Faculty of Engineering Civil Engineering

Numerical Methods

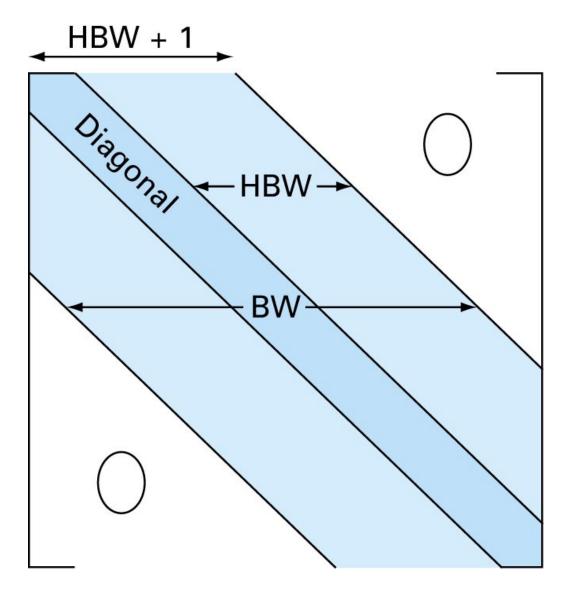


Special Matrices and Gauss-Siedel

Introduction

- Certain matrices have particular structures that can be exploited to develop efficient solution schemes.
- A **banded matrix** is a square matrix that has all elements equal to zero, with the exception of a band centered on the main diagonal. These matrices typically occur in solution of differential equations.
- The dimensions of a banded system can be quantified by two parameters: the band width BW and half-bandwidth HBW. These two values are related by BW=2HBW+1.

Banded matrix



Tridiagonal Systems

• A tridiagonal system has a bandwidth of 3:

$$\begin{bmatrix} f_1 & g_1 & & \\ e_2 & f_2 & g_2 & \\ & e_3 & f_3 & g_3 \\ & & e_4 & f_4 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \\ x_3 \\ x_4 \end{bmatrix} = \begin{cases} r_1 \\ r_2 \\ r_3 \\ r_4 \end{bmatrix}$$

• An efficient LU decomposition method, called *Thomas algorithm*, can be used to solve such an equation. The algorithm consists of three steps: decomposition, forward and back substitution, and has all the advantages of LU decomposition.

(a) Decomposition

DO
$$k = 2$$
, n
 $e_k = e_k/f_{k-1}$
 $f_k = f_k - e_k \cdot g_{k-1}$
END DO

(b) Forward substitution

 $DO \ k = 2, \ n$ $r_k = r_k - e_k \cdot r_{k-1}$ $END \ DO$

(c) Back substitution

$$x_n = r_n / f_n$$

 $DO \ k = n - 1, 1, -1$
 $x_k = (r_k - g_k \cdot x_{k+1}) / f_k$
END DO

Cholesky Decomposition

 This method is suitable for only symmetric systems where:

$$a_{ij} = a_{ji}$$
 and $A = A^T$ $[L] = \begin{bmatrix} l_{11} & 0 & 0 \\ l_{21} & l_{22} & 0 \\ l_{31} & l_{32} & l_{33} \end{bmatrix}$

$$A = L * L^{T}$$

$$\begin{bmatrix} a_{11} & a_{12} & a_{13} \\ a_{21} & a_{22} & a_{23} \\ a_{31} & a_{32} & a_{33} \end{bmatrix} = \begin{bmatrix} l_{11} & 0 & 0 \\ l_{21} & l_{22} & 0 \\ l_{31} & l_{32} & l_{33} \end{bmatrix} * \begin{bmatrix} l_{11} & l_{21} & l_{31} \\ 0 & l_{22} & l_{32} \\ 0 & 0 & l_{33} \end{bmatrix}$$

Cholesky Decomposition

$$a_{ki} - \sum_{j=1}^{i-1} l_{ij} \cdot l_{kj}$$

$$l_{ki} = \frac{l_{ki} - \sum_{j=1}^{i-1} l_{ij} \cdot l_{kj}}{l_{ii}} \quad for \quad i = 1, 2, \dots, k-1$$

$$l_{kk} = \sqrt{a_{kk} - \sum_{j=1}^{k-1} l_{kj}^2}$$

Pseudocode for Cholesky's LU Decomposition algorithm (cont'd)

```
DO k = 1.n
 DO i = 1. k - 1
   sum = 0.
   DO \ j = 1, \ i - 1
      sum = sum + a_{i,j} \cdot a_{k,j}
    END DO
   a_{ki} = (a_{ki} - sum)/a_{ii}
  END DO
  sum = 0.
  DO \ j = 1, \ k - 1
   sum = sum + a_{ki}^2
  END DO
 a_{kk} = \sqrt{a_{kk} - sum}
FND DO
```

Gauss-Siedel

- Iterative or approximate methods provide an alternative to the elimination methods. The Gauss-Seidel method is the most commonly used iterative method.
- The system [A]{X}={B} is reshaped by solving the first equation for x₁, the second equation for x₂, and the third for x₃, ...and nth equation for x_n. We will limit ourselves to a 3x3 set of equations.

Gauss-Siedel

$$\begin{aligned} a_{11}x_1 + a_{12}x_2 + a_{13}x_3 &= b_1 \\ a_{21}x_1 + a_{22}x_2 + a_{23}x_3 &= b_2 \end{aligned} \implies \begin{aligned} x_1 &= \frac{b_1 - a_{12}x_2 - a_{13}x_3}{a_{11}} \\ x_2 &= \frac{b_2 - a_{21}x_1 - a_{23}x_3}{a_{22}} \\ a_{22} \end{aligned} \\ x_1 &= \frac{b_3 - a_{31}x_1 - a_{32}x_2}{a_{33}} \end{aligned}$$

Now we can start the solution process by choosing guesses for the x's. A simple way to obtain initial guesses is to assume that they are zero. These zeros can be substituted into x_1 equation to calculate a new $x_1 = b_1/a_{11}$.

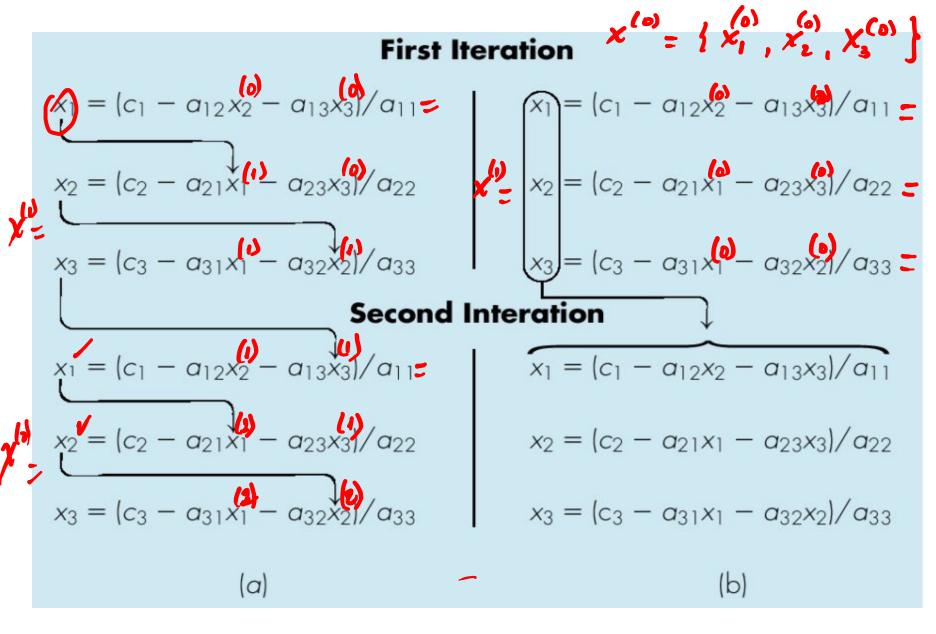
Gauss-Siedel

• New x_1 is substituted to calculate x_2 and x_3 . The procedure is repeated until the convergence criterion is satisfied:

$$\left|\varepsilon_{a,i}\right| = \left|\frac{x_i^{new} - x_i^{old}}{x_i^{new}}\right| 100\% < \varepsilon_s$$

Jacobi iteration Method

An alternative approach, called *Jacobi iteration*, utilizes a somewhat different technique. This technique includes computing a set of new x's on the basis of a set of old x's. Thus, as the new values are generated, they are not immediately used but are retained for the next iteration.



The Gauss-Seidel method

The Jacobi iteration method

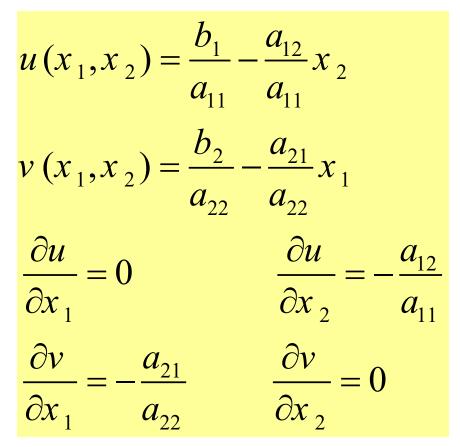
Convergence Criterion for Gauss-Seidel Method

- The gauss-siedel method is similar to the technique of fixed-point iteration.
- The Gauss-Seidel method has two fundamental problems as any iterative method:
 - 1. It is sometimes non-convergent, and
 - 2. If it converges, converges very slowly.
- Sufficient conditions for convergence of two linear equations, u(x,y) and v(x,y) are:

$$\left| \frac{\partial u}{\partial x} \right| + \left| \frac{\partial u}{\partial y} \right| < 1$$
$$\left| \frac{\partial v}{\partial x} \right| + \left| \frac{\partial v}{\partial y} \right| < 1$$

Convergence Criterion for Gauss-Seidel Method (cont'd)

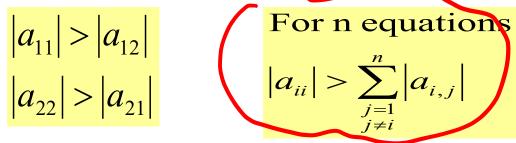
• Similarly, in case of two simultaneous equations, the Gauss-Seidel algorithm can be expressed as:



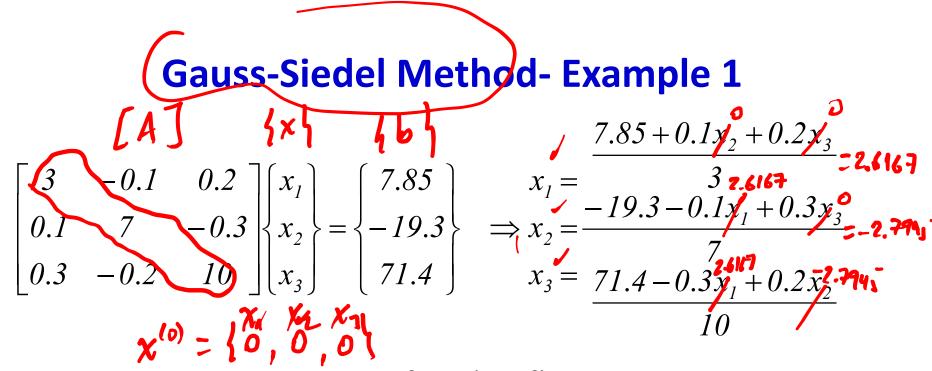
Convergence Criterion for Gauss-Seidel Method (cont'd)

Substitution into convergence criterion of two linear equations yield: $\left| \frac{a_{12}}{a_{11}} \right| < 1, \quad \left| \frac{a_{21}}{a_{22}} \right| < 1$

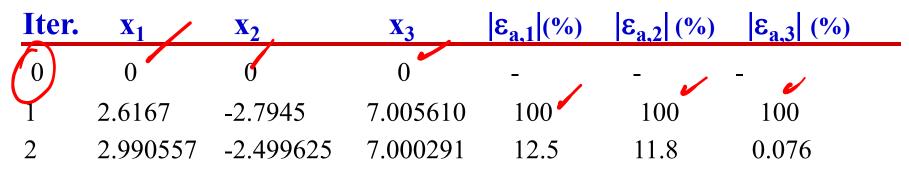
In other words, the absolute values of the slopes must be less than unity for convergence:



That is, the diagonal element must be greater than the offdiagonal element for each row.



• **Guess** x_1, x_2, x_3 = zero for the first guess



Improvement of Convergence Using Relaxation

$$x_{i}^{new} = \lambda \cdot x_{i}^{new} + (1 - \lambda) \cdot x_{i}^{old}$$

- Where λ is a weighting factor that is assigned a value between [0, 2]
- If $\lambda = 1$ the method is unmodified.
- If λ is between 0 and 1 (under relaxation) this is employed to make a non convergent system to converge.
- If λ is between 1 and 2 (over relaxation) this is employed to accelerate the convergence.

$$x_{1} + x_{2} - 2x_{3} = -20$$

$$-3x_{1} + x_{2} + 7x_{3} = -34$$

$$2x_{1} - 6x_{2} + 7x_{3} = -38$$
Rearrange so that
the equations are
diagonally dominant
$$2x_{1} - 6x_{2} - 2x_{3} = -20$$

$$2x_{1} + x_{2} - 2x_{3} = -20$$

$$2x_{1} - 6x_{2} - x_{3} = -38$$

$$-3x_{1} - x_{2} + 7x_{3} = -34$$

$$-3x_{1} - x_{2} + 7x_{3} = -34$$

$$-3x_{1} - x_{2} + 7x_{3} = -34$$

$$x_{1} = \frac{-20 - x_{2} + 2x_{3}}{-8}$$

$$x_{2} = \frac{-38 - 2x_{1} + x_{3}}{-6}$$

$$x_{3} = \frac{-34 + 3x_{1} + x_{2}}{7}$$

iteration	unknown	value	E _a	maximum ε_a
0	<i>X</i> ₁	0		
	<i>x</i> ₂	0		
	<i>X</i> ₃	0		
	<i>x</i> ₁	2.5	100.00%	
	<i>x</i> ₂	7.166667	100.00%	
	X 3	-2.7619	100.00%	100.00%
2	<i>x</i> ₁	4.08631	38.82%	
	<i>x</i> ₂	8.155754	12.13%	
	X 3	-1.94076	42.31%	42.31%
3	<i>x</i> ₁	4.004659	2.04%	
	<i>X</i> ₂	7.99168	2.05%	
	X 3	-1.99919	2.92%	2.92%

The same computation can be developed with relaxation where $\lambda = 1.2$

First iteration:

$$x_1 = \frac{-20 - x_2 + 2x_3}{-8} = \frac{-20 - 0 + 2(0)}{-8} = 2.5$$

Relaxation yields: $x_1 = 1.2(2.5) - 0.2(0) = 3$

$$x_{2} = \frac{-38 - 2x_{1} + x_{3}}{-6} = \frac{-38 - 2(3) + 0}{-6} = 7.333333$$

Relaxation yields: $x_{2} = 1.2(7.333333) - 0.2(0) = 8.8$

$$x_3 = \frac{-34 + 3x_1 + x_2}{7} = \frac{-34 + 3(3) + 8.8}{7} = -2.3142857$$

Relaxation yields: $x_3 = 1.2(-2.3142857) - 0.2(0) = -2.7771429$

lter.	unknown	value	relaxation	E _a	maximum ε_a
1	X ₁	2.5	3	100.00%	
	X ₂	7.3333333	8.8	100.00%	
	X ₃	-2.314286	-2.777143	100.00%	100.000%
2	X ₁	4.2942857	4.5531429	34.11%	
	<i>x</i> ₂	8.3139048	8.2166857	7.10%	
	X 3	-1.731984	-1.522952	82.35%	82.353%
3	X ₁	3.9078237	3.7787598	20.49%	
	<i>x</i> ₂	7.8467453	7.7727572	5.71%	
	X 3	-2.12728	-2.248146	32.26%	32.257%
4	<i>X</i> ₁	4.0336312	4.0846055	7.49%	
	<i>x</i> ₂	8.0695595	8.12892	4.38%	
	X 3	-1.945323	-1.884759	19.28%	19.280%

MATLAB M-File for Gauss-Seidel method

```
function x = GaussSeidel(A, b, es, maxit)
& Gaussfeidel(A,b): Gauss-Seidel method.
8 input:
& A = Coefficient Matrix
8 b = right hand side vector
% es = (optional) stop criterion (%) (default = 0.00001)
8
    maxit = (optional) maximum iterations (default = 50)
8 Output
% x = solution vector
if nargin < 4, maxit = 50; end
if nargin < 3, es = 0.00001; end
[m,n] = size(A);
if m ~= n, error('Matrix A must be square'); end
C = A;
for i = 1 : n
  C(i,i) = 0;
   x(i) = 0;
end
\mathbf{x} = \mathbf{x}^{\dagger}
for i = 1 : n
    C(i,1:n) = C(i,1:n) / A(i,i);
end
for i = 1: n
    d(i) = b(i) / A(i,i);
end
```

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MATLAB M-File for Gauss-Seidel method

Continued from previous page

```
disp(
                                                   x4 ...');
while (1)
    xold = x;
    for i = 1 : n
        x(i) = d(i) - C(i, :) * x;
        if x(i) ~= 0
            ea(i) = abs((x(i) - xold(i)) / x(i)) * 100;
        end
    end
    iter = iter + 1;
    disp([iter x'])
    if max(ea) <= es | iter >= maxit, break, end
end
if iter >= maxit
disp('Gauss Seidel method did not converge');
disp('results after maximum number of iterations');
else
    disp('Gauss Seidel method has converged');
end
x;
```

Gauss-Seidel Iteration

> A = [4 -1]	-1; 6 8 0;	-5 0 12];		
> b = [-2 45]	80];			
» x=Seidel(A	,b,x0,tol,	100);		
i	x 1	x 2	x 3	x4
1.0000	-0.5000	6.0000	6.4583	
2.0000	2.6146	3.6641	7.7561	
3.0000	2.3550	3.8587	7.6479	
4.0000	2.3767	3.8425	7.6569	
5.0000	2.3749	3.8439	7.6562	
6.0000	2.3750	3.8437	7.6563	
7.0000	2.3750	3.8438	7.6562	
8.0000	2.3750	3.8437	7.6563	
Gauss-Seidel	method co	nverged		

Converges faster than the Jacobi method shown in next page

Jacobi Iteration

> A = [4 -1]	-1; 6 8 0;	-5 0 12];		
> b = [-2 45]	5 80];	-		
» x=Jacobi(A	A,b,0.0001,	100);		
i	x1	x 2	x 3	x4
1.0000	-0.5000	5.6250	6.6667	
2.0000	2.5729	6.0000	6.4583	
3.0000	2.6146	3.6953	7.7387	
4.0000	2.3585	3.6641	7.7561	
5.0000	2.3550	3.8561	7.6494	
6.0000	2.3764	3.8587	7.6479	
7.0000	2.3767	3.8427	7.6568	
8.0000	2.3749	3.8425	7.6569	
9.0000	2.3749	3.8438	7.6562	
10.0000	2.3750	3.8439	7.6562	
11.0000	2.3750	3.8437	7.6563	
12.0000	2.3750	3.8437	7.6563	
13.0000	2.3750	3.8438	7.6562	
14.0000	2.3750	3.8438	7.6562	

Jacobi method converged