Viscous Damping (Sec. 1.4)



We generally assume Viscous dimping

$$F_{damper} = C J \leftarrow force is proportional to the rate of change
of the length$$



To find
$$x(t)$$
, use the same solution pacebolic as the unitamped case.
Assume $x(t) = \alpha e^{\lambda t}$ (so $\dot{x}(t) = \alpha he^{\lambda t}$ and $\ddot{x}(t) = \alpha he^{\lambda t}$)
plug there into the eq of matter
 $m(\alpha he^{\lambda t}) + c(\alpha he^{\lambda t}) + k(\alpha e^{\lambda t}) = 0$
 $(m\lambda^2 + c\lambda + k) = 0$
 $=0$ to have a nontrivial solution
 $m\lambda^2 + c\lambda + k = 0$
 $\lambda_{1,2} = \lambda m(-c \pm \sqrt{c^2 - 4}mk)$ Mot a very friendly form

Viscous Damping (cont.)

Look back at the equation of notion — $M\ddot{x} + C\dot{x} + K\dot{x} = 0$ $\ddot{x} + \frac{K}{m}\dot{x} = 0$ $Let_{m} = 25u_{n}$ $\dot{\xi} = domping rate$

Naw, the solution depends on E:

 $\chi + 2\xi\omega_{n} + \omega_{n}^{2}\chi = 0$

 $If \xi > 1: overdomped \qquad If \xi < 1 - unlerdomped \\ -no oscillation \qquad - Oscillates \\ - 2 negative real roots \qquad - complex conjugate roots \\ \lambda_{1,2} = -\xi_{n,1} \pm c_{n,1} \xi^{2} - 1 \qquad \lambda_{1,3} = -\xi_{n,1} \pm c_{n,1} \xi^{2} = -\xi_{n,2} \pm c_{n,1} \xi^{2} = -\xi_{n,2} \pm c_{n,2} \xi^{2} = -\xi_{n,2} + c_{n,2} + c_{$





Q: When might an overdamped response be desirable? Not desirable?

Desirable for no vibration or overshoot is desired

- * automatic doors
- * chemical processes

Undesirable:

- * Car suspensions (uncomfortable)
- * Controlled mechanical systems (too slow)

Underdamped Example Response



Time (s)

Energy

Kinetic Energy - energy of motion

Potential Energy - stored energy



Q: What about the spring potential?

Spring Potential

 $V_{gr} = \frac{1}{2}k d^{2}$ d = spring deflection from equil.

<u>Note</u> IF a mass is in pure rotation, we can use a simplified version of the kinetic energy $T = \frac{1}{2} \overline{\omega} \cdot \overline{H_0} \longrightarrow \frac{1}{2} \overline{I_0 \omega^2}$ Remember that this is a special case. Be careful when using

Lagrange's Equations/Method (Sec. 1.5)

- * Energy-based method
- * Allows us to ignore internal/interaction forces (if we want to)
- * Usually based around:
 - generalized coordinates an initial set of independent cards needed to describe the system
 - virtual displacements () set in derivation and for external forces

Generalized Coordinates



Lagrange's Equation (with no external forces or damping)

Define L=T-V = Kinetic Energy - Postential Energy < The Lograngian $\frac{d}{dt}\left(\frac{\partial L}{\partial \dot{q}_{i}}\right) - \frac{\partial L}{\partial \dot{q}_{i}} = 0, \quad (= 1, 2, ..., n) \quad \text{where } n \text{ is the number of generalized cosids } (= # of DOF)$

Note Be sure to include all system energies and define them consistently

Simple Linear Example of Lagrange's Method



Simple Rotational Example of Lagrange's Method

