

A non-alternating preconditioned HSS iteration method for non-Hermitian positive definite linear systems

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Received: 30 June 2014 / Revised: 9 April 2015 / Accepted: 10 April 2015 © SBMAC - Sociedade Brasileira de Matemática Aplicada e Computacional 2015

Abstract By utilizing the preconditioned Hermitian and skew-Hermitian splitting (PHSS) iteration technique, we establish a non-alternating PHSS (NPHSS) iteration method for solving large sparse non-Hermitian positive definite linear systems. The convergence analysis demonstrates that the iterative series produced by the NPHSS method converge to the unique solution of the linear system when the parameters satisfy some moderate conditions. We also give a possible optimal upper bound for the iterative spectral radius. Moreover, to reduce the computational cost, we establish an inexact variant of the NPHSS (INPHSS) iteration method whose convergence property is studied. Both theoretical and numerical results validate that the NPHSS method outperforms the PHSS method when the Hermitian part of the coefficient matrix is dominant.

Communicated by Ya-xiang Yuan.

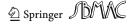
Published online: 29 April 2015

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Keywords Non-Hermitian positive definite linear systems · Hermitian and skew-Hermitian splitting · Non-alternating PHSS iteration · Spectral radius · Convergence analysis

Mathematics Subject Classification 65F10 · 65F15 · 65T10

1 Introduction

We consider the system of linear equations

$$Ax = b, \quad A \in \mathbb{C}^{n \times n}, \quad x, b \in \mathbb{C}^n,$$
 (1)

where $A \in \mathbb{C}^{n \times n}$ is a large sparse non-Hermitian and positive definite matrix. Such linear systems arise in many problems in scientific computing and engineering applications.

Since the matrix $A \in \mathbb{C}^{n \times n}$ naturally possesses the Hermitian and skew-Hermitian splitting (HSS)

$$A = H + S$$
,

where

$$H = \frac{1}{2}(A + A^*)$$
 and $S = \frac{1}{2}(A - A^*)$,

HSS (Bai et al. 2003) and preconditioned HSS (PHSS) (Bai et al. 2004, 2007a) iteration methods were proposed to compute the approximate solution of the linear system (1). In fact, the PHSS method was defined as follows.

Algorithm 1 (*The PHSS iteration method*) Let $x^{(0)} \in \mathbb{C}^n$ be an arbitrary initial guess. For $k = 0, 1, 2, \ldots$ until the sequence of iterates $\{x^{(k)}\}_{k=0}^{\infty} \subset \mathbb{C}^n$ converges, compute the next iterate $x^{(k+1)}$ according to the following procedure:

$$\begin{cases} (\alpha P + H)x^{(k+\frac{1}{2})} = (\alpha P - S)x^{(k)} + b, \\ (\alpha P + S)x^{(k+1)} = (\alpha P - H)x^{(k+\frac{1}{2})} + b, \end{cases}$$
(2)

where α is a given positive constant and $P \in \mathbb{C}^{n \times n}$ is a prescribed Hermitian positive definite matrix.

We can see that each iterate of the PHSS iteration alternates between H and S, analogously to the classical alternating direction implicit (ADI) iteration for solving partial differential equations; see Peaceman and Rachford (1955), Douglas (1962). The choice of preconditioner P is mostly dependent on the structure of the coefficient matrix A. For example, we can choose a preconditioner based on incomplete Cholesky factorization, incomplete LU factorization (Saad 2003), and incremental unknowns (Chen and Temam 1993; Yang and Wu 2014; Yang et al. 2013, 2015); see Benzi (2002), Chen (2005) for a detailed discussion.

In particular, if we choose P=I, the identity matrix, the PHSS iteration method is reduced to the HSS iteration method (Bai et al. 2003). When $P \neq I$, we can suitably choose P and α such that the induced PHSS iteration method achieves fast convergence and high computing efficiency. Theoretical analysis shows that both methods converge unconditionally to the unique solution of the linear system (1). Due to its promising performance and elegant mathematical properties, the HSS iteration has attracted many researchers' attention, resulting in numerous papers devoted to various aspects of this algorithm; see Bai et al. (2005, 2006, 2007b, 2008, 2010, 2011), Bai (2007, 2009, 2010), Benzi (2009), Bai and Golub (2007), Li et al. (2007a,b), Li et al. (2014a,b), Salkuyeh and Behnejad (2012), Yang et al. (2010), Yin and Dou (2012), Cao et al. (2012), Huang (2014).



However, the challenges of the HSS and PHSS iteration methods lie in solving the shifted skew-Hermitian sub-system of linear equations at each iteration step, which is as difficult as that of the original problem. In this work, we present a non-alternating PHSS (NPHSS) iteration method for solving non-Hermitian positive definite linear systems (1). The NPHSS iteration method can be described as follows:

Algorithm 2 (*The NPHSS iteration method*) Let $x^{(0)} \in \mathbb{C}^n$ be an arbitrary initial guess. For $\ell = 0, 1, 2, \ldots$ until the sequence of iterates $\{x^{(\ell)}\}_{\ell=0}^{\infty} \subset \mathbb{C}^n$ converges, compute the next iterate $x^{(\ell+1)}$ according to the following procedure:

$$(\alpha P + H)x^{(\ell+1)} = (\alpha P - S)x^{(\ell)} + b,$$
(3)

where α is a given positive constant and $P \in \mathbb{C}^{n \times n}$ is a prescribed Hermitian positive definite matrix.

Due to Hermitian positive definiteness of the matrix $\alpha P + H$, every sub-system in (3) can be effectively solved either exactly by a sparse Cholesky factorization, or inexactly by a preconditioned conjugate gradient (PCG) scheme (Saad 2003). Theoretical analysis shows that the iterative sequence produced by NPHSS iteration method converges to the unique solution of the linear system (1), with a loose restriction on the choice of α . The contraction factor of the NPHSS iteration can be bounded by a function that depends only on the choice of α , the smallest eigenvalue of matrix $P^{-1}H$, and the maximum modulus of eigenvalue of matrix $P^{-1}S$.

In particular, if we choose P = I, the NPHSS iteration method is reduced to the nonalternating HSS (NHSS) iteration method, which was originally proposed in Axelsson et al. (2004) and recently discussed in Li and Wu (2015). Therefore, our method is a preconditioned generalization of the method in Axelsson et al. (2004) and Li and Wu (2015). Moreover, based on the results in Axelsson et al. (2004) a more general framework of the splitting iteration method is established and its inexact variant using the preconditioned conjugate gradient method as the inner solver is discussed in our paper. The comparison of the convergence speeds of PHSS and NPHSS iteration methods is also discussed. A recent work related to this is Axelsson et al. (2014).

This paper is organized as follows. In Sect. 2, we analyze the convergence properties of the NPHSS iteration for linear system (1), including convergence condition, spectral radius of the iterative matrix, the choices of iterative parameter, etc. The comparison of the convergence speeds of PHSS and NPHSS iteration methods is discussed in Sect. 3. In Sect. 4, an inexact variant of the NPHSS (INPHSS) iteration method is presented and its convergence property is discussed. Numerical results are presented in Sect. 5 to illustrate the effectiveness of our methods. Finally, in Sect. 6, we end this work with a brief conclusion.

2 Convergence analysis of the NPHSS iteration method

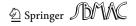
In this section, we first consider the convergence properties of the NPHSS iteration method. We can reformulate the NPHSS iteration scheme (3) as

$$x^{(\ell+1)} = M(P; \alpha)x^{(\ell)} + N(P; \alpha)b, \quad \ell = 0, 1, 2...,$$

where

$$M(P; \alpha) = (\alpha P + H)^{-1} (\alpha P - S)$$
 and $N(P; \alpha) = (\alpha P + H)^{-1}$. (4)

Here, $M(P; \alpha)$ is the iteration matrix of the NPHSS iteration method.



Note that $M(P; \alpha)$ can be reformulated as

$$M(P; \alpha) = I - (\alpha P + H)^{-1} A,$$

then the matrix $\alpha P + H$ can be viewed as a preconditioner for the coefficient matrix $A \in \mathbb{C}^{n \times n}$.

Since H is Hermitian and S is skew-Hermitian, we easily know that all eigenvalues of $P^{-1}H$ are real and positive, and all eigenvalues of $P^{-1}S$ are imaginary. Here and in the sequel, denote

$$\lambda_{\max} = \max_{\lambda_j \in sp(P^{-1}H)} \{\lambda_j\}, \quad \lambda_{\min} = \min_{\lambda_j \in sp(P^{-1}H)} \{\lambda_j\} \quad \text{and} \quad \xi_{\max} = \max_{i \notin j \in sp(P^{-1}S)} \{|\xi_j|\},$$

where sp(X) denotes the spectrum of the matrix X and $i = \sqrt{-1}$.

The following theorem gives the convergence result of the NPHSS iteration method.

Theorem 1 Let $A \in \mathbb{C}^{n \times n}$ be a positive definite matrix, let $H = \frac{1}{2}(A + A^*)$ and $S = \frac{1}{2}(A - A^*)$ be its Hermitian and skew-Hermitian parts, respectively, and let α be a positive constant. Let $P \in \mathbb{C}^{n \times n}$ be a Hermitian positive definite matrix. Then, the spectral radius $\rho(M(P;\alpha))$ of the NPHSS iteration matrix (4) satisfies $\rho(M(P;\alpha)) \leq \sigma(\alpha)$, where

$$\sigma(\alpha) = \frac{\sqrt{\alpha^2 + \xi_{\text{max}}^2}}{\alpha + \lambda_{\text{min}}}.$$
 (5)

Moreover, it holds that

- (i) If $\lambda_{min} \geq \xi_{max}$, then $\sigma(\alpha) < 1$ for any $\alpha > 0$, which means that the NPHSS iteration method is unconditionally convergent;
- (ii) if $\lambda_{\min} < \xi_{\max}$, then $\sigma(\alpha) < 1$ if and only if

$$\alpha > \frac{\xi_{\text{max}}^2 - \lambda_{\text{min}}^2}{2\lambda_{\text{min}}},\tag{6}$$

which means that the NPHSS iteration method is convergent under the condition (6).

Proof By direct computations, we have

$$\begin{split} \rho(M(P;\alpha)) &= \rho \left((\alpha P + H)^{-1} (\alpha P - S) \right) \\ &= \rho \left(\left(\alpha I + P^{-1} H \right)^{-1} \left(\alpha I - P^{-1} S \right) \right) \\ &\leq \left\| \left(\alpha I + P^{-1} H \right)^{-1} \left(\alpha I - P^{-1} S \right) \right\|_2 \\ &\leq \left\| \left(\alpha I + P^{-1} H \right)^{-1} \right\|_2 \left\| \left(\alpha I - P^{-1} S \right) \right\|_2 \\ &= \max_{\lambda_j \in sp(P^{-1} H)} \left| \frac{1}{\alpha + \lambda_j} \right| \cdot \max_{i \xi_j \in sp(P^{-1} S)} \left| \alpha - i \xi_j \right| \\ &= \max_{\lambda_j \in sp(P^{-1} H)} \frac{1}{\alpha + \lambda_j} \cdot \max_{i \xi_j \in sp(P^{-1} S)} \sqrt{\alpha^2 + \xi_j^2} \\ &= \frac{\sqrt{\alpha^2 + \xi_{\max}^2}}{\alpha + \lambda_{\min}}. \end{split}$$

Thus the upper bound of $\rho(M(P; \alpha))$ given in (5) is obtained.

By simple derivations, $\sigma(\alpha) < 1$ is equivalent to

$$2\lambda_{\min}\alpha > \xi_{\max}^2 - \lambda_{\min}^2. \tag{7}$$



If $\lambda_{\min} \geq \xi_{\max}$, then (7) holds true for any $\alpha > 0$, i.e., the NPHSS iteration converges to the unique solution of the system of linear equations (1); if $\lambda_{\min} < \xi_{\max}$, then (7) or $\sigma(\alpha) < 1$ follows if and only if α satisfies (6). Therefore, in case (ii), the sufficient and necessary condition for the convergence of the NPHSS iteration method is inequality (6).

Theorem 1 gives the convergence conditions of the NPHSS iteration method by analyzing the upper bound $\sigma(\alpha)$ of the spectral radius of the iteration matrix $M(P; \alpha)$. Since the optimal parameter α minimizing the spectral radius $\rho(M(P; \alpha))$ is hardly obtained, we instead give the parameter α^* , which minimizes the upper bound $\sigma(\alpha)$ of the spectral radius $\rho(M(P; \alpha))$, in the following corollary.

Corollary 1 *Let the conditions of Theorem* 1 *be satisfied. Then, the parameter* α^* *minimizing the upper bound* $\sigma(\alpha)$ *of the spectral radius* $\rho(M(P; \alpha))$ *is*

$$\alpha^{\star} \equiv \arg\min_{\alpha} \left\{ \sigma(\alpha) \right\} = \arg\min_{\alpha} \left\{ \frac{\sqrt{\alpha^2 + \xi_{\text{max}}^2}}{\alpha + \lambda_{\text{min}}} \right\} = \frac{\xi_{\text{max}}^2}{\lambda_{\text{min}}}$$

and

$$\sigma(\alpha^{\star}) = \frac{\xi_{\text{max}}}{\sqrt{\lambda_{\text{min}}^2 + \xi_{\text{max}}^2}}.$$
 (8)

Proof Simple calculation gives

$$\sigma'(\alpha) = \frac{\alpha \lambda_{\min} - \xi_{\max}^2}{(\alpha + \lambda_{\min})^2 \sqrt{\alpha^2 + \xi_{\max}^2}}.$$

It is obviously that $\sigma'(\alpha) > 0$ for $\alpha > \xi_{\max}^2/\lambda_{\min}$ and $\sigma'(\alpha) < 0$ for $\alpha < \xi_{\max}^2/\lambda_{\min}$. Hence, the upper bound $\sigma(\alpha)$ of the spectral radius $\rho(M(P;\alpha))$ achieves its minimum at $\alpha^* = \xi_{\max}^2/\lambda_{\min}$. Taking α^* into $\sigma(\alpha)$, the minimum value of $\sigma(\alpha)$ given in (8) is obtained.

Remark 1 For case (ii) of Theorem 1, i.e., $\lambda_{\min} < \xi_{\max}$, we see that the α^* given in Corollary 1 satisfies condition (6) since

$$\alpha^{\star} = \frac{\xi_{\text{max}}^2}{\lambda_{\text{min}}} > \frac{\xi_{\text{max}}^2 - \lambda_{\text{min}}^2}{\lambda_{\text{min}}} > \frac{\xi_{\text{max}}^2 - \lambda_{\text{min}}^2}{2\lambda_{\text{min}}}.$$

Remark 2 The parameter α^* in Corollary 1 minimizes only the upper bound $\sigma(\alpha)$ of the spectral radius of the iteration matrix. However, it is still helpful to us to choose an effective parameter α for the NPHSS iteration method. We call α^* the theoretical quasi-optimal parameter of the NPHSS iteration method.

3 Comparison of the NPHSS and PHSS methods

We first introduce a lemma briefly reviewing the convergence analysis of the PHSS method established in Bai et al. (2007a).

Lemma 1 Let the conditions of Theorem 1 be satisfied. Then, the spectral radius $\rho(L(P;\alpha))$ of the PHSS iteration matrix $L(P;\alpha) = (\alpha P + S)^{-1}(\alpha P - H)(\alpha P + H)^{-1}(\alpha P - S)$ satisfies $\rho(L(P;\alpha)) \leq \gamma(\alpha)$, where

$$\gamma(\alpha) = \max_{\lambda_j \in sp(P^{-1}H)} \left| \frac{\alpha - \lambda_j}{\alpha + \lambda_j} \right|.$$



Thus, it holds that

$$\rho(L(P;\alpha)) < \gamma(\alpha) < 1, \quad \forall \alpha > 0.$$

Moreover, the minimum point α_{\star} and the minimum value $\gamma(\alpha_{\star})$ of the upper bound $\gamma(\alpha)$ are, respectively, as

$$\alpha_{\star} \equiv \arg\min_{\alpha} \left\{ \gamma(\alpha) \right\} = \arg\min_{\alpha} \left\{ \max_{\lambda_{j} \in sp(P^{-1}H)} \left| \frac{\alpha - \lambda_{j}}{\alpha + \lambda_{j}} \right| \right\} = \sqrt{\lambda_{\min} \lambda_{\max}}$$

and

$$\gamma(\alpha_{\star}) = \frac{\sqrt{\lambda_{\text{max}}} - \sqrt{\lambda_{\text{min}}}}{\sqrt{\lambda_{\text{max}}} + \sqrt{\lambda_{\text{min}}}}.$$

Note that PHSS method is a two-step iteration. For a fair comparison, the NPHSS iteration scheme (3) can be equivalently rewritten as the following two-step iteration

$$\begin{cases} (\alpha P + H)x^{(k+\frac{1}{2})} = (\alpha P - S)x^{(k)} + b, \\ (\alpha P + H)x^{(k+1)} = (\alpha P - S)x^{(k+\frac{1}{2})} + b, \end{cases}$$
(9)

and the iteration matrix of (9) is $M(P; \alpha)^2$. Then we just need to compare the NPHSS and PHSS methods by analyzing the optimal upper bounds of the spectral radius $\rho(M(P; \alpha)^2)$ and $\rho(L(P; \alpha))$. Using Corollary 1 and Lemma 1, we give the following comparison theorem.

Theorem 2 The respective optimal upper bounds of the spectral radii of NPHSS and PHSS iteration matrices satisfy

$$\sigma^2(\alpha^*) \le \gamma(\alpha_*)$$

if and only if

$$\xi_{\text{max}} \le \sqrt{\frac{\sqrt{\lambda_{\text{max}}} - \sqrt{\lambda_{\text{min}}}}{2\sqrt{\lambda_{\text{min}}}}} \cdot \lambda_{\text{min}}.$$
 (10)

Proof From Corollary 1 and Lemma 1, inequality $\sigma^2(\alpha^*) \leq \gamma(\alpha_*)$ becomes

$$\frac{\xi_{max}^2}{\lambda_{min}^2 + \xi_{max}^2} \leq \frac{\sqrt{\lambda_{max}} - \sqrt{\lambda_{min}}}{\sqrt{\lambda_{max}} + \sqrt{\lambda_{min}}},$$

which is equivalent to

$$\frac{\xi_{\text{max}}^2}{\lambda_{\text{min}}^2} \le \frac{\sqrt{\lambda_{\text{max}}} - \sqrt{\lambda_{\text{min}}}}{2\sqrt{\lambda_{\text{min}}}}.$$
 (11)

Thus, the condition (10) follows by noticing that $\lambda_{max} \geq \lambda_{min} > 0$.

Remark 3 When inequality (10) holds, i.e., the Hermitian part of the coefficient matrix A comparing with its skew-Hermitian part is dominant, we tend to choose NPHSS method rather than PHSS method to solve the linear system (1), and vice verse.



4 Inexact NPHSS iteration method

To further improve computational efficiency of the NPHSS iteration, we develop an inexact NPHSS (INPHSS) iteration, which solves (3) by PCG iterative method. We write the INPHSS iteration scheme in the following algorithm for solving the linear system (1).

Algorithm 3 (*The INPHSS iteration method*) Given an initial guess $x^{(0)} \in \mathbb{C}^n$, then this algorithm leads to the solution of the linear system (1):

$$\ell = 0$$
:

while (not convergent)

$$r^{(\ell)} = b - Ax^{(\ell)}$$
:

approximately solve $(\alpha P + H)z^{(\ell)} = r^{(\ell)}$ by employing PCG method, such that the residual $p^{(\ell)} = r^{(\ell)} - (\alpha P + H)z^{(\ell)}$ of the iteration satisfies $||p^{(\ell)}|| \le \eta_{\ell}||r^{(\ell)}||$;

$$x^{(\ell+1)} = x^{(\ell)} + z^{(\ell)}$$
:

$$\ell = \ell + 1;$$

end

Here, $\{\eta_{\ell}\}$ is prescribed tolerances used to control the accuracies of the inner iterations.

We remark that when P=I, the INPHSS method reduces to the inexact NHSS (INHSS) method.

The convergence properties for the inexact HSS (IHSS) method have been carefully studied in Bai et al. (2003, 2008). Analogously, we can demonstrate the following convergence result about the above INPHSS iteration method.

Theorem 3 Let the conditions of Theorem 1 be satisfied. If $\{x^{(\ell)}\}_{\ell=0}^{\infty} \subseteq \mathbb{C}^n$ is an iteration sequence generated by the INPHSS iteration method and if $x^* \in \mathbb{C}^n$ is the exact solution of the linear system (1), then it holds that

$$\|x^{(\ell+1)} - x^{\star}\|_{2} \le (\sigma(\alpha) + \mu\theta\eta_{\ell})\|x^{(\ell)} - x^{\star}\|_{2}, \quad \ell = 0, 1, 2, \dots$$
 (12)

where the constants μ and θ are given by

$$\mu = \|(\alpha P + H)^{-1}\|_2, \quad \theta = \|A\|_2.$$

In particular, when

$$\sigma(\alpha) + \mu \theta \eta_{\text{max}} < 1, \tag{13}$$

then the iteration sequence $\{x^{(\ell)}\}_{\ell=0}^{\infty}\subseteq\mathbb{C}^n$ converges to $x^{\star}\in\mathbb{C}^n$, where $\eta_{max}=\max_{\ell}\{\eta_{\ell}\}$.

Proof From Algorithm 3, we have

$$x^{(\ell+1)} = x^{(\ell)} + (\alpha P + H)^{-1} (r^{(\ell)} + p^{(\ell)})$$

$$= x^{(\ell)} + (\alpha P + H)^{-1} \left(b - Ax^{(\ell)} + p^{(\ell)} \right)$$

$$= \left(I - (\alpha P + H)^{-1} A \right) x^{(\ell)} + (\alpha P + H)^{-1} b + (\alpha P + H)^{-1} p^{(\ell)}$$

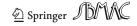
$$= (\alpha P + H)^{-1} (\alpha P - S) x^{(\ell)} + (\alpha P + H)^{-1} b + (\alpha P + H)^{-1} p^{(\ell)}. \tag{14}$$

Because $x^* \in \mathbb{C}^n$ is the exact solution of the linear system (1), it must satisfy

$$x^* = (\alpha P + H)^{-1} (\alpha P - S) x^* + (\alpha P + H)^{-1} b. \tag{15}$$

By subtracting (15) from (14), we have

$$x^{(\ell+1)} - x^* = (\alpha P + H)^{-1} (\alpha P - S)(x^{(\ell)} - x^*) + (\alpha P + H)^{-1} p^{(\ell)}.$$
 (16)



Taking norms on both sides of the identity (16), we can obtain

$$||x^{(\ell+1)} - x^{\star}||_{2} \le ||(\alpha P + H)^{-1}(\alpha P - S)||_{2}||x^{(\ell)} - x^{\star}||_{2} + ||(\alpha P + H)^{-1}||_{2}||p^{(\ell)}||_{2}$$

$$\le \sigma(\alpha)||x^{(\ell)} - x^{\star}||_{2} + \mu\eta\ell||r^{(\ell)}||_{2}.$$
(17)

Noticing that

$$||r^{(\ell)}||_2 = ||b - Ax^{(\ell)}||_2 = ||A(x^* - x^{(\ell)})||_2 \le \theta ||x^{(\ell)} - x^*||_2,$$

by (17) we can obtain (12) and (13).

We remark that Theorem 3 gives the choice of the tolerance $\{\eta_\ell\}$ for convergence. In general, to guarantee the convergence of the INPHSS iteration, the tolerance $\{\eta_\ell\}$ only needs to satisfy the condition (13).

5 Numerical examples

In this section, we are going to examine the efficiency of NHSS and NPHSS iteration methods and their inexact variants for solving linear system (1) by comparing the spectral radii,

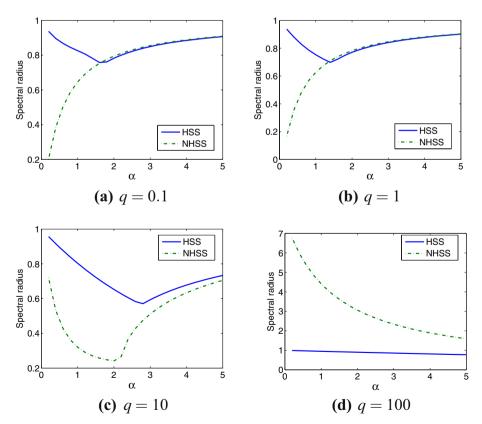


Fig. 1 Centered difference scheme. The spectral radii of the iteration matrices of NHSS and HSS methods versus α



iteration numbers (denoted as IT) and CPU time (in seconds, denoted as CPU) of these two algorithms.

We consider the three-dimensional convection-diffusion equation

$$-(u_{xx} + u_{yy} + u_{zz}) + q(u_x + u_y + u_z) = f(x, y, z)$$
(18)

on the unit cube $\Omega = [0, 1] \times [0, 1] \times [0, 1]$, with constant coefficient q and subjected to Dirichlet-type boundary conditions. Discretizing this equation with seven-point finite difference, and assuming the numbers (m) of grid points in all three directions are the same, we obtain a positive definite system (1) with the coefficient matrix

$$A = T_x \otimes I \otimes I + I \otimes T_y \otimes I + I \otimes I \otimes T_z, \tag{19}$$

where \otimes denotes the Kronecker product, and T_x , T_y , and T_z are tridiagonal matrices given by

$$T_x = \text{tridiag}(t_2, t_1, t_3), \quad T_y = \text{tridiag}(t_2, 0, t_3), \quad \text{and} \quad T_z = \text{tridiag}(t_2, 0, t_3),$$

with

$$t_1 = 6$$
, $t_2 = -1 - r$, $t_3 = -1 + r$

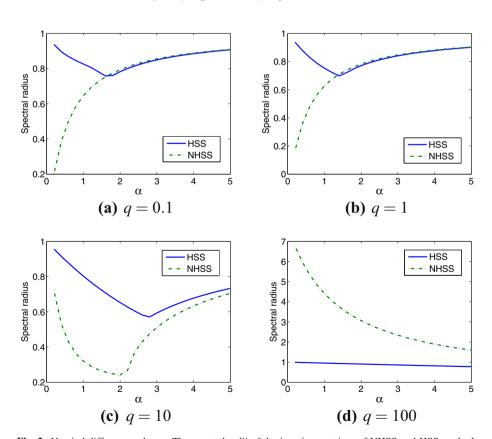


Fig. 2 Upwind difference scheme. The spectral radii of the iteration matrices of NHSS and HSS methods versus α

if the first-order derivatives are approximated by the centered difference scheme, and

$$t_1 = 6 + 6r$$
, $t_2 = -1 - 2r$, $t_3 = -1$

if the first-order derivatives are approximated by the upwind difference scheme. Here r = qh/2 is the mesh Reynolds number with h = 1/(m+1) being the step size. For details, we refer to Bai et al. (2003).

In our implementation, the initial guess is chosen to be $x^{(0)} = \mathbf{0}$ and the iteration is terminated once the current iterate $x^{(\ell)}$ satisfies $||b - Ax^{(\ell)}||_2/||b||_2 \le 10^{-6}$. In addition, we set $n = m^3 = 10^3$, and the right-hand side vector $b = A\mathbf{1}$, with $\mathbf{1}$ being the vector of all entries equal to 1. For convenience, the preconditioner P used in NPHSS and PHSS methods are chosen to be $P = \text{diag}(a_{11}, a_{22}, \ldots, a_{nn})$, where $a_{11}, a_{22}, \ldots, a_{nn}$ are main diagonal elements of the matrix A.

In IHSS, INHSS, IPHSS and INPHSS iteration methods, we set all the tolerances $\{\eta_\ell\}$ = 0.01. We solve the linear systems with coefficient matrices $\alpha P + H$ iteratively by the CG method, and solve the linear systems with the coefficient matrix $\alpha P + S$ iteratively by the GMRES method. We denote the average inner iteration numbers of CG and GMRES as IT_{CG} and IT_{GMRES}, respectively.

The comparisons of spectral radii of different iteration matrices derived by HSS, NHSS, PHSS and NPHSS iteration methods with different damping coefficients q are performed in

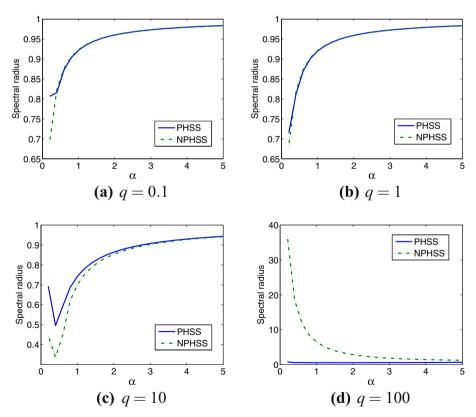


Fig. 3 Centered difference scheme. The spectral radii of the iteration matrices of NPHSS and PHSS methods



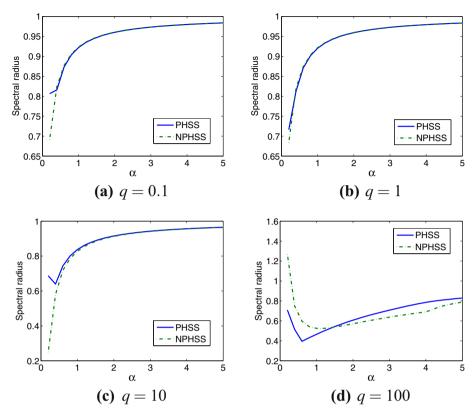


Fig. 4 Upwind difference scheme. The spectral radii of the iteration matrices of NPHSS and PHSS methods versus α

Figs. 1, 2, 3 and 4. From Figs. 1, 2, 3 and 4, we find that when q is small (the Hermitian part of the coefficient matrix is dominant), the spectral radius of the iteration matrix of NHSS method is much smaller than that of HSS method, and the spectral radius of the iteration matrix of NPHSS method is much smaller than that of PHSS method. As q becomes large (the skew-Hermitian part is dominant), HSS and PHSS methods perform better and better.

In Tables 1, 2, 3 and 4, we present iteration numbers and CPU time for HSS, NHSS, PHSS and NPHSS iteration methods with different damping coefficients q. From the results in Tables 1, 2, 3 and 4, we see that when q is small, NHSS and NPHSS methods, no matter compared with the experimental optimal parameter $\alpha_{\rm exp}$ or compared with the theoretical quasi-optimal parameter α^{\star} , perform much better than HSS and PHSS methods both in iteration numbers and in CPU time. As q becomes large, the superiorities of NHSS and NPHSS methods disappear.

In Tables 5 and 6, numerical results for IHSS, INHSS, IPHSS and INPHSS iteration methods are listed; here we adopt the iteration parameters in Tables 1 and 2 for convenience and not the experimental optimal parameters. From Tables 5 and 6, we can obtain the same conclusions as above Tables.

In addition, if Theorem 2 is applied to this numerical example, then we known that when q < 1.606 in the centered difference scheme or q < 1.725 in the upwind difference scheme,



Table 1 Numerical results for HSS and NHSS with the experimental optimal iteration parameters

Method	q	HSS			NHSS	NHSS			
Difference scheme		$\overline{\alpha_{\mathrm{exp}}}$	IT	CPU	$\alpha_{ m exp}$	IT	CPU		
Centered	0.1	1.254	35	0.2650	0.003	2	0.0160		
	1	1.458	39	0.2970	0.003	3	0.0470		
	10	2.186	23	0.1880	1.170	19	0.2030		
	100	3.945	22	0.2810	100	202	2.1570		
Upwind	0.1	1.260	35	0.2660	0.003	2	0.0310		
	1	1.515	39	0.3130	0.003	3	0.0310		
	10	2.601	28	0.2190	1.010	14	0.1560		
	100	16.010	18	0.1410	25	31	0.3280		

Table 2 Numerical results for PHSS and NPHSS with the experimental optimal iteration parameters

Method	q	PHSS			NPHSS	NPHSS			
Difference scheme		$\alpha_{ m exp}$	IT	CPU	$\alpha_{\rm exp}$	IT	CPU		
Centered	0.1	0.209	35	0.2810	0.003	2	0.0310		
	1	0.243	39	0.2970	0.003	3	0.0310		
	10	0.365	23	0.1720	0.201	19	0.2030		
	100	0.658	22	0.2650	16.600	202	2.0780		
Upwind	0.1	0.211	35	0.2650	0.003	2	0.0310		
	1	0.242	39	0.2810	0.003	3	0.0310		
	10	0.301	28	0.2190	0.091	14	0.1410		
	100	0.451	18	0.1250	0.701	31	0.3120		

Table 3 Numerical results for HSS and NHSS with the theoretical quasi-optimal iteration parameters

Method	q	HSS			NHSS			
Difference scheme		$\overline{lpha_{\star}}$	IT	CPU	α^{\star}	IT	CPU	
Centered	0.1	1.690	43	0.3280	0.003	2	0.0160	
	1	1.690	42	0.3130	0.282	10	0.1090	
	10	1.690	28	0.2190	28.175	283	2.9850	
	100	1.690	33	0.6400	0.003	3528	41.0000	
Upwind	0.1	1.698	43	0.3280	0.003	2	0.0320	
	1	1.767	42	0.3280	0.270	9	0.0940	
	10	2.459	29	0.2190	19.370	187	2.0460	
	100	9.374	30	0.2350	508.068	494	5.2500	

the NPHSS is faster than PHSS and the NHSS is faster than HSS. The above numerical results have also examined this fact.

Therefore, we tend to use NHSS or NPHSS iteration methods to solve the linear system (1) when the Hermitian part H is dominant and employ HSS or PHSS if the skew-Hermitian part S is dominant.

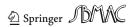


Table 4 Numerical results for PHSS and NPHSS with the theoretical quasi-optimal iteration parameters

Method	q	PHSS			NPHSS	NPHSS			
Difference scheme		α_{\star}	IT	CPU	α*	IT	CPU		
Centered	0.1	0.282	43	0.3280	0.001	2	0.0160		
	1	0.282	42	0.3290	0.047	10	0.1250		
	10	0.282	28	0.2180	4.696	283	3.1720		
	100	0.282	33	0.6400	469.578	3528	39.2340		
Upwind	0.1	0.282	43	0.3280	0.001	2	0.0160		
	1	0.282	42	0.3280	0.043	9	0.0940		
	10	0.282	29	0.2340	1.011	84	0.8750		
	100	0.282	30	0.2340	15.270	494	5.8590		

Table 5 Numerical results for IHSS and INHSS

Method	q	IHSS				INHSS		
Difference scheme		IT _{CG}	IT _{GMRES}	IT	CPU	IT _{CG}	IT	CPU
Centered	0.1	17.7	4.2	35	0.4690	20.0	2	0.0310
	1	18.0	6.7	39	0.5940	20.0	3	0.0620
	10	17.7	10.0	23	0.4220	19.1	19	0.3750
	100	15.4	10.0	22	0.3750	5.4	202	1.3280
Upwind	0.1	17.7	4.1	35	0.4690	20.0	2	0.0310
	1	18.2	6.7	39	0.6090	20.0	3	0.0620
	10	18.1	10.0	28	0.5310	19.4	14	0.2650
	100	15.3	10.0	18	0.3120	15.0	31	0.4840

Table 6 Numerical results for IPHSS and INPHSS

Method	q	IPHSS		INPHSS				
Difference scheme		IT _{CG}	IT _{GMRES}	IT	CPU	IT _{CG}	IT	CPU
Centered	0.1	17.7	4.2	35	0.4530	20.0	3	0.0470
	1	18.0	6.7	39	0.5940	20.0	4	0.0620
	10	17.7	10.0	23	0.4060	18.9	19	0.3900
	100	15.5	10.0	22	0.3590	5.4	202	1.2810
Upwind	0.1	17.5	3.9	35	0.4680	20.0	3	0.0470
	1	18.3	6.7	39	0.5940	20.0	3	0.0630
	10	18.3	10.0	28	0.5150	19.6	14	0.2970
	100	16.3	10.0	18	0.3120	15.3	31	0.5150

6 Conclusions

In this paper, we have developed a non-alternating PHSS (NPHSS) method and its inexact variant for solving non-Hermitian positive definite linear systems. Theoretical analysis demonstrates that for any initial guess and a wide range of parameter α , NPHSS method converges to the unique solution of the linear system (1). We also derive an upper bound $\sigma(\alpha)$ for the spectral radius of NPHSS iteration matrix and give the quasi-optimal parameter α^* which minimizes the upper bound $\sigma(\alpha)$. In addition, both theoretical and numerical results verify that when the Hermitian part of the coefficient matrix is dominant, NPHSS method performs better than HSS and PHSS methods. Hence, our work gives a better choice for solving the linear system (1) when the Hermitian part H of coefficient matrix A is dominant.

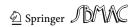
At last, we should mention that the choice of the optimal iteration parameters for the NPHSS method is an interesting but difficult topic. Some related discussions for the HSS method can be seen in Bai et al. (2006); Huang (2014). The investigation of this point may be considered in future study.

Comments: The result of recent paper (Li and Wu 2015) is the special case for P = I of our results here.

Acknowledgments The authors would like to thank the referees for their valuable comments and suggestions which greatly improve the presentation of the paper. This work is supported by the National Natural Science Foundation of China (Grant Nos. 11471150, 11401281, 11271173), and the CAPES and CNPq in Brazil.

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