

SOLUTION OF DISCRETE OSCILLATING SYSTEM ON PERSONAL COMPUTER

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SUMMARY. The article is devoted a algorithm for deriving mass matrix, stiffness matrix and damping matrix for oscillating discrete system. The algorithm is common setting equation of motion. This technique enables solving different problems of oscillating system, especially a problem of parameters optimization, by numerical methods. Comparison of different methods realized on personal computers was done.

Oscillating systems find many applications in a technical life. In this article some results of researching on computers are discussed. Most of techniques mentioned below were realized on 286X and 386X based personal computer and FORTRAN-77 is chosen as a programming language. Because of simplicity we will consider only discrete systems. It means that system could consist of elements of 3 types: mass, spring and damper and these features of system are concentrated. But the scheme is arbitrary: free or with a frame, simple or branched. Some of them are shown on the Fig. 1

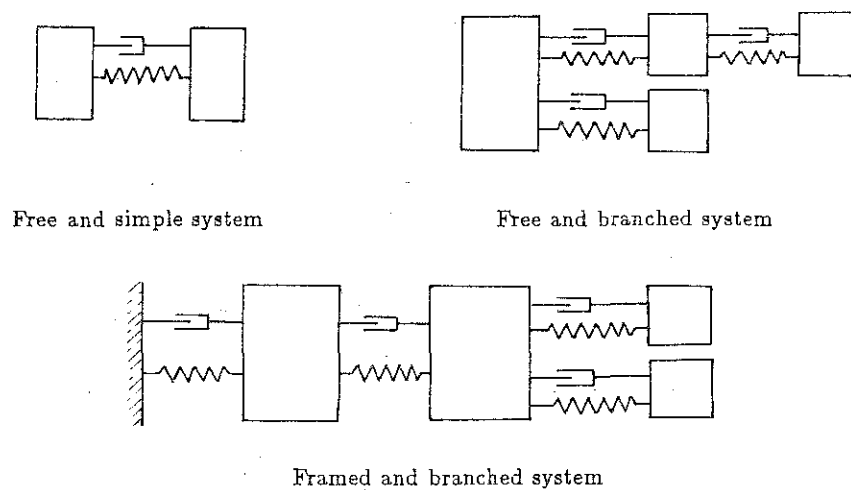


Fig. 1. Examples of discrete oscillating system

Just now we will work in a different way than normal methods when solving "by hand". We will give one common algorithm of setting equations of motion for all types of oscillating system. Let's assume that forces will affect masses which are concentrated, rigid and unsprung. Springs and dampers are massies and dampers are viscous e. g. derived force is proportional to a speed. For such a system with n degrees of freedom we have a well-known system of equations of motion:

$$\lambda \underline{M} \underline{C} = \underline{K} \underline{C} \quad (4)$$

where λ is a eigenvalue, \underline{C} is eigenvector.

Because of a symmetricity of matrix \underline{K} a QL method for a tridiagonal matrix is suitable [3]. So in advance we must transform equation (4) into standard form by using Choleski's separation:

$$\underline{M} = \underline{L} \underline{L}^T \quad (5)$$

As the mass matrix \underline{M} is diagonal, \underline{L} is diagonal again and $\underline{L} = \underline{L}^T$, so we obtain a standard equation:

$$(\underline{L}^{-1} \underline{K} \underline{L}^{-1}) (\underline{L} \underline{C}) = \lambda (\underline{L} \underline{C}) \quad (6)$$

A riding matrix $(\underline{L}^{-1} \underline{K} \underline{L}^{-1})$ is symmetric again and we can use the mentioned QL method.

2) Time-History of vibration

When damping of the system is general we should transform equation (1) into the form:

$$- \begin{bmatrix} -\underline{K} & \underline{O} \\ \underline{O} & \underline{M} \end{bmatrix} \cdot \begin{bmatrix} \underline{X} \\ \dot{\underline{X}} \end{bmatrix} + \begin{bmatrix} \underline{B} & \underline{M} \\ \underline{M} & \underline{O} \end{bmatrix} \cdot \begin{bmatrix} \dot{\underline{X}} \\ \ddot{\underline{X}} \end{bmatrix} = \begin{bmatrix} \underline{f} \\ \underline{O} \end{bmatrix} \quad (7)$$

or:

$$-\underline{P} \underline{u} + \underline{N} \dot{\underline{u}} = \underline{g}$$

where

$$\underline{P} = \begin{bmatrix} -\underline{K} & \underline{O} \\ \underline{O} & \underline{M} \end{bmatrix}; \quad \underline{u} = \begin{bmatrix} \underline{X} \\ \dot{\underline{X}} \end{bmatrix}; \quad \underline{N} = \begin{bmatrix} \underline{B} & \underline{M} \\ \underline{M} & \underline{O} \end{bmatrix}; \quad \underline{g} = \begin{bmatrix} \underline{f} \\ \underline{O} \end{bmatrix}$$

The dimension of the problem is two-times greater but \underline{F} and \underline{N} are both symmetric. A well-known method of Runge-Kutta could be used to solve equations [7].

In the cases where we could suppose the condition of proportional damping some methods of numerical intergration are better. For example when we have Rayleigh's damping:

$$\underline{B} = a_0 \underline{M} + a_1 \underline{K} \quad (8)$$

two methods of Newmark and Wilson are very elegant and convenient as a proportionality of \underline{B} enables one useful transform of matrix \underline{K} .

Another way is using main coordinates. This technique leads to a system of n single differential equations which could be solved separately by Duhamell's integral.

Two points of this domain should be treated carefully on PC: a problem of a memory and time-step for numerical integration. A size of required memory depends on used language and compiler but perhaps the size of a data segment exceeds 64 Kbyte that will complicate a storage and handling data and influence time consuming. As a time-step we recommend to choose a value:

$$\Delta t < 1/10 f_{\max} \quad (9)$$

where f_{\max} is the greatest frequency of a system

$$f = \sqrt{\lambda}/2\pi.$$

A greater step leads to accumulation of numerical errors and they can affect like a undesirable damping.

3) Optimization

Higher stage of solution is optimization of parameters (mass, stiffness and damping). Author of this article have tested a lot of methods of optimization on computers and the comparison of these methods was done in the frequency domain. We can set a following task:

Riding function

$$F = \sum_i w_i (\lambda_i - \lambda_{ib})^2 \rightarrow \min$$

where λ_{ib} $i = 1, n$ is desirable set of eigenvalues,

λ_i are eigenvalues of the system.

A similar task could be set in the amplitude domain but the form of riding function should be chosen carefully. Note that we are speaking about numerical optimization, it means that at each stage of a optimization process the system is better if it gives a smaller value of the riding function. Also, we consider only parameter optimization.

The tests on computers showed that following methods are convenient for our task:

- Conjugate gradient method (Fletcher-Reevse's method)
- Fletcher-Powell's method
- Rosenbrock's method.

The best of them is conjugate gradient method for its precision and time-consuming.

Other methods: Box's algorithm or Glass-Copper's method were tested but the results were not so good.

One difficult can appear in a optimization process: Values of stiffness of springs are greater a lot than values of masses. This leads to a "zigzag" approaching and doesn't allow to reach precise result. Some special technique of optimization should be applied in order to gain better results [2].

A simple example:

We consider an example to illustrate a problem. A torsional system in Fig. 2 has 5 inertia bodies and 6 springs. The values of parameters are following:

$m_1 = 5 \text{ kgm}^2$	$m_2 = 3 \text{ kgm}^2$	$m_3 = 5 \text{ kgm}^2$
$m_4 = 6 \text{ kgm}^2$	$m_5 = 2 \text{ kgm}^2$	
$c_1 = 10^5 \text{ Nm/rad}$	$C_2 = 2 \cdot 10^5 \text{ Nm/rad}$	$c_3 = 4 \cdot 10^5 \text{ Nm/rad}$
$C_4 = 7 \cdot 10^5 \text{ Nm/rad}$	$C_5 = 2 \cdot 10^5 \text{ Nm/rad}$	$C_6 = 10^5 \text{ Nm/rad}$

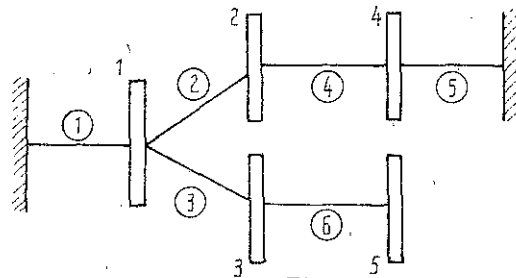


Fig. 2

The problem of eigenvalues is solved easily. The system has a spectrum:

$$104.2 \quad 201.1 \quad 269.3 \quad 449.9 \quad 643.2 \quad \text{rad/s}$$

The most difficult task is a problem of optimization. Let's suppose that retained spectrum is not convenient and a constructor want to achieve better spectrum, say for example

$$90 \quad 250 \quad 300 \quad 450 \quad 650 \quad \text{rad/s.}$$

Not all of parameters could be changed but only 1-st, 3-rd and 4-th bodies and 1-st, 5-th and 6-th springs.

In most cases the task of optimization require a lot of CPU time. For illustration we give in the Tab. 1 some value of solution on different 286X and 386X based PC. Rosenbrock's method is used and gives the result:

$$\begin{array}{lll}
 m_1 = 4.056 & m_3 = 6.584 & m_4 = 6.329 \text{ kgm}^2 \\
 C_1 = 1.063 \cdot 10^4 & C_5 = 4.832 \cdot 10^5 & C_6 = 1.34 \cdot 10^5 \text{ Nm/rad}
 \end{array}$$

Tab. 1

PC	Math. Coprocessor	CPU time	
Fujitkma 286,	8Mhz	none	11 min 32 sec
Universal Computer 286,	8Mhz	none	22 min 52 sec
Winn 286,	16Mhz	none	15 min 35 sec
Winn 286,	8Mhz	none	23 min 47 sec
Olivetti 286,	16Mhz	present	2 min 32 sec
Unitron 386,	32Mhz	none	5 min 52 sec

CONCLUSION

The algorithm for deriving mass, stiffness and damping matrix enables automatic solving the problem on personal computers. But experience showed that some difficulties could appear and we should be careful when using different methods because only some of them are convenient for our purpose.

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GIẢI HỆ DAO ĐỘNG RỜI RẠC TRÊN MÁY VI TÍNH

Bài báo đưa ra một thuật toán để xây dựng các ma trận độ cứng, ma trận hệ số cản và ma trận khối lượng đối với hệ dao động rời rạc. Thuật toán này là tổng quát cho tất cả các dạng khác nhau của hệ dao động và tạo cơ sở cho việc lập phương trình dao động trực tiếp trên máy tính. Phương pháp này cho phép giải quyết nhiều bài toán khác nhau bằng các phương pháp số, đặc biệt là bài toán tối ưu tham số. Các phương pháp khác nhau cũng được so sánh khi được sử dụng trên máy tính.