



Efficient single-step preconditioned HSS iteration methods for complex symmetric linear systems

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ARTICLE INFO

Article history:

Received 16 March 2017

Received in revised form 19 June 2017

Accepted 5 July 2017

Available online 21 July 2017

MSC 2010:

65F10

65F50

Keywords:

Complex symmetric linear system

Positive definite

HSS iteration

Spectral radius

Convergence analysis

ABSTRACT

We propose a single-step preconditioned variant of HSS (SPHSS) and an efficient parameterized SPHSS (PSPHSS) iteration method for solving a class of complex symmetric linear systems. Under suitable conditions, we analyze the convergence properties of the SPHSS and PSPHSS iteration methods. Theoretical analysis shows that the minimal upper bounds for the spectral radius of the SPHSS and PSPHSS iteration matrices are less than those of the SHSS and PSHSS iteration matrices when using the optimal parameters, respectively. Numerical results show that the PSPHSS iteration method has comparable advantage over several other iteration methods whether the experimental optimal parameters are used or not.

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1. Introduction and motivations

We consider to solve the complex symmetric system of linear equations

$$Ax \equiv (W + iT)x = b, \quad (1.1)$$

with $W, T \in \mathbb{R}^{n \times n}$ being symmetric positive semi-definite, and without loss of generality, we suppose W is positive definite. Throughout the paper, the vector $b \in \mathbb{C}^n$ is given, $x \in \mathbb{C}^n$ is unknown, and $i = \sqrt{-1}$ is the imaginary unit. This kind of complex symmetric linear systems can be found in numerical analysis and applied scientific branches, we refer to [1–5] and the references therein.

To solve the problem (1.1) iteratively, a simple Hermitian and skew-Hermitian splitting (HSS) about the coefficient matrix is commonly used. It is given by

$$A = H + S,$$

where

$$H = \frac{1}{2}(A + A^*) = W \quad \text{and} \quad S = \frac{1}{2}(A - A^*) = iT.$$

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It is apparent that A is a non-Hermitian, but positive definite matrix. Here A^* denotes the conjugate transpose of the matrix A . Bai et al. [6] first proposed the HSS iteration method to find the solution of the system (1.1) by the scheme

$$\begin{cases} (\alpha I + W)x^{(k+\frac{1}{2})} = (\alpha I - iT)x^{(k)} + b, \\ (\alpha I + iT)x^{(k+1)} = (\alpha I - W)x^{(k+\frac{1}{2})} + b, \end{cases}$$

where $\alpha > 0$ is a given constant and I is the identity matrix. For more details about the HSS iteration method and its variants, we refer to [7–13].

Later, a modification of the HSS (MHSS) iteration scheme is introduced in [2]. To speed up its convergence rate, Bai et al. [14,15] proposed a preconditioned variant of the MHSS (PMHSS) iteration method and applied it to solve distributed control problems. The PMHSS iteration method is given by

$$\begin{cases} (\alpha V + W)x^{(k+\frac{1}{2})} = (\alpha V - iT)x^{(k)} + b, \\ (\alpha V + T)x^{(k+1)} = (\alpha V + iW)x^{(k+\frac{1}{2})} - ib, \end{cases}$$

where $V \in \mathbb{R}^{n \times n}$ is a prescribed symmetric positive definite matrix. Moreover, Zheng and Ma [16] introduced the accelerated PMHSS iteration methods for complex symmetric linear systems. Li et al. [17] developed a lopsided preconditioned variant of the MHSS iteration method when the real part of the coefficient matrix is dominant. Pour and Goughery [18] proposed a new HSS (NHSS) iteration for non-Hermitian positive-definite linear systems, and a parameterized variant of NHSS iteration method is developed in [19].

Notice that all the iteration schemes mentioned above are two-step methods. Recently, Li and Wu [20] introduced the following single-step HSS (SHSS) iteration method for non-Hermitian positive definite linear systems:

$$(\alpha I + W)x^{(k+1)} = (\alpha I - iT)x^{(k)} + b.$$

The SHSS iteration method can be regarded as a general case of the shift-splitting iteration method, which has been studied for non-Hermitian positive definite linear systems in [21], for saddle point problems in [22,23] and block two-by-two linear systems in [24]. Zeng and Ma [25] proposed a parameterized variant of the SHSS (PSHSS) iteration method, which is more efficient than the SHSS iteration from their numerical experiments. The PSHSS iteration method is given by

$$(\alpha I + \omega W + T)x^{(k+1)} = [\alpha I - i(\omega T - W)]x^{(k)} + (\omega - i)b,$$

where α and ω are given positive constants.

We mention that the HSS-like iteration methods for solving complex systems of linear equations have been extended in a number of literatures. We refer to [21,26–39].

On the other hand, let $x = u + iv$ and $b = p + iq$ with $u, v, p, q \in \mathbb{R}^n$, then the complex linear system (1.1) can be rewritten as 2-by-2 block real equivalent formulation

$$\begin{bmatrix} W & -T \\ T & W \end{bmatrix} \begin{bmatrix} u \\ v \end{bmatrix} = \begin{bmatrix} p \\ q \end{bmatrix}, \quad (1.2)$$

where u, v are unknown real vectors. Chen and Ma [40] introduced the AOR-Uzawa iterative method to solve this real equivalent system. Salkuyeh et al. [41] proposed the following generalized successive overrelaxation (GSOR) iterative method

$$\begin{cases} Wu^{(k+1)} = (1 - \alpha)Wu^{(k)} + \alpha Tv^{(k)} + \alpha p, \\ Wv^{(k+1)} = -\alpha Tu^{(k+1)} + (1 - \alpha)Wv^{(k)} + \alpha q, \end{cases}$$

where $u^{(0)}$ and $v^{(0)}$ are the initial approximations of u and v , respectively. In order to accelerate the convergence rate of the GSOR method, Hezari et al. [42] developed a preconditioned variant of the GSOR (PGSOR) iterative method by the scheme

$$\begin{cases} (\omega W + T)u^{(k+1)} = (1 - \alpha)(\omega W + T)u^{(k)} + \alpha(\omega T - W)v^{(k)} + \alpha(\omega p + q), \\ (\omega W + T)v^{(k+1)} = -\alpha(\omega T - W)u^{(k+1)} + (1 - \alpha)(\omega W + T)v^{(k)} + \alpha(\omega q - p), \end{cases}$$

where α and ω are given positive constants.

In this paper, we want to make improvement on the PMHSS and the SHSS iteration methods.

This paper is organized as follows. In Section 2, two single-step preconditioned HSS iteration methods (SPHSS and PPHSS) are proposed. In Section 3, the convergence properties of the SPHSS iteration method are analyzed, including the convergence conditions, the spectral radius of the iterative matrix and the quasi-optimal parameter. In Section 4, the convergence properties of the PPHSS iteration method are considered. In Section 5, some numerical examples are presented to show the computational efficiencies of the SPHSS and PPHSS iteration methods with comparisons with several other iteration methods. Finally, a brief conclusion is made in Section 6.

2. Single-step preconditioned HSS iteration methods

It has been shown in [20] that the SHSS iteration method is convergent to the unique solution of the linear system with a loose restriction on the parameter α . Notice that the first half step of the HSS iteration method is just the SHSS method, which inspires us to take out the first half step of the PMHSS iteration method to establish the following single-step preconditioned HSS (SPHSS) iteration method.

Algorithm 1. The SPHSS Iteration Method.

Given an initial guess $x^{(0)} \in \mathbb{C}^n$, for $k = 0, 1, 2, \dots$ until the sequence of iterates $\{x^{(k)}\}$ converges, compute $x^{(k+1)}$ by the scheme

$$(\alpha V + W)x^{(k+1)} = (\alpha V - iT)x^{(k)} + b \quad (2.1)$$

where α is a given positive constant and $V \in \mathbb{R}^{n \times n}$ is a prescribed symmetric positive definite matrix.

It is obvious that the SHSS iteration method is a special case of $V = I$.

On the other hand, multiplying a parameter $\omega - i$ by the two sides of the original linear system (1.1) yields that

$$(\omega - i)Ax \equiv [(\omega W + T) + i(\omega T - W)]x = (\omega - i)b. \quad (2.2)$$

Let

$$\tilde{A} = (\omega - i)A, \quad \tilde{W} = \omega W + T, \quad \tilde{T} = \omega T - W, \quad \tilde{b} = (\omega - i)b. \quad (2.3)$$

Then we can rewrite (2.2) as

$$\tilde{A}x \equiv (\tilde{W} + i\tilde{T})x = \tilde{b}.$$

By comparing it with the original linear system (1.1), we can easily obtain the following variant of SPHSS iteration method:

$$(\alpha V + \tilde{W})x^{(k+1)} = (\alpha V - i\tilde{T})x^{(k)} + \tilde{b}. \quad (2.4)$$

Substituting (2.3) into (2.4), we get the parameterized variant of the SPHSS (PSPHSS) iteration method.

Algorithm 2. The PSPHSS Iteration Method.

Given an initial guess $x^{(0)} \in \mathbb{C}^n$, for $k = 0, 1, 2, \dots$ until the sequence of iterates $\{x^{(k)}\}$ converges, compute $x^{(k+1)}$ by the scheme

$$(\alpha V + \omega W + T)x^{(k+1)} = [\alpha V - i(\omega T - W)]x^{(k)} + (\omega - i)b, \quad (2.5)$$

where α, ω are given positive constants, and $V \in \mathbb{R}^{n \times n}$ is a prescribed symmetric positive definite matrix.

It is easy to see that the PSHSS iteration method is a special case of $V = I$.

3. Convergence analysis about the SPHSS iteration method

First, we rewrite the SPHSS iteration scheme (2.1) as

$$x^{(k+1)} = M^V(\alpha)x^{(k)} + N^V(\alpha)b, \quad k = 0, 1, 2, \dots,$$

where

$$\begin{aligned} M^V(\alpha) &= (\alpha V + W)^{-1}(\alpha V - iT) = (\alpha I + V^{-1}W)^{-1}(\alpha I - iV^{-1}T), \\ N^V(\alpha) &= (\alpha V + W)^{-1}, \end{aligned} \quad (3.1)$$

with $V \in \mathbb{R}^{n \times n}$ being a prescribed symmetric positive definite matrix.

If we define

$$F^V(\alpha) = \alpha V + W, \quad \text{and} \quad G^V(\alpha) = \alpha V - iT,$$

then it holds that

$$A = F^V(\alpha) - G^V(\alpha), \quad \text{and} \quad M^V(\alpha) = F^V(\alpha)^{-1}G^V(\alpha).$$

Hence, the splitting matrix $F^V(\alpha)$ can be taken as a preconditioner for the complex symmetric matrix $A \in \mathbb{C}^{n \times n}$, which is referred as the SPHSS preconditioner.

In particular, if we take $V = W$, then

$$M^W(\alpha) = \frac{1}{\alpha + 1}(\alpha I - iW^{-1}T), \quad N^W(\alpha) = \frac{1}{\alpha + 1}W^{-1},$$

and the SPHSS iteration scheme is now induced by the matrix splitting

$$A = F^W(\alpha) - G^W(\alpha)$$

with

$$F^W(\alpha) = (\alpha + 1)W, \quad \text{and} \quad G^W(\alpha) = \alpha W - iT.$$

In order to explore the convergence properties of the SPHSS iteration method, we use the following notations

$$\begin{aligned} \tilde{\lambda}_{\min}^V &= \min_{\tilde{\lambda}_j \in sp(V^{-\frac{1}{2}} W V^{-\frac{1}{2}})} \{\tilde{\lambda}_j\}, & \tilde{\sigma}_{\max}^V &= \max_{\tilde{\sigma}_j \in sp(V^{-\frac{1}{2}} T V^{-\frac{1}{2}})} \{\tilde{\sigma}_j\}, \\ \tilde{\lambda}_{\max}^V &= \max_{\tilde{\lambda}_j \in sp(V^{-\frac{1}{2}} W V^{-\frac{1}{2}})} \{\tilde{\lambda}_j\}, & \tilde{\sigma}_{\min}^V &= \min_{\tilde{\sigma}_j \in sp(V^{-\frac{1}{2}} T V^{-\frac{1}{2}})} \{\tilde{\sigma}_j\}, \end{aligned} \quad (3.2)$$

where $sp(X)$ is the spectral set of the matrix X .

Theorem 1. If $W, T \in \mathbb{R}^{n \times n}$ are symmetric positive semi-definite matrices with at least one of them being positive definite, then the spectral radius $\rho(M^V(\alpha))$ is bounded by

$$\delta^V(\alpha) = \frac{\sqrt{\alpha^2 + (\tilde{\sigma}_{\max}^V)^2}}{\alpha + \tilde{\lambda}_{\min}^V}. \quad (3.3)$$

Moreover,

- (i) if $\tilde{\lambda}_{\min}^V \geq \tilde{\sigma}_{\max}^V$, then $\delta^V(\alpha) < 1$ i.e. the iteration converges for any $\alpha > 0$;
- (ii) if $\tilde{\lambda}_{\min}^V < \tilde{\sigma}_{\max}^V$, then $\delta^V(\alpha) < 1$ if and only if

$$\alpha > \frac{(\tilde{\sigma}_{\max}^V)^2 - (\tilde{\lambda}_{\min}^V)^2}{2\tilde{\lambda}_{\min}^V}. \quad (3.4)$$

Proof. According to (3.1), we obtain

$$\begin{aligned} \rho(M^V(\alpha)) &= \rho((\alpha I + V^{-1}W)^{-1}(\alpha I - iV^{-1}T)) \\ &\leq \|(\alpha I + V^{-1}W)^{-1}(\alpha I - iV^{-1}T)\|_2 \\ &\leq \|(\alpha I + V^{-1}W)^{-1}\|_2 \|(\alpha I - iV^{-1}T)\|_2 \\ &= \left(\max_{\tilde{\lambda}_j \in sp(V^{-\frac{1}{2}} W V^{-\frac{1}{2}})} \frac{1}{\alpha + \tilde{\lambda}_j} \right) \left(\max_{\tilde{\sigma}_j \in sp(V^{-\frac{1}{2}} T V^{-\frac{1}{2}})} \sqrt{\alpha^2 + \tilde{\sigma}_j^2} \right) \\ &= \frac{\sqrt{\alpha^2 + (\tilde{\sigma}_{\max}^V)^2}}{\alpha + \tilde{\lambda}_{\min}^V}. \end{aligned}$$

Therefore, $\tilde{\lambda}_{\min}^V \geq \tilde{\sigma}_{\max}^V$ definitely leads to $\delta^V(\alpha) < 1$, and if $\tilde{\lambda}_{\min}^V < \tilde{\sigma}_{\max}^V$, then $\delta^V(\alpha) < 1$ if and only if

$$\alpha^2 + (\tilde{\sigma}_{\max}^V)^2 < (\alpha + \tilde{\lambda}_{\min}^V)^2,$$

which is equivalent to (3.4). \square

Corollary 1. Under the conditions of Theorem 1, the optimal parameter α_*^V which minimizes the upper bound $\delta^V(\alpha)$ of the spectral radius $\rho(M^V(\alpha))$ is given by

$$\alpha_*^V = \operatorname{argmin}_{\alpha} \left\{ \frac{\sqrt{\alpha^2 + (\tilde{\sigma}_{\max}^V)^2}}{\alpha + \tilde{\lambda}_{\min}^V} \right\} = \frac{(\tilde{\sigma}_{\max}^V)^2}{\tilde{\lambda}_{\min}^V}, \quad (3.5)$$

and

$$\delta^V(\alpha_*^V) = \frac{\tilde{\sigma}_{\max}^V}{\sqrt{(\tilde{\lambda}_{\min}^V)^2 + (\tilde{\sigma}_{\max}^V)^2}}. \quad (3.6)$$

Proof. By simple calculations, we get

$$\delta^V(\alpha)' = \frac{\alpha \tilde{\lambda}_{\min}^V - (\tilde{\sigma}_{\max}^V)^2}{(\alpha + \tilde{\lambda}_{\min}^V)^2 \sqrt{\alpha^2 + (\tilde{\sigma}_{\max}^V)^2}}.$$

Therefore, the upper bound $\delta^V(\alpha)$ of the spectral radius $\rho(M^V(\alpha))$ achieves its minimum at $\alpha_*^V = (\tilde{\sigma}_{\max}^V)^2 / \tilde{\lambda}_{\min}^V$, i.e., (3.5) holds true. Taking α_*^V into $\delta^V(\alpha)$, we obtain the minimum value of $\delta^V(\alpha)$ which is given by (3.6). \square

Moreover, according to (3.6), when $\tilde{\lambda}_{\min}^V \geq \tilde{\sigma}_{\max}^V$, if we take the optimal parameter α_*^V , then it holds that

$$\rho(M^V(\alpha_*^V)) \leq \delta^V(\alpha_*^V) \leq \frac{1}{\sqrt{2}}.$$

In particular, when $V = W$, the SPHSS iteration is given by

$$(\alpha + 1)Wx^{(k+1)} = (\alpha W - iT)x^{(k)} + b,$$

and we have $\tilde{\lambda}_{\min}^W = 1$, thus the right hand sides of (3.3)–(3.6) are related only with $\tilde{\sigma}_{\max}^W$, and we have

$$\delta^W(\alpha_*^W) = \frac{\tilde{\sigma}_{\max}^W}{\sqrt{1 + (\tilde{\sigma}_{\max}^W)^2}}. \quad (3.7)$$

Remark 1. When $V = I$, the SPHSS iteration is reduced to the SHSS iteration, and the minimal upper bound $\delta^I(\alpha)$ of the spectral radius $\rho(M^I(\alpha))$ is given by

$$\delta^I(\alpha_*^I) = \frac{\tilde{\sigma}_{\max}^I}{\sqrt{(\tilde{\lambda}_{\min}^I)^2 + (\tilde{\sigma}_{\max}^I)^2}} = \frac{\frac{\tilde{\sigma}_{\max}^I}{\tilde{\lambda}_{\min}^I}}{\sqrt{1 + \left(\frac{\tilde{\sigma}_{\max}^I}{\tilde{\lambda}_{\min}^I}\right)^2}}. \quad (3.8)$$

From (3.7) and (3.8), using the fact that $\|W^{-1}T\|_2 \leq \|W^{-1}\|_2 \|T\|_2$, we obtain

$$\delta^W(\alpha_*^W) \leq \delta^I(\alpha_*^I),$$

since $\tilde{\sigma}_{\max}^W \leq \frac{\tilde{\sigma}_{\max}^I}{\tilde{\lambda}_{\min}^I}$ and $f(x) \equiv \frac{x}{\sqrt{1+x^2}}$ is a strictly monotonic increasing function. Therefore, the SPHSS iteration using $V = W$ is more efficient than the SHSS iteration when the optimal parameters are used.

Remark 2. According to Theorem 1, the SPHSS iteration would always converge if the parameter α is properly chosen. Moreover, from (3.6), theoretically the optimal symmetric positive definite matrix V_* minimizing $\delta^V(\alpha_*^V)$ is given by

$$V_* = \operatorname{argmin}_V \left\{ \frac{\tilde{\sigma}_{\max}^V}{\sqrt{(\tilde{\lambda}_{\min}^V)^2 + (\tilde{\sigma}_{\max}^V)^2}} \right\} = \operatorname{argmin}_V \left\{ \frac{\tilde{\sigma}_{\max}^V}{\tilde{\lambda}_{\min}^V} \right\}.$$

While practically we usually take the positive definite matrix between W and T to substitute the matrix V . If both W and T are positive definite, then the better matrix V can be chosen such that

$$\frac{\tilde{\sigma}_{\max}^V}{\tilde{\lambda}_{\min}^V} = \min \left\{ \frac{\tilde{\sigma}_{\max}^W}{\tilde{\lambda}_{\min}^W}, \frac{\tilde{\sigma}_{\max}^T}{\tilde{\lambda}_{\min}^T} \right\} = \min \left\{ \tilde{\sigma}_{\max}^W, \frac{1}{\tilde{\lambda}_{\min}^T} \right\}.$$

We mention that the SPHSS iteration method has also been studied in [43] for solving a general class of non-Hermitian positive definite linear systems, and similar results about the convergence conditions and the quasi-optimal parameters are given there. However, we provide new analysis of the new SPHSS iteration method in Remarks 1 and 2.

4. Convergence analysis about the PSPHSS iteration method

In this section, we consider the convergence properties of the PSPHSS iteration method and search for the optimal parameters α_*^V and ω_*^V . First, we rewrite the PSPHSS iteration scheme (2.5) as

$$x^{(k+1)} = M^V(\alpha, \omega)x^{(k)} + N^V(\alpha, \omega)(\omega - i)b, \quad k = 0, 1, 2, \dots,$$

where

$$\begin{aligned} M^V(\alpha, \omega) &= (\alpha V + \omega W + T)^{-1}[\alpha V - i(\omega T - W)] \\ &= [\alpha I + V^{-1}(\omega W + T)]^{-1}[\alpha I - iV^{-1}(\omega T - W)] \end{aligned}$$

and

$$N^V(\alpha, \omega) = (\alpha V + \omega W + T)^{-1},$$

with α, ω being given positive constants.

As before, if we define

$$F^V(\alpha, \omega) = \alpha V + \omega W + T, \quad \text{and} \quad G^V(\alpha, \omega) = \alpha V - i(\omega T - W),$$

then it holds that

$$(\omega - i)A = F^V(\alpha, \omega) - G^V(\alpha, \omega), \quad \text{and} \quad M^V(\alpha, \omega) = F^V(\alpha, \omega)^{-1}G^V(\alpha, \omega).$$

Thus, the splitting matrix $F^V(\alpha, \omega)$ can be taken as a preconditioner for the complex symmetric matrix $(\omega - i)A \in \mathbb{C}^{n \times n}$, which is referred as the PSPHSS preconditioner.

In particular, if we take $V = W$, then the SPHSS iteration scheme is now induced by the matrix splitting

$$(\omega - i)A = F^W(\alpha, \omega) - G^W(\alpha, \omega)$$

with

$$F^W(\alpha, \omega) = (\alpha + \omega)W + T, \quad \text{and} \quad G^W(\alpha, \omega) = \alpha W - i(\omega T - W).$$

For convenience, we define

$$\hat{\lambda}_{\min}^V(\omega) = \min_{\hat{\lambda}_j \in \text{sp}(V^{-\frac{1}{2}}(\omega W + T)V^{-\frac{1}{2}})} \{\hat{\lambda}_j\} \quad \text{and} \quad \hat{\sigma}_{\max}^V(\omega) = \max_{\hat{\sigma}_j \in \text{sp}(V^{-\frac{1}{2}}(\omega T - W)V^{-\frac{1}{2}})} \{|\hat{\sigma}_j|\}. \quad (4.1)$$

Since $\tilde{W} = \omega W + T$ is also a symmetric positive definite matrix, then according to the convergence analysis about the SPHSS iteration method given in Section 3, we immediately obtain the following results.

Theorem 2. If $W, T \in \mathbb{R}^{n \times n}$ are symmetric positive semi-definite matrices with at least one of them being positive definite, then the spectral radius $\rho(M^V(\alpha, \omega))$ is bounded by

$$\delta^V(\alpha, \omega) = \frac{\sqrt{\alpha^2 + (\hat{\sigma}_{\max}^V(\omega))^2}}{\alpha + \hat{\lambda}_{\min}^V(\omega)}. \quad (4.2)$$

Furthermore, it is easy to see that

- (i) if $\hat{\lambda}_{\min}^V(\omega) \geq \hat{\sigma}_{\max}^V(\omega)$, then $\delta^V(\alpha, \omega) < 1$ i.e. the iteration converges for any $\alpha > 0$;
- (ii) if $\hat{\lambda}_{\min}^V(\omega) < \hat{\sigma}_{\max}^V(\omega)$, then $\delta^V(\alpha, \omega) < 1$ if and only if

$$\alpha > \frac{(\hat{\sigma}_{\max}^V(\omega))^2 - (\hat{\lambda}_{\min}^V(\omega))^2}{2\hat{\lambda}_{\min}^V(\omega)}. \quad (4.3)$$

Corollary 2. Under the conditions of Theorem 2, the optimal parameter $\alpha_*^V(\omega)$ minimizing the upper bound $\delta^V(\alpha, \omega)$ of the spectral radius $\rho(M^V(\alpha, \omega))$ is

$$\alpha_*^V(\omega) = \operatorname{argmin}_{\alpha} \left\{ \frac{\sqrt{\alpha^2 + (\hat{\sigma}_{\max}^V(\omega))^2}}{\alpha + \hat{\lambda}_{\min}^V(\omega)} \right\} = \frac{(\hat{\sigma}_{\max}^V(\omega))^2}{\hat{\lambda}_{\min}^V(\omega)}, \quad (4.4)$$

and

$$\delta^V(\alpha_*^V(\omega), \omega) = \frac{\hat{\sigma}_{\max}^V(\omega)}{\sqrt{(\hat{\lambda}_{\min}^V(\omega))^2 + (\hat{\sigma}_{\max}^V(\omega))^2}}. \quad (4.5)$$

According to Theorem 2, the PSPHSS iteration method converges, for any $\omega > 0$ by choosing a proper parameter α satisfying (4.3), or for any $\alpha > 0$ if the parameter ω satisfies $\hat{\lambda}_{\min}^V(\omega) \geq \hat{\sigma}_{\max}^V(\omega)$ which leads to $\delta^V(\alpha_*^V(\omega), \omega) \leq \frac{1}{\sqrt{2}}$. In addition, we have the following results by similar analysis as in [25].

Theorem 3. Let $\hat{\lambda}_{\min}^V(\omega), \hat{\sigma}_{\max}^V(\omega)$ be defined as in (4.1), $\tilde{\lambda}_{\min}^V, \tilde{\sigma}_{\max}^V, \tilde{\lambda}_{\max}^V$ and $\tilde{\sigma}_{\min}^V$ be defined as in (3.2). Denote

$$\bar{\omega}^V = \frac{\tilde{\lambda}_{\min}^V + \tilde{\lambda}_{\max}^V}{\tilde{\sigma}_{\min}^V + \tilde{\sigma}_{\max}^V}, \quad \omega_1^V = \frac{\tilde{\lambda}_{\min}^V + \tilde{\sigma}_{\min}^V}{\tilde{\sigma}_{\max}^V - \tilde{\lambda}_{\min}^V}, \quad \omega_2^V = \frac{\tilde{\lambda}_{\max}^V - \tilde{\sigma}_{\min}^V}{\tilde{\sigma}_{\min}^V + \tilde{\lambda}_{\min}^V}.$$

Then $\hat{\lambda}_{\min}^V(\omega) \geq \hat{\sigma}_{\max}^V(\omega)$, i.e. the PSPHSS iteration method converges for any $\alpha > 0$ if any of the following conditions hold:

- (1) $\tilde{\lambda}_{\min}^V \geq \tilde{\sigma}_{\max}^V$ and $\omega \geq \bar{\omega}^V$;
- (2) $\tilde{\lambda}_{\min}^V < \tilde{\sigma}_{\max}^V$ and $\bar{\omega}^V \leq \omega \leq \omega_1^V$;
- (3) $\omega_2^V \leq \omega < \bar{\omega}^V$.

Moreover, we can obtain the quasi-optimal parameter ω_*^V by similar analysis as in [25].

Theorem 4. Let $\tilde{\lambda}_{\min}^V, \tilde{\sigma}_{\max}^V, \tilde{\lambda}_{\max}^V$ and $\tilde{\sigma}_{\min}^V$ be defined as in (3.2). Then the quasi-optimal parameter ω_*^V for the PSPHSS iteration method is given by

$$\omega_*^V = \bar{\omega}^V = \frac{\tilde{\lambda}_{\min}^V + \tilde{\lambda}_{\max}^V}{\tilde{\sigma}_{\min}^V + \tilde{\sigma}_{\max}^V}.$$

Remark 3. When $V = I$, the PSPHSS iteration is reduced to the PSHSS iteration. According to [Corollary 2](#), by similar analysis as in [Remark 1](#), for comparing the PSPHSS iteration using $V = \tilde{W} = \omega W + T$ and the PSHSS iteration, we have

$$\delta^{\tilde{W}}(\alpha_*^{\tilde{W}}(\omega), \omega) \leq \delta^I(\alpha_*^I(\omega), \omega).$$

Therefore, the PSPHSS iteration using $V = \omega W + T$ is more efficient than the PSHSS iteration when the optimal parameters are used.

However, for simplicity, we consider the special case of $V = W$, and the PSPHSS iteration scheme (2.5) becomes

$$[(\alpha + \omega)W + T]x^{(k+1)} = [\alpha W - i(\omega T - W)]x^k + (\omega - i)b,$$

which can be rewritten as

$$x^{(k+1)} = M^W(\alpha, \omega)x^k + N^W(\alpha, \omega)(\omega - i)b,$$

where

$$\begin{aligned} M^W(\alpha, \omega) &= [(\alpha + \omega)W + T]^{-1}[\alpha W - i(\omega T - W)] \\ &= [(\alpha + \omega)I + W^{-1}T]^{-1}[\alpha I - i(\omega W^{-1}T - I)], \\ N^W(\alpha, \omega) &= [(\alpha + \omega)W + T]^{-1}. \end{aligned} \quad (4.6)$$

In this case, we have the following results.

Theorem 5. If we take $V = W$ for the PSPHSS iteration method, then the spectral radius $\rho(M^W(\alpha, \omega))$ is bounded by

$$\delta^W(\alpha, \omega) = \frac{\sqrt{\alpha^2 + \omega^2(\tilde{\sigma}_{\max}^W)^2 + 1}}{\alpha + \omega + \tilde{\sigma}_{\min}^W}. \quad (4.7)$$

Moreover, the PSPHSS iteration using $V = W$ converges if any of the following conditions hold:

- (1) $\omega = 1$ for any $\alpha > 0$;
- (2) $\omega \neq 1$ and

$$\alpha > \frac{\omega^2(\tilde{\sigma}_{\max}^W)^2 + 1 - (\omega + \tilde{\sigma}_{\min}^W)^2}{2(\omega + \tilde{\sigma}_{\min}^W)}, \quad (4.8)$$

in particular, $\omega \neq 1$ and $\alpha > \frac{\omega^2(\tilde{\sigma}_{\max}^W)^2 + 1 - \omega^2}{2\omega}$.

Proof. According to (4.6), it is easy to see that

$$\rho(M^W(\alpha, \omega)) \leq \max_{\tilde{\sigma}_j \in sp(W^{-\frac{1}{2}}TW^{-\frac{1}{2}})} \frac{\sqrt{\alpha^2 + (\omega\tilde{\sigma}_j - 1)^2}}{\alpha + \omega + \tilde{\sigma}_j} \leq \max_{\tilde{\sigma}_j \in sp(W^{-\frac{1}{2}}TW^{-\frac{1}{2}})} \frac{\sqrt{\alpha^2 + \omega^2\tilde{\sigma}_j^2 + 1}}{\alpha + \omega + \tilde{\sigma}_j} \leq \delta^W(\alpha, \omega). \quad (4.9)$$

Furthermore, when $\omega = 1$ for any $\alpha > 0$, we have

$$\frac{\sqrt{\alpha^2 + \omega^2\tilde{\sigma}_j^2 + 1}}{\alpha + \omega + \tilde{\sigma}_j} < 1$$

holds for any $\tilde{\sigma}_j \in sp(W^{-1}T)$, which leads to $\rho(M^W(\alpha, \omega)) < 1$. Meanwhile, when $\omega \neq 1$, the inequality (4.8) is equivalent to

$$\alpha^2 + \omega^2(\tilde{\sigma}_{\max}^W)^2 + 1 < (\alpha + \omega + \tilde{\sigma}_{\min}^W)^2,$$

which means that $\delta^W(\alpha, \omega) < 1$. \square

Similar as [Corollary 1](#), we have the following results.

Corollary 3. Under the conditions of [Theorem 5](#), the quasi-optimal parameter $\alpha_*^W(\omega)$ minimizing the upper bound $\delta^W(\alpha, \omega)$ of the spectral radius $\rho(M^W(\alpha, \omega))$ is

$$\alpha_*^W(\omega) = \operatorname{argmin}_{\alpha} \left\{ \frac{\sqrt{\alpha^2 + \omega^2(\tilde{\sigma}_{\max}^W)^2 + 1}}{\alpha + \omega + \tilde{\sigma}_{\min}^W} \right\} = \frac{\omega^2(\tilde{\sigma}_{\max}^W)^2 + 1}{\omega + \tilde{\sigma}_{\min}^W}, \quad (4.10)$$

and

$$\delta^W(\alpha_*^W(\omega), \omega) = \frac{\sqrt{\omega^2(\tilde{\sigma}_{\max}^W)^2 + 1}}{\sqrt{\omega^2(\tilde{\sigma}_{\max}^W)^2 + 1 + (\omega + \tilde{\sigma}_{\min}^W)^2}}. \quad (4.11)$$

Note that if the parameter ω is chosen such that $(\omega + \tilde{\sigma}_{\min}^W)^2 \geq \omega^2(\tilde{\sigma}_{\max}^W)^2 + 1$, then we obtain $\delta^W(\alpha_*^W(\omega), \omega) \leq \frac{1}{\sqrt{2}}$. On the other hand, according to the definitions in (3.2) and (4.1), we have

$$\tilde{\lambda}_{\min}^W = \tilde{\lambda}_{\max}^W = 1, \quad \hat{\lambda}_{\min}^W(\omega) = \omega + \tilde{\sigma}_{\min}^W,$$

and

$$\hat{\sigma}_{\max}^W(\omega) = \max_{\tilde{\sigma}_j \in \text{sp}(W^{-\frac{1}{2}}TW^{-\frac{1}{2}})} \{|\omega\tilde{\sigma}_j - 1|\} \leq \sqrt{\omega^2(\tilde{\sigma}_{\max}^W)^2 + 1}.$$

Thus, the results of Theorems 3–4 can be simplified.

5. Numerical results

In this section, we use some numerical examples to test the computational efficiencies of the SPHSS and PSPHSS iteration methods for solving the complex symmetric linear system (1.1). Numerical comparisons with the PMHSS [14], the SHSS [20], the PSHSS [25], and the PGSOR iteration methods are given to show the advantage of the SPHSS and PSPHSS iteration methods. We show the results for these methods by computations with the preconditioned GMRES methods.

We compare these methods mainly from the point of view of the number of iterations (denoted as IT) and the CPU times (denoted as CPU). Our experiments are carried out in MATLAB R2013a on Intel(R) Core(TM) CPU 1.8 × 2 Ghz and 4.00 GB of RAM. The CPU time is recorded by the command “tic-toc”.

In our computations, the initial approximation $x^{(0)} = u^{(0)} + iv^{(0)}$ is always chosen to be zero vector and the stopping criteria for all the methods are

$$\frac{\|b - Ax^{(k)}\|_2}{\|b\|_2} \leq 10^{-6},$$

where $x^{(k)} = u^{(k)} + iv^{(k)}$ is the current approximation.

Example 1. (See [2,9]) Consider a system of linear equations of the form $(W + iT)x = b$. The matrix $T \in \mathbb{R}^{n \times n}$ possesses the tensor-product form

$$T = I \otimes B_m + B_m \otimes I \quad \text{with} \quad B_m = \text{tridiag}(-1, 2, -1) \in \mathbb{R}^{m \times m},$$

so T is an $n \times n$ block-tridiagonal matrix with $n = m^2$. The matrix $W \in \mathbb{R}^{n \times n}$ is defined by

$$W = 10(I \otimes B_c + B_c \otimes I) + 9(e_1 e_m^T + e_m e_1^T) \otimes I \quad \text{with} \quad B_c = B_m - e_1 e_m^T - e_m e_1^T \in \mathbb{R}^{m \times m},$$

where e_1 and e_m are the first and m th unit basis vectors in \mathbb{R}^m , respectively. Here W is the five-point centered difference matrix which approximates the negative Laplacian operator with periodic boundary conditions, on a uniform mesh in the square $[0, 1] \times [0, 1]$ with the mesh-size $h = 1/(m + 1)$. We choose the vector $b = (1 + i)A\mathbf{1}$, where $\mathbf{1}$ is a special vector of all entries equal to 1.

Example 2. In direct frequency domain analysis of an n -degree-of-freedom linear system, researchers often need to solve the following system of linear equations

$$[(K - \varpi^2 M) + i(C_H + \varpi C_V)]x = b,$$

where K is the (real symmetric) stiffness matrix, M is the inertia (mass) matrix, which is typically real symmetric, C_V is the (real diagonal) viscous damping matrix and C_H is the (real symmetric) hysteretic damping matrix. If $C_H = \beta K$, then β is a damping coefficient, and ϖ is the driving circular frequency. For more details, we refer to [4,44].

Similar as the numerical examples given in [2,41,42], we consider to solve the following linear system as well

$$\left[(K - \varpi^2 I) + i(\beta K + 10\varpi I) \right] x = b,$$

where ϖ, β are given real constants. The matrix $K \in \mathbb{R}^{n \times n}$ possesses the tensor-product form

$$K = I \otimes B_m + B_m \otimes I \quad \text{with} \quad B_m = \frac{1}{h^2} \text{tridiag}(-1, 2, -1) \in \mathbb{R}^{m \times m}.$$

In our tests, the right-hand side vector b with its j th entry b_j is given by

$$b_j = \frac{(1 + i)j}{h^2(j + 1)^2}, \quad j = 1, 2, \dots, n.$$

In addition, we normalize the system by multiplying both sides by h^2 .

Table 1The numerical results of [Example 1](#) for different iteration methods.

Method		Grid			
		16 × 16	32 × 32	48 × 48	64 × 64
PMHSS ($V = W$)	α_*	0.7	0.7	0.7	0.7
	IT	3	3	3	3
	CPU(s)	0.0708	0.1157	0.2439	0.4796
SHSS	α_*	0.01	0.01	0.01	0.005
	IT	3	4	4	4
	CPU(s)	0.0691	0.1116	0.2391	0.4776
PSHSS	α_*	0.01	0.01	0.01	0.01
	ω_*	10	10	10	10
	IT	3	3	3	3
SPHSS ($V = W$)	CPU(s)	0.0699	0.1027	0.2169	0.4462
	α_*	0.01	0.01	0.01	0.01
	IT	3	3	3	3
PSPHSS ($V = W$)	CPU(s)	0.0665	0.0973	0.1724	0.3151
	α_*	0.01	0.01	0.01	0.01
	ω_*	10	10	10	10
PGSOR	IT	3	3	3	3
	CPU(s)	0.0671	0.0979	0.1738	0.3178
	α_*	0.99	0.99	0.99	0.99
	ω_*	10	10	10	10
	IT	2	2	2	2
	CPU(s)	0.0704	0.1541	0.3605	0.8958

Table 2Numerical results of [Example 2](#) for different iteration methods with $(\varpi, \beta) = (\pi/4, 0.02)$.

Method		Grid			
		16 × 16	32 × 32	48 × 48	64 × 64
PMHSS ($V = W$)	α_*	0.9	0.9	0.9	0.9
	IT	5	5	5	5
	CPU(s)	0.0827	0.1417	0.2814	0.5141
SHSS	α_*	0.0005	0.0005	0.0005	0.0005
	IT	5	5	5	5
	CPU(s)	0.0621	0.0865	0.1705	0.3554
PSHSS	α_*	0.005	0.005	0.005	0.005
	ω_*	5	5	5	5
	IT	5	5	5	5
SPHSS ($V = W$)	CPU(s)	0.0624	0.0863	0.1701	0.3555
	α_*	0.01	0.01	0.01	0.01
	IT	5	5	5	5
PSPHSS ($V = W$)	CPU(s)	0.0613	0.0809	0.1328	0.2597
	α_*	0.01	0.01	0.01	0.01
	ω_*	5	5	5	5
PGSOR	IT	5	5	5	5
	CPU(s)	0.0615	0.0816	0.1346	0.2647
	α_*	0.99	0.99	0.99	0.99
	ω_*	5	5	5	5
	IT	4	4	4	4
	CPU(s)	0.0801	0.2103	0.5607	1.3715

[Tables 1–2](#) show the numerical results of various iteration methods with respect to different problem sizes for [Examples 1–2](#) by using the experimental optimal parameters α_* and ω_* , respectively. The parameters $\varpi = \pi/4$ and $\beta = 0.02$ are considered in [Example 2](#). The results about iterations and CPU times are listed in the tables. It can be seen that the experimental optimal parameters of the considered methods are very stable, which brings convenience for us to find the optimal parameters for solving large complex symmetric linear systems. We can see that the SPHSS method has a little advantage over the SHSS method from the point of view of iterations and CPU times by using the experimental optimal parameters. Similarly, the PSPHSS iteration method is a little more efficient than the PSHSS iteration method by using the experimental optimal parameters from the point of view of CPU times, because it is a combination of the SPHSS and PSHSS methods, and it inherits the merits of both.

Note that if the computations are carried out without using the preconditioned GMRES methods, the SHSS method is much less efficient than the PSHSS iteration method, see [\[25\]](#). Similarly, the SPHSS method is less efficient than the PSPHSS iteration method. Therefore, according to the results of [Tables 1–2](#), we can conclude that the PSHSS and the PSHSS iteration methods outperform the other methods when the experimental optimal parameters are used.

In order to make further comparisons among the PSPHSS, the PSHSS, and the PGSOR methods, [Tables 3–4](#) show the numerical results of these three methods with respect to different problem sizes for [Example 2](#) with and without using the

Table 3Numerical results of Example 2 for the PSPHSS ($V = W$), the PSHSS and the PGSOR iteration method with $(\varpi, \beta) = (0.2, 0.5)$.

Method		Grid			
		16 × 16	32 × 32	48 × 48	64 × 64
PSPHSS	IT	4	4	4	4
$(\alpha_*, \omega_*) = (0.001, 2)$	CPU(s)	0.0595	0.0742	0.1273	0.2297
PSHSS	IT	4	4	4	4
$(\alpha_*, \omega_*) = (0.001, 2)$	CPU(s)	0.0609	0.0811	0.1566	0.3334
PGSOR	IT	3	3	3	3
$(\alpha_*, \omega_*) = (0.99, 2)$	CPU(s)	0.0707	0.1562	0.3856	0.9980
PSPHSS	IT	4	4	4	4
$(\alpha, \omega) = (0.5, 10)$	CPU(s)	0.0601	0.0745	0.1282	0.2318
PSHSS	IT	6	9	12	15
$(\alpha, \omega) = (0.5, 10)$	CPU(s)	0.0630	0.0947	0.2094	0.4626
PGSOR	IT	6	6	6	6
$(\alpha, \omega) = (0.5, 10)$	CPU(s)	0.0886	0.2724	0.7187	1.7283
PSPHSS	IT	4	4	4	4
$(\alpha, \omega) = (0.5, 0.1)$	CPU(s)	0.0612	0.0750	0.1286	0.2342
PSHSS	IT	15	28	40	52
$(\alpha, \omega) = (0.5, 0.1)$	CPU(s)	0.0711	0.1364	0.3605	1.0103
PGSOR	IT	8	8	8	8
$(\alpha, \omega) = (0.5, 0.1)$	CPU(s)	0.1024	0.2934	0.7756	1.8693

Table 4Numerical results of Example 2 for the PSPHSS ($V = W$), the PSHSS and the PGSOR iteration method with $(\varpi, \beta) = (-1, 1)$.

Method		Grid			
		16 × 16	32 × 32	48 × 48	64 × 64
PSPHSS	IT	4	4	4	4
$(\alpha_*, \omega_*) = (0.01, 1.4)$	CPU(s)	0.0687	0.0829	0.1356	0.2674
PSHSS	IT	5	5	5	5
$(\alpha_*, \omega_*) = (0.01, 5.5)$	CPU(s)	0.0842	0.0848	0.1689	0.3518
PGSOR	IT	3	3	3	3
$(\alpha_*, \omega_*) = (0.99, 1.4)$	CPU(s)	0.0985	0.2423	0.6424	1.5156
PSPHSS	IT	5	5	5	5
$(\alpha, \omega) = (0.5, 9.5)$	CPU(s)	0.0695	0.0844	0.1413	0.2763
PSHSS	IT	6	9	13	15
$(\alpha, \omega) = (0.5, 9.5)$	CPU(s)	0.0689	0.0958	0.1974	0.4439
PGSOR	IT	8	8	8	8
$(\alpha, \omega) = (0.5, 9.5)$	CPU(s)	0.1243	0.3743	0.9804	2.2316
PSPHSS	IT	4	4	4	4
$(\alpha, \omega) = (0.5, 0.05)$	CPU(s)	0.0688	0.0848	0.1405	0.2731
PSHSS	IT	13	23	32	42
$(\alpha, \omega) = (0.5, 0.05)$	CPU(s)	0.0698	0.1242	0.2938	0.8362
PGSOR	IT	8	8	8	8
$(\alpha, \omega) = (0.5, 0.05)$	CPU(s)	0.1255	0.3781	0.9769	2.2287

experimental optimal parameters, when the parameters $(\varpi, \beta) = (0.2, 0.5)$ and $(\varpi, \beta) = (-1, 1)$, respectively, where the non-optimal parameter α is always fixed, because the distances between 0.5 and the experimental optimal parameters α_* are almost the same for all these three iteration methods. We can see that the speed of convergence for the PSPHSS iteration method is affected slightly with changes in the parameters, and the PSPHSS method is the most efficient from the point of view of iterations and CPU times when the experimental optimal parameters are not used. In a word, the numerical results confirm that the PSPHSS iteration method is very powerful whether the experimental optimal parameters are used or not.

6. Conclusions

In this work, we introduce a single-step preconditioned HSS (SPHSS) and an efficient parameterized SPHSS (PSPHSS) iteration method for solving a class of complex symmetric linear systems. Theoretical analysis shows that, under a loose restriction on the parameters α and ω , the iterative sequences are convergent to the unique solution of the linear system for any initial guess. Furthermore, the quasi-optimal values of the iteration parameters for the SPHSS and PSPHSS iteration methods are also discussed. Both theoretical and numerical results show that the SPHSS and PSPHSS iteration methods have a little advantage over the SHSS and PSHSS iteration methods, respectively. Moreover, the PSPHSS iteration method outperforms several iteration methods from the point of view of the number of iterations and the CPU times when the experimental optimal parameters are used. In particular, the PSPHSS iteration method is the most efficient by comparing with the PSHSS and PGSOR methods if the experimental optimal parameters are not used.

Acknowledgments

This work is supported by NNSF with Nos. 11461046, 61563033 and 11401293, NSF of Jiangxi Province with Nos. 20161ACB21005 and 20151BAB201009, and the Scientific Research Foundation of Graduate School of Nanchang University with Nos. YC2015-S018 and CX2016143.

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