

**Indian Institute of Science  
Bangalore**

**NP-TEL  
National Programme on  
Technology Enhanced Learning**

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**Course Title**

**Finite element method for structural dynamic  
And stability analyses**

**Lecture – 05**

**Beam Elements. Reference system. Assembly of Matrices.  
Imposition of BCS. Final equation of motion.**

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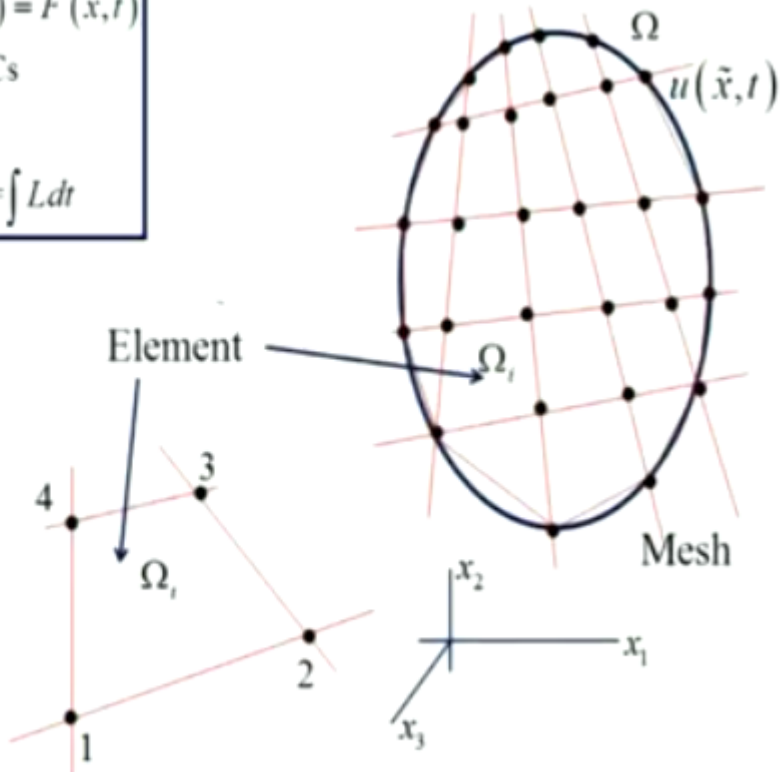
We have been talking about the development of finite element method for analyzing planar structures, so we will continue with the discussion.

$$\bar{L}u(\tilde{x},t) = F(\tilde{x},t)$$


BCs+ICs  
OR  
 $\text{Min } A = \int L dt$

**Recall**

$$\text{Mesh} \approx \Omega$$



In the last class we outline the basic framework for finite element modeling, so if  $\Omega$  is the domain that is of interest to us we partition the domain into a set of subdomains called Elements, and the union of these elements approximately equals the domain of interest and with this each domain is called an element, and what we do is we approximate the field variable that is a  $U(x_1,x_2,x_3,t)$

<b>FINITE</b>	$u(\tilde{x}, t) \text{ is approximated}$ $\text{in terms of}$ $u(\tilde{x}_i, t); i = 1, 2, \dots, N$
<b>ELEMENT</b>	$\tilde{\Omega} \approx \Omega = \bigcup_{i=1}^r \Omega_i; \Omega_i \cap \Omega_j = \emptyset \forall i \neq j$ <p>Within an element</p> $u(\tilde{x}, t) \text{ is approximated}$ $\text{as } u^e(\tilde{x}, t) = \sum_{i=1}^{N_e} N_i^e(\tilde{x}) u_i^e(\tilde{x}, t)$
	<p>An approximate numerical method to obtain solutions to PDE-s or variational problems</p>

$\Omega_i$ : Elements $\tilde{x}_i$ : Nodes $i = 1, 2, \dots, N$
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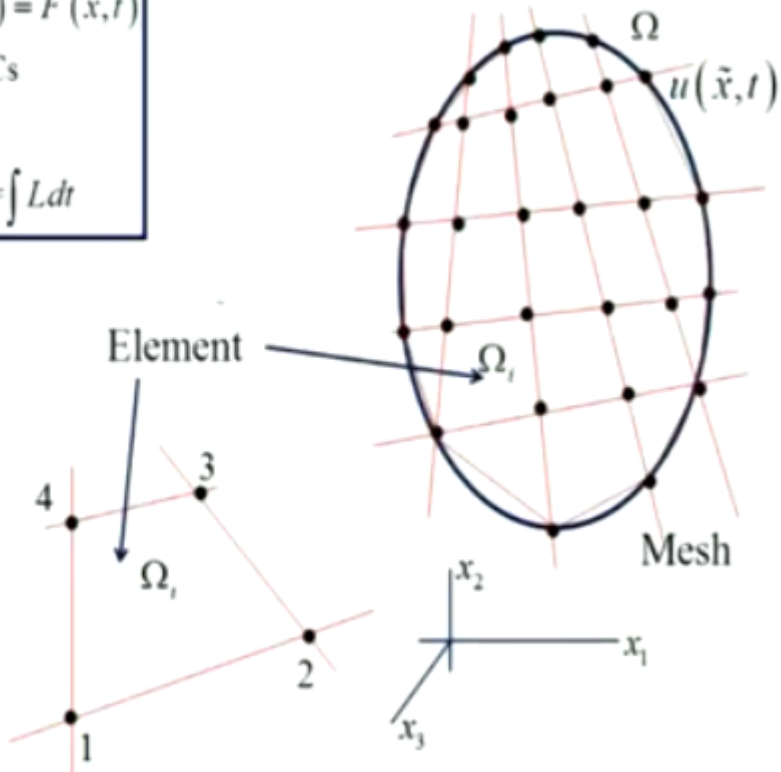
in terms of values of the field variables at the nodes, this at a set of N nodes they are these dots shown here and within an element we approximate the field variable in terms of the values at the nodes and by using interpolation. This  $NIE(\tilde{x})$  are the interpolation functions.

$$\bar{L}u(\tilde{x},t) = F(\tilde{x},t)$$

BCs+ICs  
OR  
 $\text{Min } A = \int L dt$

**Recall**

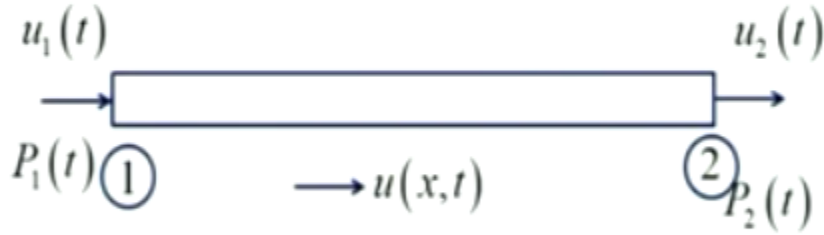
$$\text{Mesh} \approx \Omega$$



So the collection of these elements which approximates the domain of interest is known as the mesh, so this approach you know basically leads to an approximate numerical method to obtain solutions of problems posed as either in terms of a partial differential equation or in terms of a variational principle.



### Axially vibrating rod element



$$u(x,t) = u_1(t) \left(1 - \frac{x}{l}\right) + u_2(t) \frac{x}{l}$$

Now in the last class we considered the element behavior where the element was modeled as an actually vibrating rod element so it is a line element, these are the parameters of the element axial rigidity mass per unit length, length and damping and we assume that the cross section is symmetric, specifically we took that it is rectangular, and we identified two nodes  $X = 0$  and  $X = L$ , and the value of the field variable at  $X = X_1$  was  $U_1(t)$ , and the value of the field variable at  $X = L$  was  $U_2(t)$  and within this rod that the displacement field at any point  $X$  is expressed in terms of the nodal values  $U_1(t)$  and  $U_2(t)$  and these two functions  $1 - X/L$  and  $X/L$  are the interpolation functions.

Recall

Element level equation of motion

$$\frac{ml}{6} \begin{bmatrix} 2 & 1 \\ 1 & 2 \end{bmatrix} \begin{Bmatrix} \ddot{u}_1 \\ \ddot{u}_2 \end{Bmatrix} + \frac{AE}{l} \begin{bmatrix} 1 & -1 \\ -1 & 1 \end{bmatrix} \begin{Bmatrix} u_1 \\ u_2 \end{Bmatrix} = \begin{Bmatrix} P_1 \\ P_2 \end{Bmatrix}$$

$$\Rightarrow M\ddot{u} + Ku = P$$



We showed that using Lagrange's equation that means assuming this is a displacement field we can substitute this into the expression for kinetic energy and the potential energy and write the Lagrangian and apply the Lagrange's equation we get the element level equation of motion as shown here, and this  $U_1$  double dot and  $U_2$  double dot are the acceleration vector, is the acceleration vector, and  $U_1$ ,  $U_2$  this is a displacement vector, this  $M$  which is this is the element mass matrix, this is the element stiffness matrix, I also pointed out that we've use the same interpolation functions in computing elements of mass matrix as well as elements of stiffness matrix, therefore this mass matrix is called the consistent mass matrix.

## Euler-Bernoulli beam element



- Bending in the  $x$ - $y$  plane
- $x$ -axis coincides with the centroidal axis of the beam
- No coupling exists between bending and torsion

$$u(x, y, t) = -y \frac{\partial v}{\partial x}; v \equiv v(x, t)$$

$$\epsilon_{xx} = \frac{\partial u}{\partial x} = -y \frac{\partial^2 v}{\partial x^2}$$

$$2\epsilon_{xy} = \frac{\partial u}{\partial y} + \frac{\partial v}{\partial x} = -\frac{\partial v}{\partial x} + \frac{\partial v}{\partial x} = 0$$



Now we will continue with this discussion and we will now consider the behavior of an element which is modeled as an Euler-Bernoulli beam. So here we have again this is the line element this  $X$  and  $Y$  are the coordinate axis, and the displacement field along  $Y$  is denoted by  $V$ ,  $V$  is a function of  $X$  and  $T$ , and we make some of the standard assumptions that are made in Euler-Bernoulli beam theory, we are considering bending in the  $X$ - $Y$  plane and we assume that  $X$  axis coincides with the centroidal axis of the beam, and we also take in this study that there is no coupling between bending and torsion, a possibility of including this can be considered later, but right now we'll assume that there is no un-symmetric bending in the problem. So based on the standard assumption of Euler Bernoulli beam hypothesis that, for example plane sections which are normal to the neutral axis before bending remain plane and normal to the neutral axis after bending, we can postulate the form of the displacement field, so this is  $U$  in terms of  $V$  and using this we can compute the strains and that can subsequently be used in the expression for strain energy, and we get after a few simplifications the well-known expression for the strain energy due to bending.

$$\begin{aligned}
 V &= \frac{1}{2} \int_v \sigma_{xx} \varepsilon_{xx} dx \\
 &= \frac{1}{2} \int_0^L \int_A E \varepsilon_{xx}^2 dx dA \\
 &= \frac{1}{2} \int_0^L \int_A E \left( -y \frac{\partial^2 v}{\partial x^2} \right)^2 dx dA \\
 &= \frac{1}{2} \int_0^L EI_z \left( \frac{\partial^2 v}{\partial x^2} \right)^2 dx \text{ with } I_z = \int_A y^2 dA
 \end{aligned}$$

$$\begin{aligned}
 T &= \frac{1}{2} \int_0^L \int_A \rho \dot{v}^2(x,t) dx dA \\
 &= \frac{1}{2} \int_0^L m \dot{v}^2(x,t) dx \text{ with } m = \int_A \rho dA
 \end{aligned}$$



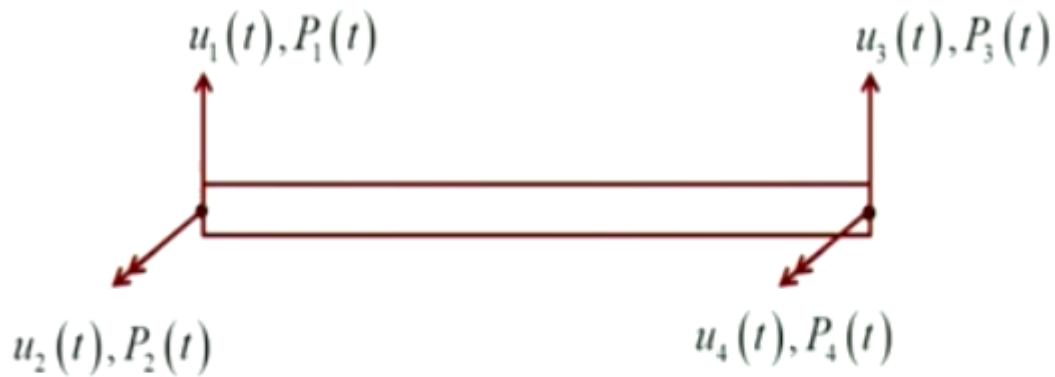
$$\begin{aligned}
 L &= T(t) - V(t) \\
 &= \frac{1}{2} \int_0^L m \dot{v}^2(x,t) dx - \frac{1}{2} \int_0^L EI_z \left( \frac{\partial^2 v}{\partial x^2} \right)^2 dx
 \end{aligned}$$

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So  $I_z$  is the moment of inertia, area moment of inertia about the Z axis this is given here. Kinetic energy is mass into velocity square integral over the domain half of that and if we use now  $\rho A$ , integral  $\rho A$  over  $A$  that is mass per unit length we get the kinetic energy in terms of mass per unit length and the velocity field.

Now based on this we can now construct the Lagrangian which is given in this expression, so this is a kinetic energy, this is the strain energy.





2-noded element with 2 dofs per node

$$v(x,t) = u_1(t)\phi_1(x) + u_2(t)\phi_2(x) + u_3(t)\phi_3(x) + u_4(t)\phi_4(x)$$



Now to obtain a finite element discretization we will now take that the element is made up of two nodes and at each node the translation and the gradient that is the slope are taken to be the nodal degrees of freedom, this is  $U_1$  and  $U_2(t)$  are the values of,  $U_1(t)$  is the field variable displacement at  $X = 0$ , this is  $\text{Dou } E / \text{Dou } X$  at  $X = 0$ , similarly  $U_3(t)$  is  $V(x,t)$  at  $X = L$ , and this  $U_4(t)$  is  $\text{Dou } V / \text{Dou } X$  at  $X = L$ . Now this element is a 2 noded element with two degrees of freedom per node and it has 4 degrees of freedom. Now the displacement field within the element as before we are going to now express in terms of the nodal degrees of freedom  $U_1(t)$ ,  $U_2(t)$ ,  $U_3(t)$ ,  $U_4(t)$  as shown here. Here  $\Phi_1$ ,  $\Phi_2$ ,  $\Phi_3$ ,  $\Phi_4$  are the interpolation functions which we need to now select.

$$v(x,t) = u_1(t)\phi_1(x) + u_2(t)\phi_2(x) + u_3(t)\phi_3(x) + u_4(t)\phi_4(x)$$

$u_i(t); i = 1, 2, 3, 4$ : Generalized coordinates

$\phi_i(x); i = 1, 2, 3, 4$ : Trial functions

$$v(0,t) = u_1(t)$$

$$\Rightarrow v(0,t) = u_1(t)\phi_1(0) + u_2(t)\phi_2(0) + u_3(t)\phi_3(0) + u_4(t)\phi_4(0)$$

Take  $\phi_1(0) = 1; \phi_2(0) = 0; \phi_3(0) = 0; \phi_4(0) = 0$

$$v'(0,t) = u_2(t)$$

$$\Rightarrow v'(0,t) = u_1(t)\phi_1'(0) + u_2(t)\phi_2'(0) + u_3(t)\phi_3'(0) + u_4(t)\phi_4'(0)$$

Take  $\phi_1'(0) = 0; \phi_2'(0) = 1; \phi_3'(0) = 0; \phi_4'(0) = 0$

Now this  $U_i(t)$  for  $i = 1, 2, 3, 4$  are the generalized coordinates and this  $\phi_i(x)$  are the trial functions. Now let us impose the boundary conditions that we already know at  $X = 0$  and  $X = L$  at  $X = 0$  we know that  $V(0,t)$  is  $U_1(t)$ , now that would mean  $V(0,t)$  is now if I substitute in this expression the  $U_1(t)$ ,  $T \phi_1(0) + U_2(t) \phi_2(0)$  etcetera. Now I know that this must be equal to  $U_1(t)$ , so that can be achieved by inspecting, if I now take  $\phi_1(0)$  to be 1, and  $\phi_2(0)$ ,  $\phi_3(0)$  and  $\phi_4(0)$  to be 0, I will be able to satisfy this requirement.

Similarly if we now consider  $\text{Dou } V / \text{Dou } X$  at  $X = 0$ , I know that it is  $U_2(t)$ , now let us again use that  $V$  prime of  $0, T$  is in terms of this, in terms of the assumed representation it is given in terms of, it is given by this equation, so again in order to satisfy this we can select  $\phi_1$  prime(0) to be 0,  $\phi_2$  prime (0) to be 1, and this  $\phi_3$  prime (0),  $\phi_4$  prime (0) to be 0, so this will ensure that this condition is met.

$$v(l,t) = u_3(t)$$

$$\Rightarrow v(l,t) = u_1(t)\phi_1(l) + u_2(t)\phi_2(l) + u_3(t)\phi_3(l) + u_4(t)\phi_4(l)$$

$$\text{Take } \phi_1(l) = 0; \phi_2(l) = 0; \phi_3(l) = 1; \phi_4(0) = 0$$

$$v'(l,t) = u_4(t)$$

$$\Rightarrow v'(l,t) = u_1(t)\phi_1'(l) + u_2(t)\phi_2'(l) + u_3(t)\phi_3'(l) + u_4(t)\phi_4'(l)$$

$$\text{Take } \phi_1'(l) = 0; \phi_2'(l) = 1; \phi_3'(l) = 0; \phi_4'(l) = 1$$

This ensures that we have satisfied the requirements on displacements at  $x = 0$  &  $x = l$



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Similarly I can impose condition at  $X = L$  where  $V(l,t)$  is  $U_3(t)$  and here we will require  $\phi_1(l)$  to be 0,  $\phi_2(l)$  to be 0,  $\phi_3(l)$  to be 1, and  $\phi_4(0)$  to be 0. Finally  $\text{Dou } V / \text{Dou } X$  at  $X = L$ , this is  $U_4(t)$  and to achieve this we have to take this conditions on  $\phi$ 's that is  $\phi_1'(l)$  is 0, this is 0, this is 0, but this has to be one. So this choice if we now select  $\phi(x)$  so that this set of 4 conditions which represents 16 conditions in total, now if we satisfy that we would ensure that we have satisfied the requirements on displacements at  $X = 0$ , and  $X = L$ .

Let us consider the equilibrium of the beam under the action of support displacements

$$u_i(t); i = 1, 2, 3, 4$$

$$\Rightarrow \phi_1(0) = 1; \phi_2(0) = 0; \phi_3(0) = 0; \phi_4(0) = 0$$

$$\Rightarrow EI \frac{d^4 \phi_1}{dx^4} = 0; \phi_1(0) = 1; \phi_2(0) = 0; \phi_3(0) = 0; \phi_4(0) = 0$$

$$\Rightarrow \phi_1(x) = ax^3 + bx^2 + cx + d \Rightarrow \phi_1(x) = 1 - 3\frac{x^2}{l^2} + 2\frac{x^3}{l^3}$$

Similarly, we get

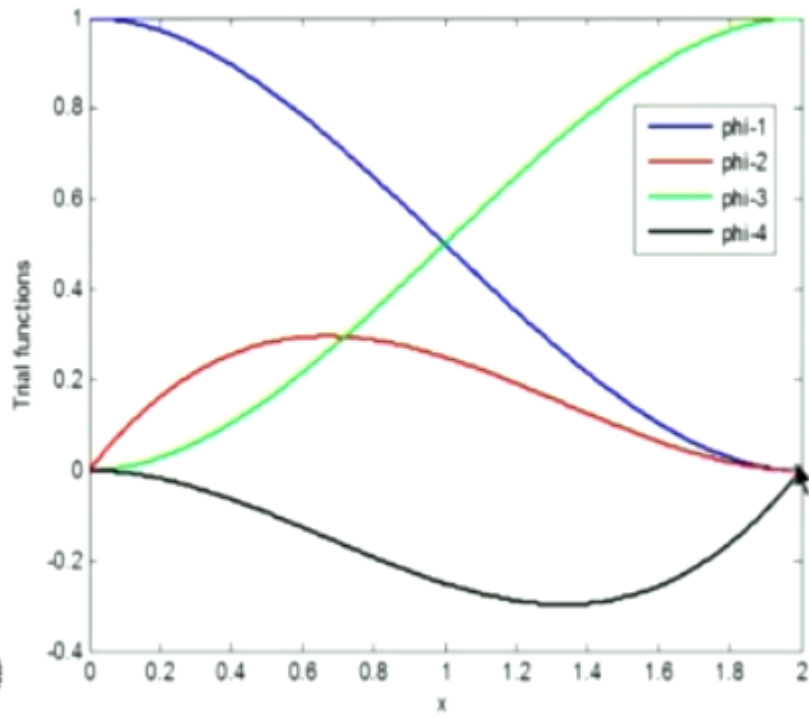
$$\phi_2(x) = x - 2\frac{x^2}{l} + \frac{x^3}{l^2};$$

$$\phi_3(x) = 3\frac{x^2}{l^2} - 2\frac{x^3}{l^3};$$

$$\phi_4(x) = -\frac{x^2}{l} + \frac{x^3}{l^2}$$



Now we can, to now make the choice for Phi(x) we can now consider the equilibrium of the beam under the action of support displacements  $U_i(t)$ ,  $i = 1, 2, 3, 4$ , so for Phi 1 for example we have 4 boundary conditions, so now if I write this equation for Phi 1 we get Phi 1(x) to be a cubic polynomial, and for Phi 1 I have 4 conditions that we have derived, now if I impose that I will get Phi 1(x) to be this 1, using the same logic that means assume a cubic polynomial with 4 constants A, B, C, D, we can derive Phi 2(x), Phi 3(x) and Phi 4(x), okay, now how do they look like so Phi 1(x) is this blue line, so at  $X = 0$  it is 1, its slope is 0, and at  $X = L$  the value is 0 is slope is also 0, similarly I have Phi 2(x) the slope here is 1 and the value is 0 and here again



the value on the slope are 0, so we define Phi 3 and Phi 4 in the same, so these are the interpolation functions that we use to represent the beam behavior.

$$\Rightarrow v(x,t) = u_1(t) \left( 1 - 3 \frac{x^2}{l^2} + 2 \frac{x^3}{l^2} \right) + u_2(t) \left( x - 2 \frac{x^2}{l} + \frac{x^3}{l^2} \right)$$

$$+ u_3(t) \left( 3 \frac{x^2}{l^2} - 2 \frac{x^3}{l^2} \right) + u_4(t) \left( -\frac{x^2}{l} + \frac{x^3}{l^2} \right)$$

$$T(t) = \frac{1}{2} \int_0^l m \dot{v}^2(x,t) dx = \frac{1}{2} \int_0^l m \left\{ \sum_{i=1}^4 \dot{u}_i(t) \phi_i(x) \right\}^2 dx$$

$$= \frac{1}{2} \frac{ml}{420} [156\dot{u}_1^2 + 4l^2\dot{u}_2^2 + 156\dot{u}_3^2 + 4l^2\dot{u}_4^2 + 44l\dot{u}_1\dot{u}_2$$

$$+ 108\dot{u}_1\dot{u}_3 - 26l\dot{u}_1\dot{u}_4 + 26l\dot{u}_2\dot{u}_3 - 6l^2\dot{u}_2\dot{u}_4 - 44l\dot{u}_3\dot{u}_4]$$

$$V(t) = \int_0^l \frac{1}{2} EI \left( \frac{\partial^2 v}{\partial x^2} \right)^2 dx = \int_0^l \frac{1}{2} EI \left( \sum_{i=1}^4 u_i(t) \phi_i''(x) \right)^2 dx$$

$$= \frac{1}{2} \frac{EI}{l^3} [12u_1^2 + 4l^2u_2^2 + 12u_3^2 + 4l^2u_4^2 + 12lu_1u_2 - 24u_1u_3$$

$$+ 12lu_1u_4 - 12lu_2u_3 + 24l^2u_2u_4 - 12lu_3u_4]$$



Now, I have now the approximation for the displacement field in terms of nodal values U1, U2, U3, U4 and these 4 interpolation functions. So now I can substitute this into the expression for kinetic energy and the potential energy. Suppose now if I substitute into the kinetic energy I have this is the expression for kinetic energy so this representation leads to this and noting that these Phi(x) are all polynomials you will easily see that in this case we will be able to integrate carry out these integrations in closed form, so the expression for kinetic energy in terms of the nodal degrees of freedom we obtain as shown here, so you have to carry out these integrations which are straightforward.

Similarly the expression for strain energy is given by this and here I substitute for Dou square V / Dou X square which is UI(t) Phi I double prime (x) summed over I = 1 to 4, so this again Phi I double prime (x) will again be polynomials and therefore carrying out this integration is a straightforward exercise and I will get this as a expression for the strain energy. So now my Lagrangian is ready T - V in terms of nodal degrees of freedom, there are 4 nodal degrees of freedom so this is a system with 4 degrees of freedom and 4 generalized coordinates,

$$L = T - V$$

$$\frac{d}{dt} \left( \frac{\partial L}{\partial \dot{u}_i} \right) - \frac{\partial L}{\partial u_i} = 0; i = 1, 2, 3, 4$$

$$\Rightarrow M\ddot{U} + KU = P$$

$$M = \frac{ml}{420} \begin{bmatrix} 156 & 22l & 54 & -13l \\ 22l & 4l^2 & 13l & -3l^2 \\ 54 & 13l & 156 & -22l \\ -13l & -3l^2 & -22l & 4l^2 \end{bmatrix} \quad \& \quad K = \frac{EI}{l^3} \begin{bmatrix} 12 & 6l & -12 & 6l \\ 6l & 4l^2 & -6l & 2l^2 \\ -12 & -6l & 12 & -6l \\ 6l & 2l^2 & -6l & 4l^2 \end{bmatrix}$$

### Remarks

- $M$  = element consistent mass matrix;  $M = M^T$
- $K$  = element stiffness matrix;  $K = K^T$
- Two noded beam element with two dofs per node
- $K$  and  $M$  are non-diagonal
- $V = \frac{1}{2} \dot{U}^T(t) K U'(t)$  &  $T = \frac{1}{2} \dot{U}^T(t) M \dot{U}'(t)$



so I need to apply the Lagrange's equation on these 4 variables, so if I do that I get the equation in the form  $M\ddot{U} + KU = P$ , where  $M\ddot{U}$  is a 4 cross 1 vector,  $U$  is 4 cross 1,  $\ddot{U}$  is 4 cross 1, so this is the vector of nodal displacements, this is a vector of nodal accelerations,  $M$  which is a 4/4 symmetric matrix is the element consistent mass matrix,  $K$  which is this 4/4 symmetric element stiffness matrix, so the  $M$  is a element consistent mass matrix it is symmetric,  $K$  is element stiffness matrix which is symmetric, so the element that we are talking about therefore is a 2 noded beam element with 2 degrees of freedom per node, and  $K$  and  $M$  we'll see that they are non-diagonal and we can express the expression for strain energy and kinetic energy can be obtained in terms of nodal displacements and nodal velocities as shown here, this can be verified. So we have now been able to derive the mass and stiffness matrix for an Euler Bernoulli beam element.

Consistent mass matrix

$$M = \frac{ml}{420} \begin{bmatrix} 156 & 22l & 54 & -13l \\ 22l & 4l^2 & 13l & -3l^2 \\ 54 & 13l & 156 & -22l \\ -13l & -3l^2 & -22l & 4l^2 \end{bmatrix}$$

Lumped mass matrix

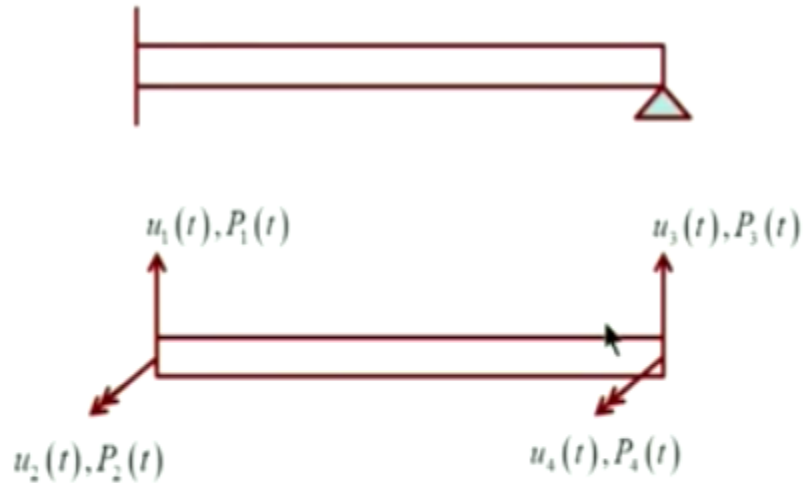
$$M = \frac{ml}{2} \begin{bmatrix} 1 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 \end{bmatrix}$$



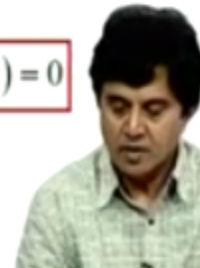
Now again let us go back to the nomenclature, I call this mass matrix as consistent mass matrix, there is an another way of forming mass matrix which is known as lumped mass matrix, for example if this is a beam element and if half of this mass is lumped here and this half is lumped here I get a mass matrix which is as shown here, the half the mass is lumped here, this is 1, and half the mass is lumped here this is 1, so this obviously doesn't take into account various cross-coupling terms that are present in the consistent mass matrix.



## Example



$$u_1(t) = 0; u_2(t) = 0; u_3(t) = 0; P_4(t) = 0$$



Now we will consider some simple examples so that we understand how to use these matrices, suppose if I have a proper cantilever, so I model this element as this beam using only one element therefore there is a 2 noded beam element with 2 degrees of freedom per node, now the imposition of the boundary conditions for this problem results in this statement  $U_1(t)$  is 0 because this displacement is 0,  $U_2(t)$  is 0 because the slope is 0 is a clamped end and similarly  $U_3(t)$  which is the displacement here which is 0,  $P_4(t)$  which is the bending moment here which is 0, so now if I write the element level equation of motion I get now for the acceleration vector the first 3 terms would be 0 and similarly for the displacement vector the first 3 terms will be 0 this  $U_4$  is the unknown associated with displacement which are 0 at the boundary, there will be a stress reaction, for example there will be shear force and bending moment here and there will be shear force here, and they are, this  $R_1, R_2, R_3$ , so this equation therefore is a set of 4 equations, 3 for the reactions, and 1 for the unknown displacement.

$$\frac{ml}{420} \begin{bmatrix} 156 & 22l & 54 & -13l \\ 22l & 4l^2 & 13l & -3l^2 \\ 54 & 13l & 156 & -22l \\ -13l & -3l^2 & -22l & 4l^2 \end{bmatrix} \begin{Bmatrix} 0 \\ 0 \\ 0 \\ \ddot{u}_4 \end{Bmatrix} + \frac{EI}{l^3} \begin{bmatrix} 12 & 6l & -12 & 6l \\ 6l & 4l^2 & -6l & 2l^2 \\ -12 & -6l & 12 & -6l \\ 6l & 2l^2 & -6l & 4l^2 \end{bmatrix} \begin{Bmatrix} 0 \\ 0 \\ 0 \\ u_4 \end{Bmatrix} = \begin{Bmatrix} R_1 \\ R_2 \\ R_3 \\ 0 \end{Bmatrix}$$

$$\frac{ml}{420} 4l^2 \ddot{u}_4 + \frac{EI}{l^3} 4l^2 u_4 = 0$$

$$\omega_1 = 20.49 \sqrt{\frac{EI}{ml^4}}; \omega_1^{\text{Exact}} = 15.41 \sqrt{\frac{EI}{ml^4}}$$

Error: 32.98%

So now we solve for the equation for the unknown displacement I get this, so you have to use the equation corresponding to the last row of this equation, so you get this single equation. Now this is a single degree freedom system in this approximation the beam is approximated as a single degree of freedom system and the generalized coordinate is  $U_4(t)$ , once I know  $U_4(t)$  I will be able to write the expression for the beam displacement, how do we do that?

$$\frac{ml}{420} \begin{bmatrix} 156 & 22l & 54 & -13l \\ 22l & 4l^2 & 13l & -3l^2 \\ 54 & 13l & 156 & -22l \\ -13l & -3l^2 & -22l & 4l^2 \end{bmatrix} \begin{Bmatrix} 0 \\ 0 \\ 0 \\ \ddot{u}_4 \end{Bmatrix} + \frac{EI}{l^3} \begin{bmatrix} 12 & 6l & -12 & 6l \\ 6l & 4l^2 & -6l & 2l^2 \\ -12 & -6l & 12 & -6l \\ 6l & 2l^2 & -6l & 4l^2 \end{bmatrix} \begin{Bmatrix} 0 \\ 0 \\ 0 \\ u_4 \end{Bmatrix} = \begin{Bmatrix} R_1 \\ R_2 \\ R_3 \\ 0 \end{Bmatrix}$$

$$\frac{ml}{420} 4l^2 \ddot{u}_4 + \frac{EI}{l^3} 4l^2 u_4 = 0$$

$$\omega_1 = 20.49 \sqrt{\frac{EI}{ml^4}}; \omega_1^{\text{Exact}} = 15.41 \sqrt{\frac{EI}{ml^4}}$$

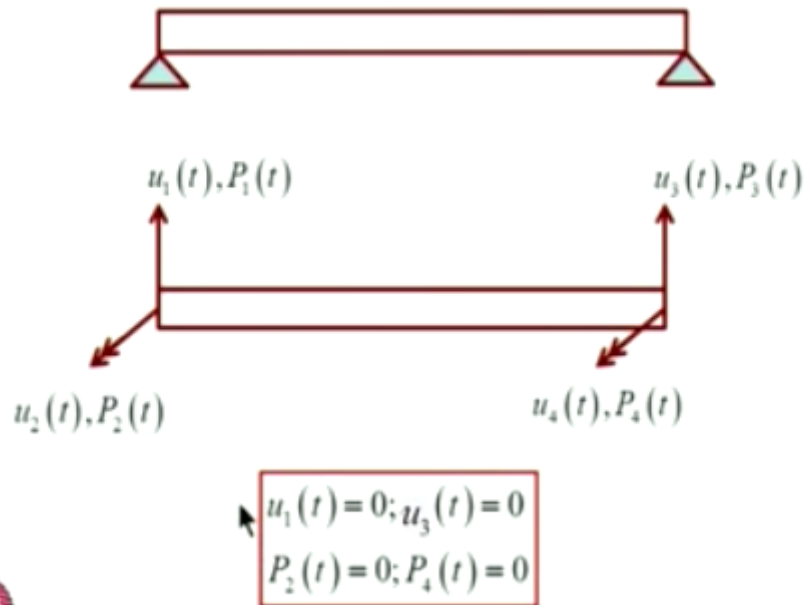
Error: 32.98%



$$\theta(x,t) = U_4(t) \Phi_4(x)$$

That is all  $\Phi_4(x)$  is a polynomial, cubic polynomial that we have already derived, so this is a displacement field for all values of  $X$  and this  $U_4(t)$  itself is governed by this differential equation, so from this you can deduce the natural frequency of the system is a single degree freedom system, square root  $K/M$  is the natural frequency if we do that we get this as answer  $20.49 \sqrt{EI/MI}$  to the power of 4. Now an exact solution to the natural frequency of this beam is available and if we reproduce here and by comparing these two we see that in this approximation the first natural frequency computation has an error of about 33%.

### Example



Now a similar exercise we can do for a simply supported beam, again this beam I approximate using only one element, so here you see that  $U_1$  is 0, because it is simply supported end,  $U_3$  is 0 and  $P_2(t)$ , what is  $P_2(t)$ ? The bending moment here which is 0 and  $P_4(t)$  which is the bending moment that is 0, so what are the unknowns in this problem?  $U_2$  and  $U_4$  and the reactions shear force at  $X = 0$  and  $X = L$ , which is  $P_3$  and  $P_1$ . So now if you write the element level equation of motion, we get this is equilibrium equation for the element, now for the beam now because we

$$\frac{ml}{420} \begin{bmatrix} 156 & 22l & 54 & -13l \\ 22l & 4l^2 & 13l & -3l^2 \\ 54 & 13l & 156 & -22l \\ -13l & -3l^2 & -22l & 4l^2 \end{bmatrix} \begin{Bmatrix} 0 \\ \ddot{u}_2 \\ 0 \\ \ddot{u}_4 \end{Bmatrix} + \frac{EI}{l^3} \begin{bmatrix} 12 & 6l & -12 & 6l \\ 6l & 4l^2 & -6l & 2l^2 \\ -12 & -6l & 12 & -6l \\ 6l & 2l^2 & -6l & 4l^2 \end{bmatrix} \begin{Bmatrix} 0 \\ u_2 \\ 0 \\ u_4 \end{Bmatrix} = \begin{Bmatrix} R_1 \\ 0 \\ R_3 \\ 0 \end{Bmatrix}$$

$$\Rightarrow \frac{ml}{420} \begin{bmatrix} 4l^2 & -3l^2 \\ -3l^2 & 4l^2 \end{bmatrix} \begin{Bmatrix} \ddot{u}_2 \\ \ddot{u}_4 \end{Bmatrix} + \frac{EI}{l^3} \begin{bmatrix} 4l^2 & 2l^2 \\ 2l^2 & 4l^2 \end{bmatrix} \begin{Bmatrix} u_2 \\ u_4 \end{Bmatrix} = \begin{Bmatrix} 0 \\ 0 \end{Bmatrix}$$

The remaining two equations govern the two unknown reactions:

$$\frac{ml}{420} \begin{bmatrix} 22l & -13l \\ 13l & -22l \end{bmatrix} \begin{Bmatrix} \ddot{u}_2 \\ \ddot{u}_4 \end{Bmatrix} + \frac{EI}{l^3} \begin{bmatrix} 6l & 6l \\ -6l & -6l \end{bmatrix} \begin{Bmatrix} u_2 \\ u_4 \end{Bmatrix}$$

NPTEL



use one element approximation, so this is U1 and U3 as 0, U2 and U4 are nonzero, so whereas R1 and R3 are the unknown reactions which we have to compute, so again this is set of 4 equations, 2 for the unknown displacements U2 and U4, and 2 for the unknown reactions R1 and R3, now if you extract from this the equation for U2 and U4 I get a 2 degree of freedom approximation to the simply supported beam structure. The remaining 2 equations govern you know they are the equations for the unknown reaction, so though if you want to find the reactions you have to first solve the equation for the displacements and substitute here.

### Natural frequencies

$$\frac{EI}{l^3} \begin{bmatrix} 4l^2 & 2l^2 \\ 2l^2 & 4l^2 \end{bmatrix} \begin{Bmatrix} u_{20} \\ u_{40} \end{Bmatrix} = \frac{ml\omega^2}{420} \begin{bmatrix} 4l^2 & -3l^2 \\ -3l^2 & 4l^2 \end{bmatrix} \begin{Bmatrix} u_{20} \\ u_{40} \end{Bmatrix}$$

$$\begin{bmatrix} 4 & 2 \\ 2 & 4 \end{bmatrix} \begin{Bmatrix} u_{20} \\ u_{40} \end{Bmatrix} = \lambda \begin{bmatrix} 4 & -3 \\ -3 & 4 \end{bmatrix} \begin{Bmatrix} u_{20} \\ u_{40} \end{Bmatrix}; \lambda = \frac{ml^4\omega^2}{420EI}$$

$$\Rightarrow \lambda = \frac{2}{7}, 6 \Rightarrow \omega_1 = 10.95 \sqrt{\frac{EI}{ml^4}} \text{ \& } \omega_2 = 50.20 \sqrt{\frac{EI}{ml^4}} \text{ (rad/s)}$$

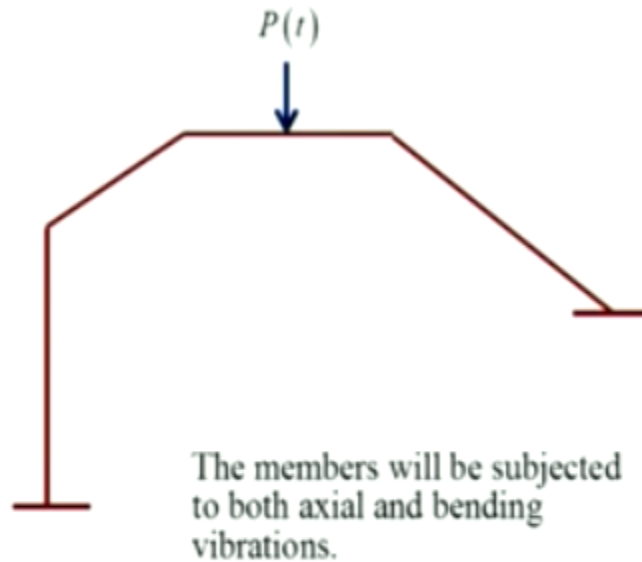
$$\omega_1^{\text{Exact}} = 9.87 \sqrt{\frac{EI}{ml^4}} \text{ \& } \omega_2^{\text{Exact}} = 39.48 \sqrt{\frac{EI}{ml^4}} \text{ (rad/s)}$$

Errors: 10.94% & 27.15%



Now let's now look at this 2 degree of freedom system, and if we are interested in finding natural frequencies we have to solve the eigenvalue problem, this is the statement of the eigenvalue problem that means we assume that the two generalized coordinates oscillate harmonically at the same frequency  $\omega$  and for that to happen this condition must be satisfied, and if we now introduce a parameter  $\lambda$ , that means I will divide by  $EI$  by  $L$  cube and call this parameter as  $\lambda$ , the eigenvalues now expressed in terms of  $\lambda$  turn out to be  $2/7$  and  $6$  from which I will compute the first and the second natural frequency, it is a 2 degree freedom system therefore I will get an approximation to first 2 natural frequencies. So these turn out to be  $10.95$  into this factor and  $50.20$  into this factor.

Now again this problem is well known to have the exact solution, and the exact solution is for the first natural frequency instead of  $10.95$  I have a factor  $9.87$ , instead of  $50.20$  I have a factor  $39.48$ , that would mean in this model the errors in the first 2, calculation of first 2 natural frequencies is about  $11\%$  and  $27\%$ . How do we improve upon these approximations? We have to introduce more nodes, and increase the size of the problem, so we will see such issues later. Now so far what we have done is we have studied an actually vibrating bar and a flexural vibrating being a single span, a single element behavior has been studied.



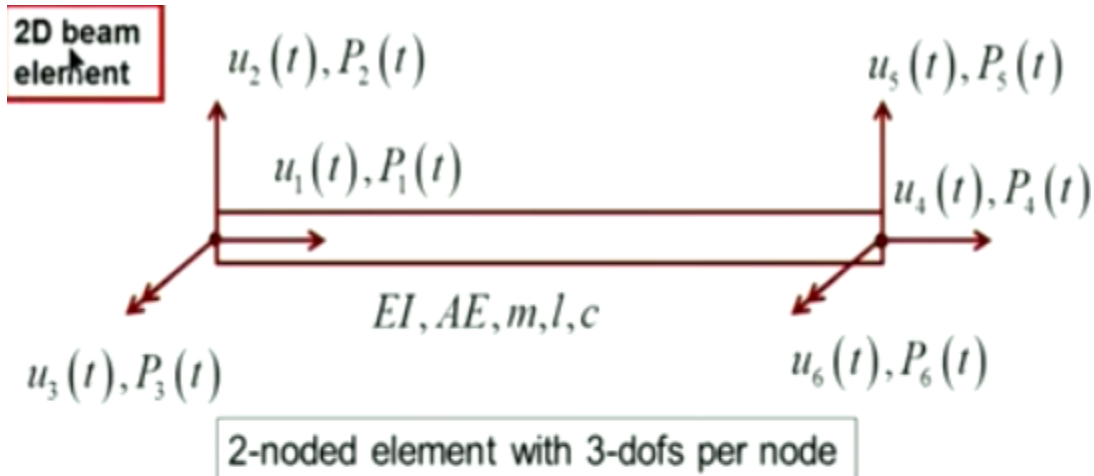
2D beam elements

**Energy:** contributions from both bending and axial oscillations.

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Now if you consider a planar structure like this carrying  $P(t)$  if you carefully you know see each member of this structure, not only deforms actually but also it flexes simultaneously that means the bending and axial vibrations coexist in each of these elements or in other words this if you compute strain energy or kinetic energy in the, each of these elements there will be contribution from actual deformation as well as contribution from flexion. So how do we treat this problem, so we will now address that issue, this leads us to the discussion of what is known as 2D beam



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element, so 2D beam element with you know is shown here again we assume that there are two nodes, now the nodal degrees of freedom for actually vibrating rod it was only the axial deformation, now for only bending it was the translation  $U_2$  and the rotation  $U_3$ . Now since both axial deformation and bending coexist at every node I should now consider  $U_1$ ,  $U_2$  and  $U_3$  one axial translation, and one transverse displacement, and one rotation, so I have  $U_1$ ,  $U_2$ ,  $U_3$  at one node, and similarly  $U_4$ ,  $U_5$ ,  $U_6$  at the other node, so therefore this beam element has 2, it is 2 noded element with 3 degrees of freedom per node, it is a 6 degree of freedom element. What are the parameters of this model? The flexural rigidity  $EI$  axial rigidity  $AE$ ,  $M$  is mass per unit length,  $L$  is length, and  $C$  is the damping parameter, we again assume that flexure is according to Euler Bernoulli beam theory, so the field variables in this case are  $U(x,t)$  which is axial deformation, and  $V(x,t)$  which is the transverse displacement, so these two are the field variables, these are the nodal values of this field variable, for example  $U_1$  and  $U_4$  are the values of  $U(x,t)$  at  $X = 0$ , and at  $X = L$ , similarly  $U_2$ ,  $U_3$ ,  $U_5$ ,  $U_6$  are nodal values associated with  $V(x,t)$  and  $\text{Dou } V / \text{Dou } X$  at  $X = 0$ , and at  $X = L$ .

$$V(t) = \frac{1}{2} \int_0^L AE \left( \frac{\partial u}{\partial x} \right)^2 dx + \frac{1}{2} \int_0^L EI \left( \frac{\partial^2 u}{\partial x^2} \right)^2 dx$$

$$T(t) = \frac{1}{2} \int_0^L m (\dot{u}^2 + \dot{v}^2) dx$$

$$u(x,t) = u_1(t) \phi_1(x) + u_4(t) \phi_4(x)$$

$$v(x,t) = u_2(t) \phi_2(x) + u_3(t) \phi_3(x) + u_5(t) \phi_5(x) + u_6(t) \phi_6(x)$$

$$\phi_1(x) = 1 - \frac{x}{l}; \phi_4(x) = \frac{x}{l}$$

$$\phi_2(x) = 1 - 3 \frac{x^2}{l^2} + 2 \frac{x^3}{l^3}; \phi_3(x) = x - 2 \frac{x^2}{l} + \frac{x^3}{l^2};$$

$$\phi_5(x) = 3 \frac{x^2}{l^2} - 2 \frac{x^3}{l^3}; \phi_6(x) = -\frac{x^2}{l} + \frac{x^3}{l^2}$$

Now we need to know again interpolate the value of these field variables V and U in terms of the nodal values, how do we do that? We will see that, now the expression for kinetic energy and potential energy if you now consider the strain energy has now contribution from axial deformation and contribution from flexure, so we have to add them, similarly the kinetic energy has contribution from axial deformation and transverse deformation, so these are the expressions for the strain energy and the kinetic energy. Now the axial displacement field is again represented like this U1 and U4 are the nodal degrees of freedom, and Phi 1 and this should be Phi 4 are the interpolation functions, and Phi 2, Phi 3, Phi 5, Phi 6 are the interpolation functions associated with bending, so Phi 1 is 1 - X / L and Phi 4(x) is X / L these are the linear interpolation functions associated with axial deformation, and these are the cubic polynomials associated with bending.



$$L(t) = V(t) - T(t) = L(u_1, u_2, \dots, u_6)$$

$$\frac{d}{dt} \left( \frac{\partial L}{\partial \dot{q}_i} \right) - \frac{\partial L}{\partial q_i} = 0; i = 1, 2, \dots, 6$$

$$\Rightarrow M\ddot{u} + Ku = 0$$

$$M = \frac{mL}{420} \begin{bmatrix} 140 & 0 & 0 & 70 & 0 & 0 \\ 0 & 156 & 22l & 0 & 54 & -13l \\ 0 & 22l & 4l^2 & 0 & 13l & -3l^2 \\ 70 & 0 & 0 & 140 & 0 & 0 \\ 0 & 54 & 13l & 0 & 156 & -22l \\ 0 & -13l & -3l^2 & 0 & -22l & 4l^2 \end{bmatrix}$$

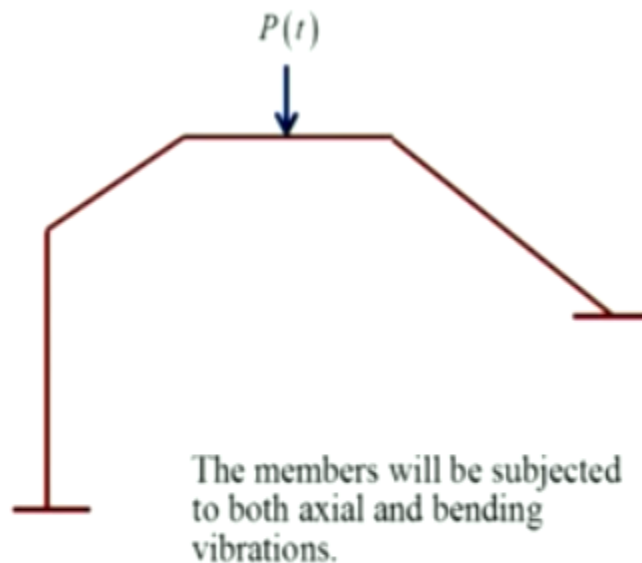


So now I can form the Lagrangian which is  $V - T$ , now this will be function of 6 generalized coordinates  $U_1, U_2$  up to  $U_6$ , so this is 6 degree of freedom model, so I can right now the Euler-Lagrange's equation and for the 6 variables  $i = 1$  to 6, and if I do that again I get a matrix form of equation of motion for the element, and this  $M$  for this case turns out to be this 6 by 6 matrix. And similarly the stiffness matrix is in terms of this 6 by 6 matrix, it is again symmetric and you can see that the contribution from axial deformation is contained here and if you partition and remove these rows and columns what remains is the consistent mass matrix for flexure, similarly here if you look at this this 2 by 2 matrix it is the stiffness matrix

$$K = \frac{EI}{l^3} \begin{bmatrix} \left(\frac{l}{r}\right)^2 & 0 & 0 & -\left(\frac{l}{r}\right)^2 & 0 & 0 \\ 0 & 12 & 6l & 0 & -12 & 6l \\ 0 & 6l & 4l^2 & 0 & -6l & 2l^2 \\ -\left(\frac{l}{r}\right)^2 & 0 & 0 & \left(\frac{l}{r}\right)^2 & 0 & 0 \\ 0 & -12 & -6l & 0 & 12 & -6l \\ 0 & 6l & 2l^2 & 0 & -6l & 4l^2 \end{bmatrix}; r = \sqrt{\frac{I}{A}}$$



corresponding to the axial deformation, this R is the radius of gyration, this is square root I/A and that is how we have written this and the remaining part suppose if you delete these rows and these columns, these rows and these columns we get the 4 by 4 matrix which is a stiffness matrix of the Euler-Bernoulli beam that we have derived.



2D beam elements

**Energy:** contributions from both bending and axial oscillations.

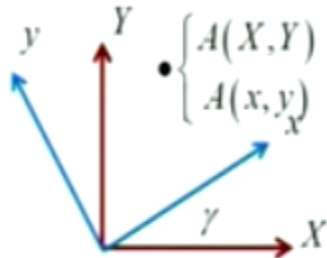
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Now so if we now return to this example one complicating issue namely that in each of these elements we have coexistence of bending and axial deformation we are able to take care of, so now if you want to discretize this beam and study this you can use for each of the element that you choose a representation of this kind. Now there is another problem here, there are few more problems that we need to address one by one, the next problem I would like to consider is now if you look at these members associated with each of these members I can think of a local coordinate system and for the entire structure I can have a global coordinate system, for example this could be the global coordinate system, whereas for this element the local coordinate system will be this.

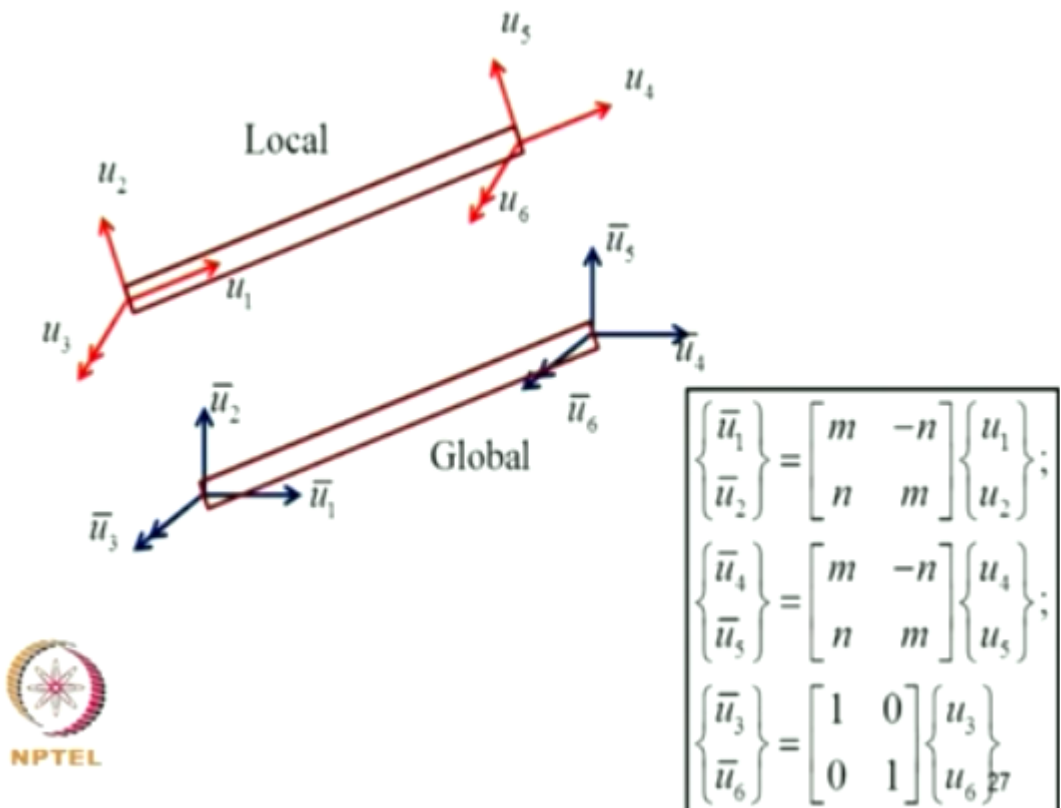
Now how do we now take care of the fact that the local coordinate system for this element does not agree with the global coordinate system, at the end of the day we should be able to describe the behavior of the structure in a common coordinate system.

## Transformation from local to global coordinate system



$$\begin{aligned}
 x &= X \cos \gamma + Y \sin \gamma \\
 y &= -X \sin \gamma + Y \cos \gamma \\
 \underbrace{\begin{Bmatrix} x \\ y \end{Bmatrix}}_{\text{Local}} &= \begin{bmatrix} \cos \gamma & \sin \gamma \\ -\sin \gamma & \cos \gamma \end{bmatrix} \underbrace{\begin{Bmatrix} X \\ Y \end{Bmatrix}}_{\text{Global}} \\
 m &= \cos \gamma; n = \sin \gamma \\
 \underbrace{\begin{Bmatrix} x \\ y \end{Bmatrix}}_{\text{Local}} &= \begin{bmatrix} m & n \\ -n & m \end{bmatrix} \underbrace{\begin{Bmatrix} X \\ Y \end{Bmatrix}}_{\text{Global}} = C \underbrace{\begin{Bmatrix} X \\ Y \end{Bmatrix}}_{\text{Global}} \\
 C &= \begin{bmatrix} m & n \\ -n & m \end{bmatrix} \quad \text{Note: } C^t = C^{-1}
 \end{aligned}$$

So now let us address that issue so that takes us to the question of transformation from local to global coordinate system, so this is you know we have to invoke the standard results from transformations of coordinates as applied to vectors, so suppose you have a coordinate system  $X$  and  $Y$  call this as global coordinate system and a local coordinate system lowercase  $x$  and  $y$  and this orientation let's say  $\gamma$ , then a point here viewed from this capital  $X, Y$  coordinate system will have coordinates capital  $X$  and  $Y$ , if you view from the local, this local coordinate system it has coordinates  $X$  and  $Y$ . Now we can ask the question how this capital  $X$  and  $Y$ , and lowercase  $x$  and  $y$  are related, so a simple consideration of the, you know geometry here leads us to the relation lower  $x$  is  $x \cos \gamma + y \sin \gamma$ , and  $y$  is  $-x \sin \gamma + y \cos \gamma$ , so in the matrix form I can write this as  $XY$  in local coordinate system is related to  $XY$  in global coordinate system through this transformation matrix, and now if we use notation  $M$  and  $N$  for  $\cos \gamma$  and  $\sin \gamma$  I get this coordinate transformation matrix, suppose if I call it as  $C$ , the  $C$  is  $MN - N$  and  $M$ , as is well known this  $C$  matrix is orthogonal that means  $C$  transpose is  $C$  inverse.



Now we will basically put to use this result to achieve what we want, so now let us consider a beam element which is like this initially we will view the element in the local coordinate system, so  $U_1, U_2, U_3, U_4, U_5, U_6$  are the degrees of freedom with respect to the local coordinate system. In the global coordinate system they are  $U_1\text{ bar}, U_2\text{ bar}, U_3\text{ bar}, U_4\text{ bar}, U_5\text{ bar}, U_6\text{ bar}$ , so now if you take the location of this node the coordinates, in local coordinate system is  $U_1, U_2$ , in global coordinate system it is  $U_1\text{ bar}, U_2\text{ bar}$ , so now you go back and use this result for this I get now I am going from local to the global so this matrix is replaced by transpose, so this is the position coordinate of this node, this is how the transformation works. Similarly if you come to this node the position coordinate is  $U_4, U_5$  and it goes to  $U_4\text{ bar}, U_5\text{ bar}$  through this relation.

$$\begin{Bmatrix} \bar{u}_1 \\ \bar{u}_2 \\ \bar{u}_3 \\ \bar{u}_4 \\ \bar{u}_5 \\ \bar{u}_6 \end{Bmatrix} = \begin{bmatrix} m & -n & 0 & 0 & 0 & 0 \\ n & m & 0 & 0 & 0 & 0 \\ 0 & 0 & 1 & 0 & 0 & 0 \\ 0 & 0 & 0 & m & -n & 0 \\ 0 & 0 & 0 & n & m & 0 \\ 0 & 0 & 0 & 0 & 0 & 1 \end{bmatrix} \begin{Bmatrix} u_1 \\ u_2 \\ u_3 \\ u_4 \\ u_5 \\ u_6 \end{Bmatrix}$$

$$\bar{u} = T_0^t u; T_0^t = \begin{bmatrix} L & 0 \\ 0 & L \end{bmatrix}; L = \begin{bmatrix} m & -n & 0 \\ n & m & 0 \\ 0 & 0 & 1 \end{bmatrix}$$


$$\Rightarrow u = T_0 \bar{u}$$



Now the rotation about Z axis remains the same because Z axis has not undergone any transformation, so U3 bar remains as U3, U6 bar remains as U6, therefore this transformation matrix is an identity matrix. Now we can collect all these results and put in a single form, the 6 degrees of freedom namely U1, U2, U3, U4, U5, U6 in local coordinate system is related to the corresponding values in global coordinate system through this transformation matrix, so this is 6 cross 1, this is also 6 cross 1, so for this transformation matrix is 6 cross 6, so we can write this in a compact form, I write this as U bar as T node transpose U, T node t transpose itself is written in this form, we notice that this matrix this is L, this is L, that repeats and this 0 is the 3 by 3 zero matrix this, so this is L, so i can write for U as T node U bar, this is local, this is global.

$$\begin{aligned}
T(t) &= \frac{1}{2} \dot{u}'(t) M \dot{u}(t) \\
&= \frac{1}{2} \dot{\bar{u}}' T_0' M T_0 \dot{\bar{u}} \\
&= \frac{1}{2} \dot{\bar{u}}' \bar{M} \dot{\bar{u}}; \\
\bar{M} &= T_0' M T_0
\end{aligned}$$

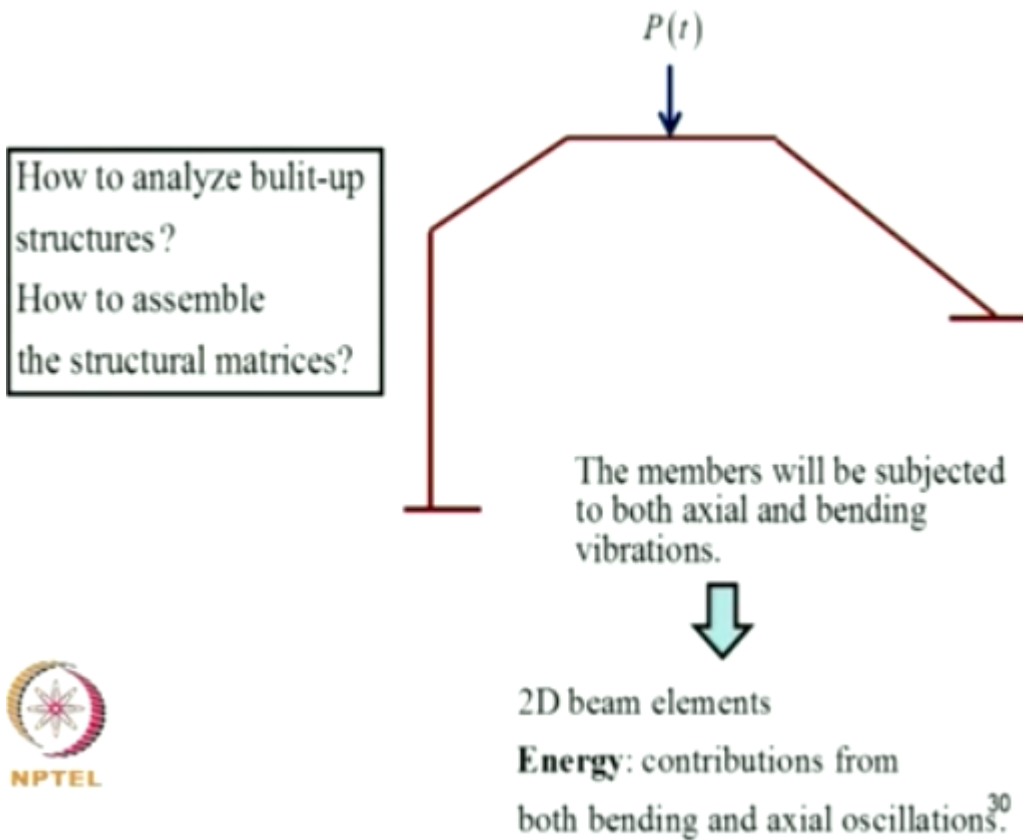
$$\begin{aligned}
V(t) &= \frac{1}{2} u'(t) K u(t) \\
&= \frac{1}{2} \bar{u}' T_0' K T_0 \bar{u} \\
&= \frac{1}{2} \bar{u}' \bar{K} \bar{u} \\
\bar{K} &= T_0' K T_0
\end{aligned}$$


 $\bar{K}$  = Element stiffness matrix in the transformed coordinate system  
 $\bar{M}$  = Element mass matrix in the transformed coordinate system

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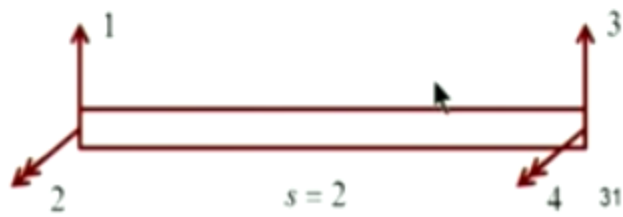
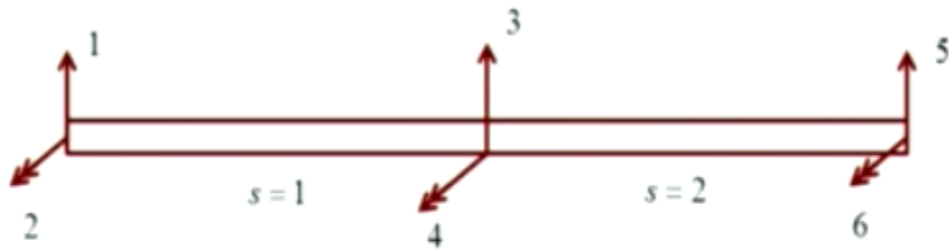
Now what we need to do is use this transformation in our expression for the kinetic energy and potential energy, so in local coordinate system the kinetic energy is given by this, the coordinate transformation doesn't change the energy, energy is a scalar but the representation for mass matrix changes so that is what we are looking at. So I know that  $U$  is  $T$  node  $U$  therefore I can write  $U$  dot transpose as  $\bar{U}$  dot transpose  $T$  node transpose  $M T$  node this, so this quantity which is  $T$  node transpose  $M T$ , I call it as  $\bar{M}$ , so this is a mass matrix in the global coordinate system for the element that is given by  $T$  node transpose empty. And a kinetic energy now I have in terms of the velocity is expressed in the global coordinate system, similar exercise we can now repeat on expression for strain energy so I have this as half  $U$  transpose  $K U$ , for  $U$  transpose now I will write, for  $U$  I will write at  $T$  node  $\bar{U}$  and  $U$  transpose will be this, and again I introduce  $\bar{K}$  which is  $T$  node transpose  $K T$  node, so this is the element stiffness matrix in the global coordinate system, and this is the expression for strain energy in the global coordinate system. So I have now therefore  $\bar{K}$  and  $\bar{M}$  which are the element structural matrices, element stiffness matrix in the transform coordinate system,  $\bar{M}$  is element mass matrix in a transform coordinate system. So we have taken care of this complication that the global and local coordinate systems could differ for different elements.



Now we will now move to the next complication that we need to deal with, now I will list what are the complication we need to deal with at this stage, we need to now take care of the issue that this structure is made up of different element which are oriented in a different direction, so if I now partition this into elements I will probably take this as one element this is at least as one element, this as one element, this is one element, we could have more elements within this it is a different issue, so the question is how to analyze a structure which is built up of different you know elements that is the first question. The next question we have to deal with is how to deal with external forces which I have not yet tackled, and how to impose the boundary condition systematically, and if there are any support motions as in case of earthquake engineering problems, how do they enter our formulation. And finally I had not talked about damping, how do we handle damping in this system, so we have to answer all these questions so I am taking up one by one. So now the next question I wish to answer is how to analyze built up structures or in other words if I now assume that this is one element, this is another element, this is one element, and so on and so forth, for each of these elements I can derive a 6 by 6 mass and stiffness matrix, but how do I now get the mass and stiffness matrix for the built-up structure? By assembling these element level structural matrices, so how do we do that?



$$p = 2; N = 6$$



To illustrate that we can consider a simple problem, suppose there is one beam which has been divided into 2 elements, say  $S=1$  and  $S=2$ .

Now let this be Euler Bernoulli beam there is no axial deformation in this illustration, so how many degrees of freedom this structure has? It has 6 degrees of freedom, this is 1, 2, 3, 4, 5, 6. Now each element has 4 degrees of freedom, so now if I take this element separately I will number these degrees of freedom as 1, 2, 3, 4, if I take the second element separately this numbering as per our scheme will be again 1, 2, 3, 4, but we should somehow impose the now the condition that what is 1 here must be 3 here, what is 2 here must be 4 here, that would ensure compatibility of displacements. How do we do that? So the basic idea is the displacement at shared nodes must match, so let  $P$  be the number of elements which was 2 in the illustration that I showed just now, let  $N$  be the total degrees of freedom which was 6 in the illustration.

## Analysis of built-up structures

- Displacements at shared nodes must match.

Let

$p$  = number of elements

$N$  = total dof-s

$\bar{U}_j; j = 1, 2, \dots, N$

Nodal displacement vector ( $N \times 1$ ) in global coordinate system for the structure.

$$\{\bar{U}\} = \{\bar{U}_j\}_{j=1}^N$$

$\{\bar{u}\}_s =$  Nodal displacement vector ( $N_s \times 1$ ) in global coordinate system for the  $s$ -th element

We write  $\{\bar{u}\}_s$  in terms of  $\{\bar{U}\}$  as



$$\{\bar{u}\}_s = [A]_s \{\bar{U}\}$$

$N_s \times N$     $N \times 1$

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Now I will denote the degrees of freedom as  $U_j$  where  $j$  runs from 1 to  $N$ , now the nodal displacement vector is a  $N \times 1$  in global coordinate system for this structure, now this is  $U_j$  where  $j = 1$  to  $N$ , but for each of the elements as I have shown here, I have this lowercase  $u$  where  $s$  which are the nodal displacement vector in global coordinate system for the  $s$ -th element,  $s = 1$ , for example I had 1, 2, 3, 4 and  $s = 2$  again I have 1, 2, 3, 4. Now we would like to write the nodal degrees of freedom at the element level in terms of the nodal degrees of freedom at the structure level, so what we do is we introduce a transformation  $U$  where  $s$  is written as  $A_s U$  into  $u$ , now  $u$  where  $s$  is some  $N_s \times 1$ , in this case  $N_s$  is for 4, each beam element has 4 degrees of freedom it is  $4 \times 1$ , whereas this  $U$  where  $s$  is  $6 \times 1$ , therefore this  $A$  matrix should be  $N_s \times N$ , so number of rows in  $A$  matrix will be equal to the number of degrees of freedom at the element level and number of columns in  $A$  matrix will be equal to the total degrees of freedom at the structure level.

**Example**

$$\begin{Bmatrix} \bar{u}_1 \\ \bar{u}_2 \\ \bar{u}_3 \\ \bar{u}_4 \end{Bmatrix}_{s=1} = \begin{bmatrix} 1 & 0 & 0 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 & 0 & 0 \\ 0 & 0 & 1 & 0 & 0 & 0 \\ 0 & 0 & 0 & 1 & 0 & 0 \end{bmatrix} \begin{Bmatrix} \bar{U}_1 \\ \bar{U}_2 \\ \bar{U}_3 \\ \bar{U}_4 \\ \bar{U}_5 \\ \bar{U}_6 \end{Bmatrix}$$

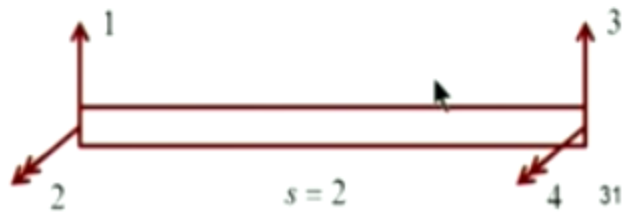
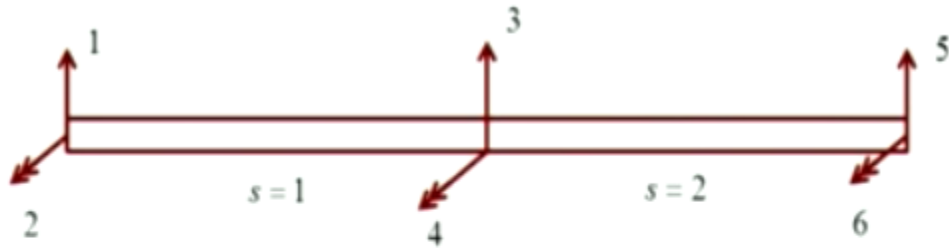
$$\begin{Bmatrix} \bar{u}_1 \\ \bar{u}_2 \\ \bar{u}_3 \\ \bar{u}_4 \end{Bmatrix}_{s=2} = \begin{bmatrix} 0 & 0 & 1 & 0 & 0 & 0 \\ 0 & 0 & 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 & 0 & 1 \end{bmatrix} \begin{Bmatrix} \bar{U}_1 \\ \bar{U}_2 \\ \bar{U}_3 \\ \bar{U}_4 \\ \bar{U}_5 \\ \bar{U}_6 \end{Bmatrix}$$



Number of rows in  $A_s$   
 = Number of rows in  $\bar{u}_s$   
 Number of columns in  $A_s$   
 = Number of rows in  $\bar{U}$

So now for this simple example that we used for this example if you know consider U1, U2, U3, U4, for the first element what is 1, 2, 3, 4 is same as 1, 2, 3, 4, for the global structure, by that I mean this 1, 2, 3, 4 and this 1, 2, 3, 4 and this 1, 2, 3, 4 are the same there is no difference so the A matrix in this case will have an identity matrix here and then 0s. For the second

$$p = 2; N = 6$$



element if you go back what is 1 is actually 3 for the global structure, and what is 2 is 4, and similarly what are 3 and 4 are respectively 5 and 6, so the A matrix now will become as shown

### Example

$$\begin{Bmatrix} \bar{u}_1 \\ \bar{u}_2 \\ \bar{u}_3 \\ \bar{u}_4 \end{Bmatrix}_{s=1} = \begin{bmatrix} 1 & 0 & 0 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 & 0 & 0 \\ 0 & 0 & 1 & 0 & 0 & 0 \\ 0 & 0 & 0 & 1 & 0 & 0 \end{bmatrix} \begin{Bmatrix} \bar{U}_1 \\ \bar{U}_2 \\ \bar{U}_3 \\ \bar{U}_4 \\ \bar{U}_5 \\ \bar{U}_6 \end{Bmatrix}$$

$$\begin{Bmatrix} \bar{u}_1 \\ \bar{u}_2 \\ \bar{u}_3 \\ \bar{u}_4 \end{Bmatrix}_{s=2} = \begin{bmatrix} 0 & 0 & 1 & 0 & 0 & 0 \\ 0 & 0 & 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 & 0 & 1 \end{bmatrix} \begin{Bmatrix} \bar{U}_1 \\ \bar{U}_2 \\ \bar{U}_3 \\ \bar{U}_4 \\ \bar{U}_5 \\ \bar{U}_6 \end{Bmatrix}$$



Number of rows in $A_s$ = Number of rows in $\bar{u}_s$ Number of columns in $A_s$ = Number of rows in $\bar{U}$
---

here, so  $\bar{U}_1$  is the third element  $\bar{U}_3$ ,  $\bar{U}_2$  is the fourth element which is  $\bar{U}_4$  and so on and so forth, so with this representation now we have the expression for kinetic energy and strain energy in terms of the local node numbering scheme. Now we will go to the built-up structure, now the kinetic energy for example in the built-up structure consists of contribution to the total kinetic energy from each of the elements, so if there are  $P$  elements this is some of  $S = 1$  to  $P$ ,  $TS(t)$ . Now  $TS(t)$  itself is given in terms of the local degrees of freedom the element level degrees of freedom as shown here this is what we have derived now, these are in the global coordinate system.



$$\begin{aligned}
 T(t) &= \sum_{s=1}^p T_s(t) \\
 &= \sum_{s=1}^p \frac{1}{2} \{\dot{\bar{u}}\}'_s [\bar{M}]_s \{\dot{\bar{u}}\}_s \\
 &= \sum_{s=1}^p \frac{1}{2} \{\dot{\bar{U}}\}' [A]_s' [\bar{M}]_s [A]_s \{\dot{\bar{U}}\} \\
 &= \frac{1}{2} \{\dot{\bar{U}}\}' \left[ \sum_{s=1}^p [A]_s' [\bar{M}]_s [A]_s \right] \{\dot{\bar{U}}\} \\
 &= \frac{1}{2} \{\dot{\bar{U}}\}' \bar{M} \{\dot{\bar{U}}\} \\
 \bar{M} &= \sum_{s=1}^p [A]_s' [\bar{M}]_s [A]_s
 \end{aligned}$$

Now I will introduce the transformation U bar is AS capital U bar, so if I substitute this I get this expression, now this summation is over the elements and you can see that the quantities inside the braces here and here are independent of S so they can be pulled out, and the summation you will be only on these terms, this summed up quantity which is written as this is the structure mass matrix in the global coordinate system, okay, so this receives contributions from each of the element, that is the element level mass matrices combined together according to this rule to get, to lead to the global level mass matrix, similarly the strain energy is again,



$$\begin{aligned}
 V(t) &= \sum_{s=1}^p V_s(t) \\
 &= \sum_{s=1}^p \frac{1}{2} \{\bar{u}\}_s' [\bar{K}]_s \{\bar{u}\}_s \\
 &= \sum_{s=1}^p \frac{1}{2} \{\bar{U}\}' [A]_s' [\bar{K}]_s [A]_s \{\bar{U}\} \\
 &= \frac{1}{2} \{\bar{U}\}' \left[ \sum_{s=1}^p [A]_s' [\bar{K}]_s [A]_s \right] \{\bar{U}\} \\
 &= \frac{1}{2} \{\bar{U}\}' \bar{K} \{\bar{U}\} \\
 \bar{K} &= \sum_{s=1}^p [A]_s' [\bar{K}]_s [A]_s
 \end{aligned}$$

the total strain energy is sum of contribution from strain energy within each elements, and this is again following the same logic it is given in terms of the element level degrees of freedom, and in terms of global degrees of freedom through this transformation like this, and again we can pull out this quantities in the brace outside the summation and I get this, and this summation is actually the assembling procedure leading to the global structure stiffness matrix, so this is summed over S = 1 to P this, so we have now therefore the expressions ready for global mass matrix and global stiffness matrix.

$$\bar{M} = \sum_{s=1}^p [A]_s^T [\bar{M}]_s [A]_s = \text{Global mass matrix}$$

$$\bar{K} = \sum_{s=1}^p [A]_s^T [\bar{K}]_s [A]_s = \text{Global stiffness matrix}$$

Global equation of motion

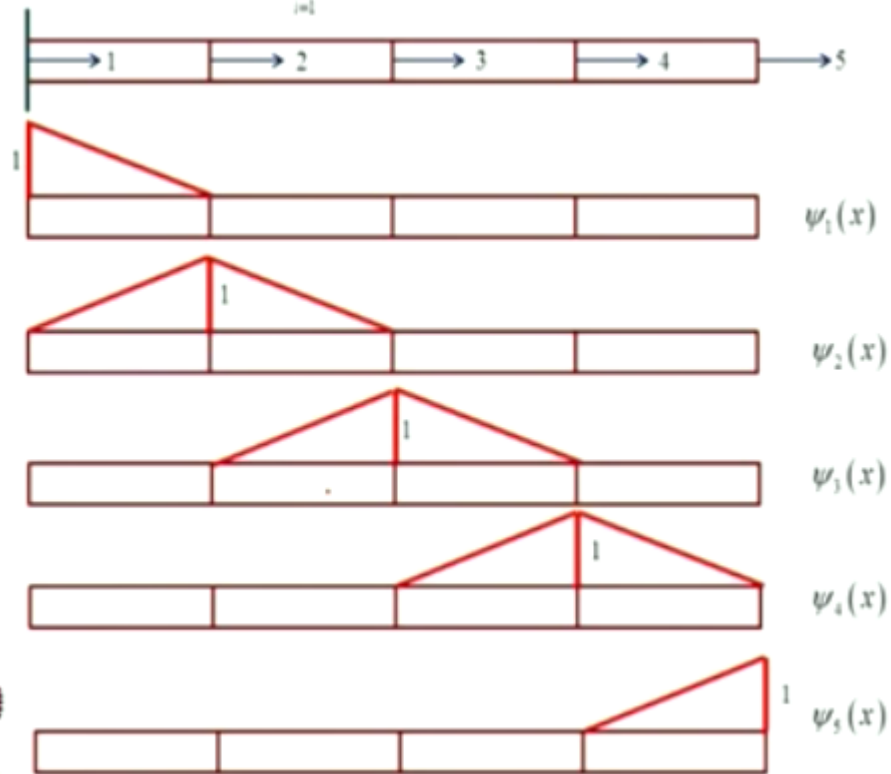
$$\bar{M}\ddot{U} + \bar{K}U = P(t)$$



Now we can write notionally the equation of motion that is a global equation of motion the right-hand side we will call it as a force vector which we are not yet you know described what it should be.



$$u(x,t) = \sum_{i=1}^n a_i(t)\psi_i(x)$$



So now let us come to that description but before that we can you know illustrate what exactly we have done by considering this simple example of a axially vibrating bar which has been discretized into a set of 4 elements 1, 2, 3, 4 and the finite element discretization implies the use of these global shape functions, this is not how we actually compute though but what we do is we carry out calculation at the element level and then sum up, so what we need to do here is we need to find for example kinetic energy in terms of the displacement field as shown here,

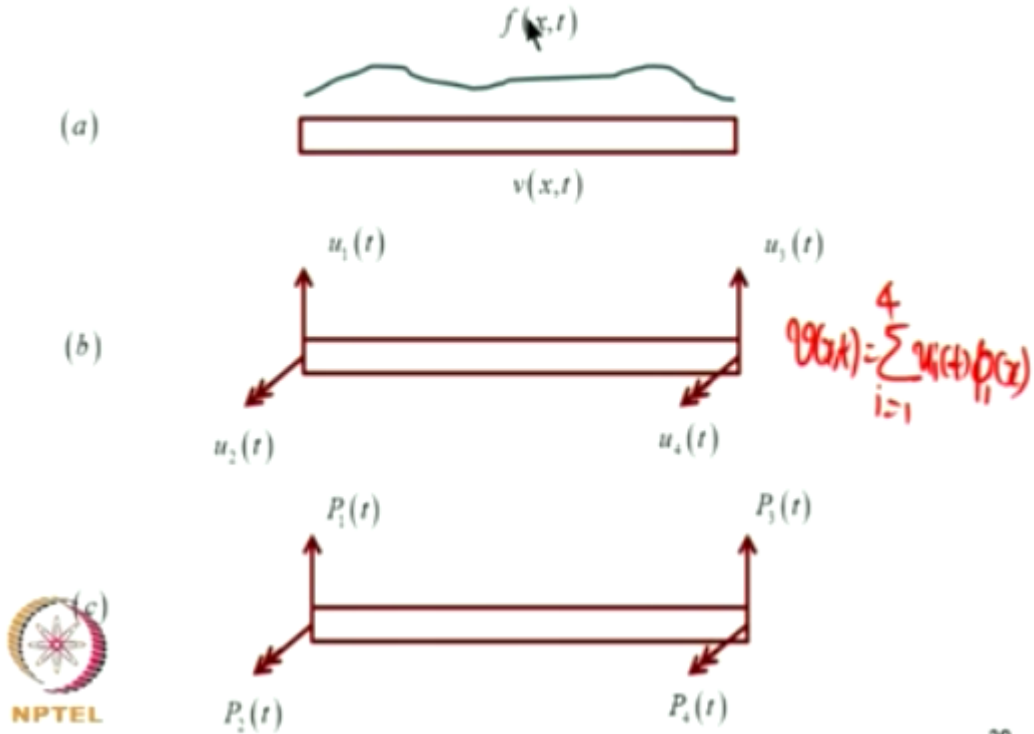


$$\begin{aligned}T(t) &= \frac{1}{2} \int_0^l m \dot{u}^2(x, t) dx \\&= \sum_{e=1}^4 \frac{1}{2} \int_{x_{e1}}^{x_{e2}} m \dot{u}_e^2(x, t) dx \\&= \sum_{e=1}^4 \frac{1}{2} \int_{x_{e1}}^{x_{e2}} m [\dot{u}_{e1}(t) \phi_{e1}(x) + \dot{u}_{e2}(t) \phi_{e2}(x)]^2 dx \\&= \sum_{e=1}^4 T_e(t) \\V(t) &= \sum_{e=1}^4 V_e(t) \\L &= \sum_{e=1}^4 [T_e(t) - V_e(t)]\end{aligned}$$

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and this we can add over the elements so I can find kinetic energy contribution from this element, this will have you know contribution from only these two trial functions, all other trial functions are 0 here similarly for this element there are two instances, for this element there are 2 instances, and this is 2 instances, what we are actually doing is we are computing these 4 situations separately, so that is what we are doing, so this is what I meant when I said total kinetic energy is sum of contribution from each of the elements, so then we can construct a Lagrangian which is again sum, and if we actually then carry out the, if you implement the Euler-Lagrange's equation on this representation we will get basically what we got here, so we are able to achieve these calculations through this simple you know in the language of matrix algebra through these arguments.

## Generalized forces



Now we will come to the question on forces, so how do we determine the nodal force? Suppose there is a beam element which is carrying a transverse load like this, so in your finite element approximation we have approximated the field variable  $V(x,t)$  in terms of  $U_1, U_2, U_3, U_4$  and the 4 interpolation functions. Now we would like to replace this distributed load in terms of these equivalent nodal forces, so we have therefore  $V(x,t)$  as  $\sum_{i=1}^4 U_i(t) \phi_i(x)$ , this is how we represent the field variable, and we would like to now replace this  $F(x,t)$  through these equivalent nodal forces, so how I can replace therefore this situation by a combination of these two situations, that is the question, so what we do is we use the principle of virtual work, we define  $\Delta V(x,t)$  as a virtual displacement, we can write this in terms of virtual displacements at the nodes and the interpolation function, because this is a representation that we are talking about.

Let  $\delta v(x, t)$  = a virtual displacement. We can write

$$\delta v(x, t) = \sum_{i=1}^4 \delta u_i(t) \phi_i(x)$$

$$\Rightarrow \sum_{i=1}^4 P_i(t) \delta u_i(t) = \int_0^l f(x, t) \delta v(x, t) dx$$

$$= \int_0^l f(x, t) \left\{ \sum_{i=1}^4 \delta u_i(t) \phi_i(x) \right\} dx$$

$$\Rightarrow \sum_{i=1}^4 \left\{ P_i(t) - \int_0^l f(x, t) \phi_i(x) dx \right\} \delta u_i(t) = 0$$

Since  $\delta u_i(t), i = 1, 2, 3, 4$  are arbitrary

$$\Rightarrow P_i(t) = \int_0^l f(x, t) \phi_i(x) dx, i = 1, 2, 3, 4$$



Now the work done by the nodal forces on the virtual displacements must be equal to the work done by these forces on these virtual displacement, so if you now write for  $\Delta U(x, t)$  this we can take the terms to the other side and we get this expression. Now since the  $\Delta U_i(t)$  are all arbitrary the only way this can be 0 for all T is that the term inside the brace must be independently equal to 0 and that leads us to the definition of the equivalent nodal forces. For example if the beam is carrying a UDL, but as a function of time I can derive the equivalent

Example:  $f(x, t) = q(t)$

$$\Rightarrow P_i(t) = \int_0^l q \phi_i(x) dx, i = 1, 2, 3, 4$$

$$P_1(t) = \frac{q(t)l}{2}; P_2(t) = \frac{q(t)l}{12}; P_3(t) = \frac{q(t)l}{2}; P_4(t) = -\frac{q(t)l}{12}$$

Example:  $f(x, t) = P(t)\delta(x - a)$

$$P_i(t) = \int_0^l P(t)\delta(x - a)\phi_i(x) dx, i = 1, 2, 3, 4$$

$$\xi = \frac{x}{l}$$

$$P_1(t) = (1 - 3\xi^2 + 2\xi^3)P(t); P_2(t) = (1 - 2\xi^2 + 3\xi^3)P(t)$$

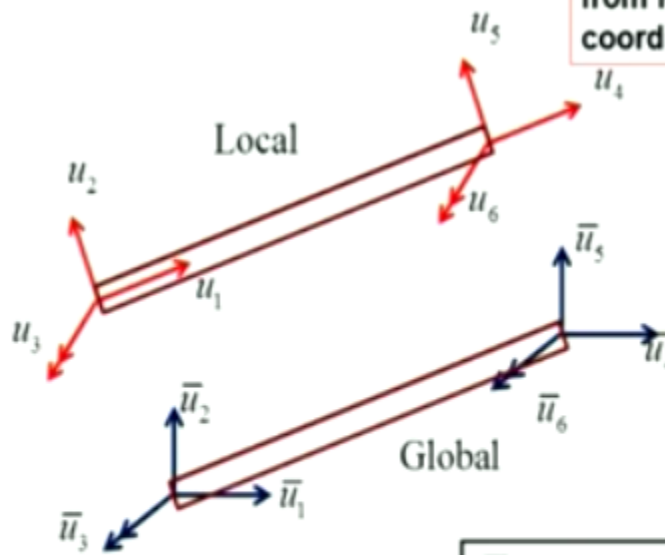
$$P_3(t) = (3\xi^2 - 2\xi^3)P(t); P_4(t) = (-\xi^2 + \xi^3)P(t)$$



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nodal forces as P1, P2, P3, P4 as shown, similarly if we have a impulsive force acting at  $X = A I$  can substitute and we can get the expression for equivalent nodal forces.

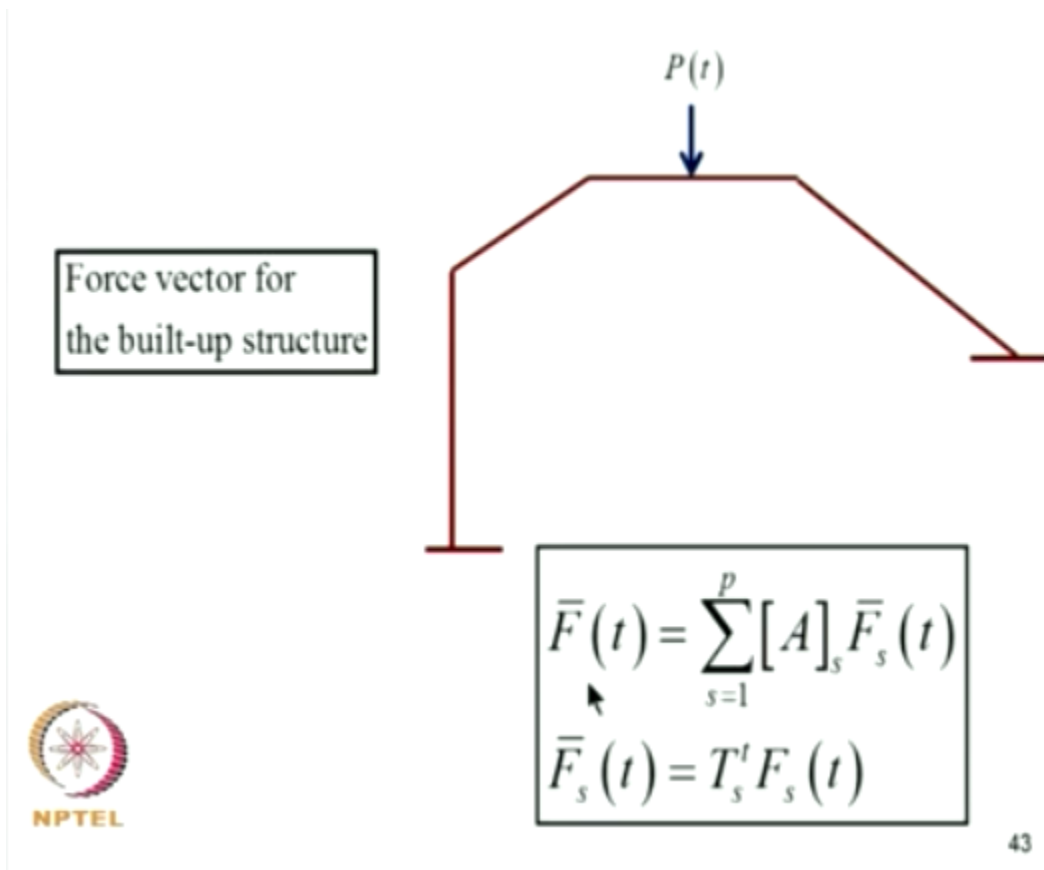
Transformation of force vector  
from local to global  
coordinate system



$$\bar{F}_s(t) = T_s^t F_s(t)$$



Now we are getting now equivalent nodal forces in the local coordinate system of the beam, how do you get in global coordinate system? It's simply the transformation, because force is a vector just like a displacement the same rule of transformation as was used for displacements also work here and this is what I get. And how do we assemble?



Suppose a force vector for a built-up structure, how do we assemble that, again we follow the same logic as we use for assembling structural matrices and the global force vector in terms of element level force vector is again through a summation using these A matrices this is as shown. So first we get, first we formulate the force in the local coordinate system and bring it to the global coordinate system and then use these A matrices and sum or  $S = 1$  to  $P$  to get the equivalent nodal forces in the global coordinate system for the built-up structure.

So now at this stage what we have obtained is the global structure equation of motion, the quantity that you are seeing in red here is a damping force so I am now arbitrarily introducing this assuming that the damping mechanism is viscous and we are going to select this damping matrix in a way that facilitates the use of normal modes which are obtained by un-damped free vibration analysis, therefore the C matrix can be introduced at a later stage, so I have arbitrarily introduced this matrix, we will return to this issue in some greater detail later.

$$\Rightarrow \bar{M} = \begin{bmatrix} M_{00} & M_{0I} \\ M_{I0} & M_{II} \end{bmatrix}; \bar{K} = \begin{bmatrix} K_{00} & K_{0I} \\ K_{I0} & K_{II} \end{bmatrix}; \bar{C} = \begin{bmatrix} C_{00} & C_{0I} \\ C_{I0} & C_{II} \end{bmatrix};$$

$$\bar{F}(t) = \begin{Bmatrix} \bar{F}_0(t) \\ \bar{F}_I(t) \end{Bmatrix} = \begin{Bmatrix} \text{Unknown reactions} \\ \text{Applied equivalent nodal forces} \end{Bmatrix}$$

### Global equation of motion

$$\begin{bmatrix} M_{00} & M_{0I} \\ M_{I0} & M_{II} \end{bmatrix} \begin{Bmatrix} 0 \\ \ddot{\bar{U}}_I \end{Bmatrix} + \begin{bmatrix} C_{00} & C_{0I} \\ C_{I0} & C_{II} \end{bmatrix} \begin{Bmatrix} 0 \\ \dot{\bar{U}}_I \end{Bmatrix} + \begin{bmatrix} K_{00} & K_{0I} \\ K_{I0} & K_{II} \end{bmatrix} \begin{Bmatrix} 0 \\ \bar{U}_I \end{Bmatrix} = \begin{Bmatrix} \bar{F}_0(t) \\ \bar{F}_I(t) \end{Bmatrix}$$

$$M_{0I} \ddot{\bar{U}}_I + C_{0I} \dot{\bar{U}}_I + K_{0I} \bar{U}_I = \bar{F}_0(t) \quad (\text{Equations for unknown reactions})$$

$$M_{II} \ddot{\bar{U}}_I + C_{II} \dot{\bar{U}}_I + K_{II} \bar{U}_I = \bar{F}_I(t) \quad (\text{Equations for unknown displacements})$$

$$\text{Number of equations} = q + (N - q) = N$$

Unknowns:  $q$  reactions ( $\bar{F}_0(t)$ ) and  $(N - q)$  displacements ( $\bar{U}_I(t)$ ).

So now the next issue that we have to consider is the boundary condition, so if you go back to this now we have tackled a few issues we have considered the fact that these elements carry both actual and bending energies so we have tackled that, and the structure itself is made up of different elements which are oriented in different directions so that we have tackled, we have a first transform all the entities in a global coordinate system and we have devised a way to assemble them. And similarly we have now tackled the issue of how to handle an external force and convert it into equivalent nodal forces in global coordinate system for the structures, at the structure level, so now the question that remains is how to deal with boundary conditions so now, so what we do is to deal with that we partition the displacement vector into a component let us call it as  $U_G(t)$  and another one let me call it  $U_I$ , this  $U_G(t)$ , the  $G$  we can take it as representing the ground, the ground displacements can be specified okay, it can be specified to be 0 or it can be specified to be a given time history, often we take that to be 0, in situation where the dynamic actions come through support motion as in case of earthquake induced vibrations  $U_G(t)$  will be specified time histories, so in any case they are known, the remaining nodal degrees of freedom are the structure degrees of freedom that we need to determine, these are the unknown nodal displacements.

Now to start with let us consider that the nodal displacements are given to be 0, so the partitioning of the displacement vector will be 0 and  $U_I$ , so if there are  $Q$  components in  $U_G(t)$  this will be  $Q \times 1$ , and this will be  $N - Q \times 1$ , now the partitioning of this displacement vector also induces a partitioning on the structural matrices and the forcing vector, so I will partition the mass matrix, stiffness matrix, damping matrix, and the force vector in this form, the size of  $M_{00}$ ,  $K_{00}$ ,  $C_{00}$  will be  $Q \times Q$  and  $M_{II}$ ,  $K_{II}$ ,  $C_{II}$  will be  $(N - Q) \times (N - Q)$ ,



so the global of equation of motion, the global equation of motion in the partition form can be written this way.

Now you see here these displacement components are specified to be 0 and they will be associated with unknown reactions, and associated with unknown displacements which is  $U_i$  bar there will be equivalent nodal forces which are externally applied, not all of them need to be nonzero it depends on the specific problem on hand, so there are now capital N number of equations and capital N number of unknowns, what are the capital N number of unknowns? Q reactions, and N - Q unknown nodal displacements, so we can write now the equation corresponding to the first row here, so  $M_{00}$  this is 0,  $M_{0i} \ddot{U}_i$  +  $C_{0i} \dot{U}_i$  +  $K_{0i} U_i$  bar is  $F_{node}(t)$ , so this set of equations represent the equations for unknown reactions, so they can be determined only after you find this  $U_i$  bar, how do I find  $U_i$  bar? I write the equation from the second row, so that is this, that is  $M_{10}$  into 0 +  $M_{22} \ddot{U}_i$  bar double dot and so on and so forth the other terms on the right-hand side I have the equivalent nodal forces. So these are the equations for unknown displacements, so the number of equations is capital N, Q reactions N - Q unknown displacements.

**Reduced structure equation of motion**

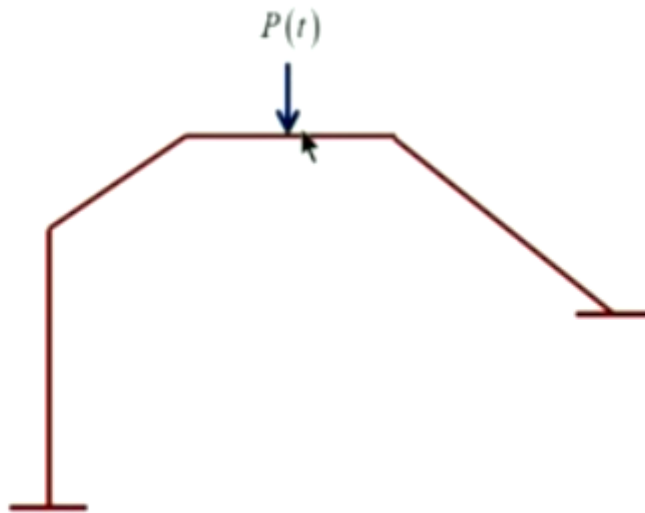
$$M_{ii} \ddot{U}_i + C_{ii} \dot{U}_i + K_{ii} U_i = \bar{F}_i(t)$$

Or simply, the governing equation of motion


$$M\ddot{U} + C\dot{U} + KU = F(t)$$

$$U(0) = U_0; \dot{U}(0) = \dot{U}_0$$

Now what we do is we avoid all this clutter of subscripts and bars and things like that, we write this equation, this set of equation which are the equations for unknown nodal displacements and velocities and accelerations simply as  $M\ddot{U} + C\dot{U} + KU = F$  which specified initial conditions, so these are the governing equilibrium equations for the structure which we need to solve, so this has taken into account the effect of boundary conditions and all the assembling as we described, so these are the unknowns that we have to solve by solving this set of coupled second order ordinary differential equations.



So now let us quickly summarize what we have been doing, we want to analyze the structure like this carrying a dynamic load like this, this structure is made up of different elements  $S = 1$ , I can have this is  $S = 2, 3, 4, 5, 6$ , so what we do for each element we first write the element level equation of motion we know the element level mass matrix, stiffness matrix, and we can

<b>Summary</b>	
Element level EOM in local coordinate system	$M_i \ddot{u}_i + C_i \dot{u}_i + K_i u_i = F_i(t)$
Element level EOM in global coordinate system	$\bar{M}_i \ddot{\bar{U}}_i + \bar{C}_i \dot{\bar{U}}_i + \bar{K}_i \bar{U}_i = \bar{F}_i(t)$ $\bar{U}_i = T_i' u_i; \bar{M}_i = T_i' M_i T_i; \bar{K}_i = T_i' K_i T_i;$ $\bar{C}_i = T_i' C_i T_i$
Global EOM after assembly of structural matrices and before imposing boundary conditions	$\bar{M} \ddot{\bar{U}} + \bar{C} \dot{\bar{U}} + \bar{K} \bar{U} = \bar{F}(t)$ $\bar{M} = \sum_{i=1}^p [A]_i' [\bar{M}]_i [A]_i; \bar{K} = \sum_{i=1}^p [A]_i' [\bar{K}]_i [A]_i$ $\bar{C} = \sum_{i=1}^p [A]_i' [\bar{C}]_i [A]_i; \bar{F}(t) = \sum_{i=1}^p [A]_i \bar{F}_i(t)$
Equations for unknown reactions	$M_{o_i} \ddot{\bar{U}}_i + C_{o_i} \dot{\bar{U}}_i + K_{o_i} \bar{U}_i = \bar{F}_{o_i}(t)$
 Equations for unknown displacements	$M_{ii} \ddot{\bar{U}}_i + C_{ii} \dot{\bar{U}}_i + K_{ii} \bar{U}_i = \bar{F}_i(t)$ $\bar{M} \ddot{\bar{U}} + \bar{C} \dot{\bar{U}} + \bar{K} \bar{U} = \bar{F}(t)$ $\bar{U}(0) = \bar{U}_0; \dot{\bar{U}}(0) = \dot{\bar{U}}_0$

notionally introduce the damping matrix and we can get the force vector also for S-th element, then we go to the global coordinate system, okay, see the local coordinate system and global coordinate system for this element it is same, suppose if I take the global coordinate system to be this, for this element 4 this local and global coordinate system coincide, not for the other elements, so you have to make a transformation, so you then get the equation of motion for elements S = 1 to P in the global coordinate system through these transformation that means for each of the element you have to find out T1, T2, T3, TP, they are the coordinate transformation matrix.

Next for each of the element we have to now define them A matrix which connects the naming of degrees of freedom at the element level to the naming of degrees of freedom at the structure level, so that leads us to the global equation of motion after assembly of structural matrices and before imposing boundary conditions, this is the equation. So this is a set of capital N number of equations, Q of them for reactions and N - Q for unknown displacements.

Then we now get the equation for unknown reactions we partition the displacement vector into Q and N - Q elements and one set of equations is equations for unknown reactions, the other set of equations is for unknown displacement, so this is the equation shown in red here is the governing equation of motion which we need to solve to obtain the, solution to the problem on hand what really happens is at the end, this is a set of ordinary differential equations T is still a continuous variable here, so these equations are known as semi discretized equations, so we have to further develop numerical methods to discretize this time and we will get a discrete map from which we will determine all the nodal displacements, velocities and accelerations,

subsequently we have to now go back to the element level descriptions through the appropriate interpolation functions and the subsequent questions will be on how to find out displacement at points which don't coincide with the nodes, how do you find, these are displacements, how do you find strains and stresses, and how do you, you know compute quantities like principle stresses, principle strains, stress metric, etcetera, etcetera, those questions need to be still addressed.

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