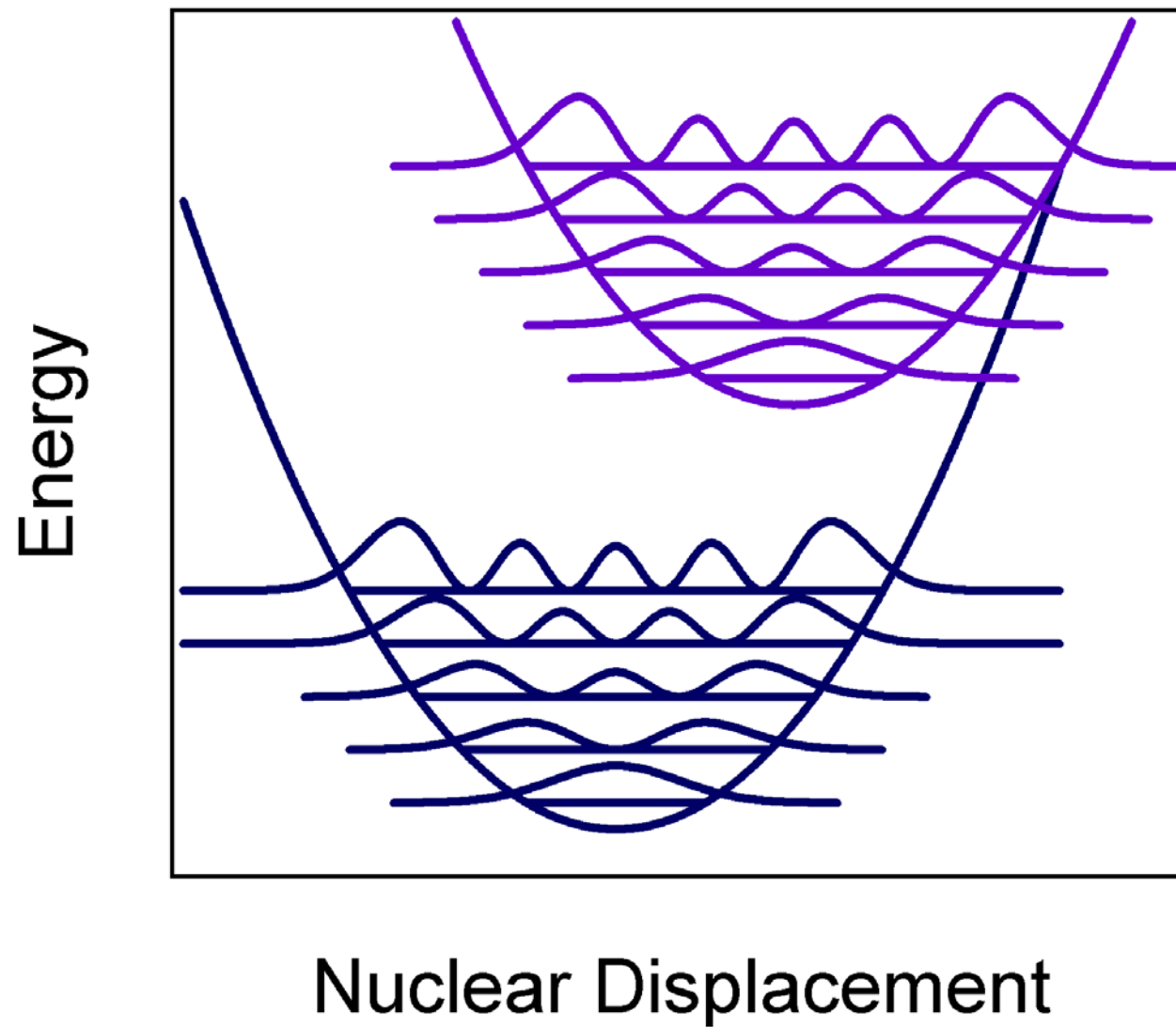


Franck-Condon factor



Franck-Condon factor

We can use a $T = 0$ K approximation to calculate the Franck-Condon factor. In this approximation the molecules are in their electronic and vibrational ground state. Therefore, the FC transitions are from $0 \rightarrow 0'$, $0 \rightarrow 1'$, $0 \rightarrow 2'$, etc. The formula for this approximation is given by:

$$FC = \sum_{n=0'}^{\infty} \frac{S^n e^{-S}}{n!} \delta(\omega - \omega_{0-0'} - n\omega_{vib})$$

Calculate the “stick spectrum” by making a table showing the relative magnitude of the first 5 vibrational lines above $0-0'$. You are given the following information.

$$\begin{aligned}\omega_{0-0'} &= 21,000 \text{ cm}^{-1} \\ \omega_{vib} &= 1600 \text{ cm}^{-1} \\ S &= 1.5\end{aligned}$$