

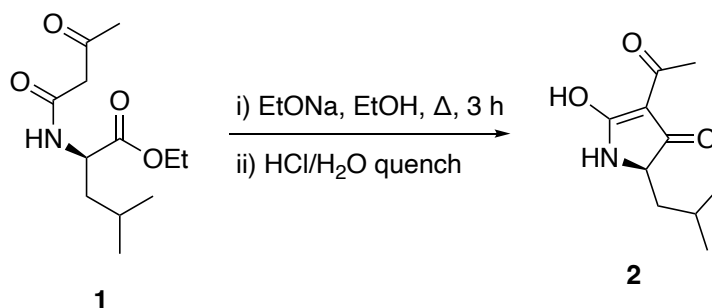
**CHEM 8321/4321****uploaded September 11, 2023****Problem Set #2****T. R. Hoye****Due in class, Monday September 18, 2023**

**Detailed Mechanism** Provide a detailed mechanism [i.e., explicitly show (using curly arrows) *EVERY* intermediate, formal charge (where relevant), equilibrium, and bond-making and -breaking step, including each proton transfer] to account for reactions **a** and **b**.

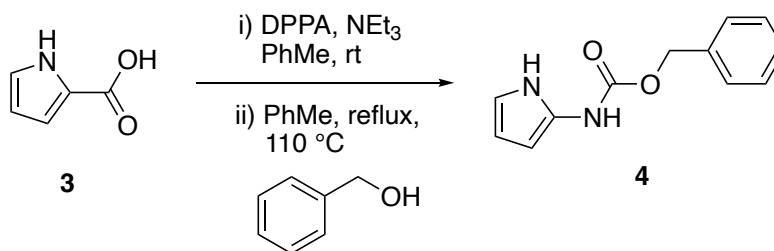
In most cases, mechanisms of reactions proceeding under acid (or cationic) conditions involve *only* neutral or positively charged intermediates,\* and mechanisms of reactions proceeding under basic (or anionic) conditions involve *only* neutral or negatively charged intermediates.\*

(\* those species derived from the organic starting material. Of course, since charge is neither created nor destroyed, there is *always* a counterion of opposite charge present to balance the cationic or anionic charged intermediate.)

(a) The Dieckmann condensation of compound **1** to form enone **2**.

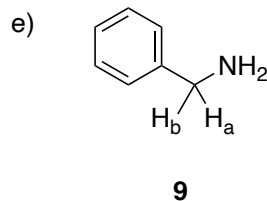
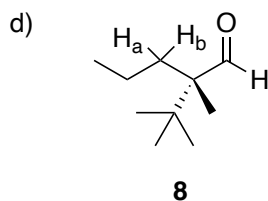
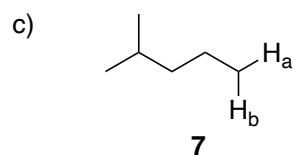
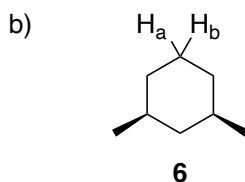
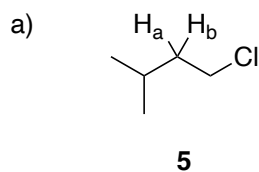


(b) The formation of carbamate **4** by first treating carboxylic acid **3** with diphenylphosphoryl azide (DPPA) then phenylmethanol (aka benzyl alcohol) in refluxing toluene. What is the name of the functional group formed after the first step? (Hint: This is the Curtius rearrangement K&C pp 116–117.)

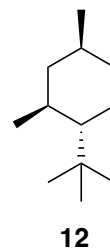
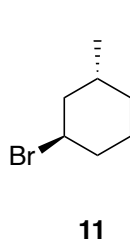
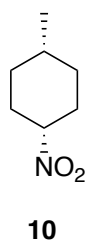


**Other Problems (Stereochemistry)**

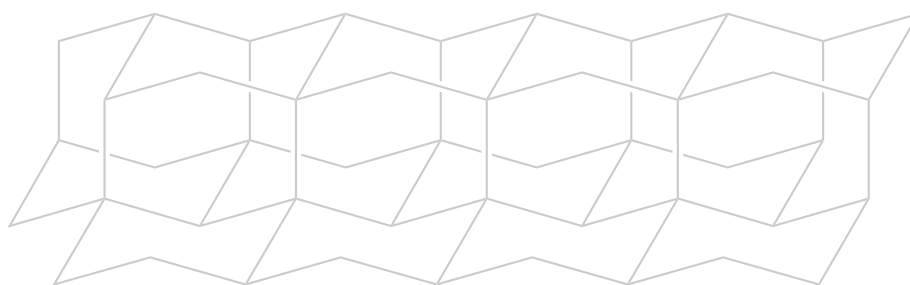
1. Indicate whether the indicated protons,  $H_a$  and  $H_b$ , in the following molecules are homotopic, diastereotopic, or enantiotopic.



2. Draw the most stable conformer of the following compounds.

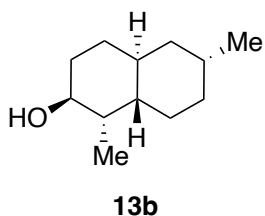
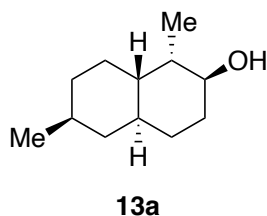


3. Draw a 3D representation of a conformation of 1-phenylhexane that contains one (and only one) syn-pentane (1,7-H,H-) interaction.



4. Indicate (circle the word) whether the two stereoisomers for the following pairs of structures are the *same*, a pair of *enantiomers*, or a pair of *diastereomers*. If they are diastereomers, indicate the number of stereogenic centers that are different in the two structures. Ignore differences in conformation.

a)

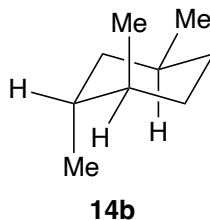
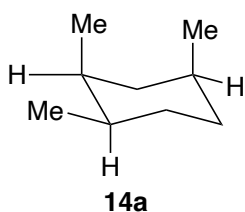


same    enantiomers    diastereomers

# of diff.  
stereocenters?

\_\_\_\_\_

b)

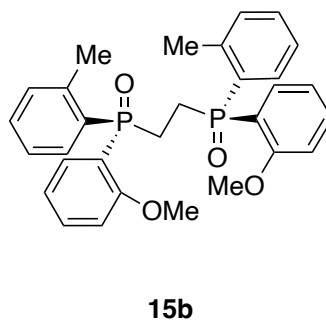
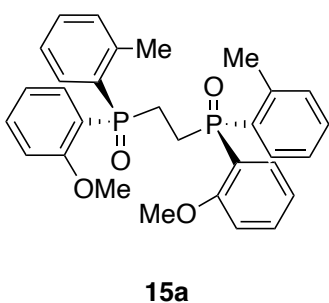


same    enantiomers    diastereomers

# of diff.  
stereocenters?

\_\_\_\_\_

c)

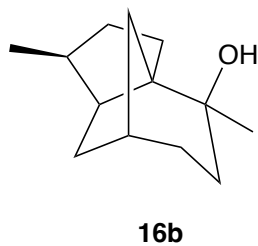
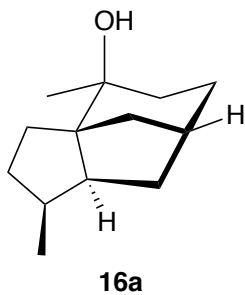


same    enantiomers    diastereomers

# of diff.  
stereocenters?

\_\_\_\_\_

d)



same    enantiomers    diastereomers

# of diff.  
stereocenters?

\_\_\_\_\_