

An Introduction to Dynamics of Structures

Giacomo Boffi

<http://intranet.dica.polimi.it/people/boffi-giacomo>

Dipartimento di Ingegneria Civile Ambientale e Territoriale
Politecnico di Milano

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Let's start with some 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

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 :

$$p(t) = p_0 \sin(\omega t - \varphi)$$

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- ▶ 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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Formulation of a Dynamical Problem

In a structural system the inertial forces depend on the time derivatives of displacements while the elastic forces, equilibrating the inertial ones, depend on the space derivatives of the displacements. ... the natural statement of the problem is hence in terms of *partial differential equations*.

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

Lumped Masses

In many structural problems, we can say that the mass is concentrated in a discrete set of *lumped masses*.

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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 component of the inertial forces/masses displacements.

Dynamic Degrees of Freedom

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

If a lumped mass can be regarded as a *point* mass then 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 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.

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.

- ▶ 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.

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.

Writing the equation 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.

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The mathematical expressions, ordinary or partial differential equations, that we are going to write express the *dynamic equilibrium* of the structural system and are known as the *Equations of Motion* (EOM).

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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.

Writing the EOM, cont.

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,

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

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.

If we write

$$\vec{p}(t) - m\ddot{\vec{u}} = 0$$

and interpret the term $-m\ddot{\vec{u}}$ as an *Inertial Force* that contrasts the acceleration of the particle, we have an equation of equilibrium for the particle.

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.

For example, considering 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*, $\mathcal{L} = \mathcal{T} - \mathcal{V}$, \mathcal{T} and \mathcal{V} being, respectively, the kinetic and the potential energy of the system expressed in terms of a vector \vec{q} of independent coordinates

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

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

1 DOF System

Structural dynamics is all about the motion of a system in the neighborhood of a point of equilibrium.

We'll start by studying the most simple of systems, a single degree of freedom system, without external forces, subjected to a perturbation of the equilibrium.

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

$$\ddot{y} = \frac{1}{m}F(y, \dot{y}) = f(y, \dot{y}).$$

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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, $y_{\text{eq.}}$, the velocity and the acceleration are zero, and hence $f(y_{\text{eq.}}, 0) = 0$.

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

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

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

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

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

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In an infinitesimal neighborhood of $y_{\text{eq.}}$, the equation of motion can be studied in terms of a linear differential equation of second order.

1 DOF System, cont.

A linear constant coefficient differential equation has the 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

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

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...

1 DOF System, cont.

If we write $s_i = r_i + \imath q_i$, then we have

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

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 s 's are imaginary, 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.

The famous box car

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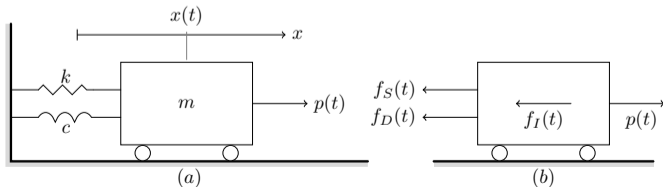
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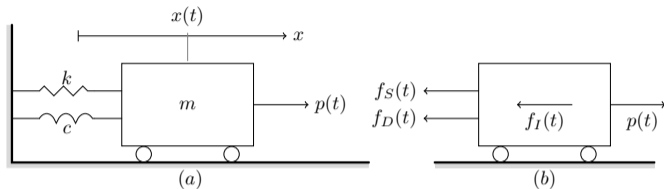


Equation of motion of the basic dynamic system

1 DOF System

Free vibrations of a SDOF system

Free vibrations of a damped system



The equation of motion can be written using the D'Alembert Principle, expressing the equilibrium of all the forces acting on the mass including the inertial force.

The forces, positive if acting in the direction of the motion, are the external force, $p(t)$, and the resisting forces due to motion, i.e., the inertial force $f_I(t)$, the damping force $f_D(t)$ and the elastic force, $f_S(t)$.

The equation of motion, merely expressing the equilibrium of these forces, is

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

EOM of the basic dynamic system, cont.

The resisting forces in

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

are functions of the displacement $x(t)$ or of one of its derivatives.
Note that the positive sense of these forces is opposite to the direction of motion.

EOM of the basic dynamic system, cont.

In accordance 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,

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

The differential equation of dynamic equilibrium

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

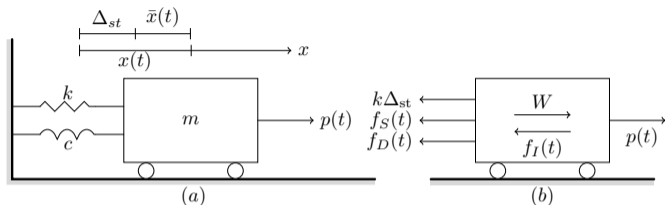
Note that this differential equation is a linear differential equation of the second order, with constant coefficients.

Influence of static forces

1 DOF System

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Considering the presence of a constant force W , the *EOM* is

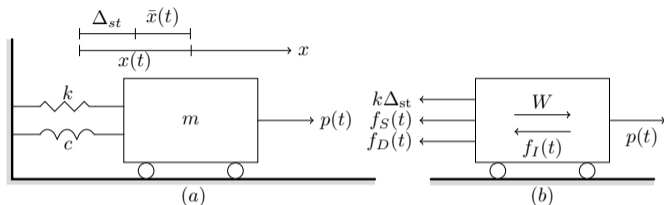
$$m\ddot{x}(t) + c\dot{x}(t) + kx(t) = p(t) + W.$$

Influence of static forces

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$$m\ddot{x}(t) + c\dot{x}(t) + kx(t) = p(t) + W.$$

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

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

Influence of static forces, cont.

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

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

Influence of static forces, cont.

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

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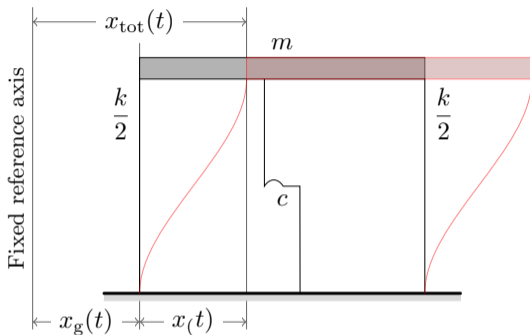
*Note that the total displacements, stresses. etc. are influenced by the static forces, and **must** be computed using the superposition of effects.*

Influence of support motion

1 DOF System

Free vibrations of a SDOF system

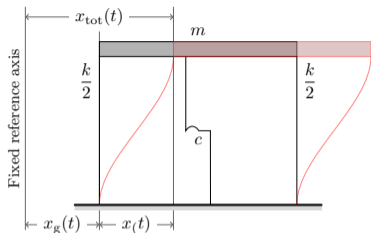
Free vibrations of a damped system



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 of a piece of equipment due to vibrations of the building in which it is housed.

Influence of support motion, cont.



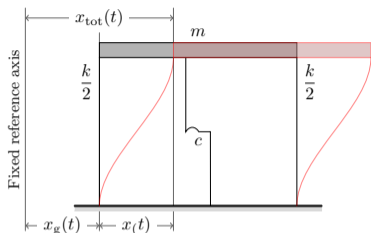
Considering a support motion $x_g(t)$, defined with respect to an inertial frame of reference, the total displacement is

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

and the total acceleration is

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

Influence of support motion, cont.



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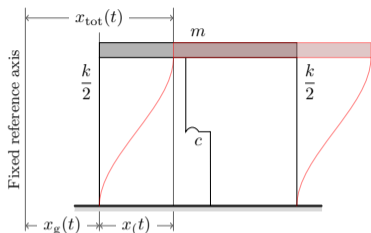
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$$\ddot{x}_{\text{tot}}(t) = \ddot{x}_g(t) + \ddot{x}(t).$$

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

$$f_1(t) = -m\ddot{x}_{\text{tot}}(t) = m\ddot{x}_g(t) + m\ddot{x}(t).$$

Influence of support motion, cont.



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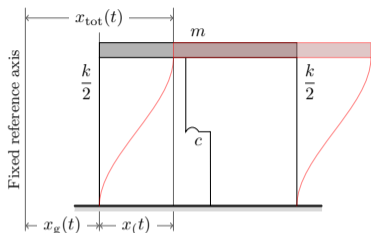
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$$f_I(t) = -m\ddot{x}_{tot}(t) = m\ddot{x}_g(t) + m\ddot{x}(t).$$

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

$$\begin{aligned} m\ddot{x}_{tot}(t) + c\dot{x}(t) + kx(t) &= 0, & \text{or,} \\ m\ddot{x}(t) + c\dot{x}(t) + kx(t) &= -m\ddot{x}_g(t) \equiv p_{\text{eff}}(t). \end{aligned}$$

Influence of support motion, cont.



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

The equation of motion,

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

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 a useful idealization. In this case, the homogeneous equation of motion is

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

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

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

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

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

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

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

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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

$$x(0) = x_0, \quad \dot{x}(0) = \dot{x}_0$$

Evaluating $x(t)$ for $t = 0$ and substituting in (39), 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

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

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

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

Undamped Free Vibrations

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Using the Euler formulas relating the imaginary argument exponentials and the trigonometric functions, can be rewritten in terms of the elementary trigonometric functions

$$x(t) = x_0 \cos(\omega_n t) + (\dot{x}_0/\omega_n) \sin(\omega_n t).$$

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Considering that for every conceivable initial conditions we can use the above representation, it is indifferent, and perfectly equivalent, to represent the general integral either in the form of exponentials of imaginary argument or 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

$$A = x_0, \quad B = \frac{\dot{x}_0}{\omega_n}.$$

Undamped Free Vibrations

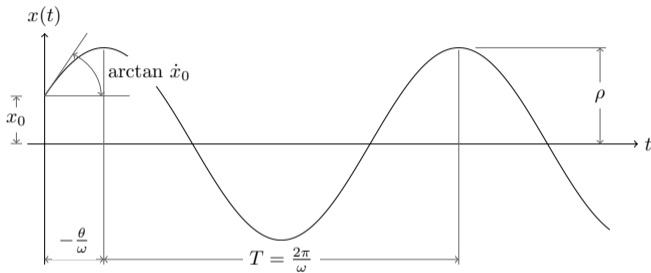
Sometimes we prefer to write $x(t)$ as a single harmonic, introducing a *phase difference* ϕ so that the amplitude of the motion, C , is put in evidence:

$$\begin{aligned}x(t) &= C \cos(\omega_n t - \phi) = C (\cos \omega_n t \cos \phi + \sin \omega_n t \sin \phi) \\ &= A \cos \omega_n t + B \sin \omega_n t\end{aligned}$$

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

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

Undamped Free Vibrations



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.

Behavior of Damped Systems

The viscous damping modifies the response of a *s dof* 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

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

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

$$m s^2 + c s + k = 0$$

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*,

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

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.

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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 ,

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

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 $\exp(-\zeta\omega_n t)$

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

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

$$x(t) = \exp(-\zeta\omega_n t) \left[\frac{\exp(i\omega_D t) + \exp(-i\omega_D t)}{2} u_0 + \frac{\exp(i\omega_D t) - \exp(-i\omega_D t)}{2i} \frac{v_0 + \zeta\omega_n u_0}{\omega_D} \right].$$

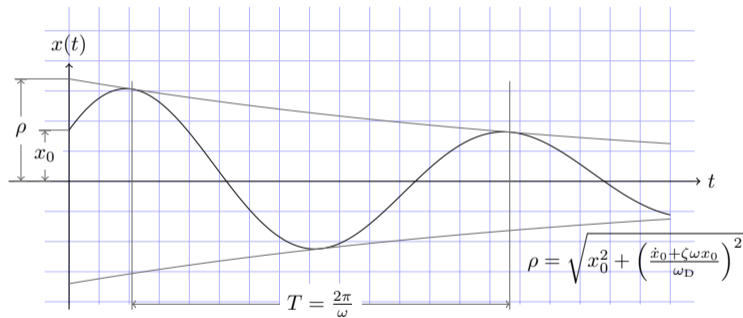
Using the Euler formulas, we finally have the preferred format of the general integral:

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

with

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

The Damped Free Response



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

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

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- ▶ as for increasing ζ we have that $\hat{\omega} \rightarrow \zeta\omega_n$, 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

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

Solving for ζ , we finally get

$$\zeta = \frac{\ln\left(\frac{u_n}{u_{n+m}}\right)}{\sqrt{(2m\pi)^2 + \left(\ln\left(\frac{u_n}{u_{n+m}}\right)\right)^2}}$$

Simplified solution

Starting from

$$\frac{u_n}{u_{n+m}} = \exp \frac{2m\pi\zeta}{\sqrt{1-\zeta^2}},$$

we note that, for the small values of ζ that are typical of structural systems, $\sqrt{1-\zeta^2} \approx 1$.

Using this approximation and the Taylor series representation of $\exp at$,

$$\frac{u_n}{u_{n+m}} \approx \exp \frac{2m\pi\zeta}{1} \approx 1 + 2m\pi\zeta + \dots$$

Solving for ζ ,

$$\zeta \approx \frac{1}{2m\pi} \frac{u_n - u_{n+m}}{u_{n+m}}.$$

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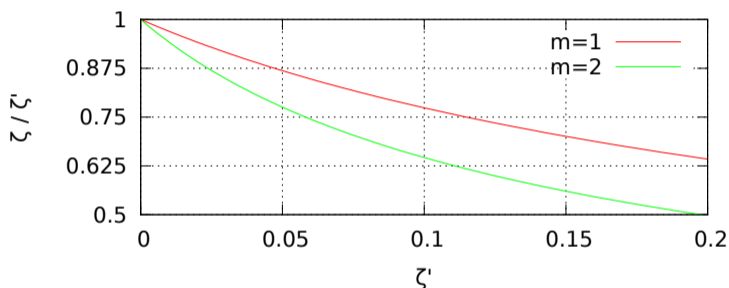
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Due to the uncertainties in the measurements and the modelization, for small values of ζ we can use the above approximation.

Errors

In abscissae the *approximate* value of ζ , in ordinates the ratio of true to approximate values



As you can see, the error grows with m , and the approximation is good only for really small values of ζ and $m = 1$.

A good alternative is to seek an iterative solution,

$$\delta = \frac{1}{2m\pi} \log \frac{u_n}{u_{n+m}}, \quad \zeta_0 = 0.0, \quad \zeta_{n+1} = \delta \sqrt{1 - \zeta_n^2}.$$