



Shift-splitting fixed point iteration method for solving generalized absolute value equations

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Abstract

Using the shift-splitting strategy, we propose a shift-splitting fixed point iteration (FPI-SS) method for solving large sparse generalized absolute value equations (GAVEs). The FPI-SS method is based on reformulating the GAVE as a two-by-two block nonlinear equation. Several different types of convergence conditions of the FPI-SS method are presented under suitable restrictions. Through numerical experiments, we demonstrate that the FPI-SS method is superior to the fixed point iteration method and the SOR-like iteration method in computing efficiency.

Keywords Generalized absolute value equation · Shift-splitting · Fixed point iteration · Convergence analysis

Mathematics Subject Classification 2010 65F10 · 65H10 · 90C05 · 90C30

1 Introduction

The generalized absolute value equation (GAVE) is formulated as:

$$Ax - B|x| = b, \quad (1)$$

where $A, B \in \mathbb{R}^{n \times n}$ are given large sparse matrices, $b \in \mathbb{R}^n$, and $|x| = (|x_1|, \dots, |x_n|)^T \in \mathbb{R}^n$ denotes the componentwise absolute value of an unknown $x \in \mathbb{R}^n$. If $B = I$, where I stands for an identity matrix of suitable dimension, the GAVE (1) can be simplified to the following absolute value equation (AVE)

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$$Ax - |x| = b. \quad (2)$$

GAVEs have arisen in various scientific and engineering fields and been presented in enormous applications since they were first introduced by Rohn [1]. Among many important applications, a well-known example is the linear complementarity problem (LCP) [2–6]. Besides LCPs, many other optimization problems can be transformed into the GAVEs (1), including linear programming and convex quadratic programming [1, 7].

Due to the existence of the nonlinear term $B|x|$, the GAVE (1) can be regarded as a weakly nonlinear system

$$Ax = G(x), \quad \text{with } G(x) = B|x| + b. \quad (3)$$

For solving the general weakly nonlinear systems

$$Ax = G(x), \quad (4)$$

where the nonlinear function $G : \mathbb{R}^n \rightarrow \mathbb{R}^n$ is B-differentiable, through the two-stage splitting $A = E - F$ and $E = M - N$, Bai for the first time introduced and studied the following two-stage iterative method [8]:

$$Mx^{(k,\ell+1)} = Nx^{(k,\ell)} + Fx^{(k)} + G(x^{(k)}), \quad \text{for } \ell = 0, 1, \dots, l_k - 1, \quad (5)$$

with $x^{(k,0)} := x^{(k)}$ and $x^{(k,l_k)} := x^{(k,l_k)}$. See also [9–11] for related methods. It is noted that the two-stage iterative method provides a general framework of matrix splitting iteration methods for solving the weakly nonlinear systems (4). For the GAVE (1), i.e., the case when $G(x) = B|x| + b$, the two-stage iterative method includes a series of existing matrix splitting iteration methods [12–16] as its special cases. For example, when $E = A$, $F = 0$, $M = E$, $N = 0$, and $l_k \equiv 1$, the two-stage iterative method reduces to the well-known Picard iteration method [12]

$$Ax^{(k+1)} = B|x^{(k)}| + b. \quad (6)$$

Recently, by reformulating the AVE (2) as a two-by-two block nonlinear equation, Ke et al. proposed an SOR-like iteration method [17] for solving the AVE (2). This method was also analyzed in [18]. The SOR-like iteration method received wide attentions and obtained considerable achievements in recent years. Using the similar technology, other SOR-like-based methods [19–21] are presented to solve the AVE (2). In order to further improve computational efficiency, Ke proposed an efficient fixed point iteration (FPI) method [22] to solve the AVE (2), which can be described as

Algorithm 1 (The FPI Method for AVE). Let $A \in \mathbb{R}^{n \times n}$ be a nonsingular matrix and $b \in \mathbb{R}^n$. Given the initial vectors $x^{(0)}, y^{(0)} \in \mathbb{R}^n$, compute $(x^{(k+1)}, y^{(k+1)})$ for $k = 0, 1, 2, \dots$ using the following iteration scheme until $\{(x^{(k)}, y^{(k)})\}_{k=0}^{+\infty}$ satisfies the stopping criterion:

$$\begin{cases} x^{(k+1)} = A^{-1}(y^{(k)} + b), \\ y^{(k+1)} = (1 - \omega)y^{(k)} + \omega|x^{(k+1)}|, \end{cases} \tag{7}$$

where ω is a positive constant.

Note that the FPI method reduces to the Picard iteration method for $\omega = 1$. Owing to the simplicity and effectiveness of FPI method for solving the AVE (2), Yu et al. developed a modified FPI (MFPI) method [23], which is a generalized version of the FPI method.

Clearly, at each step of the FPI method, a linear system $Au = f$ needs to be solved. Since A is always large and sparse, a computationally efficient way is to use matrix splitting iteration methods to obtain the approximate solution of this linear system. For solving non-Hermitian positive definite linear systems, Bai et al. first proposed the shift-splitting (SS) iteration method [24]. Motivated by its promising performance, the SS method was extended to solve many linear systems with special structure such as the saddle point problems [25], block 3×3 saddle point problems [26], and time-harmonic eddy current problems [27]. In this paper, using the shift-splitting [24] of the coefficient matrix A , we propose a shift-splitting fixed point iteration (FPI-SS) method for solving the GAVE (1). Compared with the FPI method, the coefficient matrix of the first sub-iteration scheme of our method is more diagonally dominant. Our method is more efficient than the FPI method and the SOR-like iteration method as shown in our numerical experiments.

In what follows, some notations in this work are described. For $x \in \mathbb{R}^n$, x_i stands for the i th entry of vector x for all $i = 1, 2, \dots, n$. $\text{sgn}(x) \in \mathbb{R}^n$ denotes a vector with components equal to 1, 0, or -1 depending on whether the corresponding component of the vector x is positive, zero, or negative, respectively. Let $\text{diag}(x) \in \mathbb{R}^{n \times n}$ represent a diagonal matrix with x_i as its i th diagonal entry for $i = 1, 2, \dots, n$. For matrix $M \in \mathbb{R}^{n \times n}$, $\|M\|$ denotes the spectral norm defined by $\|M\| := \max\{\|Mx\| : x \in \mathbb{R}^n, \|x\| = 1\}$, where $\|x\|$ is the 2-norm.

The organization of the remaining parts is the following. In Section 2, we present a brief introduction of the FPI method and establish the FPI-SS method for solving the GAVE (1). In Section 3, the convergence theories for the FPI-SS method are presented in detail. In Section 4, we give two numerical examples in Section 4 to verify the effectiveness of our method. Finally, the conclusions are given in Section 5.

2 The shift-splitting fixed point iteration (FPI-SS) method

Let $y = |x|$, then the GAVE (1) is equivalent to

$$\begin{cases} Ax - By = b, \\ -|x| + y = 0, \end{cases} \tag{8}$$

which can be reformulated as the following two-by-two block nonlinear equation

$$\begin{pmatrix} A & -B \\ -H(x) & I \end{pmatrix} \begin{pmatrix} x \\ y \end{pmatrix} = \begin{pmatrix} b \\ 0 \end{pmatrix}, \quad (9)$$

where $H(x) = \text{diag}(\text{sign}(x))$.

If A is a nonsingular matrix, (9) yields the following fixed point equation

$$\begin{cases} x^* = A^{-1}(By^* + b), \\ y^* = (1 - \omega)y^* + \omega|x^*|, \end{cases} \quad (10)$$

where the relaxation parameter $\omega > 0$.

Then, we can obtain the following fixed point iteration (FPI) method for the GAVE (1).

Algorithm 2 (The FPI Method for GAVE). Let $A \in \mathbb{R}^{n \times n}$ be nonsingular, $B \in \mathbb{R}^{n \times n}$ and $b \in \mathbb{R}^n$. Given the initial vectors $x^{(0)}, y^{(0)} \in \mathbb{R}^n$, compute $(x^{(k+1)}, y^{(k+1)})$ for $k = 0, 1, 2, \dots$ using the following iteration scheme until $\{(x^{(k)}, y^{(k)})\}_{k=0}^{+\infty}$ satisfies the stopping criterion:

$$\begin{cases} x^{(k+1)} = A^{-1}(By^{(k)} + b), \\ y^{(k+1)} = (1 - \omega)y^{(k)} + \omega|x^{(k+1)}|, \end{cases} \quad (11)$$

where ω is a positive constant.

It is evident that Algorithm 2 reduces to Algorithm 1 when we take $B = I$. Similarly, if we set $\omega = 1$ in Algorithm 2, the Picard iteration method (6) for solving the GAVE (1) can be obtained. Since the convergence analyses of Algorithm 2 are analogous to those of Algorithm 1 discussed in detail in [22], we do not give them here.

Importantly, by employing the following shift-splitting of the matrix A [24]

$$A = \frac{1}{2}(\alpha I + A) - \frac{1}{2}(\alpha I - A),$$

where the parameter α is a positive constant and the matrix $\alpha I + A$ is invertible, we get the following fixed point equation from (9)

$$\begin{cases} x^* = (\alpha I + A)^{-1}(\alpha I - A)x^* + 2(\alpha I + A)^{-1}(By^* + b), \\ y^* = (1 - \omega)y^* + \omega|x^*|, \end{cases} \quad (12)$$

which leads to the following FPI-SS method for the GAVE (1).

Algorithm 3 (The FPI-SS Method for GAVE). Let $A, B \in \mathbb{R}^{n \times n}$ and $b \in \mathbb{R}^n$. Let α be a positive constant such that $\alpha I + A \in \mathbb{R}^{n \times n}$ is nonsingular. Given the initial vectors $x^{(0)}, y^{(0)} \in \mathbb{R}^n$, compute $(x^{(k+1)}, y^{(k+1)})$ for $k = 0, 1, 2, \dots$ using the following iteration scheme until $\{(x^{(k)}, y^{(k)})\}_{k=0}^{+\infty}$ satisfies the stopping criterion:

$$\begin{cases} x^{(k+1)} = (\alpha I + A)^{-1}(\alpha I - A)x^{(k)} + 2(\alpha I + A)^{-1}(By^{(k)} + b), \\ y^{(k+1)} = (1 - \omega)y^{(k)} + \omega|x^{(k+1)}|, \end{cases} \quad (13)$$

where ω is a positive constant.

Remark 1 If matrix A is positive semi-definite, the condition that $\alpha I + A$ is nonsingular naturally holds. Even if matrix A is singular, we can always find some sufficiently large parameters α to ensure that $\alpha I + A$ is nonsingular. Therefore, the FPI-SS method has a broader range of application than the FPI method. In addition, owing to the positive scalar matrix αI , the matrix $\alpha I + A$ is expected to be strictly diagonally dominant and better conditioned than the matrix A . Thus, our FPI-SS method may have better computing efficiency than the FPI method.

3 Convergence of the FPI-SS method

We first give some lemmas that will be used in convergence analysis of the FPI-SS method for solving the GAVE (1).

Lemma 1 [28–30] *For any vectors $x \in \mathbb{R}^n$ and $y \in \mathbb{R}^n$, the following results hold:*

- (1) $|||x| - |y|| \leq \|x - y\|$;
- (2) if $0 \leq x \leq y$, then $\|x\|_p \leq \|y\|_p$, with $\|\cdot\|_p$ standing for p -norm of vector;
- (3) if $x \leq y$ and P is a nonnegative matrix, then $Px \leq Py$.

Lemma 2 [28, 29] *For any matrices $A, B \in \mathbb{R}^{n \times n}$, if $0 \leq A \leq B$, then $\|A\|_p \leq \|B\|_p$, with $\|\cdot\|_p$ standing for p -norm of matrix.*

Lemma 3 [28, 31] *Both roots of the real quadratic equation $x^2 - ax + b = 0$ are less than one in modulus if and only if $|b| < 1$ and $|a| < 1 + b$.*

In the remainder of this section, we assume that the GAVE (1) has a unique solution. Let (x^*, y^*) be the solution pair of (12) and $(x^{(k)}, y^{(k)})$ be generated by the FPI-SS iteration (13). The iteration errors are denoted by

$$e_k^x = x^* - x^{(k)} \quad \text{and} \quad e_k^y = y^* - y^{(k)}.$$

Then, we can get the following convergence theorem by estimating the above two iteration errors.

Theorem 1 *Let $A, B \in \mathbb{R}^{n \times n}$ and $b \in \mathbb{R}^n$. Let α be a positive constant such that $\alpha I + A \in \mathbb{R}^{n \times n}$ is nonsingular. Denote*

$$\delta = \|(\alpha I + A)^{-1}(\alpha I - A)\|, \beta = 2\|(\alpha I + A)^{-1}B\|, \gamma = |1 - \omega|,$$

and

$$E^{(k+1)} = \begin{pmatrix} \|e_{k+1}^x\| \\ \|e_{k+1}^y\| \end{pmatrix}.$$

Then, we have

$$\|E^{(k+1)}\|_\infty \leq \|L(\alpha, \omega)\|_\infty \cdot \|E^{(k)}\|_\infty, \tag{14}$$

where $\|\cdot\|_\infty$ denotes the ∞ -norm of vector or matrix and

$$L(\alpha, \omega) := \begin{pmatrix} \delta & \beta \\ \omega\delta & \omega\beta + \gamma \end{pmatrix}.$$

Furthermore, $\|L(\alpha, \omega)\|_\infty < 1$ if and only if parameters α and ω satisfy

$$\delta + \beta < 1 \text{ and } 0 < \omega < \frac{2}{1 + \delta + \beta}, \tag{15}$$

i.e., if the conditions (15) hold, the iteration sequence $\{x^{(k)}\}_{k=0}^{+\infty}$ generated by the FPI-SS iteration converges to the unique solution x^* of the GAVE (1) for any initial vector.

Proof Subtracting (13) from (12), we get

$$e_{k+1}^x = (\alpha I + A)^{-1}(\alpha I - A)e_k^x + 2(\alpha I + A)^{-1}Be_k^y, \tag{16}$$

$$e_{k+1}^y = (1 - \omega)e_k^y + \omega(|x^*| - |x^{(k+1)}|). \tag{17}$$

According to (16), we can obtain

$$\|e_{k+1}^x\| \leq \delta\|e_k^x\| + \beta\|e_k^y\|. \tag{18}$$

From (17) and Lemma 1, we have

$$\begin{aligned} \|e_{k+1}^y\| &\leq \gamma \cdot \|e_k^y\| + \omega\||x^*| - |x^{(k+1)}|\| \\ &\leq \gamma \cdot \|e_k^y\| + \omega\|x^* - x^{(k+1)}\| \\ &= \gamma \cdot \|e_k^y\| + \omega\|e_{k+1}^x\|. \end{aligned} \tag{19}$$

Rearranging (18) and (19), we find

$$\begin{pmatrix} 1 & 0 \\ -\omega & 1 \end{pmatrix} \begin{pmatrix} \|e_{k+1}^x\| \\ \|e_{k+1}^y\| \end{pmatrix} \leq \begin{pmatrix} \delta & \beta \\ 0 & \gamma \end{pmatrix} \begin{pmatrix} \|e_k^x\| \\ \|e_k^y\| \end{pmatrix}. \tag{20}$$

Let

$$P = \begin{pmatrix} 1 & 0 \\ \omega & 1 \end{pmatrix} \geq 0.$$

Multiplying (20) from left by the nonnegative matrix P and according to Lemma 1, we have

$$\begin{pmatrix} \|e_{k+1}^x\| \\ \|e_{k+1}^y\| \end{pmatrix} \leq \begin{pmatrix} \delta & \beta \\ \omega\delta & \omega\beta + \gamma \end{pmatrix} \begin{pmatrix} \|e_k^x\| \\ \|e_k^y\| \end{pmatrix}, \tag{21}$$

which can be rewritten as

$$E^{(k+1)} \leq L(\alpha, \omega) \cdot E^{(k)}. \tag{22}$$

Taking the ∞ -norm on both sides of inequality (22) and according to (2) of Lemma 1, the estimation (14) is obtained. Since

$$\|L(\alpha, \omega)\|_\infty = \max\{\delta + \beta, (\delta + \beta)\omega + \gamma\},$$

we have

$$\begin{aligned} \|L(\alpha, \omega)\|_\infty < 1 &\Leftrightarrow \begin{cases} \delta + \beta < 1 \\ (\delta + \beta)\omega + \gamma < 1 \end{cases} \Leftrightarrow \begin{cases} \delta + \beta < 1 \\ |1 - \omega| < 1 - (\delta + \beta)\omega \end{cases} \\ &\Leftrightarrow \begin{cases} \delta + \beta < 1 \\ 1 - (\delta + \beta)\omega > 0 \\ (\delta + \beta)\omega - 1 < 1 - \omega < 1 - (\delta + \beta)\omega \end{cases} \Leftrightarrow \begin{cases} \delta + \beta < 1 \\ \omega < \frac{1}{\delta + \beta} \\ 0 < \omega < \frac{2}{1 + \delta + \beta} \end{cases} \\ &\Leftrightarrow \begin{cases} \delta + \beta < 1 \\ 0 < \omega < \frac{2}{1 + \delta + \beta} \end{cases}. \end{aligned}$$

From (14), we deduce that

$$0 \leq \|E^{(k)}\|_\infty \leq \|L(\alpha, \omega)\|_\infty \cdot \|E^{(k-1)}\|_\infty \leq \dots \leq \|L(\alpha, \omega)\|_\infty^k \cdot \|E^{(0)}\|_\infty.$$

Hence if the conditions (15) are satisfied, then we have $\lim_{k \rightarrow \infty} \|E^{(k)}\|_\infty = 0$.

As

$$\|E^{(k)}\|_\infty = \max\{\|e_k^x\|, \|e_k^y\|\},$$

it follows that

$$\lim_{k \rightarrow \infty} \|e_k^x\| = 0 \quad \text{and} \quad \lim_{k \rightarrow \infty} \|e_k^y\| = 0,$$

which mean that the iteration sequence $\{(x^{(k)}, y^{(k)})\}_{k=0}^{+\infty}$ is convergent to (x^*, y^*) under the conditions (15). This proves the theorem.

Using a different error estimate by a new weighted norm, we can obtain another convergence theorem as follows.

Theorem 2 *Let the assumptions of Theorem 1 hold, $\delta, \beta,$ and γ be defined as in Theorem 1. Denote*

$$E_{\omega}^{(k+1)} = \begin{pmatrix} \|e_{k+1}^x\| \\ \omega^{-1} \|e_{k+1}^y\| \end{pmatrix}.$$

Then, we have

$$\|E_{\omega}^{(k+1)}\| \leq \|T(\alpha, \omega)\| \cdot \|E_{\omega}^{(k)}\|, \quad (23)$$

where

$$T(\alpha, \omega) := \begin{pmatrix} \delta & \omega\beta \\ \delta & \omega\beta + \gamma \end{pmatrix}.$$

Furthermore, $\|T(\alpha, \omega)\| < 1$ if and only if parameters α and ω satisfy

$$\delta\gamma < 1 \quad (24)$$

and

$$2\delta^2 + \omega^2\beta^2 + (\omega\beta + \gamma)^2 - \delta^2\gamma^2 - 1 < 0, \quad (25)$$

i.e., if the conditions (24)–(25) hold, the iteration sequence $\{x^{(k)}\}_{k=0}^{+\infty}$ generated by the FPI-SS iteration converges to the unique solution x^* of the GAVE (1) for any initial vector.

Proof Denote

$$D = \begin{pmatrix} 1 & 0 \\ 0 & \omega^{-1} \end{pmatrix} > 0.$$

According to Lemma 1, we multiply left (21) by matrix D to obtain

$$\begin{pmatrix} \|e_{k+1}^x\| \\ \omega^{-1} \|e_{k+1}^y\| \end{pmatrix} \leq \begin{pmatrix} \delta & \omega\beta \\ \delta & \omega\beta + \gamma \end{pmatrix} \begin{pmatrix} \|e_k^x\| \\ \omega^{-1} \|e_k^y\| \end{pmatrix},$$

which can be rewritten as

$$E_{\omega}^{(k+1)} \leq T(\alpha, \omega) \cdot E_{\omega}^{(k)}.$$

From the above, it follows that (23) holds.

Let λ be an eigenvalue of the matrix $Q := T(\alpha, \omega)^T T(\alpha, \omega)$. Since

$$Q = \begin{pmatrix} 2\delta^2 & 2\omega\beta\delta + \delta\gamma \\ 2\omega\beta\delta + \delta\gamma & \omega^2\beta^2 + (\omega\beta + \gamma)^2 \end{pmatrix},$$

we get

$$\text{tr}(Q) = 2\delta^2 + \omega^2\beta^2 + (\omega\beta + \gamma)^2$$

and

$$\det(Q) = \delta^2\gamma^2.$$

Thus, λ is the root of the following real quadratic equation

$$\lambda^2 - (2\delta^2 + \omega^2\beta^2 + (\omega\beta + \gamma)^2)\lambda + \delta^2\gamma^2 = 0. \tag{26}$$

From Lemma 3, it follows that $\|T(\alpha, \omega)\| < 1$ if and only if

$$\delta^2\gamma^2 < 1,$$

and

$$2\delta^2 + \omega^2\beta^2 + (\omega\beta + \gamma)^2 < 1 + \delta^2\gamma^2.$$

From (23), we conclude that

$$0 \leq \|E_\omega^{(k)}\| \leq \|T(\alpha, \omega)\| \cdot \|E_\omega^{(k-1)}\| \leq \dots \leq \|T(\alpha, \omega)\|^k \cdot \|E_\omega^{(0)}\|.$$

Hence, we have $\lim_{k \rightarrow \infty} \|E_\omega^{(k)}\| = 0$ when the conditons (24)–(25) are satisfied.

From the definition

$$\|E_\omega^{(k)}\| = \sqrt{\|e_k^x\|^2 + \omega^{-2}\|e_k^y\|^2},$$

we get

$$\lim_{k \rightarrow \infty} \|e_k^x\| = 0 \quad \text{and} \quad \lim_{k \rightarrow \infty} \|e_k^y\| = 0,$$

which mean that the iteration sequence $\{(x^{(k)}, y^{(k)})\}_{k=0}^{+\infty}$ is convergent to (x^*, y^*) under the conditions (24)–(25). This completes the proof.

Theorem 2 shows that in order to obtain the convergence of the FPI-SS method, we need to find the conditions in which $\|T(\alpha, \omega)\| < 1$ holds. Here, we give convergence conditions that are simpler than those in Theorem 2.

Corollary 1 *Let the assumptions of Theorem 1 hold, $\delta, \beta,$ and γ be defined as in Theorem 1, $E_\omega^{(k+1)}$ and $T(\alpha, \omega)$ be defined as in Theorem 2. If*

$$\delta < \frac{3 - \sqrt{5}}{2}, \beta < \frac{\sqrt{5} - 1}{2}, \tag{27}$$

and

$$\frac{\sqrt{5} - 1}{2} < \omega < \min\left\{\frac{3 - \sqrt{5}}{2\beta}, \frac{5 - \sqrt{5}}{2}\right\}, \tag{28}$$

then $\|T(\alpha, \omega)\| < 1$, i.e., the FPI-SS method is convergent when the conditions (27)–(28) hold.

Proof Let $\eta = \max\{\delta, \omega\beta, \gamma\}$, we can get

$$0 \leq T(\alpha, \omega) = \begin{pmatrix} \delta & \omega\beta \\ \delta & \omega\beta + \gamma \end{pmatrix} \leq \begin{pmatrix} \eta & \eta \\ \eta & 2\eta \end{pmatrix} = \eta \begin{pmatrix} 1 & 1 \\ 1 & 2 \end{pmatrix} := \eta K,$$

where

$$K = \begin{pmatrix} 1 & 1 \\ 1 & 2 \end{pmatrix}.$$

From lemma 2, we obtain

$$\|T(\alpha, \omega)\| \leq \|\eta K\| = \eta \|K\| = \eta \cdot \frac{3 + \sqrt{5}}{2}.$$

Let $\theta = \frac{3 - \sqrt{5}}{2}$. Hence, we have $\|T(\alpha, \omega)\| < 1$ if $\eta < \theta$. Then,

$$\begin{aligned} \eta < \theta &\Leftrightarrow \begin{cases} \delta < \theta \\ \omega\beta < \theta \\ \gamma = |1 - \omega| < \theta \end{cases} \Leftrightarrow \begin{cases} \delta < \theta \\ \omega < \frac{\theta}{\beta} \\ 1 - \theta < \omega < 1 + \theta \end{cases} \\ &\Leftrightarrow \begin{cases} \delta < \theta \\ 1 - \theta < \omega < \min\{\frac{\theta}{\beta}, 1 + \theta\} \\ 1 - \theta < \frac{\theta}{\beta} \end{cases} \\ &\Leftrightarrow \begin{cases} \delta < \theta \\ \frac{\sqrt{5}-1}{2} < \omega < \min\{\frac{3-\sqrt{5}}{2\beta}, \frac{5-\sqrt{5}}{2}\} \\ \beta < \frac{\theta}{1-\theta} = \frac{\sqrt{5}-1}{2} \end{cases} \end{aligned}$$

Therefore, if the conditions (27)–(28) are satisfied, the iteration sequence $\{(x^{(k)}, y^{(k)})\}_{k=0}^{+\infty}$ is convergent to (x^*, y^*) .

4 Numerical experiments

In this section, two examples from LCPs are presented to show the feasibility and effectiveness of the FPI-SS method. We compare the FPI-SS method with the FPI method [22] and the SOR-like iteration method [17, 18] from aspects of the numbers of iteration steps (denoted as “IT”), elapsed CPU time in seconds (denoted as “CPU”), and relative residual error (denoted as “RES”) which is defined by

$$\text{RES}(x^{(k)}) := \frac{\|Ax^{(k)} - B|x^{(k)}| - b\|_2}{\|b\|_2}.$$

In our implementation, all initial guess vectors $x^{(0)}$ and $y^{(0)}$ are chosen to zero vectors and all iterations are terminated if $\text{RES} \leq 10^{-6}$ or the maximum number

of iteration steps k_{\max} exceeds 500. All computations are performed in MATLAB R2018b on a personal computer with 2.40GHz central processing unit (Intel(R) Core(TM) i5-6200U) and 8 GB memory.

Consider the following LCP(q, M) [2]: to derive two real vectors $z, \omega \in \mathbb{R}^n$ such that

$$z \geq 0, \quad \omega = Mz + q \geq 0, \quad z^T \omega = 0, \tag{29}$$

where $M \in \mathbb{R}^{n \times n}$ and $q \in \mathbb{R}^n$ are given. From [3–6], the LCP(q, M) (29) can be formulated as the following GAVE:

$$(M + I)x - (M - I)|x| = q, \tag{30}$$

with

$$x = \frac{1}{2}((M - I)z + q).$$

Example 1 ([5, 6]) The matrix $M \in \mathbb{R}^{n \times n}$ is defined by $M = \hat{M} + \mu I \in \mathbb{R}^{n \times n}$ and $q \in \mathbb{R}^n$ is defined by $q = -Mz^*$, where

$$\hat{M} = \text{Tridiag}(-I, S, -I) = \begin{bmatrix} S & -I & 0 & \dots & 0 & 0 \\ -I & S & -I & \dots & 0 & 0 \\ 0 & -I & S & \dots & 0 & 0 \\ \vdots & \vdots & & \ddots & \vdots & \vdots \\ 0 & 0 & \dots & \dots & S & -I \\ 0 & 0 & \dots & \dots & -I & S \end{bmatrix} \in \mathbb{R}^{n \times n}$$

is a block-tridiagonal matrix,

$$S = \text{tridiag}(-1, 4, -1) = \begin{bmatrix} 4 & -1 & 0 & \dots & 0 & 0 \\ -1 & 4 & -1 & \dots & 0 & 0 \\ 0 & -1 & 4 & \dots & 0 & 0 \\ \vdots & \vdots & & \ddots & \vdots & \vdots \\ 0 & 0 & \dots & \dots & 4 & -1 \\ 0 & 0 & \dots & \dots & -1 & 4 \end{bmatrix} \in \mathbb{R}^{m \times m}$$

is a tridiagonal matrix, $n = m^2$, and $z^* = (1, 2, 1, 2, \dots, 1, 2, \dots)^T \in \mathbb{R}^n$ is the unique solution of the LCP(q, M) (29). It can be derived that $x^* = (-0.5, -1, -0.5, -1, \dots, -0.5, -1, \dots)^T \in \mathbb{R}^n$ is the exact solution after formulating the LCP(q, M) (29) as the GAVE (30).

For various problem sizes n , the optimal experimental parameters, the iteration steps, CPU time, and relative residual errors of three methods in the case of $\mu = 1$ and $\mu = 4$ are listed in Tables 1 and 2, respectively.

Table 1 Numerical results for Example 1 with $\mu = 1$

Method	n	50^2	100^2	150^2	200^2
SOR-like	ω_{exp}	1.0	1.0	1.0	1.0
	IT	65	65	65	65
	CPU	0.2185	1.3265	3.1923	6.3418
	RES	8.1731e-07	9.3586e-07	9.7624e-07	9.9659e-07
FPI	ω_{exp}	1.9	1.9	1.9	1.9
	IT	32	32	32	32
	CPU	0.1028	0.6670	1.5182	3.0045
	RES	8.6927e-07	8.1065e-07	7.8688e-07	7.7407e-07
FPI-SS	ω_{exp}	1.3	1.4	1.3	1.3
	α_{exp}	3.3	4.6	3.0	3.3
	IT	16	16	16	15
	CPU	0.0435	0.2931	0.7731	1.5064
	RES	5.5429e-07	8.3983e-07	6.0380e-07	9.9518e-07

Table 2 Numerical results for Example 1 with $\mu = 4$

Method	n	50^2	100^2	150^2	200^2
SOR-like	ω_{exp}	1.0	1.0	1.0	1.0
	IT	27	27	27	27
	CPU	0.0881	0.5479	1.3626	2.8915
	RES	7.5622e-07	8.0353e-07	8.1946e-07	8.2746e-07
FPI	ω_{exp}	1.7	1.7	1.7	1.7
	IT	13	13	13	13
	CPU	0.0410	0.2745	0.6650	1.3447
	RES	5.0722e-07	5.4099e-07	5.5243e-07	5.5818e-07
FPI-SS	ω_{exp}	1.1	1.1	1.1	1.1
	α_{exp}	4.5	6.9	7.5	7.5
	IT	10	10	9	9
	CPU	0.0324	0.1889	0.4332	0.9821
	RES	9.3877e-07	4.6551e-07	9.7719e-07	9.5782e-07

We find that each tested method converges to the exact solution and the number of iterative steps becomes smaller with the increase of μ . Notably, among these methods, the FPI-SS method requires the least iteration steps and costs the least computing time.

Example 2 ([5]) Consider the LCP(q, M) (29). The matrix $M \in \mathbb{R}^{n \times n}$ is defined by $M = \hat{M} + \mu I \in \mathbb{R}^{n \times n}$ and $q \in \mathbb{R}^n$ is defined by $q = -Mz^*$, where

$$\widehat{M} = \text{Tridiag}(-1.5I, S, -0.5I) = \begin{bmatrix} S & -0.5I & 0 & \dots & 0 & 0 \\ -1.5I & S & -0.5I & \dots & 0 & 0 \\ 0 & -1.5I & S & \dots & 0 & 0 \\ \vdots & \vdots & & \ddots & \vdots & \vdots \\ 0 & 0 & \dots & \dots & S & -0.5I \\ 0 & 0 & \dots & \dots & -1.5I & S \end{bmatrix} \in \mathbb{R}^{n \times n}$$

is a block-tridiagonal matrix,

$$S = \text{tridiag}(-1.5, 4, -0.5) = \begin{bmatrix} 4 & -0.5 & 0 & \dots & 0 & 0 \\ -1.5 & 4 & -0.5 & \dots & 0 & 0 \\ 0 & -1.5 & 4 & \dots & 0 & 0 \\ \vdots & \vdots & & \ddots & \vdots & \vdots \\ 0 & 0 & \dots & \dots & 4 & -0.5 \\ 0 & 0 & \dots & \dots & -1.5 & 4 \end{bmatrix} \in \mathbb{R}^{m \times m}$$

is a tridiagonal matrix, $n = m^2$, and $z^* = (1, 2, 1, 2, \dots, 1, 2, \dots)^T \in \mathbb{R}^n$ is the unique solution of the LCP(q, M) (29). It can be derived that $x^* = (-0.5, -1, -0.5, -1, \dots, -0.5, -1, \dots)^T \in \mathbb{R}^n$ is the exact solution after formulating the LCP(q, M) (29) as the GAVE (30).

In Tables 3 and 4, we list the numerical results of three methods by using experimental optimal parameters in the case of $\mu = 1$ and $\mu = 4$, respectively. From those results, we get the same conclusions as Example 1.

Table 3 Numerical results for Example 2 with $\mu = 1$

Method	n	50^2	100^2	150^2	200^2
SOR-like	ω_{exp}	1.0	1.0	1.0	1.0
	IT	64	65	65	65
	CPU	0.7538	3.3268	8.2104	15.4580
	RES	8.5588e-07	8.7416e-07	9.3491e-07	9.6552e-07
FPI	ω_{exp}	1.8	1.9	1.9	1.9
	IT	33	32	32	32
	CPU	0.3669	1.6016	4.5292	7.5732
	RES	7.3824e-07	9.9926e-07	9.2406e-07	8.8200e-07
FPI-SS	ω_{exp}	1.2	1.2	1.2	1.2
	α_{exp}	2.2	2.3	2.3	2.4
	IT	22	21	21	21
	CPU	0.2111	0.9085	2.6978	5.0868
	RES	9.9926e-07	8.7943e-07	7.0811e-07	6.3176e-07

Table 4 Numerical results for Example 2 with $\mu = 4$

Method	n	50^2	100^2	150^2	200^2
SOR-like	ω_{exp}	1.0	1.0	1.0	1.0
	IT	27	27	27	27
	CPU	0.3075	1.3550	3.5566	6.9559
	RES	7.4004e-07	7.9538e-07	8.1402e-07	8.2338e-07
FPI	ω_{exp}	1.7	1.7	1.7	1.7
	IT	13	13	13	13
	CPU	0.1435	0.6567	1.6867	3.2636
	RES	5.0817e-07	5.4126e-07	5.5257e-07	5.5827e-07
FPI-SS	ω_{exp}	1.1	1.1	1.1	1.1
	α_{exp}	5.7	7.0	6.0	5.8
	IT	11	10	10	10
	CPU	0.1027	0.4260	1.3000	2.5901
	RES	4.4188e-07	8.4124e-07	8.2384e-07	9.0426e-07

5 Conclusion

In this paper, by combining the shift-splitting of the coefficient matrix with the fixed point iteration (FPI) method, we proposed a shift-splitting fixed point iteration (FPI-SS) method to solve the generalized absolute value equation (GAVE). We have given several different types of convergence conditions of the FPI-SS method by introducing two different norms of the iteration error. Furthermore, using two numerical examples from linear complementarity problems, we have demonstrated that the FPI-SS method outperforms the FPI method and the SOR-like iteration method in terms of iteration steps and computing times.

Finally, we should mention that the FPI-SS method can be seen as an inexact version of the FPI method. If we replace the shift-splitting in the FPI-SS algorithm with other matrix splitting such as SOR-based splitting [28, 29, 31] and HSS-based splitting [28, 32–34], we can establish a series of inexact FPI methods which may have similar convergence results. In real applications of inexact FPI algorithms, how to choose the optimal (or quasi-optimal) parameters is an interesting and practical topic, which is left as our future work.

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Declarations

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