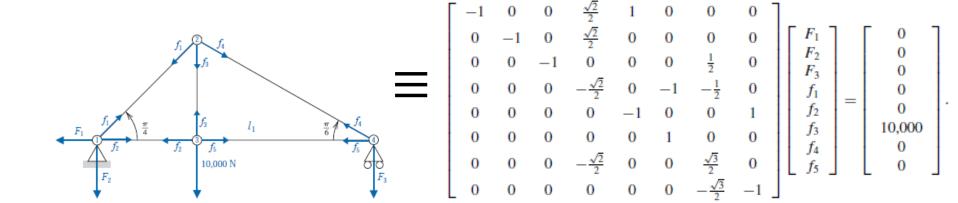
MAP 2220 – FUNDAMENTOS DE ANÁLISE NUMÉRICA 2º Semestre - 2017

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Jacobi's Method

$$x_i^{(k)} = \frac{1}{a_{ii}} \left[\sum_{\substack{j=1\\j\neq i}}^n \left(-a_{ij} x_j^{(k-1)} \right) + b_i \right], \quad \text{for } i = 1, 2, \dots, n.$$

$$\mathbf{x}^{(k)} = T_j \mathbf{x}^{(k-1)} + \mathbf{c}_j.$$

$$T_j = D^{-1}(L+U)$$
 and $\mathbf{c}_j = D^{-1}\mathbf{b}$

The Gauss-Seidel Method

$$x_i^{(k)} = \frac{1}{a_{ii}} \left[-\sum_{j=1}^{i-1} (a_{ij} x_j^{(k)}) - \sum_{j=i+1}^{n} (a_{ij} x_j^{(k-1)}) + b_i \right],$$

$$\mathbf{x}^{(k)} = T_{\varrho} \mathbf{x}^{(k-1)} + \mathbf{c}_{\varrho}.$$

$$\mathbf{x}^{(k)} = T_g \mathbf{x}^{(k-1)} + \mathbf{c}_g$$
. $T_g = (D - L)^{-1} U$ and $\mathbf{c}_g = (D - L)^{-1} \mathbf{b}$,

General Iteration Methods

To study the convergence of general iteration techniques, we need to analyze the formula

$$\mathbf{x}^{(k)} = T\mathbf{x}^{(k-1)} + \mathbf{c}$$
, for each $k = 1, 2, ...$,

where $\mathbf{x}^{(0)}$ is arbitrary. The next lemma and Theorem 7.17 on page 449 provide the key for this study.

Lemma 7.18 If the spectral radius satisfies $\rho(T) < 1$, then $(I - T)^{-1}$ exists, and

$$(I-T)^{-1} = I + T + T^2 + \dots = \sum_{j=0}^{\infty} T^j.$$

Proof Because $T\mathbf{x} = \lambda \mathbf{x}$ is true precisely when $(I - T)\mathbf{x} = (1 - \lambda)\mathbf{x}$, we have λ as an eigenvalue of T precisely when $1 - \lambda$ is an eigenvalue of I - T. But $|\lambda| \le \rho(T) < 1$, so $\lambda = 1$ is not an eigenvalue of T, and 0 cannot be an eigenvalue of I - T. Hence, $(I - T)^{-1}$ exists.

Let
$$S_m = I + T + T^2 + \cdots + T^m$$
. Then

$$(I-T)S_m = (1+T+T^2+\cdots+T^m)-(T+T^2+\cdots+T^{m+1})=I-T^{m+1},$$

and, since T is convergent, Theorem 7.17 implies that

$$\lim_{m\to\infty} (I-T)S_m = \lim_{m\to\infty} (I-T^{m+1}) = I.$$

Thus,
$$(I-T)^{-1} = \lim_{m \to \infty} S_m = I + T + T^2 + \dots = \sum_{i=0}^{\infty} T^i$$
.

Theorem 7.19 For any $\mathbf{x}^{(0)} \in \mathbb{R}^n$, the sequence $\{\mathbf{x}^{(k)}\}_{k=0}^{\infty}$ defined by

$$\mathbf{x}^{(k)} = T\mathbf{x}^{(k-1)} + \mathbf{c}, \quad \text{for each } k \ge 1,$$
 (7.11)

converges to the unique solution of $\mathbf{x} = T\mathbf{x} + \mathbf{c}$ if and only if $\rho(T) < 1$.

Proof First assume that $\rho(T) < 1$. Then,

$$\mathbf{x}^{(k)} = T\mathbf{x}^{(k-1)} + \mathbf{c}$$

$$= T(T\mathbf{x}^{(k-2)} + \mathbf{c}) + \mathbf{c}$$

$$= T^{2}\mathbf{x}^{(k-2)} + (T+I)\mathbf{c}$$

$$\vdots$$

$$= T^{k}\mathbf{x}^{(0)} + (T^{k-1} + \dots + T+I)\mathbf{c}.$$

Because $\rho(T) < 1$, Theorem 7.17 implies that T is convergent, and

$$\lim_{k\to\infty} T^k \mathbf{x}^{(0)} = \mathbf{0}.$$

Lemma 7.18 implies that

$$\lim_{k \to \infty} \mathbf{x}^{(k)} = \lim_{k \to \infty} T^k \mathbf{x}^{(0)} + \left(\sum_{j=0}^{\infty} T^j\right) \mathbf{c} = \mathbf{0} + (I - T)^{-1} \mathbf{c} = (I - T)^{-1} \mathbf{c}.$$

Hence, the sequence $\{x^{(k)}\}\$ converges to the vector $\mathbf{x} \equiv (I - T)^{-1}\mathbf{c}$ and $\mathbf{x} = T\mathbf{x} + \mathbf{c}$.

To prove the converse, we will show that for any $z \in \mathbb{R}^n$, we have $\lim_{k\to\infty} T^k z = 0$. By Theorem 7.17, this is equivalent to $\rho(T) < 1$. But $z \in \mathbb{R}^n$ was arbitrary, so by Theorem 7.17, T is convergent and $\rho(T) < 1$.

The proof of the following corollary is similar to the proofs in Corollary 2.5 on page 62. It is considered in Exercise 13.

Corollary 7.20 If ||T|| < 1 for any natural matrix norm and c is a given vector, then the sequence $\{\mathbf{x}^{(k)}\}_{k=0}^{\infty}$ defined by $\mathbf{x}^{(k)} = T\mathbf{x}^{(k-1)} + \mathbf{c}$ converges, for any $\mathbf{x}^{(0)} \in \mathbb{R}^n$, to a vector $\mathbf{x} \in \mathbb{R}^n$, with $\mathbf{x} = T\mathbf{x} + \mathbf{c}$, and the following error bounds hold:

(i)
$$\|\mathbf{x} - \mathbf{x}^{(k)}\| \le \|T\|^k \|\mathbf{x}^{(0)} - \mathbf{x}\|;$$
 (ii) $\|\mathbf{x} - \mathbf{x}^{(k)}\| \le \frac{\|T\|^k}{1 - \|T\|} \|\mathbf{x}^{(1)} - \mathbf{x}^{(0)}\|.$

We have seen that the Jacobi and Gauss-Seidel iterative techniques can be written

$$\mathbf{x}^{(k)} = T_j \mathbf{x}^{(k-1)} + \mathbf{c}_j$$
 and $\mathbf{x}^{(k)} = T_g \mathbf{x}^{(k-1)} + \mathbf{c}_g$,

using the matrices

$$T_j = D^{-1}(L+U)$$
 and $T_g = (D-L)^{-1}U$.

If $\rho(T_j)$ or $\rho(T_g)$ is less than 1, then the corresponding sequence $\{\mathbf{x}^{(k)}\}_{k=0}^{\infty}$ will converge to the solution \mathbf{x} of $A\mathbf{x} = \mathbf{b}$. For example, the Jacobi scheme has

$$\mathbf{x}^{(k)} = D^{-1}(L+U)\mathbf{x}^{(k-1)} + D^{-1}\mathbf{b},$$

and, if $\{\mathbf{x}^{(k)}\}_{k=0}^{\infty}$ converges to \mathbf{x} , then

$$\mathbf{x} = D^{-1}(L+U)\mathbf{x} + D^{-1}\mathbf{b}.$$

This implies that

$$D\mathbf{x} = (L+U)\mathbf{x} + \mathbf{b}$$
 and $(D-L-U)\mathbf{x} = \mathbf{b}$.

Since D - L - U = A, the solution x satisfies Ax = b.

We can now give easily verified sufficiency conditions for convergence of the Jacobi and Gauss-Seidel methods. (To prove convergence for the Jacobi scheme see Exercise 14, and for the Gauss-Seidel scheme see [Or2], p. 120.)

Theorem 7.21 If A is strictly diagonally dominant, then for any choice of $\mathbf{x}^{(0)}$, both the Jacobi and Gauss-Seidel methods give sequences $\{\mathbf{x}^{(k)}\}_{k=0}^{\infty}$ that converge to the unique solution of $A\mathbf{x} = \mathbf{b}$.

The relationship of the rapidity of convergence to the spectral radius of the iteration matrix T can be seen from Corollary 7.20. The inequalities hold for any natural matrix norm, so it follows from the statement after Theorem 7.15 on page 446 that

$$\|\mathbf{x}^{(k)} - \mathbf{x}\| \approx \rho(T)^k \|\mathbf{x}^{(0)} - \mathbf{x}\|.$$
 (7.12)

Thus we would like to select the iterative technique with minimal $\rho(T) < 1$ for a particular system $A\mathbf{x} = \mathbf{b}$. No general results exist to tell which of the two techniques, Jacobi or Gauss-Seidel, will be most successful for an arbitrary linear system. In special cases, however, the answer is known, as is demonstrated in the following theorem. The proof of this result can be found in [Y], pp. 120–127.

Theorem 7.22 (Stein-Rosenberg)

If $a_{ij} \le 0$, for each $i \ne j$ and $a_{ii} > 0$, for each i = 1, 2, ..., n, then one and only one of the following statements holds:

(i)
$$0 \le \rho(T_g) < \rho(T_j) < 1$$
; (ii) $1 < \rho(T_j) < \rho(T_g)$;

(iii)
$$\rho(T_j) = \rho(T_g) = 0;$$
 (iv) $\rho(T_j) = \rho(T_g) = 1.$

For the special case described in Theorem 7.22, we see from part (i) that when one method gives convergence, then both give convergence, and the Gauss-Seidel method converges faster than the Jacobi method. Part (ii) indicates that when one method diverges then both diverge, and the divergence is more pronounced for the Gauss-Seidel method.

7.4 Relaxation Techniques for Solving Linear Systems

We saw in Section 7.3 that the rate of convergence of an iterative technique depends on the spectral radius of the matrix associated with the method. One way to select a procedure to accelerate convergence is to choose a method whose associated matrix has minimal spectral radius. Before describing a procedure for selecting such a method, we need to introduce a new means of measuring the amount by which an approximation to the solution to a linear system differs from the true solution to the system. The method makes use of the vector described in the following definition.

Definition 7.23 Suppose $\tilde{\mathbf{x}} \in \mathbb{R}^n$ is an approximation to the solution of the linear system defined by $A\mathbf{x} = \mathbf{b}$. The residual vector for $\tilde{\mathbf{x}}$ with respect to this system is $\mathbf{r} = \mathbf{b} - A\tilde{\mathbf{x}}$.

In procedures such as the Jacobi or Gauss-Seidel methods, a residual vector is associated with each calculation of an approximate component to the solution vector. The true objective is to generate a sequence of approximations that will cause the residual vectors to converge rapidly to zero. Suppose we let

$$\mathbf{r}_{i}^{(k)} = (r_{1i}^{(k)}, r_{2i}^{(k)}, \dots, r_{ni}^{(k)})^{t}$$

denote the residual vector for the Gauss-Seidel method corresponding to the approximate solution vector $\mathbf{x}_i^{(k)}$ defined by

$$\mathbf{x}_i^{(k)} = (x_1^{(k)}, x_2^{(k)}, \dots, x_{i-1}^{(k)}, x_i^{(k-1)}, \dots, x_n^{(k-1)})^t.$$

The *m*th component of $\mathbf{r}_i^{(k)}$ is

$$r_{mi}^{(k)} = b_m - \sum_{j=1}^{i-1} a_{mj} x_j^{(k)} - \sum_{j=i}^n a_{mj} x_j^{(k-1)},$$
 (7.13)

or, equivalently,

$$r_{mi}^{(k)} = b_m - \sum_{j=1}^{i-1} a_{mj} x_j^{(k)} - \sum_{j=i+1}^n a_{mj} x_j^{(k-1)} - a_{mi} x_i^{(k-1)},$$

for each $m = 1, 2, \ldots, n$.

In particular, the *i*th component of $\mathbf{r}_i^{(k)}$ is

$$r_{ii}^{(k)} = b_i - \sum_{j=1}^{i-1} a_{ij} x_j^{(k)} - \sum_{j=i+1}^{n} a_{ij} x_j^{(k-1)} - a_{ii} x_i^{(k-1)},$$

SO

$$a_{ii}x_i^{(k-1)} + r_{ii}^{(k)} = b_i - \sum_{j=1}^{i-1} a_{ij}x_j^{(k)} - \sum_{j=i+1}^{n} a_{ij}x_j^{(k-1)}.$$
 (7.14)

Recall, however, that in the Gauss-Seidel method, $x_i^{(k)}$ is chosen to be

$$x_i^{(k)} = \frac{1}{a_{ii}} \left[b_i - \sum_{j=1}^{i-1} a_{ij} x_j^{(k)} - \sum_{j=i+1}^n a_{ij} x_j^{(k-1)} \right], \tag{7.15}$$

so Eq. (7.14) can be rewritten as

$$a_{ii}x_i^{(k-1)} + r_{ii}^{(k)} = a_{ii}x_i^{(k)}.$$

Consequently, the Gauss-Seidel method can be characterized as choosing $x_i^{(k)}$ to satisfy

$$x_i^{(k)} = x_i^{(k-1)} + \frac{r_{ii}^{(k)}}{a_{ii}}. (7.16)$$

We can derive another connection between the residual vectors and the Gauss-Seidel technique. Consider the residual vector $\mathbf{r}_{i+1}^{(k)}$, associated with the vector $\mathbf{x}_{i+1}^{(k)} = (x_1^{(k)}, \dots, x_i^{(k)}, x_{i+1}^{(k-1)}, \dots, x_n^{(k-1)})^t$. By Eq. (7.13) the *i*th component of $\mathbf{r}_{i+1}^{(k)}$ is

$$\begin{split} r_{i,i+1}^{(k)} &= b_i - \sum_{j=1}^i a_{ij} x_j^{(k)} - \sum_{j=i+1}^n a_{ij} x_j^{(k-1)} \\ &= b_i - \sum_{j=1}^{i-1} a_{ij} x_j^{(k)} - \sum_{j=i+1}^n a_{ij} x_j^{(k-1)} - a_{ii} x_i^{(k)}. \end{split}$$

By the manner in which $x_i^{(k)}$ is defined in Eq. (7.15) we see that $r_{i,i+1}^{(k)} = 0$. In a sense, then, the Gauss-Seidel technique is characterized by choosing each $x_{i+1}^{(k)}$ in such a way that the *i*th component of $\mathbf{r}_{i+1}^{(k)}$ is zero.

Choosing $x_{i+1}^{(k)}$ so that one coordinate of the residual vector is zero, however, is not necessarily the most efficient way to reduce the norm of the vector $\mathbf{r}_{i+1}^{(k)}$. If we modify the Gauss-Seidel procedure, as given by Eq. (7.16), to

$$x_i^{(k)} = x_i^{(k-1)} + \omega \frac{r_{ii}^{(k)}}{a_{ii}},\tag{7.17}$$

then for certain choices of positive ω we can reduce the norm of the residual vector and obtain significantly faster convergence.

Methods involving Eq. (7.17) are called **relaxation methods**. For choices of ω with $0 < \omega < 1$, the procedures are called **under-relaxation methods**. We will be interested in choices of ω with $1 < \omega$, and these are called **over-relaxation methods**. They are used to accelerate the convergence for systems that are convergent by the Gauss-Seidel technique. The methods are abbreviated **SOR**, for **Successive Over-Relaxation**, and are particularly useful for solving the linear systems that occur in the numerical solution of certain partial-differential equations.

Before illustrating the advantages of the SOR method, we note that by using Eq. (7.14), we can reformulate Eq. (7.17) for calculation purposes as

$$x_i^{(k)} = (1 - \omega)x_i^{(k-1)} + \frac{\omega}{a_{ii}} \left[b_i - \sum_{j=1}^{i-1} a_{ij} x_j^{(k)} - \sum_{j=i+1}^n a_{ij} x_j^{(k-1)} \right].$$

To determine the matrix form of the SOR method, we rewrite this as

$$a_{ii}x_i^{(k)} + \omega \sum_{j=1}^{i-1} a_{ij}x_j^{(k)} = (1 - \omega)a_{ii}x_i^{(k-1)} - \omega \sum_{j=i+1}^n a_{ij}x_j^{(k-1)} + \omega b_i,$$

so that in vector form, we have

$$(D - \omega L)\mathbf{x}^{(k)} = [(1 - \omega)D + \omega U]\mathbf{x}^{(k-1)} + \omega \mathbf{b}.$$

That is,

$$\mathbf{x}^{(k)} = (D - \omega L)^{-1} [(1 - \omega)D + \omega U] \mathbf{x}^{(k-1)} + \omega (D - \omega L)^{-1} \mathbf{b}. \tag{7.18}$$

Letting $T_{\omega} = (D - \omega L)^{-1}[(1 - \omega)D + \omega U]$ and $\mathbf{c}_{\omega} = \omega(D - \omega L)^{-1}\mathbf{b}$, gives the SOR technique the form

$$\mathbf{x}^{(k)} = T_{\omega} \mathbf{x}^{(k-1)} + \mathbf{c}_{\omega}.$$
 (7.19)

Example 1 The linear system Ax = b given by

$$4x_1 + 3x_2 = 24,$$

 $3x_1 + 4x_2 - x_3 = 30,$
 $-x_2 + 4x_3 = -24,$

has the solution $(3, 4, -5)^t$. Compare the iterations from the Gauss-Seidel method and the SOR method with $\omega = 1.25$ using $\mathbf{x}^{(0)} = (1, 1, 1)^t$ for both methods.

Solution For each k = 1, 2, ..., the equations for the Gauss-Seidel method are

$$x_1^{(k)} = -0.75x_2^{(k-1)} + 6,$$

$$x_2^{(k)} = -0.75x_1^{(k)} + 0.25x_3^{(k-1)} + 7.5,$$

$$x_3^{(k)} = 0.25x_2^{(k)} - 6,$$

and the equations for the SOR method with $\omega = 1.25$ are

$$x_1^{(k)} = -0.25x_1^{(k-1)} - 0.9375x_2^{(k-1)} + 7.5,$$

$$x_2^{(k)} = -0.9375x_1^{(k)} - 0.25x_2^{(k-1)} + 0.3125x_3^{(k-1)} + 9.375,$$

$$x_3^{(k)} = 0.3125x_2^{(k)} - 0.25x_3^{(k-1)} - 7.5.$$

The first seven iterates for each method are listed in Tables 7.3 and 7.4. For the iterates to be accurate to seven decimal places, the Gauss-Seidel method requires 34 iterations, as opposed to 14 iterations for the SOR method with $\omega = 1.25$.

Table 7.3

k	0	1	2	3	4	5	6	7
$x_1^{(k)}$	1	5.250000	3.1406250	3.0878906	3.0549316	3.0343323	3.0214577	3.0134110
$x_{2}^{(k)}$	1	3.812500	3.8828125	3.9267578	3.9542236	3.9713898	3.9821186	3.9888241
$x_3^{(k)}$	1	-5.046875	-5.0292969	-5.0183105	-5.0114441	-5.0071526	-5.0044703	-5.0027940

Table 7.4

k	0	1	2	3	4	5	6	7
$x_1^{(k)}$	1	6.312500	2.6223145	3.1333027	2.9570512	3.0037211	2.9963276	3.0000498
$x_{2}^{(k)}$	1	3.5195313	3.9585266	4.0102646	4.0074838	4.0029250	4.0009262	4.0002586
$x_3^{(k)}$	1	-6.6501465	-4.6004238	-5.0966863	-4.9734897	-5.0057135	-4.9982822	-5.0003486

An obvious question to ask is how the appropriate value of ω is chosen when the SOR method is used. Although no complete answer to this question is known for the general $n \times n$ linear system, the following results can be used in certain important situations.

Theorem 7.24 (Kahan)

If $a_{ii} \neq 0$, for each i = 1, 2, ..., n, then $\rho(T_{\omega}) \geq |\omega - 1|$. This implies that the SOR method can converge only if $0 < \omega < 2$.

The proof of this theorem is considered in Exercise 9. The proof of the next two results can be found in [Or2], pp. 123–133. These results will be used in Chapter 12.

Theorem 7.25 (Ostrowski-Reich)

If A is a positive definite matrix and $0 < \omega < 2$, then the SOR method converges for any choice of initial approximate vector $\mathbf{x}^{(0)}$.

Theorem 7.26 If A is positive definite and tridiagonal, then $\rho(T_g) = [\rho(T_j)]^2 < 1$, and the optimal choice of ω for the SOR method is

$$\omega = \frac{2}{1 + \sqrt{1 - [\rho(T_j)]^2}}.$$

With this choice of ω , we have $\rho(T_{\omega}) = \omega - 1$.

$$A = \left[\begin{array}{rrr} 4 & 3 & 0 \\ 3 & 4 & -1 \\ 0 & -1 & 4 \end{array} \right].$$

Solution This matrix is clearly tridiagonal, so we can apply the result in Theorem 7.26 if we can also who that it is positive definite. Because the matrix is symmetric, Theorem 6.24 on page 416 states that it is positive definite if and only if all its leading principle submatrices has a positive determinant. This is easily seen to be the case because

$$det(A) = 24$$
, $det \begin{pmatrix} \begin{bmatrix} 4 & 3 \\ 3 & 4 \end{bmatrix} \end{pmatrix} = 7$, and $det([4]) = 4$.

Because

$$T_j = D^{-1}(L+U) = \begin{bmatrix} \frac{1}{4} & 0 & 0 \\ 0 & \frac{1}{4} & 0 \\ 0 & 0 & \frac{1}{4} \end{bmatrix} \begin{bmatrix} 0 & -3 & 0 \\ -3 & 0 & 1 \\ 0 & 1 & 0 \end{bmatrix} = \begin{bmatrix} 0 & -0.75 & 0 \\ -0.75 & 0 & 0.25 \\ 0 & 0.25 & 0 \end{bmatrix},$$

we have

$$T_j - \lambda I = \begin{bmatrix} -\lambda & -0.75 & 0 \\ -0.75 & -\lambda & 0.25 \\ 0 & 0.25 & -\lambda \end{bmatrix},$$

SO

$$\det(T_j - \lambda I) = -\lambda(\lambda^2 - 0.625).$$

Thus

$$\rho(T_j) = \sqrt{0.625}$$

and

$$\omega = \frac{2}{1 + \sqrt{1 - [\rho(T_i)]^2}} = \frac{2}{1 + \sqrt{1 - 0.625}} \approx 1.24.$$

This explains the rapid convergence obtained in Example 1 when using $\omega = 1.25$.

SOR

To solve $A\mathbf{x} = \mathbf{b}$ given the parameter ω and an initial approximation $\mathbf{x}^{(0)}$:

INPUT the number of equations and unknowns n; the entries a_{ij} , $1 \le i, j \le n$, of the matrix A; the entries b_i , $1 \le i \le n$, of **b**; the entries XO_i , $1 \le i \le n$, of $XO = X^{(0)}$; the parameter ω ; tolerance TOL; maximum number of iterations N.

OUTPUT the approximate solution x_1, \ldots, x_n or a message that the number of iterations was exceeded.

Step 1 Set k = 1.

Step 2 While $(k \le N)$ do Steps 3–6.

Step 3 For i = 1, ..., n

set
$$x_i = (1 - \omega)XO_i + \frac{1}{a_{ii}} \left[\omega \left(-\sum_{j=1}^{i-1} a_{ij} x_j - \sum_{j=i+1}^n a_{ij} XO_j + b_i \right) \right].$$

Step 4 If $||\mathbf{x} - \mathbf{XO}|| < TOL$ then OUTPUT $(x_1, ..., x_n)$; (The procedure was successful.) STOP.

Step 5 Set k = k + 1.

Step 6 For i = 1, ..., n set $XO_i = x_i$.

Step 7 OUTPUT ('Maximum number of iterations exceeded'); (The procedure was successful.) STOP.

EXERCISE SET 7.4

Find the first two iterations of the SOR method with ω = 1.1 for the following linear systems, using x⁽⁰⁾ = 0:

a.
$$3x_1 - x_2 + x_3 = 1$$
,
 $3x_1 + 6x_2 + 2x_3 = 0$,
 $3x_1 + 3x_2 + 7x_3 = 4$.

b.
$$10x_1 - x_2 = 9,$$

 $-x_1 + 10x_2 - 2x_3 = 7,$
 $-2x_2 + 10x_3 = 6.$

$$x_i^{(k)} = \frac{1}{a_{ii}} \left[b_i - \sum_{j=1}^{i-1} a_{ij} x_j^{(k)} - \sum_{j=i+1}^n a_{ij} x_j^{(k-1)} \right],$$

$$x_i^{(k)} = (1 - \omega)x_i^{(k-1)} + \frac{\omega}{a_{ii}} \left[b_i - \sum_{j=1}^{i-1} a_{ij} x_j^{(k)} - \sum_{j=i+1}^n a_{ij} x_j^{(k-1)} \right].$$



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