

Infrared Spectra of Some Common Functional Groups

- **n-Alkanes**

- look for stretching and bending of C–H and C–C bonds

- **C–C bends**: ca. 500 cm^{-1} (out of spectral window)
- **C–C stretches**: $1200\text{--}800\text{ cm}^{-1}$, weak bands not of value for interpretation (fingerprint)

More characteristic

- **C–H stretches**: occurs from $3000\text{--}2840\text{ cm}^{-1}$
 - CH₃**: 2962 cm^{-1} , asymmetrical stretch
 2872 cm^{-1} , symmetrical stretch
 - CH₂**: 2926 cm^{-1} , asymmetrical stretch
 2853 cm^{-1} , symmetrical stretch
- **C–H bends**:
 - CH₃**: ca. 1375 cm^{-1}
 - CH₂**: ca. 1465 cm^{-1}