$$V_{avg} = \frac{1}{2} \left( V_{\min} + V_{\max} \right)$$

- $\bullet$  where  $V_{min}$  and  $V_{max}$  are the maximum and minimum values of V at the fixed nodes
- Doubts:
- Q. 1: How did we come up with the equation of alpha in slide 10?
- Answer:
- Shape functions are obtained from
- $V_e(x,y) = a + bx + cy$  where (a,b,c) are

$$\Rightarrow \begin{bmatrix} a \\ b \\ c \end{bmatrix} = \frac{1}{2A} \begin{bmatrix} x_2 y_3 - x_3 y_2 & x_3 y_1 - x_1 y_3 & x_1 y_2 - x_2 y_1 \\ y_2 - y_3 & y_3 - y_1 & y_1 - y_2 \\ x_3 - x_2 & x_1 - x_3 & x_2 - x_1 \end{bmatrix} \begin{bmatrix} V_{e1} \\ V_{e2} \\ V_{e3} \end{bmatrix}$$

• where A is the area of element e, i.e.,

$$A = \frac{1}{2} \begin{vmatrix} 1 & x_1 & y_1 \\ 1 & x_2 & y_2 \\ 1 & x_3 & y_3 \end{vmatrix} = \frac{1}{2} \{ (x_1 y_2 - x_2 y_1) + (x_3 y_1 - x_1 y_3) + (x_2 y_3 - x_3 y_2) \}$$

$$= \frac{1}{2} \{ (x_2 - x_1)(y_3 - y_1) - (x_3 - x_1)(y_2 - y_1) \}$$

We may express the below equation as

$$V_e(x,y) = a + bx + cy$$

$$V_e = \sum_{i=1}^{3} \alpha_i (x, y) V_{ei}$$

where

$$\alpha_{1} = \frac{1}{2A} \Big[ (x_{2}y_{3} - x_{3}y_{2}) + (y_{2} - y_{3})x + (x_{3} - x_{2})y \Big]$$

$$\alpha_{2} = \frac{1}{2A} \Big[ (x_{3}y_{1} - x_{1}y_{3}) + (y_{3} - y_{1})x + (x_{1} - x_{3})y \Big]$$

$$\alpha_{3} = \frac{1}{2A} \Big[ (x_{1}y_{2} - x_{2}y_{1}) + (y_{1} - y_{2})x + (x_{2} - x_{1})y \Big]$$

- Q. 2: Anticlockwise approach assigning the node numbers is not applicable for global nodes numbering?
- Ans: All the calculations for global coefficients are done from local elements coefficient matrix so no need
- Q3: Global coupling coefficient matrix (C) is always singular?
- Ans: Yes, for example, det (C)=0

$$\begin{bmatrix} C \end{bmatrix} = \begin{bmatrix} 1.2357 & -0.7786 & 0 & -0.4571 \\ -0.7786 & 1.25 & -0.4571 & -0.0143 \\ 0 & -0.4571 & 0.8238 & -0.3667 \\ -0.4571 & -0.0143 & -0.3667 & 0.8381 \end{bmatrix}$$

- Q4:
- In the last part, we are finding the partial derivative of W wrt Vk.
  - Does that mean we are finding the optimum value of V to have minimum energy?
  - If yes and if for a fixed node, if we are finding the value of Vk, does it not tamper with the original value of V, given for that particular node?
- Ans: Yes, we will find it only for free nodes

$$V_{avg} = \frac{1}{2} \left( V_{\min} + V_{\max} \right)$$

- where  $V_{min}$  and  $V_{max}$  are the maximum and minimum values of V at the fixed nodes
- (2) Band matrix method
- If all the free nodes (f) are numbered first and all prescribed nodes (p) last, we can write

$$W = \frac{1}{2} \varepsilon \begin{bmatrix} V_{f} & V_{p} \\ V_{f} & V_{p} \end{bmatrix} \begin{bmatrix} C_{ff} & C_{fp} \\ C_{pf} & C_{pp} \end{bmatrix} \begin{bmatrix} V_{f} \\ V_{p} \end{bmatrix} = \frac{1}{2} \varepsilon \begin{bmatrix} V_{f} C_{ff} + V_{p} C_{pf} \\ V_{f} C_{fp} + V_{p} C_{pp} \end{bmatrix} \begin{bmatrix} V_{f} \\ V_{p} \end{bmatrix}$$

$$W = \frac{1}{2} \varepsilon \begin{bmatrix} V_f C_{ff} V_f + V_p C_{pf} V_f + \\ V_f C_{fp} V_p + V_p C_{pp} V_p \end{bmatrix}$$

- What are free and prescribed nodes?
  - Nodes which are free
  - Nodes with the prescribed or fixed potentials

• Since V<sub>p</sub> is constant, we differentiate only w.r.t. V<sub>f</sub>, hence

$$\begin{bmatrix} C_{ff} & C_{fp} \end{bmatrix} \begin{bmatrix} V_f \\ V_p \end{bmatrix} = 0 \Rightarrow \begin{bmatrix} C_{ff} \end{bmatrix} \begin{bmatrix} V_f \end{bmatrix} = -\begin{bmatrix} C_{fp} \end{bmatrix} \begin{bmatrix} V_p \end{bmatrix}$$

This equation can be written as

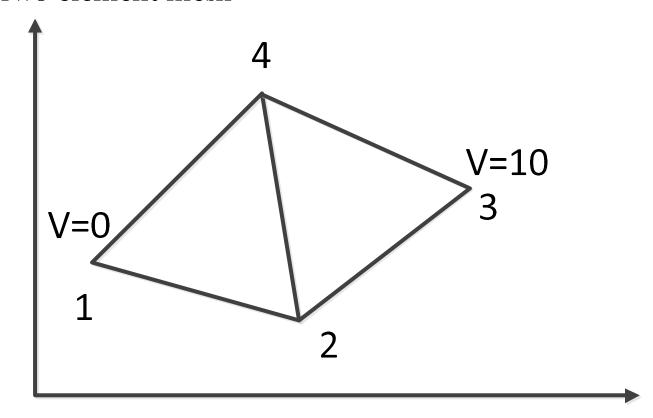
where

$$\begin{bmatrix} V \end{bmatrix} = \begin{bmatrix} V_f \end{bmatrix}; \begin{bmatrix} A \end{bmatrix} = \begin{bmatrix} C_{ff} \end{bmatrix}; \begin{bmatrix} B \end{bmatrix} = -\begin{bmatrix} C_{fp} \end{bmatrix} \begin{bmatrix} V_p \end{bmatrix}$$

- Since A is in general non-singular, the potential at the free nodes can be found out easily
- Example: Consider the two element mesh shown in Fig. Using FEM find the potentials within the mesh.

Node	$\mathbf{X}$	y
1	0.8	1.8
2	1.4	1.4
3	2.1	2.1
4	1.2	2.7

• Two element mesh



Define

$$P_{1} = (y_{2} - y_{3}), P_{2} = (y_{3} - y_{1}), P_{3} = (y_{1} - y_{2})$$

$$Q_{1} = -(x_{2} - x_{3}), Q_{2} = -(x_{3} - x_{1}), Q_{3} = -(x_{1} - x_{2})$$

• Then

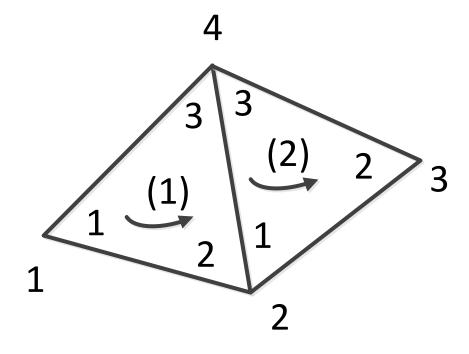
$$C_{ij}^{(e)} = \frac{1}{4A} \left( P_i P_j + Q_i Q_j \right)$$

where

$$A = \frac{1}{2} \left( P_2 Q_3 - P_3 Q_2 \right)$$

Element	1	2
$(P_1, P_2, P_3)$	(-1.3, 0.9, 0.4)	(-0.6, 1.3, -0.7)
$(Q_1,Q_2,Q_3)$	(2,4, 0.6)	(-0.9, 0.2, 0.7)
A	0.35	0.525

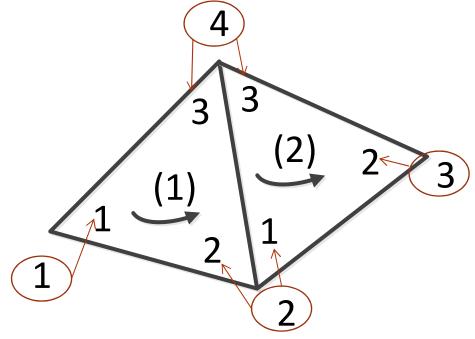
• Element numbering and local node numbering



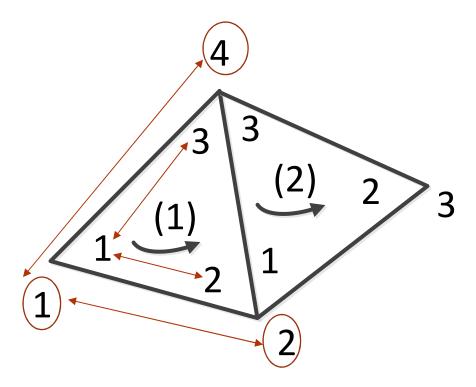
$$\begin{bmatrix} C^{(1)} \end{bmatrix} = \begin{bmatrix} 1.2357 & -0.7786 & -0.4571 \\ -0.7786 & 0.6929 & 0.0857 \\ -0.4571 & 0.0857 & 0.37114 \end{bmatrix}$$

$$\begin{bmatrix} C^{(2)} \end{bmatrix} = \begin{bmatrix} 0.5571 & -0.4571 & -0.1 \\ -0.4571 & 0.8238 & -0.3667 \\ -0.1 & -0.3667 & 0.4667 \end{bmatrix}$$

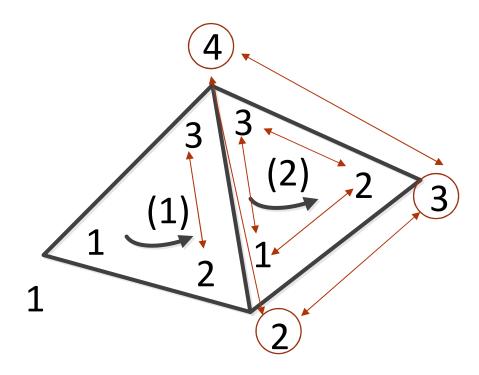
• Diagonal elements of Global coefficient matrix



$$\begin{bmatrix} C_{11} = C_{11}^{(1)} \\ C_{22} = C_{22}^{(1)} + C_{11}^{(2)} \\ C_{33} = C_{22}^{(2)} \\ C_{44} = C_{33}^{(1)} + C_{33}^{(2)} \end{bmatrix}$$



$$\begin{bmatrix} C_{11} = C_{11}^{(1)} & C_{12} = C_{12}^{(1)} & C_{13} = 0 & C_{14} = C_{13}^{(1)} \\ C_{21} = C_{12} & C_{22} = C_{22}^{(1)} + C_{11}^{(2)} \\ C_{31} = C_{13} & C_{33} = C_{22}^{(2)} \\ C_{41} = C_{14} & C_{44} = C_{33}^{(1)} + C_{33}^{(2)} \end{bmatrix}$$



$$\begin{bmatrix} C_{11} = C_{11}^{(1)} & C_{12} = C_{12}^{(1)} & C_{13} = 0 & C_{14} = C_{13}^{(1)} \\ C_{21} = C_{12} & C_{22} = C_{22}^{(1)} + C_{11}^{(2)} & C_{23} = C_{12}^{(2)} & C_{24} = C_{23}^{(1)} + C_{13}^{(2)} \\ C_{31} = C_{13} & C_{32} = C_{23} & C_{33} = C_{22}^{(2)} & C_{34} = C_{23}^{(2)} \\ C_{41} = C_{14} & C_{42} = C_{24} & C_{43} = C_{34} & C_{44} = C_{33}^{(1)} + C_{33}^{(2)} \\ \end{bmatrix}$$

$$\begin{bmatrix} C \end{bmatrix} = \begin{bmatrix} 1.2357 & -0.7786 & 0 & -0.4571 \\ -0.7786 & 1.25 & -0.4571 & -0.0143 \\ 0 & -0.4571 & 0.8238 & -0.3667 \\ -0.4571 & -0.0143 & -0.3667 & 0.8381 \end{bmatrix}$$

- Note that  $\sum_{i=1}^{4} C_{ij} = \sum_{j=1}^{4} C_{ij} = 0$
- It may be used to check if C is properly obtained
- For free nodes, we can found out as follows

First iteration (note  $V_1 = 0$ ,  $V_3 = 10$ , prescribed)  $: V_k = -\frac{1}{C_{kk}} \sum_{i=1, i \neq k}^{n} C_{ki} V_i$ 

$$V_{2} = -\frac{1}{C_{22}} \left( C_{21} V_{1} + C_{23} V_{3} + C_{24} V_{4} \right) = -\frac{1}{1.25} \left( -4.571 - 0.0143 V_{4} \right)$$

$$V_{4} = -\frac{1}{C_{44}} \left( C_{41} V_{1} + C_{42} V_{2} + C_{43} V_{3} \right) = -\frac{1}{0.831} \left( -3.667 - 0.0143 V_{2} \right)$$

Setting  $V_2 = V_4 = 0$  initially, we have,  $V_2 = 3.6568$ ,  $V_4 = 4.413$ 

Use this value for second iteration, we have,  $V_2=3.707$ ,  $V_4=4.476$ 

Band matrix method:

$$\begin{bmatrix} C_{22} & C_{24} \\ C_{42} & C_{44} \end{bmatrix} \begin{bmatrix} V_2 \\ V_4 \end{bmatrix} = - \begin{bmatrix} C_{21} & C_{23} \\ C_{41} & C_{43} \end{bmatrix} \begin{bmatrix} V_1 \\ V_3 \end{bmatrix} \qquad \because \begin{bmatrix} C_{ff} \end{bmatrix} \begin{bmatrix} V_f \end{bmatrix} = - \begin{bmatrix} C_{fp} \end{bmatrix} \begin{bmatrix} V_p \end{bmatrix}$$

$$\because \left[ C_{ff} \right] \left[ V_f \right] = - \left[ C_{fp} \right] \left[ V_p \right]$$

• Equivalently, 
$$\begin{bmatrix} 1 & 0 & 0 & 0 \\ 0 & C_{22} & 0 & C_{24} \\ 0 & 0 & 1 & 0 \\ 0 & C_{42} & 0 & C_{44} \end{bmatrix} \begin{bmatrix} V_1 \\ V_2 \\ V_3 \\ V_4 \end{bmatrix} = \begin{bmatrix} 1 & 0 \\ -C_{21} & -C_{23} \\ 0 & 1 \\ -C_{41} & -C_{43} \end{bmatrix} \begin{bmatrix} V_1 \\ V_3 \\ V_3 \end{bmatrix}$$

- Or,
- [C][V]=[B]
- where  $\begin{bmatrix} C \end{bmatrix} = \begin{bmatrix} 1 & 0 & 0 & 0 \\ 0 & 1.25 & 0 & -0.0143 \\ 0 & 0 & 1 & 0 \\ 0 & -0.0143 & 0 & 0.8381 \end{bmatrix}, \begin{bmatrix} B \end{bmatrix} = \begin{bmatrix} 0 \\ 4.571 \\ 10 \\ 3.667 \end{bmatrix}$

• Therefore,

$$\begin{bmatrix} V \end{bmatrix} = \begin{bmatrix} C \end{bmatrix}^{-1} \begin{bmatrix} B \end{bmatrix} = \begin{vmatrix} 0 \\ 3.708 \\ 10.0 \\ 4.438 \end{vmatrix}$$

- What is Functional?
- Consider a 1-D region of space in the range  $x_1 \le x \le x_2$
- A function  $\phi(x)$  and its derivative  $\phi'=d\phi/dx$  are defined in the region along with  $F(\phi,\phi',x)$
- A general expression of the functions and independent variable
- Consider the volume integral of the expression

$$I = \int_{x_1}^{x_2} F(\phi, \phi', x) dx$$

- The quantity I takes on a value for each choice of the function  $\boldsymbol{\phi}$
- It is called function of  $\phi$
- We seek a particular form of  $\phi$  that gives a maximum or minimum value of the functional
- Assume that  $\phi(x)$  has an arbitrary variation over the range but at the ends the values are clamped
- $\Phi(x_1) = \varphi_1, \ \Phi(x_2) = \varphi_2$
- To seek a minimum, we investigate variations of the functional form of  $\Phi(x)$

- Let  $\Phi_0(x)$  represent the desired minimizing function and define an arbitrary function  $\eta(x)$  that satisfies the boundary conditions
- $\eta(x_1) = \eta(x_2) = 0$
- Let  $\epsilon$  be a small number and add the quantity  $\epsilon \eta(x)$  to define the variation of  $\Phi(x)$  about the optimum point
- The functional becomes

$$I(\varepsilon) = \int_{x_1}^{x_2} F(\phi, \phi', x) dx = \int_{x_1}^{x_2} F(\phi_0 + \varepsilon \eta, \phi' + \varepsilon \eta', x) dx$$

• The condition  $dI/d\epsilon = 0$  means that  $\Phi_0(x)$  corresponds to the minimum (or maximum) of the functional

$$\frac{\partial I(\varepsilon)}{\partial \varepsilon} = \int_{x_1}^{x_2} \left( \frac{\partial F}{\partial \phi} \eta + \frac{\partial F}{\partial \phi'} \eta' \right) dx = 0$$

- Using integration by parts of the second term in the above integration
- Take

$$v = \frac{\partial F}{\partial \phi}$$
;  $u = \eta \Rightarrow du = \eta' dx$ 

Integration by parts

$$\int v du = vu - \int u dv$$

Therefore

$$\frac{\partial I(\varepsilon)}{\partial \varepsilon} = \int_{x_1}^{x_2} \left( \frac{\partial F}{\partial \phi} \eta - \eta \frac{d\left(\frac{\partial F}{\partial \phi'}\right)}{dx} \right) dx + \frac{\partial F}{\partial \phi'} \eta \Big|_{x_1}^{x_2} = 0$$

• Because of the boundary condition:  $\eta(x_1) = \eta(x_2) = 0$ 

$$\frac{\partial I(\varepsilon)}{\partial \varepsilon} = \int_{x_1}^{x_2} \left( \frac{\partial F}{\partial \phi} \eta - \eta \frac{d\left(\frac{\partial F}{\partial \phi'}\right)}{dx} \right) dx + \frac{\partial F}{\partial \phi'} \eta \Big|_{x_1}^{x_2} = 0$$

Hence,

$$\int_{x_1}^{x_2} \left( \frac{\partial F}{\partial \phi} - \frac{d \left( \frac{\partial F}{\partial \phi'} \right)}{dx} \right) \eta dx = 0$$

• The function  $\Phi(x)$  that gives minimum of the functional I therefore satisfies the equation which is called Euler equation

$$\frac{\partial F}{\partial \phi} - \frac{d\left(\frac{\partial F}{\partial \phi'}\right)}{dx} = 0; \phi' = \frac{d\phi}{dx}$$

• When the function  $\Phi(x,y,z)$  depends on three independent variables, the 3-D form of the Euler equation

$$\frac{\partial F}{\partial \phi} - \frac{\partial \left(\frac{\partial F}{\partial \phi_x}\right)}{\partial x} - \frac{\partial \left(\frac{\partial F}{\partial \phi_y}\right)}{\partial y} - \frac{\partial \left(\frac{\partial F}{\partial \phi_z}\right)}{\partial z} = 0$$

• where

$$\phi_x = \frac{d\phi}{dx}, \phi_y = \frac{d\phi}{dy}, \phi_z = \frac{d\phi}{dz}$$