

Small Oscillation (contd.-)

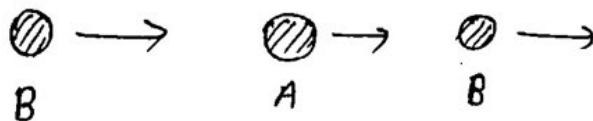
Let us now discuss the eigen vectors for three modes of vibration.

(i) For $\omega_1 = 0$

$$\begin{pmatrix} K & -K & 0 \\ -K & 2K & -K \\ 0 & -K & K \end{pmatrix} \begin{pmatrix} a_{11} \\ a_{21} \\ a_{31} \end{pmatrix} = 0$$

on solving, $a_{11} = a_{21} = a_{31} = \alpha$ (say)
so, the eigen vector is given by

$$a_1 = \begin{pmatrix} \alpha \\ \alpha \\ \alpha \end{pmatrix} \quad \text{--- (6)}$$



(Antisymmetric mode)

i-p. for $\omega_1 = 0$, all the molecules suffer same displacement in same direction.

(ii) For $\omega_2 = \sqrt{k/m}$

$$\begin{pmatrix} 0 & -K & 0 \\ -K & 2K - \frac{MK}{m} & -K \\ 0 & -K & 0 \end{pmatrix} \begin{pmatrix} a_{12} \\ a_{22} \\ a_{32} \end{pmatrix} = 0$$

on solving we have

$$a_{22} = 0$$

$$-a_{12} - a_{32} = 0$$

Therefore, $a_{22} = 0$, $a_{12} = -a_{32} = \beta$ (say)

Thus, for $\omega_2 = \sqrt{k/m}$, the eigen vector is given by

$$a_2 = \begin{pmatrix} \beta \\ 0 \\ -\beta \end{pmatrix} \quad \text{--- (7)}$$



(symmetric mode)

(iii) For $\omega_3 = \sqrt{\frac{k}{m} \left(1 + \frac{2m}{M}\right)}$

$$\begin{pmatrix} -\frac{2mK}{M} & -K & 0 \\ -K & -\frac{KM}{m} & -K \\ 0 & -K & -\frac{2mK}{M} \end{pmatrix} \begin{pmatrix} a_{13} \\ a_{23} \\ a_{33} \end{pmatrix} = 0$$

On solving we get—

$$\frac{2m}{M} a_{13} + a_{23} = 0$$

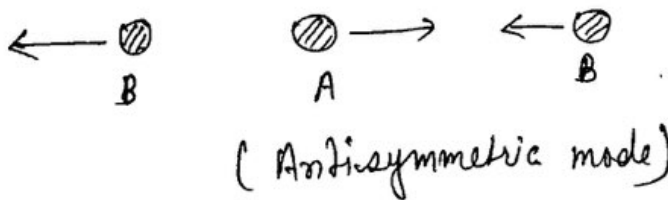
$$a_{13} + \frac{m}{M} a_{33} = 0$$

$$a_{23} + \frac{2m}{M} a_{33} = 0$$

Therefore, $a_{13} = a_{33} = \alpha$ (say) and $a_{23} = \left(-\frac{2m}{M}\right) \alpha$

Thus, the eigen vector is given by

$$a_3 = \begin{pmatrix} \alpha \\ -\frac{2m}{M} \alpha \\ \alpha \end{pmatrix} \quad \text{--- (8)}$$



Evaluation of α , β & γ

$$\text{Matrix } A = \begin{pmatrix} a_{11} & a_{12} & a_{13} \\ a_{21} & a_{22} & a_{23} \\ a_{31} & a_{32} & a_{33} \end{pmatrix} = \begin{pmatrix} \alpha & \beta & \gamma \\ \alpha & 0 & -\frac{2m}{M} \gamma \\ \alpha & -\beta & \gamma \end{pmatrix}$$

$$A^T = \begin{pmatrix} \alpha & \alpha & \alpha \\ \beta & 0 & -\beta \\ \gamma & -\frac{2m}{M} \gamma & \gamma \end{pmatrix}$$

We impose the matrix condition $A^T A = I$

$$\text{or } \begin{pmatrix} \alpha & \alpha & \alpha \\ \beta & 0 & -\beta \\ \gamma & \frac{-2m}{m}\gamma & \gamma \end{pmatrix} \begin{pmatrix} m & 0 & 0 \\ 0 & m & 0 \\ 0 & 0 & -m \end{pmatrix} \begin{pmatrix} \alpha & \beta & \gamma \\ \alpha & 0 & \frac{-2m}{m}\gamma \\ \alpha & -\beta & \gamma \end{pmatrix} = \begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix}$$

$$\text{or } \begin{pmatrix} (2m+m)\alpha^2 & 0 & 0 \\ 0 & 2m\beta^2 & 0 \\ 0 & 0 & 2m\gamma^2(1+\frac{2m}{m}) \end{pmatrix} = \begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix}$$

$$\text{or } (2m+m)\alpha^2 = 1$$

$$\therefore \alpha = \sqrt{1/(2m+m)}$$

$$\text{Similarly, } \beta = \sqrt{1/2m}$$

$$\gamma = \sqrt{\frac{1}{2m(1+\frac{2m}{m})}}$$

Eigenvectors : —

From equations (6), (7) & (8), we get—

$$a_1 = \frac{1}{\sqrt{2m+m}} \begin{pmatrix} 1 \\ 1 \\ 1 \end{pmatrix}$$

$$a_2 = \frac{1}{\sqrt{2m}} \begin{pmatrix} 1 \\ 0 \\ -1 \end{pmatrix}$$

$$a_3 = \frac{1}{\sqrt{2m(1+\frac{2m}{m})}} \begin{pmatrix} 1 \\ -2m/m \\ 1 \end{pmatrix}$$

These are called normal co-ordinates.

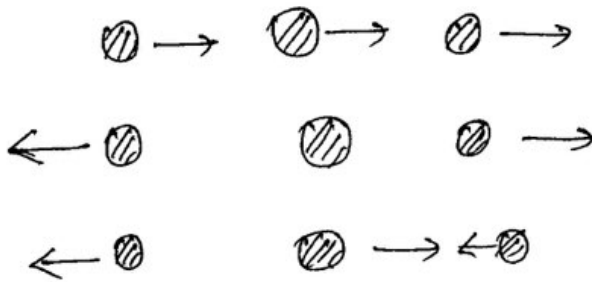
In general,

$$\begin{pmatrix} x_1 \\ x_2 \\ x_3 \end{pmatrix} = \begin{pmatrix} a_{11} & a_{12} & a_{13} \\ a_{21} & a_{22} & a_{23} \\ a_{31} & a_{32} & a_{33} \end{pmatrix} \begin{pmatrix} Q_1 \\ Q_2 \\ Q_3 \end{pmatrix}$$

$$a_1 \begin{pmatrix} x_1 \\ x_2 \\ x_3 \end{pmatrix} = \begin{pmatrix} \frac{1}{\sqrt{2m+M}} & \frac{1}{\sqrt{2m}} & \frac{1}{\sqrt{2m(1+\frac{2m}{M})}} \\ \frac{1}{\sqrt{2m+M}} & 0 & \frac{-2m}{M} \frac{1}{\sqrt{2m(1+\frac{2m}{M})}} \\ \frac{1}{\sqrt{2m+M}} & \frac{-1}{\sqrt{2m}} & \frac{1}{\sqrt{2m(1+\frac{2m}{M})}} \end{pmatrix} \begin{pmatrix} Q_1 \\ Q_2 \\ Q_3 \end{pmatrix}$$

$\therefore x_i = \sum_{k=1}^3 a_{ik} Q_k$
 where, $i=1,2,3$.

Conclusion



longitudinal normal modes of
triatomic molecules.