or  $\mathbb{R}^3$ . The vector field **u** represents the velocity in  $\Omega$ , p represents pressure, and the scalar  $\nu$  is the viscosity, which is inversely proportional to the Reynolds number. We will develop the preconditioners for the linearized version of (1) known as the *Oseen equations*, which can be written as

(2) 
$$-\nu\Delta \mathbf{u} + (\mathbf{w} \cdot \operatorname{grad}) \mathbf{u} + \operatorname{grad} p = \mathbf{f}, \\ -\operatorname{div} \mathbf{u} = 0,$$

where **w** is given such that div **w** = 0. These equations arise from a nonlinear iteration essentially of the form  $-\nu \Delta \mathbf{u}^{(m)} + (\mathbf{u}^{(m-1)} \cdot \text{grad}) \mathbf{u}^{(m)} + \text{grad} p^{(m)} = \mathbf{f}, -\text{div} \mathbf{u}^{(m)} = 0;$ see [18].

Common discretizations of (2) yield a linear system of equations

(3) 
$$\begin{pmatrix} F & B^T \\ B & 0 \end{pmatrix} \begin{pmatrix} \mathbf{u} \\ p \end{pmatrix} = \begin{pmatrix} \mathbf{f} \\ 0 \end{pmatrix},$$

where **u** and *p* now represent discrete versions of velocity and pressure, respectively. Here  $F = \nu A + N$ , where A consists of a set of uncoupled discrete Laplace operators, corresponding to diffusion, and N is a discrete convection operator. We are interested in convergence behavior of iterative methods applied to (3) as either  $\nu$  or the

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discretization mesh width h tend to zero. For small  $\nu$  it is necessary for the discretization to be fine enough to resolve features such as boundary layers [13]. Ideally, when  $\nu$  is decreased, h will also be reduced proportionally. We are concerned only with values of  $\nu$  for which stable steady-state solutions exist; for example, it is shown in [12] that values of  $1/\nu$  on the order of 1000 to several thousand are feasible.

Let  $\mathcal{A}$  denote the coefficient matrix of (3). We will consider preconditioners of the form

(4) 
$$\boldsymbol{\mathcal{Q}} = \begin{pmatrix} F & B^T \\ 0 & -X \end{pmatrix}.$$

It is easily confirmed that

(5) 
$$\mathcal{A}\mathcal{Q}^{-1} = \begin{pmatrix} I & 0\\ BF^{-1} & BF^{-1}B^TX^{-1} \end{pmatrix},$$

so that the eigenvalues of  $\mathcal{AQ}^{-1}$  are

$$\{1\} \cup \sigma(BF^{-1}B^TX^{-1}).$$

For finite element discretization, let M denote the diagonal of the pressure mass matrix; for finite differences, a natural analogue is  $M = h^d I$  for a uniform grid of width h in d dimensions. It was shown in [5] that for the choice  $X = \frac{1}{\nu}M$ , the eigenvalues of  $BF^{-1}B^TX^{-1}$  are bounded independently of h, and experimental results suggest that Krylov subspace methods such as GMRES [26] have convergence rates that are independent of the mesh size. See [19] for a rigorous analysis of this method based on the field of values and [2, 24, 27, 31] for analogous results for the Stokes equations. Similar bounds were also obtained in [11] for a different class of preconditioners based on the symmetric part of  $\mathcal{A}$ , where the preconditioning entails solution of the Stokes equations.

The convergence properties of these approaches depend on the viscosity  $\nu$ , and in general, convergence rates deteriorate as  $\nu$  decreases. For example, the results in [5] yield eigenvalues that are contained in a box in the complex plane of the form

$$[c_1\nu^2, c_2] \times i [c_3, c_4],$$

where  $\{c_j\}$  are independent of h and  $\nu$ , and in experimental results iteration counts increase roughly like  $1/\nu$ . (See also section 4.)

Our concern here is to develop alternative choices for X for which the sensitivity to  $\nu$  is less pronounced. Our starting point is an observation derived from [6]. Let G and K be two matrices of dimensions  $n_p \times n_u$  with  $n_u \ge n_p$  and such that both are of full rank  $n_p$ . The matrix  $K^T (GK^T)^{-1}G$  can then be viewed as an operator from range $(K^T)$  to itself, and it is trivial to see that this is in fact the identity operator. To apply this to (3), first assume for simplicity that B is of full rank (this assumption will be eliminated below) and let  $G = BF^{-1}$  and K = B. Our observation is then

$$B^T (BF^{-1}B^T)^{-1}BF^{-1} = I \qquad \text{on range}(B^T)$$

or, equivalently,

(6) 
$$B^T (BF^{-1}B^T)^{-1}B = F$$
 on range $(F^{-1}B^T)$ .

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Suppose for the moment that

(7) 
$$\operatorname{range}(B^T) \subset \operatorname{range}(F^{-1}B^T)$$

Then the equality (6) can be postmultiplied by  $B^T$ , and premultiplying the result by B yields

$$(BB^T)(BF^{-1}B^T)^{-1}(BB^T) = BFB^T.$$

Equivalently,

$$(BF^{-1}B^T)^{-1} = (BB^T)^{-1}(BFB^T)(BB^T)^{-1}.$$

That is, for the choice

(8) 
$$X = (BB^T) (BFB^T)^{-1} (BB^T),$$

the eigenvalues of the preconditioned operator (5) are identically 1. Moreover, it can be shown that in this case,  $\mathcal{AQ}^{-1}$  has Jordan blocks of order at most 2, so that two steps of GMRES will produce the solution.

In the rest of this paper, we examine the use of the preconditioner (4) for solving the discrete two-dimensional Oseen equations, using X defined by (8) as well as some computationally less expensive variants. Because of the presence of the high-order discrete operator  $BFB^T$  in (8), we will refer to the combination of (4) and (8) as the BFBt preconditioner; cf. [22] for other approaches to the problem of approximating the action of the inverse of  $BF^{-1}B^T$ . For our analysis, we restrict our attention to a "marker-and-cell" (MAC) finite-difference operator [15], which we outline in section 2. In section 3, we show that in the case where (2) is given with constant "wind"  $\mathbf{w}$  and periodic boundary conditions, (7) holds. (In fact the two spaces are identical.) Consequently, in this case the discussion of the previous paragraph represents a complete analysis. In section 4, we examine through a series of numerical experiments the extent to which these results reflect the behavior of the preconditioner in more realistic scenarios, that is, for Dirichlet boundary conditions or nonconstant wind. Our observations are that the convergence behavior of Krylov subspace methods is independent of  $\nu$  for the Dirichlet problem with constant wind and mildly dependent on  $\nu$ for variable wind; convergence also depends mildly on the mesh size h, with iteration counts increasing in proportion to  $h^{-1/2}$ . These conclusions also hold for variants of the BFBt preconditioner designed to keep computational costs low. In particular, use of  $\mathcal{Q}$  with (8) in an iteration entails two Poisson solutions on the pressure space and (for any X) the solution of a set of convection-diffusion equations on the velocity space. We show that these computations can be approximated using inner iterations with little degradation of performance of the outer iteration. In section 5 we make some summarizing remarks, and in an appendix we present a partial analysis of the behavior of the BFBt preconditioner for the Stokes problem.

2. Finite-difference discretization. We briefly describe the MAC finitedifference scheme. Assume that  $\Omega$  is the rectangular region  $(0,1) \times (0,1)$ , divided into a uniform  $n \times n$  grid of cells of width h = 1/n. Let  $\mathbf{u} = (u, v)^T$  denote the velocity field and  $\mathbf{w} = (a, b)^T$  the wind. The discrete velocities and pressures are defined on a staggered grid in which the discrete values of u lie in the centers of the cell boundaries orthogonal to the x-axis, the discrete values of v lie in the center of the cell boundaries orthogonal to the y-axis, and the discrete pressures lie in the cell centers. An example is shown in Figure 1.



FIG. 1. Staggered grid for the MAC finite-difference discretization.

The matrix of (3) contains three block rows, the first two of which come from the momentum equations for the individual components of the discrete velocity field and the last from the incompressibility constraint. F and B have the form

$$F = \begin{pmatrix} F_1 & 0\\ 0 & F_2 \end{pmatrix}, \qquad B = (B_1 \ B_2),$$

where  $F_i = \nu A_i + N_i$ . The submatrices are defined as follows. Let  $\phi_{jk}$  denote the value of a mesh function  $\phi$  at the point  $(jh, kh) \in \overline{\Omega}$ . The form of the indices (j, k) depends on the mesh function to which they correspond. In particular, they need not be integers. The first block row of (3) is defined by

$$\begin{array}{ll} & [-\Delta u]_{jk} \approx [A_1 u]_{jk} \equiv \frac{1}{h^2} \left( 4u_{jk} - u_{j-1,k} - u_{j+1,k} - u_{j,k+1} - u_{j,k-1} \right), \\ & [au_x]_{jk} \approx [N_1^{(x)} u]_{jk} \equiv \frac{1}{2h} \left( a_{j+1/2,k} u_{j+1,k} - a_{j-1/2,k} u_{j-1,k} \right), \\ & [bu_y]_{jk} \approx [N_1^{(y)} u]_{jk} \equiv \frac{1}{2h} \left( b_{j,k+1/2} u_{j,k+1} - b_{j,k-1/2} u_{j,k-1} \right), \\ & [p_x]_{jk} \approx [B_1^T p]_{jk} \equiv \frac{1}{h} \left( p_{j+1/2,k} - p_{j-1/2,k} \right). \end{array}$$

The second block row (associated with v) is defined analogously. The discrete convection operator  $N_1 = N_1^{(x)} + N_1^{(y)}$  represents a second-order approximation to

$$\frac{1}{2} \left[ \mathbf{w} \cdot \operatorname{grad} u + \operatorname{div} \left( u \, \mathbf{w} \right) \right];$$

this is a skew-self-adjoint version of the first convection term of (2).<sup>1</sup> The resulting matrix N is skew symmetric. The discrete incompressibility constraint is

$$(10) - [u_x + v_y]_{jk} \approx [B\mathbf{u}]_{jk} = -\left[\frac{1}{h}\left(u_{j+1/2,k} - u_{j-1/2,k}\right) + \frac{1}{h}\left(v_{j,k+1/2} - v_{j,k-1/2}\right)\right].$$

We will discuss the treatment of boundary conditions in sections 3 and 4.

**3.** Fourier analysis. Suppose that (2) is posed with constant wind **w** and periodic boundary conditions

$$u(x,0) = u(x,1), \quad u(0,y) = u(1,y)$$

For the discretization, indexing in (9) and (10) is done in mod n arithmetic. In particular, mesh points on the left and right (or top and bottom) boundaries are

<sup>&</sup>lt;sup>1</sup>This form of the convection operator leads to the skew–self-adjoint weak form used, for example, in [14, p. 53] or [28, p. 205].

identified, and if (j, k) corresponds to a pressure mesh point next to the boundary (e.g., j = 1/2), then one of its "neighboring" points is next to the opposite boundary (j = n - 1/2). The discrete versions of u, v, and p each contain  $n^2$  components, and therefore each of

(11) 
$$F_1, F_2, B_1, B_2, B_1^T, B_2^T$$

is a square matrix of order  $n^2$ . We also define a discrete convection-diffusion operator  $F_p$  on the pressure space using the first three terms of (9), that is,

$$F_p \equiv \nu A + N,$$

where A and  $N = N^{(x)} + N^{(y)}$  are specified exactly as in (9) and (j, k) now corresponds to indices for grid functions in the pressure space. An analogous idea is used in the definition of "distributive relaxation" schemes for multigrid methods applied to the Stokes and Navier–Stokes equations; see [3, 32].  $F_p$  is also defined using periodic boundary conditions for the discrete pressures. It is then straightforward to prove the following lemma by direct calculation.

LEMMA 3.1. If w is a constant vector, then  $FB^T = B^T F_p$ .

We will assume for the rest of this section that the wind **w** is constant on  $\Omega$ . The preconditioned operator can then be analyzed using Fourier techniques of the type described in [4]. It turns out that  $F_1 = F_2 = F_p$  in this case, and in some of the discussion we will refer to them collectively as  $F_*$ . Consider the discrete exponential mesh functions  $\{\psi^{(s,t)} | 0 \leq s, t < n\}$ , where

(12) 
$$\left[\psi^{(s,t)}\right]_{jk} = e^{2\pi i s x_j} e^{2\pi i t y_k}, \quad x_j = j/n, \ y_k = k/n, \qquad 0 \le j, k < n.$$

These make up an orthogonal basis for  $\mathbb{C}^{n^2}$ . The matrices associated with the periodic problem satisfy

$$F_*\psi^{(s,t)} = \left(\nu \frac{4\left(\sin^2(\pi sh) + \sin^2(\pi th)\right)}{h^2} + i \frac{w_1\sin(2\pi sh) + w_2\sin(2\pi th)}{h}\right)\psi^{(s,t)},$$
  

$$B_1\psi^{(s,t)} = \frac{1}{h}\left(1 - e^{2\pi ish}\right)\psi^{(s,t)}, \qquad B_1^T\psi^{(s,t)} = \frac{1}{h}\left(1 - e^{-2\pi ish}\right)\psi^{(s,t)},$$
  

$$B_2\psi^{(s,t)} = \frac{1}{h}\left(1 - e^{2\pi ith}\right)\psi^{(s,t)}, \qquad B_2^T\psi^{(s,t)} = \frac{1}{h}\left(1 - e^{-2\pi ith}\right)\psi^{(s,t)}.$$
  
(13)

That is, the mesh functions of (12) constitute an orthogonal basis of eigenvectors for each of these matrices. The particular choice s = t = 0 leads to  $\psi^{(0,0)} \equiv 1$ , with an associated eigenvalue equal to 0 for all the matrices of (11). Let  $S_0$  denote the space generated by  $\psi^{(0,0)}$  and let  $S_0^{\perp}$  denote the orthogonal complement of  $S_0$ . It can be verified that all eigenvalues of  $F_*$  are nonzero for other combinations of s and t, so that this matrix represents a nonsingular operator from  $S_0^{\perp}$  to itself. We can also extend  $F_*^{-1}$  to an operator on all of  $\mathbb{C}^{n^2}$  by defining it to be 0 on  $S_0$ . Therefore, it makes sense to consider the operators

(14) 
$$F^{-1}B^T, \quad B^T F_p^{-1}, \quad BF^{-1}B^T = B_1 F_1^{-1} B_1^T + B_2 F_2^{-1} B_2^T.$$

LEMMA 3.2. The Schur complement  $BF^{-1}B^T$  of (14) is a nonsingular operator from  $S_0^{\perp}$  to itself.

*Proof.* It follows immediately from (13) that

$$BF^{-1}B^T\psi^{(s,t)} = \lambda_{st}\psi^{(s,t)},$$

where

$$\lambda_{st} = \frac{4\left(\sin^2(\pi sh) + \sin^2(\pi th)\right)}{\nu \left[4\left(\sin^2(\pi sh) + \sin^2(\pi th)\right)\right] + i h \left(w_1 \sin(2\pi sh) + w_2 \sin(2\pi th)\right)}.$$

This eigenvalue is nonzero if  $s \neq 0$  or  $t \neq 0$ .  $\Box$ 

THEOREM 3.3. Let X be given by (8). If  $BF^{-1}B^T$ , X, and (the identity) I are viewed as operators from  $S_0^{\perp}$  to itself, then X is simply an alternative representation of  $BF^{-1}B^T$ , and  $BF^{-1}B^T X^{-1} = I$ .

*Proof.* It follows from Lemma 3.1 that  $F^{-1}B^T = B^T F_p^{-1}$ . Consequently,

$$\operatorname{range}(B^T) = \operatorname{range}(B^T F_p^{-1}) = \operatorname{range}(F^{-1} B^T),$$

so that (7) holds. The assertion follows from the discussion of section 1.  $\Box$ 

4. Experiments on Dirichlet problems. The results of section 3 show that (8) defines a perfect preconditioner for constant wind and periodic boundary conditions. In this section, we examine the performance of the BFBt preconditioner and some variants on more realistic problems with Dirichlet boundary conditions  $\mathbf{u} = \mathbf{g}$  on  $\partial\Omega$ .

For most of the experiments, the Oseen equations are posed on  $\Omega = (0, 1) \times (0, 1)$ and discretized with the MAC scheme. This is defined as in section 2 except when the discrete operators refer to grid indices outside  $\overline{\Omega}$  (for example, in components of the first discrete momentum equation centered at the values of u next to the bottom of  $\partial\Omega$ ). Linear extrapolation is used in these cases; see the definition of  $T_E$  in section A.2 for details. We consider two sets of coefficients, constant  $\mathbf{w} = (1, 2)$  and  $\mathbf{w} = a$  circular vortex. In the latter case,  $\mathbf{w}$  is the image of

(15) 
$$(2y(1-x^2), -2x(1-y^2))$$

under the linear mapping from  $(-1, 1) \times (-1, 1)$  to  $\Omega$ . Rather than impose explicit boundary conditions, we eliminate the discrete velocities on the boundary from the system. We use a normally distributed random vector with mean 0 and variance 1 for the right-hand side **f** of (3). We also demonstrate that the BFBt methodology is not restricted to finite differences with a set of experiments for a finite element discretization of the driven cavity problem. Both discretizations satisfy an inf-sup condition [14].

Our results are in the form of iteration counts for various combinations of  $\nu$  and h. We restrict our attention to values of  $\nu$  such that for the smallest mesh size considered,  $\frac{\mathbf{w}h}{2\nu}$  is of order 1, so that the discretizations are reasonably accurate. The iterative solver is GMRES with right-oriented preconditioning; the initial guess for all tests is  $\mathbf{u}_0 = 0, p_0 = 0$ , and the stopping criterion is

$$\frac{\|\boldsymbol{f} - \boldsymbol{\mathcal{A}} \boldsymbol{x}_{\boldsymbol{k}}\|_2}{\|\boldsymbol{f}\|_2} \le 10^{-6},$$

where f and  $x_k$  denote the right-hand side and iterate for the block system (3).

We now present the results for the BFBt preconditioner. For comparison we show analogous iteration counts for the preconditioner (4), where  $X = \frac{1}{\nu}M$  is the diagonal

Table 1

Iterations of GMRES for constant wind  $\mathbf{w} = (1, 2)$  with finite-difference discretization.

	$X = X_{BFBt}$			$X = \frac{1}{\nu}M$			
h	1/16	1/32	1/64	1/16	1/32	1/64	
$\nu = 1$	9	10	12	12	10	10	
$\nu = 1/10$	8	11	15	34	34	33	
$\nu = 1/30$	9	10	13	88	87	83	
$\nu = 1/50$	9	10	11	144	145	139	

TABLE 2 Iterations of GMRES for  $\mathbf{w} = \text{circular vortex with finite-difference discretization.}$ 

	$X = X_{BFBt}$			$X = \frac{1}{\nu}M$			
h	1/16	1/32	1/64	1/16	1/32	1/64	
$\nu = 1$	8	10	12	10	10	10	
$\nu = 1/10$	11	14	18	19	19	18	
$\nu = 1/30$	14	17	21	47	46	43	
$\nu = 1/50$	16	18	23	79	77	73	

of the scaled mass matrix as described in [5];  $X = \frac{1}{\nu}(h^2 I)$  for the MAC scheme. For the moment we are ignoring any issues of cost.<sup>2</sup>

We present three sets of results. For the MAC discretization, Table 1 shows the number of iterations for convergence of GMRES for constant wind, and Table 2 shows the results for the circular vortex. Table 3 shows analogous statistics for one example of a different discretization consisting of bilinear finite elements for both velocities and pressures, with the pressure grid of width 2h and streamline upwinding [17, p. 185] for the velocities.<sup>3</sup>

As defined, the BFBt preconditioner requires several costly subsidiary computations. For each step of GMRES, the preconditioning entails the action of  $\mathcal{Q}^{-1}$ , where  $\mathcal{Q}$  is given by (4). It can be seen from the factorization

$$\boldsymbol{\mathcal{Q}}^{-1} = \begin{pmatrix} F^{-1} & 0\\ 0 & I \end{pmatrix} \begin{pmatrix} I & B^T\\ 0 & -I \end{pmatrix} \begin{pmatrix} I & 0\\ 0 & X^{-1} \end{pmatrix}$$

that computing the action of  $\mathcal{Q}^{-1}$  entails

- computing the action of  $X^{-1}$ ;
- performing a matrix-vector product by  $B^T$ ; and
- computing the action of  $F^{-1}$ .

 $BB^T$  is a discrete Poisson operator on the pressure space (see section A.2 for a derivation for the MAC scheme), so that for the BFBt preconditioner, computing the action of  $X^{-1}$  entails solving two discrete Poisson equations. These, together with the convection-diffusion solutions, are potentially expensive operations, and the BFBt

<sup>&</sup>lt;sup>2</sup>All computations were performed in MATLAB on either a Sun SPARC-20 workstation or a DEC-Alpha 2100 4/275 workstation. For both preconditioners, the action of  $F^{-1}$  was computed using Gaussian elimination. For the BFBt preconditioner the action of  $(BB^T)^{-1}$  was computed by direct methods using the pseudoinverse of  $BB^T$ , except when h = 1/64 for the MAC discretization. In that case, this computation was done using a multigrid iteration in which the relative residual is forced to be less than  $10^{-8}$ .

<sup>&</sup>lt;sup>3</sup>The test problem is slightly different here. It is posed on  $\Omega = (-1, 1) \times (-1, 1)$  with **w** as in (15), **f** = 0 in (1), and boundary conditions u = v = 0 when  $x = \pm 1$  or y = -1, and u = 1, v = 0 when y = 1. See [5] for more details. Some of the entries in Table 3 are taken from [5].

Table 3

Iterations of GMRES for  $\mathbf{w} = circular$  vortex with bilinear finite element discretization.

	$X = X_{BFBt}$			$X = \frac{1}{\nu}M$			
h	2/16	2/32	2/64	2/16	2/32	2/64	
$\nu = 1$	7	9	11	21	22	21	
$\nu = 1/10$	10	12	15	32	36	35	
$\nu = 1/30$	13	15	17	44	56	64	
$\nu = 1/50$	15	17	19	48	72	97	

TABLE	4
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Iterations of GMRES with modified BFBt preconditioning using multigrid for the Poisson equation. Wind  $\mathbf{w} = (1, 2)$ , finite-difference discretization.

	$X = X_{BFBt/MG}$							
h	1/16	1/32	1/64	1/128				
$\nu = 1$	11	12	15	19				
$\nu = 1/10$	12	13	17	22				
$\nu = 1/30$	12	12	15	20				
$\nu = 1/50$	13	13	14	18				
$\nu = 1/100$				14				

preconditioner is significantly more costly than when  $X = \frac{1}{\nu}M$ , a diagonal matrix. We now examine what happens when less costly computations based on inner iteration are used in place of these three operations. We consider only finite differences here, although the same methodologies are applicable to other discretizations. These versions of the preconditioner also require less storage, and in particular this enabled us to further explore some trends in the data with a finer mesh size h = 1/128 and smaller viscosity parameter  $\nu = 1/100$  (used only with this fine mesh).

We first consider the effect of replacing the Poisson solutions with approximations derived from one step of the V-cycle multigrid [21, Chap. 1].<sup>4</sup> That is, the modified BFBt preconditioner uses

$$X = S_{MG} \left( BFB^T \right) S_{MG},$$

where  $S_{MG}$  is the multigrid approximation to  $(BB^T)^{-1}$ . The computational costs of this algorithm are of the same order of magnitude as when  $X = \frac{1}{\nu}M$ . (We are still solving the convection-diffusion equations exactly except when h = 1/128; in this case, we use an iterative method based on relaxation and force the relative residual norm to be less than  $10^{-8}$ .) The iteration counts for the two sets of benchmark problems are shown in Tables 4 and 5.

Next, we consider the effect of also replacing the convection-diffusion solutions with approximate solutions derived from iterative methods. We will comment on the choice of method below. The performance of any such method will depend on the relative amount of convection and diffusion in the problem, i.e., the value of  $\nu$  [7]. Therefore, rather than use a fixed number of iterations, we perform the inner iteration until the stopping criterion

(16) 
$$\frac{\|w - Fv_k\|_2}{\|w\|_2} \le \tau = 10^{-2}$$

<sup>&</sup>lt;sup>4</sup>The multigrid computation used damped Jacobi smoothing with optimal smoothing parameter  $\omega = 4/5$ , one presmoothing and one postsmoothing step, and bilinear interpolation for prolongation.

Iterations of GMRES with modified BFBt preconditioning using multigrid for the Poisson equation. Wind  $\mathbf{w} = \text{circular vortex}$ , finite-difference discretization.

	$X = X_{BFBt/MG}$					
h	1/16	1/32	1/64			
$\nu = 1$	11	12	15			
$\nu = 1/10$	14	16	20			
$\nu = 1/30$	19	21	24			
$\nu = 1/50$	21	24	27			

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Iterations of FGMRES with modified BFBt preconditioning using multigrid for the Poisson equation and iteration for the convection-diffusion equation. Wind  $\mathbf{w} = (1, 2)$ , finite-difference discretization.

	$X = X_{BFBt/MG/Iter}$				$X = \frac{1}{\nu}M$			
h	1/16	1/32	1/64	1/128	1/16	1/32	1/64	1/128
$\nu = 1$	11	13	16	20	12	11	12	12
$\nu = 1/10$	12	14	17	22	35	34	33	32
$\nu = 1/30$	12	13	15	20	111	88	85	85
$\nu = 1/50$	-	13	14	18	_	185	141	142
$\nu = 1/100$				15				

is satisfied, where w represents the right-hand side for each convection-diffusion equation and  $v_0 = 0$  is the initial guess. The number of these iterations may vary from step to step of the outer GMRES iteration, so that we are no longer using a fixed preconditioner  $\mathcal{Q}$ ; instead we have a series of operators  $\mathcal{Q}_k$  that vary with the GMRES step. This may cause difficulties for GMRES, but these can be avoided with a "flexible" variant of GMRES (FGMRES) designed for this situation [25], which we use in these tests. The two algorithms are mathematically equivalent when  $\mathcal{Q}_k$  is fixed. The results for constant wind are shown in Table 6, and for the circular vortex in Table 7. We also present the performance using the scaled mass matrix. A dash (-) indicates failure to converge; this was due to divergence of the iteration for the convection-diffusion equation and is an artifact of inaccuracy of the discretization.

We highlight some trends displayed by these data as follows.

- 1. In all tests with the constant coefficient problem (Tables 1 and 6, left, and Table 4), iteration counts with the BFBt preconditioner are independent of the viscosity  $\nu$ . Indeed, in some cases they actually decrease slightly with  $\nu$ . For the circular vortex, however, the iteration counts appear not to be independent of  $\nu$ , instead exhibiting some growth as  $\nu$  decreases (Tables 2 and 7, left, and Table 5).
- 2. In all cases, the counts with BFBt preconditioning increase slowly with  $h^{-1}$ .
- 3. The iteration counts for  $X = \frac{1}{\nu}M$  are independent of  $h^{-1}$ , but they grow roughly linearly with  $1/\nu$ ; these trends are consistent with the analysis of [5]. The counts are smaller for the circular vortex than for the constant wind; we have no explanation for this. However, the same qualitative patterns are present for both problems. In practical situations, it is often desired to compute solutions of a fixed accuracy for a variety of values of  $\nu$  by letting  $h \to 0$  and  $\nu \to 0$  simultaneously; with respect to this criterion, the BFBt preconditioner requires significantly fewer iterations as the viscosity decreases.
- 4. The results for bilinear finite elements (Table 3) are qualitatively the same as for finite differences.

TABLE	7
	•

Iterations of FGMRES with modified BFBt preconditioning using multigrid for the Poisson equation and iteration for the convection-diffusion equation. Wind  $\mathbf{w} = \text{circular vortex}$ , finite-difference discretization.

	$X = X_{BFBt/MG/Iter}$				$X = \frac{1}{\nu}M$			
h	1/16	1/32	1/64	1/128	1/16	1/32	1/64	1/128
$\nu = 1$	11	13	16	19	11	12	12	12
$\nu = 1/10$	14	16	20	25	19	19	19	18
$\nu = 1/30$	19	21	24	31	51	45	44	43
$\nu = 1/50$	—	28	27	34	_	95	73	73
$\nu = 1/100$				37				155

- 5. Comparison of Tables 4 and 5 with Tables 1 and 2 shows that replacing the exact Poisson solution with one multigrid step leads to little increase in iteration counts (less than 25% for h = 1/64). The same trends with respect to  $\nu$  and h are evident, and these carry over to h = 1/128. Note that the multigrid method used here is not necessarily an optimal choice, just a simple one.
- 6. Approximate solution of the convection-diffusion problem produces little degradation in performance of either preconditioner, even though the tolerance  $\tau$  in (16) is very mild.

We also ran some tests with the quasi-minimum residual (QMR) iteration without look-ahead [10], which (in contrast to GMRES) has a fixed cost per step. Performance for both the BFBt preconditioner and for  $X = \frac{1}{\nu}M$  was qualitatively similar to that of GMRES.

We return to the first two items on this list: the dependence of iteration counts with BFBt preconditioning on the parameters  $\nu$  (for the circular vortex) and h. To try to get a feeling for this dependence, we plot the behavior of the solver along with graphs of functions that might model this behavior. In particular, Figure 2 plots iteration counts as a function of  $\nu^{-1}$  using the entries of Table 7 for h = 1/128, together with graphs of the functions  $\nu^{-1}$  and  $\nu^{-1/2}$ ; the latter curves were shifted to make the figure easy to view. These results suggest that dependence on  $\nu$  is like  $O(\nu^{-1/2})$ , although it is difficult to make a precise statement with this limited amount of data. Similarly, Figure 3 plots iteration counts as a function of  $h^{-1}$  using the entries from Table 7 for  $\nu = 1/30$ , and these are compared with  $h^{-1}$  and  $h^{-1/2}$ . These results typify the behavior for all the values  $\nu$  in the table and suggest that the dependence on h is of order  $h^{-1/2}$ ; see also section A.2.

We comment briefly on the convection-diffusion solvers. For simple flows, it is fairly easy to construct relaxation strategies that follow the flow and converge rapidly [7], but for more complex flows, especially with recirculations, this is a more difficult task. In these tests, we used a horizontal one-line SOR iteration with relaxation parameter given by the optimal choice for the constant coefficient problem, which can be computed analytically [8]. We used this for the circular wind only for convenience of coding, and convergence was slow for these problems. Many other options for solving this problem (such as multigrid) are available. We also ran some tests with another inner iteration (using a "multidirectional relaxation") and found the number of outer iterations not to depend significantly on the choice of the method used to solve the convection-diffusion equation. This points to the importance of the convectiondiffusion problem for the BFBt preconditioner; it is critical that the approximate



FIG. 2. Comparison of iteration counts (from Table 7) with various functions of  $\nu$  for h = 1/128.



FIG. 3. Comparison of iteration counts (from Table 7) with various functions of h for  $\nu = 1/30$ .

solution to this subproblem be computed efficiently for the complete computation to be inexpensive.

5. Concluding remarks. We summarize this study as follows. The main new result is that the performance of the BFBt preconditioner for solving the steady-state Oseen equations with Dirichlet boundary conditions depends very mildly on the viscosity  $\nu$ . This stands in contrast to other preconditioning methods, where performance deteriorates more dramatically as  $\nu \to 0$ . Its performance depends on the mesh size h with iterations apparently growing in proportion to  $h^{-1/2}$ . If a series of Dirichlet problems with decreasing viscosity is to be solved in such a way that h is proportional to  $\nu$ , then we expect the iteration counts to grow like  $O(\nu^{-1/2})$ .

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This conclusion also appears to hold for "inexact" versions of the BFBt preconditioner in which the expensive subsidiary computations, solution of the Poisson equation and convection–diffusion equation, are replaced by iterative solutions with mild stopping criteria. The effectiveness of these variants of the BFBt preconditioner depends on having efficient methods for the convection–diffusion equation.

This issue is also critical for multigrid methods for the Navier–Stokes equations [3, 29, 30, 32]. It is difficult to make a simple comparison between the methods proposed here and multigrid; most reported studies show a dependence on viscosity (or Reynolds number) but not on mesh size [29, 30, 32]. An advantage of the BFBt preconditioner is that it can be applied easily to systems arising from mixed finite elements with different grids for velocities and pressure (as shown in Table 3).

If the goal is to solve a fixed Dirichlet problem on a sequence of finer meshes, then there are other methods whose performance is less dependent on the mesh size [5, 11]. This is true in particular for the Stokes equations, where several methods exist whose convergence rate is independent of h [1, 2, 24, 27].

Finally, we note that the boundary conditions are very important here. Fourier analysis is often used as a guideline in behavior of numerical methods [23], and there are numerous examples of cases where the results for periodic boundary conditions are predictive of performance for Dirichlet conditions [4]. However, here there is a qualitative difference in the two types of problems. See [20] for other examples of the effects of boundary conditions on iterative methods.

A. Appendix: Partial analysis for the preconditioned Stokes problem. We show that the minimum eigenvalue for the BFBt-preconditioned Stokes operator is no smaller than 1 and give a partial analysis that suggests why its maximum eigenvalue is proportional to  $h^{-1}$ .

**A.1. Lower bound.** Suppose the coefficient matrix of (3) is a discrete Stokes operator; i.e., it is derived from  $\nu = 1$ ,  $\mathbf{w} = 0$  in (2), with Dirichlet boundary conditions. Therefore, F = A. Let the singular value decomposition of B be denoted

$$B = U\Sigma V^T = \begin{bmatrix} U_1, u_0 \end{bmatrix} \left( \begin{pmatrix} \Sigma_1 & 0 \\ 0 & 0 \end{pmatrix} & 0 \end{pmatrix} \begin{pmatrix} V_1^T \\ V_2^T \end{pmatrix} = U_1 \Sigma_1 V_1^T$$

The columns of  $U_1$  span range(B), the columns of  $V_1$  span range $(B^T)$ , and  $u_0$  is parallel to the discrete hydrostatic pressure  $p \equiv 1$ , which determines the null space  $S_0$  of  $B^T$ . It follows that

$$BB^T = U_1 \Sigma_1^2 U_1^T.$$

As in section 3, we will treat  $B^T$  as an operator defined on  $\mathcal{S}_0^{\perp}$ ; on this space,  $BB^T$  and  $BAB^T$  represent nonsingular operators whose inverses are given by the matrix pseudoinverses

$$(BB^{T})^{-1} = U_{1}\Sigma_{1}^{-2}U_{1}^{T}, (BAB^{T})^{-1} = U_{1}\Sigma_{1}^{-1}(V_{1}^{T}AV_{1})^{-1}\Sigma_{1}^{-1}U_{1}^{T}.$$

We are interested in the minimum eigenvalue of the generalized eigenvalue problem

(17) 
$$(BA^{-1}B^T)p = \lambda (BB^T) (BAB^T)^{-1} (BB^T)p.$$

The Schur complement matrix on the left is

$$S \equiv BA^{-1}B^{T} = U_{1}\Sigma_{1}V_{1}^{T}A^{-1}V_{1}\Sigma_{1}U_{1}^{T},$$

and the preconditioning matrix on the right is

$$X = U_1 \Sigma_1 (V_1^T A V_1)^{-1} \Sigma_1 U_1^T.$$

Consequently,

(18) 
$$S = X + U_1 \Sigma_1 \left[ V_1^T A^{-1} V_1 - (V_1^T A V_1)^{-1} \right] \Sigma_1 U_1^T.$$

Using this splitting of S, we can prove the following result.

THEOREM A.1. The minimum eigenvalue of the preconditioned generalized Stokes operator with Dirichlet boundary conditions is bounded below by 1.

*Proof.* We will show that  $V_1^T A^{-1} V_1 - (V_1^T A V_1)^{-1}$  is positive semidefinite. Let

$$D = V^T A V = \begin{pmatrix} D_{11} & D_{12} \\ D_{21} & D_{22} \end{pmatrix},$$

where  $D_{ij} = V_i^T A V_j$ . D is symmetric positive definite and standard analysis gives

$$D = \begin{pmatrix} I & 0 \\ D_{21}D_{11}^{-1} & I \end{pmatrix} \begin{pmatrix} D_{11} & 0 \\ 0 & D_{22} - D_{21}D_{11}^{-1}D_{12} \end{pmatrix} \begin{pmatrix} I & D_{11}^{-1}D_{12} \\ 0 & I \end{pmatrix}.$$

Letting  $C_D \equiv D_{22} - D_{21}D_{11}^{-1}D_{12}$ , this in turn implies that

$$D^{-1} = \begin{pmatrix} D_{11}^{-1} + D_{11}^{-1} D_{12} C_D^{-1} D_{21} D_{11}^{-1} & -D_{11}^{-1} D_{12} C_D^{-1} \\ -C_D^{-1} D_{21} D_{11}^{-1} & C_D^{-1} \end{pmatrix}.$$

However,  $D^{-1} = V^T A^{-1} V$ , so that

$$V_1^T A^{-1} V_1 = [D^{-1}]_{11} = D_{11}^{-1} + D_{11}^{-1} D_{12} C_D^{-1} D_{21} D_{11}^{-1}.$$

Therefore,

$$\inf_{x \neq 0} \frac{(x, V_1^T A^{-1} V_1 x)}{(x, (V_1 A V_1)^{-1} x)} = \inf_{x \neq 0} \frac{(x, (D_{11}^{-1} + D_{11}^{-1} D_{12} C_D^{-1} D_{21} D_{11}^{-1}) x)}{(x, D_{11}^{-1} x)} \\
= \inf_{x \neq 0} \frac{(x, (D_{11} + D_{12} C_D^{-1} D_{21}) x)}{(x, D_{11} x)} \ge 1.$$

It follows that  $V_1^T A^{-1} V_1 - (V_1^T A V_1)^{-1}$  is positive semidefinite, so that (18) yields

$$\inf_{q \in \mathcal{S}_0^{\perp}} \frac{(q, Sq)}{(q, Xq)} \ge 1. \qquad \Box$$

**A.2. Largest eigenvalue.** The experiments described in section 4 show that convergence rates for the BFBt preconditioner depend on the mesh size. Figure 4 plots the maximum eigenvalue of (17) as a function of  $h^{-1}$ , where the computations were for h = 1/4, 1/8, 1/16, and 1/32. The results indicate that the maximum eigenvalue is of magnitude  $O(h^{-1})$ . By analogy with the standard convergence bounds for symmetric positive definite problems, this suggests that the iteration counts will increase in proportion to  $h^{-1/2}$ , which is consistent with the results of section 4.



FIG. 4. Maximum eigenvalue for the Stokes problem with BFBt preconditioner.

To provide some insight into this maximum eigenvalue, we write

$$\frac{(q, Sq)}{(q, Xq)} = \frac{(q, Sq)}{(q, (h^2I)q)} \frac{(q, (h^2I)q)}{(q, Xq)} \,,$$

where  $h^2 I$  is playing the role of the mass matrix for the finite-difference discretization. It is well known [9] that the first quotient on the right is bounded by a constant  $\Gamma \leq \sqrt{2}$ . We would like a bound for the second quotient of the form

$$\sup_{q \in \mathcal{S}_{0}^{\perp}} \ \frac{(q, (h^{2}I)q)}{(q, (BB^{T}) (BAB^{T})^{-1} (BB^{T})q)} \leq \frac{c}{h}$$

or, equivalently,

(19) 
$$\sup_{q \in S_{+}^{\perp}} \frac{(q, BAB^T q)}{(q, (BB^T)^2 q)} \le \frac{c}{h^3}.$$

The matrices comprising (19) have a tensor product structure derived from onedimensional operators. Let the interval [0,1] be divided into n equally spaced subintervals of width h = 1/n, where the *j*th interval is  $[x_{j-1}, x_j]$  with  $x_j = jh$  and midpoint  $\hat{x}_j = (j - 1/2)h$ . An example with n = 8 is

Consider the following three (scaled) finite-difference approximations to the onedimensional Laplace operator -u'':

$$T_D = \begin{pmatrix} 2 & -1 \\ -1 & 2 & -1 \\ & \ddots & \\ & -1 & 2 & -1 \\ & & -1 & 2 \end{pmatrix}, T_E = \begin{pmatrix} 3 & -1 \\ -1 & 2 & -1 \\ & \ddots & \\ & -1 & 2 & -1 \\ & & -1 & 3 \end{pmatrix}, T_N = \begin{pmatrix} 1 & -1 \\ -1 & 2 & -1 \\ & \ddots & \\ & -1 & 2 & -1 \\ & & -1 & 1 \end{pmatrix}$$

 $T_D$  and  $T_E$  are derived from Dirichlet boundary conditions (u(0), u(1) given), and  $T_N$  is derived from Neumann conditions (u'(0), u'(1) given).  $T_D$  is defined at the cell

boundaries  $\{x_j\}_{j=1}^{n-1}$  and is of order n-1;  $T_E$  and  $T_N$  are defined at the cell centers  $\{\hat{x}_j\}_{j=1}^n$  and are of order n. In all cases, the discrete Laplace operator at points next to the boundaries depends on the boundary conditions.  $T_E$  uses the linear extrapolation

$$u(\hat{x}_0) = -u(\hat{x}_1) + 2u(0), \quad u(\hat{x}_{n+1}) = -u(\hat{x}_n) + 2u(1);$$

 $T_N$  approximates the Neumann boundary conditions as

$$u'(0) \approx [u(\hat{x}_1) - u(\hat{x}_0)]/h, \quad u'(1) \approx [u(\hat{x}_{n+1}) - u(\hat{x}_n)]/h.$$

Note that

(20) 
$$T_E = T_N + 2E_0, \quad E_0 = \begin{pmatrix} 1 & & \\ 0 & & \\ & \ddots & \\ & & 0 \\ & & & 1 \end{pmatrix}.$$

We will also use the difference operator

$$B_D = \begin{pmatrix} -1 & & \\ 1 & -1 & & \\ & \ddots & \ddots & \\ & & 1 & -1 \\ & & & 1 \end{pmatrix},$$

which has dimensions  $n \times (n-1)$  and can be viewed as mapping cell boundary grid functions to cell centered functions. It is easily verified that

(21) 
$$B_D B_D^T = T_N, \qquad B_D T_D B_D^T = T_N^2.$$

Recall that for matrices X and Y, where X has dimensions  $r \times s$ , the tensor product of X and Y is [16, pp. 239ff]

$$X \otimes Y = \begin{pmatrix} x_{11}Y & \cdots & x_{1,s}Y \\ \vdots & \ddots & \vdots \\ x_{r,1}Y & \cdots & x_{r,s}Y \end{pmatrix}.$$

Letting  $I_r$  denote the identity matrix of order r, it is straightforward to show that

$$A = \begin{pmatrix} I_n \otimes T_D + T_E \otimes I_{n-1} & 0\\ 0 & I_{n-1} \otimes T_E + T_D \otimes I_n \end{pmatrix},$$
  
$$B = [I_n \otimes hB_D, hB_D \otimes I_n].$$

The identities (21) then imply

$$BB^{T} = h^{2}(I_{n} \otimes T_{N} + T_{N} \otimes I_{n}),$$
  

$$(BB^{T})^{2} = h^{4}(I_{n} \otimes T_{N}^{2} + 2T_{N} \otimes T_{N} + T_{N}^{2} \otimes I_{n}),$$
  

$$BAB^{T} = h^{2}(I_{n} \otimes T_{N}^{2} + T_{E} \otimes T_{N} + T_{N} \otimes T_{E} + T_{N}^{2} \otimes I_{n}).$$

Consequently, relation (19) is equivalent to

(22) 
$$\sup_{q\in\mathcal{S}_{n}^{\perp}} \frac{\left(q, \left(I_{n}\otimes T_{N}^{2}+T_{E}\otimes T_{N}+T_{N}\otimes T_{E}+T_{N}^{2}\otimes I_{n}\right)q\right)}{\left(q, \left(I_{n}\otimes T_{N}^{2}+2T_{N}\otimes T_{N}+T_{N}^{2}\otimes I_{n}\right)q\right)} \leq \frac{c}{h}$$

*Remark.* This derivation shows that  $BB^T$  is a scaled discrete Poisson operator on the pressure space with Neumann boundary conditions.

The matrices in (22) differ only by the presence of  $T_E$  in the cross-terms of the numerator; i.e., the numerator contains submatrices derived from Dirichlet conditions, whereas the denominator comes exclusively from Neumann conditions. To get a feeling for why this difference in boundary conditions leads to an inequality of the form (22), consider the generalized eigenvalue problem for the one-dimensional operators,

(23) 
$$T_N x = \lambda T_E x.$$

It is evident from (20) that there is an eigenvalue  $\lambda = 1$  of multiplicity n-2 for which every eigenvector x satisfies  $x_1 = x_n = 0$ . Moreover,  $T_N x = 0$  for the constant vector  $x \equiv 1$ , so that 0 is also an eigenvalue. If any x satisfies (23) with  $\lambda \neq 0$  and at least one of  $x_1 \neq 0$  or  $x_n \neq 0$ , then the first and last equations (of (23)) imply  $\lambda \neq 1$ . From the first equation we have

(24) 
$$x_2 - x_1 = \frac{2\lambda}{\lambda - 1} x_1.$$

The assumption  $\lambda \neq 1$  together with the (interior) equations  $k = 2, \ldots, n-1$  imply

$$x_{k+1} - x_k = x_k - x_{k-1} = \frac{2\lambda}{\lambda - 1} x_1, \quad 2 \le k \le n,$$

where the second equality comes from (24). An inductive argument then yields

(25) 
$$x_k = \frac{(2k-1)\lambda - 1}{\lambda - 1} x_1, \qquad 2 \le k \le n.$$

Finally, the last equation of (23) implies

$$x_n = \frac{1-\lambda}{2\lambda} \left( x_n - x_{n-1} \right) = -x_1.$$

Equating this expression for  $x_n$  with the one from (25) yields  $\lambda = 1/n = h$ . Therefore, we have the following result.

THEOREM A.2. The eigenvalues of the generalized eigenvalue problem (23) are  $\lambda = 1$  of multiplicity n - 2,  $\lambda = 0$ , and  $\lambda = h$ . Consequently,

$$\sup_{x} \frac{(x, T_E x)}{(x, T_N x)} = \frac{1}{h},$$

where the supremum is taken over all vectors orthogonal to the constant vector.

*Remark.* The results of section A.1 apply in general to any discretization for which the null space is spanned by constant pressures. The discussion of section A.2 applies only for the MAC scheme.

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