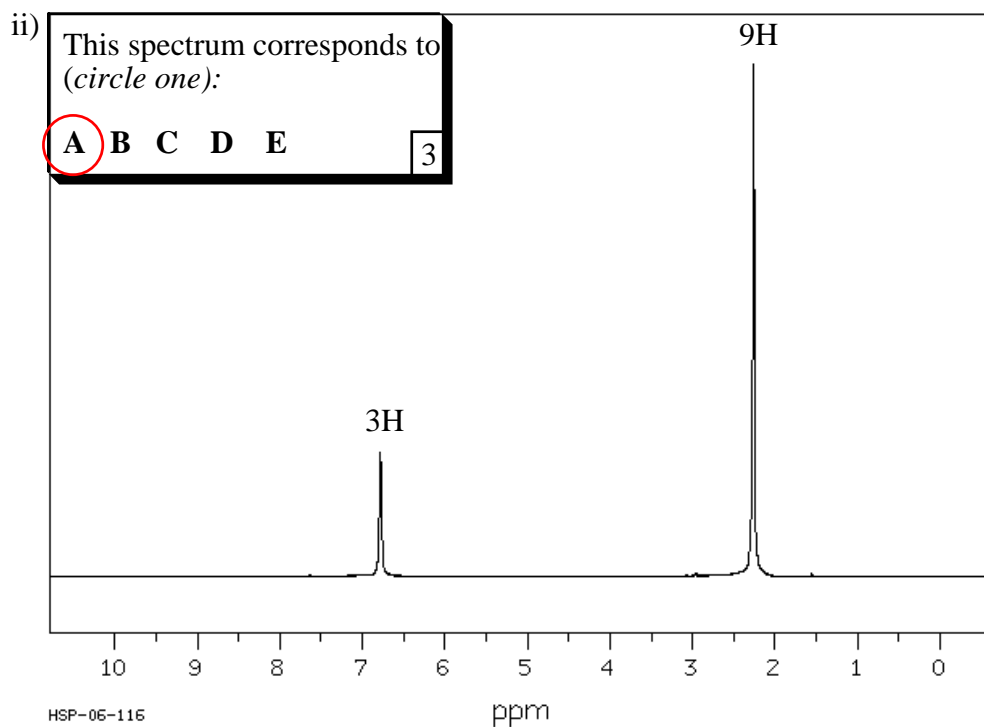
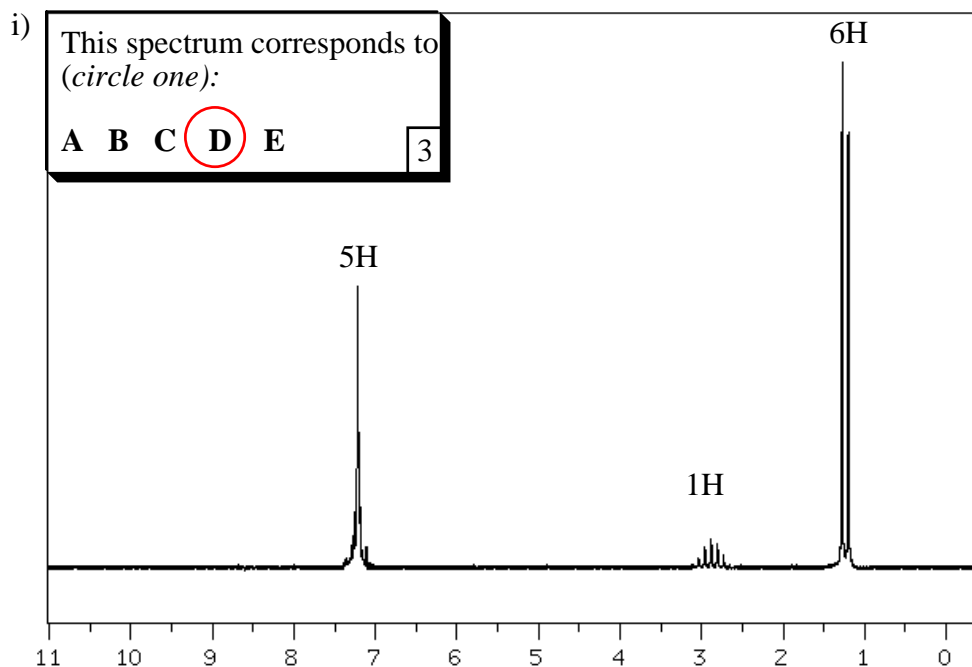
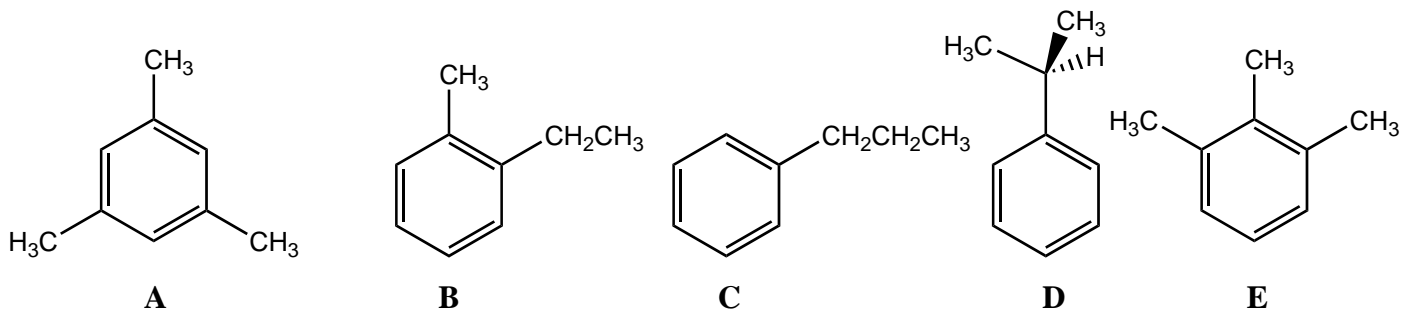
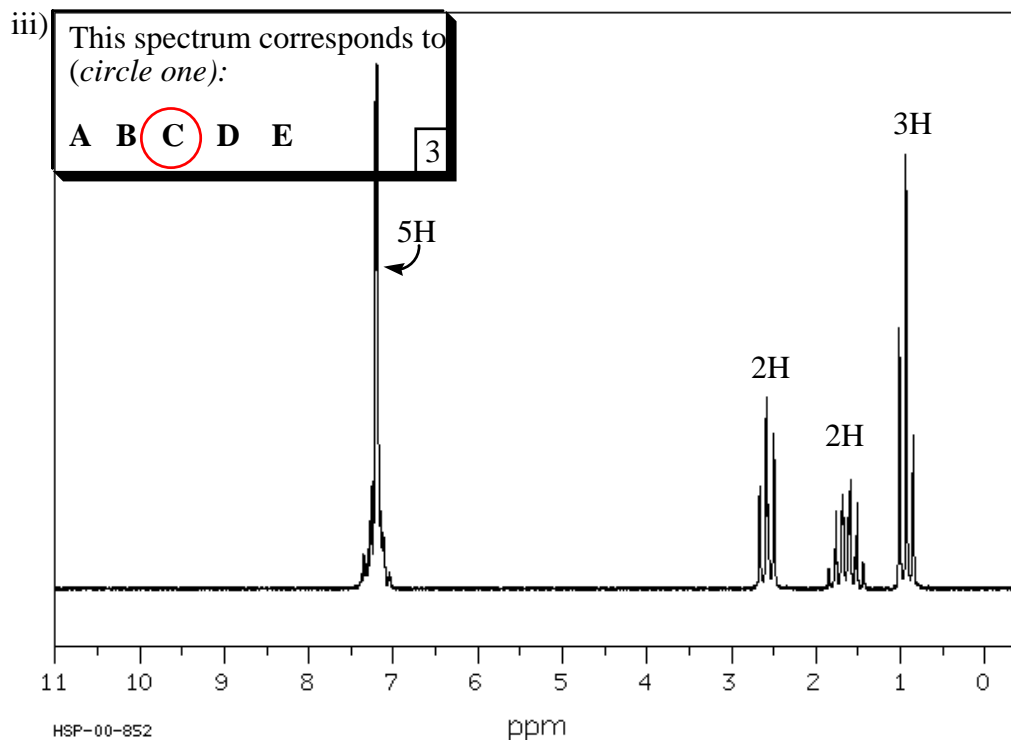
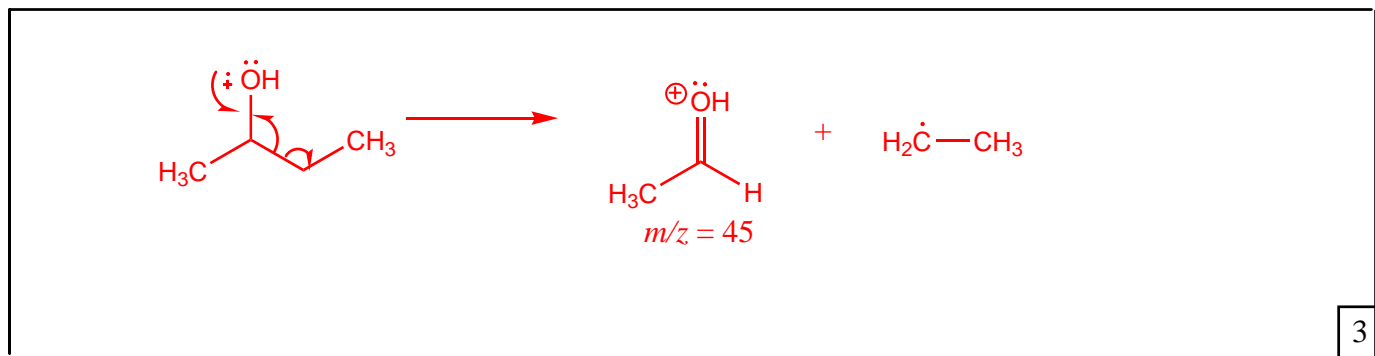


1. Match the following $^1\text{H-NMR}$ spectra to the appropriate C_9H_{12} molecule below:

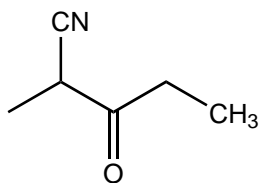




2. 2-butanol shows the base peak (highest intensity) in its mass spectrum at $m/z = 45$. Draw the molecular ion of 2-butanol (M^+) and the fragmentation mechanism that leads to this peak. Use these molecular masses in your calculations: C = 12; O = 16; H = 1.



3. Which of the following features would you **not** expect to see in the IR spectrum of the following molecule (circle one)?



a strong absorption at 1710 cm^{-1}

a medium absorption at 2250 cm^{-1}

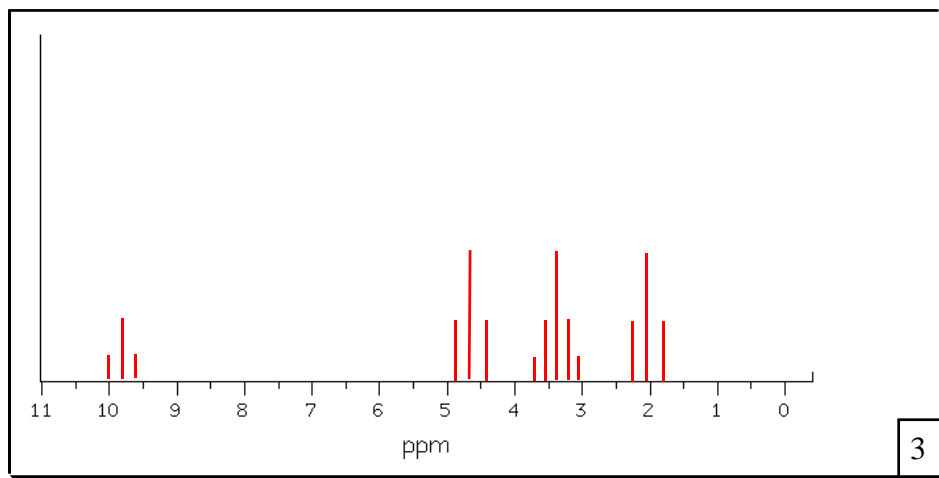
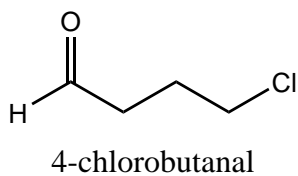
a strong absorption at 2900 cm^{-1}

a strong absorption at 3100 cm^{-1}

a medium absorption at 1465 cm^{-1}

1

4. Draw an $^1\text{H-NMR}$ spectrum for 4-chlorobutanal (shown below). Estimate chemical shifts and draw proper splitting patterns. Draw a multiplet for any splitting > 6 .



5. Use the following IR and $^1\text{H-NMR}$ spectra to deduce the structure of the proper $\text{C}_{11}\text{H}_{13}\text{BrO}_2$ isomer.

