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Computers & Numerical Analysis (STR 681) Lecture 2 Linear Algebraic Equations

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Gauss-Jordan

- The Gauss-Jordan method is a variation of Gauss elimination.
- The major difference is that when an unknown is eliminated in the Gauss-Jordan method, it is eliminated from all other equations rather than just the subsequent ones.
- In addition, all rows are normalized by dividing them by their pivot elements. Thus, the elimination step results in an identity matrix rather than a triangular matrix.
- > Have the same drawbacks of Gauss Elimination.

Gauss-Jordan



Ex.,
$$3x_1 - 0.1x_2 - 0.2x_3 = 7.85$$

 $0.1x_1 + 7x_2 - 0.3x_3 = -19.3$
 $0.3x_1 - 0.2x_2 + 10x_3 = 71.4$

$$\begin{bmatrix} 3 & -0.1 & -0.2 & 7.85 \\ 0.1 & 7 & -0.3 & -19.3 \\ 0.3 & -0.2 & 10 & 71.4 \end{bmatrix} /3$$

$$\begin{bmatrix} 1 & -0.0333333 & -0.0666667 & 2.61667 \\ 0.1 & 7 & -0.3 & -19.3 \\ 0.3 & -0.2 & 10 & 71.4 \end{bmatrix} \times 0.1 \times 0.3$$

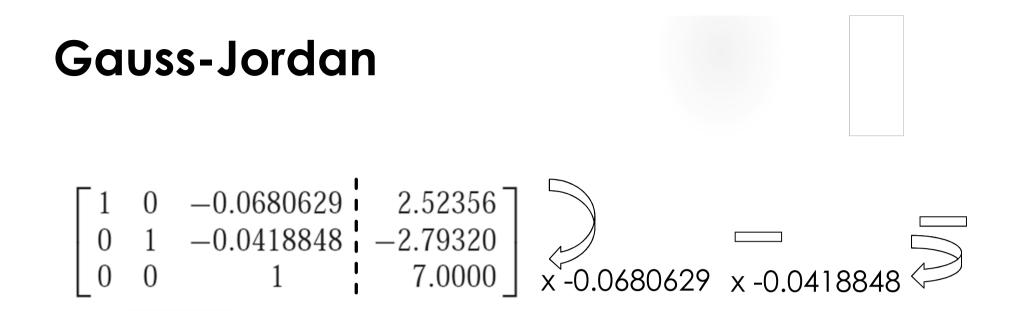
Gauss-Jordan



 $\begin{bmatrix} 1 & -0.0333333 & -0.066667 & 2.61667 \\ 0 & 7.00333 & -0.293333 & -19.5617 \\ 0 & -0.190000 & 10.0200 & 70.6150 \end{bmatrix} /7.00333$

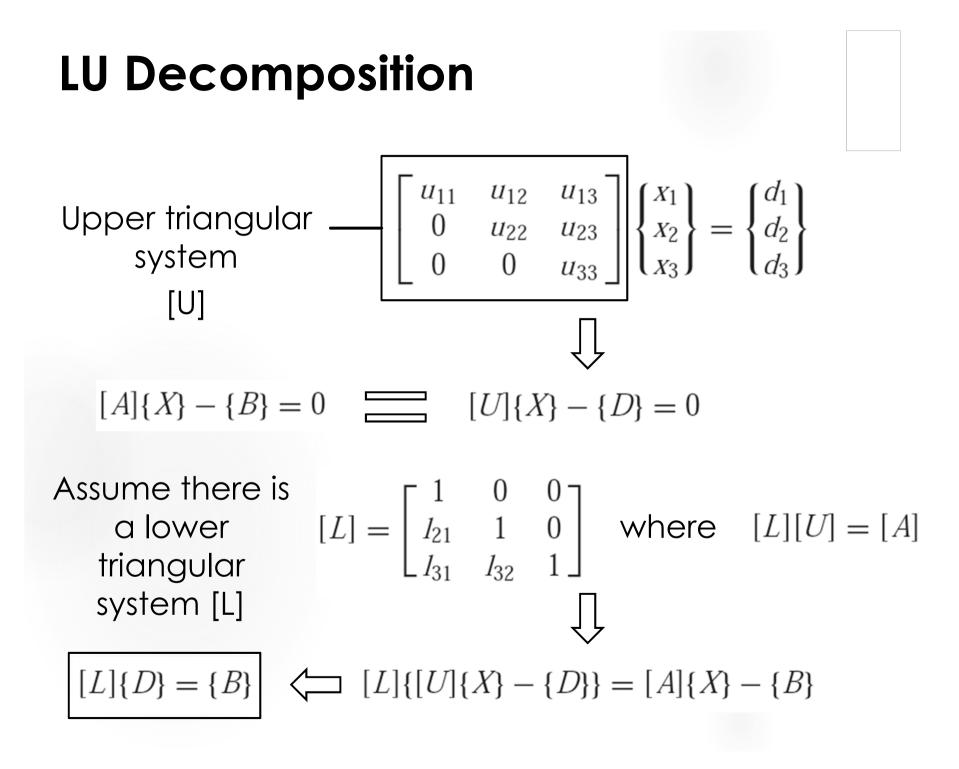
 $\begin{bmatrix} 1 & -0.0333333 & -0.066667 & 2.61667 \\ 0 & 1 & -0.0418848 & -2.79320 \\ 0 & -0.190000 & 10.0200 & 70.6150 \end{bmatrix} \times -0.0333333 \longrightarrow$

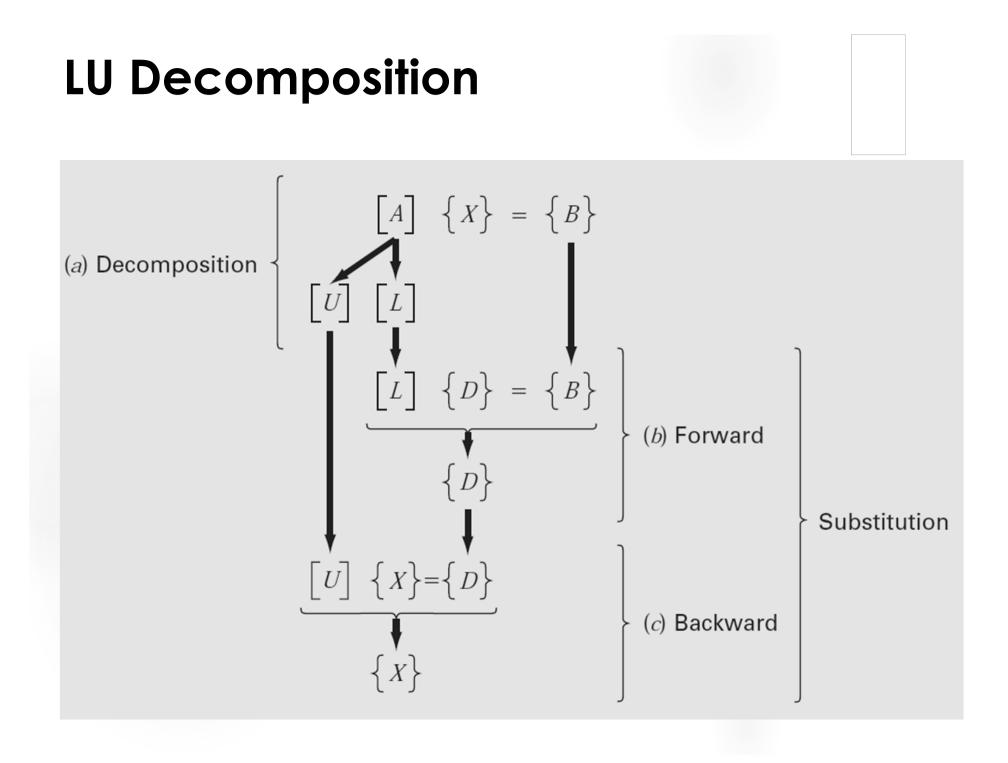
 $\begin{bmatrix} 1 & 0 & -0.0680629 & 2.52356 \\ 0 & 1 & -0.0418848 & -2.79320 \\ 0 & 0 & 10.01200 & 70.0843 \end{bmatrix} /10.012$



1	$\lceil 1 \rceil$	0	0	3.0000]
	0	1	0	-2.5000
	0	0	1	7.0000
	L			

- The time-consuming elimination step can be formulated so that it involves only operations on the matrix of coefficients,
 [A].
- It is well suited for those situations where many right-hand side vectors {B} must be evaluated for a single value of [A].
- One motive for introducing LU decomposition is that it provides an efficient means to compute the matrix inverse. The inverse has a number of valuable applications in engineering practice.
- > It also provides a means for evaluating system condition.

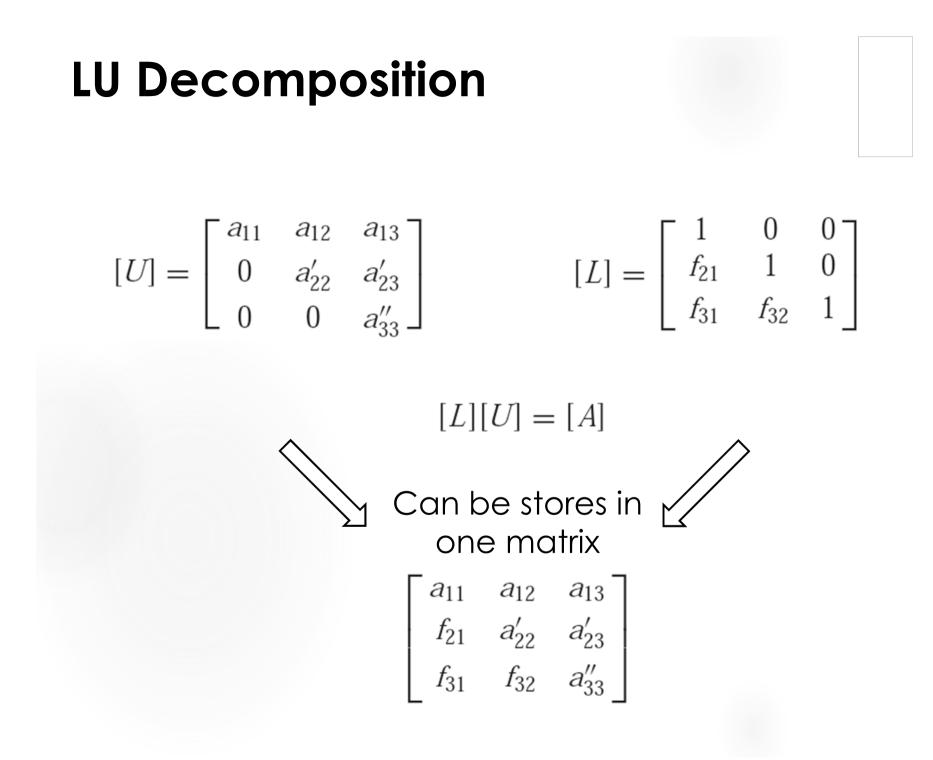






A two-step strategy for obtaining solutions:

- LU decomposition step. [A] is factored or "decomposed" into lower [L] and upper [U] triangular matrices.
- Substitution step. [L] and [U] are used to determine a solution {X} for a right-hand side {B}. This step itself consists of two steps:
 - ► Generate an intermediate vector {D} by forward substitution.
 - ▶ Then, the result can be solved by back substitution for {X}.





$$[L]{D} = {B}$$

$$d_i = b_i - \sum_{j=1}^{i-1} a_{ij} d_j \quad \text{for } i = 2, 3, \dots, n \quad \text{Forward}$$

substitution

$$\begin{bmatrix} U \end{bmatrix} \{X\} = \{D\}$$
$$x_n = d_n / a_{nn}$$
$$d_i - \sum_{j=i+1}^n a_{ij} x_j$$
$$x_i = \frac{d_i - \sum_{j=i+1}^n a_{ij} x_j}{a_{ii}}$$

Backward substitution

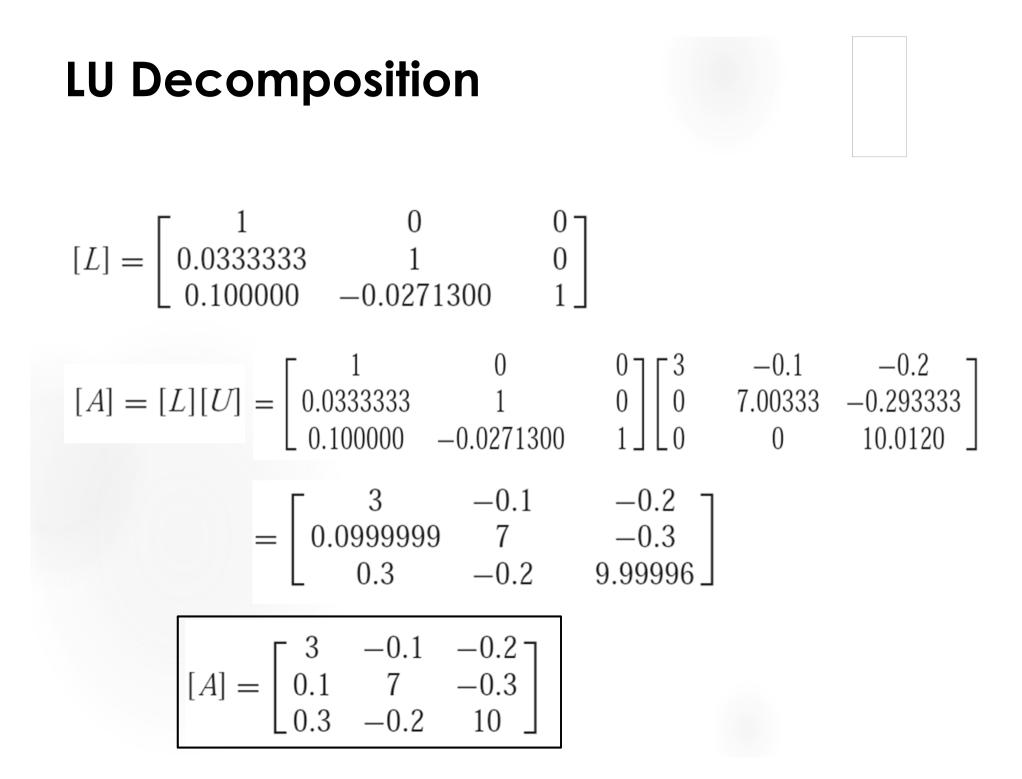
for $i = n - 1, n - 2, \dots, 1$



$$\begin{bmatrix} A \end{bmatrix} = \begin{bmatrix} 3 & -0.1 & -0.2 \\ 0.1 & 7 & -0.3 \\ 0.3 & -0.2 & 10 \end{bmatrix}$$
$$\begin{bmatrix} J \end{bmatrix}$$
$$\begin{bmatrix} 3 & -0.1 & -0.2 \\ 0 & 7.00333 & -0.293333 \\ 0 & 0 & 10.0120 \end{bmatrix}$$
After forward elimination

 \sim \sim

$$f_{21} = \frac{0.1}{3} = 0.033333333$$
$$f_{31} = \frac{0.3}{3} = 0.1000000$$
$$f_{32} = \frac{-0.19}{7.00333} = -0.0271300$$



$$\begin{bmatrix} 3 & -0.1 & -0.2 \\ 0.1 & 7 & -0.3 \\ 0.3 & -0.2 & 10 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \\ x_3 \end{bmatrix} = \begin{bmatrix} 7.85 \\ -19.3 \\ 71.4 \end{bmatrix}$$

$$[L] = \begin{bmatrix} 1 & 0 & 0 \\ 0.0333333 & 1 & 0 \\ 0.100000 & -0.0271300 & 1 \end{bmatrix} \{B\} = \begin{cases} 7.85 \\ -19.3 \\ 71.4 \end{cases}$$
$$[L]\{D\} = \{B\} \implies \begin{bmatrix} 1 & 0 & 0 \\ 0.0333333 & 1 & 0 \\ 0.100000 & -0.0271300 & 1 \end{bmatrix} \begin{cases} d_1 \\ d_2 \\ d_3 \end{cases} = \begin{cases} 7.85 \\ -19.3 \\ 71.4 \end{cases}$$

get {D} by forward substitution



 $d_1 = 7.85$

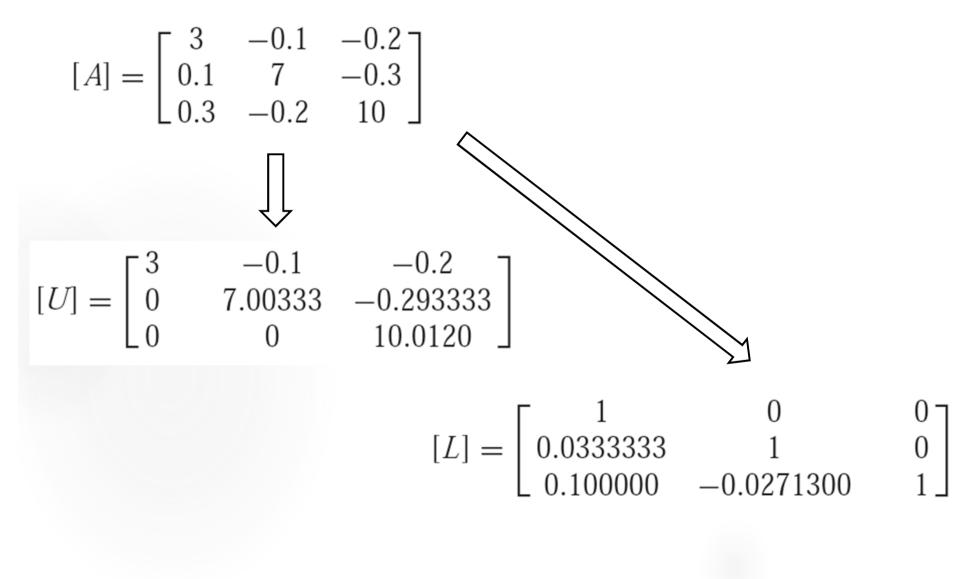
 $d_2 = -19.3 - 0.0333333(7.85) = -19.5617$

 $d_3 = 71.4 - 0.1(7.85) + 0.02713(-19.5617) = 70.0843$

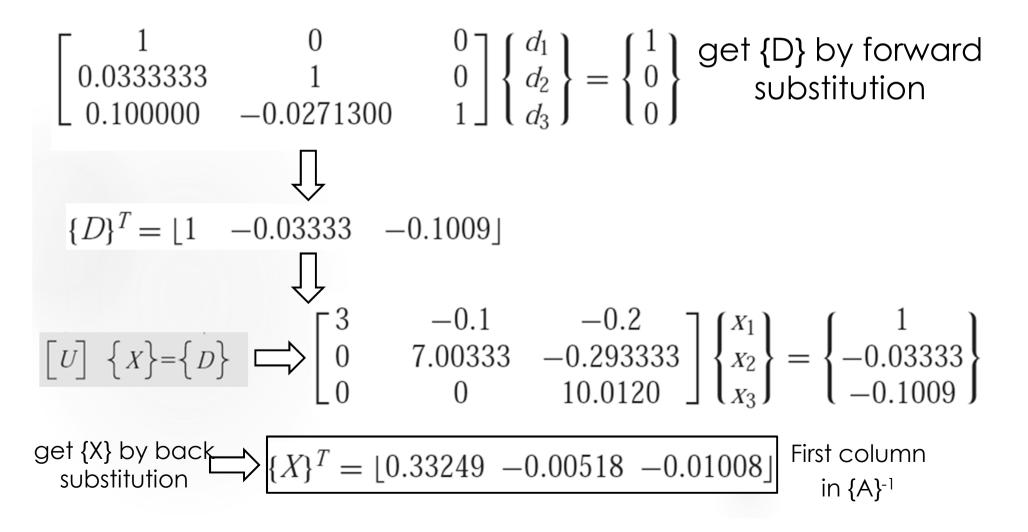
$$\{D\} = \left\{\begin{array}{c} 7.85\\ -19.5617\\ 70.0843 \end{array}\right\}$$

 \Longrightarrow Check slide # 59

$$\begin{bmatrix} U \end{bmatrix} \{X\} = \{D\} \Longrightarrow \begin{bmatrix} 3 & -0.1 & -0.2 \\ 0 & 7.00333 & -0.293333 \\ 0 & 0 & 10.0120 \end{bmatrix} \begin{bmatrix} X_1 \\ X_2 \\ X_3 \end{bmatrix} = \begin{cases} 7.85 \\ -19.5617 \\ 70.0843 \end{bmatrix}$$
$$\begin{bmatrix} X \} = \begin{cases} 3 \\ -2.5 \\ 7.00003 \end{bmatrix}$$



 $[L]{D} = {B}$



$$[A]^{-1} = \begin{bmatrix} 0.33249 & 0 & 0 \\ -0.00518 & 0 & 0 \\ -0.01008 & 0 & 0 \end{bmatrix}$$

To get the second column, repeat previous steps twice as follows:

$$\begin{bmatrix} 1 & 0 & 0 \\ 0.0333333 & 1 & 0 \\ 0.100000 & -0.0271300 & 1 \end{bmatrix} \begin{pmatrix} d_1 \\ d_2 \\ d_3 \end{pmatrix} = \begin{cases} 0 \\ 1 \\ 0 \end{pmatrix}$$
$$\begin{bmatrix} 1 & 0 & 0 \\ 0.0333333 & 1 & 0 \\ 0.100000 & -0.0271300 & 1 \end{bmatrix} \begin{pmatrix} d_1 \\ d_2 \\ d_3 \end{pmatrix} = \begin{cases} 0 \\ 0 \\ 1 \end{bmatrix}$$

$$[A]^{-1} = \begin{bmatrix} 0.33249 & 0.004944 & 0.006798 \\ -0.00518 & 0.142903 & 0.004183 \\ -0.01008 & 0.00271 & 0.09988 \end{bmatrix}$$

Check [A][A]⁻¹ = [I]

Importance of Inverse Calculation

Find out if the system is ill-conditioned:

 Scale the matrix of coefficients [A] so that the largest element in each row is 1. Invert the scaled matrix and if there are elements of [A]⁻¹ that are several orders of magnitude greater than one, it is likely that the system is ill-conditioned.

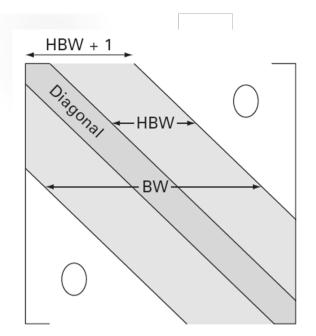
 $[A] = \begin{bmatrix} 1 & 1/2 & 1/3 \\ 1/2 & 1/3 & 1/4 \\ 1/3 & 1/4 & 1/5 \end{bmatrix} \longleftrightarrow [A]^{-1} = \begin{bmatrix} 9 & -18 & 10 \\ -36 & 96 & -60 \\ 30 & -90 & 60 \end{bmatrix}$ Hilbert Matrix

Importance of Inverse Calculation

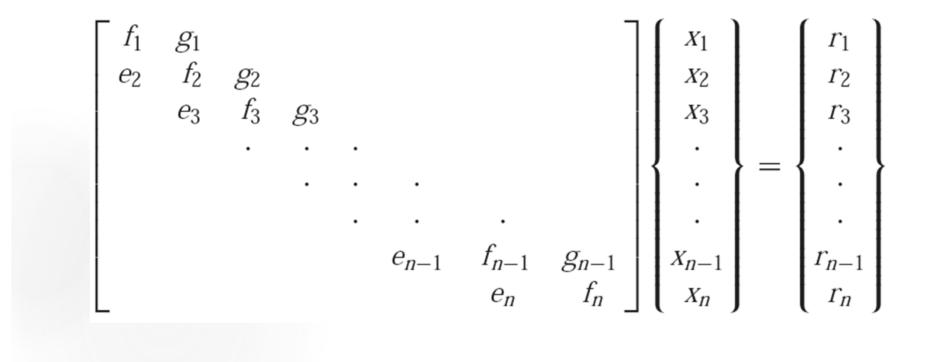
Find out if the system is ill-conditioned:

2. Multiply the inverse by the original coefficient matrix and assess whether the result is close to the identity matrix. If not, it indicates ill-conditioning.

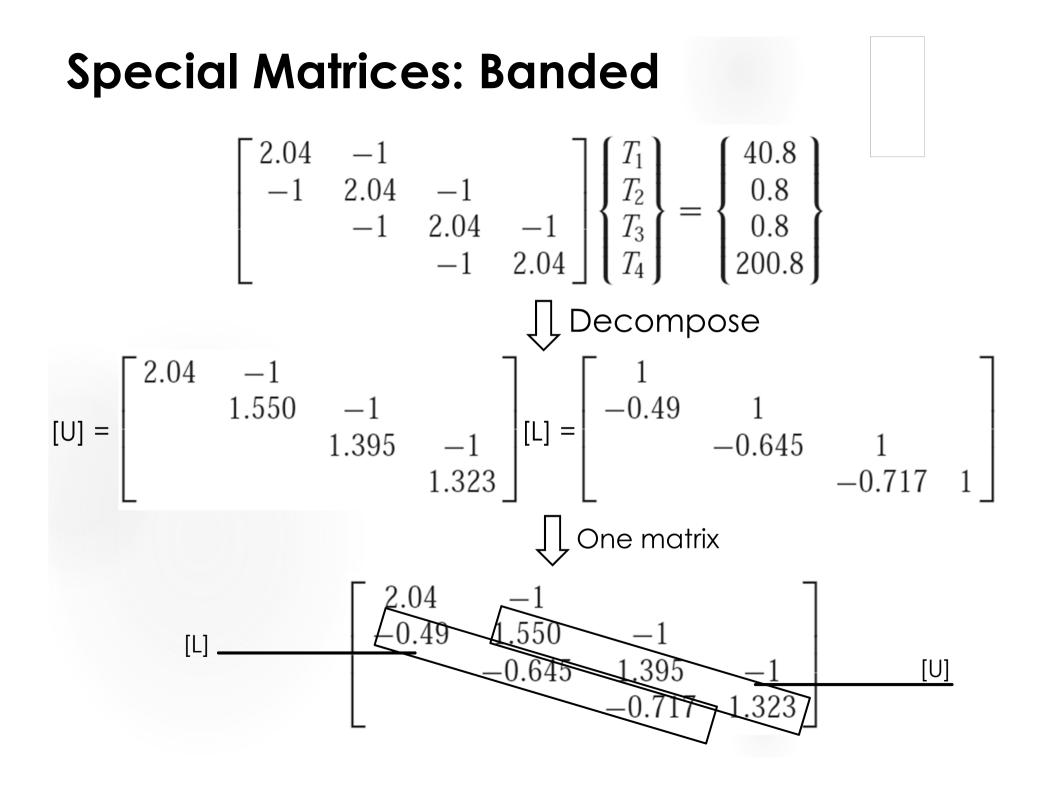
- > BW: band width
- > HBW: half band width
- ➢ BW = 2HBW + 1



- Gauss elimination or conventional LU decomposition can be employed to solve banded equations, but they are inefficient.
- None of the elements outside the band would change from their original values of zero. Thus, unnecessary space and time would be expended on the storage and manipulation of these useless zeros.
- > An alternative: Thomas Algorithm.

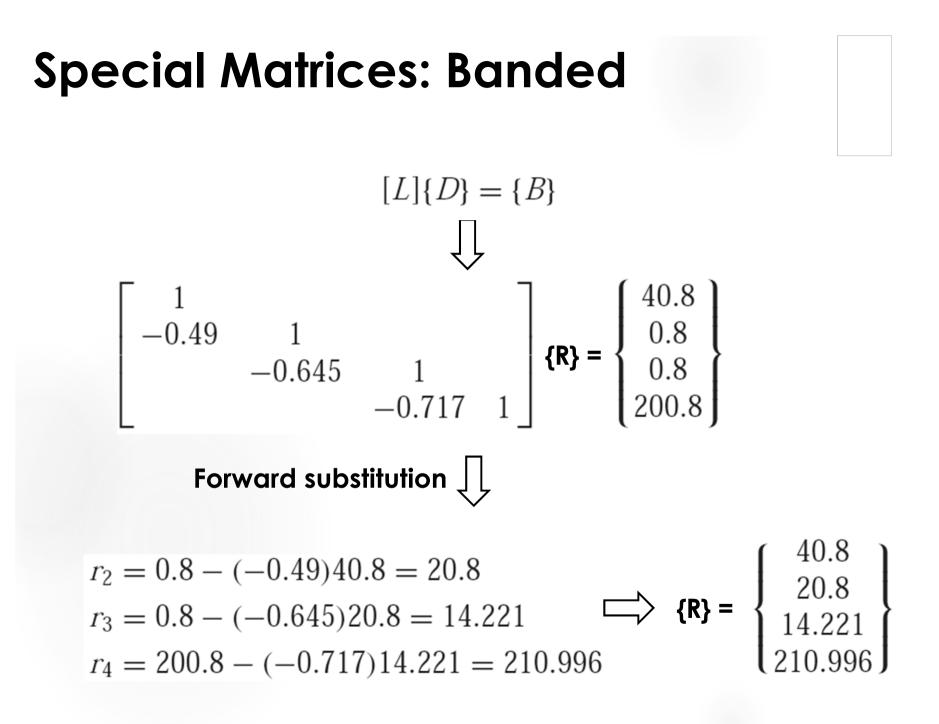


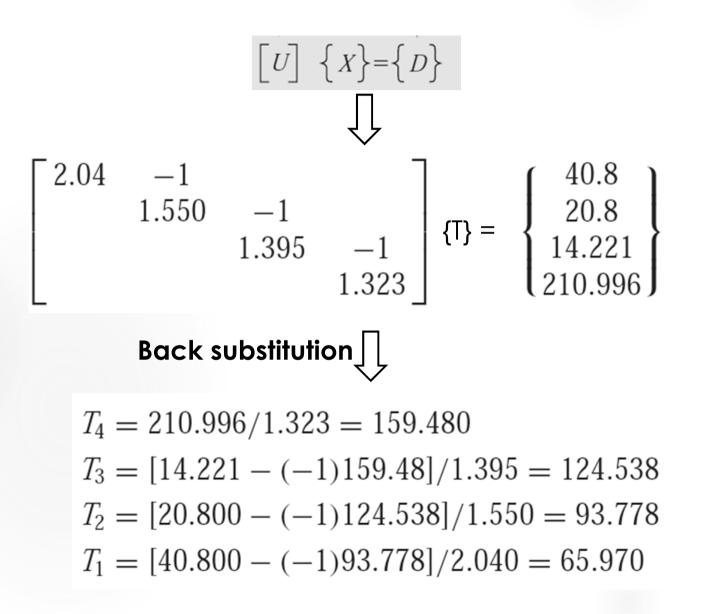




$$e_2 = -1/2.04 = -0.49$$

 $f_2 = 2.04 - (-0.49)(-1) = 1.550$
 $e_3 = -1/1.550 = -0.645$
 $f_3 = 2.04 - (-0.645)(-1) = 1.395$
 $e_4 = -1/1.395 = -0.717$
 $f_4 = 2.04 - (-0.717)(-1) = 1.323$





Gauss Seidel

- An Iterative or approximate methods provide an alternative to the elimination methods described to this point.
- Those approaches consists of guessing a value and then using a systematic method to obtain a refined estimate of the solution.

The Gauss-Seidel method is the most commonly used iterative method.

Gauss Seidel (1) $x_1 = \frac{b_1 - a_{12}x_2 - a_{13}x_3}{a_{11}}$ (2) $x_2 = \frac{b_2 - a_{21}x_1 - a_{23}x_3}{a_{22}}$ (3) $x_3 = \frac{b_3 - a_{31}x_1 - a_{32}x_2}{a_{33}}$

- 1. Assume all x's are zero.
- 2. Substitute in first equation to get x_1 .
- 3. Substitute in the second equation the new value of x_1 and the zero value of x_3 to get x_2 .
- 4. Repeat the process to calculate x_3 .
- 5. Repeat the steps 1 to 4 until:

$$|\varepsilon_{a,i}| = \left|\frac{x_i^j - x_i^{j-1}}{x_i^j}\right| 100\% < \varepsilon_s$$

Gauss Seidel

$$3x_1 - 0.1x_2 - 0.2x_3 = 7.85$$

$$0.1x_1 + 7x_2 - 0.3x_3 = -19.3$$

$$0.3x_1 - 0.2x_2 + 10x_3 = 71.4$$

$$\int_{x_1} = \frac{7.85 + 0.1x_2 + 0.2x_3}{3}$$

$$x_2 = \frac{-19.3 - 0.1x_1 + 0.3x_3}{7}$$

$$x_3 = \frac{71.4 - 0.3x_1 + 0.2x_2}{10}$$

True solution $x_1 = 3, x_2 = -2.5$, and $x_3 = 7$

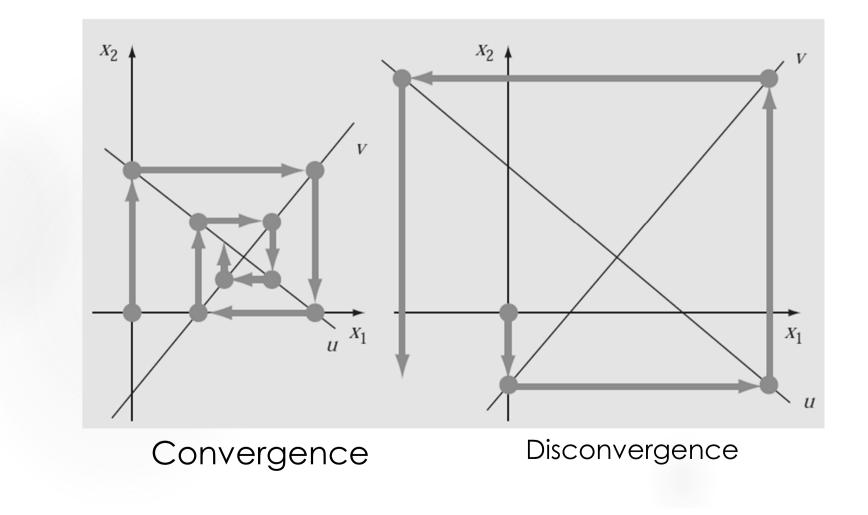
Gauss Seidel

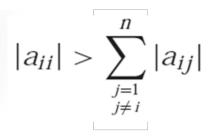


$$x_{1} = \frac{7.85 + 0.1x_{2} + 0.2x_{3}}{3}$$
$$x_{2} = \frac{-19.3 - 0.1x_{1} + 0.3x_{3}}{7}$$
$$x_{3} = \frac{71.4 - 0.3x_{1} + 0.2x_{2}}{10}$$

Trial	x ₁	X ₂	X ₃	
1	0	0	0	
2	2.616667	-2.794524	7.005610	
Error	100%	100%	100%	
3	2.990557	-2.499625	7.000291	
error	12.5%	11.8%	0.076%	

Sometimes, it is nonconvergent.





- The diagonal coefficient in each of the equations must be larger than the sum of the absolute values of the other coefficients in the equation.
- This criterion is sufficient but not necessary for convergence. That is, although the method may sometimes work if this equation is not met, convergence is guaranteed if the condition is satisfied.
- Systems where this equation holds are called diagonally dominant. Fortunately, many engineering problems of practical importance fulfill this requirement.

X2=8.5

X1=15.954

Sometimes, it is nonconvergent.

$$[A] = \begin{bmatrix} 11 & 13 \\ 11 & -9 \end{bmatrix}$$
 (Nonconvergent)
$$[A] = \begin{bmatrix} 11 & 13 \\ 11 & -9 \end{bmatrix}$$
 (Nonconvergent)
$$[11| < |13| \\ |9| < |11|$$

Trial	x ₁	x ₂		
1	0	0		
2	26	20.77778		
3	1.444442	-9.23457 34.1166		
4	36.91358			
5	-14.3196	-28.5018		
6	59.68395	61.94704		
7	-47.2101	-68.7013		
8	107.1924	120.013		

X1=15.954 X2=8.5

Sometimes, it is convergent, but very slow.

$$|x_1 - 9x_2 = 99 \qquad \implies [A] = \begin{bmatrix} 11 & -9 \\ 11 & 13 \end{bmatrix} \qquad \begin{array}{c} \text{Convergent} \\ |11| > |9| \\ |13| > |11| \end{array}$$

Trial	x ₁	x ₂	
1	0	0	
2	9	14.38462	
3	20.76923	4.426036	
4	12.6213	11.32044	But
5	18.26218	6.54739	very slow!
6	14.35696	9.851807	
7	17.06057	7.564134	
8	15.18884	9.147908	

- 2. When it converged, it often did so very slowly.
- > Relaxation represents a slight modification of the Gauss-Seidel method and is designed to enhance convergence.
- > After each new value of x is computed, that value is modified by a weighted average of the results of the previous and the present iterations: $x_i^{\text{new}} = \lambda x_i^{\text{new}} + (1 - \lambda) x_i^{\text{old}}$

where λ is a weighting factor that is assigned a value between 0 and 2.

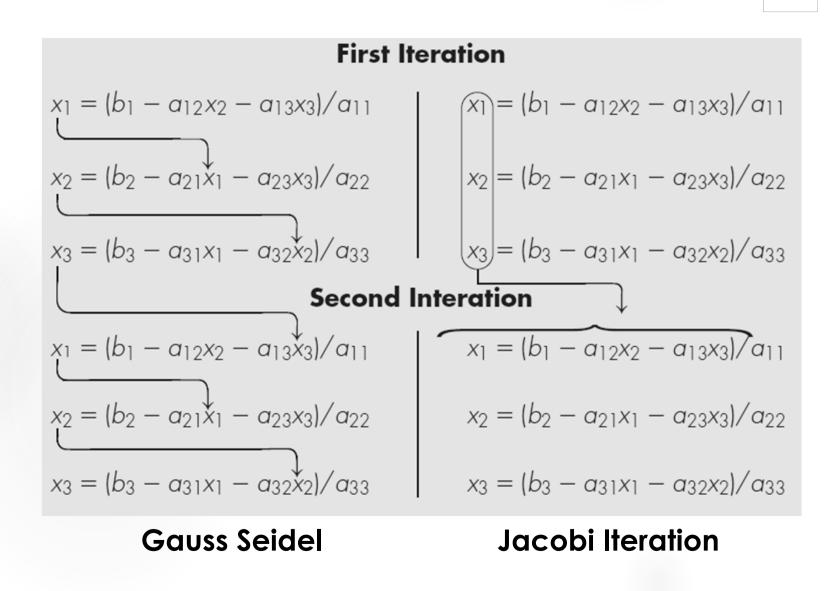
If $\lambda = 0$, the result is unmodified.

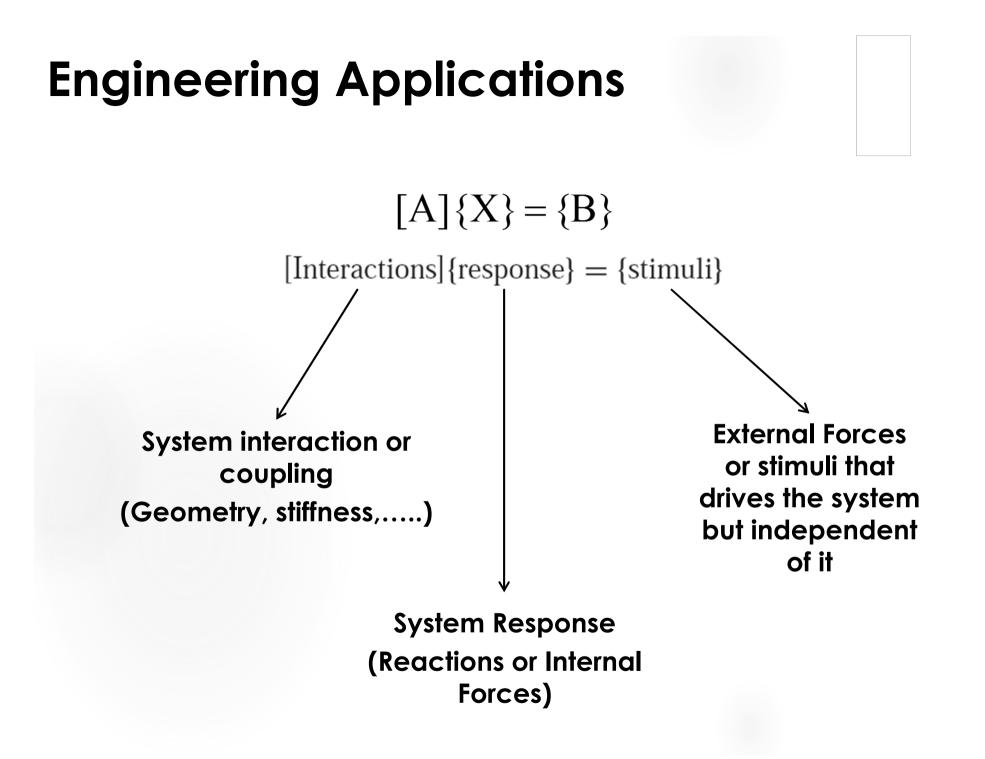
If $\lambda = 0 - 1$, "under relaxation" \rightarrow to make nonconvergent system converge If $\lambda = 1 - 2$, "successive overrelaxation," (SOR) \rightarrow to accelerate convergent system

Jacobi Iteration

- As each new x value is computed for the Gauss-Seidel method, it is immediately used in the next equation to determine another x value. Thus, if the solution is converging, the best available estimates will be employed.
- An alternative approach, called Jacobi iteration, utilizes a somewhat different tactic. Rather than using the latest available x's, this technique computes a set of new x's on the basis of a set of old x's.
- Thus, as new values are generated, they are not immediately used but rather are retained for the next iteration.

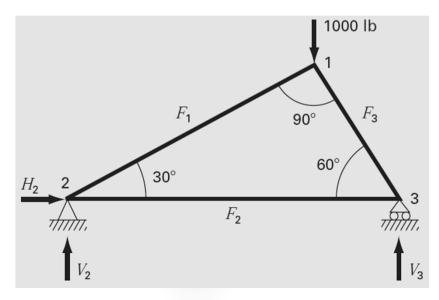
Jacobi Iteration/Gauss Seidel



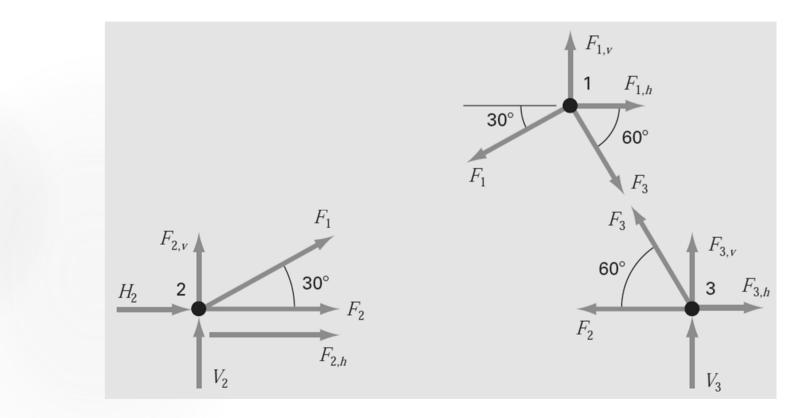


Find the forces and reactions associated with the shown truss.

- ► F₁, F₂, F₃ are the tension or compression on the members of the truss.
- ► H₂, V₂, V₃ are forces that characterize how the truss interacts with the supporting surface (reaction).
- ► 1000 lb is an external force $(F_{1,v})$



The sum of forces in both vertical and horizontal directions must be zero at each node.



Free body diagram

At node (1):

$$\Sigma F_H = 0 = -F_1 \cos 30^\circ + F_3 \cos 60^\circ + F_{1,h}$$

$$\Sigma F_V = 0 = -F_1 \sin 30^\circ - F_3 \sin 60^\circ + F_{1,v}$$

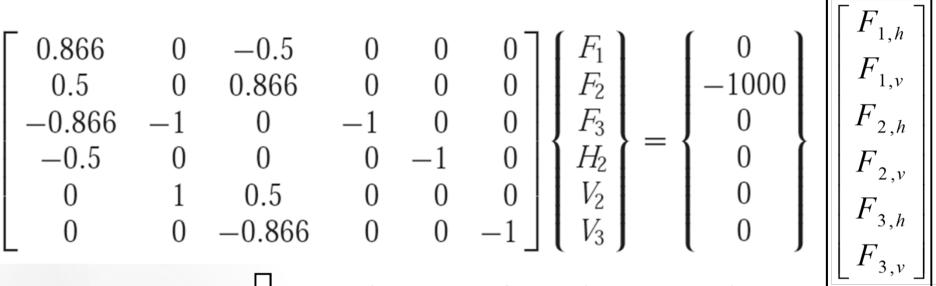
At node (2):

$$\Sigma F_H = 0 = F_2 + F_1 \cos 30^\circ + F_{2,h} + H_2$$

$$\Sigma F_V = 0 = F_1 \sin 30^\circ + F_{2,v} + V_2$$

At node (3):

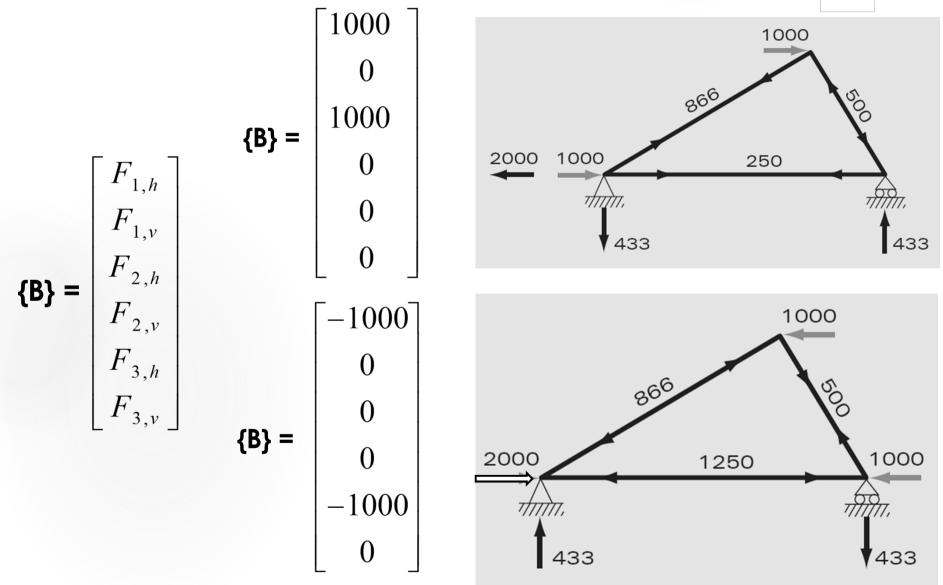
 $\Sigma F_H = 0 = -F_2 - F_3 \cos 60^\circ + F_{3,h}$ $\Sigma F_V = 0 = F_3 \sin 60^\circ + F_{3,v} + V_3$



Using any of previous techniques (LU is preferred for this problem, why?) (Partial pivoting may be required)

$F_1 = -500$	$F_2 = 433$	$F_3 = -866$
$H_2 = 0$	$V_2 = 250$	$V_3 = 750$





$$[A]^{-1} = \begin{bmatrix} 0.866 & 0.5 & 0 & 0 & 0 & 0 \\ 0.25 & -0.433 & 0 & 0 & 1 & 0 \\ -0.5 & 0.866 & 0 & 0 & 0 & 0 \\ -1 & 0 & -1 & 0 & -1 & 0 \\ -0.433 & -0.25 & 0 & -1 & 0 & 0 \\ 0.433 & -0.75 & 0 & 0 & 0 & -1 \end{bmatrix}$$

								$\begin{bmatrix} F \end{bmatrix}$
$\left(F_{1} \right)$		0.866	0.5	0	0	0	0	$\begin{bmatrix} \mathbf{I} & 1, h \\ \mathbf{F} \end{bmatrix}$
F_2		0.25	-0.433	0	0	1	0	$\Gamma_{1,v}$
F_3		-0.5	0.866	0	0	0	0	$ F_{2,h} $
H_2		-1	0	-1	0	-1	0	$ F_{2,v} $
V_2		-0.433	-0.25	0	-1	0	0	$ _{F}$
V_3		0.433	-0.75	0	0	0	-1	$\begin{bmatrix} 1 & 3, h \\ \mathbf{\Gamma} \end{bmatrix}$
()							-	$\begin{bmatrix} F_{3,v} \end{bmatrix}$