

If there is plane of symmetry within a molecule: then the molecule is **Achiral** (not chiral)

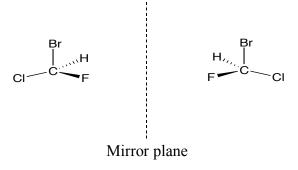
Stereogenic centre (chiral centers or asymmetric centers) - often a tetrahedral atom (carbon) with four different groups attached.

R/S Nomenclature :

R and S designation of stereoisomers

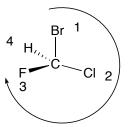
- R = Rectus - S = Sinister

Each stereogenic center analyzed separately. eg.



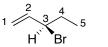
Labeling a stereogenic centre as R or S:

- identify all stereogenic centres (ie. 4 different substituents)
- assign priority based on atomic number (similar to *E* and *Z*)
- if you cannot decide, go to the next set of atoms
- with the lowest group pointing back, count 1, 2, 3:
 - clockwise \rightarrow R configuration, counterclockwise \rightarrow S configuration



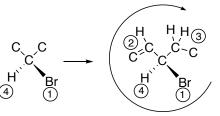


eg.



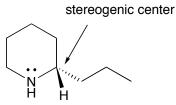
3-Bromo-1-Pentene

- #3 is stereogenic center



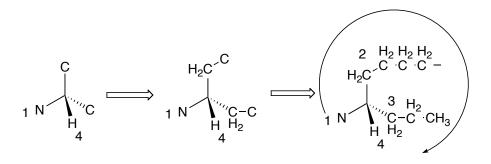
- clockwise So "R" - configuration.

eg.



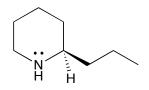
coniine: poison from Hemlock

N can invert fast



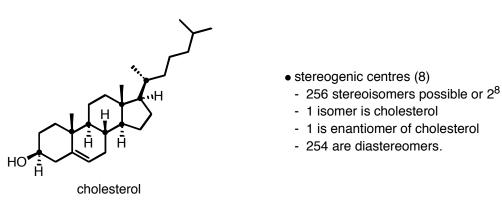
- counting 1, 2, 3 gives clockwise, BUT the smallest group is pointing forward, so the configuration is opposite of what you get if the smallest group is back
- in this case, the configuration of the stereogenic centre is "S"

To draw an enantiomer invert every stereogenic centre



Enantiomer of coniine (non toxic)

Cholesterol



- To draw enantiomer – invert every stereogenic center:

