# System Dynamics and Vibrations

Prof. Gustavo Alonso

Chapter 1: Elements of analytical dynamics Part 2

School of General Engineering Beihang University (BUAA)

#### Introduction

- >Analytical mechanics (Lagrange, variational approach)
  - Considers the <u>systems as a whole</u> → reaction and constraint forces are excluded
  - Dynamics problems formulated in terms of: kinetic energy, potential energy, virtual work of non-conservative forces
  - Equations of motion formulated in terms of generalized coordinates and generalized forces → broader and more abstract approach
  - The mathematical formulation is independent of any special system of coordinates

#### Contents

- Introduction
- Degrees of freedom and generalized coordinates
- The principle of virtual work
- The principle of D'Alembert
- Lagrange's equations

- Principle of virtual work 
   static equilibrium of systems
- Principle of D'Alembert → extension of the principle of virtual work to dynamics (for example, vibrations)

- System of N particles, of mass  $m_i$ , acted upon the applied force  $\mathbf{F}_i$  and the constraint force  $\mathbf{f}_i$ , assuming any internal force to be negligibly small:
- Rewriting Newton's second law:

$$\mathbf{F}_{i} + \mathbf{f}_{i} - m_{i}\ddot{\mathbf{r}}_{i} = \mathbf{0}, \quad i = 1, 2, ...N$$

→ Principle of D'Alembert

 $-m_i \ddot{\mathbf{r}}_i$  is the inertia force

Dynamic problems can be solved as if they were static

- The interest is to extend the principle of virtual work to the dynamical case
- The virtual work for particle  $m_i$  is:

$$(\mathbf{F}_i + \mathbf{f}_i - m_i \ddot{\mathbf{r}}_i) \cdot \delta \mathbf{r}_i = \mathbf{0}, \quad i = 1, 2, ...N$$

- Let's confine ourselves to constraint forces for which the virtual work is zero
- Summing up over the system of particles:

$$\sum_{i=1}^{N} (\mathbf{F}_{i} - m_{i} \ddot{\mathbf{r}}_{i}) \cdot \delta \mathbf{r}_{i} = 0$$

principle of virtual work + principle of D'Alembert 

generalized principle of D'Alembert

(Lagrange's version of D'Alembert's principle)

Generalized principle of D'Alembert:

The virtual work performed by the effective forces through infinitesimal virtual displacements compatible with thesystem constraints is zero

$$\sum_{i=1}^{N} (\mathbf{F}_{i} - m_{i} \ddot{\mathbf{r}}_{i}) \cdot \delta \mathbf{r}_{i} = 0$$

 $\mathbf{F}_i - m_i \ddot{\mathbf{r}}_i$   $\rightarrow$  applied force acting on particle  $m_i$ 

- The real interest is not just to derive the equations of motion, but
- To derive the extended Hamilton's principle, and then:
- To derive all system equations of motion from three scalar quantities:
  - Kinetic energy
  - Potential energy
  - Virtual work of non-conservative forces
- And then to derive the Lagrange's equations

- First case: the **position vectors**  $\mathbf{r}_i$  (i = 1, 2, ..., N) are all independent

• Starting from 
$$\sum_{i=1}^{N} (\mathbf{F}_{i} - m_{i} \ddot{\mathbf{r}}_{i}) \cdot \delta \mathbf{r}_{i} = 0$$

$$\sum_{i=1}^{N} \mathbf{F}_{i} \cdot \delta \mathbf{r}_{i} = \overline{\delta W}$$

virtual work of all applied forces

$$\sum_{i=1}^{N} \left( \mathbf{F}_{i} - m_{i} \ddot{\mathbf{r}}_{i} \right) \cdot \delta \mathbf{r}_{i} = 0$$

$$\frac{d}{dt}\left(m_{i}\dot{\mathbf{r}}_{i}\cdot\delta\mathbf{r}_{i}\right)=m_{i}\ddot{\mathbf{r}}_{i}\cdot\delta\mathbf{r}_{i}+m_{i}\dot{\mathbf{r}}_{i}\cdot\delta\dot{\mathbf{r}}_{i}=m_{i}\ddot{\mathbf{r}}_{i}\cdot\delta\mathbf{r}_{i}+\delta\left(\frac{1}{2}m_{i}\dot{\mathbf{r}}_{i}\cdot\dot{\mathbf{r}}_{i}\right)=m_{i}\ddot{\mathbf{r}}_{i}\cdot\delta\mathbf{r}_{i}+\delta T_{i}$$

$$T_i = \frac{1}{2} m_i \dot{\mathbf{r}}_i \cdot \dot{\mathbf{r}}$$
 is the kinetic energy of particle  $m_i$ 

$$-\int_{t_1}^{t_2} m_i \ddot{\mathbf{r}}_i \cdot \delta \mathbf{r}_i dt = \int_{t_1}^{t_2} \delta T_i dt - \int_{t_1}^{t_2} \frac{d}{dt} \left( m_i \dot{\mathbf{r}}_i \cdot \delta \mathbf{r}_i \right) dt = \int_{t_1}^{t_2} \delta T_i dt - m_i \dot{\mathbf{r}}_i \cdot \delta \mathbf{r}_i \Big|_{t_1}^{t_2}$$

$$-\int_{t_1}^{t_2} m_i \ddot{\mathbf{r}}_i \cdot \delta \mathbf{r}_i dt = \int_{t_1}^{t_2} \delta T_i dt - \int_{t_1}^{t_2} \frac{d}{dt} \left( m_i \dot{\mathbf{r}}_i \cdot \delta \mathbf{r}_i \right) dt = \int_{t_1}^{t_2} \delta T_i dt - m_i \dot{\mathbf{r}}_i \cdot \delta \mathbf{r}_i \Big|_{t_1}^{t_2}$$

- The virtual displacements are arbitrary
- $\rightarrow$  we choose them as to satisfy  $\delta \mathbf{r}_i = 0$  at  $t = t_1$  and  $t = t_2$

$$-\int_{t}^{t_2} m_i \ddot{\mathbf{r}}_i \cdot \delta \mathbf{r}_i dt = \int_{t}^{t_2} \delta T_i dt, \quad \delta \mathbf{r}_i = 0, \quad t = t_1, t_2; \quad i = 1, 2, \dots N$$

• Summing up over *i*:

$$-\int_{t_1}^{t_2} \sum_{i=1}^{N} m_i \ddot{\mathbf{r}}_i \cdot \delta \mathbf{r}_i dt = \int_{t_1}^{t_2} \delta T dt, \quad \delta \mathbf{r}_i = 0, \quad t = t_1, t_2; \ i = 1, 2, ...N$$

T is the kine of the following states and the states are the states a

T is the system kinetic energy

$$\int_{t_1}^{t_2} \left( \delta T + \overline{\delta W} \right) dt = 0, \quad \delta \mathbf{r}_i = 0, \quad i = 1, 2, ... N; \quad t = t_1, t_2$$

- extended Hamilton's principle
- Virtual work is  $\overline{\delta W} = \overline{\delta W_c} + \overline{\delta W_{nc}} = -\delta V + \overline{\delta W_{nc}}$  where V is the potential energy
- → then:

$$\int_{t_1}^{t_2} \left( \delta T - \delta V + \overline{\delta W_{nc}} \right) dt = 0, \quad \delta \mathbf{r}_i = 0, \quad i = 1, 2, ... N; \quad t = t_1, t_2$$

$$\int_{t_1}^{t_2} \left( \delta T - \delta V + \overline{\delta W_{nc}} \right) dt = 0, \quad \delta \mathbf{r}_i = 0, \quad i = 1, 2, ... N; \quad t = t_1, t_2$$

- → All the equations of motion can be obtained from three scalar quantities:
  - Kinetic energy
  - Potential energy
  - Virtual work of non-conservative forces

- Second case: the <u>position vectors</u> r<sub>i</sub> (i = 1, 2, ..., N) are <u>not independent</u>
- They are related by some constraint equations

$$\int_{t_1}^{t_2} \left( \delta T - \delta V + \overline{\delta W_{nc}} \right) dt = 0, \quad \delta \mathbf{r}_i = 0, \quad i = 1, 2, ... N; \quad t = t_1, t_2$$

The Kinetic energy, Potential energy, and Virtual work of non-conservative forces: all are independent of the coordinates used

→ The extended Hamilton's principle retains its form for all sets of coordinates

• Let's choose the independent generalized coordinates:

$$\int_{t_1}^{t_2} \left( \delta T - \delta V + \overline{\delta W_{nc}} \right) dt = 0, \quad \delta q_k = 0, \quad k = 1, 2, ...n; \quad t = t_1, t_2$$

- The extended Hamilton's principle in terms of generalized coordinates can be used to derive all system equations of motion (regardless the system is subjected to constraints or not)
- The only condition is that the <u>constraint forces perform no</u> <u>work</u>

- For conservative systems:  $\delta W_{nc} = 0$
- → Hamilton's principle:

$$\int_{t_1}^{t_2} \delta L dt = 0, \quad \delta q_k = 0, \quad k = 1, 2, ... n; \quad t = t_1, t_2$$

$$L = T - V$$

L is known as the Lagrangian

#### Contents

- Introduction
- Degrees of freedom and generalized coordinates
- The principle of virtual work
- The principle of D'Alembert
- Lagrange's equations

- Lagrange's equations can be derived directly from D'Alembert's principle or from the Extended Hamilton's principle (simplest)
- Kinetic energy:  $T = T(q_1, q_2, ..., q_n, \dot{q}_1, \dot{q}_2, ..., \dot{q}_n)$

$$\delta T = \sum_{k=1}^{n} \left( \frac{\delta T}{\delta q_k} \delta q_k + \frac{\delta T}{\delta \dot{q}_k} \delta \dot{q}_k \right)$$

• Potential energy:  $V = V(q_1, q_2, ..., q_n)$ 

$$\delta V = \sum_{k=1}^{n} \left( \frac{\delta V}{\delta q_k} \delta q_k \right)$$

• Virtual work of non-conservative forces:  $\overline{\delta W_{nc}} = \sum_{k=1}^{n} Q_k \delta q_k$ 

$$Q_k$$
  $(k = 1, 2, ..., n)$   $\rightarrow$  generalized non-conservative forces

$$\int_{t_1}^{t_2} \left( \delta T - \delta V + \overline{\delta W_{nc}} \right) dt = \int_{t_1}^{t_2} \sum_{k=1}^{n} \left[ \left( \frac{\delta T}{\delta q_k} - \frac{\delta V}{\delta q_k} + Q_k \right) \delta q_k + \frac{\delta T}{\delta \dot{q}_k} \delta \dot{q}_k \right] dt = 0$$

$$\delta q_k = 0, \quad k = 1, 2, ...n; \quad t = t_1, t_2$$

Integrating by parts:

$$\int_{t_{1}}^{t_{2}} \frac{\delta T}{\delta q_{k}} \delta \dot{q}_{k} dt = \int_{t_{1}}^{t_{2}} \frac{\delta T}{\delta \dot{q}_{k}} \frac{d}{dt} \delta q_{k} dt = \frac{\delta T}{\delta \dot{q}_{k}} \delta q_{k} \bigg|_{t_{1}}^{t_{2}} - \int_{t_{1}}^{t_{2}} \frac{d}{dt} \left( \frac{\delta t}{\delta \dot{q}_{k}} \right) \delta q_{k} dt$$

$$= -\int_{t_{1}}^{t_{2}} \frac{d}{dt} \left( \frac{\delta T}{\delta \dot{q}_{k}} \right) \delta q_{k} dt, \qquad k = 1, 2, ... n;$$

It has been considered  $\delta q_k = 0$  (k = 1, 2, ...n) at  $t = t_1$  and  $t = t_2$ 

$$\int_{t_1}^{t_2} \sum_{k=1}^{n} \left[ \frac{\delta T}{\delta q_k} - \frac{\delta V}{\delta q_k} + Q_k - \frac{d}{dt} \left( \frac{\delta T}{\delta \dot{q}_k} \right) \right] \delta q_k dt = 0$$

- Since the generalized virtual displacements are arbitrary
- We assign arbitrary values to  $\delta q_1$  while setting  $\delta q_2 = 0$

$$(k = 2, 3, ...n)$$

$$\int_{t_1}^{t_2} \sum_{k=1}^{n} \left[ \frac{\delta T}{\delta q_k} - \frac{\delta V}{\delta q_k} + Q_k - \frac{d}{dt} \left( \frac{\delta T}{\delta \dot{q}_k} \right) \right] \delta q_k dt = 0 \quad \Rightarrow \text{ the coefficient of } \delta q_1 \text{ is zero}$$

• Repeating the argument for all  $\delta q_2, \delta q_3, ... \delta q_n$ 

$$\frac{d}{dt} \left( \frac{\delta T}{\delta \dot{q}_k} \right) - \frac{\delta T}{\delta q_k} + \frac{\delta V}{\delta q_k} = Q_k, \quad (k = 1, 2, ...n)$$
 Lagrange equations