AN INTRODUCTION TO DYNAMICS OF STRUCTURES

Giacomo Boffi

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http://intranet.dica.polimi.it/people/boffi-giacomo
Dipartimento di Ingegneria Civile Ambientale e Territoriale

Politecnico di Milano

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INTRODUCTION

Definitions

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Dynamics of Structures all of the above, applied to a *structural system*, i.e., a system designed to stay in equilibrium

Dynamics of Structures

Our aim is to determine the stresses and deflections that a *dynamic loading* induces in a *structure* that remains in the neighborhood of a point of equilibrium.

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If we restrict ourselves to analysis of linear systems*, it is so convenient to use the principle of superposition to study the combined effects of static and dynamic loadings that different methods, of different character, are applied to these different loadings.*

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Our focus will be on *deterministic analysis*

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Dealing with deterministic dynamic loadings we will study, in order of complexity,

Harmonic Loadings a force is modulated by a harmonic function of time, characterized by a frequency ω and a phase φ:
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- the loading can be described in terms of an analytical function, e.g., $p(t) = p_o \exp(\alpha t)$
- the loading is measured experimentally, hence it is known only in a discrete set of instants; in this case, we say that we know the loading *time-history*.

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In many cases it is however possible to simplify the formulation of the structural dynamic problem to *ordinary* differential equations.

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Under this assumption, the analytical problem is greatly simplified:

1. the inertial forces are applied only at the lumped masses,
2. the only deflections that influence the inertial forces are the deflections of the lumped masses,
3. using methods of static analysis we can determine those deflections,

thus consenting the formulation of the problem in terms of a set of ordinary differential equations, one for each relevant component of the inertial forces.

Dynamic Degrees of Freedom

The *dynamic degrees of freedom* (DDOF) in a lumped mass, discretized system are the displacements components of the lumped masses associated with the relevant components of the inertial forces.

Dynamic Degrees of Freedom

If a lumped mass can be regarded as a *point* mass then at most 3 translational DDOFs will suffice to represent the associated inertial force.

On the contrary, if a lumped mass has a discrete volume its inertial force depends also on its rotations (inertial couples) and we need at most 6 DDOFs to represent the mass deflections and the inertial force.

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Of course, a continuous system has an infinite number of degrees of freedom.

Generalized Displacements

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When the masses are distributed we can simplify our problem expressing the deflections in terms of a linear combination of assigned functions of position, the coefficients of the linear combination being the *generalized coordinates* (e..g., the deflections of a rectilinear beam can be expressed in terms of a trigonometric series).

Generalized Displacements, cont.

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> *Even if the method of generalized coordinates has its beauty, we must recognize that for each different problem we have to derive an* ad hoc *formulation, with an evident loss of generality.*

Finite Element Method

The *finite elements method* (FEM) combines aspects of lumped mass and generalized coordinates methods, providing a simple and reliable method of analysis, that can be easily programmed on a digital computer.

Finite Element Method

- In the FEM, the structure is subdivided in a number of non-overlapping pieces, called the *finite elements*, delimited by common *nodes*.
- The FEM uses *piecewise approximations* (i.e., local to each element) to the field of displacements.
- In each *element* the displacement field is derived from the displacements of the *nodes* that surround each particular element, using *interpolating functions*.
- The displacement, deformation and stress fields in each element, as well as the inertial forces, can thus be expressed in terms of the unknown *nodal displacements*.
- The *nodal displacements* are the dynamical DOFs of the FEM model.

Finite Element Method

Some of the most prominent advantages of the FEM method are

1. The desired level of approximation can be achieved by further subdividing the structure.

2. The resulting equations are only loosely coupled, leading to an easier computer solution.

3. For a particular type of finite element (e.g., beam, solid, etc) the procedure to derive the displacement field and the element characteristics does not depend on the particular geometry of the elements, and can easily be implemented in a computer program.

Formulation of the equations of motion

In a deterministic dynamic analysis, given a prescribed load, we want to evaluate the displacements in each instant of time.

In many cases a limited number of DDOFs gives a sufficient accuracy; further, the dynamic problem can be reduced to the determination of the time-histories of some selected component of the response.

Writing the equation of motion

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The solution of the EOM gives the requested displacements.

The formulation of the EOM is the most important, often the most difficult part of a dynamic analysis.

We have a choice of techniques to help us in writing the EOM, namely:

- the D'Alembert Principle,
- the Principle of Virtual Displacements,
- the Variational Approach.

D'Alembert Principle

By Newton's II law of motion, for any particle the rate of change of momentum is equal to the external force,

$$\vec{p}(t) = \frac{d}{dt}(m\frac{d\vec{u}}{dt}),$$

where $\vec{u}(t)$ is the particle displacement.

In structural dynamics, we may regard the mass as a constant, and thus write

$$\vec{p}(t) = m\ddot{\vec{u}},$$

where each operation of differentiation with respect to time is denoted with a dot.

When we interpret the term $-m\ddot{\vec{u}}$ as the *inertial force* that contrasts the acceleration of the particle, we can write an equation of equilibrium for the particle, $\vec{p}(t) = m\ddot{\vec{u}}$.

D'Alembert principle, cont.

The concept that a mass develops an inertial force opposing its acceleration is known as the D'Alembert principle, and using this principle we can write the *EOM* as a simple equation of equilibrium.

The term $\vec{p}(t)$ must comprise each different force acting on the particle, including the reactions of kinematic or elastic constraints, internal forces and external, autonomous forces.

In many simple problems, D'Alembert principle is the most direct and convenient method for the formulation of the *EOM*.

Principle of virtual displacements

In a reasonably complex dynamic system, with e.g. articulated rigid bodies and external/internal constraints, the direct formulation of the *EOM* using D'Alembert principle may result difficult.

In these cases, application of the *Principle of Virtual Displacements* is very convenient, because the reactive forces do not enter the equations of motion, that are directly written in terms of the motions compatible with the restraints/constraints of the system.

Principle of Virtual Displacements, cont.

Considering, e.g., an assemblage of rigid bodies, the *pvd* states that necessary and sufficient condition for equilibrium is that, for every *virtual displacement* (i.e., any infinitesimal displacement compatible with the restraints) the total work done by all the external forces is zero.

For an assemblage of rigid bodies, writing the *EOM* requires

1. to identify all the external forces, comprising the inertial forces, and to express their values in terms of the *ddof*;
2. to compute the work done by these forces for different virtual displacements, one for each *ddof*;
3. to equate to zero all these work expressions.

The *pvd* is particularly convenient also because we have only scalar equations, even if the forces and displacements are of vectorial nature.

Variational approach

Variational approaches do not consider directly the forces acting on the dynamic system, but are concerned with the variations of kinetic and potential energy and lead, as well as the *pvd*, to a set of scalar equations.

For example, the equation of motion of a generical system can be derived in terms of the *Lagrangian function*, $\mathcal{L} = T - V$ where *T* and *V* are, respectively, the kinetic and the potential energy of the system expressed in terms of a vector $\vec{q}$ of indipendent coordinates

$$\frac{d}{dt}\left(\frac{\partial \mathcal{L}}{\partial \dot{q}_i}\right) = \frac{\partial \mathcal{L}}{\partial q_i}, \qquad i = 1, \ldots, N.$$

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The method to be used in a particular problem is mainly a matter of convenience (and, to some extent, of personal taste).

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If our system has a *constant* mass m and it's subjected to a generical, non-linear, internal force $F = F(y, \dot{y})$, where y is the displacement and $\dot{y}$ the velocity of the particle, the equation of motion is

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It is difficult to integrate the above equation in the general case, but it's easy when the motion occurs in a small neighborhood of the equilibrium position.

1 DOF System, cont.

In a position of equilibrium, $y_{eq.}$, the velocity and the acceleration are zero, and hence $f(y_{eq.}, 0) = 0$.

The force can be linearized in a neighborhood of $y_{eq.}, 0$:

$$f(y, \dot{y}) = f(y_{eq.}, 0) + \frac{\partial f}{\partial y}(y - y_{eq.}) + \frac{\partial f}{\partial \dot{y}}(\dot{y} - 0) + O(y, \dot{y}).$$

Assuming that $O(y, \dot{y})$ is small in a neighborhood of $y_{eq.}$, we can write the equation of motion

$$\ddot{x} + a\dot{x} + bx = 0$$

where $x = y - y_{eq.}$, $a = - \left. \frac{\partial f}{\partial \dot{y}} \right|_{y=y_{eq.}, \dot{y}=0}$ and $b = - \left. \frac{\partial f}{\partial y} \right|_{y=y_{eq.}, \dot{y}=0}$.

In an infinitesimal neighborhood of $y_{eq.}$ ($y_{eq.}$ being a position of static equilibrium), the equation of motion can be studied in terms of a linear differential equation of second order with constant coefficients.

1 DOF System, cont.

A linear constant coefficient differential equation has the homogeneous integral $x = A \exp(st)$, that substituted in the equation of motion gives

$$s^2 + as + b = 0$$

whose solutions are

$$s_{1,2} = -\frac{a}{2} \mp \sqrt{\frac{a^2}{4} - b}.$$

The general integral is hence

$$x(t) = A_1 \exp(s_1 t) + A_2 \exp(s_2 t).$$

Given that for a free vibration problem A_1, A_2 are given by the initial conditions, the nature of the solution depends on the sign of the real part of s_1, s_2, because $s_i = r_i + \imath q_i$ and

$$\exp(s_i t) = \exp(\imath q_i t)\exp(r_i t).$$

If one of the $r_i > 0$, the response grows infinitely over time, even for an infinitesimal perturbation of the equilibrium, so that in this case we have an *unstable equilibrium*.

If both $r_i < 0$, the response decrease over time, so we have a *stable equilibrium*.

Finally, if both $r_i = 0$ the roots s_i are purely imaginary and the response is harmonic with constant amplitude.

1 DOF System, cont.

The roots being

$$s_{1,2} = -\frac{a}{2} \mp \sqrt{\frac{a^2}{4} - b},$$

- if $a > 0$ and $b > 0$ both roots are negative or complex conjugate with negative real part, the system is asymptotically stable,
- if $a = 0$ and $b > 0$, the roots are purely imaginary, the equilibrium is indifferent, the oscillations are harmonic,
- if $a < 0$ or $b < 0$ at least one of the roots has a positive real part, and the system is unstable.

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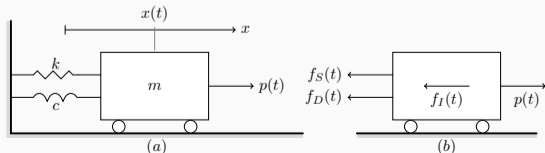
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EQUATION OF MOTION OF THE BASIC DYNAMIC SYSTEM
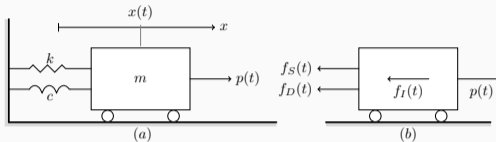
The equation of motion, written using the D'Alembert Principle, requires the equilibrium of all the forces acting on the mass *including the inertial one*.

The forces are the *external force*, $p(t)$, positive in the direction of motion and the *resisting forces*, i.e., the inertial force $f_I(t)$, the damping force $f_D(t)$ and the elastic force, $f_S(t)$, that are opposite to the direction of the acceleration, velocity and displacement.

The equation of motion, merely expressing the equilibrium of these forces, writing the resisting forces and the external force across the equal sign

$$f_I(t) + f_D(t) + f_S(t) = p(t)$$

EOM of the basic dynamic system, cont.

According to D'Alembert principle, the inertial force is the product of the mass and acceleration

$$f_I(t) = m\,\ddot{x}(t).$$

Assuming a viscous damping mechanism, the damping force is the product of the damping constant c and the velocity,

$$f_D(t) = c\,\dot{x}(t).$$

Finally, the elastic force is the product of the elastic stiffness k and the displacement,

$$f_S(t) = k\,x(t).$$

EOM OF THE BASIC DYNAMIC SYSTEM, CONT.

- The differential equation of dynamic equilibrium

$$f_I(t) + f_D(t) + f_S(t) = m\,\ddot{x}(t) + c\,\dot{x}(t) + k\,x(t) = p(t).$$

- The *resisting forces* in the EoM

$$f_I(t), f_D(t), f_S(t)$$

  are proportional to the deflection $x(t)$ or one of its time derivatives, $\dot{x}(t)$, $\ddot{x}(t)$.

- The equation of motion is a linear differential equation of the second order, with constant coefficients.

The resisting forces are, by convention, positive when opposite to the direction of motion, i.e., resisting the motion and, by convention as well, are written in the left member of the EoM.
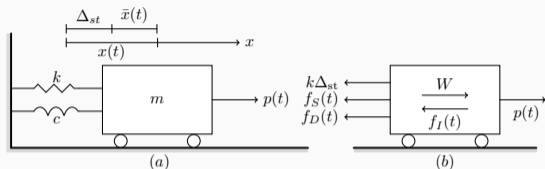
Influence of static forces

(a)   (b)
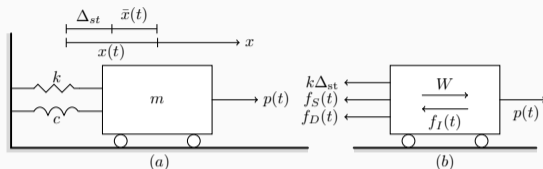
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INFLUENCE OF STATIC FORCES

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Expressing the displacement as the sum of a constant, static displacement and a dynamic displacement,

$$x(t) = \Delta_{st} + \bar{x}(t),$$

and substituting in the *EOM* we have

$$m\,\ddot{x}(t) + c\,\dot{x}(t) + k\,\Delta_{st} + k\,\bar{x}(t) = p(t) + W.$$

Recognizing that $k \Delta_{st} = W$ (so that the two terms, on opposite sides of the equal sign, cancel each other), that $\dot{x} \equiv \dot{\bar{x}}$ and that $\ddot{x} \equiv \ddot{\bar{x}}$ the *EOM* can be written as

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The equation of motion expressed with reference to the static equilibrium position is not affected by static forces.

For this reasons, all displacements in further discussions will be referenced from the equilibrium position and denoted, for simplicity, with $x(t)$.

Recognizing that $k\,\Delta_{st} = W$ (so that the two terms, on opposite sides of the equal sign, cancel each other), that $\dot{x} \equiv \dot{\bar{x}}$ and that $\ddot{x} \equiv \ddot{\bar{x}}$ the *EOM* can be written as

$$m\,\ddot{\bar{x}}(t) + c\,\dot{\bar{x}}(t) + k\,\bar{x}(t) = p(t).$$

The equation of motion expressed with reference to the static equilibrium position is not affected by static forces.

For this reasons, all displacements in further discussions will be referenced from the equilibrium position and denoted, for simplicity, with $x(t)$.

Note that the total *displacements, stresses. etc.* are influenced *by the static forces, and* must *be computed using the superposition of effects.*
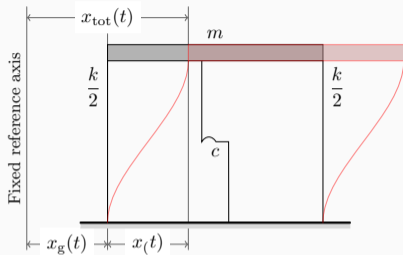
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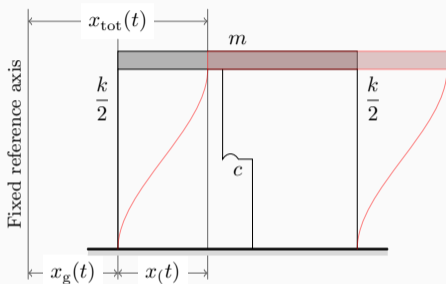
Displacements, deformations and stresses in a structure are induced also by a motion of its support.

Important examples of support motion are the motion of a building foundation due to earthquake and the motion of the base a piece of equipment due to vibrations of the building in which it is housed.

Considering a support motion $x_{\mathrm{g}}(t)$, defined with respect to a inertial frame of reference, the total displacement is

$$x_{\mathrm{tot}}(t) = x_{\mathrm{g}}(t) + x(t)$$

and the total acceleration is

$$\ddot{x}_{\mathrm{tot}}(t) = \ddot{x}_{\mathrm{g}}(t) + \ddot{x}(t).$$
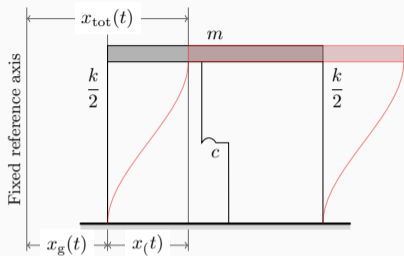
Influence of support motion, cont.

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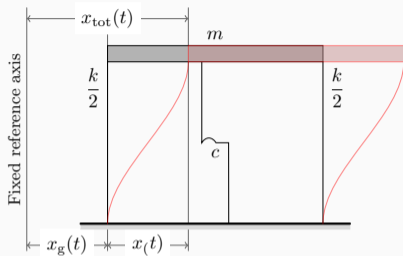
$$x_{tot}(t) = x_g(t) + x(t)$$

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$$\ddot{x}_{tot}(t) = \ddot{x}_g(t) + \ddot{x}(t).$$

While the elastic and damping forces are still proportional to *relative* displacements and velocities, the inertial force is proportional to the total acceleration,

$$f_I(t) = -m\ddot{x}_{tot}(t) = m\ddot{x}_g(t) + m\ddot{x}(t).$$

Influence of support motion, cont.

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Writing the *EOM* for a null external load, $p(t) = 0$, is hence

$$m\,\ddot{x}_{\mathrm{tot}}(t) + c\,\dot{x}(t) + k\,x(t) = 0, \qquad\qquad \text{or,}$$

$$m\,\ddot{x}(t) + c\,\dot{x}(t) + k\,x(t) = -m\,\ddot{x}_{\mathrm{g}}(t) \equiv p_{\mathrm{eff}}(t).$$

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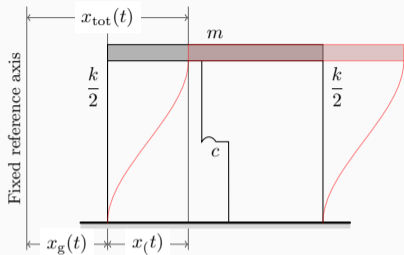$$f_{\mathrm{I}}(t) = -m\ddot{x}_{\mathrm{tot}}(t) = m\ddot{x}_{\mathrm{g}}(t) + m\ddot{x}(t).$$

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Support motion is sufficient to excite a dynamic system: $p_{\mathrm{eff}}(t) = -m\,\ddot{x}_{\mathrm{g}}(t)$.

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Free vibrations of a SDOF system

Free Vibrations

The equation of motion,

$$m\ddot{x}(t) + c\dot{x}(t) + k x(t) = p(t)$$

is a linear differential equation of the second order, with constant coefficients.

Its solution can be expressed in terms of a superposition of a *particular solution*, depending on $p(t)$, and a *free vibration* solution, that is the solution of the so called *homogeneous problem*, where $p(t) = 0$.

In the following, we will study the solution of the homogeneous problem, the so-called *homogeneous* or *complementary* solution, that is the *free vibrations* of the SDOF after a perturbation of the position of equilibrium.

Free vibrations of an undamped system

An undamped system, where $c = 0$ and no energy dissipation takes place, is just an ideal notion, as it would be a realization of *motus perpetuum*. Nevertheless, it is an useful idealization. In this case, the homogeneous equation of motion is

$$m\ddot{x}(t) + k x(t) = 0$$

which solution is of the form $\exp st$; substituting this solution in the above equation we have

$$(k + s^2 m) \exp st = 0$$

noting that $\exp st \neq 0$, we finally have

$$(k + s^2 m) = 0 \Rightarrow s = \pm\sqrt{-\frac{k}{m}}$$

As m and k are positive quantities, s must be purely imaginary.

UNDAMPED FREE VIBRATIONS

Introducing the *natural* circular frequency ω_n

$$\omega_n^2 = \frac{k}{m},$$

the solution of the algebraic equation in s is

$$s = \pm\sqrt{-\frac{k}{m}} = \pm\sqrt{-1}\sqrt{\frac{k}{m}} = \pm i\sqrt{\omega_n^2} = \pm i\omega_n$$

where $i = \sqrt{-1}$ and the general integral of the homogeneous equation is

$$x(t) = G_1 \exp(i\omega_n t) + G_2 \exp(-i\omega_n t).$$

Undamped Free Vibrations

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The solution has an imaginary part?

Undamped Free Vibrations

The solution is derived from the general integral imposing the (real) initial conditions

$$x(0) = x_0, \qquad \dot{x}(0) = \dot{x}_0 \tag{1}$$

Evaluating $x(t)$ for $t = 0$ and substituting in (1), we have

$$\begin{cases} G_1 + G_2 = x_0 \\ i\omega_n G_1 - i\omega_n G_2 = \dot{x}_0 \end{cases}$$

Solving the linear system we have

$$G_1 = \frac{ix_0 + \dot{x}_0/\omega_n}{2i}, \qquad G_2 = \frac{ix_0 - \dot{x}_0/\omega_n}{2i},$$

substituting these values in the general solution and collecting x_0 and $\dot{x}_0$, we finally find ...

Undamped Free Vibrations

$$x(t) = \frac{\exp(i\omega_n t) + \exp(-i\omega_n t)}{2} x_0 + \frac{\exp(i\omega_n t) - \exp(-i\omega_n t)}{2i} \frac{\dot{x}_0}{\omega_n}.$$

Undamped Free Vibrations

$$x(t) = \frac{\exp(i\omega_n t) + \exp(-i\omega_n t)}{2}\, x_0 + \frac{\exp(i\omega_n t) - \exp(-i\omega_n t)}{2i}\, \frac{\dot{x}_0}{\omega_n}.$$

The exponentials with imaginary argument and the trigonometric functions are connected by the Euler formulas,

$$\cos\omega_n t = \frac{\exp(i\omega_n t) + \exp(-i\omega_n t)}{2}, \qquad \sin\omega_n t = \frac{\exp(i\omega_n t) - \exp(-i\omega_n t)}{2i},$$

and the solution can be rewritten in terms of the elementary trigonometric functions

$$x(t) = x_0\,\cos(\omega_n t) + (\dot{x}_o/\omega_n)\,\sin(\omega_n t).$$

Undamped Free Vibrations

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For every real initial conditions we can use indifferently one of the above representations. We usually choose to represent the homogeneous solution as a linear combination of sine and cosine of circular frequency ω_n

Undamped Free Vibrations

Our preferred representation of the general integral of undamped free vibrations is

$$x(t) = A\cos(\omega_n t) + B\sin(\omega_n t)$$

For the usual initial conditions, we have already seen that

$$A = x_0, \qquad B = \frac{\dot{x}_0}{\omega_n}.$$
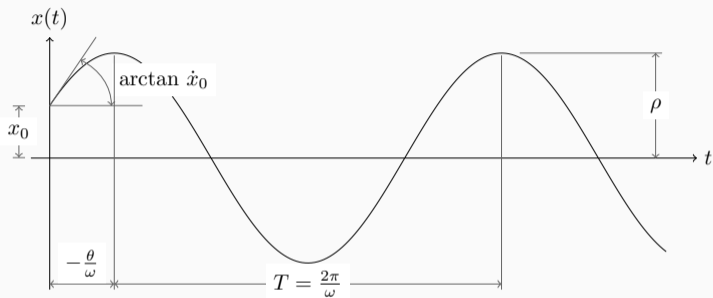
Undamped Free Vibrations

Sometimes we prefer to write $x(t)$ as a single harmonic, introducing a *phase difference* ϕ so that *the amplitude of the motion, C,* is put in evidence:

$$x(t) = A \cos \omega_n t + B \sin \omega_n t$$
$$= C \cos(\omega_n t - \varphi) = C \left(\cos \omega_n t \cos \varphi + \sin \omega_n t \sin \varphi \right)$$

From $A = C \cos \varphi$ and $B = C \sin \varphi$ we have $\tan \varphi = B/A$, from $A^2 + B^2 = C^2(\cos^2 \varphi + \sin^2 \varphi)$ we have $C = \sqrt{A^2 + B^2}$ and eventually

$$x(t) = C \cos(\omega_n t - \varphi), \quad \text{with} \begin{cases} C = \sqrt{A^2 + B^2} \\ \varphi = \arctan(B/A) \end{cases}$$

It is worth noting that the coefficients *A*, *B* and *C* have the dimension of a length, the coefficient ω_n has the dimension of the reciprocal of time and that the coefficient φ is an angle, or in other terms is adimensional.

Free vibrations of a damped system

The viscous damping modifies the response of a *sdof* system introducing a decay in the amplitude of the response. Depending on the amount of damping, the response can be oscillatory or not. The amount of damping that separates the two behaviors is denoted as *critical damping*.

THE SOLUTION OF THE EOM

The equation of motion for a free vibrating damped system is

$$m\ddot{x}(t) + c\dot{x}(t) + kx(t) = 0,$$

substituting the solution $\exp st$ in the preceding equation and simplifying, we have that the parameter s must satisfy the equation

$$ms^2 + cs + k = 0$$

or, after dividing both members by m,

$$s^2 + \frac{c}{m}s + \omega_n^2 = 0$$

whose solutions are

$$s = -\frac{c}{2m} \mp \sqrt{\left(\frac{c}{2m}\right)^2 - \omega_n^2} = \omega_n \left(-\frac{c}{2m\omega_n} \mp \sqrt{\left(\frac{c}{2m\omega_n}\right)^2 - 1}\right).$$

CRITICAL DAMPING

The behavior of the solution of the free vibration problem depends of course on the sign of the radicand $\Delta = \left(\frac{c}{2m\omega_n}\right)^2 - 1$:

$\Delta < 0$ the roots s are complex conjugate,

$\Delta = 0$ the roots are identical, double root,

$\Delta > 0$ the roots are real.

The value of c that make the radicand equal to zero is known as the *critical damping*,

$$c_{\mathrm{cr}} = 2m\omega_n = 2\sqrt{mk}.$$

Critical Damping

A single degree of freedom system is denoted as *critically damped*, *under-critically damped* or *over-critically damped* depending on the value of the damping coefficient with respect to the critical damping.

CRITICAL DAMPING

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Typical building structures are undercritically damped.

Damping Ratio

If we introduce the ratio of the damping to the critical damping, or *critical damping ratio* ζ,

$$\zeta = \frac{c}{c_{\mathrm{cr}}} = \frac{c}{2m\omega_n}, \quad c = \zeta c_{\mathrm{cr}} = 2\zeta\omega_n m$$

the equation of free vibrations can be rewritten as

$$\ddot{x}(t) + 2\zeta\omega_n\dot{x}(t) + \omega_n^2 x(t) = 0$$

and the roots $s_{1,2}$ can be rewritten as

$$s = -\zeta\omega_n \mp \omega_n\sqrt{\zeta^2 - 1}.$$

Free Vibrations of Under-critically Damped Systems

We start studying the free vibration response of under-critically damped SDOF, as this is the most important case in structural dynamics.

The determinant being negative, the roots $s_{1,2}$ are

$$s = -\zeta\omega_n \mp \omega_n\sqrt{-1}\sqrt{1 - \zeta^2} = -\zeta\omega_n \mp i\omega_D$$

with the position that

$$\omega_D = \omega_n\sqrt{1 - \zeta^2}.$$

is the *damped frequency*; the general integral of the equation of motion is, collecting the terms in $\exp(-\zeta\omega_n t)$

$$x(t) = \exp(-\zeta\omega_n t)\left[G_1\exp(-i\omega_D t) + G_2\exp(+i\omega_D t)\right]$$

Initial Conditions

By imposing the initial conditions, $u(0) = u_0$, $\dot{u}(0) = v_0$, after a bit of algebra we can write the equation of motion for the given initial conditions, namely

$$x(t) = \exp(-\zeta\omega_n t) \left[\frac{\exp(i\omega_D t) + \exp(-i\omega_D t)}{2} u_0 + \frac{\exp(i\omega_D t) - \exp(-i\omega_D t)}{2i} \frac{v_0 + \zeta\omega_n u_0}{\omega_D} \right].$$
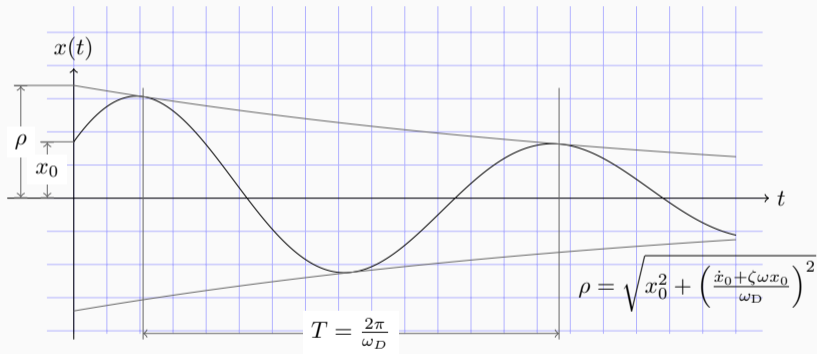
Using the Euler formulas, we finally have the preferred format of the general integral:

$$x(t) = \exp(-\zeta\omega_n t) \left[A\cos(\omega_D t) + B\sin(\omega_D t) \right]$$

with

$$A = u_0, \qquad B = \frac{v_0 + \zeta\omega_n u_0}{\omega_D}.$$

The Damped Free Response

$$\rho = \sqrt{x_0^2 + \left(\frac{\dot{x}_0 + \zeta\omega x_0}{\omega_D}\right)^2}$$

$$T = \frac{2\pi}{\omega_D}$$

Critically damped SDOF

In this case, $\zeta = 1$ and $s_{1,2} = -\omega_n$, so that the general integral must be written in the form

$$x(t) = \exp(-\omega_n t)(A + Bt).$$

The solution for given initial condition is

$$x(t) = \exp(-\omega_n t)(u_0 + (v_0 + \omega_n u_0)t),$$

note that, if $v_0 = 0$, the solution asymptotically approaches zero without crossing the zero axis.

Over-critically damped SDOF

In this case, $\zeta > 1$ and

$$s = -\zeta\omega_n \mp \omega_n\sqrt{\zeta^2 - 1} = -\zeta\omega_n \mp \hat{\omega}$$

where

$$\hat{\omega} = \omega_n\sqrt{\zeta^2 - 1}$$

and, after some rearrangement, the general integral for the over-damped SDOF can be written

$$x(t) = \exp(-\zeta\omega_n t)\left(A\cosh(\hat{\omega}t) + B\sinh(\hat{\omega}t)\right)$$

OVER-CRITICALLY DAMPED SDOF

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Note that:

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- as for increasing ζ we have that $\hat{\omega} \to \zeta\omega_n$, the velocity with which the response approaches zero slows down for increasing ζ.

Measuring damping

The real dissipative behavior of a structural system is complex and very difficult to assess.

For convenience, it is customary to express the real dissipative behavior in terms of an *equivalent viscous damping*.

In practice, we *measure* the response of a SDOF structural system under controlled testing conditions and find the value of the viscous damping (or damping ratio) for which our simplified model best matches the measurements.

For example, we could require that, under free vibrations, the real structure and the simplified model exhibit the same decay of the vibration amplitude.

LOGARITHMIC DECREMENT

Consider a SDOF system in free vibration and two positive peaks, u_n and u_{n+m}, occurring at times $t_n = n\,(2\pi/\omega_D)$ and $t_{n+m} = (n+m)\,(2\pi/\omega_D)$.

The ratio of these peaks is

$$\frac{u_n}{u_{n+m}} = \frac{\exp(-\zeta\omega_n n 2\pi/\omega_D)}{\exp(-\zeta\omega_n(n+m)2\pi/\omega_D)} = \exp(2m\pi\zeta\omega_n/\omega_D)$$

Substituting $\omega_D = \omega_n\sqrt{1-\zeta^2}$ and taking the logarithm of both members we obtain

$$\ln(\frac{u_n}{u_{n+m}}) = \delta = 2m\pi\frac{\zeta}{\sqrt{1-\zeta^2}}$$

where we have introduced δ, the logarithmic decrement; solving for ζ, we finally get

$$\zeta = \delta\left((2m\pi)^2 + \delta^2\right)^{-\frac{1}{2}}.$$

Recursive Formula for ζ

The equation of the logarithmic decrement,

$$\delta = 2m\pi \frac{\zeta}{\sqrt{1 - \zeta^2}}$$

can be formally solved for ζ:

$$\zeta = \delta 2m\pi \sqrt{1 - \zeta^2}$$

obtaining an equation that can be interpreted as generating a sequence

$$\zeta_{n+1} = \delta 2m\pi \sqrt{1 - \zeta_n^2}.$$

Starting our sequence of successive approximations with $\zeta_0 = 0$ we obtain $\zeta_1 = \delta 2m\pi$ and usually the following iterate $\zeta_2 = \delta 2m\pi \sqrt{1 - \zeta_1^2}$ has converged to the *true* value with a number of digits that exceeds the experimental accuracy.

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While the recursive formula is useful in itself, it is also useful as a first example of finding better approximations of a system's parameter using an iterative procedure.