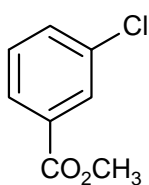


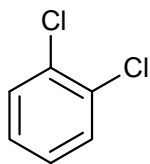
# Problem Set 3

Chem 605  
Reich

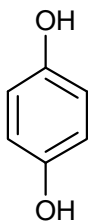
**Problem R-70C.** For each of the structures below, predict the number of carbon signals which will be seen in the room temperature  $^{13}\text{C}$  NMR spectrum. Assume that *rotation around single bonds will be fast*, and that there will be *no accidental equivalences*. Use the form: 6 peaks, 2:2:1:1:1:1. For diastereotopic effects (there are a number here) see Section 5-HMR-8.



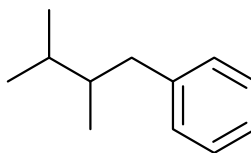
a \_\_\_\_\_



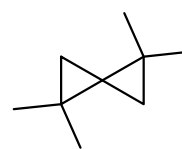
b \_\_\_\_\_



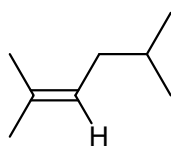
c \_\_\_\_\_



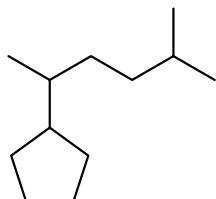
d \_\_\_\_\_



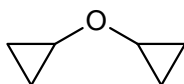
e \_\_\_\_\_



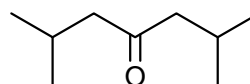
f \_\_\_\_\_



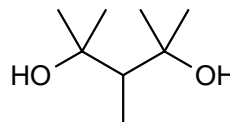
g \_\_\_\_\_



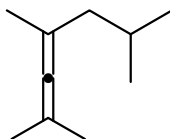
h \_\_\_\_\_



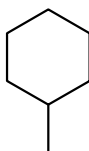
i \_\_\_\_\_



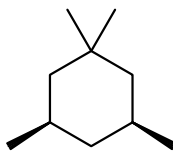
j \_\_\_\_\_



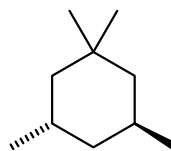
k \_\_\_\_\_



l \_\_\_\_\_



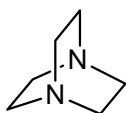
m \_\_\_\_\_



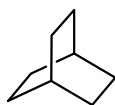
n \_\_\_\_\_



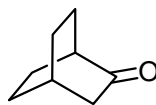
o \_\_\_\_\_



p \_\_\_\_\_



q \_\_\_\_\_



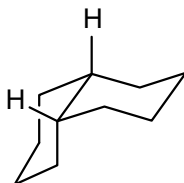
r \_\_\_\_\_



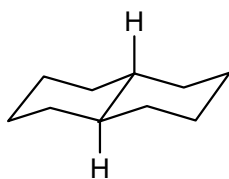
s \_\_\_\_\_



t \_\_\_\_\_



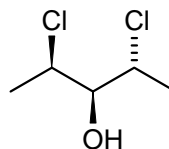
u \_\_\_\_\_



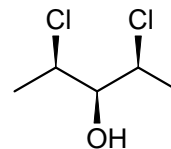
v \_\_\_\_\_



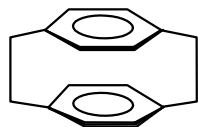
w \_\_\_\_\_



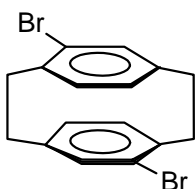
x \_\_\_\_\_



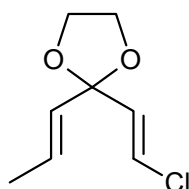
y \_\_\_\_\_



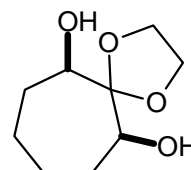
z \_\_\_\_\_



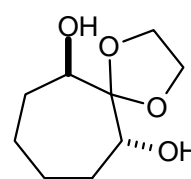
alpha \_\_\_\_\_



beta \_\_\_\_\_



gamma \_\_\_\_\_



delta \_\_\_\_\_

# Problem Set 3

Chem 605  
Reich

**Problem R-70C.** For each of the structures below, predict the number of carbon signals which will be seen in the room temperature  $^{13}\text{C}$  NMR spectrum. Assume that *rotation around single bonds will be fast*, and that there will be *no accidental equivalences*. Use the form: 6 peaks, 2:2:1:1:1:1. For diastereotopic effects (there are a number here) see Section 5-HMR-8.

