

An Improved Matrix Perturbation Method for Modified Structures

Zhong-hai XU^{1,*}, Zhong-yuan NIU², Jian-zhao ZHANG³ and
Rong-guo WANG⁴

^{1,2,3,4}National Key Laboratory of Science and Technology on Advanced Composites in
Special Environments

Harbin Institute of Technology, Harbin, 150080, PR China.

*xuzh@hit.edu.cn

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Abstract. In this paper, an improved matrix perturbation method for modified structures is presented. First we construct an iterative format on the basis of the matrix perturbation theory. By introducing a pair of initial eigensolution, we can obtain the natural vibration mode of the modified structure, and then use the Rayleigh quotient to compute the corresponding natural frequency. In this way, the computation accuracy can be greatly improved. Finally, a numerical example is included to demonstrate the validity of the proposed method.

Introduction

Structural dynamic reanalysis play an important role in structural design and optimization. Many scholars have paid their attentions to this subject.

The traditional method is the Taylor series expansion method, and later approximation method using reduced basis vectors was proposed. In the papers of Noor¹ and Nair et al.², the basis vector is the Taylor series expansion of the eigenpairs with respect to the structural parameters, and in the papers of Aktas and Moses³ and Kirsch⁴, the basis vector is the binomial expansion by solving a static problem. However, the choice of the basis vector is still a research issue. Kirsch^{5,6,7} introduced the combined approximation algorithm to the vibration analysis process, and made the equivalent treatment for the modal equations and the static equilibrium equations, then constructed the reduced basis vector in the Krylov subspace, finally, gave the error evaluation and solving process. He et al.⁸ made the modal reanalysis research for structural large topological modifications with added degrees of freedom. Rong et al.⁹, Ma et al.¹⁰, Yang et al.¹¹ and Chen et al.¹² proposed the extended combined approximation method, and improved the calculation accuracy using the Rayleigh quotient. Zhang et al.¹³ presented the modified combined approximation method based on the inverse iteration and combined approximation algorithm, and improved the calculation accuracy for structural large modifications. Liu et al.¹⁴ made the

research on the eigenvalue repeated analysis problems. Chen et al.15,16 made the research on the structural dynamic responds repeated analysis problems by introducing the Epsilon algorithm and combining with the Newmann series expansion.

Though many algorithms above are involved, the matrix perturbation method is the classical approach. Yang et al.17 proposed a method based on the Pade approximation to improve the perturbation theory. Chen et al.18 improved the calculation accuracy of the perturbation method by combining the matrix perturbation method with the Rayleigh quotient, but it isn't usually applicable to the case of large parameter modifications, as noted in the paper of Chen et al.19. So it is necessary to improve the perturbation method in the calculation accuracy and the range of application.

In this paper, we present an improved matrix perturbation method. By combining the iteration algorithm with the Rayleigh quotient, we can obtain the eigensolution of the modified structure, and the computational accuracy can be greatly improved. Numerical example is used to illustrate the validity of the proposed method.

Theoretical Background

Consider the following general structural vibration eigenvalue problem

$$[K]\{u_i\} = \lambda_i[M]\{u_i\}. \quad (1)$$

$$\{u_i\}^T[M]\{u_j\} = \delta_{ij}, \quad i, j = 1, 2, \dots, n. \quad (2)$$

where $[K]$ and $[M]$ are the structural stiffness and mass matrices, respectively, $\lambda_i = w_i^2$ is the i th eigenvalue, w_i is the i th natural frequency, $\{u_i\}$ is the eigenvector corresponding to λ_i , n is the total degrees of freedom, and δ_{ij} is the Kronecker sign. It is assumed that the eigenvalue is distinct, and Eq. (1) will be referred to as the original eigenvalue problem.

After the structure is modified, the stiffness matrix $[K]$ and the mass matrix $[M]$ are also changed, and the amounts of change are $[\Delta K]$ and $[\Delta M]$, respectively. The corresponding eigenvalue problem of the modified structure is

$$[\bar{K}]\{\bar{u}_i\} = \bar{\lambda}_i[\bar{M}]\{\bar{u}_i\}. \quad (3)$$

$$\{\bar{u}_i\}^T[\bar{M}]\{\bar{u}_j\} = \delta_{ij}, \quad i, j = 1, 2, \dots, n. \quad (4)$$

where

$$[\bar{K}] = [K] + [\Delta K] \quad (5)$$

$$[\bar{M}] = [M] + [\Delta M] \quad (6)$$

Then the i th eigenvalue $\bar{\lambda}_i$ and the corresponding eigenvector $\{\bar{u}_i\}$ can be expressed as

$$\bar{\lambda}_i = \lambda_i + \Delta\lambda_i, \quad \{\bar{u}_i\} = \{u_i\} + \{\Delta u_i\} \quad (6)$$

The First-order Perturbation Method

From Eqs. (6), the standard equations of the first-order perturbation method can be written as

$$\lambda_{1i} = \lambda_i + \Delta\lambda_i, \quad \{u_{1i}\} = \{u_i\} + \{\Delta u_{1i}\} \quad (7)$$

where

$$\Delta\lambda_{1i} = \{u_i\}^T ([\Delta K] - \lambda_i [\Delta M]) \{u_i\}, \quad \{\Delta u_{1i}\} = \sum_{j=1}^n c_{ij}^1 \{u_j\} \quad (8)$$

In the above formula, coefficients c_{ij}^1 are given as follows

$$c_{ij}^1 = \{u_j\}^T ([\Delta K] - \lambda_i [\Delta M]) \{u_i\} / (\lambda_i - \lambda_j), \quad j \neq i \quad (9)$$

$$c_{ii}^1 = -\{u_i\}^T [\Delta M] \{u_i\} / 2, \quad j = i \quad (10)$$

The second-order Perturbation Method

Similarly, the standard equations of the second-order perturbation method can be written as

$$\lambda_{2i} = \lambda_i + \Delta\lambda_{2i}, \quad \{u_{2i}\} = \{u_i\} + \{\Delta u_{2i}\} \quad (11)$$

where

$$\begin{aligned} \Delta\lambda_{2i} = & \{u_i\}^T [\Delta K] \{\Delta u_{1i}\} - \lambda_i \{u_i\}^T [\Delta M] \{\Delta u_{1i}\} \\ & - \Delta\lambda_{1i} \{u_i\}^T [M] \{\Delta u_{1i}\} - \Delta\lambda_{1i} \{u_i\}^T [\Delta M] \{u_i\} \end{aligned} \quad (12)$$

$$\{\Delta u_{2i}\} = \{\Delta u_{1i}\} + \sum_{j=1}^n c_{ij}^2 \{u_j\} \quad (13)$$

In the above formula, coefficients c_{ij}^2 are given as follows

$$c_{ij}^2 = (\{u_j\}^T [\Delta K] \{\Delta u_{1i}\} - \lambda_i \{u_j\}^T [\Delta M] \{\Delta u_{1i}\} - \Delta \lambda_{1i} \{u_j\}^T [M] \{\Delta u_{1i}\} - \Delta \lambda_{1i} \{u_j\}^T [\Delta M] \{u_i\}) / (\lambda_i - \lambda_j), \quad j \neq i. \quad (14)$$

$$c_{ii}^2 = -(\{\Delta u_{1i}\}^T [M] \{\Delta u_{1i}\} + \{u_i\}^T [\Delta M] \{\Delta u_{1i}\} + \{\Delta u_{1i}\}^T [\Delta M] \{u_i\}) / 2 \quad j = i. \quad (15)$$

William B.B. Method Mentioned by William20

Use the inner product of eigenvector with respect to the modified structure mass matrix \bar{M} to replace the inner product with respect to the original structure mass matrix M , we can obtain the William.B.B method.

Similarly, the standard equations of the William.B.B method can be written as

$$\lambda_{3i} = \lambda_i + \Delta \lambda_{3i}, \quad \{u_{3i}\} = \{u_i\} + \{\Delta u_{3i}\}. \quad (16)$$

where

$$\Delta \lambda_{3i} = \{u_i\}^T ([\Delta K] - \lambda_i [\Delta M]) \{u_i\} / (\{u_i\}^T [\bar{M}] \{u_i\}). \quad (17)$$

$$\{\Delta u_{3i}\} = \sum_{j=1}^n c_{ij}^3 \{u_j\}. \quad (18)$$

In the above formula, coefficients c_{ij}^3 are given as follows

$$c_{ij}^3 = \{u_j\}^T ([\Delta K] - \lambda_i [\Delta M]) \{u_i\} / ((\lambda_i - \lambda_j) \{u_i\}^T [\bar{M}] \{u_i\}), \quad j \neq i. \quad (19)$$

$$c_{ii}^3 = -\{u_i\}^T [\Delta M] \{u_i\} / (2 \{u_i\}^T [\bar{M}] \{u_i\}), \quad j = i. \quad (20)$$

The Perturbation Method Combining with the Rayleigh Quotient Mentioned by Chen Et Al.21

Similarly, the standard equations can be written as

$$\lambda_{4i} = \lambda_i + \Delta \lambda_{4i}, \quad \{u_{4i}\} = \{u_i\} + \{\Delta u_{4i}\}. \quad (21)$$

Consider the effects of the high-frequency modes and use the first-order perturbation method results, then have

$$\Delta \lambda_{4i} = \{u_{1i}\}^T ([\bar{K}] - \lambda_i [\bar{M}]) \{u_{1i}\} / (\{u_{1i}\}^T [\bar{M}] \{u_{1i}\}). \quad (22)$$

$$\{\Delta u_{4i}\} = \sum_{j=1}^n c_{ij}^4 \{u_j\} \quad (23)$$

In the above formula, coefficients c_{ij}^4 are given as follows:

$$c_{ij}^4 = \{u_j\}^T ([\Delta K] - \lambda_i [\Delta M] - \Delta \lambda_i [\bar{M}]) \{u_i\} / (\lambda_i - \lambda_j), \quad j \neq i \quad (24)$$

$$c_{ii}^4 = -(\{u_i\}^T [\Delta M] \{u_i\} + \{\Delta u_i\}^T [M] \{\Delta u_i\}) / 2, \quad j = i \quad (25)$$

The Proposed Method

From the previous four methods, we know that we first calculate the eigenvalues of the modified structure, and then calculate the corresponding eigenvectors. The following we first construct an iteration format on the basis of the matrix perturbation theory to compute the natural vibration mode of the modified structure, and then use the Rayleigh quotient to compute the corresponding natural frequency.

The eigenvectors of modified structure are expressed as

$$\{u_{ni}^{k+1}\} = \{u_i\} + \{\Delta u_{ni}^{k+1}\} \quad (26)$$

$$\{\Delta u_{ni}^{k+1}\} = \sum_{j=1}^n c_{ij}^{n(k+1)} \{u_j\} \quad (27)$$

Substituting Eqs. (5)-(6) into Eq. (3) and rearranging them yields

$$\begin{aligned} & -([K] - \lambda_i [M]) \{\Delta u_i\} = \\ & ([\Delta K] - \lambda_i [\Delta M]) (\{u_i\} + \{\Delta u_i\}) - \Delta \lambda_i ([M] + [\Delta M]) (\{u_i\} + \{\Delta u_i\}) \end{aligned} \quad (28)$$

Let $\Delta \lambda_i = \Delta \lambda_{ni}^k$, $\{\Delta u_i\} = \{\Delta u_{ni}^k\}$ on Eq. (28) right-hand side, and we get

$$-([K] - \lambda_i [M]) \{\Delta u_i\} = ([\Delta K] - \lambda_i [\Delta M]) \{u_{ni}^k\} - \Delta \lambda_{ni}^k [\bar{M}] \{u_{ni}^k\} \quad (29)$$

Premultiplying Eq. (29) by $\{u_j\}^T$ results in

$$\begin{aligned} & -\{u_j\}^T ([K] - \lambda_i [M]) \{\Delta u_i\} = \\ & \{u_j\}^T ([\Delta K] - \lambda_i [\Delta M]) \{u_{ni}^k\} - \Delta \lambda_{ni}^k \{u_j\}^T [\bar{M}] \{u_{ni}^k\} \end{aligned} \quad (30)$$

In Eq. (30), let $\{\Delta u_i\} = \{\Delta u_{ni}^{k+1}\}$. Substituting Eq. (27) into Eq. (30) and noting that $\{u_j\}$ is M-normalized, we have the following iterative formula

$$c_{ij}^{n(k+1)} = \{u_j\}^T ([\Delta K] - \lambda_i [\Delta M] - \Delta \lambda_{ni}^k \bar{M}) \{u_{ni}^k\} / (\lambda_i - \lambda_j) \quad j \neq i \quad (31)$$

For the case of $j=i$, the coefficient $c_{ii}^{n(k+1)}$ can be obtained by the \bar{M} -normalized condition

$$(\{u_i\} + \{\Delta u_i\})^T [\bar{M}] (\{u_i\} + \{\Delta u_i\}) = 1 \quad (32)$$

Expanding Eq. (32) and rearranging them yields

$$-2\{u_i\}^T [M] \{\Delta u_i\} = (\{u_i\} + \{\Delta u_i\})^T [\Delta M] (\{u_i\} + \{\Delta u_i\}) + \{\Delta u_i\}^T [M] \{\Delta u_i\} \quad (33)$$

Similarly, noting Eqs. (2) and (27), we can obtain

$$c_{ii}^{n(k+1)} = -(\{u_{ni}^k\}^T [\Delta M] \{u_{ni}^k\} + \{\Delta u_{ni}^k\} [M] \{\Delta u_{ni}^k\}) / 2 \quad (34)$$

Based on the obtained $c_{ii}^{n(k+1)}$ and $c_{ij}^{n(k+1)}$, the eigenvectors $\{u_{ni}^{k+1}\}$ can be obtained from Eqs. (26) and (27).

Then we use the Rayleigh quotient and get the eigenvalues

$$\lambda_{ni}^{k+1} = (\{u_{ni}^{k+1}\}^T [\bar{K}] \{u_{ni}^{k+1}\}) / (\{u_{ni}^{k+1}\}^T [\bar{M}] \{u_{ni}^{k+1}\}) \quad (35)$$

and further get

$$\Delta \lambda_{ni}^{k+1} = \{u_{ni}^{k+1}\}^T ([\bar{K}] - \lambda_i [\bar{M}]) \{u_{ni}^{k+1}\} / (\{u_{ni}^{k+1}\}^T [\bar{M}] \{u_{ni}^{k+1}\}) \quad (36)$$

Throughout the iterative process, we first select the iteration-based initial values $\Delta \lambda_{ni}^0$, $\{\Delta u_{ni}^0\}$ and $\{u_{ni}^0\}$. After obtaining the $\Delta \lambda_{ni}^{k+1}$, $\{\Delta u_{ni}^{k+1}\}$ and $\{u_{ni}^{k+1}\}$, we regard them as the new iteration-based and substitute them into Eqs. (31) and (34), and make the repeated iteration of the loop equation until the results converge to the required accuracy.

Numerical Example

In this section, we will use a numerical example to demonstrate the validity of the proposed method.

We consider the following truss with 8 joints which is shown in Fig. 1. The cross section of all members is $A = 1.2 \times 10^{-3} m^2$, the modulus of elasticity is $E = 2.1 \times 10^{11} Pa$, and the density is $\rho = 7.85 \times 10^3 kg/m^3$. We will adopt the lumped mass in the finite element computation. We assume that the cross section of the bar ① is increased by 20% and 60%, respectively.

To compare the accuracies, we adopt the first-order perturbation method, the second-order perturbation method, the William.B.B method, the perturbation method combining with the Rayleigh quotient and the proposed method respectively to compute the eigenvalues and eigenvectors of the modified structure. In the proposed method, we choose the results of first-order perturbation method as the initial iteration base, and make iteration one time, which makes the cost of computing minimum. For convenience, we denote the above five methods as method 1-method 5 sequentially in the following Figures.

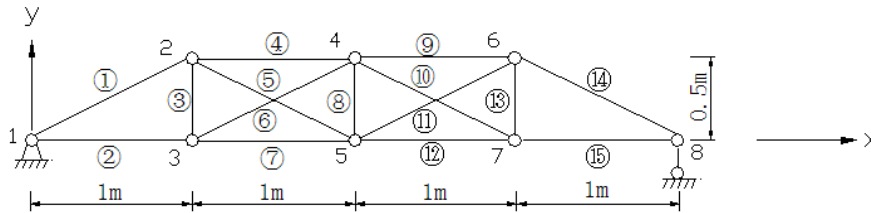


Fig. 1 The truss with 8 joints

The error of eigenvalue is calculated by

$$tr_i = |(\lambda_{Ei} - \lambda_{Ii}) / \lambda_{Ei}| \quad (37)$$

λ_{Ei} is the exact eigenvalue of the modified structure, λ_{Ii} is the approximate eigenvalue computed by the previous several methods. The errors are denoted by tr1, tr2, tr3, tr4 and trn1, respectively.

The error of eigenvector is calculated by

$$ur_i = \text{abs}(\{u_{Ei}\}^T * \{u_{Ii}\} / \text{sqrt}(\{u_{Ei}\}^T * \{u_{Ei}\}) / \text{sqrt}(\{u_{Ii}\}^T * \{u_{Ii}\})) \quad (38)$$

u_{Ei} is the exact eigenvector of the modified structure, u_{Ii} is the approximate eigenvector computed by the previous several methods. The errors are denoted by ur1, ur2, ur3, ur4 and urn1, respectively.

The error results of the first five eigensolutions are listed in Tables 1-4. We use the first-order perturbation solution as the initial iteration base, and make iteration one time in the proposed method, at this time Eqs. (31) and (34) are degenerated into Eqs. (24) and (25), respectively, so Eq. (26) is identical to Eq. (21), as we have seen in Tables 3-4, the values of urn1 and ur4 are the same, i.e. the eigenvector accuracy of the proposed method is consistent with that obtained by the perturbation method combining with the Rayleigh quotient, but more accurate than the other three methods, as we have seen in Figs. 4-5. The eigenvalue accuracy of the proposed method is much higher than the other four methods, as shown in Tables 1-2 and Figs. 2-3. We also know when a large amount of structure modification, such as modification of 60%, as shown in Tables 2 and 4, Figs. 3 and 5, the proposed method also has high accuracy, and the validity of the proposed method is demonstrated.

Table 1 The eigenvalue accuracies comparison (cross section is increased by 20%)

i	tr1	tr2	tr3	tr4	trn1
1	0.0016	0.0003	0.0015	5.57e-5	1.91e-6
2	0.0004	0.0001	0.0004	4.32e-5	3.84e-6
3	0.0040	0.0003	0.0037	1.54e-5	1.37e-6
4	0.0003	4.4e-5	0.0002	3.04e-6	2.96e-8
5	7.81e-5	4.75e-7	1.08e-5	1.93e-7	1.64e-10

Table 2 The eigenvalue accuracies comparison (cross section is increased by 60%)

i	tr1	tr2	tr3	tr4	trn1
1	0.0107	0.0061	0.0102	0.0034	0.0011
2	0.0024	0.0027	0.0026	0.0024	0.0020
3	0.0275	0.0075	0.0257	0.0014	0.0005
4	0.0031	0.0009	0.0022	0.0002	2.01e-5
5	0.0007	6.48e-6	0.0001	1.42e-5	1.08e-7

Table 3 The eigenvector accuracies comparison (cross section is increased by 20%)

i	ur1	ur2	ur3	ur4	urn1
1	0.999998	0.999999	0.999998	1.000000	1.000000
2	0.999964	0.999997	0.999967	0.999997	0.999997
3	0.999961	0.999997	0.999964	0.999997	0.999997
4	0.999993	1.000000	0.999995	1.000000	1.000000
5	1.000000	1.000000	1.000000	1.000000	1.000000

Table 4 The eigenvector accuracies comparison (cross section is increased by 60%)

i	ur1	ur2	ur3	ur4	urn1
1	0.999885	0.999964	0.999899	0.999965	0.999965
2	0.998633	0.998883	0.998781	0.998879	0.998879
3	0.998357	0.998924	0.998543	0.998937	0.998937
4	0.999511	0.999992	0.999646	0.999992	0.999992
5	0.999973	1.000000	0.999991	1.000000	1.000000

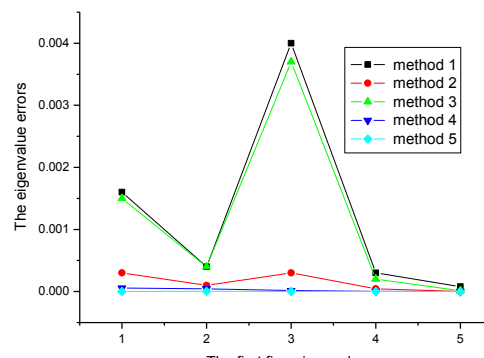


Fig. 2 The eigenvalue accuracies comparison (cross section is increased by 20%)

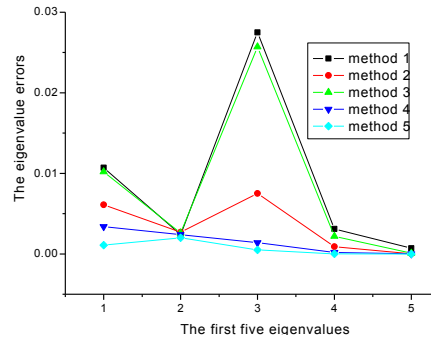


Fig. 3 The eigenvalue accuracies comparison (cross section is increased by 60%)

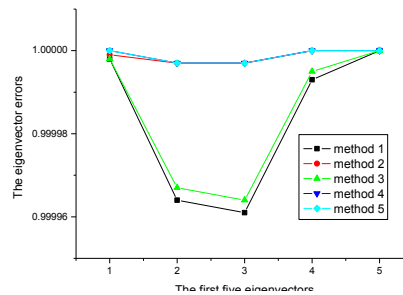


Fig.4 The eigenvector accuracies comparison (cross section is increased by 20%)

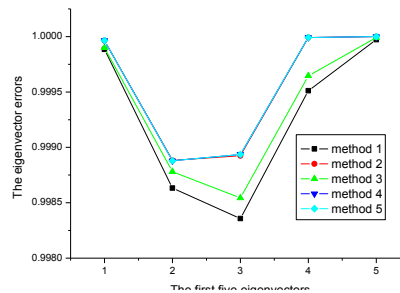


Fig. 5 The eigenvector accuracies comparison (cross section is increased by 60%)

Conclusions

An improved matrix perturbation method for modified structures has been proposed. By constructing an iterative format and combining with the Rayleigh quotient, we can obtain the eigensolution of the modified structure, and the calculation accuracy can be greatly improved. A numerical example has demonstrated the validity of the proposed method.

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