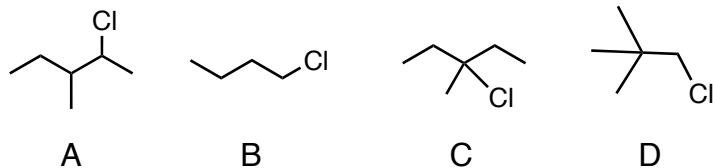
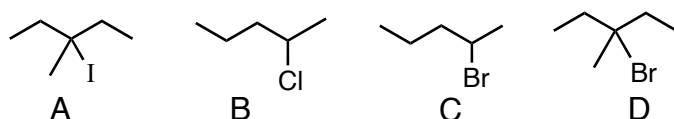


**DUE in class: Wednesday November 15 @ 8am**

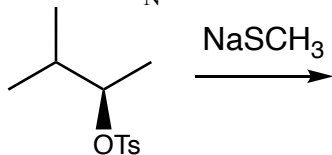
1. Rank the following substrates according to their expected rate of reaction with a nucleophile in an **S<sub>N</sub>2 reaction** from fastest >>> slowest.



2. Rank the following substrates according to their expected rate of reaction with a nucleophile in an **S<sub>N</sub>1 reaction** from fastest >>> slowest.



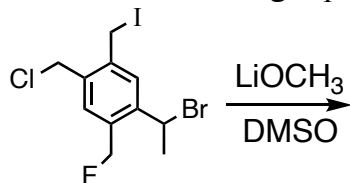
3. Draw the product and determine the configuration (R or S) for the starting material AND the product resulting from the S<sub>N</sub>2 reaction below:



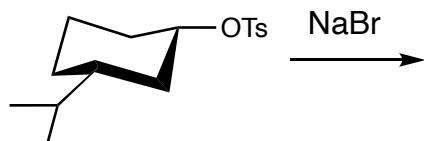
4. Which one of the following statements is true about the substrate 3-bromo-3-methylhexane?

- A. It can react via both the S<sub>N</sub>1 and S<sub>N</sub>2 substitution reaction mechanisms
- B. It cannot react by either the S<sub>N</sub>1 or S<sub>N</sub>2 substitution reaction mechanisms
- C. It can react via the S<sub>N</sub>1 but not the S<sub>N</sub>2 substitution reaction mechanism
- D. It can react via the S<sub>N</sub>2 but not the S<sub>N</sub>1 substitution reaction mechanism

5. The S<sub>N</sub>2 reaction below was conducted with only ONE equivalent of LiOCH<sub>3</sub> (i.e. only sufficient LiOCH<sub>3</sub> to substitute at **a single** position). What is the predicted major product?

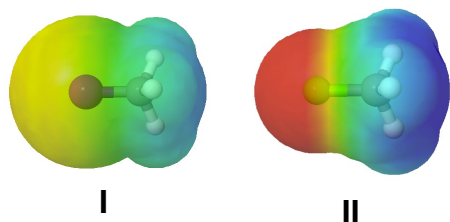


6. The name of the product formed from the following  $S_N2$  reaction is (HINT: redraw the chair conformation as a planar 6-membered ring):



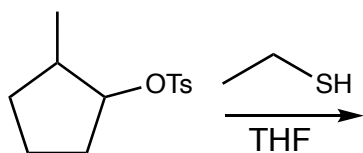
- A. trans-1-bromo-3-isopropylcyclohexane
- B. cis-1-bromo-3-isopropylcyclohexane
- C. trans-1-isopropyl-3-bromocyclohexane
- D. cis-1-isopropyl-3-bromocyclohexane

7. The electrostatic potential map for  $\text{CH}_3\text{I}$  (see I, below) shows much less deep red and blue colors than the electrostatic potential map for  $\text{CH}_3\text{F}$  (see II below). This is because:

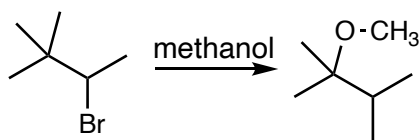


- A. The C-I bond is more polar than C-F so there is more  $\delta^+$  on I and more  $\delta^-$  on C
- B. The C-I bond is more polar than C-F so there is more  $\delta^-$  on I and more  $\delta^+$  on C
- C. The C-F bond is more polar than C-I so there is more  $\delta^-$  on F and more  $\delta^+$  on C
- D. The C-F bond is more polar than C-I so there is more  $\delta^+$  on F and more  $\delta^-$  on C

8. Predict whether the following reaction is likely to proceed by an  $S_N1$  or  $S_N2$  reaction and *justify your answer* (i.e., consider the nature of the nucleophile, nature of the substrate, nature of the leaving group, and the solvent):



9. Draw the individual steps (including curved arrows!!) for the reaction mechanism to explain the following  $S_N1$  reaction:



10. Jimmy says the following reaction will proceed via an  $S_N2$  mechanism because "OTs is an excellent leaving group and  $\text{CH}_3\text{OH}$  is a strong nucleophile". Do you agree with Jimmy? Why or why not? What major final product is expected to form from this reaction (HINT: consider possible rearrangements)?



Jimmy

