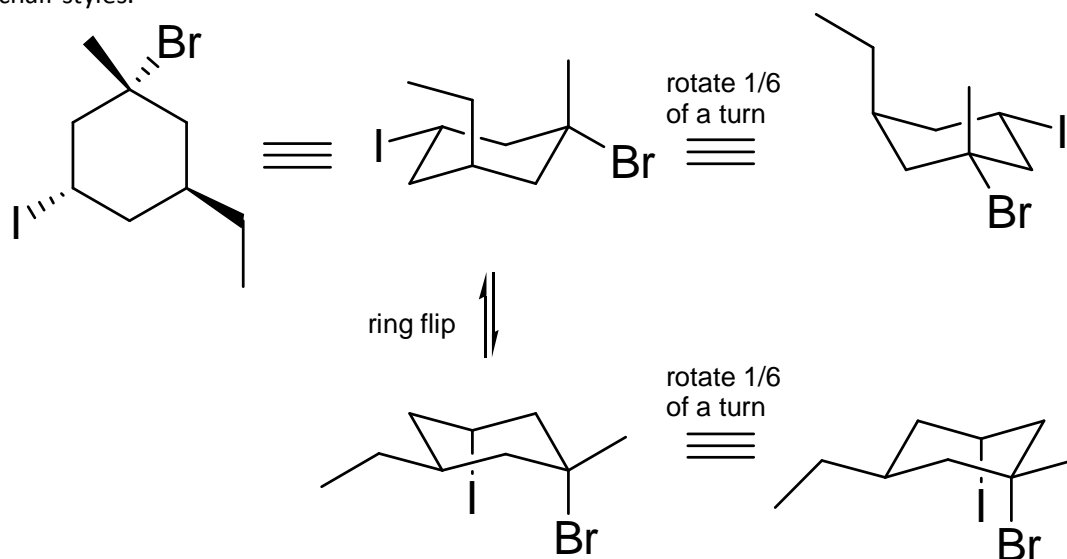
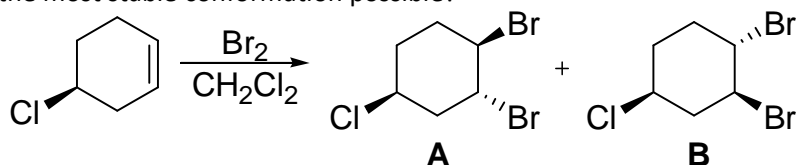


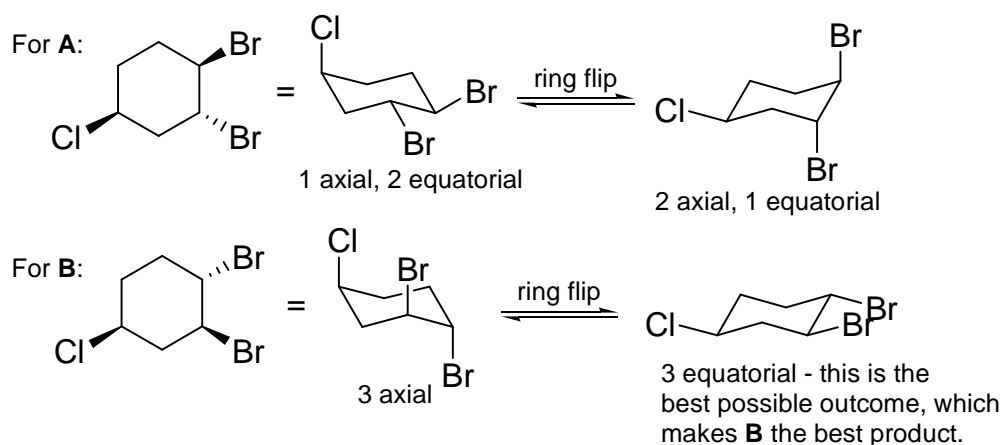
- 1) Using the templates below, draw the flat, top-down view of (1R,3R,5S)-1-bromo-3-ethyl-5-iodo-1-methylcyclohexane. Convert this to one chair conformer, then the other. Show each ring-flip isomer in both chair styles.



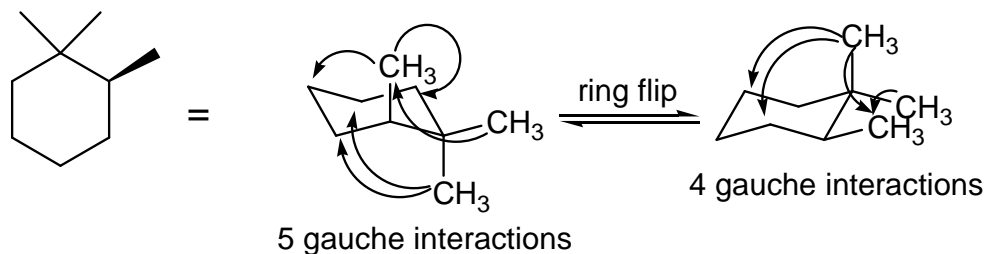
- 2) Reacting (S)-4-chlorocyclohex-1-ene with Br_2 and CH_2Cl_2 gives two products. Are these products enantiomers or diastereomers? Based on the two chair conformations of each of these molecules, which product has the most stable conformation possible?



These are diastereomers, since some chiral centers are the same and some are flipped.



- 3) For the molecule (S)-1,1,2-trimethylcyclohexane, how many total gauche interactions exist in each of its two ring-flip forms?



Both of these forms have at least one gauche interaction between two groups that are on adjacent carbons. You can see this more clearly with a Newman projection. In the long run, it's still true that the form with more axial groups has more gauche interactions and is less stable. However, it's not by as big a margin as you might initially think.