

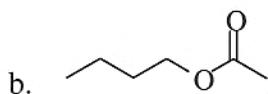
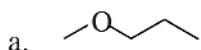
NOTE: This Version is Longer than the Real Test Will Be

- Predict the ^1H NMR spectra for the following molecules. Include predicted:
 - chemical shifts
 - integration
 - splitting pattern (singlet, doublet, triplet, quartet, etc., multiplet)

Example



3's, 2H, t
1's (or 2's), 2H, pentet (or multiplet)
3's, 2H, t



- Assign the dimethylbenzene isomer for which the ^{13}C NMR spectrum has:

a. 3 signals (q, s, d)

b. 4 signals (q, s, d, d)

c. 5 signals (q, s, d, d, d)

- Match the circled proton or protons in the following compounds with the correct chemical shift.

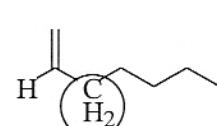
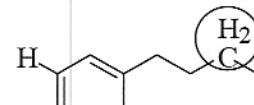
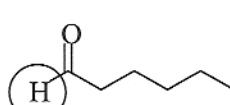
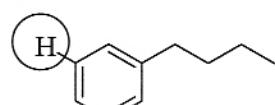
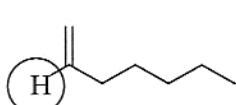
1.20

2.05

5.70

7.17

9.55

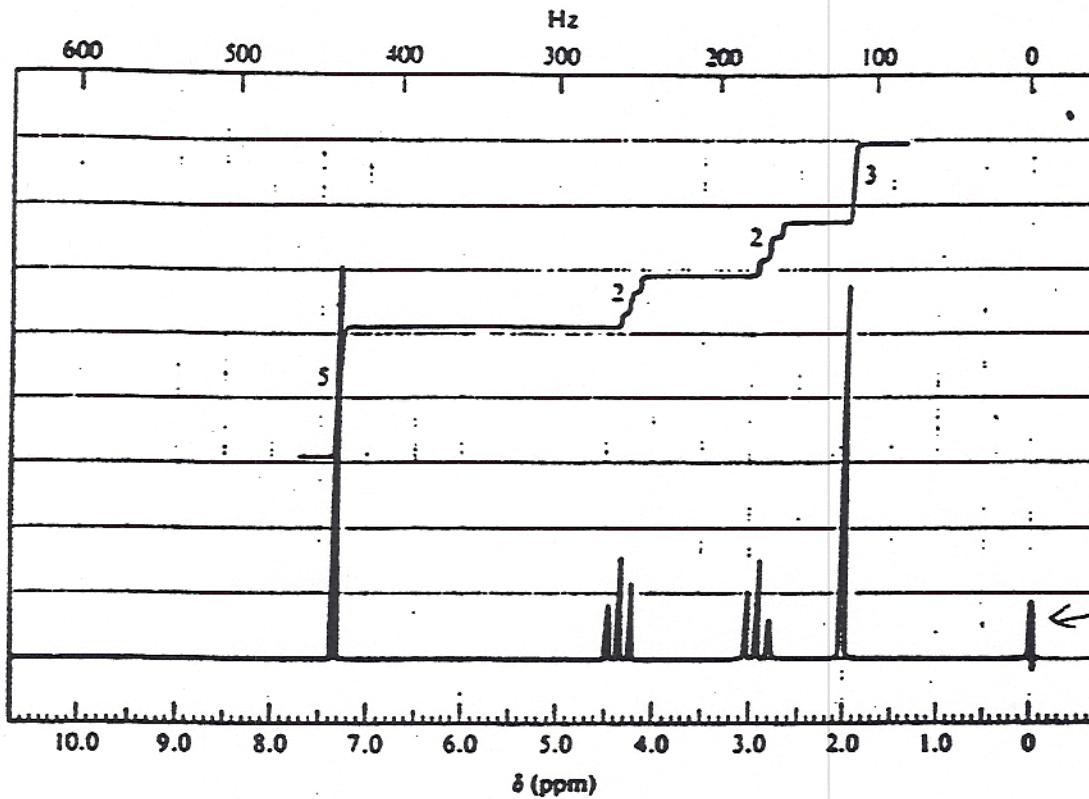


DRAW STRUCTURES FOR THE MOLECULES IN PROBLEMS 3-9

3. $C_{10}H_{12}O_2$

IR: 1740 (strong), 750 (strong), 700 (strong)

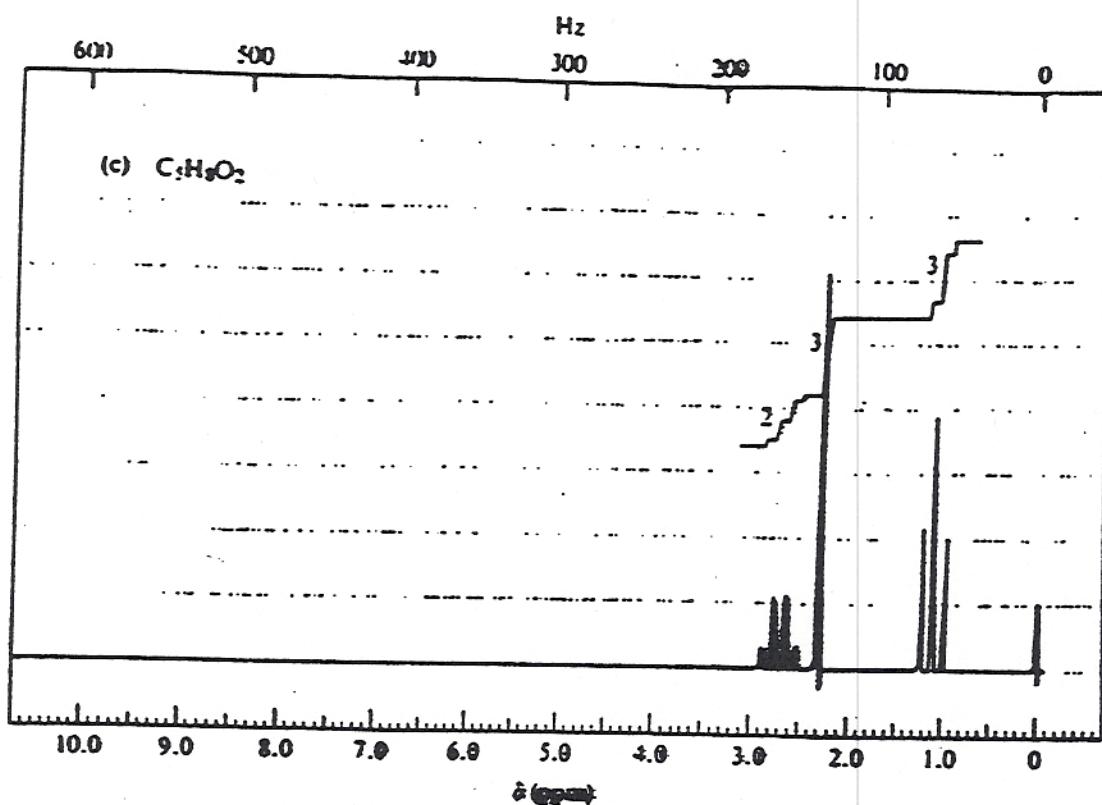
^{13}C NMR: 185 (s), 155 (s), 135 (d), 130 (d), 128 (d), 35 (t), 28 (t), 20 (q)



Note:
Signal
at zero
is reference
signal, nothing
in actual
molecule.

(4) $C_5H_8O_2$

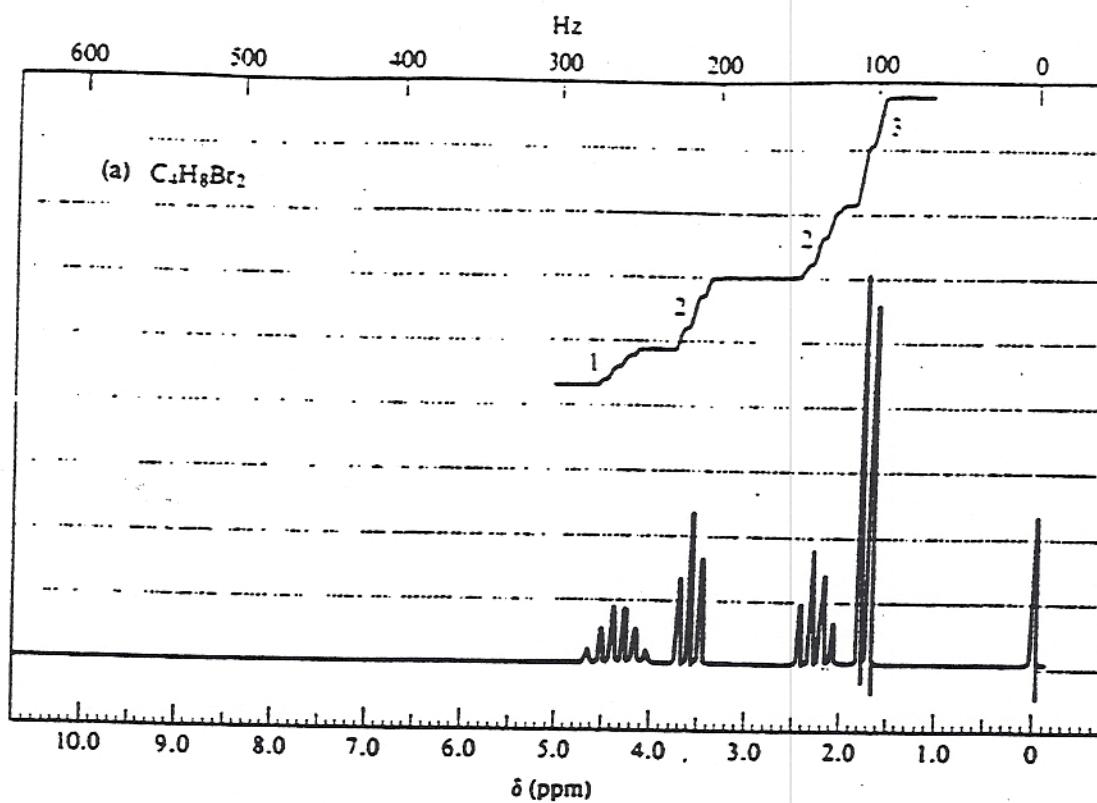
IR: 1720 (s), 1725 (s)



5. $C_4H_8Br_2$

IR: nothing interesting

^{13}C NMR: 45 (d), 37 (t), 24 (t), 18 (q)



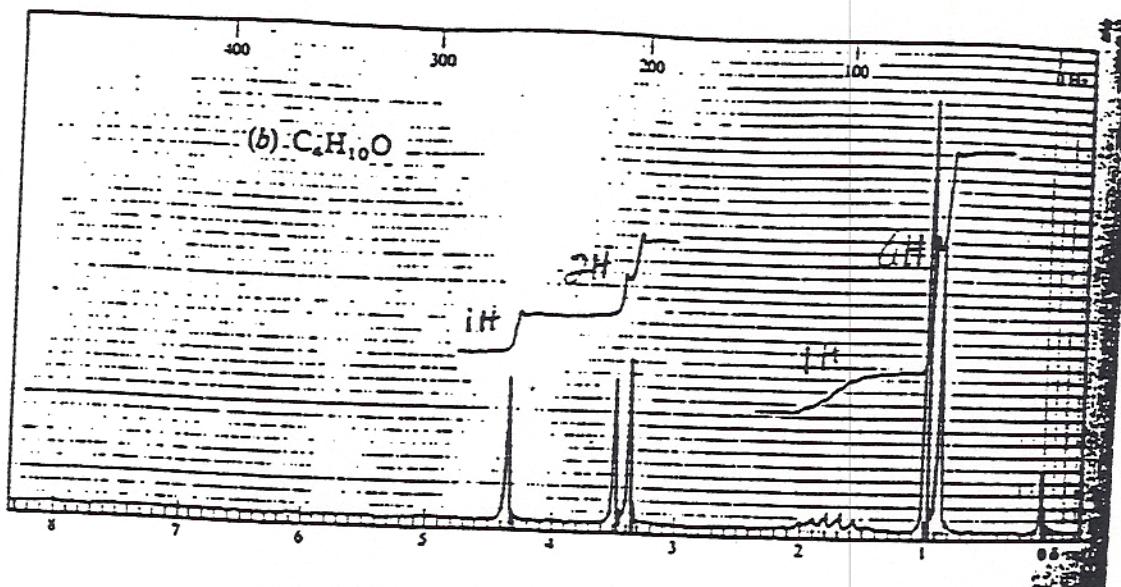
(6) C₁₁H₁₆O

IR: 820 (strong)

¹³C: 145 (s), 132 (s), 128 (d), 120 (d), 75 (t), 35 (d), 20 (q), 18 (q)

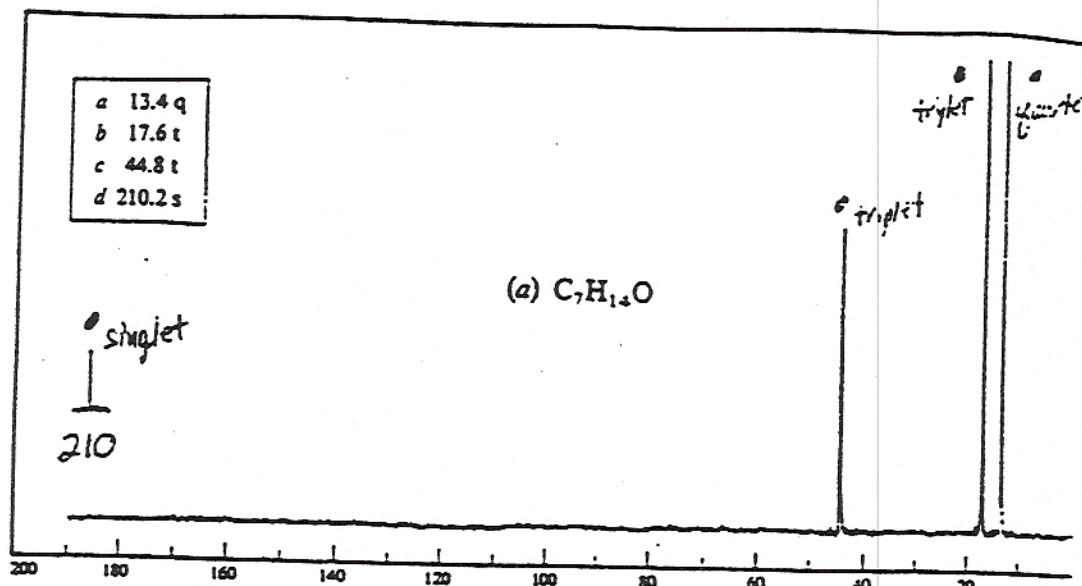
¹H NMR: 1.25 (6H, d), 1.30 (3H, t), 2.90 (m, 1H), 4.15 (2H, q), 6.66 (2H, d), 6.97 (2H, d)

7c C₄H₁₀O

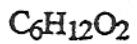


(8c) C₇H₁₄O

IR: 1710 (strong)

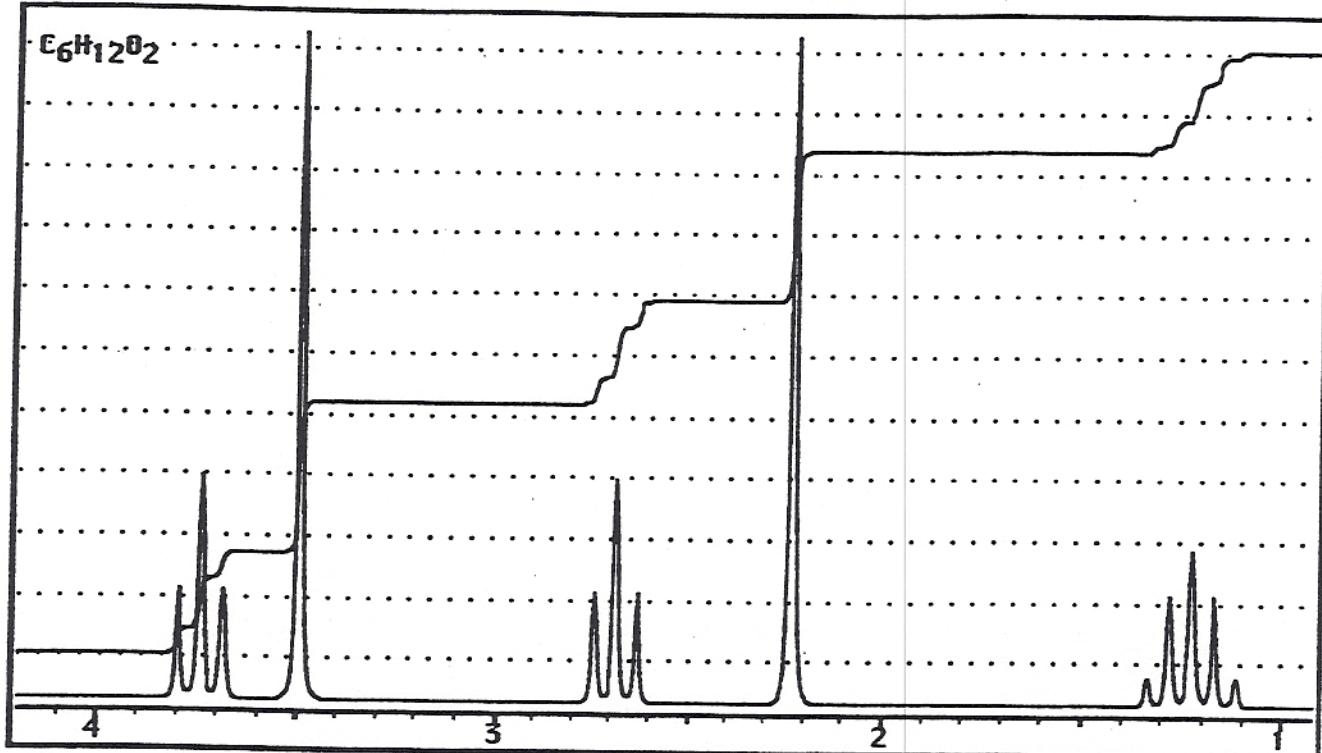


Q
18. Show the structures for the following molecule, based on the spectroscopic information provided. (10 points)



IR: 1710, strong

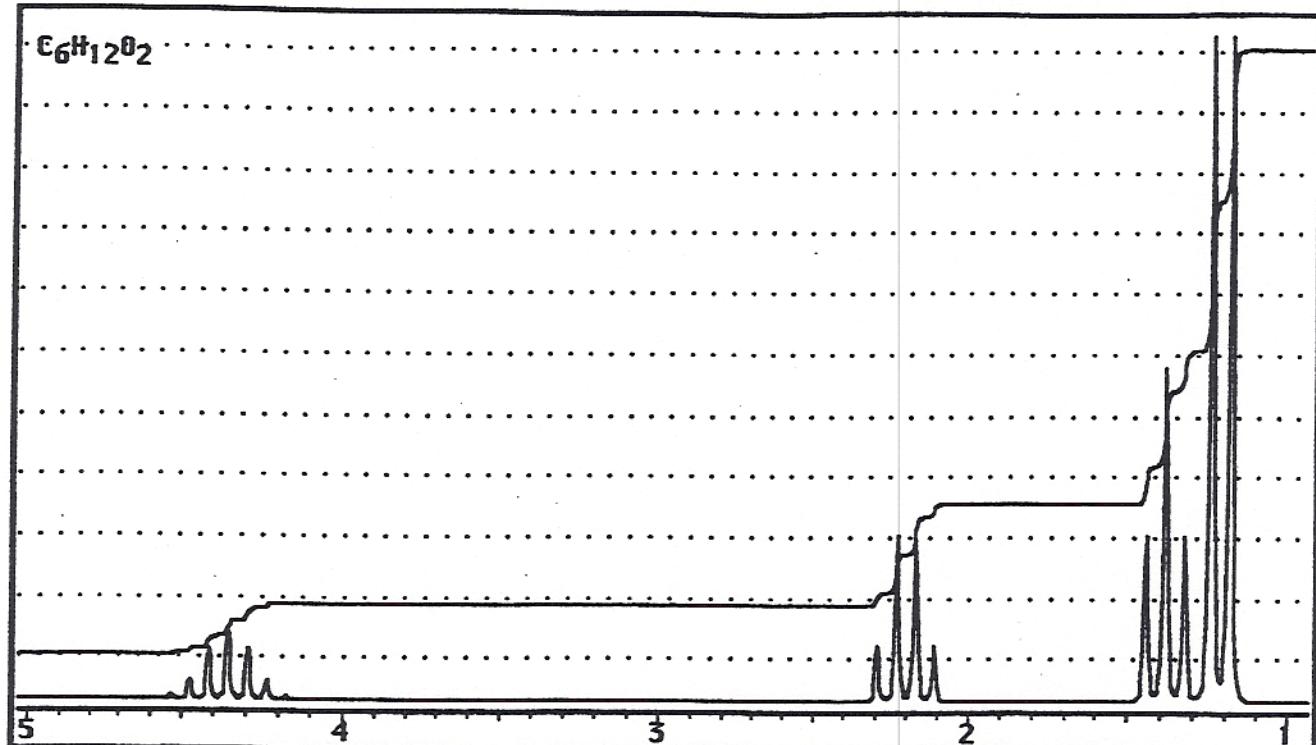
^{13}C NMR: 200 (s), 75 (t), 65 (q), 40 (t), 30 (t), 20 (q)



10. Show the structures for the following molecule, based on the spectroscopic information provided. (10 points)

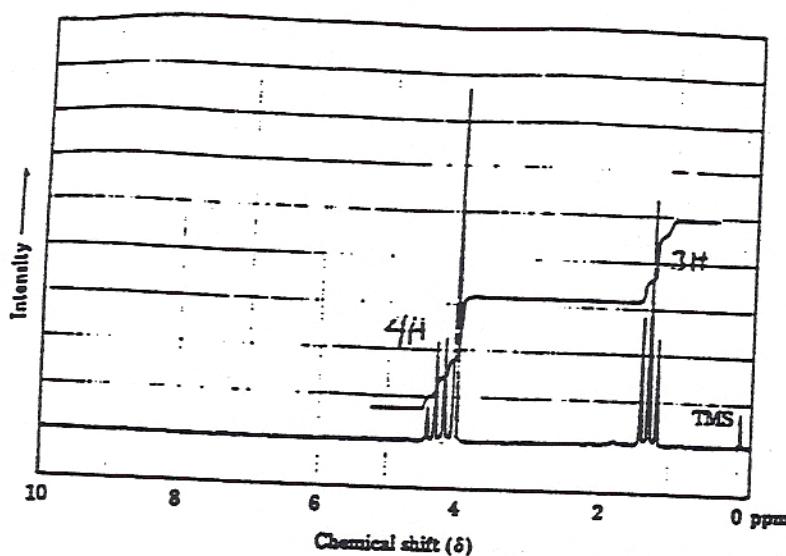
C₆H₁₂O₂ IR: 1745, strong

¹³C NMR: 20 (q), 30 (q), 48 (t), 78 (d), 185 (s)





IR: 1740



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