## **CONCEPT:** R AND S CONFIGURATION

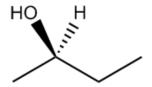
According to IUPAC protocol, each molecule must have a unique, unambiguous name – even stereoisomers.

Step 1. Assign priorities to the four atoms on the chiral center according to their atomic mass on the periodic table.

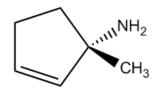
Step 2. When there is a tie between atomic weights, compare the next set of adjacent atoms (playoffs!).

Step 3. Double bonds count twice. Triple bonds count three times.

**EXAMPLE:** Determine priorities for the following chiral center:



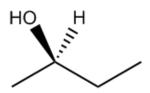
**EXAMPLE:** Determine priorities for the following chiral center:



Step 4. IF the last priority group is in the back, then trace a path from \_\_\_\_\_\_ to \_\_\_\_\_ priority.

- Clockwise = \_\_\_\_\_, Counterclockwise = \_\_\_\_ 

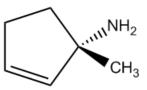
  □ Always Ignore group 4



Step 5. If the last priority group is **NOT in the back**, \_\_\_\_\_ that group with the group that is on the dash.

- Trace path as always, but this time \_\_\_\_\_ the sign since you \_\_\_\_\_ groups.

  - \_\_ > \_\_



**PRACTICE:** Provide the full name for the following molecules, taking stereochemistry into account.

a.

b.

c.