Advanced Dynamics

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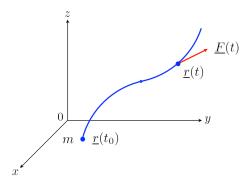
1 Dynamics

1.1 Coordinate frames

1.1.1 Inertial coordinate frame

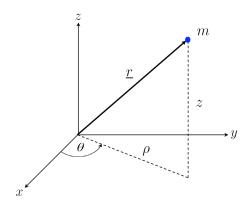
A frame of reference (coordinate system) that is not rotating and not accelerated.

1.1.2 Cartesian coordinate frame



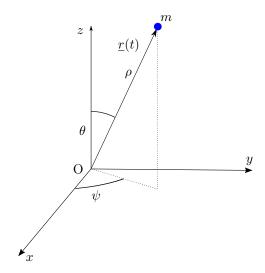
$$\underline{r} = \begin{pmatrix} x(t) \\ y(t) \\ z(t) \end{pmatrix} \quad \underline{v} = \begin{pmatrix} \dot{x}(t) \\ \dot{y}(t) \\ \dot{z}(t) \end{pmatrix} \quad \underline{a} = \begin{pmatrix} \ddot{x}(t) \\ \ddot{y}(t) \\ \ddot{z}(t) \end{pmatrix}$$

1.1.3 Cylindrical coordinate frame



$$\underline{r} = \begin{pmatrix} \rho(t) \\ \theta(t) \\ z(t) \end{pmatrix} \ \underline{v} = \begin{pmatrix} \dot{\rho}(t) \\ \rho(t)\dot{\theta}(t) \\ \dot{z}(t) \end{pmatrix} \ \underline{a} = \begin{pmatrix} \ddot{\rho}(t) - \rho(t)\dot{\theta}(t)^2 \\ 2\dot{\rho}(t)\dot{\theta}(t) + \rho(t)\ddot{\theta}(t) \\ \ddot{z}(t) \end{pmatrix}$$

1.1.4 Spherical coordinate frame



$$\underline{r} = \begin{pmatrix} \rho(t) \\ \theta(t) \\ \psi(t) \end{pmatrix} \underline{v} = \begin{pmatrix} \dot{\rho} \\ \rho \dot{\theta} cos(\psi) \\ \rho \dot{\psi} \end{pmatrix} \underline{a} = \begin{pmatrix} \ddot{\rho} - \rho \dot{\theta}^2 \cos^2 \psi - \rho \dot{\psi}^2 \\ 2\dot{\rho} \dot{\theta} \cos(\psi) + \rho \ddot{\theta} \cos(\psi) - 2\rho \dot{\theta} \dot{\psi} \sin(\psi) \\ 2\dot{\rho} \dot{\psi} + \rho \dot{\theta} \sin(\psi) \cos(\psi) + \rho \ddot{\psi} \end{pmatrix}$$

1.2 Constraints

A constraint is a scalar relation that limits the possible displacement of particles.

(a) Rigid link

$$|\underline{r}_{AB}| = \sqrt{(x_A - x_B)^2 + (y_A - y_B)^2 + (z_A - z_B)^2} = const$$

(b) Rotating tube

$$\begin{pmatrix} x_A \\ y_A \\ z_A \end{pmatrix} = \rho_A(t)\underline{e}(t)$$

where $\underline{e}(t)$ denotes the direction of the rotating tube.

(c) Planar motion

$$z_A = const$$

(d) Rolling without slipping

$$\dot{x} = R\dot{\phi}$$

Holonomic constraint:

A holonomic constraint is a constraint that only depends on the current position of the particles involved and the current time. A non-holonomic constraint also depends on time-derivatives of the motion.

1.3 Generalized coordinates

A set of variables used to uniquely describe the configuration of a system are called generalized coordinates. This can comprise, e.g., distances or angles.

1.4 Degree of Freedom

The degree of freedom (DOF) of a system is defined as the number of independent generalized coordinates necessary to describe the motion of the system completely and uniquely.

$$DOF = dim \cdot n - k$$

k number of holonomic constraints applied to the system

• There are always different possibilities of interpreting the DOF of a multi body system. Be sure that you define the constraints such that each body you consider is only moving in the exact way it is supposed to.

1.5 GIVENS ROTATION

Givens rotation matrices define a rotation about one of the three coordinate directions about some angle α , β , or γ referring to a rotation about the x-axis, y-axis or z-axis respectively.

$$\underline{\underline{R}}_{x} = \begin{pmatrix} 1 & 0 & 0 \\ 0 & \cos \alpha & \sin \alpha \\ 0 & -\sin \alpha & \cos \alpha \end{pmatrix}$$

$$\underline{\underline{R}}_{y} = \begin{pmatrix} \cos \beta & 0 & -\sin \beta \\ 0 & 1 & 1 \\ \sin \beta & 0 & \cos \beta \end{pmatrix}$$

$$\underline{\underline{R}}_{z} = \begin{pmatrix} \cos \gamma & \sin \gamma & 0 \\ -\sin \gamma & \cos \gamma & 0 \\ 0 & 0 & 1 \end{pmatrix}$$

1.6 Taylor series

The Taylor series of a function f(x), which is infinitely differentiable at some number a is given by,

$$f(a) + \frac{f'(a)}{1!}(x-a) + \frac{f''(a)}{2!}(x-a)^2 + \frac{f'''(a)}{3!}(x-a)^2 + \cdots,$$

or equivalently,

$$\sum_{n=0}^{\infty} \frac{f^{(n)}(a)}{n!} (x-a)^n.$$

1.6.1 Small angle approximation

Based on a Taylor series of the second order some trigonometrical functions can be approximated well for small angles.

$$\sin \theta \approx \theta$$

$$\cos \theta \approx 1 - \frac{\theta^2}{2}$$

$$\tan \theta \approx \theta$$

2 Single-particle dynamics

2.1 Linear momentum

$$\underline{P} := m\underline{v}$$

$$\underline{P} \quad \text{linear momentum} \quad \text{in} \quad \text{kg m s}^{-1}$$

$$m \quad \text{mass} \dots \dots \quad \text{in} \quad \text{kg}$$

$$v \quad \text{velocity} \dots \dots \quad \text{in} \quad \text{m s}^{-1}$$

2.2 Principle of linear momentum for a single particle (LMP)

The linear momentum principle is equivalent to Newton's second axiom. It states that a resultant forces acting on a particle causes an acceleration of the particle.

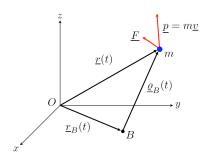
- Use for problems based on linear motion.
- Use to find equation of motion based on the acting forces or find the acting forces based on a defined motion.
- If no resultant force is acting on the particle, the acceleration is zero and the linear momentum of the particle is conserved.
- The knowledge of all forces acting on a particle leads directly to its acceleration. If the initial velocity is known, the particle's velocity can be derived as a function of time. If the initial position is known additionally, the particle's position can be calculated as well.
- The LMP is described by a vectorial equation. Therefore you can set up a scalar equation for each coordinate direction.
- Only valid in an inertial frame!

Remark on LMP in cylindrical coordinates:

In case of a curvilinear motion where cylindrical coordinates are helpful, you need to express the vector $m\underline{a}$ as a function of the cylindrical coordinates (ρ, θ) and project it to the radial and azimuthal unit vectors $\underline{e}_{\rho},\underline{e}_{\theta}$ afterwards. With this trick you get the linear momentum principle in cylindrical coordinates (forces acting on the particle are then expressed in the $\underline{e}_{\alpha},\underline{e}_{\theta}$ system).

A special case of this is what has been derived in Problem 2.2, where we treat a circular motion with a constant velocity.

2.3 Angular momentum



$$\underline{H}_B(t) := \underline{\varrho}_B(t) \times \underline{P}(t)$$

$\underline{H}_B(t)$	angular momentum w.r.t. point B	$_{ m in}$	$\mathrm{kg}\mathrm{m}^2\mathrm{s}^{-1}$
$\rho_{_{B}}(t)$	vector between point B and the particle m	$_{ m in}$	m
	linear momentum of particle m		

2.4 Principle of angular momentum for a single particle w.r.t. B (AMP)

The angular momentum principle can be seen as Newton's second axiom for rotation. It states that the angular momentum of a particle is influenced by the torque acting on the particle. The angular momentum itself can be calculated w.r.t an arbitrary reference point B which can be fixed or moving.

$$\underline{\dot{H}}_B + \underline{v}_B \times \underline{P} = \underline{M}_B$$

$\underline{\dot{H}}_{B}$	change angular momentum w.r.t. B	in	$\mathrm{kg}\mathrm{m}^2\mathrm{s}^{-2}$
\underline{v}_B	velocity of B	in	$\mathrm{m}\mathrm{s}^{-1}$
$\underline{\underline{P}}$	linear momentum		
M_{B}	resultant torque w.r.t. $B \dots \dots$	in	$N m = kg m^2 s^{-2}$

- \bullet Often used if a particle is in curvilinear motion.
- If $\underline{v}_B = 0$ or $\underline{v}_B || \underline{P}$ and $\underline{M}_B = 0$, the angular momentum w.r.t point B is conserved.
- If the angular momentum is conserved and the particle's initial velocity is given, it is often useful to equate the initial angular momentum with the particle's angular momentum at a later time.

- Choose points with $\underline{v}_B = 0$ to avoid calculating the terms of the cross product.
- Although the AMP is a vectorial equation, it is convenient to write it in a scalar form if a 2D problem needs to be solved. In the 2D case, the angular momentum and therefore it's derivative are always normal to the plane of projection.

2.5 Work-energy principle for a single particle

A force acting on a moving particle will change the particle's kinetic energy T_i .

$$W_{12} = T_2 - T_1$$

 W_{12} work done on particle between point 1 and 2 in $J = kg m^2 s^{-2}$ T_i Kinetic energy at point i in $kg m^2 s^{-2}$

• The kinetic energy of a particle is given by,

$$T_i = \frac{1}{2}mv_i^2.$$

- Used if only the initial and end state of a particle is of interest and not the states in between.
- Can easily be calculated if the force acting on the particle is parallel to the particle's velocity.
- The work done by a force is zero, if the force is normal to the particle's velocity.
- The work done on a particle by a force is given by,

$$W = \int_{P_1}^{P_2} \underline{F} \cdot \underline{dr} = \int_{P_1}^{P_2} |\underline{F}| \cos(\alpha) dS.$$

2.6 Definition of a potential force

A force is called potential force if a potential functions exists, such that the negative gradient of this function is equal to the force.

$$\underline{F} = -\frac{\partial V}{\partial r} = -\nabla V(\underline{r})$$

 $\begin{array}{lll} \underline{F} & \text{potential force...} & \text{in} & N = \text{kg m s}^{-2} \\ V & \text{potential function} & \text{in} & N \, \text{m} = \text{kg m}^2 \, \text{s}^{-2} \\ r & \text{position vector...} & \text{in} & \text{m} \end{array}$

If a force has a potential, the work done by this force between two points does not depend on the path between the points. In this case, the work done by the force can be calculated by the difference of the values of the potential function in this points, i.e., $W_{12} = V_1 - V_2 = V(x_1, y_1, z_1) - V(x_2, y_2, z_2)$.

An example for a potential force is the gravitational force. It's potential is given by,

$$V(\underline{r}) = mgz,$$

because,

$$-\nabla V(\underline{r}) = \begin{pmatrix} 0 \\ 0 \\ -mg \end{pmatrix}.$$

2.7 Work-energy principle for a conservative system

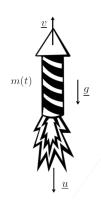
The work energy principle for a conservative system holds true, if all forces acting on a system are either potential (=conservative) or do no work. In this case, the work done by potential forces can be expressed as the difference of the potentials at the start and the end point of the desired motion.

$$T_1 + V_1 = T_2 + V_2 \quad \Leftrightarrow \quad \frac{d}{dt}[T(t) + V(t)] = 0$$

- T_i Kinetic energy in point i. in $J = kg m^2 s^{-2}$
- V_i Potential energy in point i in $J = kg m^2 s^{-2}$
- T(t) Kinetic energy at time t.. in $J = kg m^2 s^{-2}$
- V(t) Potential energy at time t in $J = kg m^2 s^{-2}$
- Always check if the system of interest is conservative
- A conservative system has the advantage that you do not need to know the path of the particle to calculate the work done by the forces acting on it.
- If some of the forces are potential and some not, express the work done by the potential forces as a difference of the potentials and calculate the work done by the non-conservative forces with the standard procedure, if you know the path of the particle.
- Some common examples for potentials are given by:

Force	Potential
Gravitation	mgz
Linear spring	$-\frac{1}{2}kx^{2}$
Torsional spring	$-\frac{1}{2}k\theta^2$

2.8 Mass accretion equation



The mass accretion equation describes the motion of a particle that loses or gains mass and is under the influence of some external force. The loss or gain of mass is described by a change of momentum $\dot{m}u$ where u denotes the velocity of the gained or ejected mass.

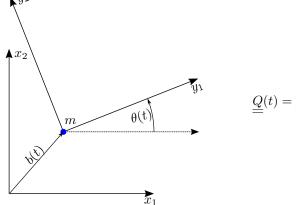
$$\dot{m}\underline{v} + m\dot{\underline{v}} - \dot{m}\underline{u} = \underline{F} \quad \Leftrightarrow \quad m\dot{\underline{v}} = \underline{F} + \dot{m}\underline{v}_{rel}$$

m	mass depending on time	in	kg
\dot{m}	change of mass	in	$\mathrm{kg}\mathrm{s}^{-1}$
\underline{v}	velocity of the particle	in	$\mathrm{m}\mathrm{s}^{-1}$
\underline{u}	velocity of the ejected or gained mass	in	$\mathrm{m}\mathrm{s}^{-1}$
$\overline{\dot{v}}$	acceleration	in	$\mathrm{m}\mathrm{s}^{-2}$
\underline{F}	resultant force	in	$N = kg m s^{-2}$
$\underline{v}_{rel} = \underline{u} - \underline{v}$	relative velocity		

An exemplary application of this principle is a rocket that fights gravity while it is loosing mass. In that case the acting force is the gravitational force, v represents the velocity of the rocket, \underline{u} represents the velocity of the exhaust gases. Note that the gravitational force varies as well since the mass of the rocket varies.

2.9 General transformation to a moving, rotating y-frame

Such a frame can be described as follows:



$$\underline{\underline{Q}}(t) = \underline{\underline{R}}_z = \begin{pmatrix} \cos \theta & -\sin \theta & 0\\ \sin \theta & \cos \theta & 0\\ 0 & 0 & 1 \end{pmatrix}$$

$$\underline{\underline{x}} = \underline{\underline{\underline{Q}}}(t)\underline{\underline{y}} + \underline{\underline{b}}(t)$$

\underline{x}	position vector in inertial x-frame	in	m
\overline{y}	position vector in non-inertial y-frame	in	m
$\overline{Q}(t)$	rotation matrix that aligns the y axes with the \underline{x} axes		
	$\underline{\underline{Q}}^T = \underline{\underline{Q}}^{-1}$ and $\det\left(\underline{\underline{Q}}\right) = 1$		
b(t)	translational vector between the two origins of the coordinate systems	$_{ m in}$	m

2.10 Linear momentum principle in a planar, rotating frame

Considering particles in non-inertial frames we have to consider the influence of non-inertial forces

$$\begin{split} m \underline{\ddot{y}} &= \underline{\tilde{F}} + \underline{F}_{centrifugal} + \underline{F}_{Coriolis} + \underline{F}_{Euler} \\ \underline{F}_{centrifugal} &= m \underline{\Omega}^2 \underline{y} \\ \underline{F}_{Coriolis} &= -2m\underline{\Omega} \times \underline{\dot{y}} \\ \underline{F}_{Euler} &= -m\underline{\dot{\Omega}} \times \underline{y} \end{split}$$

m	mass	$_{ m in}$	kg
\dot{y}	velocity in y-frame	$_{ m in}$	$\mathrm{m}\mathrm{s}^{-1}$
$\overline{\ddot{y}}$	acceleration in y-frame	$_{ m in}$	$\mathrm{m}\mathrm{s}^{-2}$
$\overline{ ilde{F}}$	active and constraint force w.r.t. the moving frame	$_{ m in}$	$N = kg m^2 s^{-2}$
$\underline{\Omega}$	angular velocity of the y-frame		
$\dot{\Omega}$	angular acceleration of the y-frame	$_{ m in}$	$\rm rads^{-2}$

• Note that the y-frame denotes the moving frame.

3 Multi-particle dynamics

3.1 Center of mass for a system of particles

The center of mass represents the location in which the resultant gravitational force of all included particles attacks. Or in other words it is the point in which the resultant moment of all gravitational forces is zero.

$$\underline{r}_C = \frac{1}{M} \sum_{i=1}^n m_i \underline{r}_i, \text{ where } M = \sum_{i=1}^n m_i$$

\underline{r}_C	center of mass	in	m
\check{M}	total mass of the system	in	kg
m_i	individual masses of particles	in	kg
\underline{r}_i	locations of particles	in	\mathbf{m}
n	number of particles		

3.2 Linear momentum of a system of particles

$$\underline{P} := \sum_{i=1}^{n} \underline{P}_{i} = \sum_{i=1}^{n} m_{i} \underline{v}_{i}$$

3.3 Resultant force on a system of particles

$$\underline{F} := \sum_{i=1}^{n} \underline{F}_{i}^{int} + \sum_{i=1}^{n} \underline{F}_{i}^{ext}$$

• Note that the sum of the inner forces is zero by definition.

3.4 Linear momentum principle for a system of particles

The linear momentum principle for a system of particles is equivalent to Newton's second axiom. It states that a resultant forces acting on such a system causes an acceleration of the center of mass.

$$\underline{\dot{P}} = M\underline{a}_C = \sum_{i=1}^{n} \underline{F}_i^{ext} = \underline{F}^{ext}$$

$\underline{\dot{P}}$	change of linear momentum	$_{ m in}$	$ m kgms^{-2}$
M	total mass of the system	in	kg
a_C	acceleration of the center of mass	in	$\mathrm{m}\mathrm{s}^{-2}$
$\frac{F_i^{ext}}{F^{ext}}$	individual resultant forces on particles	in	$N = kg m s^{-2}$
\underline{F}^{ext}	total resultant force on system	in	$N = kg m s^{-2}$
n	number of particles		

- If no resultant force is acting on the system, the acceleration of the center of mass is zero and the linear momentum of the system is conserved.
- If all external forces on a system are known the acceleration of the center of mass can be determined. Vice versa if the acceleration of the center of mass is known the sum of the external forces can be calculated.
- Since the linear momentum principle is a vectorial equation it can be split up and considered in each coordinate direction separately. This means also that it can also be conserved in one direction while it changes in the other direction.
- Only valid in an inertial frame!

$3.5\,$ Angular momentum principle for a system of particles w.r.t. a point B

The angular momentum principle is the same as in the case of a single particle. The only difference is, that \underline{P} is equal to the linear momentum of the system of particles, $\underline{P} = m\underline{v}_C$.

$$\underline{\dot{H}}_B + \underline{v}_B \times \underline{P} = \underline{M}_B^{ext}$$

$$\begin{array}{lll} \underline{\dot{H}}_B & \text{change of angular momentum w.r.t. } B & \text{in} & \log \mathrm{m}^2 \, \mathrm{s}^{-2} \\ \underline{v}_B & \text{velocity of point } B \dots & & \text{in} & \mathrm{m} \, \mathrm{s}^{-1} \\ \underline{P} & \text{linear momentum of the system} \dots & \text{in} & \log \mathrm{m} \, \mathrm{s}^{-1} \\ \underline{M}_B^{ext} & \text{external torque acting on the system} & & \text{in} & \mathrm{N} \, \mathrm{m} = \log \mathrm{m}^2 \, \mathrm{s}^{-2} \end{array}$$

• The angular momentum is conserved if and only if:

$$-\underline{M}_{B}^{ext} \equiv 0$$
 AND either $\underline{v}_{B} \equiv 0$ or $\underline{P} \equiv 0$.

- Alternatively it is possible that: $\underline{M}_{B}^{ext} \underline{v}_{B} \times \underline{P} \equiv 0$.
- When considering impulsive forces it makes sense to assess the change of angular momentum over an infinitesimally short time, therefore allowing negligence of non-impulsive forces.

3.6 Collision of two particles: coefficient of restitution

 $e = -\frac{\text{relative normal velocity after impact}}{\text{relative normal velocity before impact}} = -\frac{u_2^n - u_1^n}{v_2^n - v_1^n}$

- e coefficient of restitution.....
- u_i^n individual normal velocities after impact. in ms⁻¹
- v_i^n individual normal velocities before impact in ms⁻¹
- e=1 implies that we are dealing with a perfectly elastic collision.
- e = 0 implies that we are dealing with a perfectly inelastic collision.
- Due to energy conservation e must satisfy 0 < e < 1.
- e can be interpreted as a measure of the kinetic energy preserved during the collision.

 The amount of energy lost goes into deformation of the bodies.
- The linear momentum of the individual particles is always conserved in tangential direction.
- The linear momentum of the system is always conserved, since no forces act in tangential direction and the forces in normal direction are opposite in direction and equal in magnitude and therefore cancel out.

3.7 VELOCITY RELATION BETWEEN VELOCITIES BEFORE AND AFTER IMPACT

$$u_1^n = \frac{m_1 - em_2}{m_1 + m_2} v_1^n + \frac{(1+e)m_2}{m_1 + m_2} v_2^n \tag{1}$$

$$u_2^n = \frac{(1+e)m_1}{m_1 + m_2} v_1^n + \frac{m_2 - em_1}{m_1 + m_2} v_2^n$$
 (2)

3.8 Work-energy principle for a system of particles

$$W_{12} = T_2 - T_1$$

$$T := \sum_{i=1}^{n} \frac{1}{2} m_i |\underline{v}_i|^2$$

• Always check if the system is conservative. If so you can apply conservation of mechanical energy:

$$T_1 + V_1 = T_2 + V_2$$

- If the system is not conservative split the work done in a conservative and a non-conservative part.
- To calculate the non-conservative part of the work keep in mind the definition of work. Only the component of the force that is parallel to the velocity of its point of attack does work!

$$W_{12}^{nc} = \int_{P_1}^{P_2} \underline{F} \cdot \underline{dr} = \int_{P_1}^{P_2} |\underline{F}| \cos(\alpha) dS$$

• Check if the system is a rigid system of particles! If yes the work done by the internal forces is zero. \rightarrow 3.8.1, 3.9

3.8.1 RIGID SYSTEM OF PARTICLES

$$\frac{d}{dt} \left| \underline{r}_i - \underline{r}_j \right|^2 \equiv 0 \ \forall \ i, j = 1, \dots, n$$

 \underline{i} location of particle i in m number of particles

3.9 Work-energy principle for a rigid system of particles

The difference compared to the work-energy principle for a general system of particles is that in the case of a rigid system of particles the work done by the internal forces equals zero.

$$W_{12}^{ext} = T_2 - T_1$$

 $W_{12}^{\rm ext}$ total work done by external forces on the particles between point 1 and 2 in $J = kg\,m\,s^{-2}$ total kinetic energy of the system at t_i in $J = kg\,m\,s^{-2}$

4 RIGID BODY DYNAMICS

4.0.1 Connection vector

$$\underline{r}_{AB} = \underline{r}_B - \underline{r}_A$$

4.1 Velocity transfer formula for a rigid body

$$\underline{v}_B = \underline{v}_A + \underline{\omega} \times \underline{r}_{AB}$$

- Remember that angular velocities can be added up.
- It is not always necessary to do vector calculations. Clever usage of the SDPG (Theorem of projected velocities) and the instantaneous center of rotation can be a solution too.

4.1.1 Instantaneous center of rotation

The instantaneous center of rotation O of a planar rigid body \mathcal{B} is defined as the point $O \in \mathcal{B}'$ for which

$$\underline{v}_{O} = 0$$

Hence the extended rigid body \mathcal{B}' includes the original body \mathcal{B} and the instantaneous center of rotation O.

• If the instantaneous center of rotation and the angular velocity of a body is known the velocity of any point that is part of the body can be easily calculated.

4.1.2 Instantaneous axis of rotation

As seen in Mechanics I, in any moving rigid body there can be found an axis whose points only have translational velocity. This axis is called the instantaneous axis of rotation.

4.2 Center of mass of a rigid body ${\cal B}$

$$\underline{r}_C = \frac{1}{M} \int_{\mathcal{B}} \underline{r} dm = \frac{1}{M} \int_{\mathcal{B}} \underline{r} \rho dV$$

• Normally this integral can be split up into a simple sum.

4.3 Linear momentum principle for a rigid body ${\cal B}$

$$\underline{\dot{P}} = M\underline{a}_C = \underline{F}^{ext}$$

- If no resultant force is acting on the body, the acceleration of the center of mass is zero and the linear momentum of the system is conserved.
- If all external forces on a system are known the acceleration of the center of mass can be determined. Vice versa if the acceleration of the center of mass is known the sum of the external forces can be calculated.
- Since the linear momentum principle is a vectorial equation it can be split up and considered in each direction separately. This means also that it can be conserved in one direction while it changes in the other direction.
- Only valid in an inertial frame!

4.4 The angular momentum of a rigid body \mathcal{B} w.r.t. a fixed point B, belonging to the body, or w.r.t. its center of mass $B \equiv C$

$$\underline{H}_B = \int_{\mathcal{B}} \underline{\varrho}_B \times (\underline{\omega} \times \underline{\varrho}_B) dm = \underline{\underline{I}}_B \underline{\omega}$$

 $\begin{array}{llll} \underline{H}_B & \text{angular momentum of } \mathcal{B} \text{ w.r.t. } B \dots & \text{in} & \text{kg m}^2 \, \text{s}^{-1} \\ \underline{\varrho}_B & \text{distance from location of dm to point } B & \text{in} & \text{m} \\ \underline{\omega} & \text{rotational velocity of } \mathcal{B} \dots & \text{in} & \text{rad s}^{-1} \\ \underline{I}_B & \text{moment of inertia of } \mathcal{B} \dots & \text{in} & \text{kg m}^2 \end{array}$

- Remember that angular momenta w.r.t. the same point can be added up.
- Do not forget to use the angular momentum transfer formula whenever necessary!

4.5 Mass moments of inertia of a rigid body \mathcal{B}

 $\underline{\underline{I}}_{B}$ is the **moment of intertia tensor** of the body \mathcal{B} w.r.t. the point B.

$$\underline{\underline{I}}_B := \begin{pmatrix} I_{xx} & I_{xy} & I_{xz} \\ I_{xy} & I_{yy} & I_{yz} \\ I_{xz} & I_{yz} & I_{zz} \end{pmatrix}$$

where

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$$I_{xx} = \int_{\mathcal{B}} (y^2 + z^2) dm \qquad I_{xy} = -\int_{\mathcal{B}} xy dm \qquad I_{xz} = -\int_{\mathcal{B}} xz dm$$
$$I_{yy} = \int_{\mathcal{B}} (x^2 + z^2) dm \qquad I_{yz} = -\int_{\mathcal{B}} yz dm \qquad I_{zz} = \int_{\mathcal{B}} (x^2 + y^2) dm$$

- The mass moments of inertia are always given at the exam, either as an equation or directly as a value.
- Remember that moments of inertia formulated w.r.t. to the same point can simply be added up.
- Remember to apply Steiner's theorem wherever necessary.

4.6 Angular momentum principle for a rigid body ${\mathcal B}$ with respect to an arbitrary point B

$$\underline{\dot{H}}_B + \underline{v}_B \times \underline{P} = \underline{M}_B^{ext}$$

- $\underline{H}_B = \underline{\underline{I}}_B \cdot \underline{\omega}$ if and only if Point B is fixed or coincides with the center of mass!
- The angular momentum is conserved if and only if:
 - $-\underline{M}_{B}^{ext} \equiv 0$ AND either $\underline{v}_{B} \equiv 0$ or $\underline{P} \equiv 0$.
 - Alternatively it is possible that: $\underline{M}_{B}^{ext} \underline{v}_{B} \times \underline{P} \equiv 0$.
- When considering impulsive forces it makes sense to assess the change of angular momentum over an infinitesimally short time, therefore allowing negligence of nonimpulsive forces.
- When applying the angular momentum principle to a rolling cylinder it is possible to choose the contact point with the surface as a point of respect, since that way you get rid of the external moment of the contact force and $\underline{v_B} \times \underline{P}$, because in that case $\underline{v_B} \parallel \underline{v_C}$. However if the cylinder still slides consider the next tip:
- When applying the angular momentum principle in a point that is not the center of mass or that is not fixed you must use the angular momentum transfer formula!

4.7 Angular momentum transfer formula for a rigid body ${\cal B}$ between two arbitrary points A and B

$$\underline{H}_B = \underline{H}_A + \underline{P} \times \underline{r}_{AB}$$

 $\begin{array}{llll} \underline{H}_B & \text{angular momentum w.r.t. point } B \dots & \text{in} & \text{kg m}^2 \, \text{s}^{-1} \\ \underline{H}_A & \text{angular momentum w.r.t. point } A \dots & \text{in} & \text{kg m}^2 \, \text{s}^{-1} \\ \underline{P} & \text{linear momentum of } \mathcal{B} \dots & \text{in} & \text{kg m} \, \text{s}^{-1} \\ \underline{T}_{AB} & \text{connection vector between } A \text{ and } B \end{array}$

4.8 Parallel axis theorem (Steiner's Theorem)

$$\underline{\underline{I}}_{CO} = \underline{\underline{I}}_{CC} + M \left(|\underline{r}_{CO}|^2 \underline{\mathbb{I}}_{CC} - \underline{r}_{CO} \underline{r}_{CO}^T \right) = \underline{\underline{I}}_{CC} + M \begin{pmatrix} b^2 + c^2 & -ab & -ac \\ -ab & a^2 + c^2 & -bc \\ -ac & -bc & a^2 + b^2 \end{pmatrix}$$

$$\underline{\underline{I}}_{CC} \quad \text{moment of inertia tensor of } \mathcal{B} \text{ w.r.t. some point } O \dots \qquad \text{in} \quad \text{kg m}^2$$

$$\underline{\underline{I}}_{CC} \quad \text{moment of inertia tensor of } \mathcal{B} \text{ w.r.t. the center of mass of } \mathcal{B} \text{ in} \quad \text{kg m}^2$$

$$\underline{M} \quad \text{total mass of } \mathcal{B} \dots \qquad \text{in} \quad \text{kg}$$

$$\underline{r}_{CO} \quad \text{connection vector between } C \text{ and } O \dots \qquad \text{in} \quad \text{in} \quad \text{m}$$

$$\underline{\mathbf{I}} \quad \text{unity matrix in } \mathbb{R}^3 \dots \qquad \text{in} \quad \text{m}$$

4.9 Parallel axis theorem (Steiner's theorem) for planar bodies

$$I_A = I_C + M|\underline{r}_{CA}|^2$$

4.10 Kinetic energy T of a rigid body \mathcal{B}

$$T = \frac{1}{2}M|\underline{v}_C|^2 + \frac{1}{2}\underline{\omega}^T\underline{\underline{I}}_C\underline{\omega}$$

• For a fixed point B:

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$$T = \frac{1}{2} \underline{\omega} \underline{I}_B \underline{\omega}$$

4.11 Work-energy principle for a rigid body

$$W_{12}^{ext} = T_2 - T_1$$

 W_1^{ext} total work done by all forces on \mathcal{B} in $J = kg m^2 s^{-2}$ T_i total kinetic energy of the system at t_i in $J = kg m^2 s^{-2}$

4.12 Differentiation of quantities in moving frames

$$\underline{\dot{u}} = \underline{\mathring{u}} + \underline{\Omega} \times \underline{u}$$

- Differentiation in moving frame makes sense since aligning the chosen coordinate system with the principle axes of the body in question simplifies calculations. The vector-matrix multiplications are much simpler if you have the principal moments of inertia instead of a general inertia tensor.
- The rotational velocity of the moving frame can coincide with the rotational velocity of the body but in general does not have to.

4.12.1 Principal moments of inertia

Consider a basis $[\underline{u}_1 \quad \underline{u}_2 \quad \underline{u}_3]$ in which the moment of inertia becomes diagonal.

$$\underline{\underline{I}}_B = \begin{pmatrix} I_1 & 0 & 0 \\ 0 & I_2 & 0 \\ 0 & 0 & I_3 \end{pmatrix}$$

 $[\underline{u}_1 \quad \underline{u}_2 \quad \underline{u}_3]$ are then called the principal axes of \mathcal{B} and I_i are the principal moments of inertia

4.13 Euler's equations of motion for a spinning top

$$I_1\dot{\omega}_1 + (I_3 - I_2)\omega_3\omega_2 = M_1$$

$$I_2\dot{\omega}_2 + (I_1 - I_3)\omega_1\omega_3 = M_2$$

$$I_3\dot{\omega}_3 + (I_2 - I_1)\omega_2\omega_1 = M_3$$

 I_i i_{TH} principal moment of inertia . . . in kg m² $\dot{\omega}_i$ rotational acceleration in i-direction in rad s⁻² M_i torque on \mathcal{B} in i-direction in N m = kg m² s⁻²

- The Euler equations are nothing else than the AMP applied to a fixed reference point or the center of mass of a body. A rotating, principal frame is used.
- As a result, they are only valid for a fixed point or the center of mass of a body.
- They can only be applied if a body fixed, principal frame is used.
- You can modify the equations to get a general formulation if the angular velocity of the coordinate frame is not equal to the angular velocity of the body.

4.14 TSP-RULE

$$\underline{\omega}_{\varphi} \times \underline{\omega}_{\psi} \approx \frac{1}{I_3} \underline{M}_{O}^{ext}$$

 $\begin{array}{lllll} \underline{\omega}_{\varphi} & \text{precession} & & \text{in} & \text{rad s}^{-1} \\ \underline{\omega}_{\psi} & \text{spin} & & \text{in} & \text{rad s}^{-1} \\ \underline{M}_{O}^{ext} & \text{external torque acting on gyroscope} & \text{in} & \text{N m} = \text{kg m}^2 \, \text{s}^{-2} \end{array}$

- The TSP-Rule holds true for a fast spinning gyroscope.
- The external torque is aligned with the middle finger.
- Spin is aligned with the index finger.
- Precession axis is aligned with the thumb.

5 VIBRATIONS

5.1 Equation of motion of a dampened, forced one degree of freedom oscillator

$$\ddot{x} + 2\delta\dot{x} + \omega_0^2 x = f_0(t)$$

where

$$\delta = \frac{c}{2m} \qquad \omega_0 = \sqrt{\frac{k}{m}}$$

c damper constant in kg s⁻¹
m vibrating mass in kg δ characteristic damping in s⁻¹
k spring constant in kg s⁻² ω_0 natural frequency in rad s⁻¹

5.2 General solution for free vibrations (one DOF) with distinct roots $\lambda_1 \neq \lambda_2$

The solution in the time domain is

$$x(t) = A_1 e^{\lambda_1 t} + A_2 e^{\lambda_2 t}$$

while in the phase space is

$$\underline{\dot{x}}(t) = A_1 \underline{s}_1 e^{\lambda_1 t} + A_2 \underline{s}_2 e^{\lambda_2 t}$$

where $\lambda_{1,2} = -\delta \pm \sqrt{\delta^2 - \omega_0^2}$ and $\underline{s}_i = (1, \lambda_i)$

 $\begin{array}{lllll} \delta & \mbox{characteristic damping} & \mbox{in} & \mbox{s}^{-1} \\ \omega_0 & \mbox{natural frequency}.... & \mbox{in} & \mbox{rad}\,\mbox{s}^{-1} \end{array}$

5.3 General solution for free vibrations (one DOF) with repeated roots $\lambda=\lambda_1=\lambda_2$

The solution in the time domain is

$$x(t) = A_1 e^{\lambda t} + A_2 t e^{\lambda t}$$

while in the phase space is

$$\underline{\dot{x}}(t) = (A_1 \underline{s}_1 + A_2 \overline{\underline{s}}_1)e^{\lambda t} + A_2 \underline{s}_2 t e^{\lambda t}$$

where $\lambda_{1,2} = -\delta = -\omega_0$, $\underline{s}_i = (1, \lambda)$ and $\underline{\tilde{s}}_1 = (0, 1)$

5.4 Forced Harmonic Response

$$\ddot{x} + 2\delta \dot{x} + \omega_0^2 x = f_0 \sin(\omega t)$$

the steady state solution is

$$x_n(t) = A\sin(\omega t - \phi)$$

where

$$A = \frac{f_0}{\sqrt{(\omega_0^2 - \omega^2) + 4\delta\omega^2}} \quad \text{and} \quad \phi = \arctan\left(\frac{2\delta\omega}{\omega_0^2 - \omega^2}\right)$$

or in the form

$$x_p(t) = \frac{f_0}{\omega_0^2} V(\eta, D) \sin(\omega t - \phi(\eta, D))$$

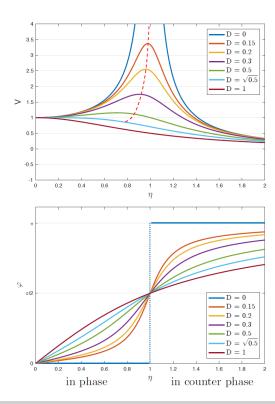
where

$$V = \frac{1}{\sqrt{(1 - \eta^2)^2 + (2D\eta)^2}} \quad \text{and} \quad \phi = \arctan\left(\frac{2D\eta}{1 - \eta^2}\right)$$

and

$$\eta = \frac{\omega}{\omega_0} \quad \text{and} \quad D = \frac{\delta}{\omega_0}$$

D Lehr's damping



$5.5\,$ Equation for vibrations with n degrees of freedom

$$\underline{M\ddot{x}} + \underline{C\dot{x}} + \underline{Kx} = \underline{F}(t)$$

 $\begin{array}{llll} \underline{x}(t) & \text{deviation from the stable equilibrium} & & \text{in} & \text{m} \\ \underline{\underline{M}} & \text{positive definite, symmetric mass matrix} & & \text{in} & \text{kg} \\ \underline{\underline{C}} & \text{positive semidefinite, symmetric damping matrix} & & \text{in} & \text{kg s}^{-1} \\ \underline{\underline{K}} & \text{positive semidefinite, symmetric stiffness matrix} & & \text{in} & \text{kg s}^{-2} \\ \underline{\underline{F}}(t) & \text{external force acting on the system} & & & \text{in} & \text{kg m s}^{-2} \end{array}$

5.6 STRUCTURAL DAMPING

$$\underline{C} = a\underline{M} + b\underline{K}$$

5.7 Generalized eigenvalue problem

$$(\lambda^2 \underline{\underline{M}} + \underline{\underline{K}})\underline{u} = \underline{0}$$

5.8 Purely imaginary eigenvalue pairs

 $\lambda = \pm i \ \omega_j, \ \omega_j > 0, \ \underline{u}_j \in \mathbb{R}^n$

The solution in the time domain is

$$\underline{x}_{i}(t) = \underline{u}_{i} A_{j} \sin(\omega_{j} t + \phi_{j})$$

5.9 Zero eigenvalue pairs

 $\lambda = 0^2, \ u \in \mathbb{R}^n$

The solution in the time domain is

$$\underline{x}_j(t) = u_j(B_j + C_j t)$$

5.10 Modal decomposition of free, undampened vibrations

$$\underline{\underline{U}}^T \underline{\underline{M}} \underline{\underline{U}} = \operatorname{diag}(m_1, m_2, \dots, m_n)$$
(3)

$$\underline{\underline{U}}^T \underline{K} \underline{U} = \operatorname{diag}(k_1, k_2, \dots, k_n)$$
(4)

5.11 Modal equation of motion

$$m_j \ddot{y}_j + c_j \dot{y}_j + k_j y_j = f_j(t), \ j = 1, \dots n$$
 $c_j = a m_j + b k_j \ge 0$

5.12 Forced response to harmonic forcing of a n degrees of freedom system

For harmonic forcing

$$\underline{F}(t) = \underline{F}_0 \sin(\omega t)$$

the steady-state solution of a generic modal amplitude $[y_P(t)]_i$ is given by

$$[y_P(t)]_j = A_i^P \sin(\omega t + \phi_j)$$

with

$$A_j^P = \frac{f_j^0}{\sqrt{(\omega_j^2 - \omega^2)^2 + c_j^2 \omega^2}}, \quad \phi_j = \arctan\left(\frac{c_j \omega}{\omega_j^2 - \omega^2}\right) \quad \text{and} \quad f_j^0 = (\underline{\underline{U}}^T \underline{F}_0)_j$$

- Note that this solution is the particular solution to $\underline{\underline{M}}\underline{\ddot{x}} + \underline{\underline{C}}\underline{\dot{x}} + \underline{\underline{K}}\underline{x} = \underline{F}(t)$, transformed to modal coordinates such that the system is decoupled.
 - Like that the system is described by a series of one DOF oscillators, one for each modal coordinate. For those the solutions for forced, damped vibrations are applicable again.
- Each mode has its resonance frequency $\omega_j = \sqrt{\frac{k_j}{m_j}}$.
- When the external forcing is given as above we only care about the particular solution as given above, since the transient response defined by the homogeneous solution decays after some time.
- Remember that from modal coordinates one needs to transform back to the original x-coordinate frame.

$$\underline{x}_p = \underline{U}\underline{y}_p(t) = \begin{pmatrix} \underline{u}_1 & \underline{u}_1 & \cdots & \underline{u}_n \end{pmatrix} \begin{pmatrix} A_1^P \sin(\omega t + \phi_1) \\ \vdots \\ A_n^P \sin(\omega t + \phi_n) \end{pmatrix}$$

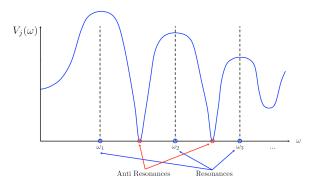
which can be written as:

$$\underline{x}_p(t) = \sum_{k=1}^n A_k^P \underline{u}_k \sin(\omega t + \phi_k)$$

Based on that an amplification factor can be established where $B_j(\omega) = \sum_{k=1}^n A_k^P(\underline{u}_k)_j$ is the amplitude of the particular response of the j^{th} degree of freedom.

$$V_j(\omega) = \frac{|B_j|}{|B_j(0)|}$$

That is the relation of the amplification at any frequency ω w.r.t. the amplification at a frequency of $\omega = 0$.



• As seen in figure ?? the amplification factor has maxima at each of the modal resonance frequencies and for the example shown here it has anti resonances between the maxima. However this depends on the underlying system. See Exercise 13.3 for a counterexample.

6 Analytic Dynamics

6.0.1 Generalized Coordinates

The generalized coordinates are variables required to define the position of a body uniquely. They are not unique themselves.

6.0.2 Constraints

Definition 1. A constraint is a scalar function of generalized coordinates, generalized velocities and time. This constraint is non-holonomic. Also accelerations, and higher derivatives?

$$f(\underline{q},\underline{\dot{q}},t) = 0$$

Definition 2. A sceleronomic constraint is only a function of generalized coordinates.

$$f(\underline{q}) = 0$$

Definition 3. A **rheonomic** constraint is only a function of generalized coordinates and time.

$$f(\underline{q},t)$$

Definition 4. If non-holonomic constraints can be integrated, they will become holonomic and also eliminate DoFs.

Every holonomic constraint can be given in **Pfaffian form** (exact differential):

$$f(\underline{q},t) = 0 \quad \to \quad df = \sum_{i=1}^{n} \frac{\partial f(\underline{q},t)}{\partial q_i} + \frac{\partial f(\underline{q},t)}{\partial t} dt = 0$$

Definition 5. It can also be given in **velocity form** by dividing by dt:

$$\frac{df}{dt} = \sum_{i=1}^{n} \frac{\partial f(\underline{q},t)}{\partial \dot{q}_{i}} + \frac{\partial f(\underline{q},t)}{\partial t} = 0$$

Definition 6. In practice many constraints are given in a linear velocity form:

$$\sum_{i=1}^{n} a_i(\underline{q}, t)\dot{q}_i + b(\underline{q}, t) = 0$$

6.0.3 Showing that a constraint is holonomic

A constraint in **velocity form** is holonomic if its coefficients a_i and b_i can be mapped to a **pfaffian form**. If so then the velocity form is an exact differential

$$\frac{df}{dt} = \sum_{i=1}^{n} a_i(\underline{q}, t)\dot{q}_i + b(\underline{q}, t)$$

What remains to be showed is that:

$$\frac{\partial f(\underline{q},t)}{\partial a_i} = Ca_i \quad , \quad \frac{\partial f(\underline{q},t)}{\partial t} = Cb, \quad \text{where } C = C(\underline{q},t)$$

The condition for partial derivatives to be integrable

$$\frac{\partial^2 f(\underline{q},t)}{\partial a_k \partial a_i} = \frac{\partial^2 f(\underline{q},t)}{\partial a_i \partial a_k} \Rightarrow \frac{\partial C a_i}{\partial a_k = \partial C a_i} \partial q_i \quad ; \quad \frac{\partial C b}{\partial a_k} = \frac{\partial C a_i}{\partial t}$$

If C = 0 the integration constant does not exist, the trivial solution of the problem is found. Thus the constraint is non-holonomic.

6.1 Virtual displacements

6.1.1 Sceleronomic constraint

$$\hat{\underline{r}}(t) = q_1 \hat{\underline{e}}_1 + \dots + q_n \hat{\underline{e}}_n$$
 Actual path

$$\underline{\tilde{r}}(t) = \tilde{q}_1 \underline{\hat{e}}_1 + \dots + \tilde{q}_n \underline{\hat{e}}_n$$
 Variational path

$$\delta \underline{\hat{r}} = \underline{\tilde{r}} - \underline{\hat{r}} = \delta q_1 \underline{\hat{e}}_1 + \dots + \delta q_n \underline{\hat{e}}_n$$
 Virtual displacement

- The virtual displacement is the difference between the actual and the virtual path.
- It is equivalent to freezing time and moving to a possible, close path.
- Possible means that all constraints are satisfied, i.e. $\tilde{r}(t)$ and $\hat{r}(t)$ must lie on f(q) = 0.

$$\boxed{d\hat{\underline{r}} = \frac{\partial \hat{\underline{r}}}{\partial q_1} dq_1 + \dots + \frac{\partial \hat{\underline{r}}}{\partial q_n} dq_n} \text{ Infinitesimal displacement}$$

$$\boxed{\hat{\underline{v}} = \frac{\partial \hat{\underline{r}}}{\partial q_1} \dot{q}_1 + \dots + \frac{\partial \hat{\underline{r}}}{\partial q_n} \dot{q}_n} \text{ Generalized velocity}$$

• $d\hat{r}$ is tangent to the trajectory, $\delta \hat{r}$ is not.

$$\boxed{\underline{n} = \frac{\nabla f}{|\nabla f|} = \frac{1}{|\nabla f|} \left(\frac{\partial f}{\partial q_1} \hat{\underline{e}}_1 + \dots + \frac{\partial f}{\partial q_n} \hat{\underline{e}}_2 \right)} \text{ Normal vector}$$

• Both are perpendicular to the gradient of the constraint plane. Thus the following conditions must hold.

$$\nabla f \cdot d\underline{\hat{r}} = 0 \Rightarrow \frac{\partial f}{\partial q_1} dq_1 + \dots + \frac{\partial f}{\partial q_1} dq_n = 0$$
$$\nabla f \cdot \partial \underline{\hat{r}} = 0 \Rightarrow \frac{\partial f}{\partial q_1} \delta q_1 + \dots + \frac{\partial f}{\partial q_n} \delta q_n = 0$$

6.1.2 Rheonomic constraint

- $d\hat{r}$ and \hat{v} do not lay on the constraint surface $\nabla f \cdot d\hat{r} \neq 0$
- $\delta \hat{\underline{r}}$ does lay on the tangent plane $\nabla f \cdot \delta \hat{\underline{r}} = \delta f = 0$

6.1.3 Kinematical method to compute virtual displacements

Recall the definition of velocity:

$$\underline{v} = \frac{d\underline{r}}{dt} = \sum_{i=1}^{n} \frac{\partial \underline{r}}{\partial q_i} \dot{q}_i + \frac{\partial \underline{r}}{\partial t}$$

and of virtual displacements:

$$\delta \underline{r} = \sum_{i=1}^{n} \frac{\partial \underline{r}}{\partial q_i} \delta q_i$$

- 1. Take the velocity formula.
- 2. Identify the linear terms in \dot{q}_i .
- 3. Discard the rest.
- 4. Replace \dot{q}_i with δq_i .

Use this method when only the velocity is known.

6.2 D'ALEMBERTS PRINCIPLE

6.2.1 Reaction forces

A constraint force keeps the system on the surface f(q,t) = 0 in the configuration space.

$$\boxed{\underline{R} = R\underline{n}}$$
 Constraint force

Definition 7. The virtual work is the work done by a force onto a virtual displacement.

The virtual work of a constraint force is zero.

Definition 8. An ideal constraint is such that the associated constraint force does not do work along virtual displacements compatible with the constraint.

6.2.2 Projected Newton's equations

1. Assume a system of N particles of mass m_i , subjected to holonomic constraints. The LMP writes:

$$m_i \ddot{\underline{r}}_i = \underline{F}_i + \underline{R}_i$$

2. Choose generalized coordinates q compatible with the constraints (admissible):

$$\underline{r}_i = \underline{r}_i(q,t)$$

and the corresponding virtual displacement:

$$\delta \underline{r}_i = \frac{\partial \underline{r}_i}{\partial \underline{q}} \delta \underline{q}$$

3. Bring all terms of the LMP to the LHS (dynamic equilibrium)

$$m_i \ddot{\underline{r}}_i - \underline{F}_i - \underline{R}_i = \underline{0}$$

4. Project the dynamic equilibrium onto the virtual displacement:

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

5. Reaction forces do not do work on an admissible virtual displacement:

$$\sum_{i=1}^{N} \delta \underline{r}_i \cdot \underline{R}_i = 0$$

$$\sum_{i=1}^{N} \delta \underline{r}_i \cdot (m_i \underline{\ddot{r}}_i - \underline{F}_i) = 0$$
 D'alembers principle

5. Recall the definition of the virtual displacements

$$\delta \underline{r}_i = \frac{\partial \underline{r}_i}{\partial \underline{q}} \delta \underline{q} = \sum_{k=1}^p \frac{\partial \underline{r}_i}{\partial q_k} \delta q_k$$

6. And insert it into d'alemberts principle

$$\sum_{i=1}^{N} (m_i \underline{\ddot{r}}_i - \underline{F}_i) \cdot \sum_{k=1}^{p} \frac{\partial \underline{r}_i}{\partial q_k} \delta q_k = 0$$

7. Since all general coordinates can be varied independently (admissible) we can write p independent equations:

$$\left| \sum_{i=1}^{N} (m_i \underline{\ddot{r}}_i - \underline{F}_i) \cdot \frac{\partial \underline{r}_i}{\partial q_k} = 0, \quad \forall k = 1, \cdots, p \right|$$

6.2.3 Generalized forces

Definition 9. The projection of the active forces onto the admissible direction dictated by the variation of q_k is called **generalized force**.

$$Q_k = \sum_{i=1}^{N} \underline{F}_i \cdot \frac{\partial \underline{r}_i}{\partial q_k}$$

6.3 Lagrange equations

Definition 10. Consider a system of N particles. The **kinetic energy** is defined as:

$$\mathcal{T} = \sum_{i=1}^{N} \frac{1}{2} m_i \dot{\underline{r}}_i \cdot \dot{\underline{r}}_i = \sum_{i=1}^{N} \frac{1}{2} m_i |\dot{\underline{r}}_i|^2$$

Note that:

$$\frac{d}{dt} \left(\frac{\partial \mathcal{T}}{\partial \dot{r}_i} \right) = m_i \ddot{\underline{r}}_i$$

And therefore d'Alembert's priciple can be written as:

$$\textstyle\sum_{i=1}^{N}\left[\frac{d}{dt}\left(\frac{\partial\mathcal{R}}{\partial\dot{\underline{r}}_{i}}\right)-\underline{F}_{i}\right]\cdot\delta\underline{r}_{i}=0\Rightarrow\sum_{k=1}^{p}\sum_{i=1}^{N}\left[\frac{d}{dt}\left(\frac{\partial\mathcal{T}}{\partial\dot{\underline{r}}_{i}}\right)-\underline{F}_{i}\right]\cdot\frac{\partial\underline{r}_{i}}{\partial q_{k}}\delta q_{k}=0$$

If the GC's are kinematically admissible, then

$$\sum_{i=1}^{N} \left[\frac{d}{dt} \left(\frac{\partial \mathcal{T}}{\partial \dot{\underline{r}}_i} \right) - \underline{F}_i \right] \cdot \frac{\partial \underline{r}_i}{\partial q_k} = 0 \quad k = 1, \dots, p$$

Note that:

$$\frac{d}{dt} \left(\frac{\partial \mathcal{T}}{\partial \dot{\underline{r}}_i} \cdot \frac{\partial \underline{r}_i}{\partial q_k} \right) = \frac{d}{dt} \left(\frac{\partial \mathcal{T}}{\partial \dot{\underline{r}}_i} \cdot \frac{\partial \underline{r}_i}{\partial q_k} \right) + \frac{\partial \mathcal{T}}{\partial \dot{\underline{r}}_i} \cdot \frac{d}{dt} \left(\frac{\partial \underline{r}_i}{\partial q_k} \right)$$

and therefore:

$$\frac{d}{dt} \left(\frac{\partial \mathcal{T}}{\partial \dot{\underline{r}}} \right) \cdot \frac{\partial \underline{r}_i}{\partial q_k} = -\frac{\partial \mathcal{T}}{\partial \dot{\underline{r}}_i} \cdot \frac{d}{dt} \left(\frac{\partial \underline{r}_i}{\partial q_k} \right) + \frac{d}{dt} \left(\frac{\partial \mathcal{T}}{\partial \dot{\underline{r}}_i} \cdot \frac{\partial \underline{r}_i}{\partial q_k} \right)$$

Substitution yields:

$$\sum_{i=1}^{N} \left[-\frac{\partial \mathcal{T}}{\partial \dot{\underline{r}}_{i}} \cdot \frac{d}{dt} \left(\frac{\partial \underline{r}_{i}}{\partial q_{k}} \right) + \frac{d}{dt} \left(\frac{\partial \mathcal{T}}{\partial \dot{\underline{r}}_{i}} \cdot \frac{\partial \underline{r}_{i}}{\partial q_{k}} \right) - \underline{F}_{i} \cdot \frac{\partial \underline{r}_{i}}{\partial q_{k}} \right] = 0 \quad k = 1, \dots, p$$

Now consider red:

$$\frac{\partial \mathcal{T}}{\partial \dot{\underline{r}}_i} \cdot \frac{d}{dt} \left(\frac{\partial \underline{r}_i}{\partial q_k} \right) = \frac{\partial \mathcal{T}}{\partial \dot{\underline{r}}_i} \cdot \frac{\partial \dot{\underline{r}}_i}{\partial q_k} = \frac{\partial \mathcal{T}}{\partial q_k}$$

Then consider blue:

$$\underline{\dot{r}}_i = \sum_{i=1}^p \frac{\partial \underline{r}_i}{\partial q_k} \dot{q}_k + \frac{\partial \underline{r}_i}{\partial t} \quad \Rightarrow \quad \frac{\partial \dot{\underline{r}}_i}{\partial \dot{q}_k} = \frac{\partial \underline{r}_i}{\partial q_k}$$

and therefore

$$\frac{\partial \mathcal{T}}{\partial \dot{\underline{r}}_i} \cdot \frac{\partial \underline{r}_i}{\partial q_k} = \frac{\partial \mathcal{T}}{\partial \dot{\underline{r}}_i} \cdot \frac{\partial \dot{\underline{r}}_i}{\partial \dot{q}_i} = \frac{\partial \mathcal{T}}{\partial q_k}$$

$$\frac{d}{dt} \left(\frac{\partial \mathcal{T}}{\partial \dot{q}_k} \right) - \frac{\partial \mathcal{T}}{\partial q_k} - Q_k = 0, \quad k = 1, \dots, p$$

- LEs are nothing else than d'Alembert's principle, written in terms of kinetic energies.
- As such, the chosen generalized coordinates are supposed to identically satisfy the constraints (i.e. the are unconstrained coordinates.
- By working with energy, the inertial terms involve less algebra than the d'Alembert's principle.

6.3.1 Potential active forces

If all active forces are potential:

$$\underline{F}_i = -\frac{\partial \mathcal{V}}{\partial \underline{r}_i} + \underline{F}_i^{nc}$$

Therefore the generalized forces become:

$$Q_k = \sum_{i=1}^N \underline{F}_i \cdot \frac{\partial_{\underline{r}_i}}{\partial q_k} = -\sum_{i=1}^N \left[\frac{\partial \mathcal{V}}{\partial \underline{r}_i} \cdot \frac{\partial \underline{r}_i}{\partial q_k} + \underline{F}_i^{nc} \cdot \frac{\partial \underline{r}_i}{\partial q_k} \right] = -\frac{\partial \mathcal{V}}{\partial q_k} + Q_k^{nc}$$

And then the LEs could be written as:

$$\frac{d}{dt} \left(\frac{\partial \mathcal{T}}{\partial \dot{q}_k} \right) - \frac{\partial \mathcal{T}}{\partial q_k} + \frac{\partial \mathcal{V}}{\partial q_k} - Q_k^{nc} = 0, \quad k = 1, \dots, p$$

$$\frac{d}{dt} \left(\frac{\partial \mathcal{T}}{\partial \underline{\dot{q}}} \right) - \frac{\partial \mathcal{T}}{\partial \underline{q}} + \frac{\partial \mathcal{V}}{\partial \underline{q}} = 0$$

- 6.3.2 Lagrange equations for constrained coordinates
- 6.4 Hamiltons principle
- 6.5 Non-holonomic dynamics

7 Linearization

- 7.1 Equilibrium and stability
- 7.2 Mass, stiffness and damping matrices
- 7.3 Imposed motion

8 Vibrations

- 8.1 Vibration modes, Mode orthogonality
- 8.2 Modal superposition
- 8.3 Frequency response function
- 8.4 Damped Vibrations

9 Advanced 3D Rigid Body Dynamics

- 9.1 Parametrization of rotations
- 9.2 Euler equations
- 9.3 Lagrangian Dynamics for 3D multibody systems