

# *Group Theory in Quantum Mechanics*

## *Lecture 17* (4.2.13)

### *Vibrational modes and symmetry reciprocity: Induced reps*

(Int.J.Mol.Sci, 14, 714(2013) p.755-774 , QTCA Unit 5 Ch. 15 )  
(PSDS - Ch. 4)

*Review: Hamiltonian local-symmetry eigensolution in global and local  $|\mathbf{P}^{(\mu)}\rangle$ -basis*

*Molecular vibrational modes vs. Hamiltonian eigenmodes*

*Molecular K-matrix construction*

*$D_3 \supset C_2(i_3)$  local-symmetry K-matrix eigensolutions*

*$D_3$ -direct-connection K-matrix eigensolutions*

*$D_3 \supset C_3(\mathbf{r}^{\pm l})$  local symmetry K-matrix eigensolutions*

*Applied symmetry reduction and splitting*

*Subduced irep  $D^\alpha(D^3) \downarrow C_2 = d^{0_2} \oplus d^{l_2} \oplus \dots$  correlation*

*Subduced irep  $D^\alpha(D^3) \downarrow C_3 = d^{0_3} \oplus d^{l_3} \oplus \dots$  correlation*

*Spontaneous symmetry breaking and clustering: Frobenius Reciprocity , band structure*

*Induced rep  $d^a(C_2) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus \dots$  correlation*

*Induced rep  $d^a(C_3) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus \dots$  correlation*

*$D_6$  symmetry and Hexagonal Bands*

*Cross product of the  $C_2$  and  $D_3$  characters gives all  $D_6 = D_3 \times C_2$  characters and irreps*

→ Review: *Hamiltonian local-symmetry eigensolution in global and local  $|P^{(\mu)}\rangle$ -basis* ←

Molecular vibrational modes vs. Hamiltonian eigenmodes

Molecular K-matrix construction

$D_3 \supset C_2(i_3)$  local-symmetry K-matrix eigensolutions

$D_3$ -direct-connection K-matrix eigensolutions

$D_3 \supset C_3(\mathbf{r}^{\pm l})$  local symmetry K-matrix eigensolutions

Applied symmetry reduction and splitting

Subduced irep  $D^\alpha(D^3) \downarrow C_2 = d^{0_2} \oplus d^{l_2} \oplus \dots$  correlation

Subduced irep  $D^\alpha(D^3) \downarrow C_3 = d^{0_3} \oplus d^{l_3} \oplus \dots$  correlation

Spontaneous symmetry breaking and clustering: Frobenius Reciprocity , band structure

Induced rep  $d^a(C_2) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus \dots$  correlation

Induced rep  $d^a(C_3) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus \dots$  correlation

$D_6$  symmetry and Hexagonal Bands

Cross product of the  $C_2$  and  $D_3$  characters gives all  $D_6 = D_3 \times C_2$  characters and irreps

# Compare Global vs Local $|g\rangle$ -basis vs. Global vs Local $|P^{(\mu)}\rangle$ -basis

$D_3$  global group product table

1	$r^2$	$r$	$i_1$	$i_2$	$(i_3)$
$r$	1	$r^2$	$(i_3)$	$i_1$	$i_2$
$r^2$	$r$	1	$i_2$	$(i_3)$	$i_1$
$i_1$	$(i_3)$	$i_2$	1	$r$	$r^2$
$i_2$	$i_1$	$(i_3)$	$r^2$	1	$r$
$(i_3)$	$i_2$	$i_1$	$r$	$r^2$	1

$D_3$  global projector product table

$D_3$	$P_{xx}^{A_1}$	$P_{yy}^{A_2}$	$P_{xx}^E$	$P_{xy}^E$	$P_{yx}^E$	$P_{yy}^E$
$P_{xx}^{A_1}$	$P_{xx}^{A_1}$	.	.	.	.	.
$P_{yy}^{A_2}$	.	$P_{yy}^{A_2}$	.	.	.	.
$P_{xx}^E$	.	.	$P_{xx}^E$	$P_{xy}^E$	.	.
$P_{yx}^E$	.	.	$P_{yx}^E$	$P_{yy}^E$	.	.
$P_{xy}^E$	.	.	.	.	$P_{xx}^E$	$P_{xy}^E$
$P_y^E$	.	.	.	.	$P_y^E$	$P_y^E$

$$P_{ab}^{(m)} P_{cd}^{(n)} = \delta^{mn} \delta_{bc} P_{ad}^{(m)}$$

Change Global to Local by switching

...column-P with column- $P^\dagger$

....and row-P with row- $P^\dagger$

Just switch  $r$  with  $r^\dagger = r^2$ . (all others are self-conjugate)

$D_3$  local group table

1	$r$	$r^2$	$i_1$	$i_2$	$(i_3)$
$r^2$	1	$r$	$i_2$	$(i_3)$	$i_1$
$r$	$r^2$	1	$(i_3)$	$i_1$	$i_2$
$i_1$	$i_2$	$(i_3)$	1	$r$	$r^2$
$i_2$	$(i_3)$	$i_2$	$r^2$	1	$r$
$(i_3)$	$i_1$	$i_2$	$r$	$r^2$	1

$D_3$  local projector product table

$D_3$	$P_{xx}^{A_1}$	$P_{yy}^{A_2}$	$P_{xx}^E$	$P_{yx}^E$	$P_{xy}^E$	$P_{yy}^E$
$P_{xx}^{A_1}$	$P_{xx}^{A_1}$	.	.	.	.	.
$P_{yy}^{A_2}$	.	$P_{yy}^{A_2}$	.	.	.	.
$P_{xx}^E$	.	.	$P_{xx}^E$	0	$P_{xy}^E$	0
$P_{xy}^E$	.	.	0	$P_{xx}^E$	0	$P_{xy}^E$
$P_{yx}^E$	.	.	$P_{yx}^E$	0	$P_{yy}^E$	0
$P_{yy}^E$	.	.	0	$P_{yx}^E$	0	$P_{yy}^E$

$$\bar{P}_{ab}^{(m)} \bar{P}_{cd}^{(n)} = \delta^{mn} \delta_{bc} \bar{P}_{ad}^{(m)}$$

(Just switch  $P_{yx}^E$  with  $P_{yx}^E = P_{xy}^E$ .)

## *D*<sub>3</sub> global-g group matrices in |P<sup>(μ)</sup>⟩-basis

$$R^P(\mathbf{g}) = TR^G(\mathbf{g})T^\dagger =$$

$$\left| \mathbf{P}_{xx}^{A_1} \right\rangle \quad \left| \mathbf{P}_{yy}^{A_2} \right\rangle \quad \left| \mathbf{P}_{\textcolor{blue}{xx}}^{E_1} \right\rangle \quad \left| \mathbf{P}_{\textcolor{blue}{yx}}^{E_1} \right\rangle \quad \left| \mathbf{P}_{\textcolor{blue}{xy}}^{E_1} \right\rangle \quad \left| \mathbf{P}_{yy}^{E_1} \right\rangle$$

$D^{A_1}(\mathbf{g})$	.	.	.	.	.
.	$D^{A_2}(\mathbf{g})$	.	.	.	.
.	.	$D_{xx}^{E_1}(\mathbf{g})$	$D_{xy}^{E_1}$	.	.
.	.	$D_{yx}^{E_1}(\mathbf{g})$	$D_{yy}^{E_1}$	.	.
.	.	.	.	$D_{xx}^{E_1}(\mathbf{g})$	$D_{xy}^{E_1}$
.	.	.	.	$D_{yx}^{E_1}(\mathbf{g})$	$D_{yy}^{E_1}$

|P<sup>(μ)</sup>⟩-base  
ordering to  
concentrate  
global-g  
D-matrices

$$\bar{R}^P(\mathbf{g}) = \bar{T}R^G(\mathbf{g})\bar{T}^\dagger =$$

$$\left| \mathbf{P}_{xx}^{A_1} \right\rangle \quad \left| \mathbf{P}_{yy}^{A_2} \right\rangle \quad \left| \mathbf{P}_{\textcolor{blue}{xx}}^{E_1} \right\rangle \quad \left| \mathbf{P}_{\textcolor{blue}{xy}}^{E_1} \right\rangle \quad \left| \mathbf{P}_{\textcolor{blue}{yx}}^{E_1} \right\rangle \quad \left| \mathbf{P}_{yy}^{E_1} \right\rangle$$

$D^{A_1}(\mathbf{g})$	.	.	.	.	.
.	$D^{A_2}(\mathbf{g})$	.	.	.	.
.	.	$D_{xx}^{E_1}(\mathbf{g})$	.	$D_{xy}^{E_1}(\mathbf{g})$	.
.	.	.	$D_{xx}^{E_1}$	.	$D_{xy}^{E_1}$
.	.	$D_{yx}^{E_1}(\mathbf{g})$	.	$D_{yy}^{E_1}(\mathbf{g})$	.
.	.	.	$D_{yx}^{E_1}$	.	$D_{yy}^{E_1}$

Global g-matrix component

$$\left\langle \begin{array}{c} \mu \\ m'n \end{array} \middle| \mathbf{g} \middle| \begin{array}{c} \mu \\ mn \end{array} \right\rangle = D_{m'm}^\mu(\mathbf{g})$$

## *D*<sub>3</sub> local- $\bar{\mathbf{g}}$ group matrices in |P<sup>(μ)</sup>⟩-basis

$$R^P(\bar{\mathbf{g}}) = TR^G(\bar{\mathbf{g}})T^\dagger =$$

$$\left| \mathbf{P}_{xx}^{A_1} \right\rangle \quad \left| \mathbf{P}_{yy}^{A_2} \right\rangle \quad \left| \mathbf{P}_{\textcolor{red}{xx}}^{E_1} \right\rangle \quad \left| \mathbf{P}_{\textcolor{red}{yx}}^{E_1} \right\rangle \quad \left| \mathbf{P}_{\textcolor{red}{xy}}^{E_1} \right\rangle \quad \left| \mathbf{P}_{yy}^{E_1} \right\rangle$$

$D^{A_1}(\mathbf{g})$	.	.	.	.	.
.	$D^{A_2}(\mathbf{g})$	.	.	.	.
.	.	$D_{xx}^{E_1}(\mathbf{g})$	.	$D_{xy}^{E_1}(\mathbf{g})$	.
.	.	.	$D_{xx}^{E_1}$	.	$D_{xy}^{E_1}$
.	.	$D_{yx}^{E_1}(\mathbf{g})$	.	$D_{yy}^{E_1}(\mathbf{g})$	.
.	.	.	$D_{yx}^{E_1}$	.	$D_{yy}^{E_1}$

|P<sup>(μ)</sup>⟩-base  
ordering to  
concentrate  
local- $\bar{\mathbf{g}}$   
D-matrices  
and  
H-matrices

$$\bar{R}^P(\bar{\mathbf{g}}) = \bar{T}R^G(\bar{\mathbf{g}})\bar{T}^\dagger =$$

$$\left| \mathbf{P}_{xx}^{A_1} \right\rangle \quad \left| \mathbf{P}_{yy}^{A_2} \right\rangle \quad \left| \mathbf{P}_{\textcolor{red}{xx}}^{E_1} \right\rangle \quad \left| \mathbf{P}_{\textcolor{red}{xy}}^{E_1} \right\rangle \quad \left| \mathbf{P}_{\textcolor{red}{yx}}^{E_1} \right\rangle \quad \left| \mathbf{P}_{yy}^{E_1} \right\rangle$$

$D^{A_1}(\mathbf{g})$	.	.	.	.	.
.	$D^{A_2}(\mathbf{g})$	.	.	.	.
.	.	$D_{xx}^{E_1}(\mathbf{g})$	$D_{xy}^{E_1}(\mathbf{g})$	.	.
.	.	.	$D_{xx}^{E_1}$	$D_{xy}^{E_1}$	.
.	.	$D_{yx}^{E_1}(\mathbf{g})$	$D_{yy}^{E_1}(\mathbf{g})$	.	.
.	.	.	$D_{yx}^{E_1}$	$D_{yy}^{E_1}$	.

Local  $\bar{\mathbf{g}}$ -matrix component

$$\left\langle \begin{array}{c} \mu \\ mn' \end{array} \middle| \bar{\mathbf{g}} \middle| \begin{array}{c} \mu \\ mn \end{array} \right\rangle = D_{nn'}^\mu(\mathbf{g}^{-1}) = D_{n'n}^{\mu*}(\mathbf{g})$$

### D<sub>3</sub> Hamiltonian local- H matrices in |P<sup>(μ)</sup>⟩-basis

**H matrix in |g⟩-basis:**

$$(\mathbf{H})_G = \sum_{g=1}^{\circ G} r_g \bar{\mathbf{g}} = \begin{pmatrix} r_0 & r_2 & r_1 & i_1 & i_2 & i_3 \\ r_1 & r_0 & r_1 & i_3 & i_1 & i_2 \\ r_2 & r_1 & r_0 & i_2 & i_3 & i_1 \\ i_i & i_3 & i_2 & r_0 & r_1 & r_2 \\ i_2 & i_1 & i_3 & r_2 & r_0 & r_1 \\ i_3 & i_2 & i_1 & r_1 & r_2 & r_0 \end{pmatrix}$$

**H matrix in |P<sup>(μ)</sup>⟩-basis:**

$$(\mathbf{H})_P = \bar{T} (\mathbf{H})_G \bar{T}^\dagger =$$

$$\begin{array}{c|c|c|c|c} \left| \mathbf{P}_{xx}^{A_1} \right\rangle & \left| \mathbf{P}_{yy}^{A_2} \right\rangle & \left| \mathbf{P}_{xx}^{E_1} \right\rangle \left| \mathbf{P}_{xy}^{E_1} \right\rangle & \left| \mathbf{P}_{yx}^{E_1} \right\rangle \left| \mathbf{P}_{yy}^{E_1} \right\rangle \\ \hline H^{A_1} & \cdot & \cdot & \cdot & \cdot \\ \cdot & H^{A_2} & \cdot & \cdot & \cdot \\ \cdot & \cdot & H_{xx}^{E_1} & H_{xy}^{E_1} & \cdot \\ \cdot & \cdot & H_{yx}^{E_1} & H_{yy}^{E_1} & \cdot \\ \cdot & \cdot & \cdot & \cdot & H_{xx}^{E_1} & H_{xy}^{E_1} \\ \cdot & \cdot & \cdot & \cdot & H_{yx}^{E_1} & H_{yy}^{E_1} \end{array}$$

$$H_{ab}^\alpha = \langle \mathbf{P}_m^\mu | \mathbf{H} | \mathbf{P}_n^\mu \rangle = \frac{\langle \mathbf{1} | \mathbf{P}_{am}^\mu \mathbf{H} \mathbf{P}_{nb}^\mu | \mathbf{1} \rangle}{(norm)^2} = \langle \mathbf{1} | \mathbf{H} \mathbf{P}_{am}^\mu \mathbf{P}_{nb}^\mu | \mathbf{1} \rangle = \delta_{mn} \langle \mathbf{1} | \mathbf{H} \mathbf{P}_{ab}^\mu | \mathbf{1} \rangle = \sum_{g=1}^{\circ G} \langle \mathbf{1} | \mathbf{H} | \mathbf{g} \rangle D_{ab}^{\alpha*}(g) = \sum_{g=1}^{\circ G} r_g D_{ab}^{\alpha*}(g)$$

$$H^{A_1} = r_0 D^{A_1*}(1) + r_1 D^{A_1*}(r^1) + r_1^* D^{A_1*}(r^2) + i_1 D^{A_1*}(i_1) + i_2 D^{A_1*}(i_2) + i_3 D^{A_1*}(i_3) = r_0 + r_1 + r_1^* + i_1 + i_2 + i_3$$

$$H^{A_2} = r_0 D^{A_2*}(1) + r_1 D^{A_2*}(r^1) + r_1^* D^{A_2*}(r^2) + i_1 D^{A_2*}(i_1) + i_2 D^{A_2*}(i_2) + i_3 D^{A_2*}(i_3) = r_0 + r_1 + r_1^* - i_1 - i_2 - i_3$$

$$H_{xx}^{E_1} = r_0 D_{xx}^{E_1*}(1) + r_1 D_{xx}^{E_1*}(r^1) + r_1^* D_{xx}^{E_1*}(r^2) + i_1 D_{xx}^{E_1*}(i_1) + i_2 D_{xx}^{E_1*}(i_2) + i_3 D_{xx}^{E_1*}(i_3) = (2r_0 - r_1 - r_1^* - i_1 - i_2 + 2i_3)/2$$

$$H_{xy}^{E_1} = r_0 D_{xy}^{E_1*}(1) + r_1 D_{xy}^{E_1*}(r^1) + r_1^* D_{xy}^{E_1*}(r^2) + i_1 D_{xy}^{E_1*}(i_1) + i_2 D_{xy}^{E_1*}(i_2) + i_3 D_{xy}^{E_1*}(i_3) = \sqrt{3}(-r_1 + r_1^* - i_1 + i_2)/2 = H_{yx}^{E_1*}$$

$$H_{yy}^{E_1} = r_0 D_{yy}^{E_1*}(1) + r_1 D_{yy}^{E_1*}(r^1) + r_1^* D_{yy}^{E_1*}(r^2) + i_1 D_{yy}^{E_1*}(i_1) + i_2 D_{yy}^{E_1*}(i_2) + i_3 D_{yy}^{E_1*}(i_3) = (2r_0 - r_1 - r_1^* + i_1 + i_2 - 2i_3)/2$$

$$= r_0 + 2r_1 + 2i_{12} + i_3$$

$$= r_0 + 2r_1 - 2i_{12} - i_3$$

$$= r_0 - r_1 - i_{12} + i_3$$

$$= 0$$

$$= r_0 - r_1 + i_{12} - i_3$$

$$C_2 = \{\mathbf{1}, \mathbf{i}_3\}$$

*Local symmetry determines all levels and eigenvectors with just 4 real parameters*

$$\begin{pmatrix} H_{xx}^{E_1} & H_{xy}^{E_1} \\ H_{yx}^{E_1} & H_{yy}^{E_1} \end{pmatrix} = \frac{1}{2} \begin{pmatrix} 2r_0 - r_1 - r_1^* - i_1 - i_2 + 2i_3 & \sqrt{3}(-r_1 + r_1^* - i_1 + i_2) \\ \sqrt{3}(-r_1^* + r_1 - i_1 + i_2) & 2r_0 - r_1 - r_1^* + i_1 + i_2 - 2i_3 \end{pmatrix}$$

$$= \begin{pmatrix} r_0 - r_1 - i_{12} + i_3 & 0 \\ 0 & r_0 - r_1 - i_{12} - i_3 \end{pmatrix}$$

*Choosing local C<sub>2</sub> = {1, i<sub>3</sub>} symmetry with local constraints r<sub>1</sub> = r<sub>1</sub><sup>\*</sup> = r<sub>2</sub> and i<sub>1</sub> = i<sub>2</sub> for: r<sub>1</sub> = r<sub>1</sub><sup>\*</sup> and i<sub>1</sub> = i<sub>2</sub>*

$$\mathbf{P}_{mn}^{(\mu)} = \sum_{\mathbf{g}}^{\ell(\mu)} \mathbf{D}_{mn}^{(\mu)} \mathbf{g}^* \mathbf{g}$$

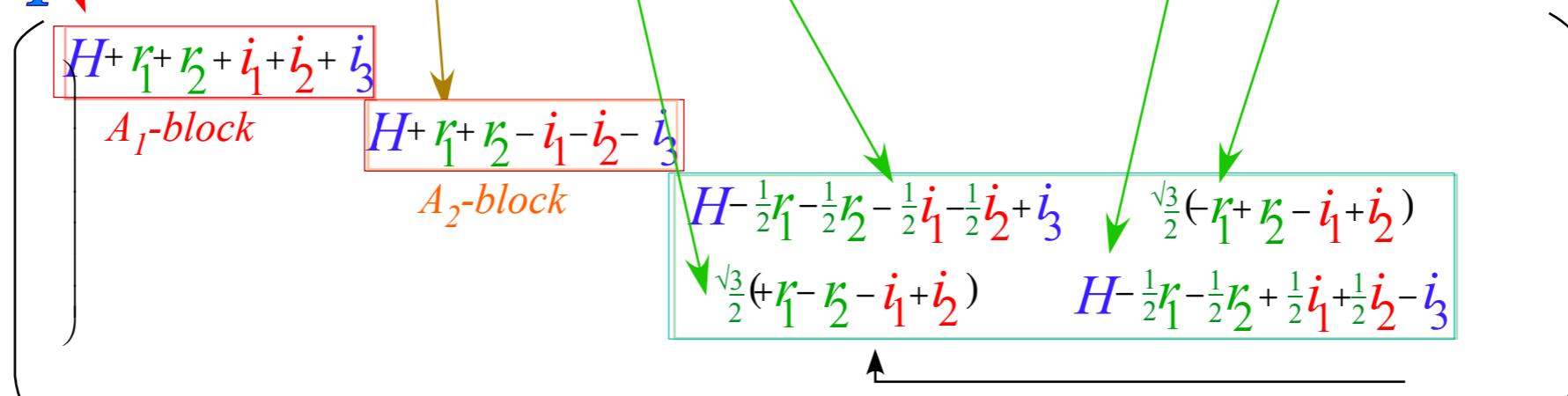
*Spectral Efficiency: Same  $D(a)_{mn}$  projectors give a lot!*

$$\begin{array}{ccccccc} 1 & \mathbf{r}^1 & \mathbf{r}^2 & \mathbf{i}_1 & \mathbf{i}_2 & \mathbf{i}_3 \\ \mathbf{P}_{x,x}^{A_1} = & \frac{(1 \ 1 \ 1 \ 1 \ 1 \ 1)/6}{(1 \ 1 \ 1 \ -1 \ -1 \ -1)/6} & & & & & \\ \mathbf{P}_{y,y}^{A_2} = & & & & & & \end{array}$$

$$\begin{array}{ccccccc} 1 & \mathbf{r}^1 & \mathbf{r}^2 & \mathbf{i}_1 & \mathbf{i}_2 & \mathbf{i}_3 \\ \mathbf{P}_{x,x}^E = & \frac{(2 \ -1 \ -1 \ -1 \ -1 \ +2)/6}{(0 \ 1 \ -1 \ -1 \ +1 \ 0)/\sqrt{3}/2} & & & & & \\ \mathbf{P}_{y,x}^E = & & & & & & \end{array}$$

$$\begin{array}{ccccccc} 1 & \mathbf{r}^1 & \mathbf{r}^2 & \mathbf{i}_1 & \mathbf{i}_2 & \mathbf{i}_3 \\ \mathbf{P}_{x,y}^E = & \frac{(0 \ -1 \ 1 \ -1 \ +1 \ 0)/\sqrt{3}/2}{(2 \ -1 \ -1 \ +1 \ +1 \ -2)/6} & & & & & \\ \mathbf{P}_{y,y}^E = & & & & & & \end{array}$$

- Eigenstates (shown before)
- Complete Hamiltonian



- Local symmetry eigenvalue formulae (L.S. => off-diagonal zero.)

$$C_2 = \{1, \mathbf{i}_3\}$$

Local symmetry determines all levels and eigenvectors with just 4 real parameters

$$r_1 = r_2 = r_1^* = r, \quad i_1 = i_2 = i_1^* = i$$

$$\begin{aligned} & A_1\text{-level: } H + 2r + 2i + i_3 \\ & \text{gives: } A_1\text{-level: } H + 2r - 2i - i_3 \\ & E_x\text{-level: } H - r - i + i_3 \\ & E_y\text{-level: } H - r + i - i_3 \end{aligned}$$

*Global (LAB) symmetry*

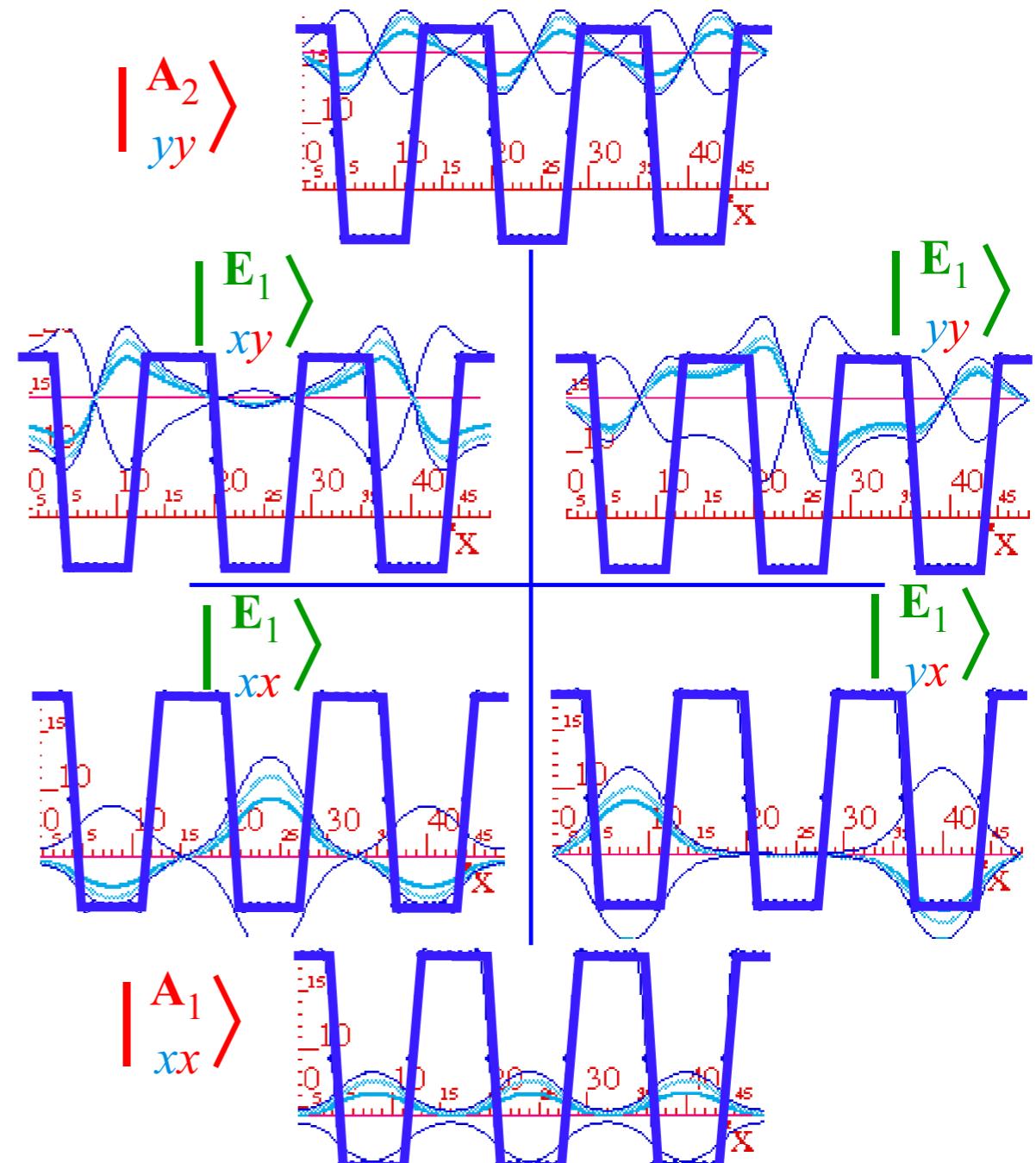
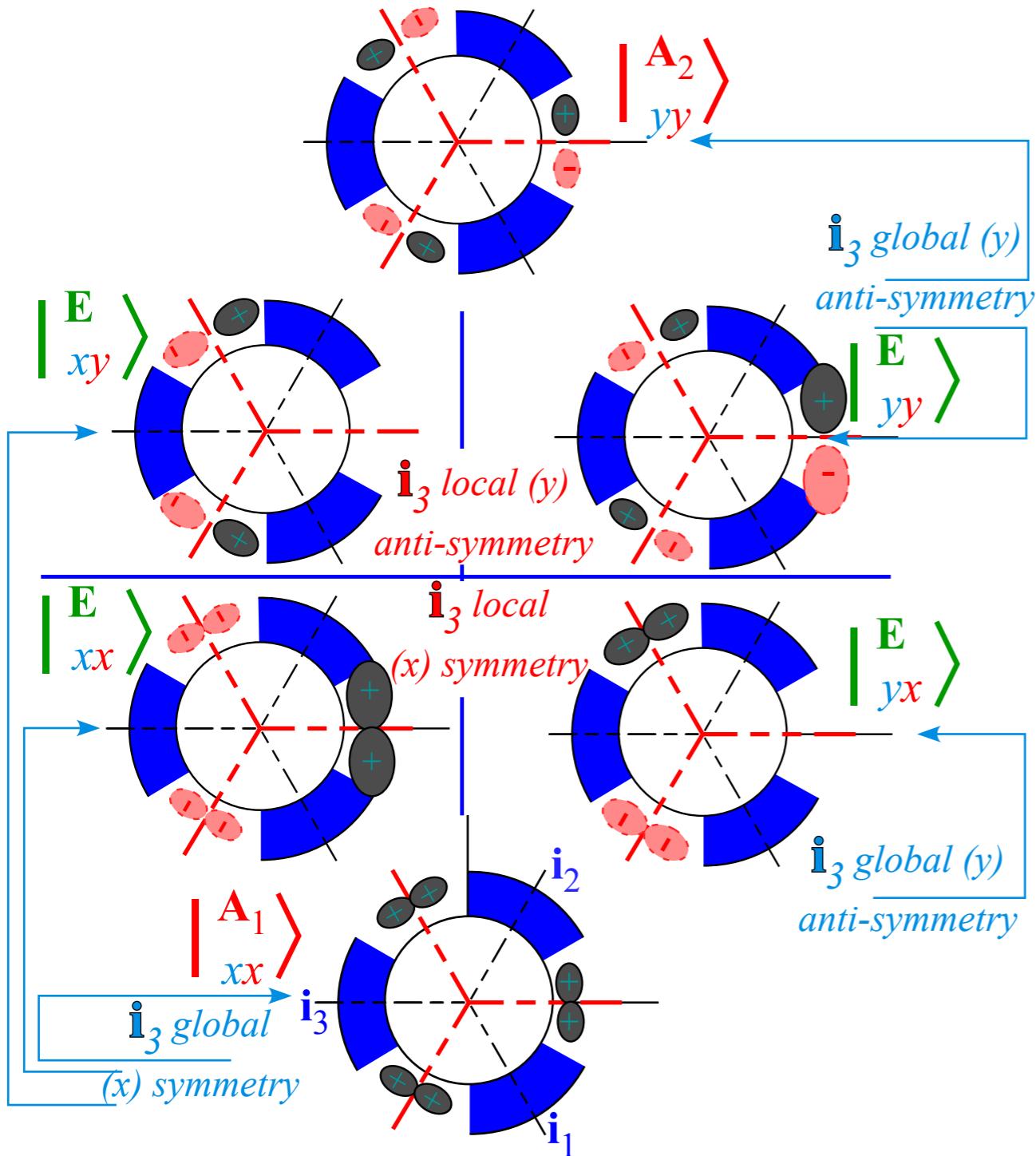
$$\mathbf{i}_3|_{eb}^{(m)}\rangle = \mathbf{i}_3 \mathbf{P}_{eb}^{(m)} |1\rangle \\ = (-1)^e |^{(m)}\rangle$$

$D_3 > C_2$   $\mathbf{i}_3$  projector states

$$|_{eb}^{(m)}\rangle = \mathbf{P}_{eb}^{(m)} |1\rangle$$

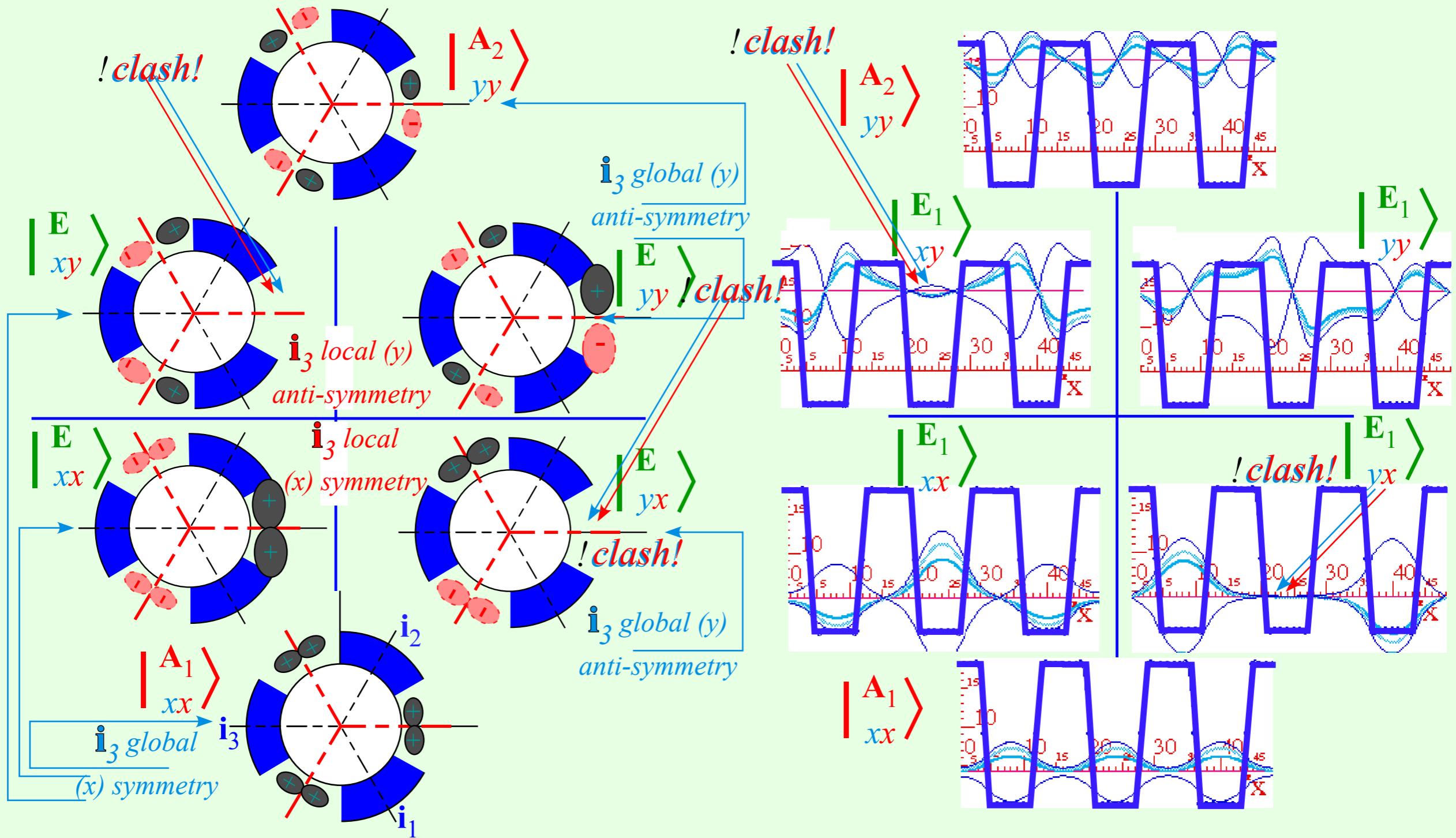
*Local (BOD) symmetry*

$$\bar{\mathbf{i}}_3|_{eb}^{(m)}\rangle = \bar{\mathbf{i}}_3 \mathbf{P}_{eb}^{(m)} |1\rangle = \mathbf{P}_{eb}^{(m)} \bar{\mathbf{i}}_3 |1\rangle \\ = \mathbf{P}_{eb}^{(m)} \mathbf{i}_3^\dagger |1\rangle = (-1)^b |^{(m)}\rangle$$



# *When there is no there, there...*

Nobody Home  
where **LOCAL**  
and **GLOBAL**



Review: Hamiltonian local-symmetry eigensolution in global and local  $|\mathbf{P}^{(\mu)}\rangle$ -basis

→ Molecular vibrational modes vs. Hamiltonian eigenmodes ←

Molecular K-matrix construction

$D_3 \supset C_2(i_3)$  local-symmetry K-matrix eigensolutions

$D_3$ -direct-connection K-matrix eigensolutions

$D_3 \supset C_3(\mathbf{r}^{\pm l})$  local symmetry K-matrix eigensolutions

Applied symmetry reduction and splitting

Subduced irep  $D^\alpha(D^3) \downarrow C_2 = d^{0_2} \oplus d^{l_2} \oplus \dots$  correlation

Subduced irep  $D^\alpha(D^3) \downarrow C_3 = d^{0_3} \oplus d^{l_3} \oplus \dots$  correlation

Spontaneous symmetry breaking and clustering: Frobenius Reciprocity , band structure

Induced rep  $d^a(C_2) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus \dots$  correlation

Induced rep  $d^a(C_3) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus \dots$  correlation

$D_6$  symmetry and Hexagonal Bands

Cross product of the  $C_2$  and  $D_3$  characters gives all  $D_6 = D_3 \times C_2$  characters and ireps

## *Molecular vibrational modes vs. Hamiltonian eigenmodes*

Classical equations of coupled harmonic motion are Newtonian  $\mathbf{F}=\mathbf{M}\cdot\mathbf{a}$  relations of  $n$ -dimensional force vector  $\mathbf{F}$ , acceleration vector  $\mathbf{a}$ , and mass operator  $\mathbf{M}=M\cdot\mathbf{1}$  for  $D_3$ -symmetry. Force  $\mathbf{F}$  is a (-)derivative of potential  $V(x)$  that becomes a  $\mathbf{F}=-\mathbf{K}\cdot\mathbf{x}$  matrix expression.

$$-M\partial_t^2 x^a = \frac{\partial V}{\partial x^a} = \sum_b K_{ab} x^b$$

## *Molecular vibrational modes vs. Hamiltonian eigenmodes*

Classical equations of coupled harmonic motion are Newtonian  $\mathbf{F}=\mathbf{M}\cdot\mathbf{a}$  relations of  $n$ -dimensional force vector  $\mathbf{F}$ , acceleration vector  $\mathbf{a}$ , and mass operator  $\mathbf{M}=M\cdot\mathbf{1}$  for  $D_3$ -symmetry. Force  $\mathbf{F}$  is a (-)derivative of potential  $V(x)$  that becomes a  $\mathbf{F}=-\mathbf{K}\cdot\mathbf{x}$  matrix expression.

$$-M\partial_t^2 x^a = \frac{\partial V}{\partial x^a} = \sum_b K_{ab} x^b$$

## Molecular vibrational modes vs. Hamiltonian eigenmodes

Classical equations of coupled harmonic motion are Newtonian  $\mathbf{F}=\mathbf{M}\cdot\mathbf{a}$  relations of  $n$ -dimensional force vector  $\mathbf{F}$ , acceleration vector  $\mathbf{a}$ , and mass operator  $\mathbf{M}=M\cdot\mathbf{1}$  for  $D_3$ -symmetry. Force  $\mathbf{F}$  is a (-)derivative of potential  $V(x)$  that becomes a  $\mathbf{F}=-\mathbf{K}\cdot\mathbf{x}$  matrix expression.

$$-M\partial_t^2 x^a = \frac{\partial V}{\partial x^a} = \sum_b K_{ab} x^b$$

Compare classical equation to Schrodinger's equation for quantum motion.

$$i\hbar\partial_t \psi^a = \sum_b H_{ab} \psi^b$$

## Molecular vibrational modes vs. Hamiltonian eigenmodes

Classical equations of coupled harmonic motion are Newtonian  $\mathbf{F}=\mathbf{M}\cdot\mathbf{a}$  relations of  $n$ -dimensional force vector  $\mathbf{F}$ , acceleration vector  $\mathbf{a}$ , and mass operator  $\mathbf{M}=M\cdot\mathbf{1}$  for  $D_3$ -symmetry. Force  $\mathbf{F}$  is a (-)derivative of potential  $V(x)$  that becomes a  $\mathbf{F}=-\mathbf{K}\cdot\mathbf{x}$  matrix expression.

$$-M\partial_t^2 x^a = \frac{\partial V}{\partial x^a} = \sum_b K_{ab} x^b$$

Compare classical equation to Schrodinger's equation for quantum motion.

$$i\hbar\partial_t\psi^a = \sum_b H_{ab}\psi^b$$

Squared time generator ( $i\hbar\partial_t=\mathbf{H})^2$  has classical form with  $K=H^2$  and  $M=\hbar^2$ .

$$-\hbar^2\partial_t^2\psi^a = \sum_b K_{ab}\psi^b \text{ where: } K = H^2$$

## Molecular vibrational modes vs. Hamiltonian eigenmodes

Classical equations of coupled harmonic motion are Newtonian  $\mathbf{F}=\mathbf{M}\cdot\mathbf{a}$  relations of  $n$ -dimensional force vector  $\mathbf{F}$ , acceleration vector  $\mathbf{a}$ , and mass operator  $\mathbf{M}=M\cdot\mathbf{1}$  for  $D_3$ -symmetry. Force  $\mathbf{F}$  is a (-)derivative of potential  $V(x)$  that becomes a  $\mathbf{F}=-\mathbf{K}\cdot\mathbf{x}$  matrix expression.

$$-M\partial_t^2 x^a = \frac{\partial V}{\partial x^a} = \sum_b K_{ab} x^b$$

Compare classical equation to Schrodinger's equation for quantum motion.

$$i\hbar\partial_t\psi^a = \sum_b H_{ab}\psi^b$$

Squared time generator ( $i\hbar\partial_t=\mathbf{H})^2$  has classical form with  $K=H^2$  and  $M=\hbar^2$ .

$$-\hbar^2\partial_t^2\psi^a = \sum_b K_{ab}\psi^b \text{ where: } K = H^2$$

$(\mathbf{H}/\hbar)$ -eigenvalues are quantum angular frequencies  $\epsilon_m/\hbar=\omega_m$ . (Like Planck axiom:  $\epsilon=\hbar\omega$ .)

## Molecular vibrational modes vs. Hamiltonian eigenmodes

Classical equations of coupled harmonic motion are Newtonian  $\mathbf{F}=\mathbf{M}\cdot\mathbf{a}$  relations of  $n$ -dimensional force vector  $\mathbf{F}$ , acceleration vector  $\mathbf{a}$ , and mass operator  $\mathbf{M}=M\cdot\mathbf{1}$  for  $D_3$ -symmetry. Force  $\mathbf{F}$  is a (-)derivative of potential  $V(x)$  that becomes a  $\mathbf{F}=-\mathbf{K}\cdot\mathbf{x}$  matrix expression.

$$-M\partial_t^2 x^a = \frac{\partial V}{\partial x^a} = \sum_b K_{ab} x^b$$

Compare classical equation to Schrodinger's equation for quantum motion.

$$i\hbar\partial_t\psi^a = \sum_b H_{ab}\psi^b$$

Squared time generator ( $i\hbar\partial_t=\mathbf{H})^2$  has classical form with  $K=H^2$  and  $M=\hbar^2$ .

$$-\hbar^2\partial_t^2\psi^a = \sum_b K_{ab}\psi^b \text{ where: } K = H^2$$

$(\mathbf{H}/\hbar)$ -eigenvalues are quantum angular frequencies  $\epsilon_m/\hbar=\omega_m$ . (Like Planck axiom:  $\epsilon=\hbar\omega$ .)

$(\mathbf{K}/M)$ -eigenvalues are classical *squared* frequencies  $k_m/M=\omega_m^2$ . (Like Hooke's law:  $k/M=\omega^2$ .)

## Molecular vibrational modes vs. Hamiltonian eigenmodes

Classical equations of coupled harmonic motion are Newtonian  $\mathbf{F}=\mathbf{M}\cdot\mathbf{a}$  relations of  $n$ -dimensional force vector  $\mathbf{F}$ , acceleration vector  $\mathbf{a}$ , and mass operator  $\mathbf{M}=M\cdot\mathbf{1}$  for  $D_3$ -symmetry. Force  $\mathbf{F}$  is a (-)derivative of potential  $V(x)$  that becomes a  $\mathbf{F}=-\mathbf{K}\cdot\mathbf{x}$  matrix expression.

$$-M\partial_t^2 x^a = \frac{\partial V}{\partial x^a} = \sum_b K_{ab} x^b$$

Compare classical equation to Schrodinger's equation for quantum motion.

$$i\hbar\partial_t\psi^a = \sum_b H_{ab}\psi^b$$

Squared time generator ( $i\hbar\partial_t=\mathbf{H})^2$  has classical form with  $K=H^2$  and  $M=\hbar^2$ .

$$-\hbar^2\partial_t^2\psi^a = \sum_b K_{ab}\psi^b \text{ where: } K = H^2$$

$(\mathbf{H}/\hbar)$ -eigenvalues are quantum angular frequencies  $\epsilon_m/\hbar=\omega_m$ . (Like Planck axiom:  $\epsilon=\hbar\omega$ .)

$(\mathbf{K}/M)$ -eigenvalues are classical *squared* frequencies  $k_m/M=\omega_m^2$ . (Like Hooke's law:  $k/M=\omega^2$ .)

Apart from normalization, eigenvectors of  $\mathbf{H}$  and  $\mathbf{K}$  are the same.

## Molecular vibrational modes vs. Hamiltonian eigenmodes

Classical equations of coupled harmonic motion are Newtonian  $\mathbf{F}=\mathbf{M}\cdot\mathbf{a}$  relations of  $n$ -dimensional force vector  $\mathbf{F}$ , acceleration vector  $\mathbf{a}$ , and mass operator  $\mathbf{M}=M\cdot\mathbf{1}$  for  $D_3$ -symmetry. Force  $\mathbf{F}$  is a (-)derivative of potential  $V(x)$  that becomes a  $\mathbf{F}=-\mathbf{K}\cdot\mathbf{x}$  matrix expression.

$$-M\partial_t^2 x^a = \frac{\partial V}{\partial x^a} = \sum_b K_{ab} x^b$$

Compare classical equation to Schrodinger's equation for quantum motion.

$$i\hbar\partial_t\psi^a = \sum_b H_{ab}\psi^b$$

Squared time generator ( $i\hbar\partial_t=\mathbf{H})^2$  has classical form with  $K=H^2$  and  $M=\hbar^2$ .

$$-\hbar^2\partial_t^2\psi^a = \sum_b K_{ab}\psi^b \text{ where: } K = H^2$$

$(\mathbf{H}/\hbar)$ -eigenvalues are quantum angular frequencies  $\epsilon_m/\hbar=\omega_m$ . (Like Planck axiom:  $\epsilon=\hbar\omega$ .)

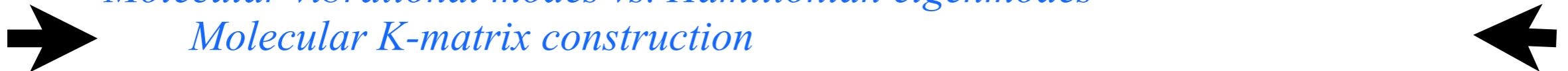
$(\mathbf{K}/M)$ -eigenvalues are classical *squared* frequencies  $k_m/M=\omega_m^2$ . (Like Hooke's law:  $k/M=\omega^2$ .)

Apart from normalization, eigenvectors of  $\mathbf{H}$  and  $\mathbf{K}$  are the same.

And, each eigenvalue set corresponds to its respective energy spectrum.

*Review: Hamiltonian local-symmetry eigensolution in global and local  $|\mathbf{P}^{(\mu)}\rangle$ -basis*

*Molecular vibrational modes vs. Hamiltonian eigenmodes*



*Molecular K-matrix construction*

*$D_3 \supset C_2(i_3)$  local-symmetry K-matrix eigensolutions*

*$D_3$ -direct-connection K-matrix eigensolutions*

*$D_3 \supset C_3(\mathbf{r}^{\pm l})$  local symmetry K-matrix eigensolutions*



*Applied symmetry reduction and splitting*

*Subduced irep  $D^\alpha(D^3) \downarrow C_2 = d^{0_2} \oplus d^{l_2} \oplus \dots$  correlation*

*Subduced irep  $D^\alpha(D^3) \downarrow C_3 = d^{0_3} \oplus d^{l_3} \oplus \dots$  correlation*

*Spontaneous symmetry breaking and clustering: Frobenius Reciprocity , band structure*

*Induced rep  $d^a(C_2) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus \dots$  correlation*

*Induced rep  $d^a(C_3) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus \dots$  correlation*

*$D_6$  symmetry and Hexagonal Bands*

*Cross product of the  $C_2$  and  $D_3$  characters gives all  $D_6 = D_3 \times C_2$  characters and ireps*

## Molecular K-matrix construction

Classical modes are eigenvectors of force-field matrix  $K$  or operator  $\mathbf{K}$ .

Harmonic potential  $V(\mathbf{x})$  is a quadratic  $K$ -form of coordinates  $x_a$  based on six  $D_3$ -labeled axes  $\hat{\mathbf{x}}^a$  or  $|a\rangle$ .

$$V(x) = \sum_{(k)} \frac{1}{2} \langle x | \mathbf{K} | x \rangle \quad \text{where: } |x\rangle = \sum_a x_a |a\rangle, \quad (a, b) = (1, r^1, r^2, i_1, i_2, i_3)$$

## Molecular K-matrix construction

Classical modes are eigenvectors of force-field matrix  $K$  or operator  $\mathbf{K}$ .

Harmonic potential  $V(\mathbf{x})$  is a quadratic  $K$ -form of coordinates  $x_a$  based on six  $D_3$ -labeled axes  $\hat{\mathbf{x}}^a$  or  $|a\rangle$ .

$$V(x) = \sum_{(k)} \frac{1}{2} \langle x | \mathbf{K} | x \rangle \quad \text{where: } |x\rangle = \sum_a x_a |a\rangle, \quad (a, b) = (1, r^1, r^2, i_1, i_2, i_3)$$

Each  $\mathbf{K}$  component  $K_{ab} = \langle a | \mathbf{K} | b \rangle$  is a sum over spring  $k$ -constants that connect axis- $\mathbf{x}^a$  to axis- $\mathbf{x}^b$  multiplied by factor  $(\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)(\hat{\mathbf{k}}_b \bullet \hat{\mathbf{x}}^b)$  for projecting spring- $k$ 's end vectors  $\hat{\mathbf{k}}_a$  and  $\hat{\mathbf{k}}_b$  onto  $\hat{\mathbf{x}}^a$  and  $\hat{\mathbf{x}}^b$  at respective connections.

## Molecular K-matrix construction

Classical modes are eigenvectors of force-field matrix  $K$  or operator  $\mathbf{K}$ .

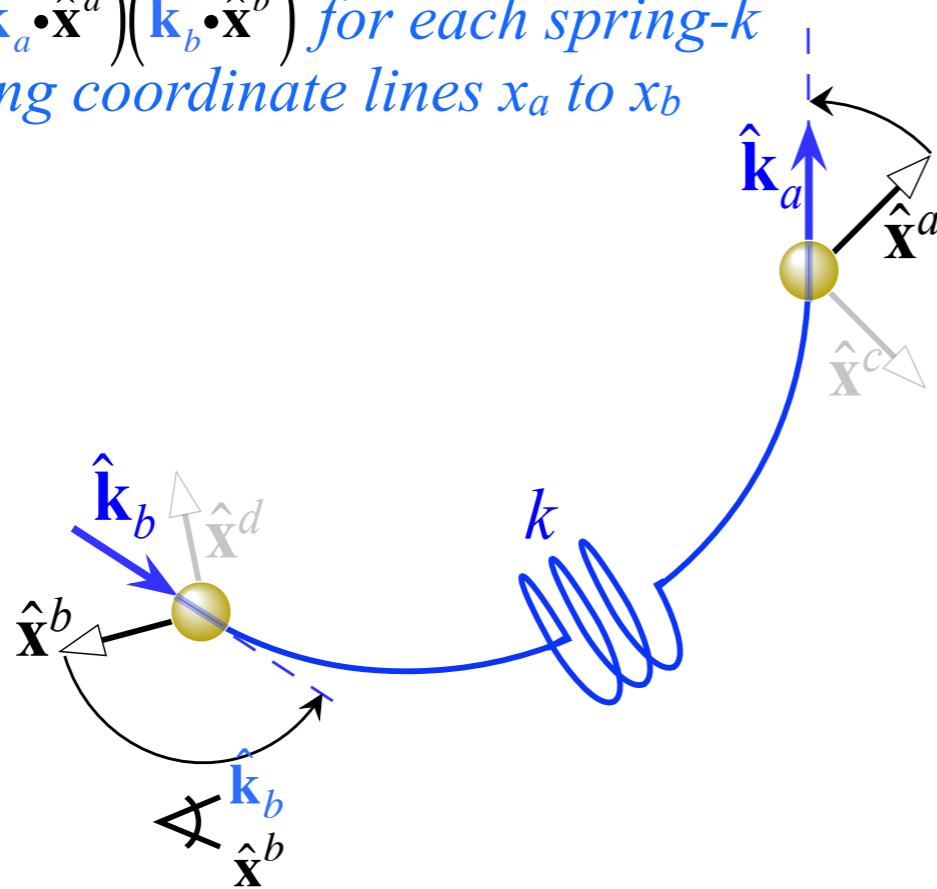
Harmonic potential  $V(\mathbf{x})$  is a quadratic  $K$ -form of coordinates  $x_a$  based on six  $D_3$ -labeled axes  $\hat{\mathbf{x}}^a$  or  $|a\rangle$ .

$$V(x) = \sum_{(k)} \frac{1}{2} \langle x | \mathbf{K} | x \rangle \quad \text{where: } |x\rangle = \sum_a x_a |a\rangle, \quad (a, b) = (1, r^1, r^2, i_1, i_2, i_3)$$

Each  $\mathbf{K}$  component  $K_{ab} = \langle a | \mathbf{K} | b \rangle$  is a sum over spring  $k$ -constants that connect axis- $\mathbf{x}^a$  to axis- $\mathbf{x}^b$  multiplied by factor  $(\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)(\hat{\mathbf{k}}_b \bullet \hat{\mathbf{x}}^b)$  for projecting spring- $k$ 's end vectors  $\hat{\mathbf{k}}_a$  and  $\hat{\mathbf{k}}_b$  onto  $\hat{\mathbf{x}}^a$  and  $\hat{\mathbf{x}}^b$  at respective connections.

*Coupling  $K_{ab} = \langle a | \mathbf{K} | b \rangle$*

*Sum  $k \cdot (\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)(\hat{\mathbf{k}}_b \bullet \hat{\mathbf{x}}^b)$  for each spring- $k$  connecting coordinate lines  $x_a$  to  $x_b$*



## Molecular K-matrix construction

Classical modes are eigenvectors of force-field matrix  $K$  or operator  $\mathbf{K}$ .

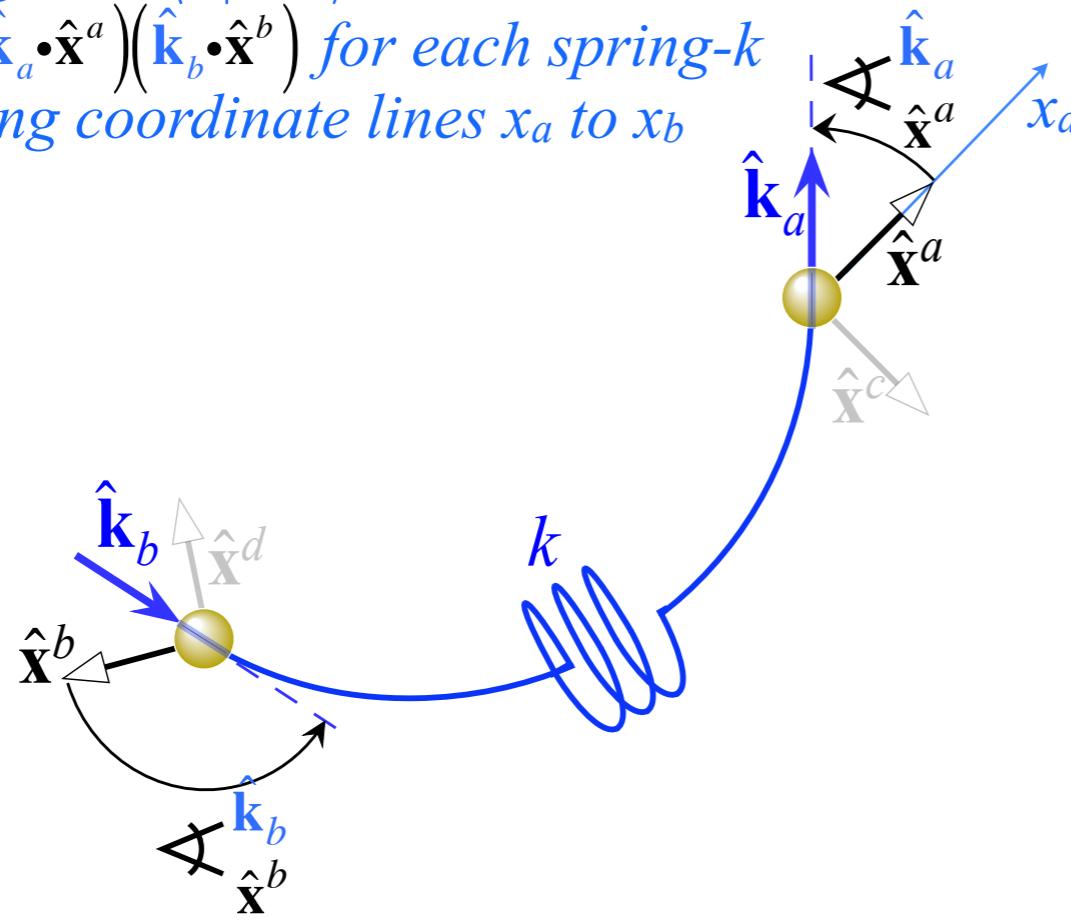
Harmonic potential  $V(\mathbf{x})$  is a quadratic  $K$ -form of coordinates  $x_a$  based on six  $D_3$ -labeled axes  $\hat{\mathbf{x}}^a$  or  $|a\rangle$ .

$$V(x) = \sum_{(k)} \frac{1}{2} \langle x | \mathbf{K} | x \rangle \quad \text{where: } |x\rangle = \sum_a x_a |a\rangle, \quad (a, b) = (1, r^1, r^2, i_1, i_2, i_3)$$

Each  $\mathbf{K}$  component  $K_{ab} = \langle a | \mathbf{K} | b \rangle$  is a sum over spring  $k$ -constants that connect axis- $\mathbf{x}^a$  to axis- $\mathbf{x}^b$  multiplied by factor  $(\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)(\hat{\mathbf{k}}_b \bullet \hat{\mathbf{x}}^b)$  for projecting spring- $k$ 's end vectors  $\hat{\mathbf{k}}_a$  and  $\hat{\mathbf{k}}_b$  onto  $\hat{\mathbf{x}}^a$  and  $\hat{\mathbf{x}}^b$  at respective connections.

*Coupling  $K_{ab} = \langle a | \mathbf{K} | b \rangle$*

*Sum  $k \cdot (\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)(\hat{\mathbf{k}}_b \bullet \hat{\mathbf{x}}^b)$  for each spring- $k$  connecting coordinate lines  $x_a$  to  $x_b$*



*Diagonal  $K_{aa} = \langle a | \mathbf{K} | a \rangle$*

*Sum  $-\frac{1}{2}k(\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)^2$  for each spring- $k$  connected to coordinate line  $x_a$*

## Molecular K-matrix construction

Classical modes are eigenvectors of force-field matrix  $K$  or operator  $\mathbf{K}$ .

Harmonic potential  $V(\mathbf{x})$  is a quadratic  $K$ -form of coordinates  $x_a$  based on six  $D_3$ -labeled axes  $\hat{\mathbf{x}}^a$  or  $|a\rangle$ .

$$V(x) = \sum_{(k)} \frac{1}{2} \langle x | \mathbf{K} | x \rangle \quad \text{where: } |x\rangle = \sum_a x_a |a\rangle, \quad (a, b) = (1, r^1, r^2, i_1, i_2, i_3)$$

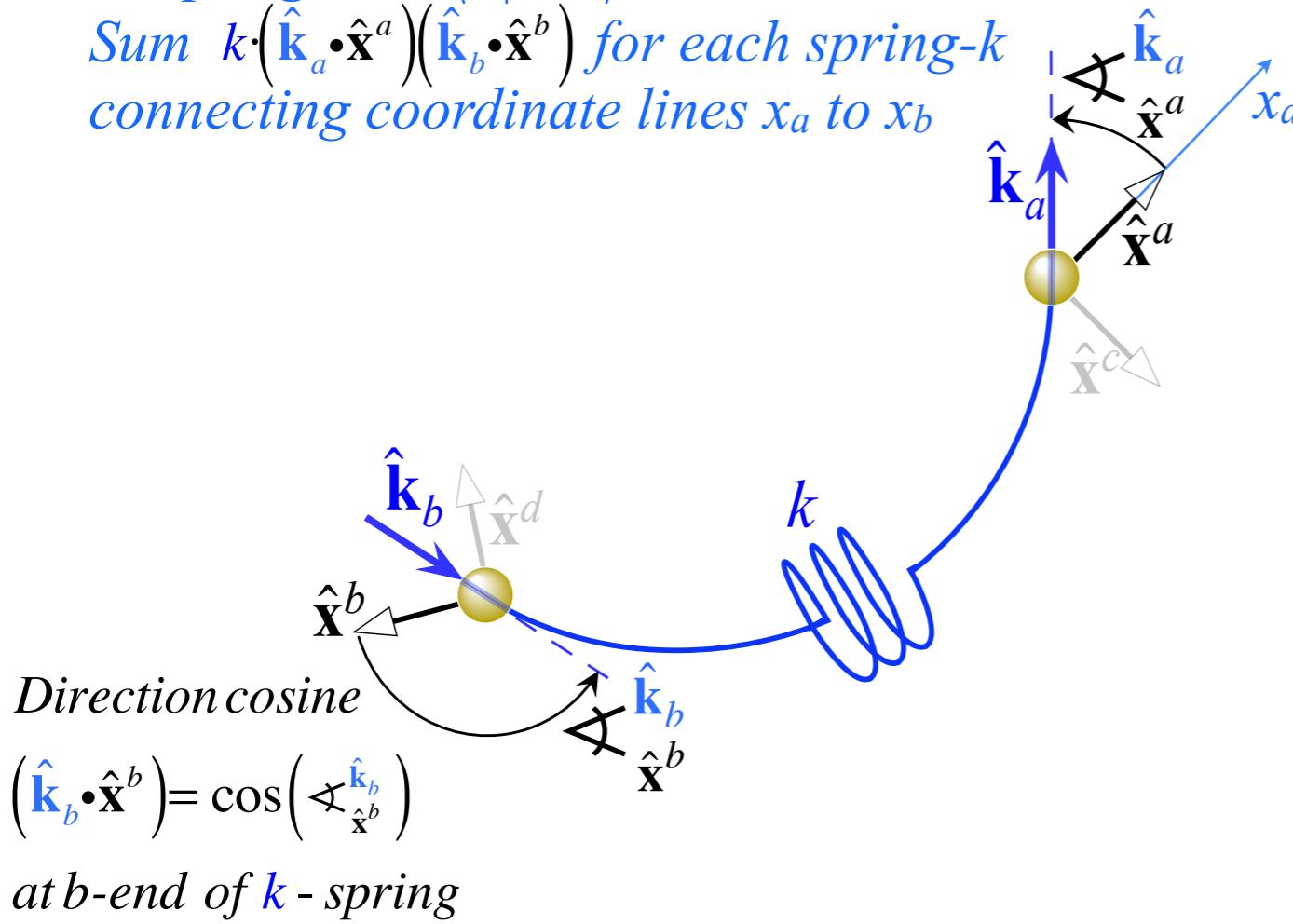
Each  $\mathbf{K}$  component  $K_{ab} = \langle a | \mathbf{K} | b \rangle$  is a sum over spring  $k$ -constants that connect axis- $\mathbf{x}^a$  to axis- $\mathbf{x}^b$  multiplied by factor  $(\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)(\hat{\mathbf{k}}_b \bullet \hat{\mathbf{x}}^b)$  for projecting spring- $k$ 's end vectors  $\hat{\mathbf{k}}_a$  and  $\hat{\mathbf{k}}_b$  onto  $\hat{\mathbf{x}}^a$  and  $\hat{\mathbf{x}}^b$  at respective connections.

*Coupling  $K_{ab} = \langle a | \mathbf{K} | b \rangle$*

*Sum  $k \cdot (\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)(\hat{\mathbf{k}}_b \bullet \hat{\mathbf{x}}^b)$  for each spring- $k$  connecting coordinate lines  $x_a$  to  $x_b$*

*Diagonal  $K_{aa} = \langle a | \mathbf{K} | a \rangle$*

*Sum  $-\frac{1}{2} k (\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)^2$  for each spring- $k$  connected to coordinate line  $x_a$*



## Molecular K-matrix construction

Classical modes are eigenvectors of force-field matrix  $K$  or operator  $\mathbf{K}$ .

Harmonic potential  $V(\mathbf{x})$  is a quadratic  $K$ -form of coordinates  $x_a$  based on six  $D_3$ -labeled axes  $\hat{\mathbf{x}}^a$  or  $|a\rangle$ .

$$V(x) = \sum_{(k)} \frac{1}{2} \langle x | \mathbf{K} | x \rangle \quad \text{where: } |x\rangle = \sum_a x_a |a\rangle, \quad (a, b) = (1, r^1, r^2, i_1, i_2, i_3)$$

Each  $\mathbf{K}$  component  $K_{ab} = \langle a | \mathbf{K} | b \rangle$  is a sum over spring  $k$ -constants that connect axis- $\mathbf{x}^a$  to axis- $\mathbf{x}^b$  multiplied by factor  $(\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)(\hat{\mathbf{k}}_b \bullet \hat{\mathbf{x}}^b)$  for projecting spring- $k$ 's end vectors  $\hat{\mathbf{k}}_a$  and  $\hat{\mathbf{k}}_b$  onto  $\hat{\mathbf{x}}^a$  and  $\hat{\mathbf{x}}^b$  at respective connections.

## Molecular K-matrix construction

Classical modes are eigenvectors of force-field matrix  $K$  or operator  $\mathbf{K}$ .

Harmonic potential  $V(\mathbf{x})$  is a quadratic  $K$ -form of coordinates  $x_a$  based on six  $D_3$ -labeled axes  $\hat{\mathbf{x}}^a$  or  $|a\rangle$ .

$$V(x) = \sum_{(k)} \frac{1}{2} \langle x | \mathbf{K} | x \rangle \quad \text{where: } |x\rangle = \sum_a x_a |a\rangle, \quad (a, b) = (1, r^1, r^2, i_1, i_2, i_3)$$

Each  $\mathbf{K}$  component  $K_{ab} = \langle a | \mathbf{K} | b \rangle$  is a sum over spring  $k$ -constants that connect axis- $\mathbf{x}^a$  to axis- $\mathbf{x}^b$  multiplied by factor  $(\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)(\hat{\mathbf{k}}_b \bullet \hat{\mathbf{x}}^b)$  for projecting spring- $k$ 's end vectors  $\hat{\mathbf{k}}_a$  and  $\hat{\mathbf{k}}_b$  onto  $\hat{\mathbf{x}}^a$  and  $\hat{\mathbf{x}}^b$  at respective connections.

$$V(x) = \frac{1}{2} \sum_{a,b} K_{ab} x_a x_a \quad \text{where: } K_{ab} = \begin{cases} \sum_{(k)} \frac{k}{2} (\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)^2 & \text{if } a = b \\ - \sum_{(k)} k (\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)(\hat{\mathbf{k}}_b \bullet \hat{\mathbf{x}}^b) & \text{if } a \neq b \end{cases}$$

## Molecular K-matrix construction

Classical modes are eigenvectors of force-field matrix  $K$  or operator  $\mathbf{K}$ .

Harmonic potential  $V(\mathbf{x})$  is a quadratic  $K$ -form of coordinates  $x_a$  based on six  $D_3$ -labeled axes  $\hat{\mathbf{x}}^a$  or  $|a\rangle$ .

$$V(x) = \sum_{(k)} \frac{1}{2} \langle x | \mathbf{K} | x \rangle \quad \text{where: } |x\rangle = \sum_a x_a |a\rangle, \quad (a, b) = (1, r^1, r^2, i_1, i_2, i_3)$$

Each  $\mathbf{K}$  component  $K_{ab} = \langle a | \mathbf{K} | b \rangle$  is a sum over spring  $k$ -constants that connect axis- $\mathbf{x}^a$  to axis- $\mathbf{x}^b$  multiplied by factor  $(\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)(\hat{\mathbf{k}}_b \bullet \hat{\mathbf{x}}^b)$  for projecting spring- $k$ 's end vectors  $\hat{\mathbf{k}}_a$  and  $\hat{\mathbf{k}}_b$  onto  $\hat{\mathbf{x}}^a$  and  $\hat{\mathbf{x}}^b$  at respective connections.

$$V(x) = \frac{1}{2} \sum_{a,b} K_{ab} x_a x_b \quad \text{where: } K_{ab} = \begin{cases} \sum_{(k)} \frac{k}{2} (\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)^2 & \text{if } a = b \\ - \sum_{(k)} k (\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)(\hat{\mathbf{k}}_b \bullet \hat{\mathbf{x}}^b) & \text{if } a \neq b \end{cases}$$

$$\begin{aligned} V(x) &= \sum_{(k)} \frac{k}{2} (\Delta \ell_k)^2 = \sum_{(k)} \frac{k}{2} \sum_{a,b} (\hat{\mathbf{k}}_a \bullet \mathbf{x}^a - \hat{\mathbf{k}}_b \bullet \mathbf{x}^b)^2 \\ &= \sum_{(k)} \frac{k}{2} \sum_a (\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)^2 x_a^2 - \sum_{(k)} k \sum_{a \neq b} (\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)(\hat{\mathbf{k}}_b \bullet \hat{\mathbf{x}}^b) x_a x_b \end{aligned}$$

## Molecular K-matrix construction

Classical modes are eigenvectors of force-field matrix  $K$  or operator  $\mathbf{K}$ .

Harmonic potential  $V(\mathbf{x})$  is a quadratic  $K$ -form of coordinates  $x_a$  based on six  $D_3$ -labeled axes  $\hat{\mathbf{x}}^a$  or  $|a\rangle$ .

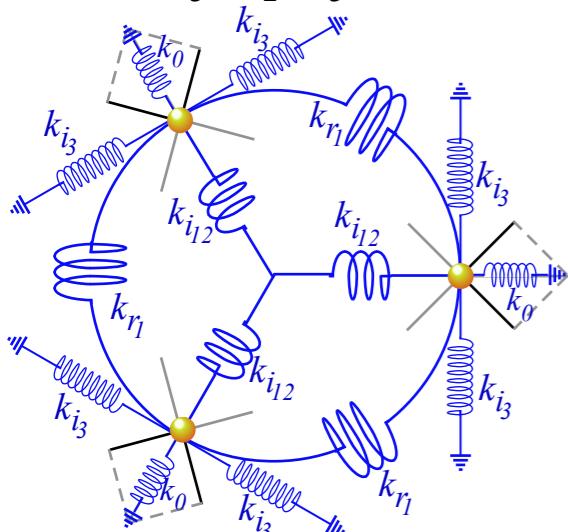
$$V(x) = \sum_{(k)} \frac{1}{2} \langle x | \mathbf{K} | x \rangle \quad \text{where: } |x\rangle = \sum_a x_a |a\rangle, \quad (a, b) = (1, r^1, r^2, i_1, i_2, i_3)$$

Each  $\mathbf{K}$  component  $K_{ab} = \langle a | \mathbf{K} | b \rangle$  is a sum over spring  $k$ -constants that connect axis- $\mathbf{x}^a$  to axis- $\mathbf{x}^b$  multiplied by factor  $(\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)(\hat{\mathbf{k}}_b \bullet \hat{\mathbf{x}}^b)$  for projecting spring- $k$ 's end vectors  $\hat{\mathbf{k}}_a$  and  $\hat{\mathbf{k}}_b$  onto  $\hat{\mathbf{x}}^a$  and  $\hat{\mathbf{x}}^b$  at respective connections.

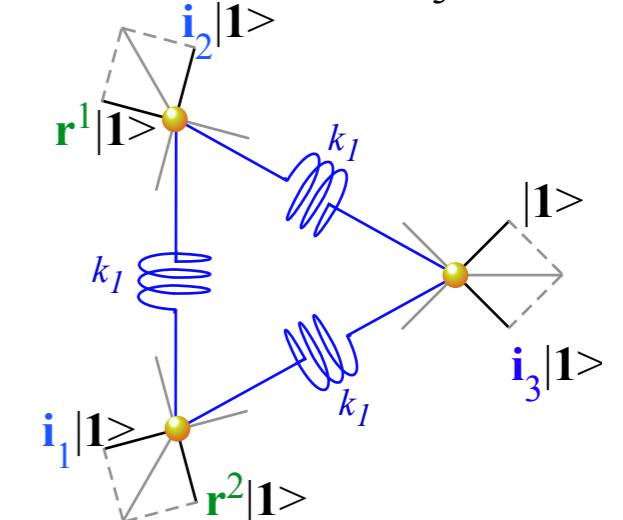
$$V(x) = \frac{1}{2} \sum_{a,b} K_{ab} x_a x_b \quad \text{where: } K_{ab} = \begin{cases} \sum_{(k)} \frac{k}{2} (\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)^2 & \text{if } a = b \\ - \sum_{(k)} k (\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)(\hat{\mathbf{k}}_b \bullet \hat{\mathbf{x}}^b) & \text{if } a \neq b \end{cases}$$

$$\begin{aligned} V(x) &= \sum_{(k)} \frac{k}{2} (\Delta \ell_k)^2 = \sum_{(k)} \frac{k}{2} \sum_{a,b} (\hat{\mathbf{k}}_a \bullet \mathbf{x}^a - \hat{\mathbf{k}}_b \bullet \mathbf{x}^b)^2 \\ &= \sum_{(k)} \frac{k}{2} \sum_a (\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)^2 x_a^2 - \sum_{(k)} k \sum_{a \neq b} (\hat{\mathbf{k}}_a \bullet \hat{\mathbf{x}}^a)(\hat{\mathbf{k}}_b \bullet \hat{\mathbf{x}}^b) x_a x_b \end{aligned}$$

Local  $D_3$   $C_2 \supset (i_3)$  model



Direct connection  $D_3$  model



*Review: Hamiltonian local-symmetry eigensolution in global and local  $|\mathbf{P}^{(\mu)}\rangle$ -basis*

### *Molecular vibrational modes vs. Hamiltonian eigenmodes*

*Molecular K-matrix construction*



*$D_3 \supset C_2(i_3)$  local-symmetry K-matrix eigensolutions*



*$D_3$ -direct-connection K-matrix eigensolutions*

*$D_3 \supset C_3(\mathbf{r}^{\pm l})$  local symmetry K-matrix eigensolutions*

### *Applied symmetry reduction and splitting*

*Subduced irep  $D^\alpha(D^3) \downarrow C_2 = d^{0_2} \oplus d^{l_2} \oplus \dots$  correlation*

*Subduced irep  $D^\alpha(D^3) \downarrow C_3 = d^{0_3} \oplus d^{l_3} \oplus \dots$  correlation*

### *Spontaneous symmetry breaking and clustering: Frobenius Reciprocity , band structure*

*Induced rep  $d^a(C_2) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus \dots$  correlation*

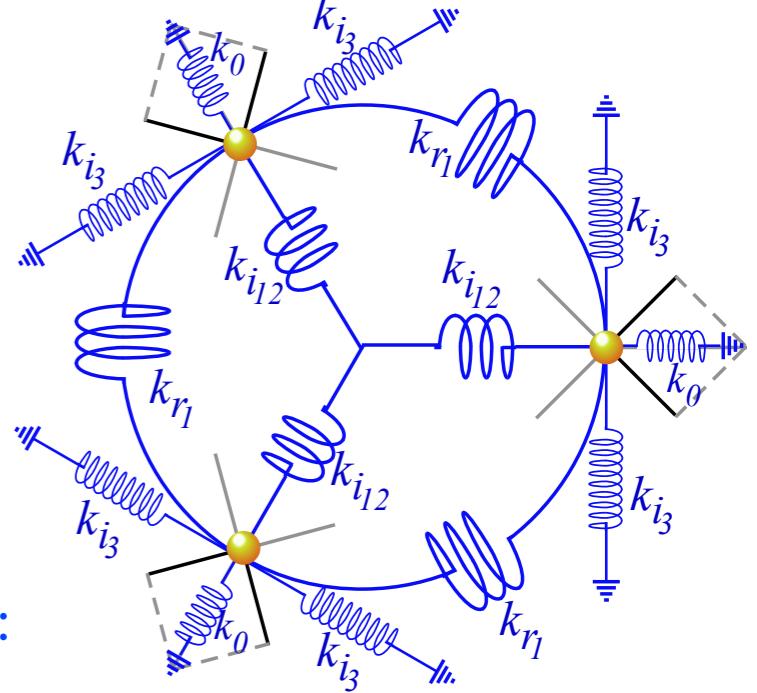
*Induced rep  $d^a(C_3) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus \dots$  correlation*

### *$D_6$ symmetry and Hexagonal Bands*

*Cross product of the  $C_2$  and  $D_3$  characters gives all  $D_6 = D_3 \times C_2$  characters and ireps*

*D<sub>3</sub>⊃C<sub>2</sub>(i<sub>3</sub>) local-symmetry vibrational K-matrix eigensolutions*

*Local D<sub>3</sub> C<sub>2</sub>▷(i<sub>3</sub>) model*



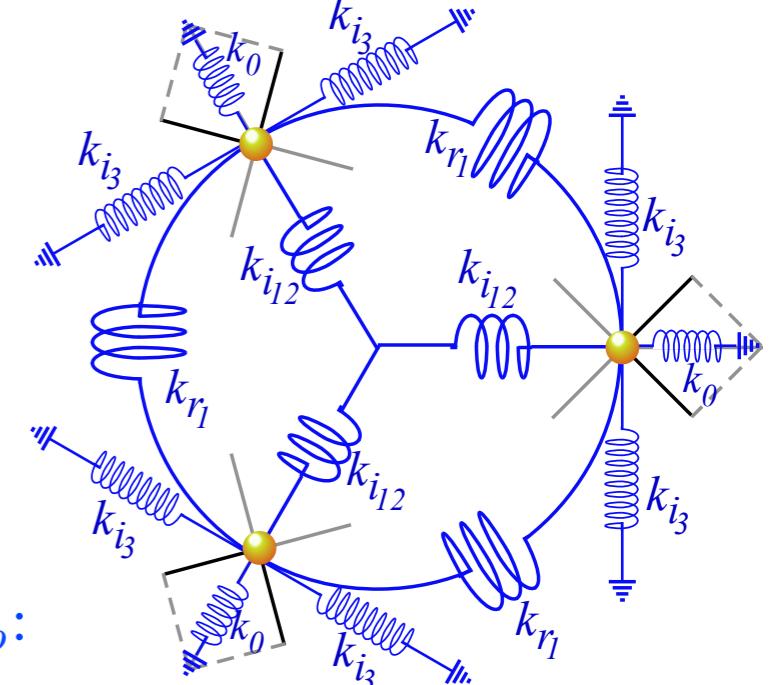
*D<sub>3</sub>⊃C<sub>2</sub>(i<sub>3</sub>) local-symmetry vibrational K-matrix*

1<sup>st</sup>-row parameters  $g_b = \langle \mathbf{1} | \mathbf{K} | \mathbf{g}_b \rangle = K_{1b}$  of the force matrix  $K_{ab}$ :

$D_3 \supset C_2(i_3)$  model has internal [  $k_r$  (angular),  $k_i$  (radial) ] and external [  $k_3$  (angular),  $k_0$  (radial) ] constants between masses and lab frame.

## $D_3 \supset C_2(i_3)$ local-symmetry vibrational K-matrix eigensolutions

## Local $D_3$ $C_2^{\supset}(i_3)$ model



## $D_3 \supset C_2(i_3)$ local-symmetry vibrational K-matrix

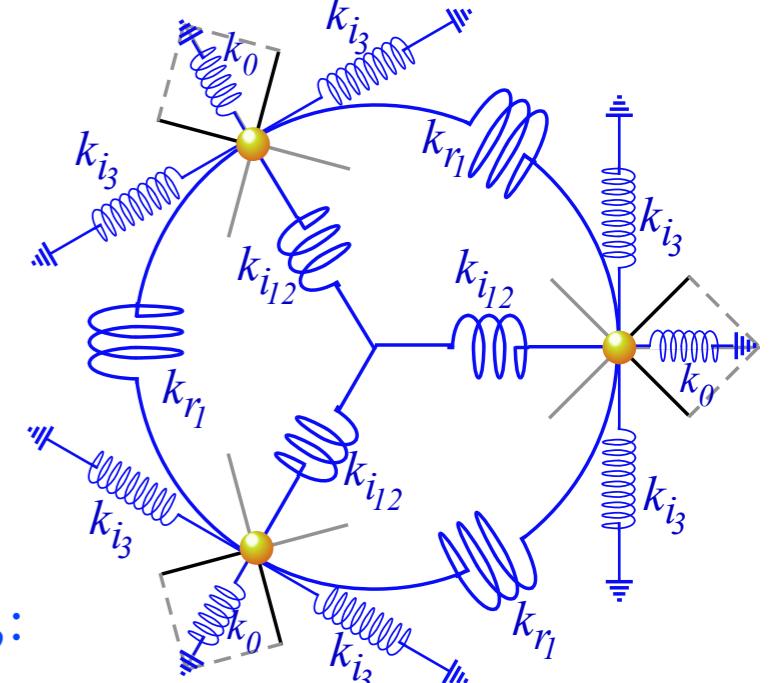
1<sup>st</sup>-row parameters  $g_b = \langle \mathbf{1} | \mathbf{K} | \mathbf{g}_b \rangle = K_{1b}$  of the force matrix  $K_{ab}$ :

$D_3 \supset C_2(i_3)$  model has internal [ $k_r$ (angular), $k_i$ (radial)] and external [ $k_3$ (angular), $k_0$ (radial)] constants between masses and lab frame.

$ g_b\rangle$	$ \mathbf{1}\rangle$	$ \mathbf{r}^1\rangle$	$ \mathbf{r}^2\rangle$	$ \mathbf{i}_1\rangle$	$ \mathbf{i}_2\rangle$	$ \mathbf{i}_3\rangle$
$\langle \mathbf{1}   \mathbf{K}   g_b \rangle =$	$k_i/2$ $+k_r$ $+k_3$ $+k_0/2$	$k_i/2$ $-k_r/2$ $+0$ $+0$	$k_i/2$ $-k_r/2$ $+0$ $+0$	$k_i/2$ $+k_r/2$ $+0$ $+0$	$k_i/2$ $+k_r/2$ $+0$ $+0$	$k_i/2$ $-k_r$ $-k_3$ $+k_0/2$

## $D_3 \supset C_2(i_3)$ local-symmetry vibrational K-matrix eigensolutions

## Local $D_3$ $C_2^{\supset}(i_3)$ model



## $D_3 \supset C_2(i_3)$ local-symmetry vibrational K-matrix

1<sup>st</sup>-row parameters  $g_b = \langle \mathbf{1} | \mathbf{K} | \mathbf{g}_b \rangle = K_{1b}$  of the force matrix  $K_{ab}$ :

$D_3 \supset C_2(i_3)$  model has internal [ $k_r$ (angular), $k_i$ (radial)] and external [ $k_3$ (angular), $k_0$ (radial)] constants between masses and lab frame.

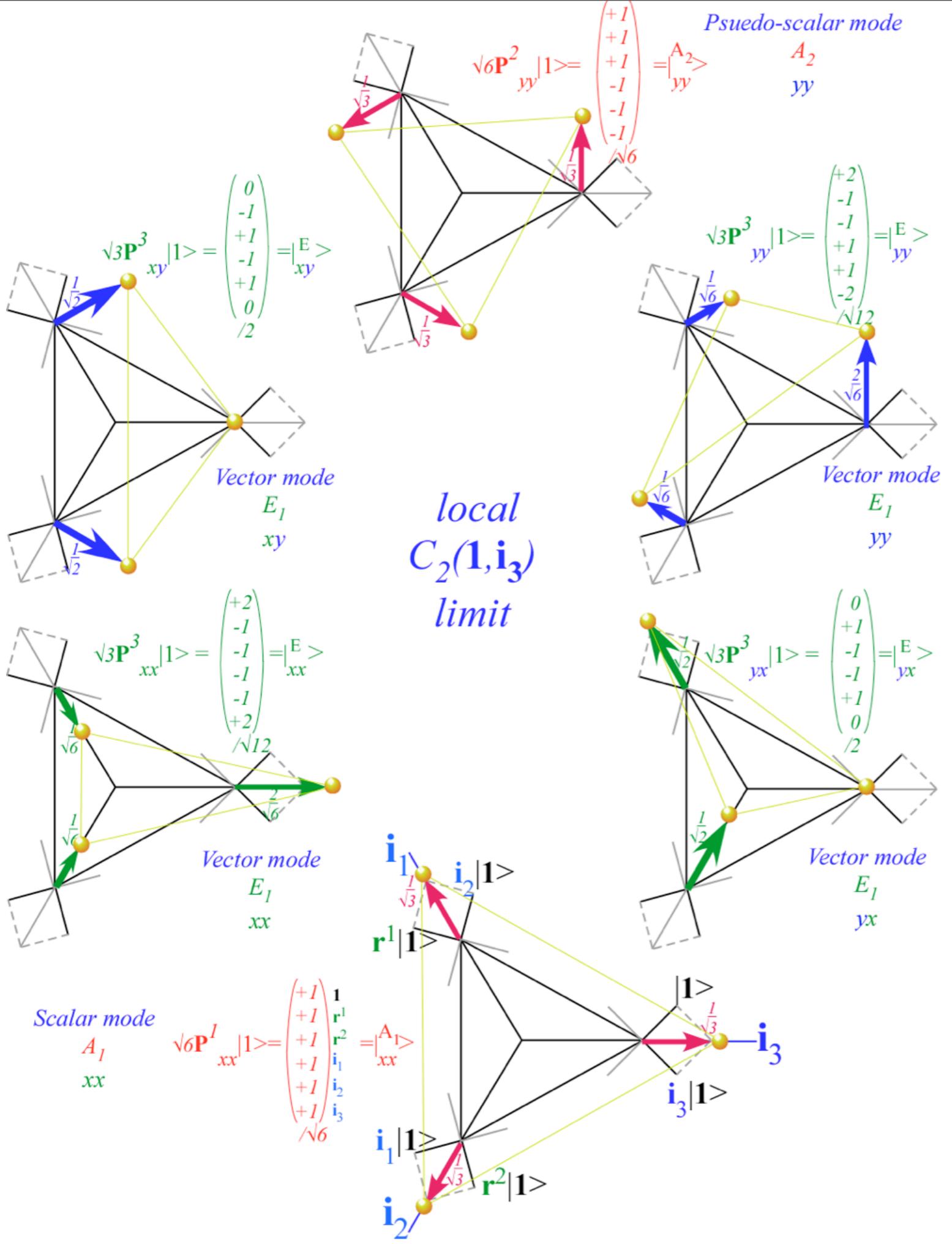
$ g_b\rangle$	$ \mathbf{1}\rangle$	$ \mathbf{r}^1\rangle$	$ \mathbf{r}^2\rangle$	$ \mathbf{i}_1\rangle$	$ \mathbf{i}_2\rangle$	$ \mathbf{i}_3\rangle$
$\langle \mathbf{1}   \mathbf{K}   g_b \rangle =$	$k_i/2$ $+k_r$ $+k_3$ $+k_0/2$	$k_i/2$ $-k_r/2$ $+0$ $+0$	$k_i/2$ $-k_r/2$ $+0$ $+0$	$k_i/2$ $+k_r/2$ $+0$ $+0$	$k_i/2$ $+k_r/2$ $+0$ $+0$	$k_i/2$ $-k_r$ $-k_3$ $+k_0/2$

## $D_3 \supset C_2(i_3)$ local-symmetry vibrational K-matrix eigenvalues $K_m/M = \omega_m^2$

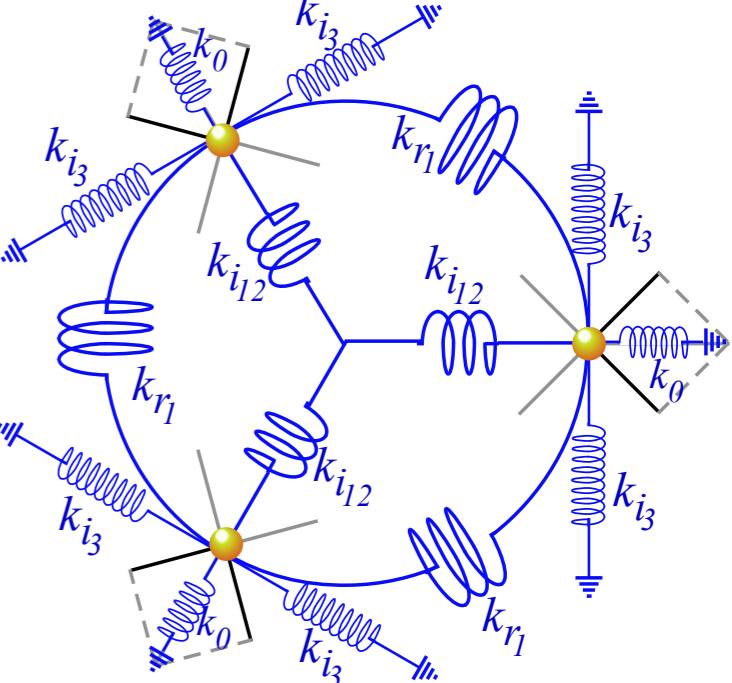
$$K_{xx}^{A_1} = r_0 + r_1 + r_1^* + i_1 + i_2 + i_3 = k_0 + 3k_i$$

$$K_{yy}^{A_2} = r_0 + r_1 + r_1^* - i_1 - i_2 - i_3 = 3k_3$$

$$\begin{pmatrix} K_{xx}^{E_1} & K_{xy}^{E_1} \\ K_{yx}^{E_1} & K_{yy}^{E_1} \end{pmatrix} = \frac{1}{2} \begin{pmatrix} 2r_0 - r_1 - r_1^* - i_1 - i_2 + 2i_3 & \sqrt{3}(-r_1 + r_1^* - i_1 + i_2) \\ \sqrt{3}(-r_1^* + r_1 - i_1 + i_2) & 2r_0 - r_1 - r_1^* + i_1 + i_2 - 2i_3 \end{pmatrix} = \begin{pmatrix} k_0 & 0 \\ 0 & k_3 + 2k_r \end{pmatrix}$$



*Local  $D_3$   $C_2(i_3)$  model*



*Review: Hamiltonian local-symmetry eigensolution in global and local  $|\mathbf{P}^{(\mu)}\rangle$ -basis*

### *Molecular vibrational modes vs. Hamiltonian eigenmodes*

*Molecular K-matrix construction*

*$D_3 \supset C_2(i_3)$  local-symmetry K-matrix eigensolutions*

*$D_3$ -direct-connection K-matrix eigensolutions*

*$D_3 \supset C_3(\mathbf{r}^{\pm l})$  local symmetry K-matrix eigensolutions*



### *Applied symmetry reduction and splitting*

*Subduced irep  $D^\alpha(D^3) \downarrow C_2 = d^{0_2} \oplus d^{l_2} \oplus ..$  correlation*

*Subduced irep  $D^\alpha(D^3) \downarrow C_3 = d^{0_3} \oplus d^{l_3} \oplus ..$  correlation*

### *Spontaneous symmetry breaking and clustering: Frobenius Reciprocity , band structure*

*Induced rep  $d^a(C_2) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus ..$  correlation*

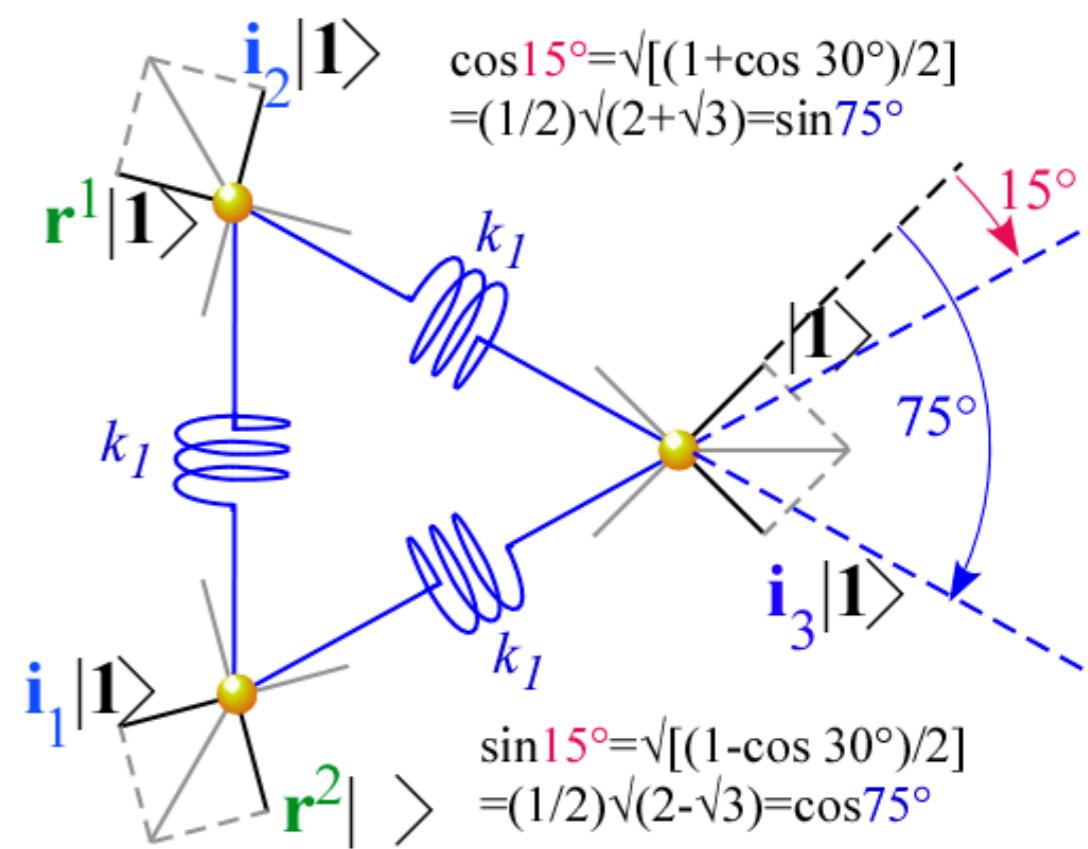
*Induced rep  $d^a(C_3) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus ..$  correlation*

### *$D_6$ symmetry and Hexagonal Bands*

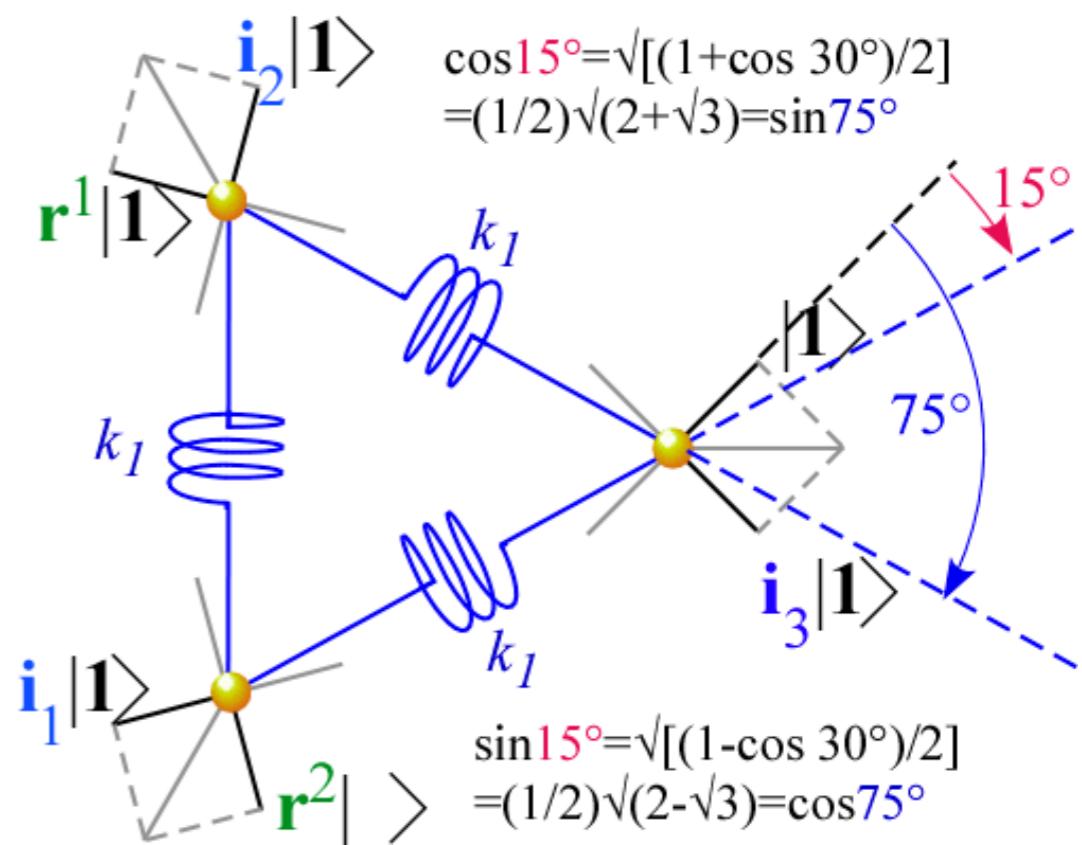
*Cross product of the  $C_2$  and  $D_3$  characters gives all  $D_6 = D_3 \times C_2$  characters and ireps*

## *D<sub>3</sub>-direct-connection K-matrix eigensolutions*

## *D<sub>3</sub>-direct-connection vibrational K-matrix*



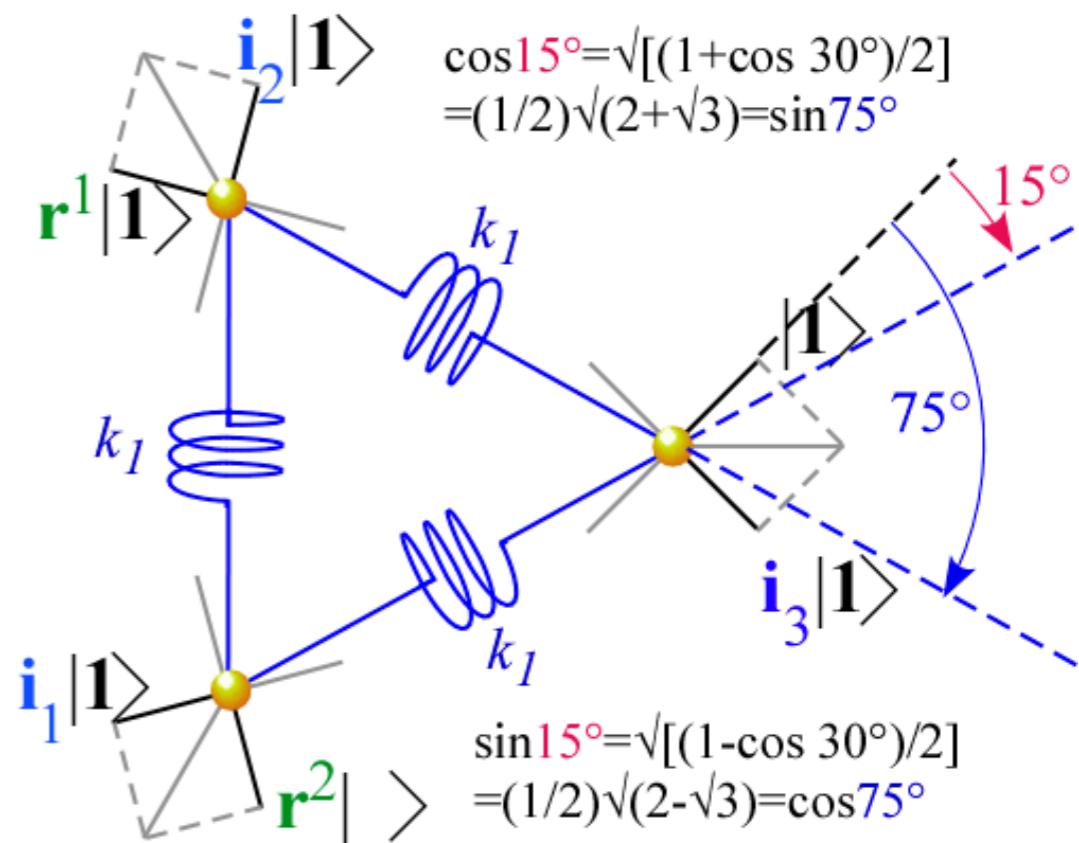
## D<sub>3</sub>-direct-connection K-matrix eigensolutions



## D<sub>3</sub>-direct-connection vibrational K-matrix

$ g_b\rangle$	$ \mathbf{1}\rangle$	$ \mathbf{r}^1\rangle$	$ \mathbf{r}^2\rangle$	$ \mathbf{i}_1\rangle$	$ \mathbf{i}_2\rangle$	$ \mathbf{i}_3\rangle$
$\langle \mathbf{1}   \mathbf{K}   g_b \rangle =$	$k_1(\cos^2 75^\circ + \cos^2 15^\circ) = k_1$	$k_1 \cos 75^\circ \cdot \cos 15^\circ = \frac{k_1}{4}$	$k_1 \cos 15^\circ \cdot \cos 75^\circ = \frac{k_1}{4}$	$k_1 \cos 15^\circ \cdot \cos 15^\circ = \frac{k_1(2 - \sqrt{3})}{4}$	$k_1 \cos 75^\circ \cdot \cos 75^\circ = \frac{k_1(2 + \sqrt{3})}{4}$	$k_1(\cos^2 75^\circ - \cos^2 15^\circ) = \frac{k_1}{2}$

## D<sub>3</sub>-direct-connection K-matrix eigensolutions



## D<sub>3</sub>-direct-connection vibrational K-matrix

$ g_b\rangle$	$ 1\rangle$	$ r^1\rangle$	$ r^2\rangle$	$ i_1\rangle$	$ i_2\rangle$	$ i_3\rangle$
$\langle 1  \mathbf{K}  g_b\rangle =$	$k_1(\cos^2 75^\circ + \cos^2 15^\circ) = k_1$	$k_1 \cos 75^\circ \cdot \cos 15^\circ = \frac{k_1}{4}$	$k_1 \cos 15^\circ \cdot \cos 75^\circ = \frac{k_1}{4}$	$k_1 \cos 15^\circ \cdot \cos 15^\circ = \frac{k_1(2 - \sqrt{3})}{4}$	$k_1 \cos 75^\circ \cdot \cos 75^\circ = \frac{k_1(2 + \sqrt{3})}{4}$	$k_1(\cos^2 75^\circ - \cos^2 15^\circ) = \frac{k_1}{2}$

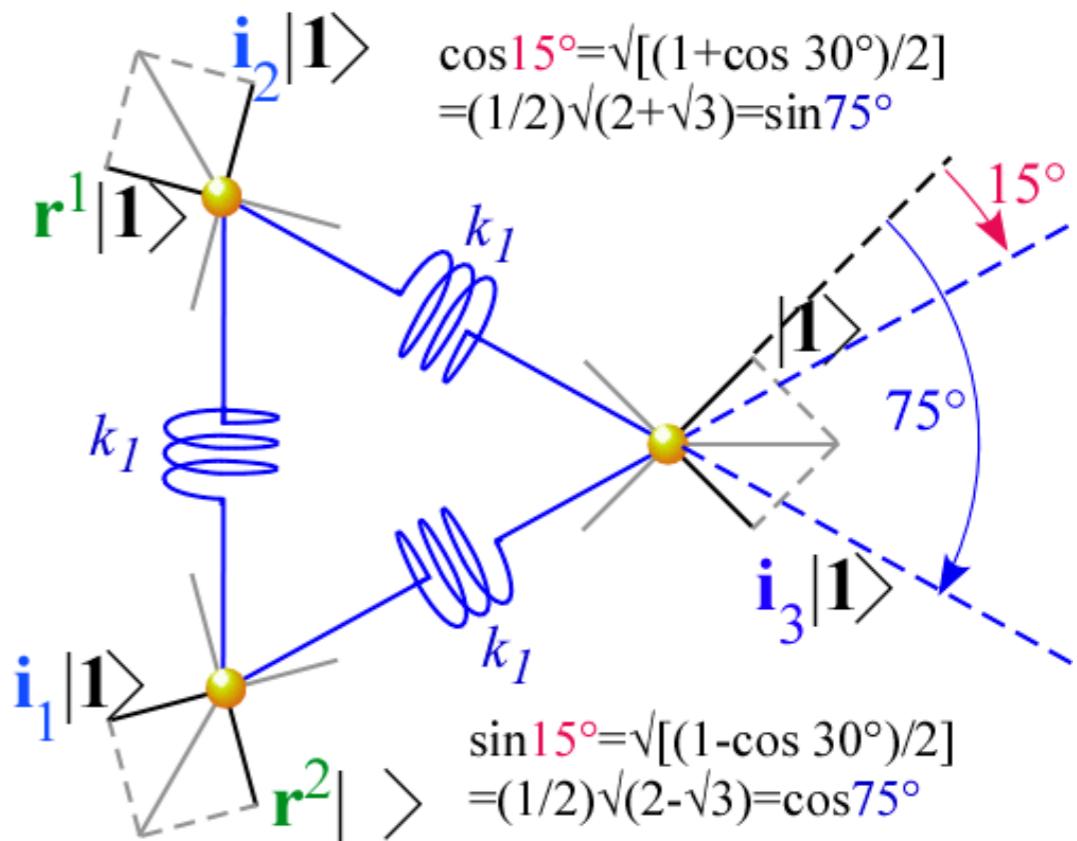
## D<sub>3</sub>-direct-connection vibrational K-matrix eigenvalues $K_m/M = \omega_m^2$

$$K_{xx}^{A_1} = 3k_1$$

$$K_{yy}^{A_2} = 0$$

$$\begin{pmatrix} K_{xx}^{E_1} & K_{xy}^{E_1} \\ K_{yx}^{E_1} & K_{yy}^{E_1} \end{pmatrix} = \begin{pmatrix} \frac{3k_1}{4} & \frac{3k_1}{4} \\ \frac{3k_1}{4} & \frac{3k_1}{4} \end{pmatrix}$$

## D<sub>3</sub>-direct-connection K-matrix eigensolutions



## D<sub>3</sub>-direct-connection vibrational K-matrix

$ g_b\rangle$	$ \mathbf{1}\rangle$	$ \mathbf{r}^1\rangle$	$ \mathbf{r}^2\rangle$	$ \mathbf{i}_1\rangle$	$ \mathbf{i}_2\rangle$	$ \mathbf{i}_3\rangle$
$\langle \mathbf{1}   \mathbf{K}   g_b \rangle =$	$k_1(\cos^2 75^\circ + \cos^2 15^\circ) = k_1$	$k_1 \cos 75^\circ \cdot \cos 15^\circ = \frac{k_1}{4}$	$k_1 \cos 15^\circ \cdot \cos 75^\circ = \frac{k_1}{4}$	$k_1 \cos 15^\circ \cdot \cos 15^\circ = \frac{k_1(2 - \sqrt{3})}{4}$	$k_1 \cos 75^\circ \cdot \cos 75^\circ = \frac{k_1(2 + \sqrt{3})}{4}$	$k_1(\cos^2 75^\circ - \cos^2 15^\circ) = \frac{k_1}{2}$

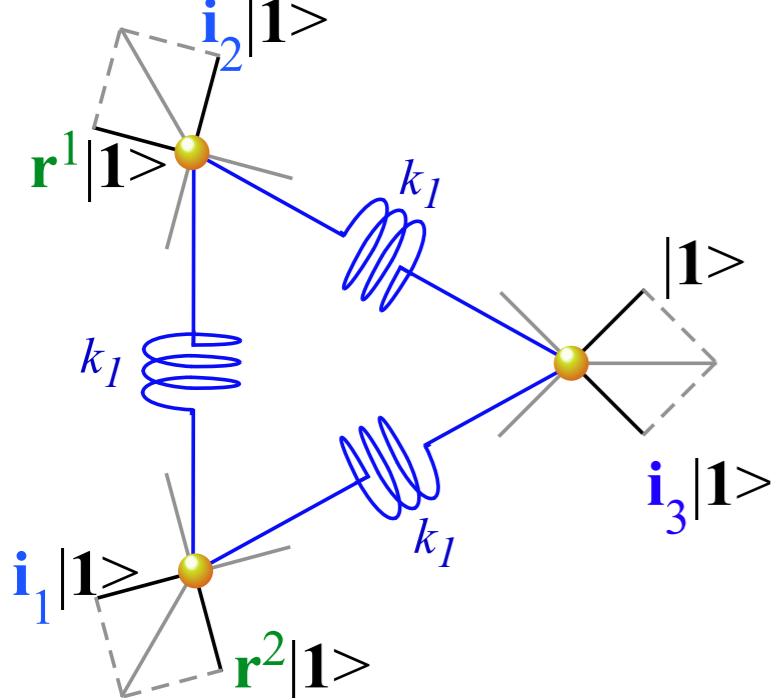
## D<sub>3</sub>-direct-connection vibrational K-matrix eigenvalues $K_m/M = \omega_m^2$

$$\begin{aligned}
 K_{xx}^{A_1} &= 3k_1 \\
 K_{yy}^{A_2} &= 0 \\
 \begin{pmatrix} K_{xx}^{E_1} & K_{xy}^{E_1} \\ K_{yx}^{E_1} & K_{yy}^{E_1} \end{pmatrix} &= \begin{pmatrix} \frac{3k_1}{4} & \frac{3k_1}{4} \\ \frac{3k_1}{4} & \frac{3k_1}{4} \end{pmatrix}
 \end{aligned}$$

$E_1$  Eigenvectors in terms of  $D_3 \supset C_2(i_3)$   $E_1$ -vectors

$$\begin{aligned}
 \mathbf{K} \begin{pmatrix} E_1 \\ g(+ \ ) \end{pmatrix} &= \mathbf{K} \left( \begin{pmatrix} E_1 \\ g\mathbf{x} \end{pmatrix} + \begin{pmatrix} E_1 \\ g\mathbf{y} \end{pmatrix} \right) \frac{1}{\sqrt{2}} = \frac{3k_1}{2} \begin{pmatrix} E_1 \\ g(+ \ ) \end{pmatrix}, \\
 \mathbf{K} \begin{pmatrix} E_1 \\ g(- \ ) \end{pmatrix} &= \mathbf{K} \left( \begin{pmatrix} E_1 \\ g\mathbf{x} \end{pmatrix} - \begin{pmatrix} E_1 \\ g\mathbf{y} \end{pmatrix} \right) \frac{1}{\sqrt{2}} = 0 \begin{pmatrix} E_1 \\ g(- \ ) \end{pmatrix}, \quad g = (\mathbf{x} \text{ or } \mathbf{y}).
 \end{aligned}$$

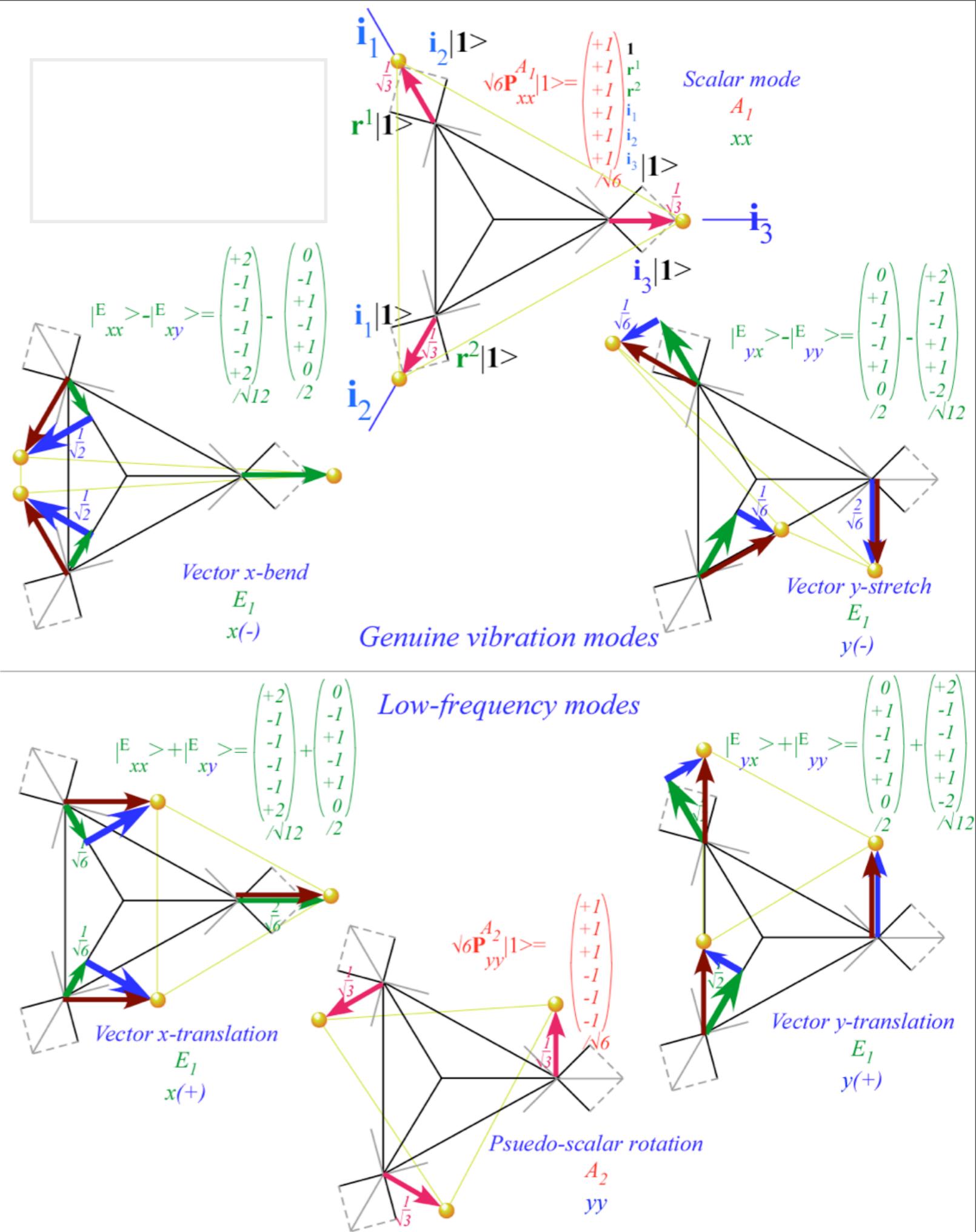
# Direct connection $D_3$ model



$$K_{xx}^{A_1} = 3k_1$$

$$K_{yy}^{A_2} = 0$$

$$\begin{pmatrix} K_{xx}^{E_1} & K_{xy}^{E_1} \\ K_{yx}^{E_1} & K_{yy}^{E_1} \end{pmatrix} = \begin{pmatrix} \frac{3k_1}{4} & \frac{3k_1}{4} \\ \frac{3k_1}{4} & \frac{3k_1}{4} \end{pmatrix}$$



*Review: Hamiltonian local-symmetry eigensolution in global and local  $|\mathbf{P}^{(\mu)}\rangle$ -basis*

### *Molecular vibrational modes vs. Hamiltonian eigenmodes*

*Molecular K-matrix construction*

*$D_3 \supset C_2(i_3)$  local-symmetry K-matrix eigensolutions*

*$D_3$ -direct-connection K-matrix eigensolutions*

*$D_3 \supset C_3(\mathbf{r}^{\pm l})$  local symmetry K-matrix eigensolutions*



### *Applied symmetry reduction and splitting*

*Subduced irep  $D^\alpha(D^3) \downarrow C_2 = d^{0_2} \oplus d^{l_2} \oplus \dots$  correlation*

*Subduced irep  $D^\alpha(D^3) \downarrow C_3 = d^{0_3} \oplus d^{l_3} \oplus \dots$  correlation*

### *Spontaneous symmetry breaking and clustering: Frobenius Reciprocity , band structure*

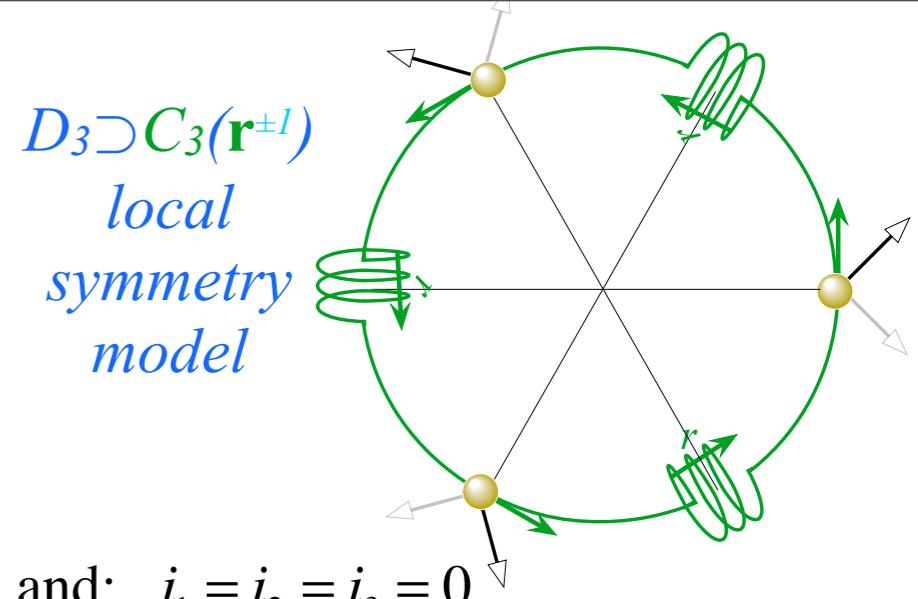
*Induced rep  $d^a(C_2) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus \dots$  correlation*

*Induced rep  $d^a(C_3) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus \dots$  correlation*

### *$D_6$ symmetry and Hexagonal Bands*

*Cross product of the  $C_2$  and  $D_3$  characters gives all  $D_6 = D_3 \times C_2$  characters and irreps*

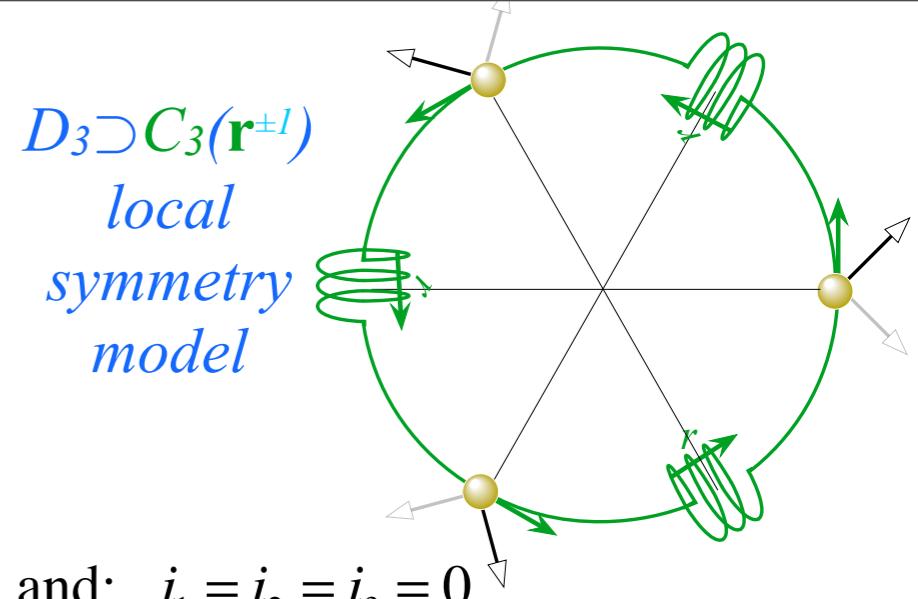
# $D_3 \supset C_3(\mathbf{r}^{\pm l})$ local symmetry K-matrix eigensolutions



$D_3 \supset C_3(\mathbf{r}^{\pm l})$  local symmetry vibrational K-matrix Set:  $r_1 = \mathbf{r} = -r_2^*$ , and:  $i_1 = i_2 = i_3 = 0$

$$\begin{aligned}
 K_{xx}^{A_1} &= r_0 + \mathbf{r}_1 + \mathbf{r}_1^* + i_1 + i_2 + i_3 & = r_0 \\
 K_{yy}^{A_2} &= r_0 + \mathbf{r}_1 + \mathbf{r}_1^* - i_1 - i_2 - i_3 & = r_0 \\
 \begin{pmatrix} K_{xx}^{E_1} & K_{xy}^{E_1} \\ K_{yx}^{E_1} & K_{yy}^{E_1} \end{pmatrix} &= \frac{1}{2} \begin{pmatrix} 2r_0 - \mathbf{r}_1 - \mathbf{r}_1^* - i_1 - i_2 + 2i_3 & \sqrt{3}(-\mathbf{r}_1 + \mathbf{r}_1^* - i_1 + i_2) \\ \sqrt{3}(-\mathbf{r}_1^* + \mathbf{r}_1 - i_1 + i_2) & 2r_0 - \mathbf{r}_1 - \mathbf{r}_1^* + i_1 + i_2 - 2i_3 \end{pmatrix}_{\substack{r_1 = \mathbf{r} = -r_2^* \\ i_1 = i_2 = i_3 = 0}} &= \begin{pmatrix} r_0 & -i\mathbf{r} \frac{\sqrt{3}}{2} \\ +i\mathbf{r} \frac{\sqrt{3}}{2} & r_0 \end{pmatrix}
 \end{aligned}$$

# $D_3 \supset C_3(\mathbf{r}^{\pm l})$ local symmetry K-matrix eigensolutions



$D_3 \supset C_3(\mathbf{r}^{\pm l})$  local symmetry vibrational K-matrix Set:  $r_1 = \mathbf{r} = -r_2^*$ , and:  $i_1 = i_2 = i_3 = 0$

$$\begin{aligned}
 K_{xx}^{A_1} &= r_0 + \mathbf{r}_1 + \mathbf{r}_1^* + i_1 + i_2 + i_3 & = r_0 \\
 K_{yy}^{A_2} &= r_0 + \mathbf{r}_1 + \mathbf{r}_1^* - i_1 - i_2 - i_3 & = r_0 \\
 \begin{pmatrix} K_{xx}^{E_1} & K_{xy}^{E_1} \\ K_{yx}^{E_1} & K_{yy}^{E_1} \end{pmatrix} &= \frac{1}{2} \begin{pmatrix} 2r_0 - \mathbf{r}_1 - \mathbf{r}_1^* - i_1 - i_2 + 2i_3 & \sqrt{3}(-\mathbf{r}_1 + \mathbf{r}_1^* - i_1 + i_2) \\ \sqrt{3}(-\mathbf{r}_1^* + \mathbf{r}_1 - i_1 + i_2) & 2r_0 - \mathbf{r}_1 - \mathbf{r}_1^* + i_1 + i_2 - 2i_3 \end{pmatrix} &= \begin{pmatrix} r_0 & -i\mathbf{r} \frac{\sqrt{3}}{2} \\ +i\mathbf{r} \frac{\sqrt{3}}{2} & r_0 \end{pmatrix} \\
 &\quad \underset{i_1=i_2=i_3=0}{=} &
 \end{aligned}$$

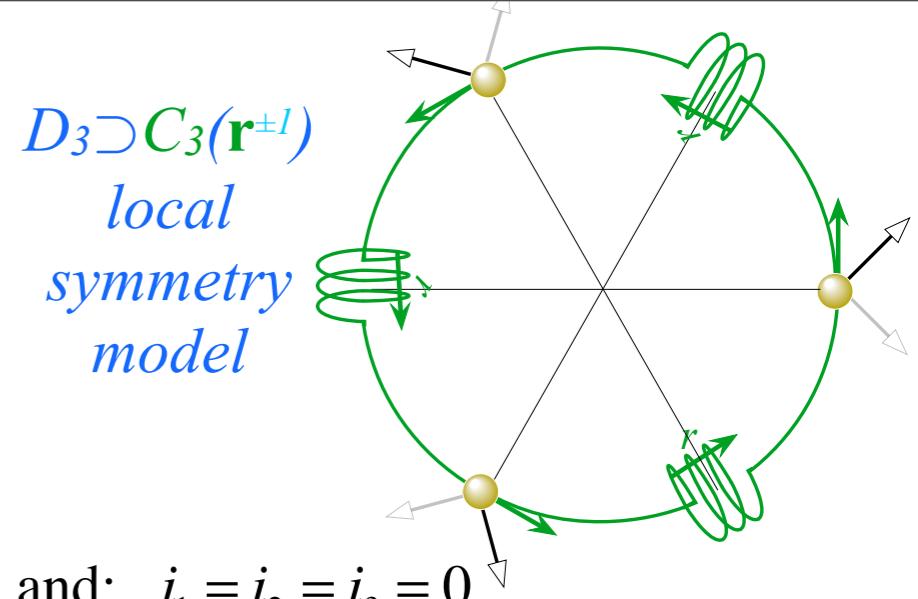
# $D_3 \supset C_3(\mathbf{r}^{\pm l})$ local symmetry vibrational K-matrix eigenvalues $K_m/M = \omega_m^2$

$$K_{xx}^{A_1} = r_0$$

$$K_{yy}^{A_2} = r_0$$

$$\begin{pmatrix} K_{xx}^{E_1} & K_{xy}^{E_1} \\ K_{yx}^{E_1} & K_{yy}^{E_1} \end{pmatrix} = \begin{pmatrix} r_0 & -i\mathbf{r} \frac{\sqrt{3}}{2} \\ +i\mathbf{r} \frac{\sqrt{3}}{2} & r_0 \end{pmatrix} \Rightarrow \begin{pmatrix} r_0 + \mathbf{r} \frac{\sqrt{3}}{2} & 0 \\ 0 & r_0 - \mathbf{r} \frac{\sqrt{3}}{2} \end{pmatrix}$$

# $D_3 \supset C_3(\mathbf{r}^{\pm l})$ local symmetry K-matrix eigensolutions



$D_3 \supset C_3(\mathbf{r}^{\pm l})$  local symmetry vibrational K-matrix Set:  $r_1 = \mathbf{r} = -r_2^*$ , and:  $i_1 = i_2 = i_3 = 0$

$$\begin{aligned}
 K_{xx}^{A_1} &= r_0 + \mathbf{r}_1 + \mathbf{r}_1^* + i_1 + i_2 + i_3 & = r_0 \\
 K_{yy}^{A_2} &= r_0 + \mathbf{r}_1 + \mathbf{r}_1^* - i_1 - i_2 - i_3 & = r_0 \\
 \begin{pmatrix} K_{xx}^{E_1} & K_{xy}^{E_1} \\ K_{yx}^{E_1} & K_{yy}^{E_1} \end{pmatrix} &= \frac{1}{2} \begin{pmatrix} 2r_0 - \mathbf{r}_1 - \mathbf{r}_1^* - i_1 - i_2 + 2i_3 & \sqrt{3}(-\mathbf{r}_1 + \mathbf{r}_1^* - i_1 + i_2) \\ \sqrt{3}(-\mathbf{r}_1^* + \mathbf{r}_1 - i_1 + i_2) & 2r_0 - \mathbf{r}_1 - \mathbf{r}_1^* + i_1 + i_2 - 2i_3 \end{pmatrix} &= \begin{pmatrix} r_0 & -i\mathbf{r} \frac{\sqrt{3}}{2} \\ +i\mathbf{r} \frac{\sqrt{3}}{2} & r_0 \end{pmatrix} \\
 &\quad \underset{i_1=i_2=i_3=0}{=} &
 \end{aligned}$$

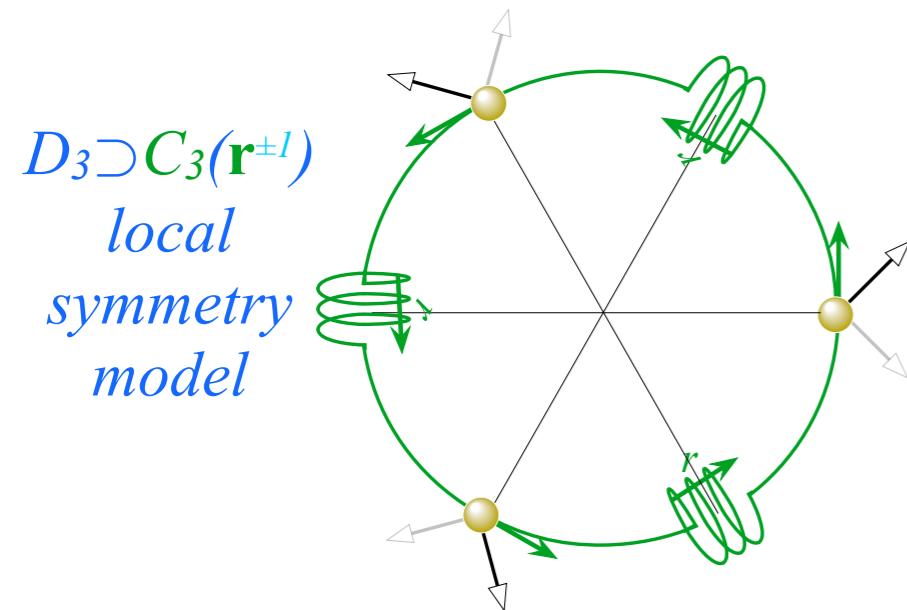
## $D_3 \supset C_3(\mathbf{r}^{\pm l})$ local symmetry vibrational K-matrix eigenvalues $K_m/M = \omega_m^2$

$E_1$  Eigenvectors in terms of  $D_3 \supset C_2(i_3)$   $E_1$ -vectors

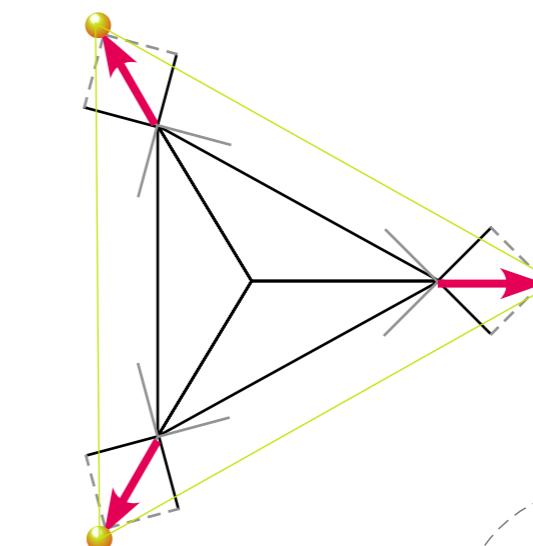
$$\begin{aligned}
 K_{xx}^{A_1} &= r_0 \\
 K_{yy}^{A_2} &= r_0 \\
 \begin{pmatrix} K_{xx}^{E_1} & K_{xy}^{E_1} \\ K_{yx}^{E_1} & K_{yy}^{E_1} \end{pmatrix} &= \begin{pmatrix} r_0 & -i\mathbf{r} \frac{\sqrt{3}}{2} \\ +i\mathbf{r} \frac{\sqrt{3}}{2} & r_0 \end{pmatrix} \Rightarrow \begin{pmatrix} r_0 + \mathbf{r} \frac{\sqrt{3}}{2} & 0 \\ 0 & r_0 - \mathbf{r} \frac{\sqrt{3}}{2} \end{pmatrix}
 \end{aligned}$$

$$\begin{aligned}
 \mathbf{K} \begin{pmatrix} E_1 \\ g(1)_3 \end{pmatrix} &= \mathbf{K} \left( \begin{pmatrix} E_1 \\ gx \end{pmatrix} + i \begin{pmatrix} E_1 \\ gy \end{pmatrix} \right) \frac{1}{\sqrt{2}} = +\mathbf{r} \frac{\sqrt{3}}{2} \begin{pmatrix} E_1 \\ g(1)_3 \end{pmatrix}, \\
 \mathbf{K} \begin{pmatrix} E_1 \\ g(2)_3 \end{pmatrix} &= \mathbf{K} \left( \begin{pmatrix} E_1 \\ gx \end{pmatrix} - i \begin{pmatrix} E_1 \\ gy \end{pmatrix} \right) \frac{1}{\sqrt{2}} = -\mathbf{r} \frac{\sqrt{3}}{2} \begin{pmatrix} E_1 \\ g(2)_3 \end{pmatrix}.
 \end{aligned}$$

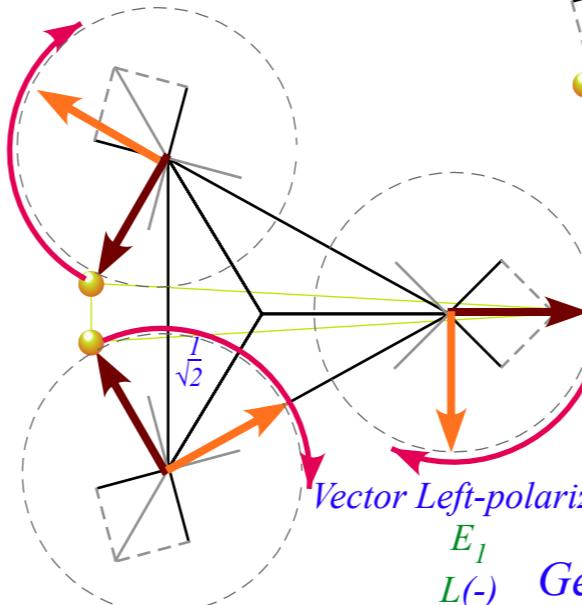
# $D_3 \supset C_3(\mathbf{r}^{\pm 1})$ local symmetry K-matrix eigensolutions



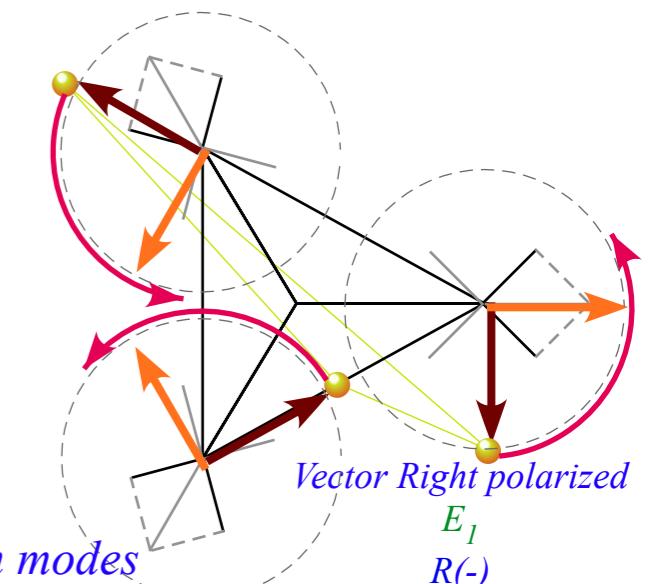
Strong  
 $C_3$  coupling  
limit



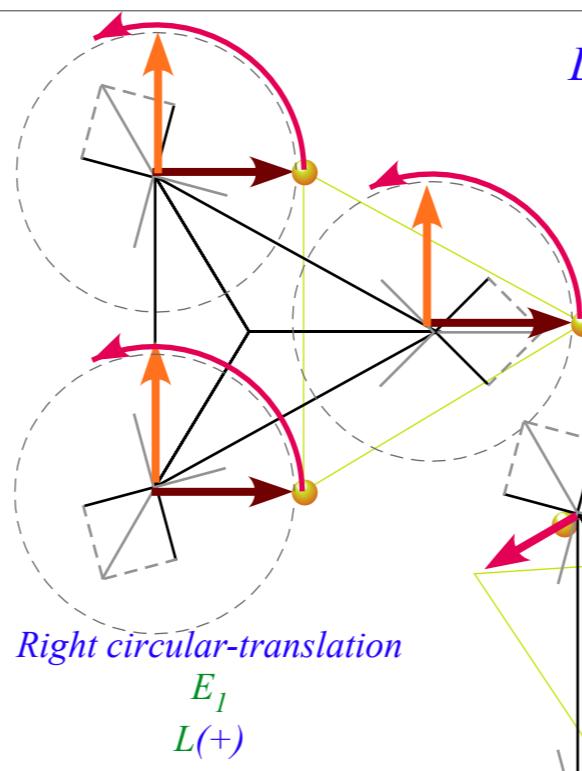
Scalar mode  
 $A_1$   
 $xx$



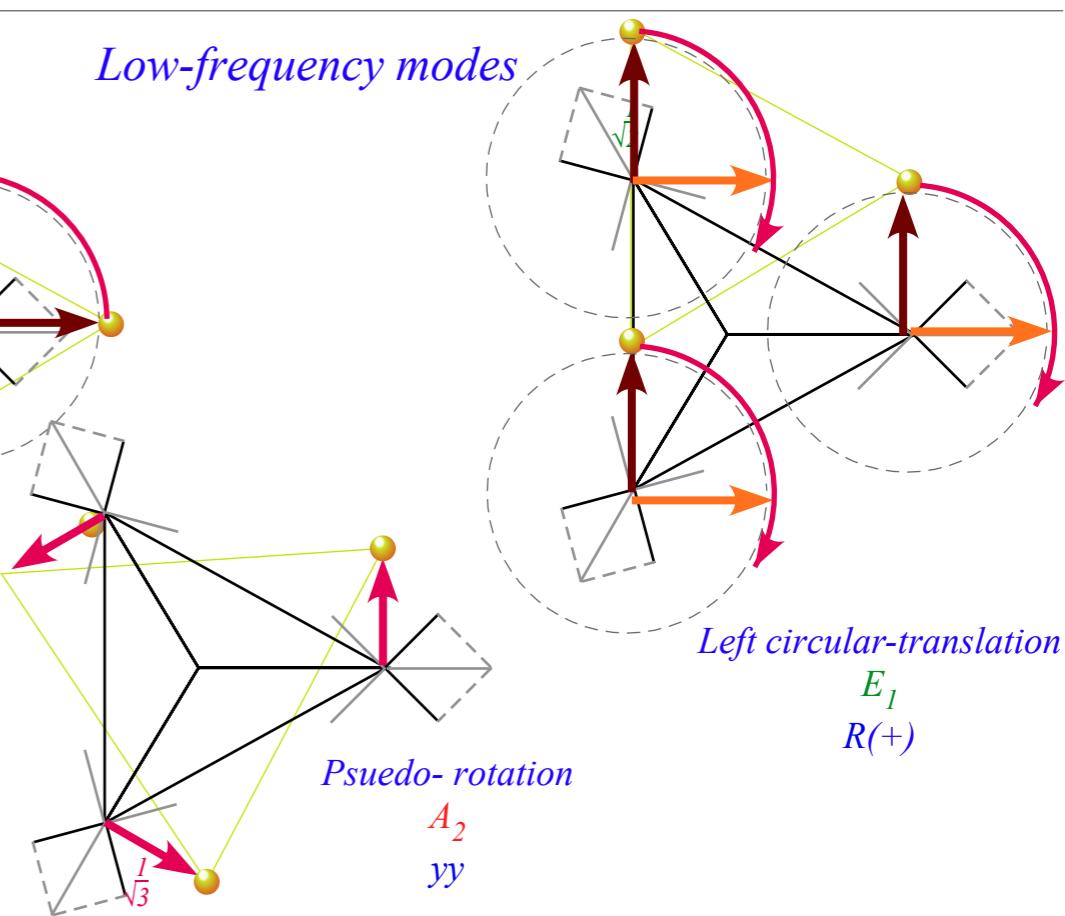
Vector Left-polarized  
 $E_1$   
 $L(-)$  Genuine vibration modes



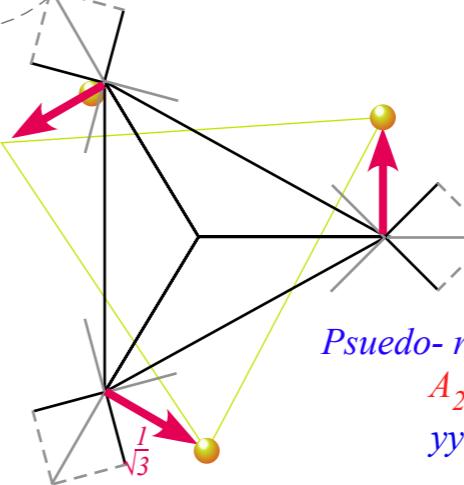
Vector Right polarized  
 $E_1$   
 $R(-)$



Right circular-translation  
 $E_1$   
 $L(+)$



Left circular-translation  
 $E_1$   
 $R(+)$



Pseudo- rotation  
 $A_2$   
 $yy$

*Review: Hamiltonian local-symmetry eigensolution in global and local  $|\mathbf{P}^{(\mu)}\rangle$ -basis*

*Molecular vibrational modes vs. Hamiltonian eigenmodes*

*Molecular K-matrix construction*

*$D_3 \supset C_2(i_3)$  local-symmetry K-matrix eigensolutions*

*$D_3$ -direct-connection K-matrix eigensolutions*

*$D_3 \supset C_3(\mathbf{r}^{\pm l})$  local symmetry K-matrix eigensolutions*

→ *Applied symmetry reduction and splitting* ←

*Subduced irep  $D^\alpha(D^3) \downarrow C_2 = d^{0_2} \oplus d^{l_2} \oplus ..$  correlation*

*Subduced irep  $D^\alpha(D^3) \downarrow C_3 = d^{0_3} \oplus d^{l_3} \oplus ..$  correlation*

*Spontaneous symmetry breaking and clustering: Frobenius Reciprocity , band structure*

*Induced rep  $d^a(C_2) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus ..$  correlation*

*Induced rep  $d^a(C_3) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus ..$  correlation*

*$D_6$  symmetry and Hexagonal Bands*

*Cross product of the  $C_2$  and  $D_3$  characters gives all  $D_6 = D_3 \times C_2$  characters and ireps*

*Applied symmetry reduction and splitting: Subduced irep  $D^\alpha(D^3) \downarrow C_2 = d^{0_2} \oplus d^{l_2} \oplus \dots$  correlation*

*Applied symmetry reduction and splitting: Subduced irep  $D^\alpha(D^3) \downarrow C_2 = d^{0_2} \oplus d^{1_2} \oplus \dots$  correlation*

$$\begin{array}{llll}
 D_3 \supset C_2 & \xrightarrow{\mathbf{P}^\alpha \text{ relabel/split}} & \xrightarrow{D^\alpha \text{ relabel/reduce}} & \xrightarrow{\omega^\alpha \text{ relabel/split}} \\
 A_1 & \mathbf{P}^{A_1} = \mathbf{P}^{A_1} \mathbf{P}^{0_2} = \mathbf{P}_{0_2 0_2}^{A_1} & \Rightarrow D^{A_1} \downarrow C_2 \sim d^{0_2} & \Rightarrow \omega^{A_1} \rightarrow \omega^{0_2} \\
 A_2 & \mathbf{P}^{A_2} = \mathbf{P}^{A_2} \mathbf{P}^{1_2} = \mathbf{P}_{1_2 1_2}^{A_2} & \Rightarrow D^{A_2} \downarrow C_2 \sim d^{1_2} & \Rightarrow \omega^{A_2} \rightarrow \omega^{1_2} \\
 E_1 & \mathbf{P}^{E_1} = \mathbf{P}^{E_1} \mathbf{P}^{0_2} + \mathbf{P}^{E_1} \mathbf{P}^{1_2} & \Rightarrow D^{E_1} \downarrow C_2 \sim & \Rightarrow \omega^{E_1} \rightarrow \omega^{0_2} \\
 & = \mathbf{P}_{0_2 0_2}^{E_1} + \mathbf{P}_{1_2 1_2}^{E_1} & d^{0_2} \oplus d^{1_2} & \searrow \omega^{1_2}
 \end{array}$$

Applied symmetry reduction and splitting: Subduced irep  $D^\alpha(D^3) \downarrow C_2 = d^{0_2} \oplus d^{1_2} \oplus \dots$  correlation

$D_3 \supset C_2$	$\mathbf{P}^\alpha$ relabel/split	$D^\alpha$ relabel/reduce	$\omega^\alpha$ relabel/split	$D_3 \supset C_2$	$0_2$	$1_2$
$A_1$	$\mathbf{P}^{A_1} = \mathbf{P}^{A_1} \mathbf{P}^{0_2} = \mathbf{P}_{0_2 0_2}^{A_1}$	$\Rightarrow D^{A_1} \downarrow C_2 \sim d^{0_2}$	$\Rightarrow \omega^{A_1} \rightarrow \omega^{0_2}$	$A_1$	1	.
$A_2$	$\mathbf{P}^{A_2} = \mathbf{P}^{A_2} \mathbf{P}^{1_2} = \mathbf{P}_{1_2 1_2}^{A_2}$	$\Rightarrow D^{A_2} \downarrow C_2 \sim d^{1_2}$	$\Rightarrow \omega^{A_2} \rightarrow \omega^{1_2}$	$A_2$	.	1
$E_1$	$\mathbf{P}^{E_1} = \mathbf{P}^{E_1} \mathbf{P}^{0_2} + \mathbf{P}^{E_1} \mathbf{P}^{1_2}$ $= \mathbf{P}_{0_2 0_2}^{E_1} + \mathbf{P}_{1_2 1_2}^{E_1}$	$\Rightarrow D^{E_1} \downarrow C_2 \sim$ $d^{0_2} \oplus d^{1_2}$	$\Rightarrow \omega^{E_1} \rightarrow \omega^{0_2}$ $\searrow \omega^{1_2}$	$E_1$	1	1

*Applied symmetry reduction and splitting: Subduced irep  $D^\alpha(D^3) \downarrow C_2 = d^{0_2} \oplus d^{l_2} \oplus \dots$  correlation*

$D_3 \supset C_2$	$\mathbf{P}^\alpha$ relabel/split	$D^\alpha$ relabel/reduce	$\omega^\alpha$ relabel/split	$D_3 \supset C_2$	$0_2$	$1_2$
$A_1$	$\mathbf{P}^{A_1} = \mathbf{P}^{A_1} \mathbf{P}^{0_2} = \mathbf{P}_{0_2 0_2}^{A_1}$	$\Rightarrow D^{A_1} \downarrow C_2 \sim d^{0_2}$	$\Rightarrow \omega^{A_1} \rightarrow \omega^{0_2}$	$A_1$	1	.
$A_2$	$\mathbf{P}^{A_2} = \mathbf{P}^{A_2} \mathbf{P}^{1_2} = \mathbf{P}_{1_2 1_2}^{A_2}$	$\Rightarrow D^{A_2} \downarrow C_2 \sim d^{1_2}$	$\Rightarrow \omega^{A_2} \rightarrow \omega^{1_2}$	$A_2$	.	1
$E_1$	$\mathbf{P}^{E_1} = \mathbf{P}^{E_1} \mathbf{P}^{0_2} + \mathbf{P}^{E_1} \mathbf{P}^{1_2}$ $= \mathbf{P}_{0_2 0_2}^{E_1} + \mathbf{P}_{1_2 1_2}^{E_1}$	$\Rightarrow D^{E_1} \downarrow C_2 \sim$ $d^{0_2} \oplus d^{1_2}$	$\Rightarrow \omega^{E_1} \rightarrow \omega^{0_2}$ $\searrow \omega^{1_2}$	$E_1$	1	1

*Applied symmetry reduction and splitting: Subduced irep  $D^\alpha(D^3) \downarrow C_3 = d^{0_3} \oplus d^{l_3} \oplus \dots$  correlation*

Applied symmetry reduction and splitting: Subduced irep  $D^\alpha(D^3) \downarrow C_2 = d^{0_2} \oplus d^{l_2} \oplus \dots$  correlation

$D_3 \supset C_2$	$\underline{P^\alpha \text{ relabel/split}}$	$\underline{D^\alpha \text{ relabel/reduce}}$	$\underline{\omega^\alpha \text{ relabel/split}}$	$D_3 \supset C_2$	$0_2$	$1_2$
$A_1$	$P^{A_1} = P^{A_1} P^{0_2} = P_{0_2 0_2}^{A_1}$	$\Rightarrow D^{A_1} \downarrow C_2 \sim d^{0_2}$	$\Rightarrow \omega^{A_1} \rightarrow \omega^{0_2}$	$A_1$	1	.
$A_2$	$P^{A_2} = P^{A_2} P^{1_2} = P_{1_2 1_2}^{A_2}$	$\Rightarrow D^{A_2} \downarrow C_2 \sim d^{1_2}$	$\Rightarrow \omega^{A_2} \rightarrow \omega^{1_2}$	$A_2$	.	1
$E_1$	$P^{E_1} = P^{E_1} P^{0_2} + P^{E_1} P^{1_2}$ $= P_{0_2 0_2}^{E_1} + P_{1_2 1_2}^{E_1}$	$\Rightarrow D^{E_1} \downarrow C_2 \sim d^{0_2} \oplus d^{1_2}$	$\Rightarrow \omega^{E_1} \rightarrow \omega^{0_2}$ $\searrow \omega^{1_2}$	$E_1$	1	1

Applied symmetry reduction and splitting: Subduced irep  $D^\alpha(D^3) \downarrow C_3 = d^{0_3} \oplus d^{l_3} \oplus \dots$  correlation

$D_3 \supset C_3$	$\underline{P^\alpha \text{ relabel/split}}$	$\underline{D^\alpha \text{ relabel/reduce}}$	$\underline{\omega^\alpha \text{ relabel/split}}$
$A_1$	$P^{A_1} = P^{A_1} P^{0_3} = P_{0_3 0_3}^{A_1}$	$\Rightarrow D^{A_1} \downarrow C_3 \sim d^{0_3}$	$\Rightarrow \omega^{A_1} \rightarrow \omega^{0_3}$
$A_2$	$P^{A_2} = P^{A_2} P^{0_3} = P_{0_3 0_3}^{A_2}$	$\Rightarrow D^{A_2} \downarrow C_3 \sim d^{0_3}$	$\Rightarrow \omega^{A_2} \rightarrow \omega^{0_3}$
$E_1$	$P^{E_1} = P^{E_1} P^{1_3} + P^{E_1} P^{2_3}$ $= P_{1_3 1_3}^{E_1} + P_{2_3 2_3}^{E_1}$	$\Rightarrow D^{E_1} \downarrow C_3 \sim d^{1_3} \oplus d^{2_3}$	$\Rightarrow \omega^{E_1} \rightarrow \omega^{1_3}$ $\searrow \omega^{2_3}$

Applied symmetry reduction and splitting: Subduced irep  $D^\alpha(D^3) \downarrow C_2 = d^{0_2} \oplus d^{l_2} \oplus \dots$  correlation

$D_3 \supset C_2$	$\mathbf{P}^\alpha$ relabel/split	$D^\alpha$ relabel/reduce	$\omega^\alpha$ relabel/split	$D_3 \supset C_2$	$0_2$	$1_2$
$A_1$	$\mathbf{P}^{A_1} = \mathbf{P}^{A_1} \mathbf{P}^{0_2} = \mathbf{P}_{0_2 0_2}^{A_1}$	$\Rightarrow D^{A_1} \downarrow C_2 \sim d^{0_2}$	$\Rightarrow \omega^{A_1} \rightarrow \omega^{0_2}$	$A_1$	1	.
$A_2$	$\mathbf{P}^{A_2} = \mathbf{P}^{A_2} \mathbf{P}^{l_2} = \mathbf{P}_{l_2 l_2}^{A_2}$	$\Rightarrow D^{A_2} \downarrow C_2 \sim d^{l_2}$	$\Rightarrow \omega^{A_2} \rightarrow \omega^{l_2}$	$A_2$	.	1
$E_1$	$\mathbf{P}^{E_1} = \mathbf{P}^{E_1} \mathbf{P}^{0_2} + \mathbf{P}^{E_1} \mathbf{P}^{l_2}$ $= \mathbf{P}_{0_2 0_2}^{E_1} + \mathbf{P}_{l_2 l_2}^{E_1}$	$\Rightarrow D^{E_1} \downarrow C_2 \sim d^{0_2} \oplus d^{l_2}$	$\Rightarrow \omega^{E_1} \rightarrow \omega^{0_2}$ $\searrow \omega^{l_2}$	$E_1$	1	1

Applied symmetry reduction and splitting: Subduced irep  $D^\alpha(D^3) \downarrow C_3 = d^{0_3} \oplus d^{l_3} \oplus \dots$  correlation

$D_3 \supset C_3$	$\mathbf{P}^\alpha$ relabel/split	$D^\alpha$ relabel/reduce	$\omega^\alpha$ relabel/split	$D_3 \supset C_3$	$0_3$	$1_3$	$2_3$
$A_1$	$\mathbf{P}^{A_1} = \mathbf{P}^{A_1} \mathbf{P}^{0_3} = \mathbf{P}_{0_3 0_3}^{A_1}$	$\Rightarrow D^{A_1} \downarrow C_3 \sim d^{0_3}$	$\Rightarrow \omega^{A_1} \rightarrow \omega^{0_3}$	$A_1$	1	.	.
$A_2$	$\mathbf{P}^{A_2} = \mathbf{P}^{A_2} \mathbf{P}^{0_3} = \mathbf{P}_{0_3 0_3}^{A_2}$	$\Rightarrow D^{A_2} \downarrow C_3 \sim d^{0_3}$	$\Rightarrow \omega^{A_2} \rightarrow \omega^{0_3}$	$A_2$	1	.	.
$E_1$	$\mathbf{P}^{E_1} = \mathbf{P}^{E_1} \mathbf{P}^{1_3} + \mathbf{P}^{E_1} \mathbf{P}^{2_3}$ $= \mathbf{P}_{1_3 1_3}^{E_1} + \mathbf{P}_{2_3 2_3}^{E_1}$	$\Rightarrow D^{E_1} \downarrow C_3 \sim d^{1_3} \oplus d^{2_3}$	$\Rightarrow \omega^{E_1} \rightarrow \omega^{1_3}$ $\searrow \omega^{2_3}$	$E_1$	.	1	1

Review: Hamiltonian local-symmetry eigensolution in global and local  $|\mathbf{P}^{(\mu)}\rangle$ -basis

Molecular vibrational modes vs. Hamiltonian eigenmodes

Molecular K-matrix construction

$D_3 \supset C_2(i_3)$  local-symmetry K-matrix eigensolutions

$D_3$ -direct-connection K-matrix eigensolutions

$D_3 \supset C_3(\mathbf{r}^{\pm l})$  local symmetry K-matrix eigensolutions

Applied symmetry reduction and splitting

Subduced irep  $D^\alpha(D^3) \downarrow C_2 = d^{02} \oplus d^{l_2} \oplus \dots$  correlation

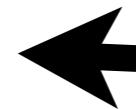
Subduced irep  $D^\alpha(D^3) \downarrow C_3 = d^{03} \oplus d^{l_3} \oplus \dots$  correlation



Spontaneous symmetry breaking and clustering: Frobenius Reciprocity , band structure

Induced rep  $d^a(C_2) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus \dots$  correlation

Induced rep  $d^a(C_3) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus \dots$  correlation



$D_6$  symmetry and Hexagonal Bands

Cross product of the  $C_2$  and  $D_3$  characters gives all  $D_6 = D_3 \times C_2$  characters and irreps

Applied symmetry reduction and splitting: Subduced irep  $D^\alpha(D^3) \downarrow C_2 = d^{0_2} \oplus d^{1_2} \oplus \dots$  correlation

$D_3 \supset C_2$	$\frac{\mathbf{P}^\alpha \text{ relabel/split}}{\mathbf{P}^{A_1} = \mathbf{P}^{A_1} \mathbf{P}^{0_2} = \mathbf{P}_{0_2 0_2}^{A_1}}$	$\frac{D^\alpha \text{ relabel/reduce}}{\Rightarrow D^{A_1} \downarrow C_2 \sim d^{0_2}}$	$\frac{\omega^\alpha \text{ relabel/split}}{\Rightarrow \omega^{A_1} \rightarrow \omega^{0_2}}$	$\frac{D_3 \supset C_2}{\begin{array}{c cc} & \mathbf{0}_2 & \mathbf{1}_2 \\ \hline A_1 & 1 & \cdot \\ A_2 & \cdot & 1 \\ E_1 & 1 & 1 \end{array}}$	$D^{A_1}(D_3) \downarrow C_2 \sim d^{0_2}$
$A_1$					$D^{A_2}(D_3) \downarrow C_2 \sim d^{1_2}$
$A_2$	$\mathbf{P}^{A_2} = \mathbf{P}^{A_2} \mathbf{P}^{1_2} = \mathbf{P}_{1_2 1_2}^{A_2}$	$\Rightarrow D^{A_2} \downarrow C_2 \sim d^{1_2}$	$\Rightarrow \omega^{A_2} \rightarrow \omega^{1_2}$		
$E_1$	$\mathbf{P}^{E_1} = \mathbf{P}^{E_1} \mathbf{P}^{0_2} + \mathbf{P}^{E_1} \mathbf{P}^{1_2}$ $= \mathbf{P}_{0_2 0_2}^{E_1} + \mathbf{P}_{1_2 1_2}^{E_1}$	$\Rightarrow D^{E_1} \downarrow C_2 \sim d^{0_2} \oplus d^{1_2}$	$\Rightarrow \omega^{E_1} \rightarrow \omega^{0_2}$ $\searrow \omega^{1_2}$		$D^{E_1}(D_3) \downarrow C_2 \sim d^{0_2} \oplus d^{1_2}$
				$d^{0_2}(C_2) \uparrow D_3$	
				$\sim D^{A_1} \oplus D^{E_1}$	
				$d^{1_2}(C_2) \uparrow D_3$	
				$\sim D^{A_2} \oplus D^{E_1}$	

Spontaneous symmetry breaking

and clustering: Induced rep  $d^a(C_2) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus \dots$  correlation

Applied symmetry reduction and splitting: Subduced irep  $D^\alpha(D^3) \downarrow C_3 = d^{0_3} \oplus d^{1_3} \oplus \dots$  correlation

$D_3 \supset C_3$	$\frac{\mathbf{P}^\alpha \text{ relabel/split}}{\mathbf{P}^{A_1} = \mathbf{P}^{A_1} \mathbf{P}^{0_3} = \mathbf{P}_{0_3 0_3}^{A_1}}$	$\frac{D^\alpha \text{ relabel/reduce}}{\Rightarrow D^{A_1} \downarrow C_3 \sim d^{0_3}}$	$\frac{\omega^\alpha \text{ relabel/split}}{\Rightarrow \omega^{A_1} \rightarrow \omega^{0_3}}$	$\frac{D_3 \supset C_3}{\begin{array}{c ccc} & \mathbf{0}_3 & \mathbf{1}_3 & \mathbf{2}_3 \\ \hline A_1 & 1 & \cdot & \cdot \\ A_2 & 1 & \cdot & \cdot \\ E_1 & \cdot & 1 & 1 \end{array}}$	$D^{A_1}(D_3) \downarrow C_3 \sim d^{0_3}$
$A_1$					$D^{A_2}(D_3) \downarrow C_3 \sim d^{0_3}$
$A_2$	$\mathbf{P}^{A_2} = \mathbf{P}^{A_2} \mathbf{P}^{0_3} = \mathbf{P}_{0_3 0_3}^{A_2}$	$\Rightarrow D^{A_2} \downarrow C_3 \sim d^{0_3}$	$\Rightarrow \omega^{A_2} \rightarrow \omega^{0_3}$		
$E_1$	$\mathbf{P}^{E_1} = \mathbf{P}^{E_1} \mathbf{P}^{1_3} + \mathbf{P}^{E_1} \mathbf{P}^{2_3}$ $= \mathbf{P}_{1_3 1_3}^{E_1} + \mathbf{P}_{2_3 2_3}^{E_1}$	$\Rightarrow D^{E_1} \downarrow C_3 \sim d^{1_3} \oplus d^{2_3}$	$\Rightarrow \omega^{E_1} \rightarrow \omega^{1_3}$ $\searrow \omega^{2_3}$		$D^{E_1}(D_3) \downarrow C_3 \sim d^{1_3} \oplus d^{2_3}$
				$d^{0_3}(C_3) \uparrow D_3$	
				$\sim D^{A_1} \oplus D^{A_2}$	
				$d^{1_3}(C_3) \uparrow D_3$	
				$\sim D^{E_1}$	
				$d^{2_3}(C_3) \uparrow D_3$	
				$\sim D^{E_1}$	

Spontaneous symmetry breaking

and clustering: Induced rep  $d^a(C_3) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus \dots$  correlation

### *Frobenius Reciprocity Theorem*

$$\text{Number of } D^\alpha \text{ in } d^k(K) \uparrow G = \text{Number of } d^k \text{ in } D^\alpha(G) \downarrow K$$

### *Frobenius Reciprocity Theorem*

$$\text{Number of } D^\alpha \text{ in } d^k(K) \uparrow G = \text{Number of } d^k \text{ in } D^\alpha(G) \downarrow K$$

*..and regular representation*

$D_3 \supset C_1$	$0_1 = 1_1$
$A_1$	1
$A_2$	1
$E_1$	2

## Frobenius Reciprocity Theorem

$$\text{Number of } D^\alpha \text{ in } d^k(K) \uparrow G = \text{Number of } d^k \text{ in } D^\alpha(G) \downarrow K$$

..and regular representation

$D_3 \supset C_1$	$0_1 = 1_1$
$A_1$	1
$A_2$	1
$E_1$	2

$D_3 \supset C_2$	$0_2$	$1_2$
$A_1$	1	.
$A_2$	.	1
$E_1$	1	1

$D_3 \supset C_3$	$0_3$	$1_3$	$2_3$
$A_1$	1	.	.
$A_2$	1	.	.
$E_1$	.	1	1

*Review: Hamiltonian local-symmetry eigensolution in global and local  $|\mathbf{P}^{(\mu)}\rangle$ -basis*

*Molecular vibrational modes vs. Hamiltonian eigenmodes*

*Molecular K-matrix construction*

*$D_3 \supset C_2(i_3)$  local-symmetry K-matrix eigensolutions*

*$D_3$ -direct-connection K-matrix eigensolutions*

*$D_3 \supset C_3(\mathbf{r}^{\pm l})$  local symmetry K-matrix eigensolutions*

*Applied symmetry reduction and splitting*

*Subduced irep  $D^\alpha(D^3) \downarrow C_2 = d^{0_2} \oplus d^{l_2} \oplus \dots$  correlation*

*Subduced irep  $D^\alpha(D^3) \downarrow C_3 = d^{0_3} \oplus d^{l_3} \oplus \dots$  correlation*

*Spontaneous symmetry breaking and clustering: Frobenius Reciprocity , band structure*

*Induced rep  $d^a(C_2) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus \dots$  correlation*

*Induced rep  $d^a(C_3) \uparrow D^3 = D^\alpha \oplus D^\beta \oplus \dots$  correlation*

→  *$D_6$  symmetry and Hexagonal Bands*

*Cross product of the  $C_2$  and  $D_3$  characters gives all  $D_6 = D_3 \times C_2$  characters and ireps*

## *D<sub>6</sub>* symmetry and Hexagonal Bands

*D<sub>6</sub>* is the *outer product* ( $\times$ ) product  $D_3 \times C_2$  of  $D_3$  and  $C_2$ . (Requires  $C_2$  to commute with all of  $D_3$ .)

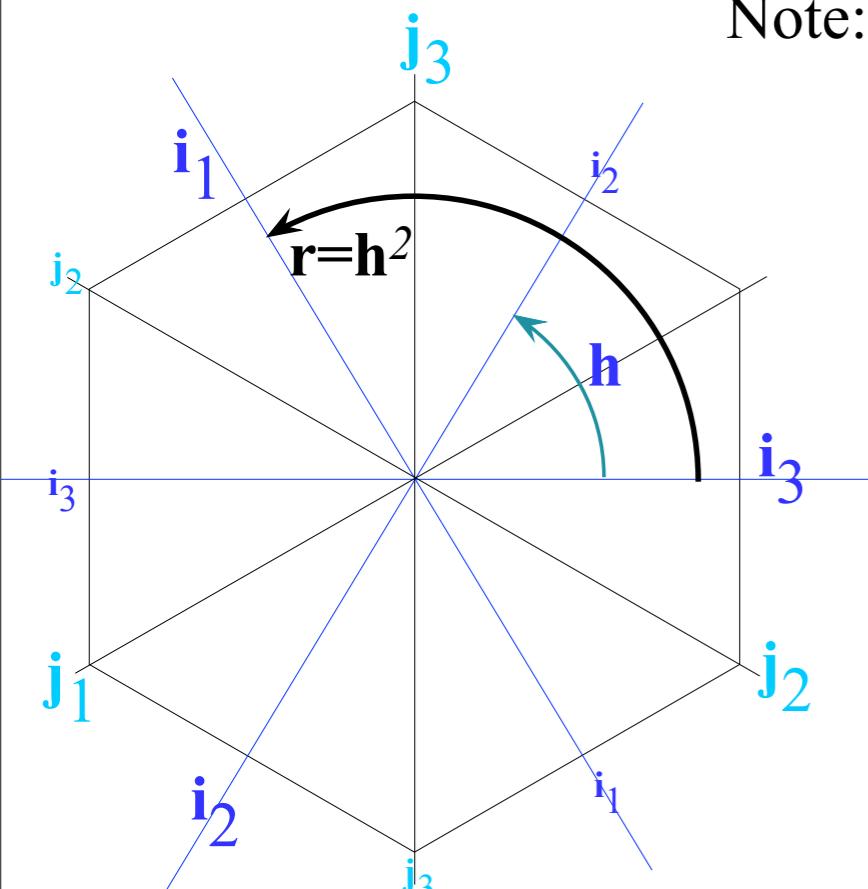
$$D_6 = D_3 \times C_2 = \{1, \mathbf{r}, \mathbf{r}^2, \mathbf{i}_1, \mathbf{i}_2, \mathbf{i}_3\} \times \{1, \mathbf{R}_z\}$$

$\times$  product and  $D_6$  operators. Define *hexagonal generator*  $\mathbf{h}$  of subgroup  $C_6 = \{1, \mathbf{h}, \mathbf{h}^2, \mathbf{h}^3, \mathbf{h}^4, \mathbf{h}^5\}$

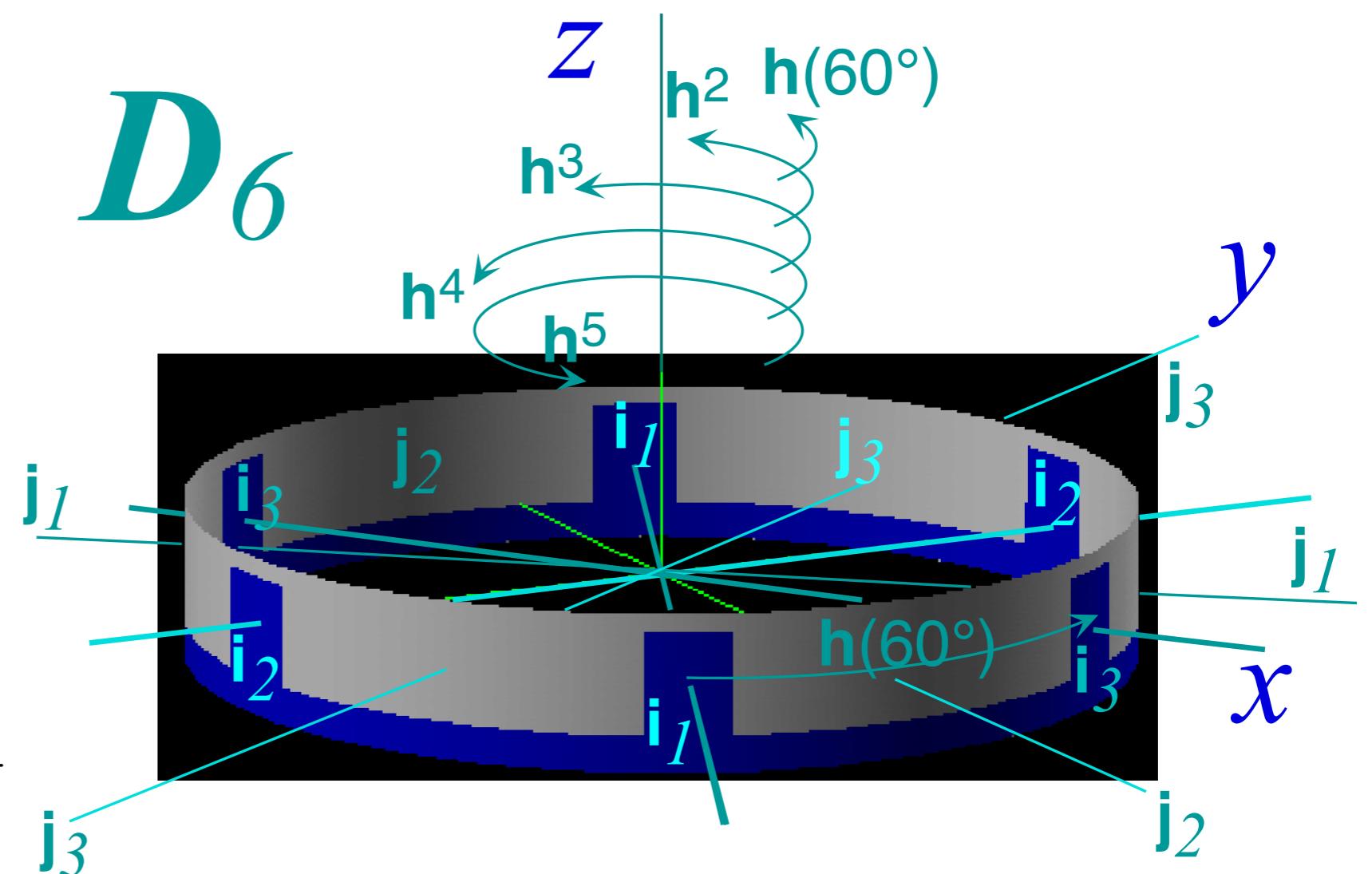
$$D_6 = D_3 \times C_2 = \{1, \mathbf{r}, \mathbf{r}^2, \mathbf{i}_1, \mathbf{i}_2, \mathbf{i}_3, 1 \cdot \mathbf{R}_z, \mathbf{r} \cdot \mathbf{R}_z, \mathbf{r}^2 \cdot \mathbf{R}_z, \mathbf{i}_1 \cdot \mathbf{R}_z, \mathbf{i}_2 \cdot \mathbf{R}_z, \mathbf{i}_3 \cdot \mathbf{R}_z\}$$

$$D_6 = D_3 \times C_2 = \{1, \mathbf{h}^2, \mathbf{h}^4, \mathbf{i}_1, \mathbf{i}_2, \mathbf{i}_3, \mathbf{h}^3, \mathbf{h}^5, \mathbf{h}, \mathbf{j}_1, \mathbf{j}_2, \mathbf{j}_3\}$$

Note:  $\mathbf{h}^2 = \mathbf{r}$  and  $\mathbf{h}^3 = \mathbf{R}_z$  and  $\mathbf{h}^4 = \mathbf{r}^2$  and  $\mathbf{h}^5 = \mathbf{r} \cdot \mathbf{R}_z$



NOTE:  
The  $\mathbf{i}_a$  and  $\mathbf{j}_b$  do not flip over the potential plot.



Electrostatic potential  $V(\phi)$  doesn't care which way is "up." Wells remain wells, and barriers remain barriers under all  $D_6$  operations.

Cross product of the  $C_2$  and  $D_3$  characters gives all  $D_6 = D_3 \times C_2$  characters.

$D_3$	$\begin{array}{ c c c } \hline 1 & \{\mathbf{r}, \mathbf{r}^2\} & \{\mathbf{i}_1, \mathbf{i}_2, \mathbf{i}_3\} \\ \hline \end{array}$			$\begin{array}{ c c } \hline C_2^z & 1 & \mathbf{R}_z \\ \hline (A) & 1 & 1 \\ (B) & 1 & -1 \\ \hline \end{array}$			$D_3 \times C_2^z$		
$\chi^{A_1}(\mathbf{g})$	1	1	1	$A_1 \cdot (A)$	1·1	1·1	1·1	1·1	1·1
$\chi^{A_2}(\mathbf{g})$	1	1	-1	$A_2 \cdot (A)$	1·1	1·1	-1·1	1·1	1·1
$\chi^{E_1}(\mathbf{g})$	2	-1	0	$E_1 \cdot (A)$	2·1	-1·1	0·1	2·1	-1·1
				$A_1 \cdot (B)$	1·1	1·1	1·1	1·(-1)	1·(-1)
				$A_2 \cdot (B)$	1·1	1·1	-1·1	1·(-1)	1·(-1)
				$E_1 \cdot (B)$	2·1	-1·1	0·1	2·(-1)	-1·(-1)
									0·(-1)
$\chi_g^\mu(D_6)$	$\begin{array}{ c c c } \hline 1 & \{\mathbf{h}^2, \mathbf{h}^4\} & \{\mathbf{i}_1, \mathbf{i}_2, \mathbf{i}_3\} \\ \hline \end{array}$			$\begin{array}{ c c c } \hline \mathbf{h}^3 & \{\mathbf{h}, \mathbf{h}^5\} & \{\mathbf{j}_1, \mathbf{j}_2, \mathbf{j}_3\} \\ \hline \end{array}$			$D_3 \times C_2^z$		
	$A_1$	1	1	1	1	1	1		
	$A_2$	1	1	-1	1	1		-1	
	$E_2$	2	-1	0	2	-1		0	
	$B_2$	1	1	1	-1	-1		-1	
	$B_1$	1	1	-1	-1	-1		1	
	$E_1$	2	-1	0	-2	1		0	

(Recall  $C_2 \times C_2 = D_2$  characters made of two  $C_2$  groups)

$$\chi_g^\mu(D_6) =$$

Unit translation  
or

60° hex rotation  $\mathbf{h}$   
determines  
 $A_p$  vs  $B_p$   
(+1) vs (-1)

Y-rotation  
or  
180° flip  $\mathbf{j}_3$   
determines  
 $X_1$  vs  $X_2$   
(+1) vs (-1)

Cross product of the  $C_2$  and  $D_3$  ireps gives all  $D_6 = D_3 \times C_2$  ireps.

$\mathbf{g} =$	$1$	$\mathbf{r} = \mathbf{h}^2$	$\mathbf{r}^2 = \mathbf{h}^4$	$\mathbf{i}_1$	$\mathbf{i}_2$	$\mathbf{i}_3$	$\mathbf{h}^3$	$\mathbf{h}^3\mathbf{r} = \mathbf{h}^5$	$\mathbf{h}^3\mathbf{r}^2 = \mathbf{h}^1$	$\mathbf{h}^3\mathbf{i}_1 = \mathbf{j}_1$	$\mathbf{h}^3\mathbf{i}_2 = \mathbf{j}_2$	$\mathbf{h}^3\mathbf{i}_3 = \mathbf{j}_3$
$D^{A_1}(g) =$	1	1	1	1	1	1	1	1	1	1	1	1
$D^{A_2}(g) =$	1	1	1	-1	-1	-1	1	1	1	-1	-1	-1
$D^{E_2}(g) =$	$\begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix}$	$\begin{pmatrix} -1 & -\sqrt{3} \\ \frac{\sqrt{3}}{2} & \frac{-1}{2} \end{pmatrix}$	$\begin{pmatrix} -1 & \sqrt{3} \\ \frac{\sqrt{3}}{2} & \frac{1}{2} \end{pmatrix}$	$\begin{pmatrix} -1 & -\sqrt{3} \\ \frac{\sqrt{3}}{2} & \frac{1}{2} \end{pmatrix}$	$\begin{pmatrix} -1 & \sqrt{3} \\ \frac{\sqrt{3}}{2} & \frac{1}{2} \end{pmatrix}$	$\begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$	$\begin{pmatrix} -1 & -\sqrt{3} \\ \frac{\sqrt{3}}{2} & \frac{-1}{2} \end{pmatrix}$	$\begin{pmatrix} -1 & \sqrt{3} \\ \frac{\sqrt{3}}{2} & \frac{-1}{2} \end{pmatrix}$	$\begin{pmatrix} -1 & -\sqrt{3} \\ \frac{\sqrt{3}}{2} & \frac{1}{2} \end{pmatrix}$	$\begin{pmatrix} -1 & \sqrt{3} \\ \frac{\sqrt{3}}{2} & \frac{1}{2} \end{pmatrix}$	$\begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$	
$D^{B_2}(g) =$	1	1	1	1	1	1	-1	-1	-1	-1	-1	-1
$D^{B_1}(g) =$	1	1	1	-1	-1	-1	-1	-1	-1	1	1	1
$D^{E_1}(g) =$	$\begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix}$	$\begin{pmatrix} -1 & -\sqrt{3} \\ \frac{\sqrt{3}}{2} & \frac{-1}{2} \end{pmatrix}$	$\begin{pmatrix} -1 & \sqrt{3} \\ \frac{\sqrt{3}}{2} & \frac{1}{2} \end{pmatrix}$	$\begin{pmatrix} -1 & -\sqrt{3} \\ \frac{\sqrt{3}}{2} & \frac{1}{2} \end{pmatrix}$	$\begin{pmatrix} -1 & \sqrt{3} \\ \frac{\sqrt{3}}{2} & \frac{1}{2} \end{pmatrix}$	$\begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$	$\begin{pmatrix} 1 & \sqrt{3} \\ \frac{-\sqrt{3}}{2} & \frac{1}{2} \end{pmatrix}$	$\begin{pmatrix} 1 & -\sqrt{3} \\ \frac{\sqrt{3}}{2} & \frac{1}{2} \end{pmatrix}$	$\begin{pmatrix} 1 & \sqrt{3} \\ \frac{\sqrt{3}}{2} & \frac{-1}{2} \end{pmatrix}$	$\begin{pmatrix} 1 & -\sqrt{3} \\ \frac{\sqrt{3}}{2} & \frac{-1}{2} \end{pmatrix}$	$\begin{pmatrix} -1 & 0 \\ 0 & 1 \end{pmatrix}$	

Unit translation

or

$60^\circ$  hex rotation  $\mathbf{h}$

determines

$A_p$  vs  $B_p$

(+1) vs (-1)

$Y$ -rotation

or

$180^\circ$  flip  $\mathbf{j}_3$

determines

$X_1$  vs  $X_2$

(+1) vs (-1)

Cross product of the  $C_2$  and  $D_3$  ireps gives all  $D_6 = D_3 \times C_2$  ireps.

$\mathbf{g} =$	$1$	$\mathbf{r} = \mathbf{h}^2$	$\mathbf{r}^2 = \mathbf{h}^4$	$\mathbf{i}_1$	$\mathbf{i}_2$	$\mathbf{i}_3$	$\mathbf{h}^3$	$\mathbf{h}^3\mathbf{r} = \mathbf{h}^5$	$\mathbf{h}^3\mathbf{r}^2 = \mathbf{h}^1$	$\mathbf{h}^3\mathbf{i}_1 = \mathbf{j}_1$	$\mathbf{h}^3\mathbf{i}_2 = \mathbf{j}_2$	$\mathbf{h}^3\mathbf{i}_3 = \mathbf{j}_3$
$D^{A_1}(g) =$	1	1	1	1	1	1	1	1	1	1	1	1
$D^{A_2}(g) =$	1	1	1	-1	-1	-1	1	1	1	-1	-1	-1
$D^{E_2}(g) =$	$\begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix}$	$\begin{pmatrix} -1 & -\sqrt{3} \\ \frac{1}{2} & \frac{\sqrt{3}}{2} \end{pmatrix}$	$\begin{pmatrix} -1 & \sqrt{3} \\ \frac{1}{2} & \frac{-\sqrt{3}}{2} \end{pmatrix}$	$\begin{pmatrix} -1 & -\sqrt{3} \\ \frac{1}{2} & \frac{\sqrt{3}}{2} \end{pmatrix}$	$\begin{pmatrix} -1 & \sqrt{3} \\ \frac{1}{2} & \frac{-\sqrt{3}}{2} \end{pmatrix}$	$\begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$	$\begin{pmatrix} -1 & -\sqrt{3} \\ \frac{1}{2} & \frac{\sqrt{3}}{2} \end{pmatrix}$	$\begin{pmatrix} -1 & \sqrt{3} \\ \frac{1}{2} & \frac{-\sqrt{3}}{2} \end{pmatrix}$	$\begin{pmatrix} -1 & -\sqrt{3} \\ \frac{1}{2} & \frac{\sqrt{3}}{2} \end{pmatrix}$	$\begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$	$\begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$	
$D^{B_2}(g) =$	1	1	1	1	1	1	-1	-1	-1	-1	-1	-1
$D^{B_1}(g) =$	1	1	1	-1	-1	-1	-1	-1	-1	1	1	1
$D^{E_1}(g) =$	$\begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix}$	$\begin{pmatrix} -1 & -\sqrt{3} \\ \frac{1}{2} & \frac{\sqrt{3}}{2} \end{pmatrix}$	$\begin{pmatrix} -1 & \sqrt{3} \\ \frac{1}{2} & \frac{-\sqrt{3}}{2} \end{pmatrix}$	$\begin{pmatrix} -1 & \sqrt{3} \\ \frac{1}{2} & \frac{-\sqrt{3}}{2} \end{pmatrix}$	$\begin{pmatrix} -1 & \sqrt{3} \\ \frac{1}{2} & \frac{-\sqrt{3}}{2} \end{pmatrix}$	$\begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$	$\begin{pmatrix} 1 & \sqrt{3} \\ \frac{1}{2} & \frac{-\sqrt{3}}{2} \end{pmatrix}$	$\begin{pmatrix} 1 & \sqrt{3} \\ \frac{1}{2} & \frac{-\sqrt{3}}{2} \end{pmatrix}$	$\begin{pmatrix} 1 & \sqrt{3} \\ \frac{1}{2} & \frac{-\sqrt{3}}{2} \end{pmatrix}$	$\begin{pmatrix} 1 & -\sqrt{3} \\ \frac{1}{2} & \frac{\sqrt{3}}{2} \end{pmatrix}$	$\begin{pmatrix} 1 & -\sqrt{3} \\ \frac{1}{2} & \frac{\sqrt{3}}{2} \end{pmatrix}$	$\begin{pmatrix} -1 & 0 \\ 0 & 1 \end{pmatrix}$

Unit translation

or

60° hex rotation  $\mathbf{h}$

determines

$A_p$  vs  $B_p$

(+1) vs (-1)

Y-rotation

or

180° flip  $\mathbf{j}_3$

determines

$X_1$  vs  $X_2$

(+1) vs (-1)

$D_3 \supset C_2(j_3)$	$0_2$	$1_2$
$A_1$	1	.
$A_2$	.	1
$E_2$	1	1
$B_2$	.	1
$B_1$	1	.
$E_1$	1	1

$D_6 \supset C_3(h)$	$0_6$	$1_6$	$2_6$	$3_6$	$4_6$	$5_6$
$A_1$	1	.	.	.	.	.
$A_2$	1	.	.	.	.	.
$E_2$	.	.	1	.	1	.
$B_2$	.	.	.	1	.	.
$B_1$	.	.	.	1	.	.
$E_1$	.	1	.	.	.	1

*D<sub>6</sub> Band structure  
and related  
induced  
representations*

