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EVALUATION OF EIGEN VALUES AND EIGEN VECTORS FOR FREE VIBRATION ANALYSIS USING FINITE ELEMENT PROCEDURES R R GADPAL

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Abstract: The generalised problem in free vibration analysis of any structure is that of evaluating an eigenvalue λ (= ω^2), which is a measure of the frequency of vibration together with the corresponding eigenvector x indicating the mode shape. This presentation consists of determination of fundamental vibration frequencies of a general structure. The frequencies can be determined by the finite element method using characteristic polynomial technique, Vector iteration technique and transformation methods.. Power iteration, inverse iteration and subspace iteration methods use the property of the Rayleigh quotient. Power iteration leads to evaluation of the largest eigenvalue. Subspace iteration technique is suitable for large scale problems and used in several codes. The inverse iteration scheme can be used for evaluating the lowest eigenvalues. The basic approach in transformation method is to transform the matrices to a simpler form and then to determine the eigenvalues and eigenvectors. In this presentation the main focus is to discuss the transformation method, to develop the related computer program and validate the results. The results show good agreement with those determined by classical method.

Keywords: Free vibration analysis, Fundamental natural frequencies, Mode shapes, Eigenvalues and Eigenvectors , Generalised eigenproblem, Standard eigenproblem, Stiffness and mass matrices.



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INTRODUCTION

Free vibration analysis is often required for most of the important structures / structural elements in the field of civil, mechanical, automobile, aerospace, optical, marine, nuclear and structural engineering. The differential characteristics in free vibration analysis enable engineers to design better and lighter structures. The Study of their free vibration behaviour is very important to the structural engineers when these structures are subjected to external complicated dynamic loads such as earthquake, wind, impact and wave forces. An understanding of the free vibration frequencies of any system (especially, the fundamental frequency) is the prerequisite to the understanding of its response to forced vibration. In civil engineering, buildings, horizontal floors, beams, columns are directly exposed to static and dynamic loadings. To determine the eigenvalues and eigenvectors which are the measure of the frequency of vibration and mode shapes the suitable method is to be selected by the analyst as per the requirements in design of structures.

II LITERATURE REVIEW

The literature of eigenproblems is quite large, both for theoretical aspects and numerical algorithms, and only a small fraction of it can be cited. Hughes T.J.R. [1], Bathe K. J. [2] and Kardestuncer H. [3] have written extensive discussions about eigenproblems in their Finite Element Books. Wilkinson J.H. [4] wrote a book about the algebraic eigenvalue problem in 1965. Bathe K. J. and Wilson E.L. [5] published a paper on solution methods for eigenvalue problems in structural mechanics in 1973. In 1980, Parlett B.N. [6] introduced the method for solution of the symmetric eigenvalue problem. In 1984, Jennings A. [7] wrote about eigenvalue methods for vibration analysis. Sehmi N.S. [8] gave large order structural eigenvalue techniques in 1989. Cheung Y.K. and Leung A.Y.T. [9]. published a book on finite element methods in dynamics in 1991 In 1994, Tichler V.A. and Venk ayya V.B. [10] evaluated eigenvalue routines for large scale applications. Bertolini A.F. [11] reviewed eigensolution procedures for linear dynamic finite element analysis in 1998.

Undamped free vibration analysis of the entire building is performed as per established methods of mechanics using appropriate masses and elastic stiffness of the structural system to obtain natural period (T) and mode shape $\{\phi\}$ of those of its modes of vibration that need to be considered as per I.S.1893-2002(part 1) clause No. 7.8.4.2.[12]

Analytical solutions for dynamic response of structures are available for very few cases i.e. for structural elements with simple geometry and boundary conditions. But for elements with complex boundary and boundary conditions, solutions are possible only with the help of numerical methods. The most commonly used numerical method is finite element method.

Thus the present study evaluates the first few dominant modes of vibration frequencies of structures. The frequencies estimated by the proposed formulation and program coincide well with those obtained by the finite element method, which can serve as a design aid for structural engineers.

III. FINITE ELEMENT FORMULATION

For a positive definite symmetric stiffness matrix [K] of size n x n, there are n real eigenvalues and corresponding eigenvectors satisfying equation (1).

$$[K] - \lambda [M] \{x\} = \{0\} \qquad \dots (1)$$

The above eigenproblem asks for the values of a scalar λ such that the matrix equation (1) has solutions other than trivial solution $\{x\} = \{0\}$. There are at most n nonzero roots λ_i , not necessarily all distinct. The λ_i are called eigenvalues, characteristic values, latent roots or principal values. The eigenvalues may be arranged in ascending order such that

$$0 \le \lambda_1 \le \lambda_2 \le \dots \le \lambda_n$$
 ... (2)

The vector corresponding to each λ_i is an $\{x\}_i$ called as an eigenvector, characteristic vector, principal vector, normal mode or natural mode. The eigenvectors possess the property of being orthogonal with respect to both the stiffness and mass matrices:

$$\mathbf{x}_i^{\mathsf{T}} \mathbf{M} \mathbf{x}_j = 0$$
 if $i \neq j$...(3)

The lengths of eigenvectors are generally normalized so that

$$\mathbf{x}_{i}^{\mathsf{T}} \mathbf{M} \mathbf{x}_{i} = 1$$
 ... (4)

The foregoing normalization of the eigenvectors leads to the relation

$$\mathbf{x}_{i}^{\mathsf{T}} \mathbf{K} \mathbf{x}_{i} = \lambda_{i}$$
 ... (5)

In many codes, other normalization schemes are also used. The length of an eigenvector may be fixed by setting its largest component to a preset value, say, unity.

The eqn (1) in the form ($[K] - \lambda [M]$) $\{x\} = \{0\}$ is called *generalised eigenproblem* or simply eigenproblem and further can be simplified as

$$[K] \{x\} = \lambda [M] \{x\} \qquad \dots (6)$$

If , however, matrix [M] happens to be identity matrix [I] or premultiplying both sides of above equation by [M -1], we get,

[M
$$^{-1}$$
] [K] {x} = λ [M $^{-1}$] [M]) {x}
[A] {x} = λ [I] {x}
or [A] {x} = λ {x} ...(7)
where, [A] = [M $^{-1}$] [K]

Equation (6) is called standard eigenproblem in which matrix [A] will be un-symmetric and moreover it is necessary to find [M -1].

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The eigenvalue and eigenvector evaluation procedures fall into the following basic categories.

- 1. Characteristic polynomial technique
- 2. Vector iteration techniques
- 3. Transformation methods

Amongst these methods, transformation methods are suitable for large scale problems and will be discussed in details.

Characteristic Polynomial

From eqn.(6), we have $([K] - \lambda [M])\{x\} = \{0\}$. If the eigenvector is to be non-trivial, the required condition is

$$det ([K] - \lambda [M]) = \{0\}$$
 ... (8)

This represents the characteristic polynomial in λ .

The characteristic polynomial method can solve 2 x 2 problems by hand calculations. However it is also found uneconomical for computer usage because it is rather tedious and requires further mathematical considerations. We now discuss the other two categories.

Vector Iteration Methods

Various vector iteration methods use the properties of *Rayleigh Quotient*. For the generalised eigenvalue problem given in eqn. (6), Rayleigh quotient can be defined as ,

$$Q(\mathbf{v}) = \frac{\mathbf{v}^{\mathrm{T}} \mathbf{K} \mathbf{v}}{\mathbf{v}^{\mathrm{T}} \mathbf{M} \mathbf{v}} \dots$$

(9)

where, **v** is an arbitrary vector. A fundamental property of the Rayleigh quotient is that it lies between the smallest and largest eigenvalue.

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$$\lambda_1 \leq Q(\mathbf{v}) \leq \lambda_n$$
 ... (10)

In the inverse iteration scheme [14], we start with a trial vector \mathbf{x}^{o} and obtain the eigenvector \mathbf{x}^{k} after normalization and satisfying the requirements of tolerance. This scheme converges to the lowest eigenvalue, provided the trial vector does not coincide with the one of the eigenvectors. Other eigenvalues can be obtained by shifting, or by taking the trial vector from a space that is M – orthogonal to the calculated eigenvectors.

Transformation Methods

The basic approach in this method is to transform the matrices to a simpler form and then to determine the eigenvalues and eigenvectors. The major methods in this category are the generalised Jacobi Method and the QR method. These methods are suitable for large-scale problems. In the QR method, the matrices are firs reduced to tridigonalization form using Householder matrices. The generalised Jacobi Method uses the transformation to simultaneously diagonalize the stiffness and mass matrices. This method needs the full matrix locations and is quite efficient for calculating all eigenvalues and eigenvectors for small problems.

If all the eigenvectors are arranged as columns of a square matrix \mathbf{X} and all eigenvalues as the diagonal elements of a square matrix $\mathbf{\Lambda}$, then the generalised eigenvalue problem can be written in the form

(13)

Using M- orthonormality of eigenvectors, we have,

$$[X]^{T}[M][X] = [\Lambda]$$
 ... (14)

and

$$[X]^T[K][X] = [I]$$
 ... (15)

where [I] = Identity matrix

Generalised Jacobi Method

In the generalised Jacobi method a series of transformations P_1 , P_2 ,...., P_l are used such that if P represents the product

$$P = P_1 P_2 P_l$$
 ... (16)

Then the off diagonal terms of P^TKP and P^TMP are zero. In practice, the off-diagonal terms are set to be less than a value smaller than tolerance.

$$[K^*] = [P_I]^T \dots [P_2]^T [P_1]^T [K] [P_1] [P_2] \dots [P_I]$$
 ... (17a)

$$[M^*] = [P_I]^T \dots [P_2]^T [P_1]^T [M] [P_1] [P_2] \dots [P_I]$$
 ... (17b)

[K^*] and [M^*] are the diagonal matrices. To eliminate [K_{ii}] and [M_{ii}] simultaneously we have to let transformation matrix [P] in such a way that at step k,

$$\begin{bmatrix} 1 & 0 & 0 & 0 & 0 \\ 0 & 1 & 0 & \alpha & 0 \\ 0 & 0 & 1 & 0 & 0 \\ 0 & \beta & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 & 1 \end{bmatrix} \text{ and thus, } \begin{bmatrix} P_k \end{bmatrix}^T = \begin{bmatrix} 1 & 0 & 0 & 0 & 0 \\ 0 & 1 & 0 & \beta & 0 \\ 0 & 0 & 1 & 0 & 0 \\ 0 & \alpha & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 & 1 \end{bmatrix} \dots (18)$$

[P_k] has all diagonal elements equal to 1, has a value of α at row i and column j and β at row j and column i, and all other elements equal to zero. The scalars α and β are chosen so that the ij locations of P_k^TKP and P_k^TMP are simultaneously zero.

$$\begin{bmatrix} 1 & 0 & 0 & 0 & 0 \\ 0 & 1 & 0 & \beta & 0 \\ 0 & 0 & 1 & 0 & 0 \\ 0 & \alpha & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 & 1 \end{bmatrix} \begin{bmatrix} K_{11} & K_{12} & K_{13} & K_{14} & K_{15} \\ K_{21} & K_{22} & K_{23} & K_{24} & K_{25} \\ K_{31} & K_{32} & K_{33} & K_{34} & K_{35} \\ K_{41} & K_{42} & K_{43} & K_{44} & K_{45} \\ K_{51} & K_{52} & K_{53} & K_{54} & K_{55} \end{bmatrix} \begin{bmatrix} 1 & 0 & 0 & 0 & 0 \\ 0 & 1 & 0 & \alpha & 0 \\ 0 & 0 & 1 & 0 & 0 \\ 0 & \beta & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 & 1 \end{bmatrix}$$

Considering only 2 and 4 rows and 2 and 4 columns are affected

$$\begin{aligned} & \text{Considering only 2 and 4 rows and 2 and 4 columns are affected} \\ & = \begin{bmatrix} K_{11} & K_{12} & K_{13} & K_{14} & K_{15} \\ K_{21} + \beta K_{41} & K_{22} + \beta K_{42} & K_{23} + \beta K_{43} & K_{24} + \beta K_{44} & K_{25} + \beta K_{45} \\ K_{31} & K_{32} & K_{33} & K_{34} & K_{35} \\ K_{41} + \alpha K_{21} & K_{42} + \alpha K_{22} & K_{43} + \alpha K_{23} & K_{44} + \alpha K_{24} & K_{45} + \alpha K_{25} \\ K_{51} & K_{52} & K_{53} & K_{54} & K_{55} \end{bmatrix} \begin{bmatrix} 1 & 0 & 0 & 0 & 0 \\ 0 & 1 & 0 & \alpha & 0 \\ 0 & 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 & 1 \end{bmatrix} \\ & = \begin{bmatrix} K_{11} & K_{12} + \beta K_{14} & K_{13} & K_{14} + \alpha K_{12} & K_{15} \\ K_{21} + \beta K_{41} & K_{22} + \beta K_{42} + \beta K_{24} + \beta^2 K_{44} & K_{23} + \beta K_{43} & K_{24} + \beta K_{44} + \alpha K_{22} + \alpha \beta K_{42} & K_{25} + \beta K_{45} \\ K_{31} & K_{32} + \beta K_{34} & K_{33} & K_{34} + \alpha K_{34} & K_{35} \\ K_{41} + \alpha K_{21} & K_{42} + \alpha K_{22} + \beta K_{44} + \alpha \beta K_{24} & K_{43} + \alpha K_{23} & K_{44} + \alpha K_{24} + \alpha K_{42} + \alpha^2 K_{22} & K_{45} + \alpha K_{25} \\ K_{51} & K_{52} + \beta K_{54} & K_{53} & \alpha K_{52} + K_{54} & K_{55} \end{bmatrix} \end{bmatrix}$$

Thus new $K_{24} = K_{24} + \beta K_{44} + \alpha K_{22} + \alpha \beta K_{42} = (1 + \alpha \beta) K_{24} + \alpha K_{22} + \beta K_{44}$

Now in general in order to make the non-diagonal elements equal to zero we can write $P_k^T KP$ and P_k^TMP as,

$$(1+ \alpha \beta)^{K_{ij}} + \alpha K_{ii} + \beta K_{jj} = 0$$
 ... (19) and
$$(1+ \alpha \beta)^{M_{ij}} + \alpha M_{ii} + \beta M_{jj} = 0$$
 ... (20)

Premultiplying eqn (19) by M_{ij} and eq. (20) by K_{ij} and solving the simultaneous equations, we get

$$\alpha \left({^{K_{ii}}M_{ij}}_{-} {^{M_{ii}}K_{ij}} \right) + \beta \left({^{K_{jj}}M_{ij}}_{-} {^{M_{jj}}K_{ij}} \right) = 0$$
or $\alpha \left({^{A}} \right) + \beta \left({^{B}} \right) = 0$... (21)

where, $A = (K_{ii}M_{ij} - M_{ii}K_{ij})$ and $B = (K_{jj}M_{ij} - M_{jj}K_{ij})$

Solving Eqn (21), $\beta = -\frac{\alpha A}{B}$ and after substituting in eqn (19), we get

$$K_{ij} = \alpha^2 \frac{A}{B} K_{ij} + \alpha K_{ii} + \beta K_{jj} = 0$$
 ... (22)

Multiplying eqn (22) by B and dividing by K_{ij} , we get

$$B - \frac{\alpha^2}{A} + \frac{\alpha K_{ii}(B)}{K_{ij}} - \frac{\alpha K_{jj}(A)}{K_{ij}} = 0$$

$$\alpha^2_{A + \alpha} \left[\frac{K_{jj}(A)}{K_{ij}} - \frac{K_{ii}(B)}{K_{ij}} \right] - B = 0 \qquad ... (23)$$
Introducing $C = \left[\frac{K_{jj}(A)}{K_{ij}} - \frac{K_{ii}(B)}{K_{ij}} \right]$ and substituting values of A and B, we get

 $C = M_{jj}K_{ii} - M_{ii}K_{jj}$ Multiplying eqn (23) by 0.5, we get

$$0.5^{\alpha^2} A + 0.5 \quad \alpha C - 0.5 B = 0$$
 ... (24)

Solving eqn (24), we get

$$\alpha = \frac{-0.5 C \pm \sqrt{0.25 C^2 + AB}}{A} \qquad ... (25)$$

Particularly,

When,
$$A \neq 0$$
, $B \neq 0$, $\alpha = \frac{-0.5 \ C \pm \sqrt{0.25 \ C^2 + AB}}{A}$ and $\beta = -\frac{\alpha A}{B}$;
$$A = 0, \ \beta = 0 \ \text{and} \ \alpha = -\frac{\frac{K_{ij}}{K_{ii}}}{\frac{K_{ij}}{K_{jj}}};$$

$$B = 0, \ \alpha = 0 \ \text{and} \ \beta = -\frac{K_{ij}}{K_{jj}} \qquad \dots (26)$$

When both A and B are zero, any one of the two values listed can be chosen.

Adopting the above procedure for determination of matrices $[K^*]$ and $[M^*]$ in eqs. (17), the eigenvalues and eigenvectors are given by,

$$\lambda \text{ii} = \frac{\frac{K_{ii}}{M_{ii}}}{M_{ii}} \text{ and } \mathbf{x}_{ij} = \frac{P_{ij}}{\sqrt{Mi}}, \text{ or in general}$$

$$\Lambda = \frac{K^*}{M^*} \text{ and } \mathbf{X} = \frac{P}{\sqrt{M}^*}$$
... (27)
$$\begin{bmatrix} \frac{1}{M_{11}} & \cdots & 0 & 0 \\ 0 & \frac{1}{M_{22}} & 0 & 0 \\ \vdots & & \ddots & \vdots \\ 0 & \cdots & 0 & \frac{1}{M_{nn}} \end{bmatrix} \text{ and } \frac{1}{\sqrt{M}^*} = \begin{bmatrix} \frac{1}{M_{11}} & \cdots & 0 & 0 \\ 0 & \frac{1}{M_{22}} & 0 & 0 \\ \vdots & & \ddots & \vdots \\ 0 & \cdots & 0 & \frac{1}{M_{nn}} \end{bmatrix}$$
where, $\frac{1}{M^*} = \frac{P_{ij}}{\sqrt{M}}$ and $\frac{1}{\sqrt{M}} = \frac{P_{ij}}{\sqrt{M}} = \frac{P_{ij}}{\sqrt{M}} = \frac{P_{ij}}{\sqrt{M}}$ and $\frac{1}{\sqrt{M}} = \frac{P_{ij}}{\sqrt{M}} = \frac{P_$

In the generalized Jacobi program, the elements of K and M are zeroed out in the order indicated in fig.1. Once P_k is defined by determining α and β , P_k^T [] P_k can be performed on Kand M as shown in fig.2. Also by starting with P = I, the product PP_k is computed after each step. When all elements are covered as shown in fig.1, one pass is completed. After operations at step k, some of the previously zeroed elements are altered. Another pass is conducted to check for the value of the diagonal elements. The transformation is performed if the elements at ij is larger than a tolerance value. A tolerance 10^{-6} smallest Kii is used for stiffness, and 10^{-6} X largest Mii is used for the mass. The tolerance can be redefined for higher accuracy. The process stops when all off-diagonal elements are less than the tolerance.

Fig.1 Diagonalization

Fig .2 Multiplication of $P_k^T[\]P_k$ Column Row $row i + \beta (row j)$ Symmetric $row j + \alpha(row i)$

If the diagonal masses are less than the tolerance, the diagonal value is replaced by the tolerance value, thus, a large eigenvalue will be obtained. In this method **K** need not be positive definite. On the basis of above formulation a computer program is developed and results are compared with the standard available results.

IV NUMERICAL EXAMPLE

Example 1: A cantilever having geometrical and material properties is as shown in fig. 3 Determine all the eigenvalues and eigenvectors for the beam shown in fig. 3 using the program developed on the basis of above formulation.

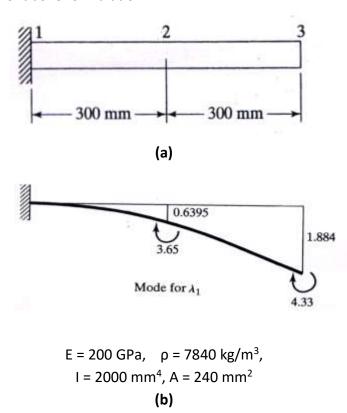


Fig. 3 Cantilever Beam Model

Solution:

The degree of freedom at each node is deflection (amplitude) w and slope $\partial w/\partial x$. Using standard finite element procedure [15], we obtain the stiffness and mass matrices by hand calculations:

The input data for the developed program is same as that for inverse iteration program. However, the program coverts to full matrices in calculations. Convergence occurs at the fourth sweep. The solution is presented in Table 1. and compared with the standard results.

Table 1. Natural frequencies & Mode shapes

Results reported in literature in full form		Results obtained with present analysis	
Natural	Mode shapes (eigen vectors)	Natural	Mode shapes (eigen vectors) U
frequency	U	frequency	
(eigenvalue)		(eigen value) λ	
λ			
2.0304 x 10 ⁴	0.64, 3.65, 1.88, 4.32	2.032x 10 ⁴	0.6395 , _{3.652} , 1.884 , 4.326
8.0987 x 10 ⁵	-1.37, 1.39, 1.901, 15.27	8.183x 10 ⁵	-1.375, 1.390, 1.906, 15.25
9.2651×10^6	-0.20, 27.16, -2.12,-33.84	9.35×10^6	-0.2043, 27.17, -2.121,-33.81
7.7974×10^7	0.8986, 30.89, 3.546, 119.15	8.268×10^7	0.9055, 30.89, 3.776, 119.2

The natural fundamental frequencies of the structure / component can be determined using classical methods as well as using finite element technique which is widely reported in the literature. It is well known that the numerical analysis results are valid only for particular values of the parameters considered in the analysis. The structural engineers concerned with dynamic analysis or design of structures need a design formula or program for rapid determination of the governing natural frequency. The numerical values obtained by running the program are quite gratifying with those reported in literature[16] and reproduced in Table 1.

VI CONCLUSIONS:

The fundamental frequencies and mode shapes were determined by finite element method. The devised program is quite useful for determining Fundamental natural frequencies, Mode shapes, by evaluating eigenvalues and eigenvectors in the Generalised eigenproblem.,

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