

"Easy" ways to identify chiral molecules:

- * presence of just one chirality center -
if there is only one, then the molecule is definitely chiral.

- * presence of an internal mirror plane (plane of symmetry) - even if chirality centers are present, if the molecule has an internal mirror plane, it will be superimposable on its mirror image + therefore achiral.

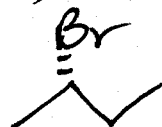
Nomenclature (ZUPAC)

consider:



(R)

+



(S)

* 2-bromobutane

* ~~enantiomers~~

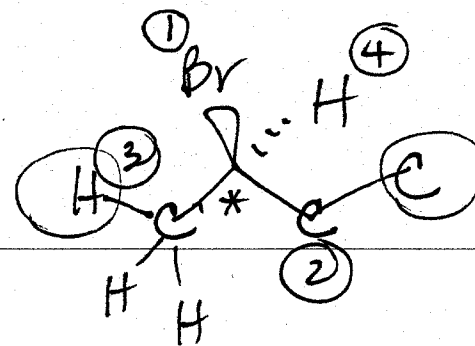
* how can we name them so as to tell them apart?

Use Cahn-Ingold-Prelog rules of prioritization.

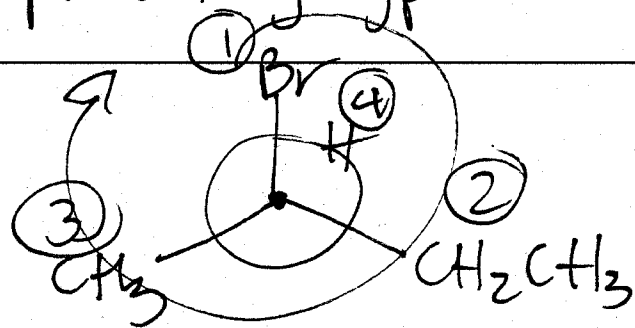
1. Each of the 4 gps. attached to the chirality center is assigned a priority. 1 \rightarrow 4 (high \rightarrow low)

a. Go by atomic #. Higher atomic # = higher priority.

b. If you can't choose based on those atoms, go one atom out.



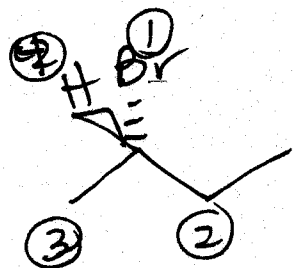
2. Align the molecule so that the lowest priority gp. is pointing away.



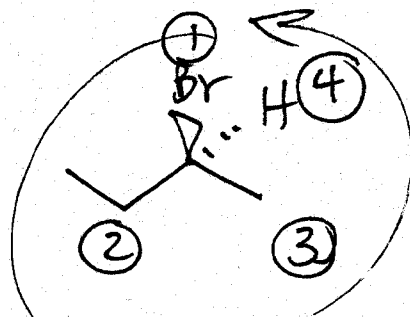
3. Draw a circle 1 → 2 → 3

a. Clockwise - (R) rectus (right)

b. Counterclockwise - (S) sinister (left)



≡



(S)

What if there are multiple bonds?

