

## CONCEPT: $^1\text{H}$ NMR – SPIN-SPLITTING WITHOUT J-VALUES

Also known as *spin-spin coupling*, or *J-coupling*, this describes the distances between different protons.

**Note:** This topic can be taught with or without J-values. Check with your professor to determine how much detail you should learn. For now, we will start with the *simplest explanation*, (should suffice for 90% of professors), which is without J-values.

- Adjacent, \_\_\_\_\_ - \_\_\_\_\_ protons will split each other's magnetic response to the NMR
  - ☐ We use the \_\_\_\_\_ rule to determine how many splits we will achieve
  - ☐ *Pascal's Triangle* predicts the shape of the splits we will get

Pascal's Triangle with the  $(n + 1)$  Rule

n									
0	singlet								
1	doublet								
2	triplet								
3	quartet								
4	quintet								
5	sextet								
6	septet								
7	octet								

1

1 1

1 2 1

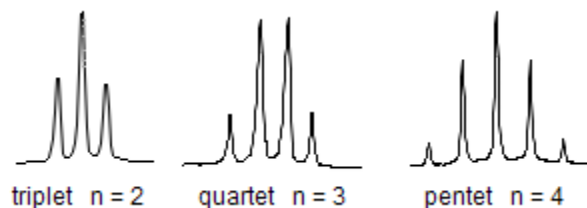
1 3 3 1

1 4 6 4 1

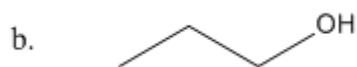
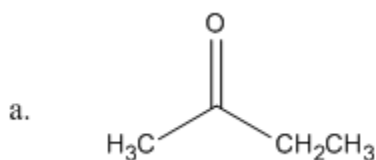
1 5 10 10 5 1

1 6 15 20 15 6 1

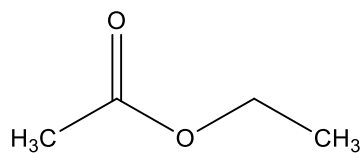
1 7 21 35 35 21 7 1



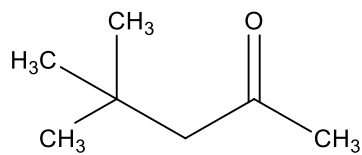
**EXAMPLE:** How will the following protons be split?



PRACTICE: Predict the splitting pattern (multiplicity) for the following molecule:



PRACTICE: Predict the splitting pattern (multiplicity) for the following molecule:



PRACTICE: Which of the following compounds gives a  $^1\text{H}$  NMR spectrum consisting of only a singlet, a triplet, and a pentet?

