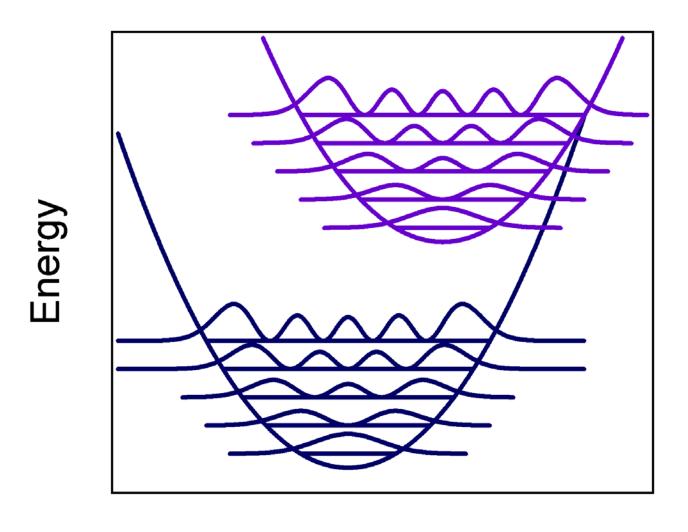
Franck-Condon factor



Nuclear Displacement

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 10 vibrational lines above 0-0'. You are given the follow information.

$$\omega_{0-0'} = 17,000 \ cm^{-1}$$
 $\omega_{vib} = 300 \ cm^{-1}$
 $S = 3.2$