

ME667 Course Notes

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1 Iterative Methods for Matrix Inversion

1. Basic Concept
2. Point-Jacobi, Gauss-Seidel, SOR

Inversion by iterative schemes:

$$\textcircled{1} \quad A\vec{x} = \vec{b}$$

$$\text{Let } A = A_1 - A_2$$

$$A_1\vec{x} = A_2\vec{x} + \vec{b}$$

One can construct an iterative method by writing

$$A_1\vec{x}^{(k+1)} = A_2\vec{x}^{(k)} + \vec{b}$$

k is the iteration index. Starting with some guess $\vec{x}^{(0)}$ this scheme can be used to obtain solution to the system $A\vec{x} = \vec{b}$

The idea is as $k \rightarrow \infty$ $\vec{x}^{(k+1)} \rightarrow \vec{x}^{(k)}$

$$\begin{aligned} \text{Then we get } A_1\vec{x}^{(k+1)} - A_2\vec{x}^{(k)} &\approx (A_1 - A_2)\vec{x}^{(k)} \\ &= A\vec{x}^{(k)} - \vec{b} \end{aligned}$$

Requirements:

① A_1 should be easily invertable

② Iterations should converge

$$\lim_{k \rightarrow \infty} \vec{x}^{(k)} = \vec{x}$$

Let the error at iteration k be $\epsilon^{(k)} = \vec{x} - \vec{x}^{(k)}$

we can write

$$A_1(\vec{x} - \vec{x}^{(k)}) = A_2(\vec{x} - \vec{x}^{(k)}) + (\vec{b} - \vec{b})$$

$$\Rightarrow A_1\epsilon^{(k+1)} = A_2\epsilon^{(k)}$$

$$\epsilon^{(k+1)} = A_1^{-1}A_2\epsilon^{(k)}$$

$$\epsilon^{(k)} = (A_1^{-1}A_2)^k \epsilon^{(0)}$$

so we need
 $\lim_{k \rightarrow \infty} \epsilon^{(k)} = 0$

$A_1^{-1} A_2 = B^K$ will tend to zero

iff $|\lambda_i|_{\max} \leq 1$ λ_i are eigenvalues

$\rho = |\lambda|_{\max}$ is the spectral radius

Point-Jacobi

To have easy inversion of a matrix a diagonal matrix is preferred.

So let A_1 be Diagonal matrix

$$A_1 = \begin{bmatrix} -4 & & & \\ & -4 & & \\ & & -4 & \\ & & & \ddots \end{bmatrix} \quad \text{in the Laplace problem}$$

Then $A_2 = A_1 - A = \begin{bmatrix} 0 & & -1 & \\ -1 & 0 & -1 & -1 \\ -1 & 0 & -1 & \\ & & \ddots & \ddots \end{bmatrix}$ The other elements

$$A_1^{-1} = \begin{bmatrix} -1/4 & & & \\ & -1/4 & & \\ & & \ddots & \end{bmatrix}$$

$$A_1 \phi^{(k+1)} = A_2 \phi^{(k)} + \vec{b} \quad \text{where } \phi \text{ is the unknown vector}$$

For each ϕ

$$\boxed{\phi^{(k+1)} = -\frac{1}{4} A_2 \phi^{(k)} - \frac{1}{4} \vec{b}}$$

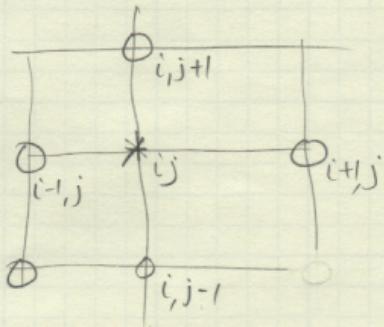
Using A_2 and the index notation

$$\boxed{\phi_{ij}^{(k+1)} = \frac{1}{4} [\phi_{i-1,j}^{(k)} + \phi_{i+1,j}^{(k)} + \phi_{i,j-1}^{(k)} + \phi_{i,j+1}^{(k)}] - \frac{1}{4} b_{ij}}$$

i, j are used in the same order as the matrix eqⁿ in the Laplace eqⁿ ①

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22-139 100 sheets



ϕ_{ij} depends one values of ϕ at "immediate" neighbors. No matrix storage involve

So starting with $\phi_{ij}^{(0)}$ guess (could be = 0) one just updates $\phi_{ij}^{(k+1)}$ using above eqⁿ. ①.

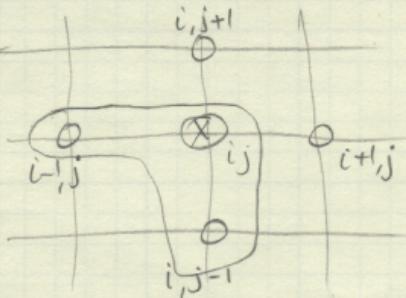
eigenvalues of $A_1^{-1} A_2 = -\frac{1}{4} A_2$ should be such that $| \lambda_{i,j} |_{\max} < 1$

From Linear Algebra for (uniform grids) eigenvalues are

$$\lambda_{mn} = \frac{1}{2} \left[\cos \frac{\pi m}{M} + \cos \frac{\pi n}{N} \right]$$

$M = 1, 2, 3, \dots, M-1$ & $N = 1, 2, 3, \dots, N-1$ are grid points in x & y

Gauss-Seidel



In point-Jacobi iterative scheme notice that we solve ϕ_{ij} in an orderly fashion. In other words $\phi_{i-1,j}$ & $\phi_{i,j-1}$ will be solved before ϕ_{ij} .

So we will have $\phi_{i-1,j}^{(k+1)}$, $\phi_{i,j-1}^{(k+1)}$ available before we try to obtain $\phi_{ij}^{(k+1)}$. Gauss-Seidel makes use of "latest" or most updated values of ϕ_{ij} .

Then

$$\phi_{ij}^{(k+1)} = \frac{1}{4} \left[\phi_{i-1,j}^{(k+1)} + \phi_{i+1,j}^{(k)} + \phi_{i,j-1}^{(k+1)} + \phi_{i,j+1}^{(k)} \right] - \frac{1}{4} b_{ij}$$

In the matrix form $A = A_1 - A_2$

$$\text{Here } A_1 = D - L \quad , \quad A_2 = U$$

$$A_1 = \begin{pmatrix} -4 & & & \\ -1 & -4 & & \\ & -1 & -4 & \\ & & \ddots & \ddots \end{pmatrix}$$

$$A_2 = \begin{pmatrix} 0 & -1 & & -1 \\ 0 & -1 & & -1 \\ 0 & 0 & -1 & \\ & & & 0 \end{pmatrix}$$

lower triangular
easy to invert

for Gauss-Seidel

$$\lambda_{mn} = \frac{1}{4} \left[\cos \frac{\pi m}{M} + \cos \frac{n\pi}{N} \right]^2$$

Square of the eigenvalues of point-Jacobi
Twice as fast

SOR (Successive over relaxation)

Note $A_1 \vec{x}^{(k+1)} = A_2 \vec{x}^{(k)} + \vec{b}$

$\phi^{(k+1)} = \phi^{(k)} + \vec{d}$ or change in ϕ at
each iteration is $\phi^{(k+1)} - \phi^{(k)} = \vec{d}$.

The idea of SOR is to "accelerate" this change
at each iteration in order to reduce # of iterations.
This is done by saying

$$\phi^{(k+1)} - \phi^{(k)} = w \vec{d}$$

w is the relaxation factor

$w > 1$ for accelerated convergence. Hence over-relaxation

$w = 1$ Gauss-Seidel

$w < 1$ slower convergence

We first use Gauss-Seidel, which can be written as

$$D \tilde{\phi}^{(k+1)} = L \phi^{(k+1)} + U \phi^{(k)} + \vec{b}$$

$\tilde{\phi}^{(k+1)}$ is not considered as the "accepted" value
at $k+1$.

$$\phi^{(k+1)} = \phi^{(k)} + w(\tilde{\phi}^{(k+1)} - \phi^{(k)}) \text{ is used for}$$

weighted convergence

$$\phi^{(k+1)} = (1-w) \phi^{(k)} + w \tilde{\phi}^{(k+1)}$$

$w=0 \Rightarrow$ Gauss-Seidel (GS)

$w > 1$ over-relaxation

So we weight the predicted value for GS more than the old iteration value

Combining the two eqns

$$\tilde{\phi}^{(k+1)} = D^{-1}L\phi^{(k+1)} + D^{-1}U\phi^{(k)} + D^{-1}b$$

$$\phi^{(k+1)} = (1-w)I\phi^{(k)} + wD^{-1}U\phi^{(k)} + wD^{-1}b + wD^{-1}L\phi^{(k+1)}$$

$$\Rightarrow (I - wD^{-1}L)\phi^{(k+1)} = [(1-w)I + wD^{-1}U]\phi^{(k)} + wD^{-1}b$$

$$\phi^{(k+1)} = \underbrace{[I - wD^{-1}L]^{-1}}_{GSOR} \underbrace{[(1-w)I + wD^{-1}U]\phi^{(k)}}_{+ (I - wD^{-1}L)^{-1}wD^{-1}b}$$

GSOR

To optimize convergence means selecting w to minimize the largest eigenvalue

For rectangular regions with uniform mesh, w_{opt} can be obtained analytically

$$w_{opt} = 2 \left\{ \frac{1 - \sqrt{1 - \lambda_{max}^2}}{\lambda_{max}^2} \right\}, \quad \begin{array}{l} \lambda_{max} \text{ is max. eigen} \\ \text{value of point-Jacob} \end{array}$$