Structural Matrices in MDOF Systems

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Today we will study the properties of structural matrices, that is the operators that relate the vector of system coordinates \mathbf{x} and its time derivatives $\dot{\mathbf{x}}$ and $\ddot{\mathbf{x}}$ to the forces acting on the system nodes, \mathbf{f}_{S} , \mathbf{f}_{D} and \mathbf{f}_{I} , respectively.

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In the end, we will see again the solution of a MDOF problem by superposition, and in general today we will revisit many of the subjects of our previous class, but you know

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We already met the mass and the stiffness matrix. M and K, and tangentially we introduced also the dampig matrix \boldsymbol{C} .

We have seen that these matrices express the linear relation that holds between the vector of system coordinates x and its time derivatives \dot{x} and \ddot{x} to the forces acting on the system nodes, f_S , f_D and f_I , elastic, damping and inertial force vectors.

$$\mathbf{M} \ddot{\mathbf{x}} + \mathbf{C} \dot{\mathbf{x}} + \mathbf{K} \mathbf{x} = \mathbf{p}(t)$$

 $\mathbf{f}_1 + \mathbf{f}_D + \mathbf{f}_S = \mathbf{p}(t)$

Also, we know that **M** and **K** are symmetric and definite positive, and that it is possible to uncouple the equation of motion expressing the system coordinates in terms of the eigenvectors, $\mathbf{x}(t) = \sum q_i \psi_i$, where the q_i are the modal coordinates and the eigenvectors ψ_i are the non-trivial solutions to the equation of free vibrations,

$$\left(\textbf{\textit{K}}-\omega^{2}\textbf{\textit{M}}\right)\psi=\textbf{0}$$

From the homogeneous, undamped problem

$$\mathbf{M}\ddot{\mathbf{x}} + \mathbf{K}\mathbf{x} = \mathbf{0}$$

introducing separation of variables

$$\mathbf{x}(t) = \mathbf{\psi} \left(A \sin \omega t + B \cos \omega t \right)$$

we wrote the homogeneous linear system

$$(\mathbf{K} - \omega^2 \mathbf{M}) \, \mathbf{\psi} = \mathbf{0}$$

whose non-trivial solutions ψ_i for ω_i^2 such that $\|\mathbf{K} - \omega_i^2 \mathbf{M}\| = 0$ are the eigenvectors.

It was demonstrated that, for each pair of distint eigenvalues ω_r^2 and ω_s^2 , the corresponding eigenvectors obey the ortogonality condition,

$$\psi_s^T \mathbf{M} \psi_r = \delta_{rs} M_r, \quad \psi_s^T \mathbf{K} \psi_r = \delta_{rs} \omega_r^2 M_r.$$

Additional Orthogonality Relationships

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From

$$\mathbf{K} \mathbf{\psi}_s = \mathbf{\omega}_s^2 \mathbf{M} \mathbf{\psi}_s$$

premultiplying by $\psi_r^T KM^{-1}$ we have

$$\boldsymbol{\psi}_{\textit{r}}^{\textit{T}} \textit{K} \textit{M}^{-1} \textit{K} \, \boldsymbol{\psi}_{\textit{s}} = \boldsymbol{\omega}_{\textit{s}}^{2} \boldsymbol{\psi}_{\textit{r}}^{\textit{T}} \textit{K} \, \boldsymbol{\psi}_{\textit{s}}$$

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,

premultiplying the first equation by $\psi_r^T KM^{-1}KM^{-1}$

$$\psi_r^T \mathbf{K} \mathbf{M}^{-1} \mathbf{K} \mathbf{M}^{-1} \mathbf{K} \psi_s = \omega_s^2 \psi_r^T \mathbf{K} \mathbf{M}^{-1} \mathbf{K} \psi_s =$$

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and, generalizing,

$$\mathbf{\psi}_{r}^{T}\left(\mathbf{K}\mathbf{M}^{-1}\right)^{b}\mathbf{K}\,\mathbf{\psi}_{s}=\delta_{rs}\left(\omega_{r}^{2}\right)^{b+1}M_{r}.$$

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premultiplying the first eq. by $\psi_r^T \left(\emph{M} \emph{K}^{-1} \right)^2$ we have

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Choice of Property Formulati Defining $X_{rs}(k) = \psi_r^T \boldsymbol{M} (\boldsymbol{M}^{-1} \boldsymbol{K})^k \psi_s$ we have

$$\begin{cases} X_{rs}(0) = \psi_r^T \mathbf{M} \psi_s &= \delta_{rs} \left(\omega_s^2\right)^0 M_s \\ X_{rs}(1) = \psi_r^T \mathbf{K} \psi_s &= \delta_{rs} \left(\omega_s^2\right)^1 M_s \\ X_{rs}(2) = \psi_r^T \left(\mathbf{K} \mathbf{M}^{-1}\right)^1 \mathbf{K} \psi_s &= \delta_{rs} \left(\omega_s^2\right)^2 M_s \\ \dots \\ X_{rs}(n) = \psi_r^T \left(\mathbf{K} \mathbf{M}^{-1}\right)^{n-1} \mathbf{K} \psi_s &= \delta_{rs} \left(\omega_s^2\right)^n M_s \end{cases}$$

Observing that $(\mathbf{M}^{-1}\mathbf{K})^{-1}=(\mathbf{K}^{-1}\mathbf{M})^1$

$$\begin{cases} X_{rs}(-1) = \psi_r^T \left(\mathbf{M} \mathbf{K}^{-1} \right)^1 \mathbf{M} \psi_s &= \delta_{rs} \left(\omega_s^2 \right)^{-1} M_s \\ \dots \\ X_{rs}(-n) = \psi_r^T \left(\mathbf{M} \mathbf{K}^{-1} \right)^n \mathbf{M} \psi_s &= \delta_{rs} \left(\omega_s^2 \right)^{-n} M_s \end{cases}$$

finally

$$X_{rs}(k) = \delta_{rs} \omega_s^{2k} M_s$$
 for $k = -\infty, ..., \infty$.

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Choice of Property

Given a system whose state is determined by the generalized displacements x_j of a set of nodes, we define the flexibility coefficient f_{jk} as the deflection, in direction of x_j , due to the application of a unit force in correspondance of the displacement x_k .

The matrix $\mathbf{F} = [f_{jk}]$ is the *flexibility matrix*.

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In general, the dynamic degrees of freedom correspond to the points where there is

- application of external forces and/or
- presence of inertial forces.

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Given a load vector $\mathbf{p} = \{p_k\}$, the displacementent x_i is

$$x_j = \sum f_{jk} p_k$$

or, in vector notation,

$$x = F p$$

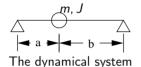
Flexibility Matrix

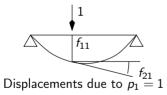
Example Stiffness Matrix Mass Matrix Damping Matrix Geometric Stiffness External Loading

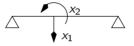
Flexibility Matrix Example

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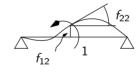








The degrees of freedom



and due to $p_2 = 1$.

Momentarily disregarding inertial effects, each node shall be in equilibrium under the action of the external forces and the elastic forces, hence taking into accounts all the nodes, all the external forces and all the elastic forces it is possible to write the vector equation of equilibrium

$$p = f_{\mathsf{S}}$$

and, substituting in the previos vector expression of the displacements

$$x = F f_{S}$$

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The *stiffness matrix* \boldsymbol{K} can be simply defined as the inverse of the flexibility matrix \boldsymbol{F} ,

$$K = F^{-1}$$
.

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Stiffness

The *stiffness matrix* \boldsymbol{K} can be simply defined as the inverse of the flexibility matrix \boldsymbol{F} ,

$$K = F^{-1}$$
.

As an alternative definition, consider an unary vector of displacements,

$$oldsymbol{e}^{(i)} = \left\{ \delta_{ij}
ight\}$$
 , $j = 1, \ldots, N$,

and the vector of nodal forces \mathbf{k}_i that, applied to the structure, produces the displacements $\mathbf{e}^{(i)}$

$$\mathbf{F} \mathbf{k}_i = \mathbf{e}^{(i)}, \qquad i = 1, \ldots, N.$$

Collecting all the ordered $\mathbf{e}^{(i)}$ in a matrix \mathbf{E} , it is clear that $\mathbf{E} \equiv \mathbf{I}$ and we have, writing all the equations at once,

$$F[k_i] = [e^{(i)}] = E = I.$$

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Collecting all the ordered $\mathbf{e}^{(i)}$ in a matrix \mathbf{E} , it is clear that $\mathbf{E} \equiv \mathbf{I}$ and we have, writing all the equations at once,

$$F[k_i] = [e^{(i)}] = E = I.$$

Collecting the ordered force vectors in a matrix $oldsymbol{K} = \left[ec{k}_i
ight]$ we have

$$FK = I$$
, $\Rightarrow K = F^{-1}$,

giving a physical interpretation to the columns of the stiffness matrix. Finally, writing the nodal equilibrium, we have

$$p = f_S = K x$$
.

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The elastic strain energy V can be written in terms of displacements and external forces.

$$V = \frac{1}{2} \boldsymbol{p}^T \boldsymbol{x} = \frac{1}{2} \begin{cases} \boldsymbol{p}^T \boldsymbol{F} \boldsymbol{p}, \\ \boldsymbol{x}^T \boldsymbol{K} \boldsymbol{x}. \end{cases}$$

Because the elastic strain energy of a stable system is always greater than zero, ${\pmb K}$ is a positive definite matrix.

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Choice of Property

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Because the elastic strain energy of a stable system is always greater than zero, \boldsymbol{K} is a positive definite matrix.

On the other hand, for an unstable system, think of a compressed beam, there are displacement patterns that are associated to zero strain energy.

Symmetry

Two sets of loads \mathbf{p}^A and \mathbf{p}^B are applied, one after the other, to an elastic system; the work done is

$$V_{AB} = \frac{1}{2} \boldsymbol{p}^{AT} \boldsymbol{x}^{A} + \boldsymbol{p}^{AT} \boldsymbol{x}^{B} + \frac{1}{2} \boldsymbol{p}^{BT} \boldsymbol{x}^{B}.$$

If we revert the order of application the work is

$$V_{BA} = \frac{1}{2} \boldsymbol{p}^{BT} \boldsymbol{x}^{B} + \boldsymbol{p}^{BT} \boldsymbol{x}^{A} + \frac{1}{2} \boldsymbol{p}^{AT} \boldsymbol{x}^{A}.$$

The total work being independent of the order of loading,

$$\boldsymbol{p}^{AT}\boldsymbol{x}^{B}=\boldsymbol{p}^{BT}\boldsymbol{x}^{A}.$$

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Expressing the displacements in terms of \mathbf{F} .

$${m p}^{A}{}^T{m F}\,{m p}^B={m p}^B{}^T{m F}{m p}^A,$$

both terms are scalars so we can write

$$\boldsymbol{p}^{A^T} \boldsymbol{F} \, \boldsymbol{p}^B = \left(\boldsymbol{p}^{B^T} \boldsymbol{F} \boldsymbol{p}^A \right)^T = \boldsymbol{p}^{A^T} \boldsymbol{F}^T \, \boldsymbol{p}^B.$$

Because this equation holds for every \boldsymbol{p} , we conclude that

$$\mathbf{F} = \mathbf{F}^T$$
.

The inverse of a symmetric matrix is symmetric, hence

$$K = K^T$$
.

Choice of Property Formulation

For the kind of *structures* we mostly deal with in our examples, problems, exercises and assignments, that is *simple structures*, it is usually convenient to compute first the flexibility matrix applying the Principle of Virtual Displacements and later the stiffness matrix, using inversion,

 $K = F^{-1}$.

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On the other hand, the PVD approach cannot work in practice for *real structures*, because the number of degrees of freedom necessary to model the structural behaviour exceeds our ability to apply the PVD...

The stiffness matrix for large, complex structures to construct different methods required are.

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E.g., the Finite Element Method.

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The most common procedure to construct the matrices that describe the behaviour of a complex system is the *Finite Element Method*, or *FEM*. The procedure can be sketched in the following terms:

▶ the structure is subdivided in non-overlapping portions, the *finite elements*, bounded by *nodes*, connected by the same nodes,

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- ▶ the structure is subdivided in non-overlapping portions, the *finite elements*, bounded by *nodes*, connected by the same nodes,
- ▶ the state of the structure can be described in terms of a vector **x** of generalized *nodal displacements*,

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- ▶ the element stiffness matrix, K_{el} establishes a linear relation between an element nodal displacements and forces,

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- ▶ the structure is subdivided in non-overlapping portions, the *finite elements*, bounded by *nodes*, connected by the same nodes,
- ▶ the state of the structure can be described in terms of a vector **x** of generalized *nodal displacements*,
- lacktriangle there is a mapping between element and structure DOFs, $i_{\mathsf{el}} \mapsto r$,
- ▶ the element stiffness matrix, K_{el} establishes a linear relation between an element nodal displacements and forces,
- ▶ for each FE, all local k_{ij} 's are contributed to the global stiffness k_{rs} 's, with $i \mapsto r$ and $j \mapsto s$, taking in due consideration differences between local and global systems of reference.

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The most common procedure to construct the matrices that describe the behaviour of a complex system is the *Finite Element Method*, or *FEM*. The procedure can be sketched in the following terms:

- ▶ the structure is subdivided in non-overlapping portions, the *finite elements*, bounded by *nodes*, connected by the same nodes,
- ▶ the state of the structure can be described in terms of a vector **x** of generalized *nodal displacements*,
- ▶ there is a mapping between element and structure *DOF*'s, $i_{el} \mapsto r$,
- ▶ the *element stiffness matrix*, **K**_{el} establishes a linear relation between an element nodal displacements and forces,
- ▶ for each FE, all local k_{ij} 's are contributed to the global stiffness k_{rs} 's, with $i \mapsto r$ and $j \mapsto s$, taking in due consideration differences between local and global systems of reference.

Note that in the *r*-th *global* equation of equilibrium we have internal forces caused by the nodal displacements of the *FE* that have nodes i_{el} such that $i_{el} \mapsto r$, thus implying that global K is a *banded* matrix.

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Consider a 2-D inextensible beam element, that has 4 *DOF*, namely two transverse end displacements x_1 , x_2 and two end rotations, x_3 , x_4 . The element stiffness is computed using 4 shape functions ϕ_i , the transverse displacement being $v(s) = \sum_i \phi_i(s) x_i$, $0 \le s \le L$, the different ϕ_i are such all end displacements or rotation are zero, except the one corresponding to index i.

The shape functions for a beam are

$$\begin{aligned} & \phi_1(s) = 1 - 3\left(\frac{s}{L}\right)^2 + 2\left(\frac{s}{L}\right)^3, & \phi_2(s) = 3\left(\frac{s}{L}\right)^2 - 2\left(\frac{s}{L}\right)^3, \\ & \phi_3(s) = \left(\frac{s}{L}\right) - 2\left(\frac{s}{L}\right)^2 + \left(\frac{s}{L}\right)^3 & \phi_4(s) = -\left(\frac{s}{L}\right)^2 + \left(\frac{s}{L}\right)^3. \end{aligned}$$

The element stiffness coefficients can be computed using, what else, the PVD: we compute the external virtual work done by a variation δx_i by the force due to a unit displacement x_j , that is k_{ij} ,

$$\delta W_{\rm ext} = \delta x_i k_{ij}$$
,

the virtual internal work is the work done by the variation of the curvature, $\delta x_i \phi_i''(s)$ by the bending moment associated with a unit x_j , $\phi_j''(s) EJ(s)$,

$$\delta W_{\rm int} = \int_0^L \delta x_i \varphi_i''(s) \varphi_j''(s) EJ(s) \, \mathrm{d}s.$$

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The equilibrium condition is the equivalence of the internal and external virtual works, so that simplifying δx_i we have

$$k_{ij} = \int_0^L \varphi_i''(s) \varphi_j''(s) EJ(s) ds.$$

For EJ = const,

$$\mathbf{f}_{S} = \frac{EJ}{L^{3}} \begin{bmatrix} 12 & -12 & 6L & 6L \\ -12 & 12 & -6L & -6L \\ 6L & -6L & 4L^{2} & 2L^{2} \\ 6L & -6L & 2L^{2} & 4L^{2} \end{bmatrix} \mathbf{x}$$

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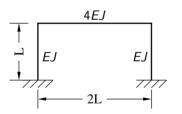
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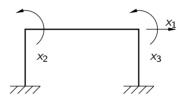
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The mass matrix maps the nodal accelerations to nodal inertial forces, and the most common assumption is to concentrate all the masses in nodal point masses, without rotational inertia, computed *lumping* a fraction of each element mass (or a fraction of the supported mass) on all its bounding nodes.

This procedure leads to a so called *lumped* mass matrix, a diagonal matrix with diagonal elements greater than zero for all the translational degrees of freedom and diagonal elements equal to zero for angular degrees of freedom.

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The mass matrix is definite positive *only* if all the structure *DOF*'s are translational degrees of freedom, otherwise *M* is semi-definite positive and the eigenvalue procedure is not directly applicable. This problem can be overcome either by using a *consistent* mass matrix or using the *static condensation* procedure.

Using our beam example as a reference, consider the inertial forces associated with a single nodal acceleration \ddot{x}_j , $f_{l,j}(s)=m(s)\varphi_j(s)\ddot{x}_j$ and denote with $m_{ij}\ddot{x}_j$ the reaction associated with the i-nth degree of freedom of the element, by the PVD

$$\delta x_i m_{ij} \ddot{x}_j = \int \delta x_i \varphi_i(s) m(s) \varphi_j(s) ds \ddot{x}_j$$

simplifying

$$m_{ij} = \int m(s) \varphi_i(s) \varphi_j(s) ds.$$

For $m(s) = \overline{m} = \text{const.}$

$$\mathbf{f_l} = \frac{\overline{m}L}{420} \begin{bmatrix} 156 & 54 & 22L & -13L \\ 54 & 156 & 13L & -22L \\ 22L & 13L & 4L^2 & -3L^2 \\ -13L & -22L & -3L^2 & 4L^2 \end{bmatrix} \ddot{\mathbf{x}}$$

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Pro

- some convergence theorem of FEM theory holds only if the mass matrix is consistent,
- sligtly more accurate results,
- no need for static condensation.

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Pro

- some convergence theorem of *FEM* theory holds only if the mass matrix is consistent.
- sligtly more accurate results,
- no need for static condensation.

Contra

- ▶ **M** is no more diagonal, heavy computational aggravation.
- static condensation is computationally beneficial, inasmuch it reduces the global number of degrees of freedom.

Damping Matrix

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For each element $c_{ij} = \int c(s) \phi_i(s) \phi_j(s) ds$ and the damping matrix \boldsymbol{C} can be assembled from element contributions.

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For each element $c_{ij} = \int c(s) \varphi_i(s) \varphi_j(s) \, \mathrm{d}s$ and the damping matrix \boldsymbol{C} can be assembled from element contributions.

However, using the FEM $\mathbf{C}^{\star} = \mathbf{\Psi}^{T} \mathbf{C} \mathbf{\Psi}$ is not diagonal and the modal equations are no more uncoupled!

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For each element $c_{ij} = \int c(s) \phi_i(s) \phi_j(s) ds$ and the damping matrix C can be assembled from element contributions.

However, using the FEM $\mathbf{C}^{\star} = \mathbf{\Psi}^{T} \mathbf{C} \mathbf{\Psi}$ is not diagonal and the modal equations are no more uncoupled!

The alternative is to write directly the global damping matrix, in terms of the underdetermined coefficients c_b ,

$$C = \sum_b c_b M \left(M^{-1} K \right)^b$$
.

With our definition of **C**.

$$oldsymbol{\mathcal{C}} = \sum_b \mathfrak{c}_b oldsymbol{M} \left(oldsymbol{M}^{-1} oldsymbol{K}
ight)^b$$
 ,

assuming normalized eigenvectors, we can write the individual component of ${\pmb C}^\star = {\pmb \Psi}^{{\cal T}} {\pmb C} \, {\pmb \Psi}$

$$c_{ij}^{\star} = \boldsymbol{\psi}_{i}^{T} \boldsymbol{C} \, \boldsymbol{\psi}_{j} = \delta_{ij} \sum_{b} \mathfrak{c}_{b} \omega_{j}^{2b}$$

due to the additional orthogonality relations, we recognize that now ${\pmb C}^{\star}$ is a diagonal matrix.

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$$oldsymbol{\mathcal{C}} = \sum_b \mathfrak{c}_b oldsymbol{M} \left(oldsymbol{M}^{-1} oldsymbol{K}
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due to the additional orthogonality relations, we recognize that now C^* is a diagonal matrix.

Introducing the modal damping C_j we have

$$C_j = \boldsymbol{\psi}_j^T \boldsymbol{C} \, \boldsymbol{\psi}_j = \sum_b \mathfrak{c}_b \omega_j^{2b} = 2\zeta_j \omega_j$$

and we can write a system of linear equations in the c_h .

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$$\ddot{q}_i + 2\zeta_i \omega_i \dot{q}_i + \omega_i^2 q_i = p_i^*$$

Using

$$\mathbf{C} = \mathfrak{c}_0 \mathbf{M} + \mathfrak{c}_1 \mathbf{K} + \mathfrak{c}_2 \mathbf{K} \mathbf{M}^{-1} \mathbf{K}$$

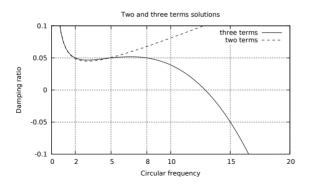
we have

$$2 \times 0.05 \begin{Bmatrix} \omega_1 \\ \omega_2 \\ \omega_3 \end{Bmatrix} = \begin{bmatrix} 1 & \omega_1^2 & \omega_1^4 \\ 1 & \omega_2^2 & \omega_2^4 \\ 1 & \omega_3^2 & \omega_3^4 \end{bmatrix} \begin{Bmatrix} \mathfrak{c}_0 \\ \mathfrak{c}_1 \\ \mathfrak{c}_2 \end{Bmatrix}$$

Solving for the c's and substituting above, the resulting damping matrix is orthogonal to every eigenvector of the system, for the first three modes, leads to a modal damping ratio that is equal to 5%.

Computing the coefficients \mathfrak{c}_0 , \mathfrak{c}_1 and \mathfrak{c}_2 to have a 5% damping at frequencies $\omega_1=2,\ \omega_2=5$ and $\omega_3=8$ we have $\mathfrak{c}_0=1200/9100,\ \mathfrak{c}_1=159/9100$ and $\mathfrak{c}_2=-1/9100.$

Writing $\zeta(\omega)=\frac{1}{2}\left(\frac{\mathfrak{c}_0}{\omega}+\mathfrak{c}_1\omega+\mathfrak{c}_2\omega^3\right)$ we can plot the above function, along with its two term equivalent $(\mathfrak{c}_0=10/70,\mathfrak{c}_1=1/70)$.



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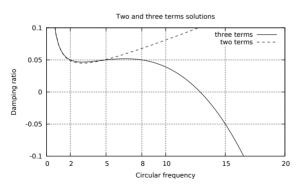
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Negative damping? No. thank you: use only an even number of terms.

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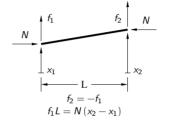
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Choice of Property

A common assumption is based on a linear approximation, for a beam element



It is possible to compute the geometrical stiffness matrix using *FEM*, shape functions and PVD,

$$k_{\mathsf{G},ij} = \int N(s) \varphi_i'(s) \varphi_j'(s) \,\mathrm{d}s,$$

for constant N

$$K_{G} = \frac{N}{30L} \begin{bmatrix} 36 & -36 & 3L & 3L \\ -36 & 36 & -3L & -3L \\ 3L & -3L & 4L^{2} & -L^{2} \\ 3L & -3L & -L^{2} & 4L^{2} \end{bmatrix}$$

Following the same line of reasoning that we applied to find nodal inertial forces, by the PVD and the use of shape functions we have

$$p_i(t) = \int p(s, t) \phi_i(s) ds.$$

For a constant, uniform load $p(s,t)=\overline{p}={
m const.}$ applied on a beam element,

$$\boldsymbol{p} = \overline{p}L \left\{ \frac{1}{2} \quad \frac{1}{2} \quad \frac{L}{12} \quad -\frac{L}{12} \right\}^T$$

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Simplified Approach

Some structural parameter is approximated, only translational *DOF*'s are retained in dynamic analysis.

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Simplified Approach

Some structural parameter is approximated, only translational *DOF*'s are retained in dynamic analysis.

Consistent Approach

All structural parameters are computed according to the FEM, and all DOF's are retained in dynamic analysis.

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All structural parameters are computed according to the *FEM*, and all *DOF*'s are retained in dynamic analysis.

If we choose a simplified approach, we must use a procedure to remove unneeded structural *DOF*'s from the model that we use for the dynamic analysis.

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If we choose a simplified approach, we must use a procedure to remove unneeded structural *DOF*'s from the model that we use for the dynamic analysis.

Enter the Static Condensation Method.

Static Condensation

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We have, from a *FEM* analysis, a stiffnes matrix that uses all nodal *DOF*'s, and from the lumped mass procedure a mass matrix were only translational (and maybe a few rotational) *DOF*'s are blessed with a non zero diagonal term.

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We have, from a *FEM* analysis, a stiffnes matrix that uses all nodal *DOF*'s, and from the lumped mass procedure a mass matrix were only translational (and maybe a few rotational) *DOF*'s are blessed

and partition the displacement vector \mathbf{x} in two subvectors: a) \mathbf{x}_A , all the DOF's that are associated with inertial forces and b) \mathbf{x}_B , all the remaining DOF's not associated with inertial forces.

with a non zero diagonal term. In this case, we can always rearrange

$$\mathbf{x} = \left\{ \mathbf{x}_{A} \quad \mathbf{x}_{B} \right\}^{T}$$

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Static Condensation Example After rearranging the *DOF*'s, we must rearrange also the rows (equations) and the columns (force contributions) in the structural matrices, and eventually partition the matrices so that

$$egin{aligned} \left\{ egin{aligned} \mathbf{f}_I \\ \mathbf{0} \end{matrix}
ight\} &= \left[egin{aligned} \mathbf{M}_{AA} & \mathbf{M}_{AB} \\ \mathbf{M}_{BA} & \mathbf{M}_{BB} \end{matrix}
ight] \left\{ egin{aligned} \ddot{\mathbf{x}}_A \\ \ddot{\mathbf{x}}_B \end{matrix}
ight\} \end{aligned} \\ \mathbf{f}_S &= \left[egin{aligned} \mathbf{K}_{AA} & \mathbf{K}_{AB} \\ \mathbf{K}_{BA} & \mathbf{K}_{BB} \end{matrix}
ight] \left\{ egin{aligned} \mathbf{x}_A \\ \mathbf{x}_B \end{matrix}
ight\} \end{aligned}$$

with

$$\mathbf{\textit{M}}_{BA} = \mathbf{\textit{M}}_{AB}^{\mathsf{T}} = \mathbf{0}, \quad \mathbf{\textit{M}}_{BB} = \mathbf{0}, \quad \mathbf{\textit{K}}_{BA} = \mathbf{\textit{K}}_{AB}^{\mathsf{T}}$$

After rearranging the DOF's, we must rearrange also the rows (equations) and the columns (force contributions) in the structural matrices, and eventually partition the matrices so that

$$\begin{cases} \mathbf{f}_I \\ \mathbf{0} \end{cases} = \begin{bmatrix} \mathbf{M}_{AA} & \mathbf{M}_{AB} \\ \mathbf{M}_{BA} & \mathbf{M}_{BB} \end{bmatrix} \begin{cases} \ddot{\mathbf{x}}_A \\ \ddot{\mathbf{x}}_B \end{cases}$$

$$\mathbf{f}_S = \begin{bmatrix} \mathbf{K}_{AA} & \mathbf{K}_{AB} \\ \mathbf{K}_{BA} & \mathbf{K}_{BB} \end{bmatrix} \begin{cases} \mathbf{x}_A \\ \mathbf{x}_B \end{cases}$$

with

$$\mathbf{\textit{M}}_{BA} = \mathbf{\textit{M}}_{AB}^{\mathsf{T}} = \mathbf{0}, \quad \mathbf{\textit{M}}_{BB} = \mathbf{0}, \quad \mathbf{\textit{K}}_{BA} = \mathbf{\textit{K}}_{AB}^{\mathsf{T}}$$

Finally we rearrange the loadings vector and write...

Static Condensation, 3

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... the equation of dynamic equilibrium,

$$egin{aligned} oldsymbol{
ho}_A &= oldsymbol{M}_{AA}\ddot{oldsymbol{x}}_A + oldsymbol{M}_{AB}\ddot{oldsymbol{x}}_B + oldsymbol{K}_{AA}oldsymbol{x}_A + oldsymbol{K}_{AB}oldsymbol{x}_B \\ oldsymbol{
ho}_B &= oldsymbol{M}_{BA}\ddot{oldsymbol{x}}_A + oldsymbol{M}_{BB}\ddot{oldsymbol{x}}_B + oldsymbol{K}_{BA}oldsymbol{x}_A + oldsymbol{K}_{BB}oldsymbol{x}_B \end{aligned}$$

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... the equation of dynamic equilibrium,

$$egin{aligned} oldsymbol{p}_A &= oldsymbol{M}_{AA}\ddot{f x}_A + oldsymbol{M}_{AB}\ddot{f x}_B + oldsymbol{K}_{AA}oldsymbol{x}_A + oldsymbol{K}_{AB}oldsymbol{x}_B \\ oldsymbol{p}_B &= oldsymbol{M}_{BA}\ddot{f x}_A + oldsymbol{M}_{BB}\ddot{f x}_B + oldsymbol{K}_{BA}oldsymbol{x}_A + oldsymbol{K}_{BB}oldsymbol{x}_B \end{aligned}$$

The terms in red are zero, so we can simplify

$$oldsymbol{M}_{AA}\ddot{oldsymbol{x}}_A + oldsymbol{K}_{AA}oldsymbol{x}_A + oldsymbol{K}_{AB}oldsymbol{x}_B = oldsymbol{p}_A \ oldsymbol{K}_{BA}oldsymbol{x}_A + oldsymbol{K}_{BB}oldsymbol{x}_B = oldsymbol{p}_B$$

solving for x_B in the 2nd equation and substituting

$$oldsymbol{x}_B = oldsymbol{K}_{BB}^{-1} oldsymbol{p}_B - oldsymbol{K}_{BB}^{-1} oldsymbol{K}_{BA} oldsymbol{x}_A \ oldsymbol{p}_A - oldsymbol{K}_{AB} oldsymbol{K}_{BB}^{-1} oldsymbol{p}_B = oldsymbol{M}_{AA} \ddot{oldsymbol{x}}_A + ig(oldsymbol{K}_{AA} - oldsymbol{K}_{AB} oldsymbol{K}_{BB}^{-1} oldsymbol{K}_{BA}ig) oldsymbol{x}_A$$

Matrices Science Be

Going back to the homogeneous problem, with obvious positions we can write

$$\left(\overline{\textbf{\textit{K}}}-\omega^2\overline{\textbf{\textit{M}}}\right)\psi_{\textit{A}}=\textbf{0}$$

but the ψ_A are only part of the structural eigenvectors, because in essentially every application we must consider also the other DOF's, so we write

$$\psi_i = egin{cases} \psi_{A,i} \ \psi_{B,i} \end{cases}$$
 , with $\psi_{B,i} = extbf{\emph{K}}_{BB}^{-1} extbf{\emph{K}}_{BA} \psi_{A,i}$

Example

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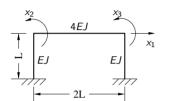
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Static Condensation Example



$$\mathbf{K} = \frac{2EJ}{L^3} \begin{bmatrix} 12 & 3L & 3L \\ 3L & 6L^2 & 2L^2 \\ 3L & 2L^2 & 6L^2 \end{bmatrix}$$

Disregarding the factor $2EJ/L^3$,

$$\mathbf{K}_{BB} = L^2 \begin{bmatrix} 6 & 2 \\ 2 & 6 \end{bmatrix}$$
, $\mathbf{K}_{BB}^{-1} = \frac{1}{32L^2} \begin{bmatrix} 6 & -2 \\ -2 & 6 \end{bmatrix}$, $\mathbf{K}_{AB} = \begin{bmatrix} 3L & 3L \end{bmatrix}$

The matrix \overline{K} is

$$\overline{\mathbf{K}} = \frac{2EJ}{L^3} \left(12 - \mathbf{K}_{AB} \mathbf{K}_{BB}^{-1} \mathbf{K}_{AB}^T \right) = \frac{39EJ}{2L^3}$$