Sparse Matrices — Basic Iterative Methods Matrix Computations — CPSC 5006 E

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Sparse Matrices — Basic Iterative Methods

- Jacobi.
- Gauss-Seidel.
- SOR.
- SSOR.
- Saad. Iterative Methods for Sparse Linear System sections 4.1 and 4.2.
- Golub-Van Loan, section 10.1.
- Barrett et al. Templates for the Solution of Linear Systems, section 2.2.



Jacobi (p. 510)

Perhaps the simplest iterative scheme is the **Jacobi iteration**. It is defined for matrices that have non zero diagonal elements¹. The method can be motivated by rewritting the 3-by-3 system Ax = b as follows:

$$x_1 = (b_1 - a_{12}x_2 - a_{13}x_3)/a_{11}$$

 $x_2 = (b_2 - a_{21}x_1 - a_{23}x_3)/a_{22}$
 $x_3 = (b_3 - a_{31}x_1 - a_{32}x_2)/a_{33}$

Suppose $x^{(k)}$ is an approximation to $x = A^{-1}b$. A natural way to generate a new approximation $x^{(k+1)}$ is to compute:

$$x_1^{(k+1)} = (b_1 - a_{12}x_2^{(k)} - a_{13}x_3^{(k)})/a_{11}$$

$$x_2^{(k+1)} = (b_2 - a_{21}x_1^{(k)} - a_{23}x_3^{(k)})/a_{22}$$

$$x_3^{(k+1)} = (b_3 - a_{31}x_1^{(k)} - a_{32}x_2^{(k)})/a_{33}$$

¹This assumption can be reduced

Jacobi (p. 510)

The general equation is

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

The sum can be divided in two parts, the first part up to i-1 and the second part starting à i+1

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

Jacobi Update

Note that in the Jacobi iteration one does not use the most recently available information when computing $x_i^{(k+1)}$. For example $x_1^{(k)}$ is used in the calculation of $x_2^{(k+1)}$ even though component $x_1^{(k+1)}$ is known.

This allows to update the vector $x^{(k+1)}$, from the vector $x^{(k)}$, looping over the component in any order. It also allows easy parallelization on multiple processors.

Jacobi (p. 510)

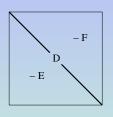
For general n, we have the Jacobi algorithm:

Algorithm 1 Jacobi. $A \in \mathbb{R}^{n \times n}$ with non zero diagonal elements. $b \in \mathbb{R}^n$, $x^{(0)}$, an initial guess, ε a stopping criterion and *MaxIter*, the maximum number of iterations if it does not converge.

```
1: for k=1: MaxIter
2: for i=1: n
3: x_i^{(k+1)} = \left(b_i - \sum_{j=1}^{i-1} a_{ij} x_j^{(k)} - \sum_{j=i+1}^{n} a_{ij} x_j^{(k)}\right) a_{ii}^{-1}
4: end
5: if \|x^{(k+1)} - x^{(k)}\| < \varepsilon then
6: Break
7: end
8: end
```

Matrix Form of the Jacobi Method (p. 511)

Relaxation schemes: based on the decomposition



$$A = D - E - F$$
 $D = \text{diag}(A);$
 $E = \text{negative of the strict lower}$

part of $A;$
 $F = \text{negative of the strict upper}$

part of A .

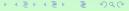
Simplest method for solving Ax = b: Jacobi iteration

$$Dx^{(k+1)} = (E+F)x^{(k)} + b$$

$$x^{(k+1)} = D^{-1}((E+F)x^{(k)} + b)$$

$$x^{(k+1)} = D^{-1}(E+F)x^{(k)} + D^{-1}b$$

Analysed using iteration matrix $M_{Jac} = D^{-1}(E + F)$.



Jacobi (Matrix Form) (p. 510)

For general n, we have the matrix form of the Jacobi algorithm:

Algorithm 2 Jacobi. $M_{Jac} = D^{-1}(E+F) \in \mathbb{R}^{n \times n}$. $b_{Jac} = D^{-1}b \in \mathbb{R}^n$, $x^{(0)}$, an initial guess, ε a stopping criterion and MaxIter, the maximum number of iterations if it does not converge.

- 1: **for** k = 1 : *MaxIter*
- 2: $x^{(k+1)} = M_{Jac}x^{(k)} + b_{Jac}$
- 3: **if** $||x^{(k+1)} x^{(k)}|| < \varepsilon$ **then**
- 4: Break
- 5: **end**
- 6: **end**

Note: This algorithm contains only a matrix-vector update.



Gauss-Seidel

Jacobi:

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

Idea: Use the newest value of x_i^{k+1} as soon as it is available.

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

This gives the Gauss-Seidel method.

Gauss-Seidel Update

Because the Gauss-Seidel method uses the most recently available information when computing $x_i^{(k+1)}$, it usually converge faster than the Jacobi method.

However, because we cannot compute $x_i^{(k+1)}$ before $x_{i-1}^{(k+1)}$ is computed, parallelization on multiple processors is not easy. Also, the Gauss-Seidel method depends on the traversal order of the unknowns. The classics are the forward loop from 1 to n, the reverse loop from n to 1, a combination of both, a random traversal, etc.

Gauss-Seidel (p. 510)

For general *n*, we have the Gauss-Seidel algorithm:

Algorithm 3 Gauss-Seidel. $A \in \mathbb{R}^{n \times n}$ with non zero diagonal elements. $b \in \mathbb{R}^n$, $x^{(0)}$, an initial guess, ε a stopping criterion and *MaxIter*, the maximum number of iterations if it does not converge.

```
1: for k = 1 : MaxIter
2: for i = 1 : n
3: x_i^{(k+1)} = \left(b_i - \sum_{j=1}^{i-1} a_{ij} x_j^{(k+1)} - \sum_{j=i+1}^{n} a_{ij} x_j^{(k)}\right) a_{ii}^{-1}
4: end
5: if ||x^{(k+1)} - x^{(k)}|| < \varepsilon then
6: Break
7: end
8: end
```

Gauss-Seidel

<u>Idea:</u> correct the *i*-th component of the current approximate solution, i = 1, 2, ...n, to zero out i - th component of residual.

Gauss-Seidel: $x_i^{new} = \frac{1}{a_{ii}} \left[b_i - \sum_{j < i} a_{ij} x_j^{new} - \sum_{j > i} a_{ij} x_j^{old} \right]$

Matrix form of Gauss-Seidel:

$$(D-E)x^{(k+1)} = Fx^{(k)} + b$$

Analysed using iteration matrix $M_{GS} = (D - E)^{-1}(F)$. Can also define a backward Gauss-Seidel Iteration:

$$(D-F)x^{(k+1)} = Ex^{(k)} + b$$

and a Symmetric Gauss-Seidel Iteration: forward sweep followed by backward sweep.

Relaxation

<u>Relaxation</u> is based on relaxing the Gauss-Seidel iteration:

•
$$x_i^{(k+1)} = \xi_i^{(k)} + \omega(\xi_i^{GS} - \xi_i^{(k)})$$

- $0 < \omega < 1 \Leftrightarrow Under-relaxation$.
- $\omega = 1 \Leftrightarrow \mathsf{Gauss}\text{-}\mathsf{Seidel}.$
- $1 < \omega < 2 \Leftrightarrow Over-relaxation$.

Over-relaxation is based on the decomposition:

$$\omega A = (D - \omega E) - (\omega F + (1 - \omega)D)$$

→ successive overrelaxation, (SOR):

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

corresponding to iteration matrix

$$M_{\omega SOR} = (D - \omega E)^{-1} (\omega F + (1 - \omega)D).$$



Iteratives Matrices

Iteration matrices

In Jacobi, Gauss-Seidel, or SOR, the iteration is of the form,

$$x^{(k+1)} = Mx^{(k)} + f$$

where

•
$$M_{Jac} = D^{-1}(E+F) = I - D^{-1}A$$

•
$$M_{GS}(A) = (D - E)^{-1}F = I - (D - E)^{-1}A$$

•
$$M_{\omega SOR}(A) = (D - \omega E)^{-1} (\omega F + (1 - \omega)D)$$

= $I - (\omega^{-1}D - E)^{-1}A$

Convergence

Jacobi and Gauss-Seidel converge for diagonal dominant matrices

SOR converges for $0<\omega<2$ for SPD matrices

Optimal ω known for 'consistently ordered matrices' (eig-vals of $\alpha^{-1}D^{-1}E + \alpha D^{-1}F$ indep. of α):

$$\omega_{\text{optimal}} = \frac{2}{1 + \sqrt{1 - \rho(M_{Jac})^2}}.$$