

# Internal Coordinates

If the potential energy of a system depends only on the internal coordinates of the system, then the motion of the centre of mass can always be separated from the internal motion.

Consider two point masses  $m_1$  and  $m_2$ , both in motion and interacting with each other.

$$E = \frac{1}{2}m_1(\dot{x}_1^2 + \dot{y}_1^2 + \dot{z}_1^2) + \frac{1}{2}m_2(\dot{x}_2^2 + \dot{y}_2^2 + \dot{z}_2^2) + V(x_1, y_1, z_1, x_2, y_2, z_2)$$

Define centre of mass coordinates:

$$X = \frac{m_1x_1 + m_2x_2}{m_1 + m_2} \quad Y = \frac{m_1y_1 + m_2y_2}{m_1 + m_2} \quad Z = \frac{m_1z_1 + m_2z_2}{m_1 + m_2}$$

and internal coordinates:

$$x = x_1 - x_2 \quad y = y_1 - y_2 \quad z = z_1 - z_2$$

then

$$E = \underbrace{\frac{1}{2}(m_1 + m_2)(\dot{X}^2 + \dot{Y}^2 + \dot{Z}^2)}_{\text{translational energy}} + \underbrace{\frac{1}{2}\mu(\dot{x}^2 + \dot{y}^2 + \dot{z}^2) + V(x, y, z)}_{\text{internal energy}}$$

where **reduced mass**  $\mu = \frac{m_1m_2}{m_1 + m_2}$

In terms of wave functions:

$$\Psi = \psi_{\text{trans}}(X, Y, Z) \cdot \psi_{\text{int}}(x, y, z)$$

# Rotation as Translational Motion

For a particle of mass  $m$  moving in 3-D space

$$\begin{aligned}\hat{H} = \hat{T} + \hat{V} &= -\frac{\hbar^2}{2m} \nabla^2 + V(r, \theta, \phi) \\ &= -\frac{\hbar^2}{2m} \left( \frac{\partial^2}{\partial r^2} + \frac{2}{r} \frac{\partial}{\partial r} + \frac{1}{r^2} \Lambda^2 \right) + V(r, \theta, \phi)\end{aligned}$$

where  $\Lambda^2 = \frac{\partial^2}{\partial \theta^2} + \frac{\cos \theta}{\sin \theta} \frac{\partial}{\partial \theta} + \frac{1}{\sin^2 \theta} \frac{\partial^2}{\partial \phi^2}$  Legendrian

Suppose the particle is confined to the surface of a sphere, i.e.  $r = R$ .

$$\hat{H} = -\frac{\hbar^2}{2mR^2} \Lambda^2 \quad \text{a function of } \theta \text{ and } \phi \text{ only}$$

A **rigid rotator** is a pair of masses at a fixed distance apart ( $R$ ), freely rotating about the **centre of mass**.

$$\begin{aligned}\hat{H} &= -\frac{\hbar^2}{2\mu R^2} \Lambda^2 & \mu &= \frac{m_1 m_2}{m_1 + m_2} \\ &= -\frac{\hbar^2}{2I} \Lambda^2 & I &= \mu R^2\end{aligned}$$

If the particle is confined to a ring (the equator),  $\theta = \pi/2$ .

$$\hat{H} = -\frac{\hbar^2}{2I} \frac{d^2}{d\phi^2} = -\frac{\hbar^2}{2\mu} \frac{d^2}{ds^2} \quad \text{s is the distance along the circumference}$$

# The Particle on a Ring

The Schrödinger Equation looks like that of the free particle, so the solutions are similar:

$$\Phi_m = \frac{1}{\sqrt{2\pi}} e^{im\phi} \quad E_m = \frac{\hbar^2}{2I} m^2 \quad m = 0, \pm 1, \pm 2, \dots$$

Quantization is due to a **cyclic boundary**

$$\Phi(\phi) = \Phi(\phi + 2m\pi)$$

**condition:**

Except for  $m = 0$  the states are twofold **degenerate**.

Real functions can be constructed by taking **linear**

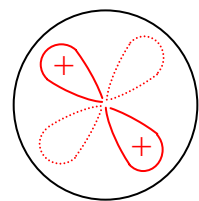
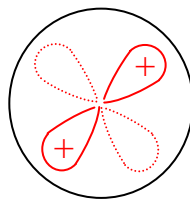
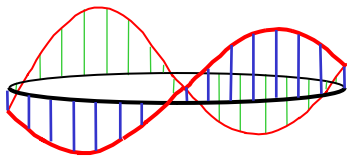
**combinations:**

$$\Phi_m^+ = \frac{1}{\sqrt{2}} [\Phi_m + \Phi_{-m}] = \frac{1}{\sqrt{\pi}} \cos(|m|\phi)$$

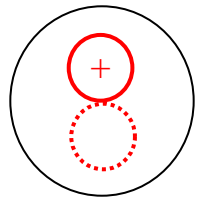
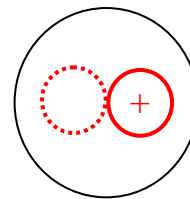
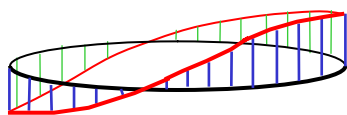
$$\Phi_m^- = \frac{-i}{\sqrt{2}} [\Phi_m - \Phi_{-m}] = \frac{1}{\sqrt{\pi}} \sin(|m|\phi)$$

$m$

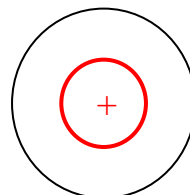
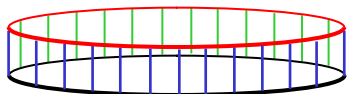
$\pm 2$



$\pm 1$



0



# The Particle on a Sphere

$$\Lambda^2 \Psi(\theta, \phi) = -\frac{2I}{\hbar^2} E \Psi(\theta, \phi)$$

This type of equation is well known (to applied mathematicians):

$$\Lambda^2 Y_{lm}(\theta, \phi) = -l(l+1) Y_{lm}(\theta, \phi) \quad \begin{cases} l = 0, 1, 2, \dots \\ m = 0, \pm 1, \pm 2, \dots, \pm l \end{cases}$$

The solutions are the spherical harmonics:

$$Y_{lm}(\theta, \phi) = \frac{1}{\sqrt{2\pi}} \Theta_{lm}(\theta) e^{im\phi}$$

where  $\Theta_{lm}$  are the associated Legendre polynomials.

$l$	$m$	$\Theta_{lm}$
0	0	$\sqrt{1/2}$
1	0	$\sqrt{3/2} \cos \theta$
1	$\pm 1$	$\sqrt{3/4} \sin \theta$
2	0	$\sqrt{5/8} (3 \cos^2 \theta - 1)$
2	$\pm 1$	$\sqrt{15/4} \sin \theta \cos \theta$
2	$\pm 2$	$\sqrt{15/16} \sin^2 \theta$

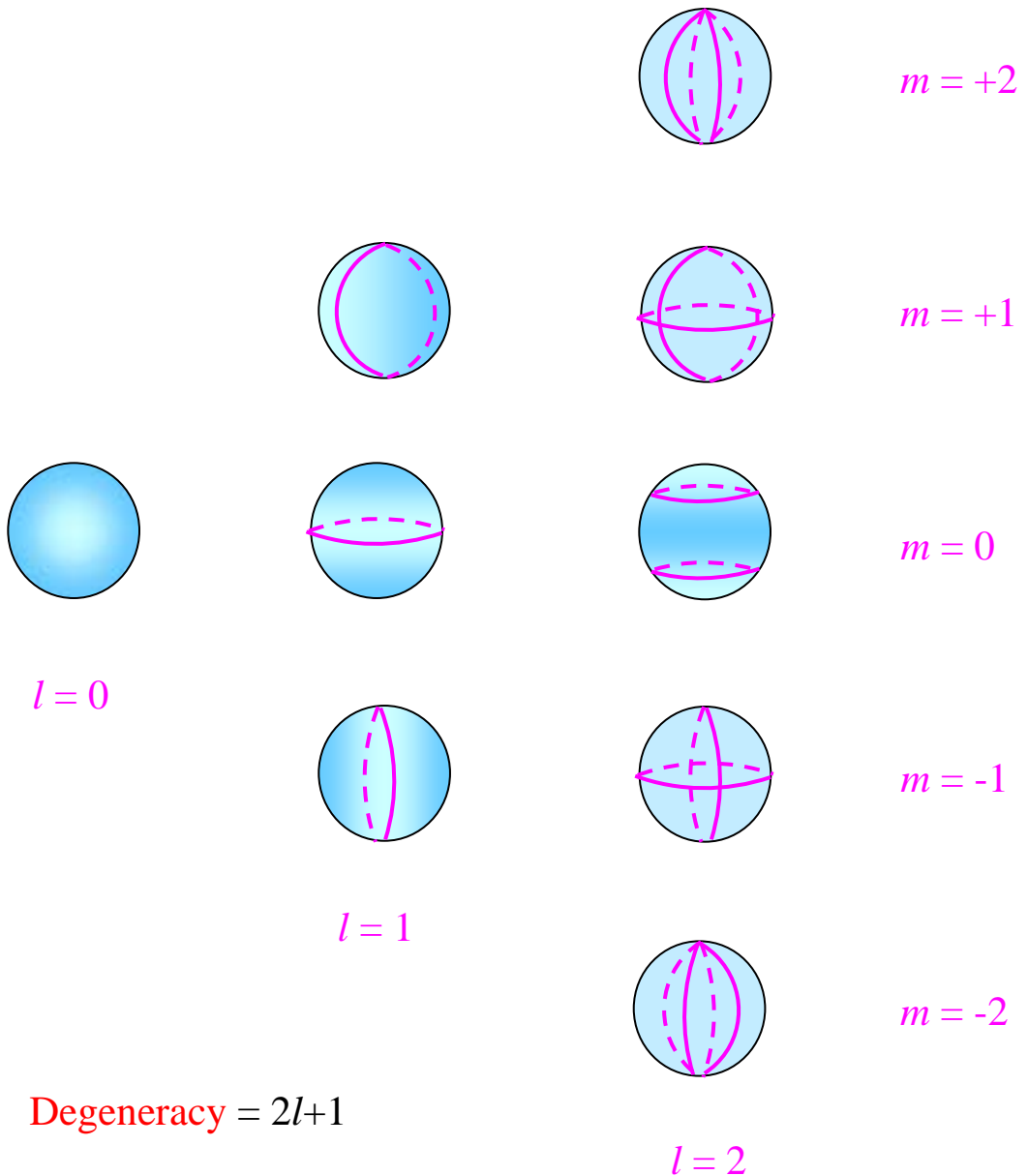
Comparing equations,

$$E_{lm} = \frac{\hbar^2}{2I} l(l+1)$$

# Spherical Harmonics: Real Wavefunctions

$$Z_{l,m}^+ = (1/\sqrt{2})[Y_{l,m} + Y_{l,-m}] = (1/\sqrt{\pi})\cos(|m|\phi)\Theta_{lm}(\theta)$$

$$Z_{l,m}^- = (-i/\sqrt{2})[Y_{l,m} - Y_{l,-m}] = (1/\sqrt{\pi})\sin(|m|\phi)\Theta_{lm}(\theta)$$



# Rotational/Orbital Angular Momentum

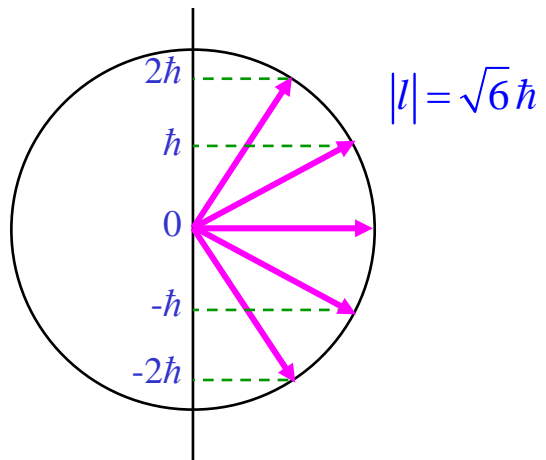
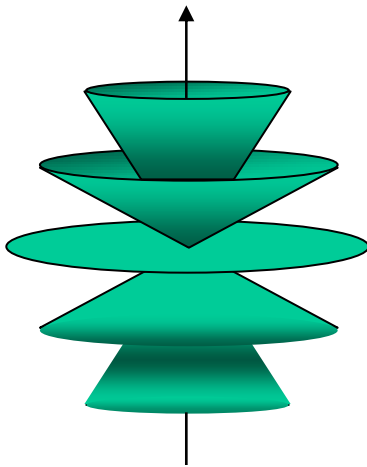
$$E_{lm} = \frac{\hbar^2}{2I} l(l+1) \quad \begin{cases} l = 0, 1, 2, \dots \\ m = 0, \pm 1, \pm 2, \dots, \pm l \end{cases}$$

The energy of a rotating body (or particle in orbit) is quantized.

There are  $(2l+1)$  degenerate states which have the same energy determined by the quantum number  $l$ .

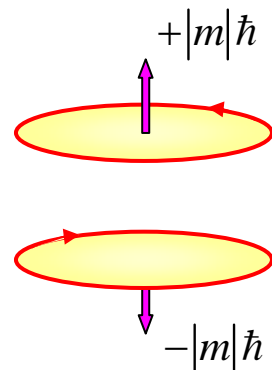
The different states, labelled by quantum numbers  $m$ , are related by simple symmetry transformations, i.e. they correspond to different orientations in space.

The orientation of a rotating body is quantized.



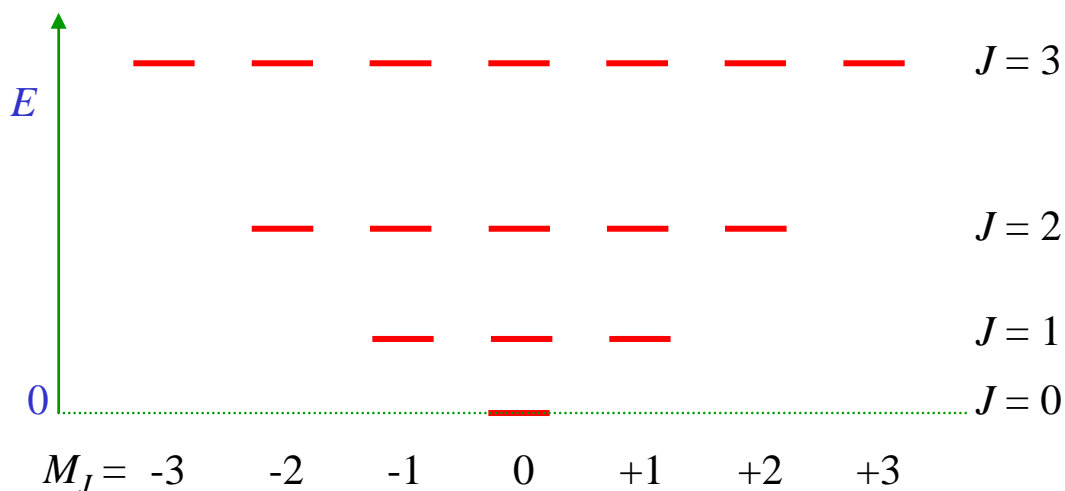
For the particle on a ring

$$E_m = \frac{\hbar^2}{2I} m^2 \quad m = 0, \pm 1, \pm 2, \dots$$



# Rotational Spectra of Diatomic Molecules

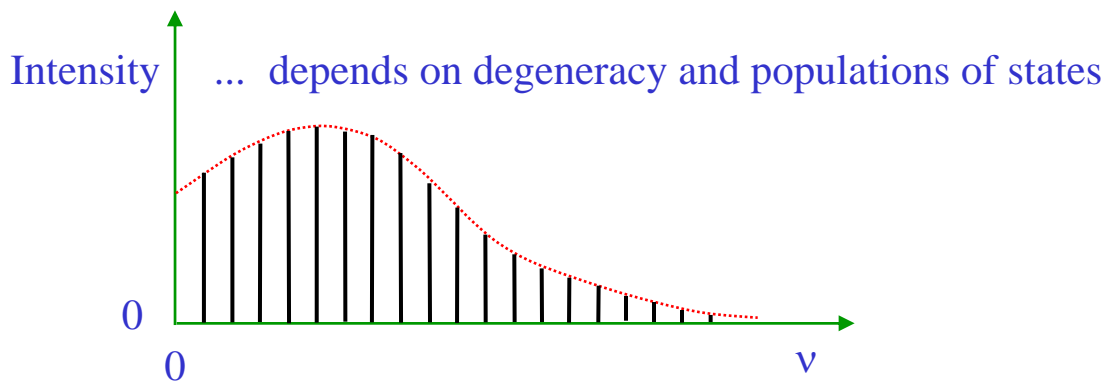
$$E_J = B J(J + 1) \quad B = \frac{\hbar^2}{2I} \quad J = 0, 1, 2, \dots$$



**Selection Rules:**  $\Delta J = \pm 1$        $\Delta M_J = 0, \pm 1$

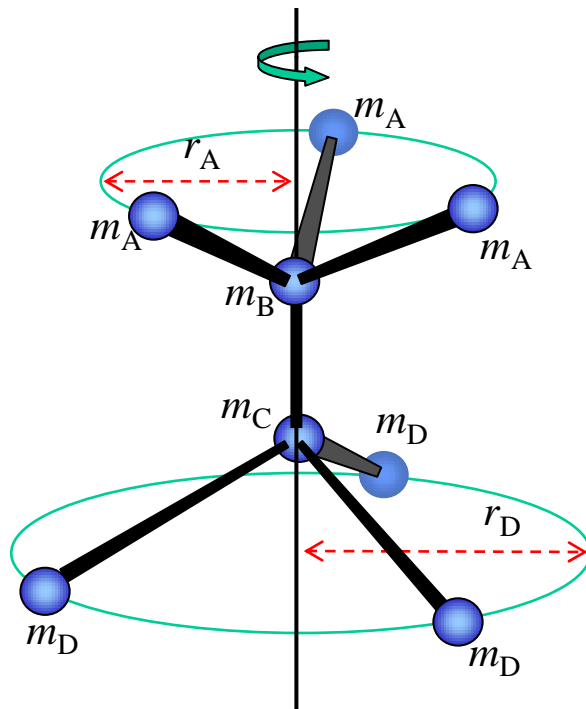
and the molecule must have a dipole moment.

**Transitions**  $\Delta E = E_{J+1} - E_J = 2B(J + 1) = 2B, 4B, 6B, \dots$



# The Moment of Inertia of a Rotating Molecule

$$I = 3m_A r_A^2 + 3m_D r_D^2$$





# Moments of Inertia – Principal Axes

Consider a molecule as a system of point masses whose positions are fixed relative to each other.

Centre of gravity:

$$\vec{r}_0 = \frac{\sum_k m_k \vec{r}_k}{\sum_k m_k}$$

Put a Cartesian coordinate system at this centre and define the three moments of inertia.

$$I_x = \sum_k m_k r_{kx}^2 \text{ etc.}$$

$r_{kx}$  is the perpendicular distance of nucleus  $k$  from the  $x$  axis

If  $I_{xy} = m_k r_{kx} r_{ky} \neq 0 \text{ etc.}$

rotate the coordinate system until  $I_{x'y'} = I_{y'z'} = I_{z'x'} = 0 \text{ etc.}$

It is always possible to find unique principal axes and thus calculate principal moments of inertia ( $I_a I_b I_c$ ).

Linear Rotator  $I_a = I_b \neq 0 \quad I_c = 0$

Spherical Top  $I_a = I_b = I_c \quad I_c \neq 0$

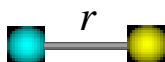
Symmetric Top  $I_a = I_b \neq I_c \quad I_c \neq 0$

prolate top  $I_a = I_b > I_c$

oblate top  $I_a = I_b < I_c$

Asymmetric Top  $I_a \neq I_b \neq I_c$

## Linear Molecules



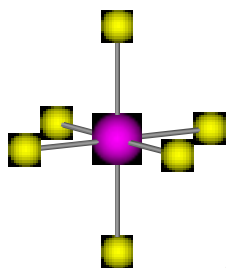
$$I = \frac{m_1 m_2}{m_1 + m_2} r^2 = \mu r^2$$



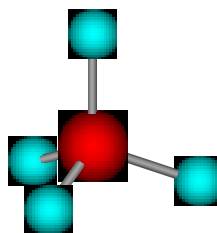
$$I = \frac{m_1 m_2 r_{12}^2 + m_2 m_3 r_{23}^2 + m_1 m_3 r_{13}^2}{m_1 + m_2 + m_3}$$

$$= 2m_1 r_{12}^2 \quad \text{if } m_1 = m_3$$

## Spherical Tops

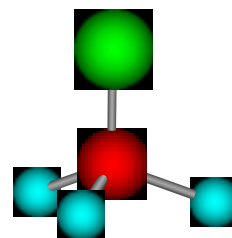
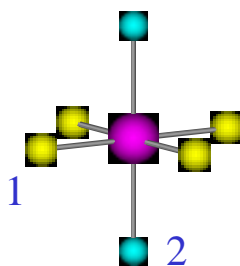
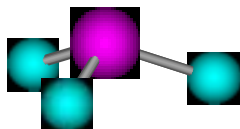


$$I = 4m_1 r^2$$



$$I = \frac{8}{3} m_1 r^2$$

## Symmetric Tops



$$I_{\parallel} = 4m_1 r_1^2$$

$$I_{\perp} = 2m_1 r^2 + 2m_2 r_2^2$$