

Numerical Analysis II

Iterative techniques in matrix algebra

Instructor: Wei-Cheng Wang ¹

Department of Mathematics
National TsingHua University

Spring 2011



¹These slides are based on Prof. Tsung-Ming Huang(NTNU)'s original slides

Outline

- 1 Norms of vectors and matrices
- 2 Eigenvalues and eigenvectors
- 3 Iterative techniques for solving linear systems
- 4 Error bounds and iterative refinement
- 5 The conjugate gradient method



Definition

$\|\cdot\| : \mathbb{R}^n \rightarrow \mathbb{R}$ is a vector norm if

- (i) $\|x\| \geq 0, \forall x \in \mathbb{R}^n,$
- (ii) $\|x\| = 0$ if and only if $x = 0,$
- (iii) $\|\alpha x\| = |\alpha| \|x\| \forall \alpha \in \mathbb{R}$ and $x \in \mathbb{R}^n,$
- (iv) $\|x + y\| \leq \|x\| + \|y\| \forall x, y \in \mathbb{R}^n.$

Definition

The ℓ_2 and ℓ_∞ norms for $x = [x_1, x_2, \dots, x_n]^T$ are defined by

$$\|x\|_2 = (x^T x)^{1/2} = \left\{ \sum_{i=1}^n x_i^2 \right\}^{1/2} \quad \text{and} \quad \|x\|_\infty = \max_{1 \leq i \leq n} |x_i|.$$

The ℓ_2 norm is also called the Euclidean norm.



Theorem (Cauchy-Bunyakovsky-Schwarz inequality)

For each $x = [x_1, x_2, \dots, x_n]^T$ and $y = [y_1, y_2, \dots, y_n]^T$ in \mathbb{R}^n ,

$$x^T y = \sum_{i=1}^n x_i y_i \leq \left\{ \sum_{i=1}^n x_i^2 \right\}^{1/2} \left\{ \sum_{i=1}^n y_i^2 \right\}^{1/2} = \|x\|_2 \cdot \|y\|_2.$$

Proof: If $x = 0$ or $y = 0$, the result is immediate.

Suppose $x \neq 0$ or $y \neq 0$. For each $\alpha \in \mathbb{R}$,

$$0 \leq \|x - \alpha y\|_2^2 = \sum_{i=1}^n (x_i - \alpha y_i)^2 = \sum_{i=1}^n x_i^2 - 2\alpha \sum_{i=1}^n x_i y_i + \alpha^2 \sum_{i=1}^n y_i^2,$$

and

$$2\alpha \sum_{i=1}^n x_i y_i \leq \sum_{i=1}^n x_i^2 + \alpha^2 \sum_{i=1}^n y_i^2 = \|x\|_2^2 + \alpha^2 \|y\|_2^2.$$



Since $\|x\|_2 > 0$ and $\|y\|_2 > 0$, we can let

$$\alpha = \frac{\|x\|_2}{\|y\|_2}$$

to give

$$\left(2 \frac{\|x\|_2}{\|y\|_2}\right) \left(\sum_{i=1}^n x_i y_i\right) \leq \|x\|_2^2 + \frac{\|x\|_2^2}{\|y\|_2^2} \|y\|_2^2 = 2\|x\|_2^2.$$

Thus

$$x^T y = \sum_{i=1}^n x_i y_i \leq \|x\|_2 \|y\|_2.$$



For each $x, y \in \mathbb{R}^n$,

$$\begin{aligned}\|x + y\|_\infty &= \max_{1 \leq i \leq n} |x_i + y_i| \leq \max_{1 \leq i \leq n} (|x_i| + |y_i|) \\ &\leq \max_{1 \leq i \leq n} |x_i| + \max_{1 \leq i \leq n} |y_i| = \|x\|_\infty + \|y\|_\infty\end{aligned}$$

and

$$\begin{aligned}\|x + y\|_2^2 &= \sum_{i=1}^n (x_i + y_i)^2 = \sum_{i=1}^n x_i^2 + 2 \sum_{i=1}^n x_i y_i + \sum_{i=1}^n y_i^2 \\ &\leq \|x\|_2^2 + 2\|x\|_2\|y\|_2 + \|y\|_2^2 = (\|x\|_2 + \|y\|_2)^2,\end{aligned}$$

which gives

$$\|x + y\|_2 \leq \|x\|_2 + \|y\|_2.$$

Definition

A sequence $\{x^{(k)} \in \mathbb{R}^n\}_{k=1}^\infty$ is convergent to x with respect to the norm $\|\cdot\|$ if $\forall \varepsilon > 0, \exists$ an integer $N(\varepsilon)$ such that

$$\|x^{(k)} - x\| < \varepsilon, \quad \forall k \geq N(\varepsilon).$$

Theorem

$\{x^{(k)} \in \mathbb{R}^n\}_{k=1}^{\infty}$ converges to x with respect to $\|\cdot\|_{\infty}$ if and only if

$$\lim_{k \rightarrow \infty} x_i^{(k)} = x_i, \quad \forall i = 1, 2, \dots, n.$$

Proof: “ \Rightarrow ” Given any $\varepsilon > 0$, \exists an integer $N(\varepsilon)$ such that

$$\max_{1 \leq i \leq n} |x_i^{(k)} - x_i| = \|x^{(k)} - x\|_{\infty} < \varepsilon, \quad \forall k \geq N(\varepsilon).$$

This result implies that

$$|x_i^{(k)} - x_i| < \varepsilon, \quad \forall i = 1, 2, \dots, n.$$

Hence

$$\lim_{k \rightarrow \infty} x_i^{(k)} = x_i, \quad \forall i.$$



“ \Leftarrow ” For a given $\varepsilon > 0$, let $N_i(\varepsilon)$ represent an integer with

$$|x_i^{(k)} - x_i| < \varepsilon, \quad \text{whenever } k \geq N_i(\varepsilon).$$

Define

$$N(\varepsilon) = \max_{1 \leq i \leq n} N_i(\varepsilon).$$

If $k \geq N(\varepsilon)$, then

$$\max_{1 \leq i \leq n} |x_i^{(k)} - x_i| = \|x^{(k)} - x\|_\infty < \varepsilon.$$

This implies that $\{x^{(k)}\}$ converges to x with respect to $\|\cdot\|_\infty$. □



Theorem

For each $x \in \mathbb{R}^n$,

$$\|x\|_\infty \leq \|x\|_2 \leq \sqrt{n}\|x\|_\infty.$$

Proof: Let x_j be a coordinate of x such that

$$\|x\|_\infty^2 = |x_j|^2 \leq \sum_{i=1}^n x_i^2 = \|x\|_2^2,$$

so $\|x\|_\infty \leq \|x\|_2$ and

$$\|x\|_2^2 = \sum_{i=1}^n x_i^2 \leq \sum_{i=1}^n x_j^2 = nx_j^2 = n\|x\|_\infty^2,$$

so $\|x\|_2 \leq \sqrt{n}\|x\|_\infty$.



Definition

A matrix norm $\|\cdot\|$ on the set of all $n \times n$ matrices is a real-valued function satisfying for all $n \times n$ matrices A and B and all real number α :

- (i) $\|A\| \geq 0$;
- (ii) $\|A\| = 0$ if and only if $A = 0$;
- (iii) $\|\alpha A\| = |\alpha| \|A\|$;
- (iv) $\|A + B\| \leq \|A\| + \|B\|$;
- (v) $\|AB\| \leq \|A\| \|B\|$;

Theorem

If $\|\cdot\|$ is a vector norm on \mathbb{R}^n , then

$$\|A\| = \max_{\|x\|=1} \|Ax\|$$

is a matrix norm.

For any $z \neq 0$, we have $x = z/\|z\|$ as a unit vector. Hence

$$\|A\| = \max_{\|x\|=1} \|Ax\| = \max_{z \neq 0} \left\| A \left(\frac{z}{\|z\|} \right) \right\| = \max_{z \neq 0} \frac{\|Az\|}{\|z\|}.$$

Corollary

$$\|Az\| \leq \|A\| \cdot \|z\|.$$

Theorem

If $A = [a_{ij}]$ is an $n \times n$ matrix, then

$$\|A\|_{\infty} = \max_{1 \leq i \leq n} \sum_{j=1}^n |a_{ij}|.$$



Proof: Let x be an n -dimension vector with

$$1 = \|x\|_\infty = \max_{1 \leq i \leq n} |x_i|.$$

Then

$$\begin{aligned} \|Ax\|_\infty &= \max_{1 \leq i \leq n} \left| \sum_{j=1}^n a_{ij} x_j \right| \\ &\leq \max_{1 \leq i \leq n} \sum_{j=1}^n |a_{ij}| \max_{1 \leq j \leq n} |x_j| = \max_{1 \leq i \leq n} \sum_{j=1}^n |a_{ij}|. \end{aligned}$$

Consequently,

$$\|A\|_\infty = \max_{\|x\|_\infty=1} \|Ax\|_\infty \leq \max_{1 \leq i \leq n} \sum_{j=1}^n |a_{ij}|.$$

On the other hand, let p be an integer with

$$\sum_{j=1}^n |a_{pj}| = \max_{1 \leq i \leq n} \sum_{j=1}^n |a_{ij}|,$$



and x be the vector with

$$x_j = \begin{cases} 1, & \text{if } a_{pj} \geq 0, \\ -1, & \text{if } a_{pj} < 0. \end{cases}$$

Then

$$\|x\|_\infty = 1 \quad \text{and} \quad a_{pj}x_j = |a_{pj}|, \quad \forall j = 1, 2, \dots, n,$$

so

$$\|Ax\|_\infty = \max_{1 \leq i \leq n} \left| \sum_{j=1}^n a_{ij}x_j \right| \geq \left| \sum_{j=1}^n a_{pj}x_j \right| = \left| \sum_{j=1}^n |a_{pj}| \right| = \max_{1 \leq i \leq n} \sum_{j=1}^n |a_{ij}|.$$

This result implies that

$$\|A\|_\infty = \max_{\|x\|_\infty=1} \|Ax\|_\infty \geq \max_{1 \leq i \leq n} \sum_{j=1}^n |a_{ij}|.$$

which gives

$$\|A\|_\infty = \max_{1 \leq i \leq n} \sum_{j=1}^n |a_{ij}|. \quad \square$$



Eigenvalues and eigenvectors

Definition (Characteristic polynomial)

If A is a square matrix, the characteristic polynomial of A is defined by

$$p(\lambda) = \det(A - \lambda I).$$

Definition (Eigenvalue and eigenvector)

If p is the characteristic polynomial of the matrix A , the zeros of p are eigenvalues of the matrix A . If λ is an eigenvalue of A and $x \neq 0$ satisfies $(A - \lambda I)x = 0$, then x is an eigenvector of A corresponding to the eigenvalue λ .

Definition (Spectrum and Spectral Radius)

The set of all eigenvalues of a matrix A is called the spectrum of A . The spectral radius of A is

$$\rho(A) = \max\{|\lambda|; \lambda \text{ is an eigenvalue of } A\}.$$

Theorem

If A is an $n \times n$ matrix, then

- (i) $\|A\|_2 = \sqrt{\rho(A^T A)}$;
- (ii) $\rho(A) \leq \|A\|$ for any matrix norm.

Proof: Proof for the second part. Suppose λ is an eigenvalue of A and $x \neq 0$ is a corresponding eigenvector such that $Ax = \lambda x$ and $\|x\| = 1$. Then

$$|\lambda| = |\lambda|\|x\| = \|\lambda x\| = \|Ax\| \leq \|A\|\|x\| = \|A\|,$$

that is, $|\lambda| \leq \|A\|$. Since λ is arbitrary, this implies that $\rho(A) = \max |\lambda| \leq \|A\|$. □

Theorem

For any A and any $\varepsilon > 0$, there exists a matrix norm $\|\cdot\|$ such that

$$\rho(A) < \|A\| < \rho(A) + \varepsilon.$$

Definition

We call an $n \times n$ matrix A convergent if

$$\lim_{k \rightarrow \infty} (A^k)_{ij} = 0 \quad \forall i = 1, 2, \dots, n \quad \text{and} \quad j = 1, 2, \dots, n.$$

Theorem

The following statements are equivalent.

- 1 A is a *convergent matrix*;
- 2 $\lim_{k \rightarrow \infty} \|A^k\| = 0$ for *some* matrix norm;
- 3 $\lim_{k \rightarrow \infty} \|A^k\| = 0$ for *all* matrix norm;
- 4 $\rho(A) < 1$;
- 5 $\lim_{k \rightarrow \infty} A^k x = 0$ for *any* x .



Iterative techniques for solving linear systems

- For small dimension of linear systems, it requires for direct techniques.
- For large systems, iterative techniques are efficient in terms of both computer storage and computation.

The basic idea of iterative techniques is to split the coefficient matrix A into

$$A = M - (M - A),$$

for some matrix M , which is called the **splitting matrix**. Here we assume that A and M are both **nonsingular**. Then the original problem is rewritten in the equivalent form

$$Mx = (M - A)x + b.$$

This suggests an iterative process

$$x^{(k)} = (I - M^{-1}A)x^{(k-1)} + M^{-1}b \equiv Tx^{(k-1)} + c,$$

where T is usually called the **iteration matrix**. The initial vector $x^{(0)}$ can be arbitrary or be chosen according to certain conditions.



Two criteria for choosing the splitting matrix M are

- $x^{(k)}$ is easily computed. More precisely, the system $Mx^{(k)} = y$ is easy to solve;
- the sequence $\{x^{(k)}\}$ converges rapidly to the exact solution.

Note that one way to achieve the second goal is to choose M so that M^{-1} approximate A^{-1} ,

In the following subsections, we will introduce some of the mostly commonly used classic iterative methods.



Jacobi Method

If we decompose the coefficient matrix A as

$$A = L + D + U,$$

where D is the **diagonal part**, L is the **strictly lower triangular part**, and U is the **strictly upper triangular part**, of A , and choose $M = D$, then we derive the iterative formulation for **Jacobi method**:

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

With this method, the iteration matrix $T_J = -D^{-1}(L + U)$ and $c = D^{-1}b$. Each component $x_i^{(k)}$ can be computed by

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



$$\begin{aligned}
 a_{11}x_1^{(k)} + a_{12}x_2^{(k-1)} + a_{13}x_3^{(k-1)} + \cdots + a_{1n}x_n^{(k-1)} &= b_1 \\
 a_{21}x_1^{(k-1)} + a_{22}x_2^{(k)} + a_{23}x_3^{(k-1)} + \cdots + a_{2n}x_n^{(k-1)} &= b_2 \\
 &\vdots \\
 a_{n1}x_1^{(k-1)} + a_{n2}x_2^{(k-1)} + a_{n3}x_3^{(k-1)} + \cdots + a_{nn}x_n^{(k)} &= b_n.
 \end{aligned}$$

Algorithm (Jacobi Method)

Given $x^{(0)}$, tolerance TOL , maximum number of iteration M .

Set $k = 1$.

While $k \leq M$ and $\|x - x^{(0)}\|_2 \geq TOL$

Set $k = k + 1$, $x^{(0)} = x$.

For $i = 1, 2, \dots, n$

$$x_i = \left(b_i - \sum_{j=1}^{i-1} a_{ij}x_j^{(0)} - \sum_{j=i+1}^n a_{ij}x_j^{(0)} \right) / a_{ii}$$

End For

End While

Example

Consider the linear system $Ax = b$ given by

$$\begin{aligned}E_1: & 10x_1 - x_2 + 2x_3 & = & 6, \\E_2: & -x_1 + 11x_2 - x_3 + 3x_4 & = & 25, \\E_3: & 2x_1 - x_2 + 10x_3 - x_4 & = & -11, \\E_4: & & 3x_2 - x_3 + 8x_4 & = & 15\end{aligned}$$

which has the unique solution $x = [1, 2, -1, 1]^T$.

Solving equation E_i for x_i , for $i = 1, 2, 3, 4$, we obtain

$$\begin{aligned}x_1 & = & & & 1/10x_2 & - & 1/5x_3 & & + & 3/5, \\x_2 & = & 1/11x_1 & & & + & 1/11x_3 & - & 3/11x_4 & + & 25/11, \\x_3 & = & -1/5x_1 & + & 1/10x_2 & & & + & 1/10x_4 & - & 11/10, \\x_4 & = & & - & 3/8x_2 & + & 1/8x_3 & & + & 15/8.\end{aligned}$$



Then $Ax = b$ can be rewritten in the form $x = Tx + c$ with

$$T = \begin{bmatrix} 0 & 1/10 & -1/5 & 0 \\ 1/11 & 0 & 1/11 & -3/11 \\ -1/5 & 1/10 & 0 & 1/10 \\ 0 & -3/8 & 1/8 & 0 \end{bmatrix} \quad \text{and} \quad c = \begin{bmatrix} 3/5 \\ 25/11 \\ -11/10 \\ 15/8 \end{bmatrix}$$

and the iterative formulation for Jacobi method is

$$x^{(k)} = Tx^{(k-1)} + c \quad \text{for } k = 1, 2, \dots$$

The numerical results of such iteration is list as follows:



k	x_1	x_2	x_3	x_4
0	0.0000	0.0000	0.0000	0.0000
1	0.6000	2.2727	-1.1000	1.8750
2	1.0473	1.7159	-0.8052	0.8852
3	0.9326	2.0533	-1.0493	1.1309
4	1.0152	1.9537	-0.9681	0.9738
5	0.9890	2.0114	-1.0103	1.0214
6	1.0032	1.9922	-0.9945	0.9944
7	0.9981	2.0023	-1.0020	1.0036
8	1.0006	1.9987	-0.9990	0.9989
9	0.9997	2.0004	-1.0004	1.0006
10	1.0001	1.9998	-0.9998	0.9998



Matlab code of Example

```
clear all; delete rslt.dat; diary rslt.dat; diary on;
n = 4; xold = zeros(n,1); xnew = zeros(n,1); T = zeros(n,n);
T(1,2) = 1/10; T(1,3) = -1/5; T(2,1) = 1/11;
T(2,3) = 1/11; T(2,4) = -3/11; T(3,1) = -1/5;
T(3,2) = 1/10; T(3,4) = 1/10; T(4,2) = -3/8; T(4,3) = 1/8;
c(1,1) = 3/5; c(2,1) = 25/11; c(3,1) = -11/10; c(4,1) = 15/8;
xnew = T * xold + c; k = 0;
fprintf(' k      x1      x2      x3      x4      \n');
while ( k <= 100 & norm(xnew-xold) > 1.0d-14 )
    xold = xnew; xnew = T * xold + c; k = k + 1;
    fprintf('%3.0f ',k);
    for jj = 1:n
        fprintf('%5.4f ',xold(jj));
    end
    fprintf('\n');
end
```

Gauss-Seidel Method

When computing $x_i^{(k)}$ for $i > 1$, $x_1^{(k)}, \dots, x_{i-1}^{(k)}$ have already been computed and are likely to be better approximations to the exact x_1, \dots, x_{i-1} than $x_1^{(k-1)}, \dots, x_{i-1}^{(k-1)}$. It seems reasonable to compute $x_i^{(k)}$ using these most recently computed values. That is

$$\begin{aligned} a_{11}x_1^{(k)} + a_{12}x_2^{(k-1)} + a_{13}x_3^{(k-1)} + \cdots + a_{1n}x_n^{(k-1)} &= b_1 \\ a_{21}x_1^{(k)} + a_{22}x_2^{(k)} + a_{23}x_3^{(k-1)} + \cdots + a_{2n}x_n^{(k-1)} &= b_2 \\ a_{31}x_1^{(k)} + a_{32}x_2^{(k)} + a_{33}x_3^{(k)} + \cdots + a_{3n}x_n^{(k-1)} &= b_3 \\ &\vdots \\ a_{n1}x_1^{(k-1)} + a_{n2}x_2^{(k-1)} + a_{n3}x_3^{(k-1)} + \cdots + a_{nn}x_n^{(k)} &= b_n. \end{aligned}$$

This improvement induce the Gauss-Seidel method.

The Gauss-Seidel method sets $M = D + L$ and defines the iteration as

$$x^{(k)} = -(D + L)^{-1}Ux^{(k-1)} + (D + L)^{-1}b.$$



That is, Gauss-Seidel method uses $T_G = -(D + L)^{-1}U$ as the iteration matrix. The formulation above can be rewritten as

$$x^{(k)} = -D^{-1} \left(Lx^{(k)} + Ux^{(k-1)} - b \right).$$

Hence each component $x_i^{(k)}$ can be computed by

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

- For Jacobi method, only the components of $x^{(k-1)}$ are used to compute $x^{(k)}$. Hence $x_i^{(k)}, i = 1, \dots, n$, can be computed in parallel at each iteration k .
- At each iteration of Gauss-Seidel method, since $x_i^{(k)}$ can not be computed until $x_1^{(k)}, \dots, x_{i-1}^{(k)}$ are available, the method is not a parallel algorithm in nature.



Algorithm (Gauss-Seidel Method)

Given $x^{(0)}$, tolerance TOL , maximum number of iteration M .

Set $k = 1$.

For $i = 1, 2, \dots, n$

$$x_i = \left(b_i - \sum_{j=1}^{i-1} a_{ij}x_j - \sum_{j=i+1}^n a_{ij}x_j^{(0)} \right) / a_{ii}$$

End For

While $k \leq M$ and $\|x - x^{(0)}\|_2 \geq TOL$

Set $k = k + 1$, $x^{(0)} = x$.

For $i = 1, 2, \dots, n$

$$x_i = \left(b_i - \sum_{j=1}^{i-1} a_{ij}x_j - \sum_{j=i+1}^n a_{ij}x_j^{(0)} \right) / a_{ii}$$

End For

End While



Example

Consider the linear system $Ax = b$ given by

$$\begin{aligned}E_1: & 10x_1 - x_2 + 2x_3 & = & 6, \\E_2: & -x_1 + 11x_2 - x_3 + 3x_4 & = & 25, \\E_3: & 2x_1 - x_2 + 10x_3 - x_4 & = & -11, \\E_4: & 3x_2 - x_3 + 8x_4 & = & 15\end{aligned}$$

which has the unique solution $x = [1, 2, -1, 1]^T$.

Gauss-Seidel method gives the equation

$$\begin{aligned}x_1^{(k)} &= \frac{1}{10}x_2^{(k-1)} - \frac{1}{5}x_3^{(k-1)} + \frac{3}{5}, \\x_2^{(k)} &= \frac{1}{11}x_1^{(k)} + \frac{1}{11}x_3^{(k-1)} - \frac{3}{11}x_4^{(k-1)} + \frac{25}{11}, \\x_3^{(k)} &= -\frac{1}{5}x_1^{(k)} + \frac{1}{10}x_2^{(k)} + \frac{1}{10}x_4^{(k-1)} - \frac{11}{10}, \\x_4^{(k)} &= -\frac{3}{8}x_2^{(k)} + \frac{1}{8}x_3^{(k)} + \frac{15}{8}.\end{aligned}$$



The numerical results of such iteration is list as follows:

k	x_1	x_2	x_3	x_4
0	0.0000	0.0000	0.0000	0.0000
1	0.6000	2.3273	-0.9873	0.8789
2	1.0302	2.0369	-1.0145	0.9843
3	1.0066	2.0036	-1.0025	0.9984
4	1.0009	2.0003	-1.0003	0.9998
5	1.0001	2.0000	-1.0000	1.0000

- The results of Example appear to imply that the Gauss-Seidel method is superior to the Jacobi method.
- This is almost always true, but there are linear systems for which the Jacobi method converges and the Gauss-Seidel method does not.
- See Exercises 17 and 18.



Matlab code of Example

```
clear all; delete rslt.dat; diary rslt.dat; diary on;
n = 4; xold = zeros(n,1); xnew = zeros(n,1); A = zeros(n,n);
A(1,1)=10; A(1,2)=-1; A(1,3)=2; A(2,1)=-1; A(2,2)=11; A(2,3)=-1; A(2,4)=3; A(3,1)=2; A(3,2)=-1;
A(3,3)=10; A(3,4)=-1; A(4,2)=3; A(4,3)=-1; A(4,4)=8; b(1)=6; b(2)=25; b(3)=-11; b(4)=15;
for ii = 1:n
    xnew(ii) = b(ii);
    for jj = 1:ii-1
        xnew(ii) = xnew(ii) - A(ii,jj) * xnew(jj);
    end
    for jj = ii+1:n
        xnew(ii) = xnew(ii) - A(ii,jj) * xold(jj);
    end
    xnew(ii) = xnew(ii) / A(ii,ii);
end
k = 0; fprintf(' k      x1      x2      x3      x4      \n');
while ( k <= 100 & norm(xnew-xold) > 1.0d-14 )
    xold = xnew; k = k + 1;
    for ii = 1:n
        xnew(ii) = b(ii);
        for jj = 1:ii-1
            xnew(ii) = xnew(ii) - A(ii,jj) * xnew(jj);
        end
        for jj = ii+1:n
            xnew(ii) = xnew(ii) - A(ii,jj) * xold(jj);
        end
        xnew(ii) = xnew(ii) / A(ii,ii);
    end
    fprintf('%3.0f ',k);
    for jj = 1:n
        fprintf('%5.4f ',xold(jj));
    end
    fprintf('\n');
end
diary off
```

Lemma (20)

If $\rho(T) < 1$, then $(I - T)^{-1}$ exists and

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

Proof: Let λ be an eigenvalue of T , then $1 - \lambda$ is an eigenvalue of $I - T$. But $|\lambda| \leq \rho(A) < 1$, so $1 - \lambda \neq 0$ and 0 is not an eigenvalue of $I - T$, which means $(I - T)$ is nonsingular.

Next we show that $(I - T)^{-1} = I + T + T^2 + \dots$. Since

$$(I - T) \left(\sum_{i=0}^m T^i \right) = I - T^{m+1},$$

and $\rho(T) < 1$ implies $\|T^m\| \rightarrow 0$ as $m \rightarrow \infty$, we have

$$(I - T) \left(\lim_{m \rightarrow \infty} \sum_{i=0}^m T^i \right) = (I - T) \left(\sum_{i=0}^{\infty} T^i \right) = I. \quad \square$$



Theorem

For *any* $x^{(0)} \in \mathbb{R}^n$, the sequence produced by

$$x^{(k)} = Tx^{(k-1)} + c, \quad k = 1, 2, \dots,$$

converges to the *unique* solution of $x = Tx + c$ if and only if

$$\rho(T) < 1.$$

Proof: Suppose $\rho(T) < 1$. The sequence of vectors $x^{(k)}$ produced by the iterative formulation are

$$x^{(1)} = Tx^{(0)} + c$$

$$x^{(2)} = Tx^{(1)} + c = T^2x^{(0)} + (T + I)c$$

$$x^{(3)} = Tx^{(2)} + c = T^3x^{(0)} + (T^2 + T + I)c$$

\vdots

In general

$$x^{(k)} = T^k x^{(0)} + (T^{k-1} + T^{k-2} + \dots + T + I)c.$$



Since $\rho(T) < 1$, $\lim_{k \rightarrow \infty} T^k x^{(0)} = 0$ for any $x^{(0)} \in \mathbb{R}^n$. By Lemma 20,

$$(T^{k-1} + T^{k-2} + \cdots + T + I)c \rightarrow (I - T)^{-1}c, \quad \text{as } k \rightarrow \infty.$$

Therefore

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

Conversely, suppose $\{x^{(k)}\} \rightarrow x = (I - T)^{-1}c$. Since

$$\begin{aligned} x - x^{(k)} &= Tx + c - Tx^{(k-1)} - c = T(x - x^{(k-1)}) = T^2(x - x^{(k-2)}) \\ &= \cdots = T^k(x - x^{(0)}). \end{aligned}$$

Let $z = x - x^{(0)}$. Then

$$\lim_{k \rightarrow \infty} T^k z = \lim_{k \rightarrow \infty} (x - x^{(k)}) = 0.$$

It follows from theorem $\rho(T) < 1$.



Theorem

If $\|T\| < 1$, then the sequence $x^{(k)}$ converges to x for any initial $x^{(0)}$ and

$$\textcircled{1} \quad \|x - x^{(k)}\| \leq \|T\|^k \|x - x^{(0)}\|$$

$$\textcircled{2} \quad \|x - x^{(k)}\| \leq \frac{\|T\|^k}{1 - \|T\|} \|x^{(1)} - x^{(0)}\|.$$

Proof: Since $x = Tx + c$ and $x^{(k)} = Tx^{(k-1)} + c$,

$$\begin{aligned} x - x^{(k)} &= Tx + c - Tx^{(k-1)} - c \\ &= T(x - x^{(k-1)}) \\ &= T^2(x - x^{(k-2)}) = \dots = T^k(x - x^{(0)}). \end{aligned}$$

The first statement can then be derived

$$\|x - x^{(k)}\| = \|T^k(x - x^{(0)})\| \leq \|T\|^k \|x - x^{(0)}\|.$$

For the second result, we first show that

$$\|x^{(n)} - x^{(n-1)}\| \leq \|T\|^{n-1} \|x^{(1)} - x^{(0)}\| \quad \text{for any } n \geq 1.$$



Since

$$\begin{aligned}x^{(n)} - x^{(n-1)} &= Tx^{(n-1)} + c - Tx^{(n-2)} - c \\&= T(x^{(n-1)} - x^{(n-2)}) \\&= T^2(x^{(n-2)} - x^{(n-3)}) = \dots = T^{n-1}(x^{(1)} - x^{(0)}),\end{aligned}$$

we have

$$\|x^{(n)} - x^{(n-1)}\| \leq \|T\|^{n-1} \|x^{(1)} - x^{(0)}\|.$$

Let $m \geq k$,

$$\begin{aligned}&x^{(m)} - x^{(k)} \\&= (x^{(m)} - x^{(m-1)}) + (x^{(m-1)} - x^{(m-2)}) + \dots + (x^{(k+1)} - x^{(k)}) \\&= T^{m-1}(x^{(1)} - x^{(0)}) + T^{m-2}(x^{(1)} - x^{(0)}) + \dots + T^k(x^{(1)} - x^{(0)}) \\&= (T^{m-1} + T^{m-2} + \dots + T^k)(x^{(1)} - x^{(0)}),\end{aligned}$$



hence

$$\begin{aligned} & \|x^{(m)} - x^{(k)}\| \\ \leq & \left(\|T\|^{m-1} + \|T\|^{m-2} + \dots + \|T\|^k \right) \|x^{(1)} - x^{(0)}\| \\ = & \|T\|^k \left(\|T\|^{m-k-1} + \|T\|^{m-k-2} + \dots + 1 \right) \|x^{(1)} - x^{(0)}\|. \end{aligned}$$

Since $\lim_{m \rightarrow \infty} x^{(m)} = x$,

$$\begin{aligned} & \|x - x^{(k)}\| \\ = & \lim_{m \rightarrow \infty} \|x^{(m)} - x^{(k)}\| \\ \leq & \lim_{m \rightarrow \infty} \|T\|^k \left(\|T\|^{m-k-1} + \|T\|^{m-k-2} + \dots + 1 \right) \|x^{(1)} - x^{(0)}\| \\ = & \|T\|^k \|x^{(1)} - x^{(0)}\| \lim_{m \rightarrow \infty} \left(\|T\|^{m-k-1} + \|T\|^{m-k-2} + \dots + 1 \right) \\ = & \|T\|^k \frac{1}{1 - \|T\|} \|x^{(1)} - x^{(0)}\|. \end{aligned}$$

This proves the second result.



Theorem

If A is *strictly diagonal dominant*, then both the *Jacobi* and *Gauss-Seidel* methods *converges* for *any* initial vector $x^{(0)}$.

Proof: By assumption, A is strictly diagonal dominant, hence $a_{ii} \neq 0$ (otherwise A is singular) and

$$|a_{ii}| > \sum_{j=1, j \neq i}^n |a_{ij}|, \quad i = 1, 2, \dots, n.$$

For Jacobi method, the iteration matrix $T_J = -D^{-1}(L + U)$ has entries

$$[T_J]_{ij} = \begin{cases} -\frac{a_{ij}}{a_{ii}}, & i \neq j, \\ 0, & i = j. \end{cases}$$

Hence

$$\|T_J\|_{\infty} = \max_{1 \leq i \leq n} \sum_{j=1, j \neq i}^n \left| \frac{a_{ij}}{a_{ii}} \right| = \max_{1 \leq i \leq n} \frac{1}{|a_{ii}|} \sum_{j=1, j \neq i}^n |a_{ij}| < 1,$$

and this implies that the Jacobi method converges.



For Gauss-Seidel method, the iteration matrix $T_G = -(D + L)^{-1}U$. Let λ be any eigenvalue of T_G and y , $\|y\|_\infty = 1$, is a corresponding eigenvector. Thus

$$T_G y = \lambda y \quad \implies \quad -Uy = \lambda(D + L)y.$$

Hence for $i = 1, \dots, n$,

$$-\sum_{j=i+1}^n a_{ij}y_j = \lambda a_{ii}y_i + \lambda \sum_{j=1}^{i-1} a_{ij}y_j.$$

This gives

$$\lambda a_{ii}y_i = -\lambda \sum_{j=1}^{i-1} a_{ij}y_j - \sum_{j=i+1}^n a_{ij}y_j$$

and

$$|\lambda| |a_{ii}| |y_i| \leq |\lambda| \sum_{j=1}^{i-1} |a_{ij}| |y_j| + \sum_{j=i+1}^n |a_{ij}| |y_j|.$$

Choose the index k such that $|y_k| = 1 \geq |y_j|$ (this index can always be found since $\|y\|_\infty = 1$). Then



Successive over-relaxation (SOR) method

Definition

Suppose $\tilde{x} \in \mathbb{R}^n$ is an approximated solution of $Ax = b$. The **residual vector** r for \tilde{x} is $r = b - A\tilde{x}$.

Let the approximate solution $\mathbf{x}^{(k,i)}$ produced by Gauss-Seidel method be defined by

$$\mathbf{x}^{(k,i)} = \left[x_1^{(k)}, \dots, x_{i-1}^{(k)}, x_i^{(k-1)}, \dots, x_n^{(k-1)} \right]^T$$

and

$$r_i^{(k)} = \left[r_{1i}^{(k)}, r_{2i}^{(k)}, \dots, r_{ni}^{(k)} \right]^T = b - A\mathbf{x}^{(k,i)}$$

be the corresponding residual vector. Then the m th component of $r_i^{(k)}$ is

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



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 $m = 1, 2, \dots, n$.

In particular, the i th component of $r_i^{(k)}$ 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

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



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

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

Relaxation method is modified the Gauss-Seidel procedure to

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

for certain choices of positive ω such that the norm of the residual vector is reduced and the convergence is significantly faster.



These methods are called for

$\omega < 1$: under relaxation,

$\omega = 1$: Gauss-Seidel method,

$\omega > 1$: over relaxation.

Over-relaxation methods are called **SOR (Successive over-relaxation)**. To determine the matrix of the SOR method, we rewrite (1) 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 if $A = L + D + U$, then we have

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

or

$$\begin{aligned} x^{(k)} &= (D + \omega L)^{-1} [(1 - \omega)D - \omega U]x^{(k-1)} + \omega(D + \omega L)^{-1}b \\ &\equiv T_\omega x^{(k-1)} + c_\omega. \end{aligned}$$



Example

The linear system $Ax = b$ given by

$$\begin{aligned}4x_1 + 3x_2 &= 24, \\3x_1 + 4x_2 - x_3 &= 30, \\-x_2 + 4x_3 &= -24,\end{aligned}$$

has the solution $[3, 4, -5]^T$.

- Numerical results of Gauss-Seidel method with $x^{(0)} = [1, 1, 1]^T$:

k	x_1	x_2	x_3
0	1.0000000	1.0000000	1.0000000
1	5.2500000	3.8125000	-5.0468750
2	3.1406250	3.8828125	-5.0292969
3	3.0878906	3.9267578	-5.0183105
4	3.0549316	3.9542236	-5.0114441
5	3.0343323	3.9713898	-5.0071526
6	3.0214577	3.9821186	-5.0044703
7	3.0134110	3.9888241	-5.0027940



- Numerical results of SOR method with $\omega = 1.25$ and $x^{(0)} = [1, 1, 1]^T$:

k	x_1	x_2	x_3
0	1.0000000	1.0000000	1.0000000
1	6.3125000	3.5195313	-6.6501465
2	2.6223145	3.9585266	-4.6004238
3	3.1333027	4.0102646	-5.0966863
4	2.9570512	4.0074838	-4.9734897
5	3.0037211	4.0029250	-5.0057135
6	2.9963276	4.0009262	-4.9982822
7	3.0000498	4.0002586	-5.0003486



- Numerical results of SOR method with $\omega = 1.6$ and $x^{(0)} = [1, 1, 1]^T$:

k	x_1	x_2	x_3
0	1.0000000	1.0000000	1.0000000
1	7.8000000	2.4400000	-9.2240000
2	1.9920000	4.4560000	-2.2832000
3	3.0576000	4.7440000	-6.3324800
4	2.0726400	4.1334400	-4.1471360
5	3.3962880	3.7855360	-5.5975040
6	3.0195840	3.8661760	-4.6950272
7	3.1488384	4.0236774	-5.1735127



Matlab code of SOR

```
clear all; delete rslt.dat; diary rslt.dat; diary on;
n = 3; xold = zeros(n,1); xnew = zeros(n,1); A = zeros(n,n); DL = zeros(n,n); DU = zeros(n,n);
A(1,1)=4; A(1,2)=3; A(2,1)=3; A(2,2)=4; A(2,3)=-1; A(3,2)=-1; A(3,3)=4;
b(1,1)=24; b(2,1)=30; b(3,1)=-24; omega=1.25;
for ii = 1:n
    DL(ii,ii) = A(ii,ii);
    for jj = 1:ii-1
        DL(ii,jj) = omega * A(ii,jj);
    end
    DU(ii,ii) = (1-omega)*A(ii,ii);
    for jj = ii+1:n
        DU(ii,jj) = - omega * A(ii,jj);
    end
end
c = omega * (DL \ b); xnew = DL \ ( DU * xold ) + c;
k = 0; fprintf(' k      x1      x2      x3      \n');
while ( k <= 100 & norm(xnew-xold) > 1.0d-14 )
    xold = xnew; k = k + 1; xnew = DL \ ( DU * xold ) + c;
    fprintf('%3.0f ',k);
    for jj = 1:n
        fprintf('%5.4f ',xold(jj));
    end
    fprintf('\n');
end
diary off
```



Theorem (Kahan)

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

Theorem (Ostrowski-Reich)

If A is *positive definite* and the relaxation parameter ω satisfying $0 < \omega < 2$, then the SOR iteration *converges* for *any* initial vector $x^{(0)}$.

Theorem

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 iteration is

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

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

Example

The matrix

$$A = \begin{bmatrix} 4 & 3 & 0 \\ 3 & 4 & -1 \\ 0 & -1 & 4 \end{bmatrix},$$

given in previous example, is positive definite and tridiagonal.

Since

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



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

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

This explains the rapid convergence obtained in previous example when using $\omega = 0.125$



Symmetric Successive Over Relaxation (SSOR) Method

Let A be symmetric and $A = D + L + L^T$. The idea is in fact to implement the SOR formulation **twice**, **one forward** and **one backward**, at each iteration. That is, SSOR method defines

$$(D + \omega L)x^{(k-\frac{1}{2})} = [(1 - \omega)D - \omega L^T]x^{(k-1)} + \omega b, \quad (2)$$

$$(D + \omega L^T)x^{(k)} = [(1 - \omega)D - \omega L]x^{(k-\frac{1}{2})} + \omega b. \quad (3)$$

Define

$$\begin{cases} M_\omega: = D + \omega L, \\ N_\omega: = (1 - \omega)D - \omega L^T. \end{cases}$$

Then from the iterations (2) and (3), it follows that

$$\begin{aligned} x^{(k)} &= (M_\omega^{-T} N_\omega^T M_\omega^{-1} N_\omega) x^{(k-1)} + \omega (M_\omega^{-T} N_\omega^T M_\omega^{-1} + M_\omega^{-T}) b \\ &\equiv T(\omega)x^{(k-1)} + M(\omega)^{-1}b. \end{aligned}$$



But

$$\begin{aligned} & ((1 - \omega)D - \omega L)(D + \omega L)^{-1} + I \\ &= (-\omega L - D - \omega D + 2D)(D + \omega L)^{-1} + I \\ &= -I + (2 - \omega)D(D + \omega L)^{-1} + I \\ &= (2 - \omega)D(D + \omega L)^{-1}. \end{aligned}$$

Thus

$$M(\omega)^{-1} = \omega (D + \omega L^T)^{-1} (2 - \omega)D(D + \omega L)^{-1},$$

then the splitting matrix is

$$M(\omega) = \frac{1}{\omega(2 - \omega)} (D + \omega L)D^{-1} (D + \omega L^T).$$

The iteration matrix is

$$T(\omega) = (D + \omega L^T)^{-1} [(1 - \omega)D - \omega L](D + \omega L)^{-1} [(1 - \omega)D - \omega L^T]$$



Error bounds and iterative refinement

Example

The linear system $Ax = b$ given by

$$\begin{bmatrix} 1 & 2 \\ 1.0001 & 2 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \end{bmatrix} = \begin{bmatrix} 3 \\ 3.0001 \end{bmatrix}$$

has the unique solution $x = [1, 1]^T$.

The poor approximation $\tilde{x} = [3, 0]^T$ has the residual vector

$$r = b - A\tilde{x} = \begin{bmatrix} 3 \\ 3.0001 \end{bmatrix} - \begin{bmatrix} 1 & 2 \\ 1.0001 & 2 \end{bmatrix} \begin{bmatrix} 3 \\ 0 \end{bmatrix} = \begin{bmatrix} 0 \\ -0.0002 \end{bmatrix},$$

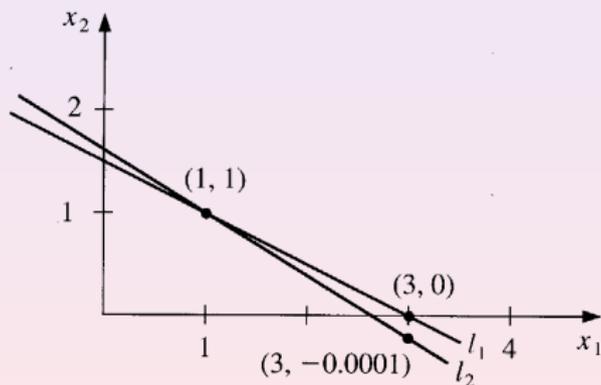
so $\|r\|_\infty = 0.0002$. Although the norm of the residual vector is small, the approximation $\tilde{x} = [3, 0]^T$ is obviously quite poor; in fact, $\|x - \tilde{x}\|_\infty = 2$.



The solution of above example represents the intersection of the lines

$$l_1 : x_1 + 2x_2 = 3 \quad \text{and} \quad l_2 : 1.0001x_1 + 2x_2 = 3.0001.$$

l_1 and l_2 are nearly parallel. The point $(3, 0)$ lies on l_1 which implies that $(3, 0)$ also lies close to l_2 , even though it differs significantly from the intersection point $(1, 1)$.



Theorem

Suppose that \tilde{x} is an approximate solution of $Ax = b$, A is nonsingular matrix and $r = b - A\tilde{x}$. Then

$$\|x - \tilde{x}\| \leq \|r\| \cdot \|A^{-1}\|$$

and if $x \neq 0$ and $b \neq 0$,

$$\frac{\|x - \tilde{x}\|}{\|x\|} \leq \|A\| \cdot \|A^{-1}\| \frac{\|r\|}{\|b\|}.$$

Proof: Since

$$r = b - A\tilde{x} = Ax - A\tilde{x} = A(x - \tilde{x})$$

and A is nonsingular, we have

$$\|x - \tilde{x}\| = \|A^{-1}r\| \leq \|A^{-1}\| \cdot \|r\|. \quad (4)$$

Moreover, since $b = Ax$, we have

$$\|b\| \leq \|A\| \cdot \|x\|.$$



It implies that

$$\frac{1}{\|x\|} \leq \frac{\|A\|}{\|b\|}. \quad (5)$$

Combining Equations (4) and (5), we have

$$\frac{\|x - \tilde{x}\|}{\|x\|} \leq \frac{\|A\| \cdot \|A^{-1}\|}{\|b\|} \|r\|.$$



Definition (Condition number)

The condition number of nonsingular matrix A is

$$\kappa(A) = \|A\| \cdot \|A^{-1}\|.$$

For any nonsingular matrix A ,

$$1 = \|I\| = \|A \cdot A^{-1}\| \leq \|A\| \cdot \|A^{-1}\| = \kappa(A).$$



Definition

A matrix A is **well-conditioned** if $\kappa(A)$ is close to **1**, and is **ill-conditioned** when $\kappa(A)$ is significantly greater than 1.

In previous example,

$$A = \begin{bmatrix} 1 & 2 \\ 1.0001 & 2 \end{bmatrix}.$$

Since

$$A^{-1} = \begin{bmatrix} -10000 & 10000 \\ 5000.5 & -5000 \end{bmatrix},$$

we have

$$\kappa(A) = \|A\|_{\infty} \cdot \|A^{-1}\|_{\infty} = 3.0001 \times 20000 = 60002 \gg 1.$$



How to estimate the effective condition number in t -digit arithmetic without having to invert the matrix A ?

- If the approximate solution \tilde{x} of $Ax = b$ is being determined using t -digit arithmetic and Gaussian elimination, then

$$\|r\| = \|b - A\tilde{x}\| \approx 10^{-t}\|A\| \cdot \|\tilde{x}\|.$$

- All the arithmetic operations in Gaussian elimination technique are performed using t -digit arithmetic, but the residual vector r are done in double-precision (i.e., $2t$ -digit) arithmetic.
- Use the Gaussian elimination method which has already been calculated to solve

$$Ay = r.$$

Let \tilde{y} be the approximate solution. Then

$$\tilde{y} \approx A^{-1}r = A^{-1}(b - A\tilde{x}) = x - \tilde{x}$$

and

$$x \approx \tilde{x} + \tilde{y}.$$



Moreover,

$$\begin{aligned}\|\tilde{y}\| &\approx \|x - \tilde{x}\| = \|A^{-1}r\| \\ &\leq \|A^{-1}\| \cdot \|r\| \approx \|A^{-1}\|(10^{-t}\|A\| \cdot \|\tilde{x}\|) = 10^{-t}\|\tilde{x}\|\kappa(A).\end{aligned}$$

It implies that

$$\kappa(A) \approx \frac{\|\tilde{y}\|}{\|\tilde{x}\|} 10^t.$$

Iterative refinement

In general, $\tilde{x} + \tilde{y}$ is a more accurate approximation to the solution of $Ax = b$ than \tilde{x} .



Algorithm (Iterative refinement)

Given tolerance TOL , maximum number of iteration M , number of digits of precision t .

Solve $Ax = b$ by using Gaussian elimination in t -digit arithmetic.

Set $k = 1$

while ($k \leq M$)

 Compute $r = b - Ax$ in $2t$ -digit arithmetic.

 Solve $Ay = r$ by using Gaussian elimination in t -digit arithmetic.

 If $\|y\|_{\infty} < TOL$, then stop.

 Set $k = k + 1$ and $x = x + y$.

End while



Example

The linear system given by

$$\begin{bmatrix} 3.3330 & 15920 & -10.333 \\ 2.2220 & 16.710 & 9.6120 \\ 1.5611 & 5.1791 & 1.6852 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \\ x_3 \end{bmatrix} = \begin{bmatrix} 15913 \\ 28.544 \\ 8.4254 \end{bmatrix}$$

has the exact solution $x = [1, 1, 1]^T$.

Using Gaussian elimination and five-digit rounding arithmetic leads successively to the augmented matrices

$$\left[\begin{array}{ccc|c} 3.3330 & 15920 & -10.333 & 15913 \\ 0 & -10596 & 16.501 & -10580 \\ 0 & -7451.4 & 6.5250 & -7444.9 \end{array} \right]$$

and

$$\left[\begin{array}{ccc|c} 3.3330 & 15920 & -10.333 & 15913 \\ 0 & -10596 & 16.501 & -10580 \\ 0 & 0 & -5.0790 & -4.7000 \end{array} \right].$$



The approximate solution is

$$\tilde{x}^{(1)} = [1.2001, 0.99991, 0.92538]^T.$$

The residual vector corresponding to \tilde{x} is computed in double precision to be

$$\begin{aligned} r^{(1)} &= b - A\tilde{x}^{(1)} \\ &= \begin{bmatrix} 15913 \\ 28.544 \\ 8.4254 \end{bmatrix} - \begin{bmatrix} 3.3330 & 15920 & -10.333 \\ 2.2220 & 16.710 & 9.6120 \\ 1.5611 & 5.1791 & 1.6852 \end{bmatrix} \begin{bmatrix} 1.2001 \\ 0.99991 \\ 0.92538 \end{bmatrix} \\ &= \begin{bmatrix} 15913 \\ 28.544 \\ 8.4254 \end{bmatrix} - \begin{bmatrix} 15913.00518 \\ 28.26987086 \\ 8.611560367 \end{bmatrix} = \begin{bmatrix} -0.00518 \\ 0.27412914 \\ -0.186160367 \end{bmatrix}. \end{aligned}$$

Hence the solution of $Ay = r^{(1)}$ to be

$$\tilde{y}^{(1)} = [-0.20008, 8.9987 \times 10^{-5}, 0.074607]^T$$

and the new approximate solution $x^{(2)}$ is

$$x^{(2)} = x^{(1)} + \tilde{y}^{(1)} = [1.0000, 1.0000, 0.99999]^T.$$



Using the suggested stopping technique for the algorithm, we compute $r^{(2)} = b - A\tilde{x}^{(2)}$ and solve the system $Ay^{(2)} = r^{(2)}$, which gives

$$\tilde{y}^{(2)} = [1.5002 \times 10^{-9}, 2.0951 \times 10^{-10}, 1.0000 \times 10^{-5}]^T.$$

Since

$$\|\tilde{y}^{(2)}\|_{\infty} \leq 10^{-5},$$

we conclude that

$$\tilde{x}^{(3)} = \tilde{x}^{(2)} + \tilde{y}^{(2)} = [1.0000, 1.0000, 1.0000]^T$$

is sufficiently accurate. □

In the linear system

$$Ax = b,$$

A and b can be represented exactly. Realistically, the matrix A and vector b will be perturbed by δA and δb , respectively, causing the linear system

$$(A + \delta A)x = b + \delta b$$

to be solved in place of $Ax = b$.



Theorem

Suppose A is nonsingular and

$$\|\delta A\| < \frac{1}{\|A^{-1}\|}.$$

Then the solution \tilde{x} of $(A + \delta A)\tilde{x} = b + \delta b$ approximates the solution x of $Ax = b$ with the error estimate

$$\frac{\|x - \tilde{x}\|}{\|x\|} \leq \frac{\kappa(A)}{1 - \kappa(A)(\|\delta A\|/\|A\|)} \left(\frac{\|\delta b\|}{\|b\|} + \frac{\|\delta A\|}{\|A\|} \right).$$

- If A is well-conditioned, then small changes in A and b produce correspondingly small changes in the solution x .
- If A is ill-conditioned, then small changes in A and b may produce large changes in x .



The conjugate gradient method

Consider the linear systems

$$Ax = b$$

where A is large sparse and symmetric positive definite. Define the inner product notation

$$\langle x, y \rangle = x^T y \quad \text{for any } x, y \in \mathbb{R}^n.$$

Theorem

Let A be symmetric positive definite. Then x^ is the solution of $Ax = b$ if and only if x^* minimizes*

$$g(x) = \langle x, Ax \rangle - 2 \langle x, b \rangle .$$



Proof: (“ \Rightarrow ”) Rewrite $g(x)$ as

$$\begin{aligned}g(x) &= \langle x - x^*, A(x - x^*) \rangle + \langle x, Ax^* \rangle + \langle x^*, Ax \rangle \\ &\quad - \langle x^*, Ax^* \rangle - 2 \langle x, b \rangle \\ &= \langle x - x^*, A(x - x^*) \rangle - \langle x^*, Ax^* \rangle \\ &\quad + 2 \langle x, Ax^* \rangle - 2 \langle x, b \rangle \\ &= \langle x - x^*, A(x - x^*) \rangle - \langle x^*, Ax^* \rangle + 2 \langle x, Ax^* - b \rangle .\end{aligned}$$

Suppose that x^* is the solution of $Ax = b$, i.e., $Ax^* = b$. Then

$$g(x) = \langle x - x^*, A(x - x^*) \rangle - \langle x^*, Ax^* \rangle$$

which minimum occurs at $x = x^*$.



(“ \Leftarrow ”) Fixed vectors x and v , for any $\alpha \in \mathbb{R}$,

$$\begin{aligned} f(\alpha) &\equiv g(x + \alpha v) \\ &= \langle x + \alpha v, Ax + \alpha Av \rangle - 2 \langle x + \alpha v, b \rangle \\ &= \langle x, Ax \rangle + \alpha \langle v, Ax \rangle + \alpha \langle x, Av \rangle + \alpha^2 \langle v, Av \rangle \\ &\quad - 2 \langle x, b \rangle - 2\alpha \langle v, b \rangle \\ &= \langle x, Ax \rangle - 2 \langle x, b \rangle + 2\alpha \langle v, Ax \rangle - 2\alpha \langle v, b \rangle + \alpha^2 \langle v, Av \rangle \\ &= g(x) + 2\alpha \langle v, Ax - b \rangle + \alpha^2 \langle v, Av \rangle. \end{aligned}$$

Because f is a quadratic function of α and $\langle v, Av \rangle$ is positive, f has a minimal value when $f'(\alpha) = 0$. Since

$$f'(\alpha) = 2 \langle v, Ax - b \rangle + 2\alpha \langle v, Av \rangle,$$

the minimum occurs at

$$\hat{\alpha} = -\frac{\langle v, Ax - b \rangle}{\langle v, Av \rangle} = \frac{\langle v, b - Ax \rangle}{\langle v, Av \rangle}.$$



and

$$\begin{aligned}g(x + \hat{\alpha}v) &= f(\hat{\alpha}) = g(x) - 2 \frac{\langle v, b - Ax \rangle}{\langle v, Av \rangle} \langle v, b - Ax \rangle \\ &\quad + \left(\frac{\langle v, b - Ax \rangle}{\langle v, Av \rangle} \right)^2 \langle v, Av \rangle \\ &= g(x) - \frac{\langle v, b - Ax \rangle^2}{\langle v, Av \rangle}.\end{aligned}$$

So, for any nonzero vector v , we have

$$g(x + \hat{\alpha}v) < g(x) \quad \text{if} \quad \langle v, b - Ax \rangle \neq 0 \quad (6)$$

and

$$g(x + \hat{\alpha}v) = g(x) \quad \text{if} \quad \langle v, b - Ax \rangle = 0. \quad (7)$$

Suppose that x^* is a vector that minimizes g . Then

$$g(x^* + \hat{\alpha}v) \geq g(x^*) \quad \text{for any } v. \quad (8)$$



From (6), (7) and (8), we have

$$\langle v, b - Ax^* \rangle = 0 \text{ for any } v,$$

which implies that $Ax^* = b$. □

Let

$$r = b - Ax.$$

Then

$$\alpha = \frac{\langle v, b - Ax \rangle}{\langle v, Av \rangle} = \frac{\langle v, r \rangle}{\langle v, Av \rangle}.$$

If $r \neq 0$ and if v and r are not orthogonal, then

$$g(x + \alpha v) < g(x)$$

which implies that $x + \alpha v$ is closer to x^* than is x .



Let $x^{(0)}$ be an initial approximation to x^* and $v^{(1)} \neq 0$ be an initial search direction. For $k = 1, 2, 3, \dots$, we compute

$$\alpha_k = \frac{\langle v^{(k)}, b - Ax^{(k-1)} \rangle}{\langle v^{(k)}, Av^{(k)} \rangle},$$
$$x^{(k)} = x^{(k-1)} + \alpha_k v^{(k)}$$

and choose a new search direction $v^{(k+1)}$.

Question: How to choose $\{v^{(k)}\}$ such that $\{x^{(k)}\}$ converges rapidly to x^* ?

Let $\Phi : \mathbb{R}^n \rightarrow \mathbb{R}$ be a differential function on x . Then it holds

$$\frac{\Phi(x + \varepsilon p) - \Phi(x)}{\varepsilon} = \nabla \Phi(x)^T p + O(\varepsilon).$$

The right hand side takes minimum at

$$p = -\frac{\nabla \Phi(x)}{\|\nabla \Phi(x)\|} \quad (\text{i.e., the largest descent})$$

for all p with $\|p\| = 1$ (neglect $O(\varepsilon)$).



Denote $x = [x_1, x_2, \dots, x_n]^T$. Then

$$g(x) = \langle x, Ax \rangle - 2 \langle x, b \rangle = \sum_{i=1}^n \sum_{j=1}^n a_{ij} x_i x_j - 2 \sum_{i=1}^n x_i b_i.$$

It follows that

$$\frac{\partial g}{\partial x_k}(x) = 2 \sum_{i=1}^n a_{ki} x_i - 2b_k, \quad \text{for } k = 1, 2, \dots, n.$$

Therefore, the gradient of g is

$$\nabla g(x) = \left[\frac{\partial g}{\partial x_1}(x), \frac{\partial g}{\partial x_2}(x), \dots, \frac{\partial g}{\partial x_n}(x) \right]^T = 2(Ax - b) = -2r.$$



Steepest descent method (gradient method)

Given an initial $x_0 \neq 0$.

For $k = 1, 2, \dots$

$$r_{k-1} = b - Ax_{k-1}$$

If $r_{k-1} = 0$, then stop;

$$\text{else } \alpha_k = \frac{r_{k-1}^T r_{k-1}}{r_{k-1}^T A r_{k-1}}, \quad x_k = x_{k-1} + \alpha_k r_{k-1}.$$

End for

Theorem

If x_k, x_{k-1} are two approximations of the steepest descent method for solving $Ax = b$ and $\lambda_1 \geq \lambda_2 \geq \dots \geq \lambda_n > 0$ are the eigenvalues of A , then it holds:

$$\|x_k - x^*\|_A \leq \left(\frac{\lambda_1 - \lambda_n}{\lambda_1 + \lambda_n} \right) \|x_{k-1} - x^*\|_A,$$

where $\|x\|_A = \sqrt{x^T A x}$. Thus the gradient method is convergent.

- If the condition number of A ($= \lambda_1/\lambda_n$) is large, then $\frac{\lambda_1 - \lambda_n}{\lambda_1 + \lambda_n} \approx 1$. The gradient method converges very slowly. Hence this method is not recommendable.
- It is favorable to choose that the search directions $\{v^{(i)}\}$ as mutually A -conjugate, where A is symmetric positive definite.

Definition

Two vectors p and q are called A -conjugate (A -orthogonal), if $p^T A q = 0$.



Lemma

Let $v_1, \dots, v_n \neq 0$ be pairwise A -conjugate. Then they are *linearly independent*.

Proof: From

$$0 = \sum_{j=1}^n c_j v_j$$

follows that

$$0 = (v_k)^T A \left(\sum_{j=1}^n c_j v_j \right) = \sum_{j=1}^n c_j (v_k)^T A v_j = c_k (v_k)^T A v_k,$$

so $c_k = 0$, for $k = 1, \dots, n$.



Theorem

Let A be *symm. positive definite* and $v_1, \dots, v_n \in \mathbb{R}^n \setminus \{0\}$ be *pairwisely A -orthogonal*. Give x_0 and let $r_0 = b - Ax_0$. For $k = 1, \dots, n$, let

$$\alpha_k = \frac{\langle v_k, b - Ax_{k-1} \rangle}{\langle v_k, Av_k \rangle} \quad \text{and} \quad x_k = x_{k-1} + \alpha_k v_k.$$

Then $Ax_n = b$ and

$$\langle b - Ax_k, v_j \rangle = 0, \quad \text{for each } j = 1, 2, \dots, k-1.$$

Proof: Since, for each $k = 1, 2, \dots, n$,

$$x_k = x_{k-1} + \alpha_k v_k,$$

we have

$$\begin{aligned} Ax_n &= Ax_{n-1} + \alpha_n Av_n = (Ax_{n-2} + \alpha_{n-1} Av_{n-1}) + \alpha_n Av_n \\ &\vdots \\ &= Ax_0 + \alpha_1 Av_1 + \alpha_2 Av_2 + \cdots + \alpha_n Av_n. \end{aligned}$$



It implies that

$$\begin{aligned} & \langle Ax_n - b, v_k \rangle \\ = & \langle Ax_0 - b, v_k \rangle + \alpha_1 \langle Av_1, v_k \rangle + \cdots + \alpha_n \langle Av_n, v_k \rangle \\ = & \langle Ax_0 - b, v_k \rangle + \alpha_1 \langle v_1, Av_k \rangle + \cdots + \alpha_n \langle v_n, Av_k \rangle \\ = & \langle Ax_0 - b, v_k \rangle + \alpha_k \langle v_k, Av_k \rangle \\ = & \langle Ax_0 - b, v_k \rangle + \frac{\langle v_k, b - Ax_{k-1} \rangle}{\langle v_k, Av_k \rangle} \langle v_k, Av_k \rangle \\ = & \langle Ax_0 - b, v_k \rangle + \langle v_k, b - Ax_{k-1} \rangle \\ = & \langle Ax_0 - b, v_k \rangle \\ & + \langle v_k, b - Ax_0 + Ax_0 - Ax_1 + \cdots - Ax_{k-2} + Ax_{k-2} - Ax_{k-1} \rangle \\ = & \langle Ax_0 - b, v_k \rangle + \langle v_k, b - Ax_0 \rangle + \langle v_k, Ax_0 - Ax_1 \rangle \\ & + \cdots + \langle v_k, Ax_{k-2} - Ax_{k-1} \rangle \\ = & \langle v_k, Ax_0 - Ax_1 \rangle + \cdots + \langle v_k, Ax_{k-2} - Ax_{k-1} \rangle . \end{aligned}$$



For any i

$$x_i = x_{i-1} + \alpha_i v_i \quad \text{and} \quad Ax_i = Ax_{i-1} + \alpha_i Av_i,$$

we have

$$Ax_{i-1} - Ax_i = -\alpha_i Av_i.$$

Thus, for $k = 1, \dots, n$,

$$\begin{aligned} & \langle Ax_n - b, v_k \rangle \\ &= -\alpha_1 \langle v_k, Av_1 \rangle - \dots - \alpha_{k-1} \langle v_k, Av_{k-1} \rangle = 0 \end{aligned}$$

which implies that $Ax_n = b$.

Suppose that

$$\langle r_{k-1}, v_j \rangle = 0 \quad \text{for } j = 1, 2, \dots, k-1. \quad (9)$$

By the result

$$r_k = b - Ax_k = b - A(x_{k-1} + \alpha_k v_k) = r_{k-1} - \alpha_k Av_k$$



it follows that

$$\begin{aligned}\langle r_k, v_k \rangle &= \langle r_{k-1}, v_k \rangle - \alpha_k \langle Av_k, v_k \rangle \\ &= \langle r_{k-1}, v_k \rangle - \frac{\langle v_k, b - Ax_{k-1} \rangle}{\langle v_k, Av_k \rangle} \langle Av_k, v_k \rangle \\ &= 0.\end{aligned}$$

From assumption (9) and A -orthogonality, for $j = 1, \dots, k-1$

$$\langle r_k, v_j \rangle = \langle r_{k-1}, v_j \rangle - \alpha_k \langle Av_k, v_j \rangle = 0$$

which is completed the proof by the mathematic induction. □

Method of conjugate directions:

Let A be symmetric positive definite, $b, x_0 \in \mathbb{R}^n$. Given $v_1, \dots, v_n \in \mathbb{R}^n \setminus \{0\}$ pairwise A -orthogonal.

$$r_0 = b - Ax_0,$$

For $k = 1, \dots, n$,

$$\alpha_k = \frac{\langle v_k, r_{k-1} \rangle}{\langle v_k, Av_k \rangle}, \quad x_k = x_{k-1} + \alpha_k v_k,$$

$$r_k = r_{k-1} - \alpha_k Av_k = b - Ax_k.$$

End For



Practical Implementation

- In k -th step a direction v_k which is A -orthogonal to v_1, \dots, v_{k-1} must be determined.
- It allows for orthogonalization of r_k against v_1, \dots, v_k .
- Let $r_k \neq 0$, $g(x)$ decreases strictly in the direction $-r_k$. For $\varepsilon > 0$ small, we have $g(x_k - \varepsilon r_k) < g(x_k)$.

If $r_{k-1} = b - Ax_{k-1} \neq 0$, then we use r_{k-1} to generate v_k by

$$v_k = r_{k-1} + \beta_{k-1}v_{k-1}. \quad (10)$$

Choose β_{k-1} such that

$$\begin{aligned} 0 &= \langle v_{k-1}, Av_k \rangle = \langle v_{k-1}, Ar_{k-1} + \beta_{k-1}Av_{k-1} \rangle \\ &= \langle v_{k-1}, Ar_{k-1} \rangle + \beta_{k-1} \langle v_{k-1}, Av_{k-1} \rangle. \end{aligned}$$



That is

$$\beta_{k-1} = -\frac{\langle v_{k-1}, Ar_{k-1} \rangle}{\langle v_{k-1}, Av_{k-1} \rangle}. \quad (11)$$

Theorem

Let v_k and β_{k-1} be defined in (10) and (11), respectively. Then r_0, \dots, r_{k-1} are mutually orthogonal and

$$\langle v_k, Av_i \rangle = 0, \quad \text{for } i = 1, 2, \dots, k-1.$$

That is $\{v_1, \dots, v_k\}$ is an A -orthogonal set.

Having chosen v_k , we compute

$$\begin{aligned} \alpha_k &= \frac{\langle v_k, r_{k-1} \rangle}{\langle v_k, Av_k \rangle} = \frac{\langle r_{k-1} + \beta_{k-1}v_{k-1}, r_{k-1} \rangle}{\langle v_k, Av_k \rangle} \\ &= \frac{\langle r_{k-1}, r_{k-1} \rangle}{\langle v_k, Av_k \rangle} + \beta_{k-1} \frac{\langle v_{k-1}, r_{k-1} \rangle}{\langle v_k, Av_k \rangle} \\ &= \frac{\langle r_{k-1}, r_{k-1} \rangle}{\langle v_k, Av_k \rangle}. \end{aligned}$$



Since

$$r_k = r_{k-1} - \alpha_k Av_k,$$

we have

$$\langle r_k, r_k \rangle = \langle r_{k-1}, r_k \rangle - \alpha_k \langle Av_k, r_k \rangle = -\alpha_k \langle r_k, Av_k \rangle.$$

Further, from (12),

$$\langle r_{k-1}, r_{k-1} \rangle = \alpha_k \langle v_k, Av_k \rangle,$$

so

$$\begin{aligned} \beta_k &= -\frac{\langle v_k, Ar_k \rangle}{\langle v_k, Av_k \rangle} = -\frac{\langle r_k, Av_k \rangle}{\langle v_k, Av_k \rangle} \\ &= \frac{(1/\alpha_k) \langle r_k, r_k \rangle}{(1/\alpha_k) \langle r_{k-1}, r_{k-1} \rangle} = \frac{\langle r_k, r_k \rangle}{\langle r_{k-1}, r_{k-1} \rangle}. \end{aligned}$$



Algorithm (Conjugate Gradient method (CG-method))

Let A be s.p.d., $b \in \mathbb{R}^n$, choose $x_0 \in \mathbb{R}^n$, $r_0 = b - Ax_0 = v_0$.

If $r_0 = 0$, then $N = 0$ stop, otherwise for $k = 0, 1, \dots$

(a). $\alpha_k = \frac{\langle r_k, r_k \rangle}{\langle v_k, Av_k \rangle},$

(b). $x_{k+1} = x_k + \alpha_k v_k,$

(c). $r_{k+1} = r_k - \alpha_k Av_k,$

(d). If $r_{k+1} = 0$, let $N = k + 1$, stop.

(e). $\beta_k = \frac{\langle r_{k+1}, r_{k+1} \rangle}{\langle r_k, r_k \rangle},$

(f). $v_{k+1} = r_{k+1} + \beta_k v_k.$

- Theoretically, the exact solution is obtained in n steps.
- If A is well-conditioned, then approximate solution is obtained in about \sqrt{n} steps.
- If A is ill-conditioned, then the number of iterations may be greater than n .



Select a nonsingular matrix C so that

$$\tilde{A} = C^{-1}AC^{-T}$$

is better conditioned.

Consider the linear system

$$\tilde{A}\tilde{x} = \tilde{b},$$

where

$$\tilde{x} = C^T x \quad \text{and} \quad \tilde{b} = C^{-1}b.$$

Then

$$\tilde{A}\tilde{x} = (C^{-1}AC^{-T})(C^T x) = C^{-1}Ax.$$

Thus,

$$Ax = b \Leftrightarrow \tilde{A}\tilde{x} = \tilde{b} \quad \text{and} \quad x = C^{-T}\tilde{x}.$$



Since

$$\tilde{x}_k = C^T x_k,$$

we have

$$\begin{aligned}\tilde{r}_k &= \tilde{b} - \tilde{A}\tilde{x}_k = C^{-1}b - (C^{-1}AC^{-T})C^T x_k \\ &= C^{-1}(b - Ax_k) = C^{-1}r_k.\end{aligned}$$

Let

$$\tilde{v}_k = C^T v_k \quad \text{and} \quad w_k = C^{-1}r_k.$$

Then

$$\begin{aligned}\tilde{\beta}_k &= \frac{\langle \tilde{r}_k, \tilde{r}_k \rangle}{\langle \tilde{r}_{k-1}, \tilde{r}_{k-1} \rangle} = \frac{\langle C^{-1}r_k, C^{-1}r_k \rangle}{\langle C^{-1}r_{k-1}, C^{-1}r_{k-1} \rangle} \\ &= \frac{\langle w_k, w_k \rangle}{\langle w_{k-1}, w_{k-1} \rangle}.\end{aligned}$$



Thus,

$$\begin{aligned}\tilde{\alpha}_k &= \frac{\langle \tilde{r}_{k-1}, \tilde{r}_{k-1} \rangle}{\langle \tilde{v}_k, \tilde{A}\tilde{v}_k \rangle} = \frac{\langle C^{-1}r_{k-1}, C^{-1}r_{k-1} \rangle}{\langle C^T v_k, C^{-1}AC^{-T}C^T v_k \rangle} \\ &= \frac{\langle w_{k-1}, w_{k-1} \rangle}{\langle C^T v_k, C^{-1}Av_k \rangle}\end{aligned}$$

and, since

$$\begin{aligned}\langle C^T v_k, C^{-1}Av_k \rangle &= (v_k)^T CC^{-1}Av_k = (v_k)^T Av_k \\ &= \langle v_k, Av_k \rangle,\end{aligned}$$

we have

$$\tilde{\alpha}_k = \frac{\langle w_{k-1}, w_{k-1} \rangle}{\langle v_k, Av_k \rangle}.$$

Further,

$$\tilde{x}_k = \tilde{x}_{k-1} + \tilde{\alpha}_k \tilde{v}_k, \quad \text{so } C^T x_k = C^T x_{k-1} + \tilde{\alpha}_k C^T v_k$$

and

$$x_k = x_{k-1} + \tilde{\alpha}_k v_k.$$



Continuing,

$$\tilde{r}_k = \tilde{r}_{k-1} - \tilde{\alpha}_k \tilde{A} \tilde{v}_k,$$

so

$$C^{-1}r_k = C^{-1}r_{k-1} - \tilde{\alpha}_k C^{-1}AC^{-T}C^T v_k$$

and

$$r_k = r_{k-1} - \tilde{\alpha}_k A v_k.$$

Finally,

$$\tilde{v}_{k+1} = \tilde{r}_k + \tilde{\beta}_k \tilde{v}_k \quad \text{and} \quad C^T v_{k+1} = C^{-1}r_k + \tilde{\beta}_k C^T v_k,$$

so

$$v_{k+1} = C^{-T}C^{-1}r_k + \tilde{\beta}_k v_k = C^{-T}w_k + \tilde{\beta}_k v_k.$$



Algorithm (Preconditioned CG-method (PCG-method))

Choose C and x_0 . Set $r_0 = b - Ax_0$, solve $Cw_0 = r_0$ and $C^T v_1 = w_0$.

If $r_0 = 0$, then $N = 0$ stop, otherwise for $k = 1, 2, \dots$

(a). $\alpha_k = \langle w_{k-1}, w_{k-1} \rangle / \langle v_k, Av_k \rangle,$

(b). $x_k = x_{k-1} + \alpha_k v_k,$

(c). $r_k = r_{k-1} - \alpha_k Av_k,$

(d). If $r_k = 0$, let $N = k + 1$, stop.

Otherwise, solve $Cw_k = r_k$ and $C^T z_k = w_k,$

(e). $\beta_k = \langle w_k, w_k \rangle / \langle w_{k-1}, w_{k-1} \rangle,$

(f). $v_{k+1} = z_k + \beta_k v_k.$

