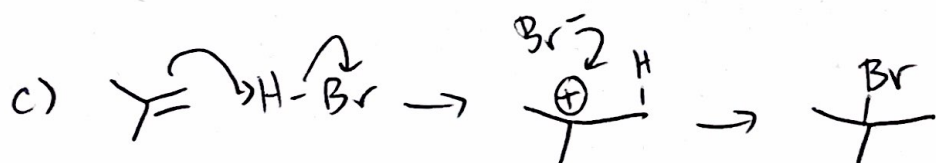
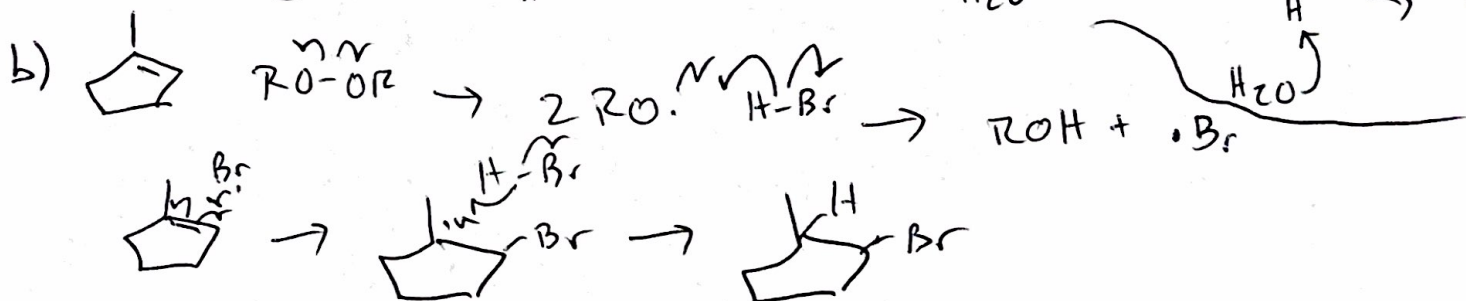
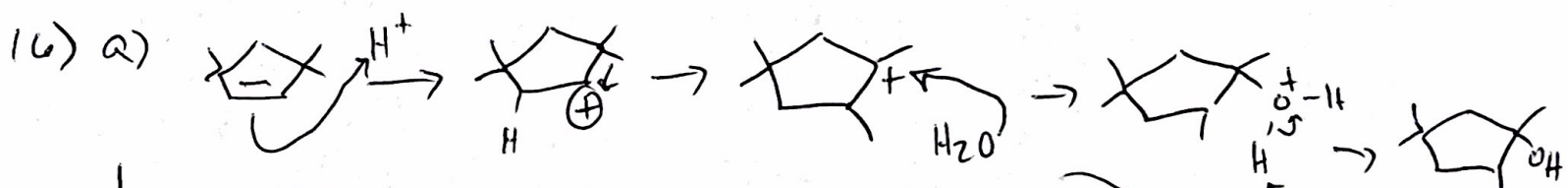
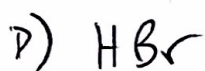
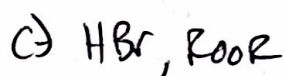
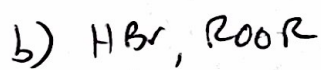
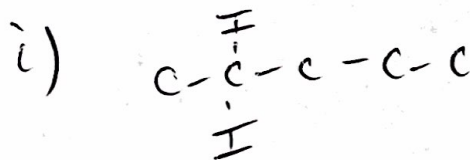
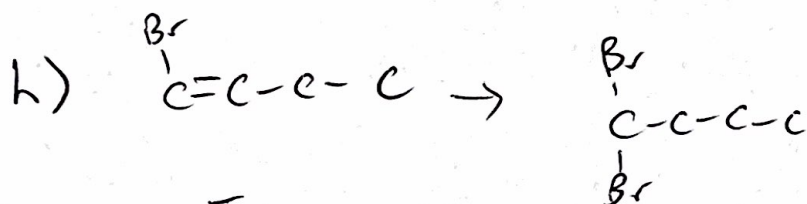
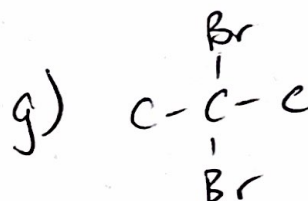
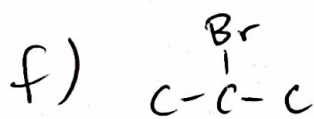
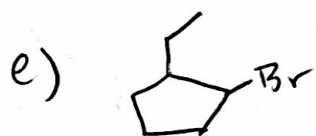
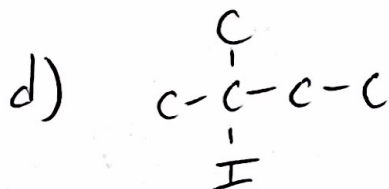
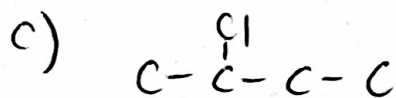
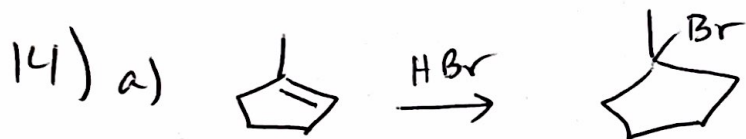
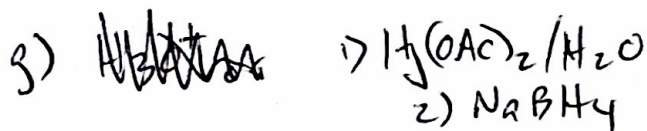
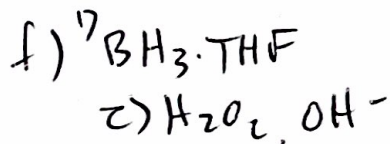
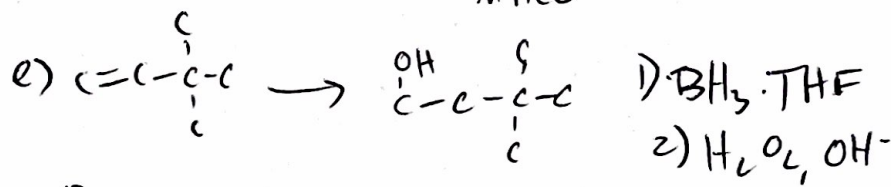
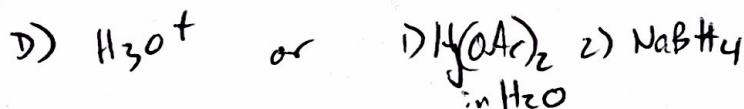
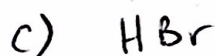
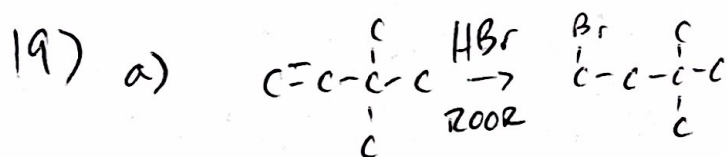
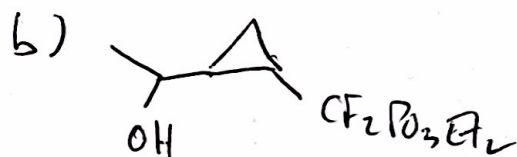
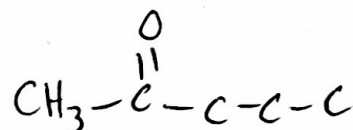
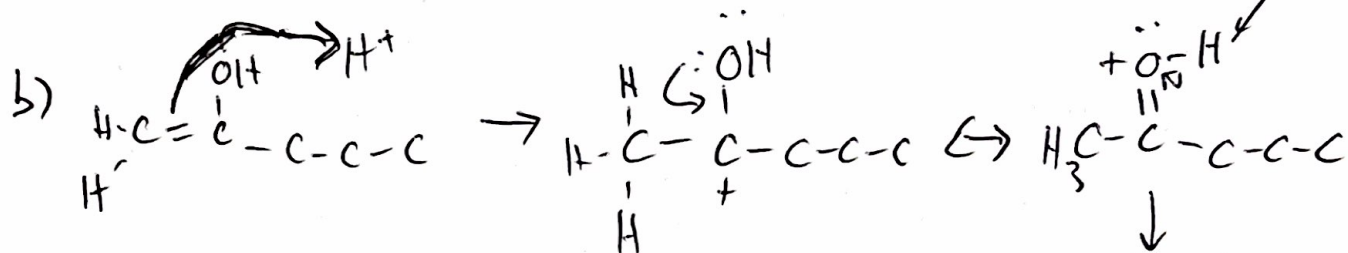
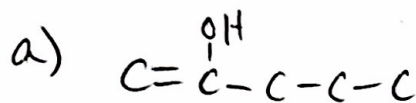
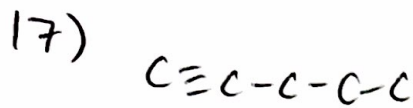
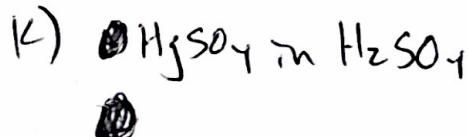
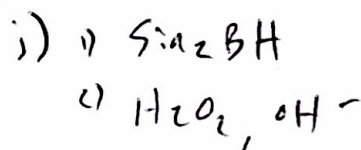
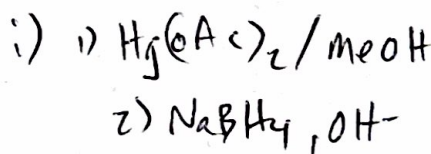


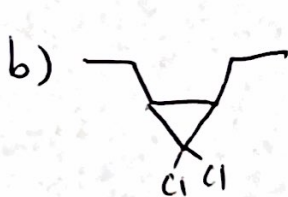
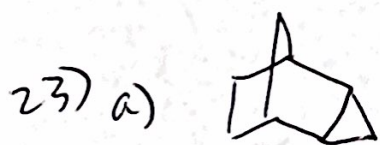
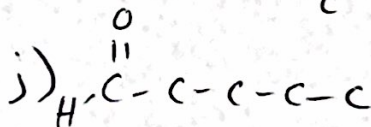
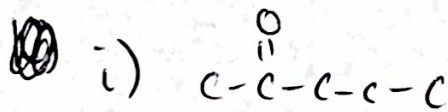
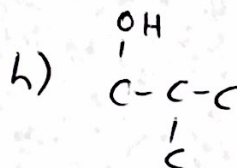
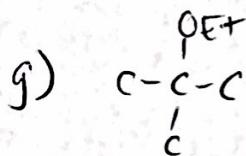
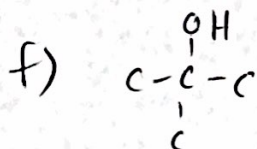
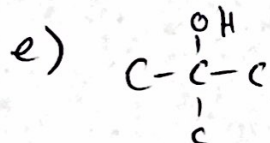
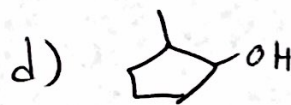
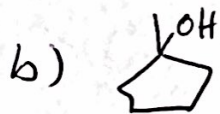
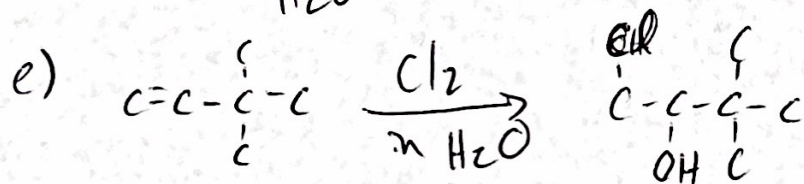
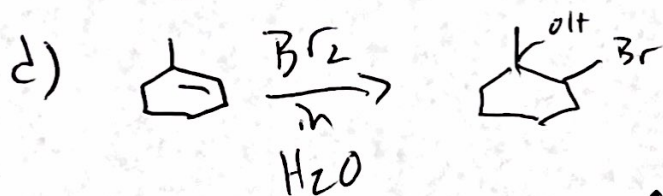
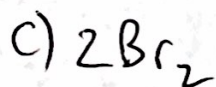
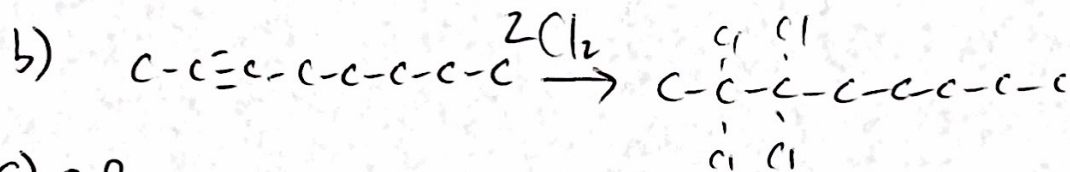
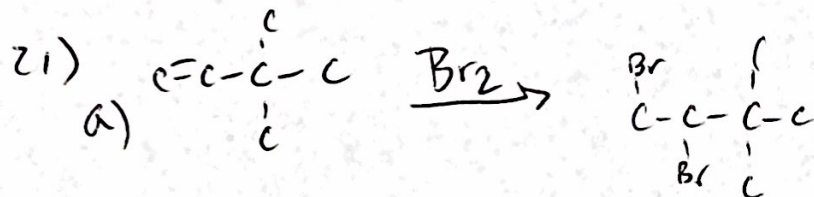
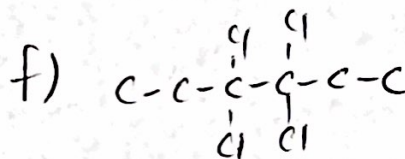
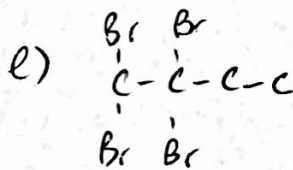
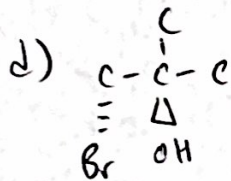
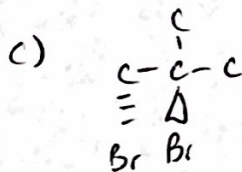
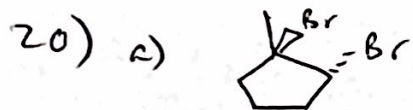
Chapter 11



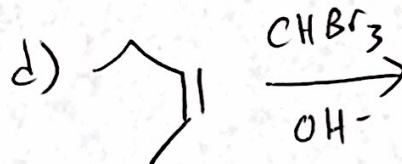
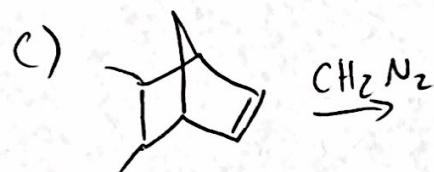
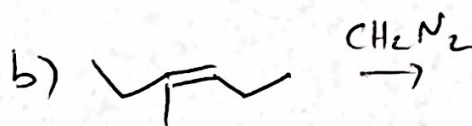


h) yes, if form carbocation it could rearrange w/ a methyl shift



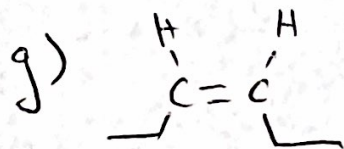
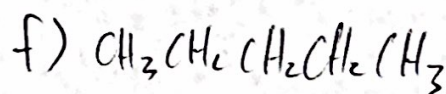
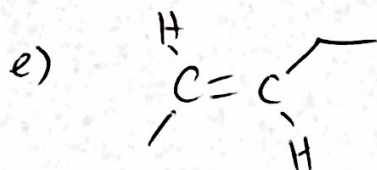
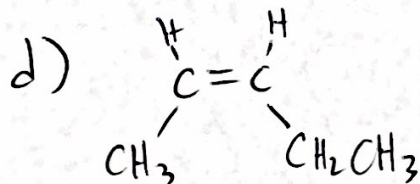
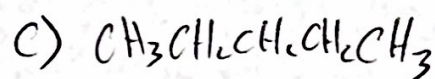
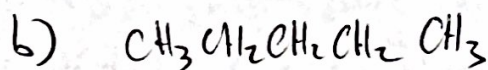
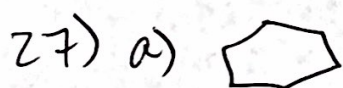
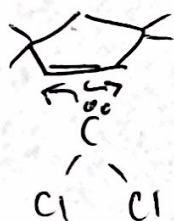
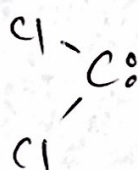
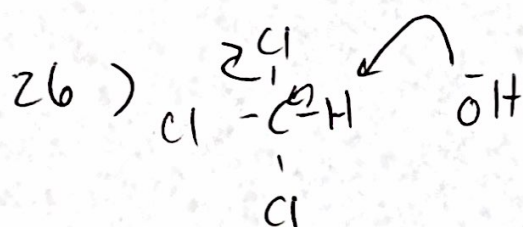


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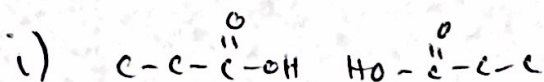
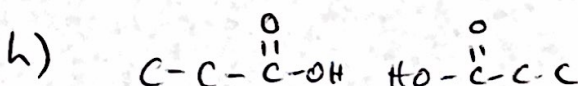
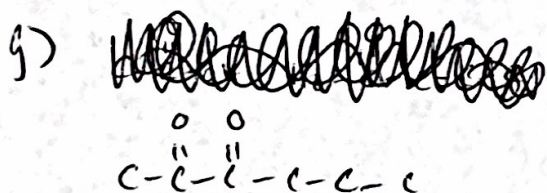
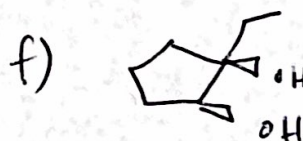
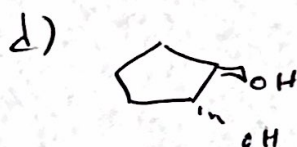
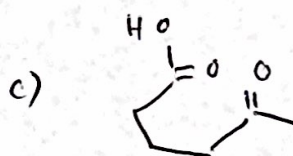
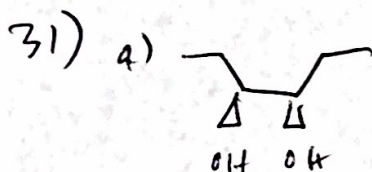
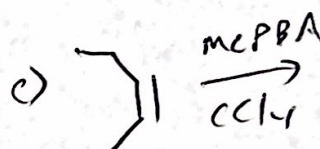
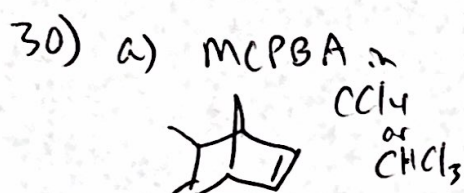
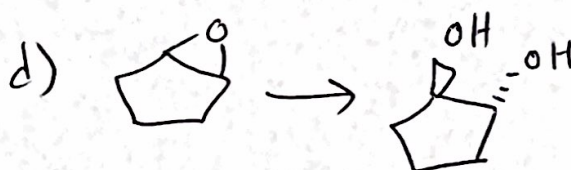
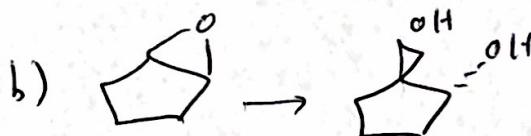
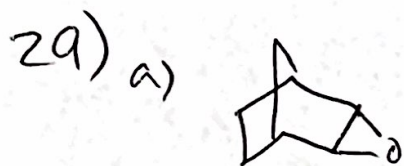


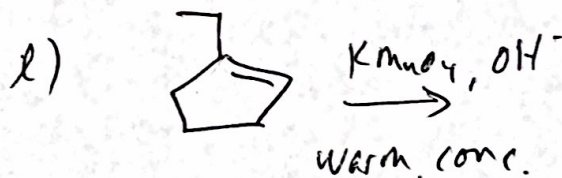
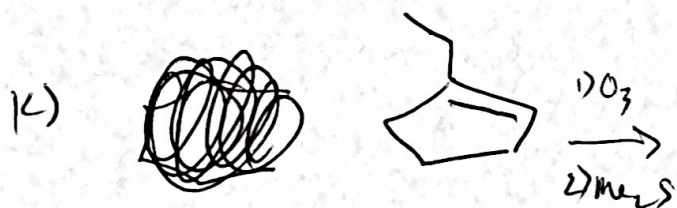
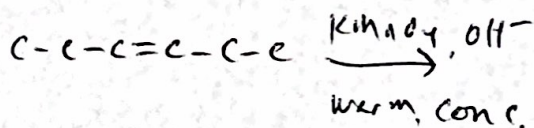
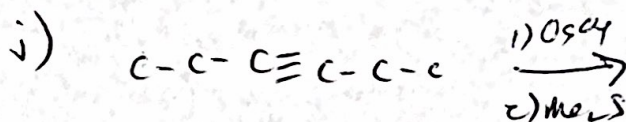
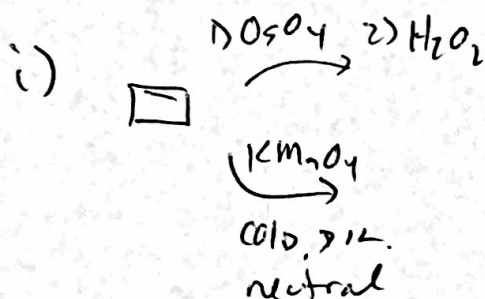
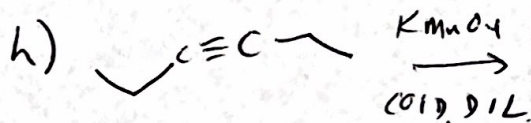
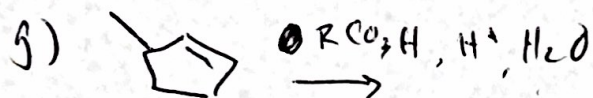
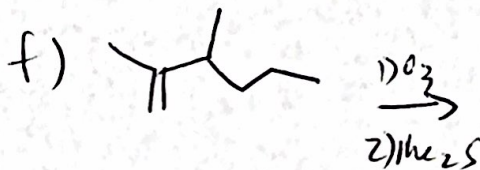
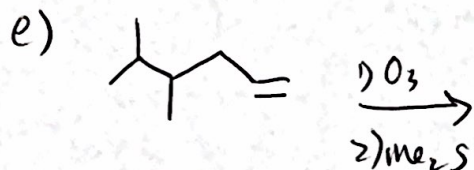
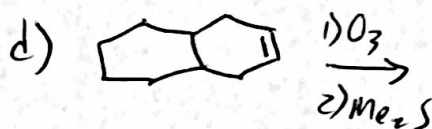
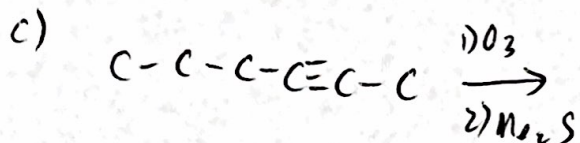
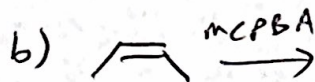
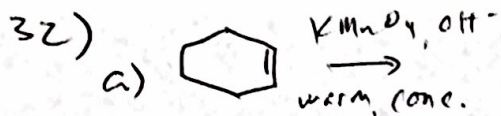
25) a) Simmons Smith Reagent or CH_2N_2

b) carbene

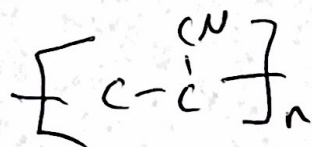
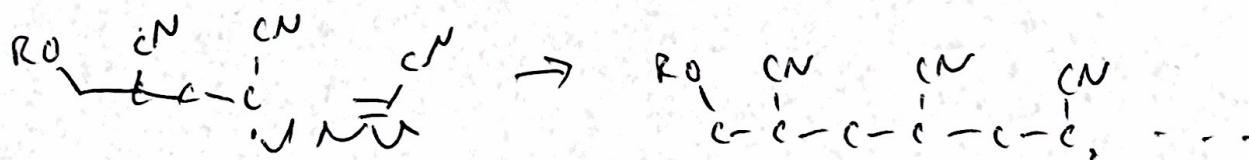
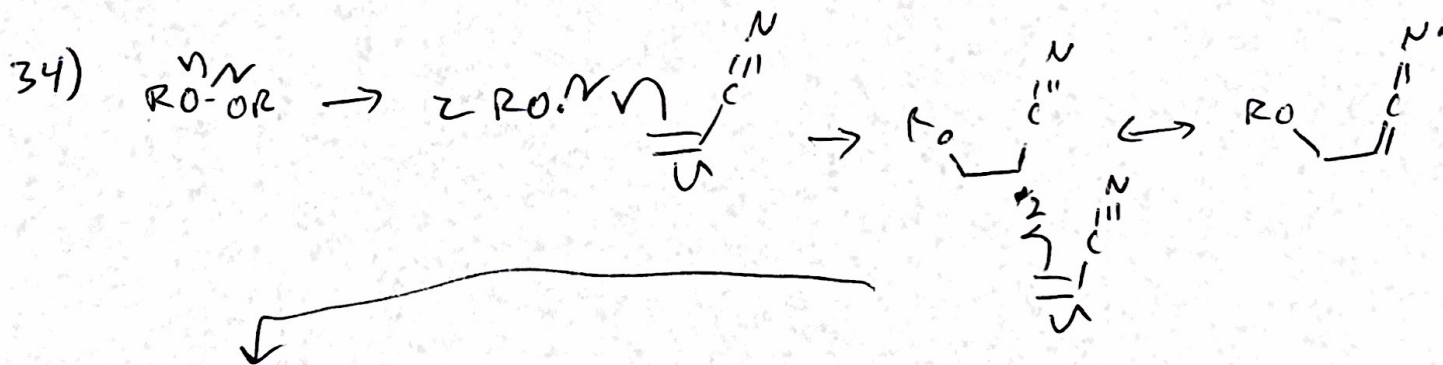
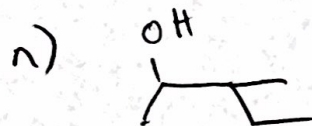
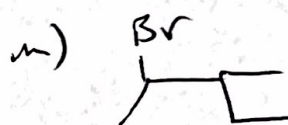
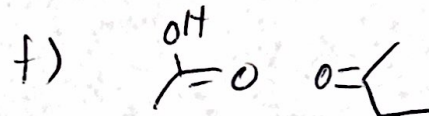
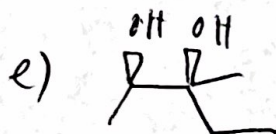
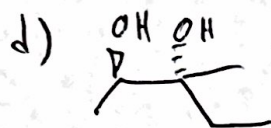
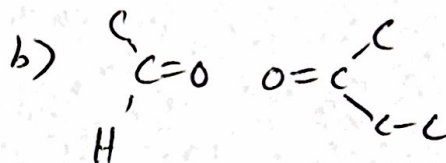
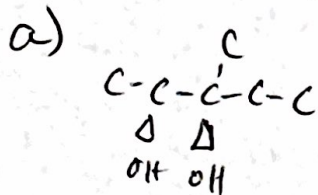
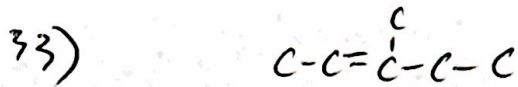


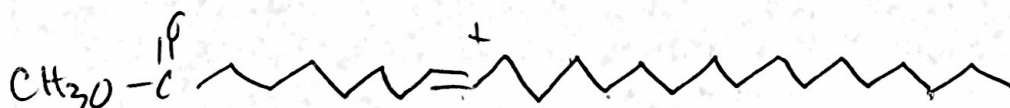
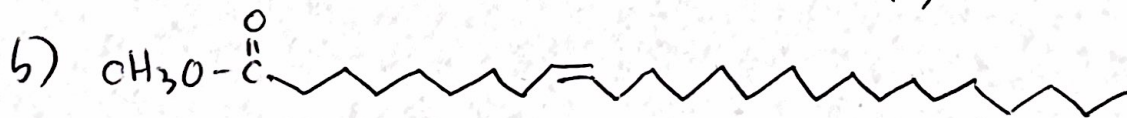
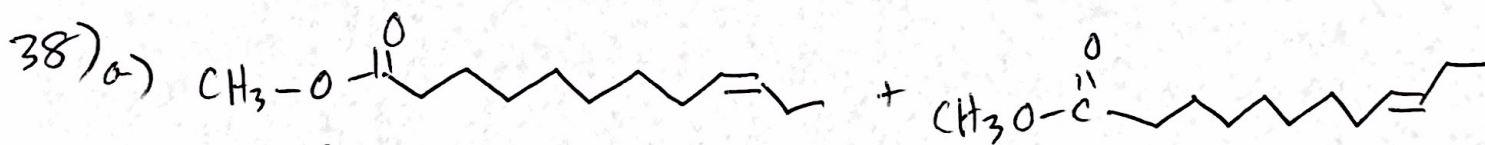
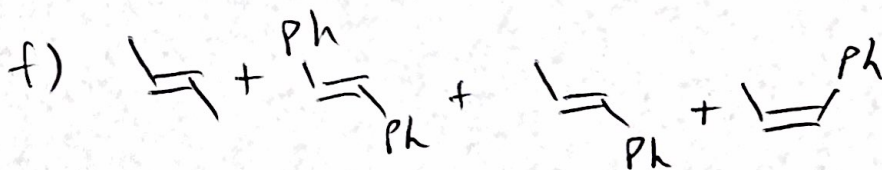
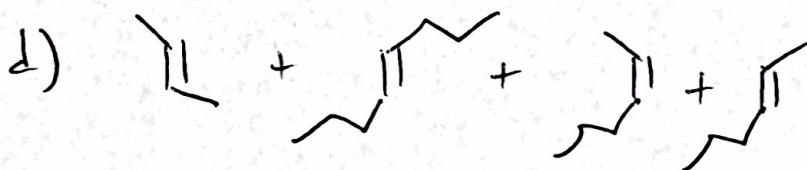
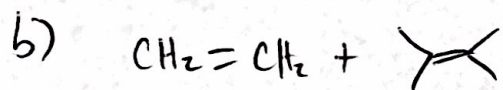
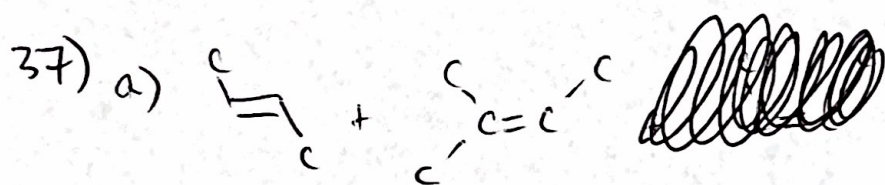
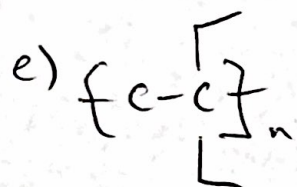
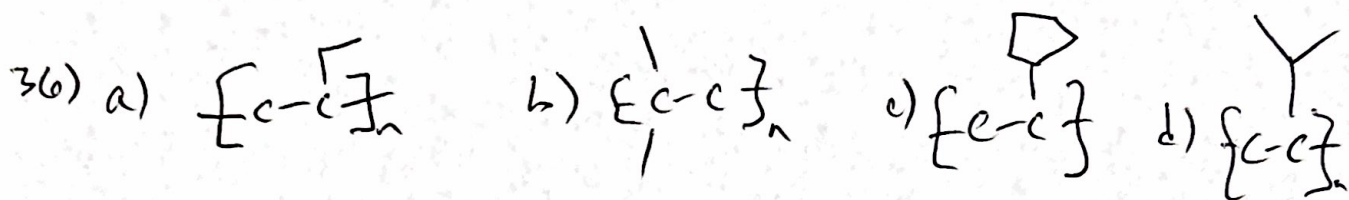
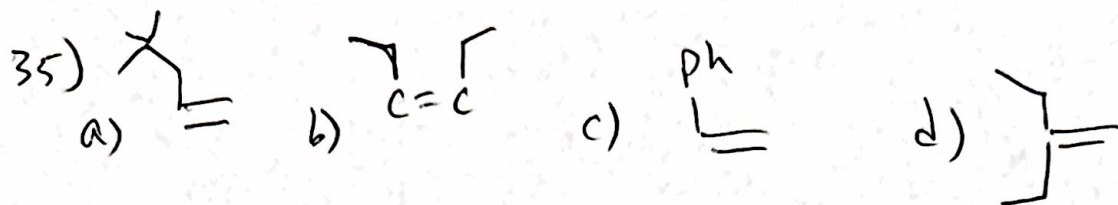
28) The top side of the alkene is too crowded for the H_2/Pt to ATTACK so it can only ATTACK from underneath.



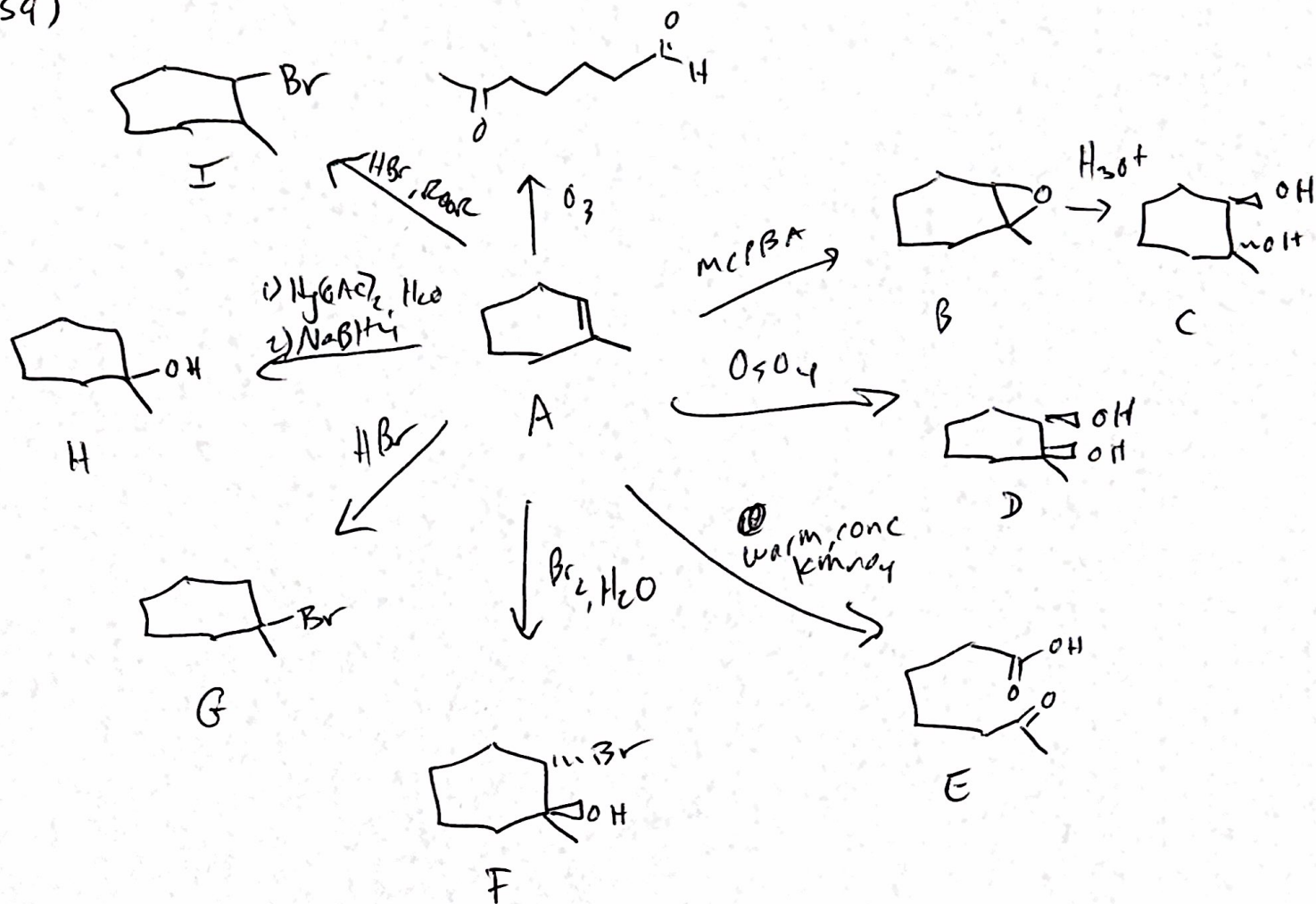


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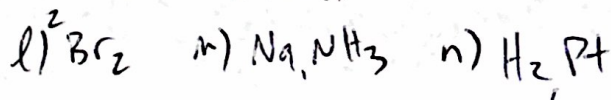
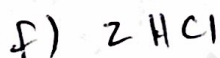
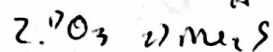
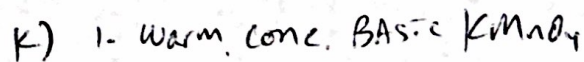
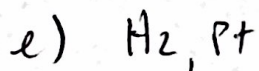
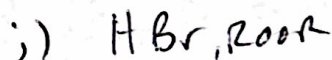
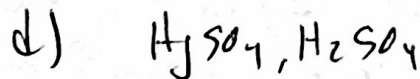
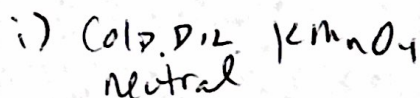
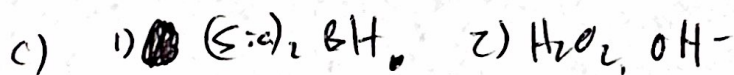
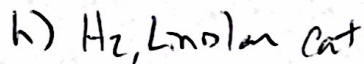
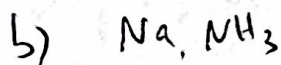
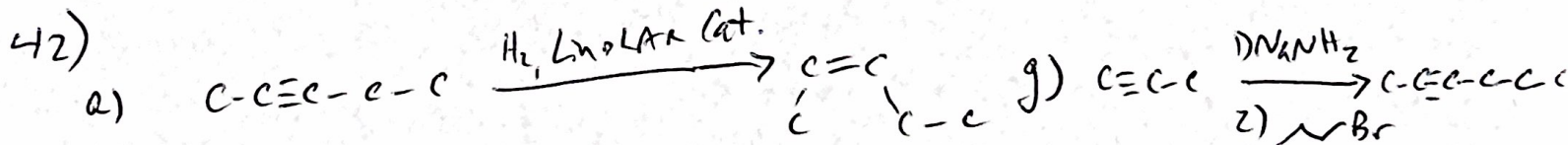
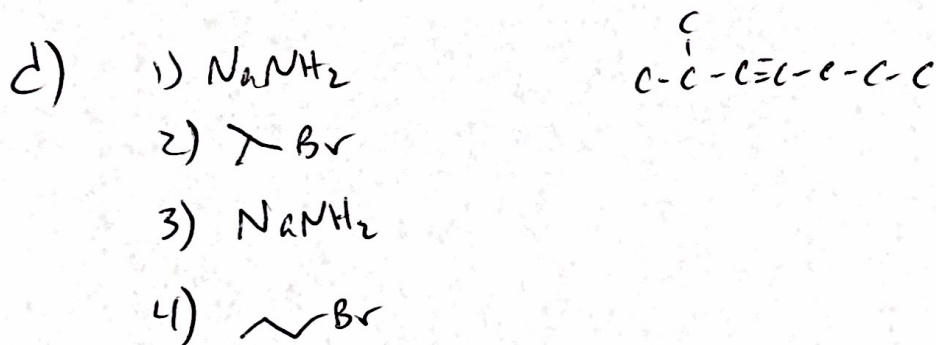
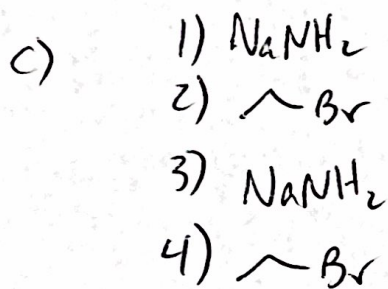
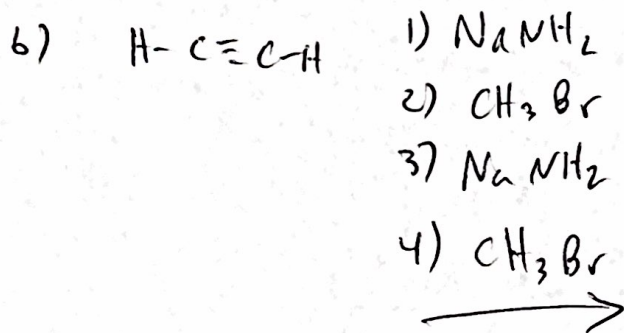
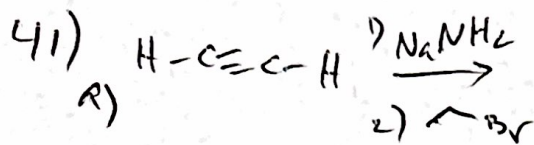




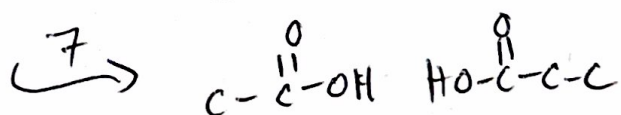
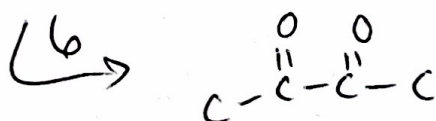
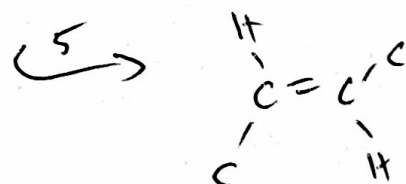
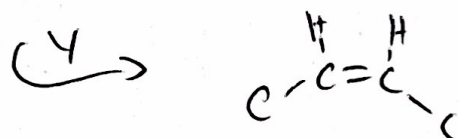
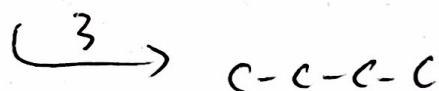
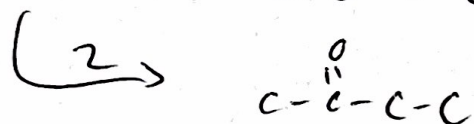
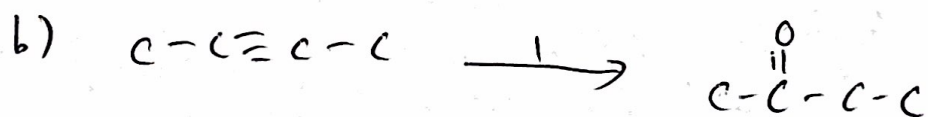
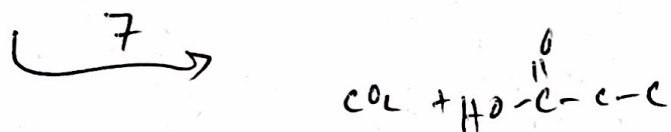
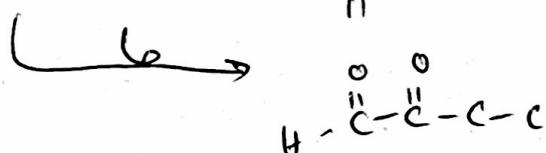
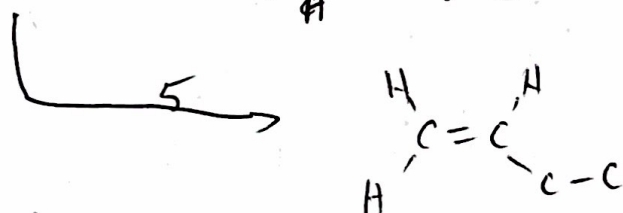
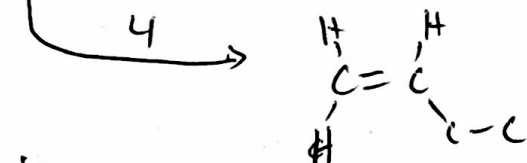
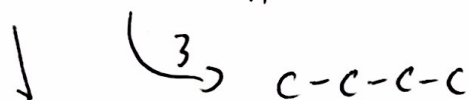
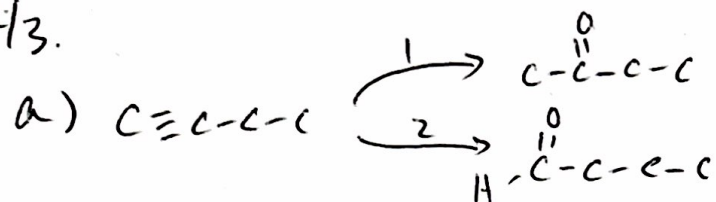
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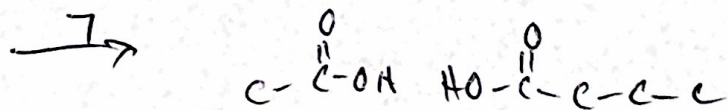
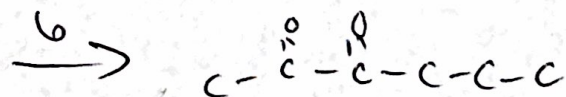
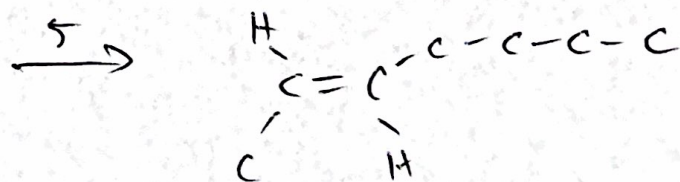
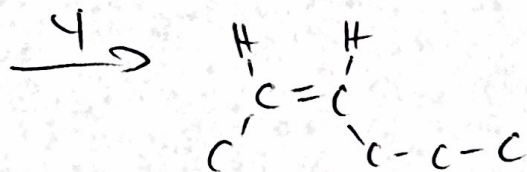
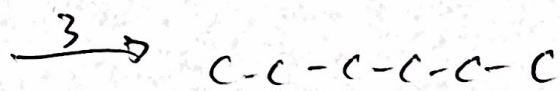
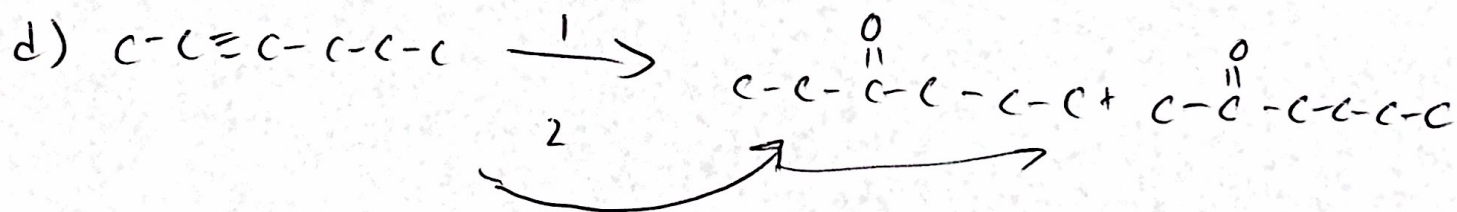
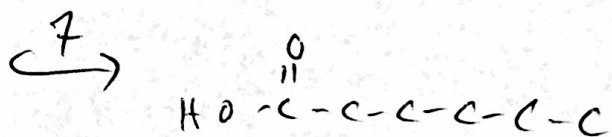
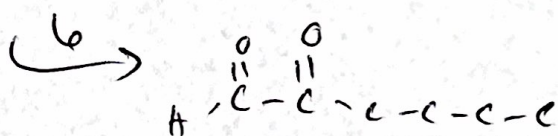
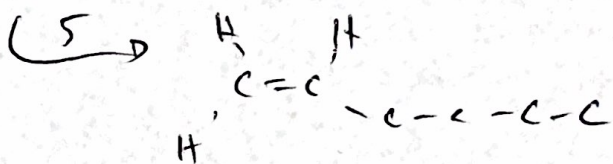
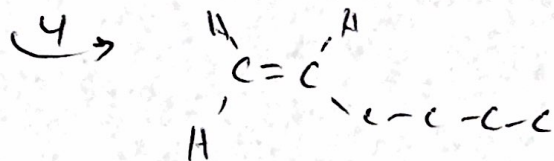
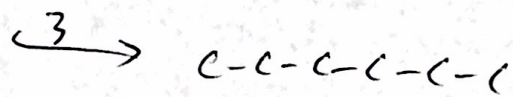
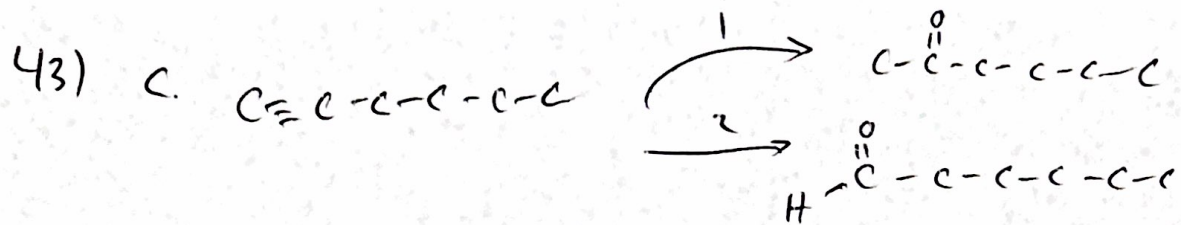


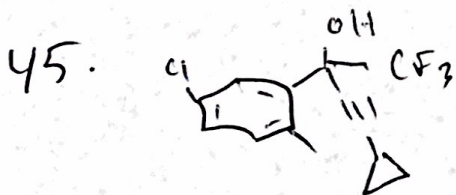
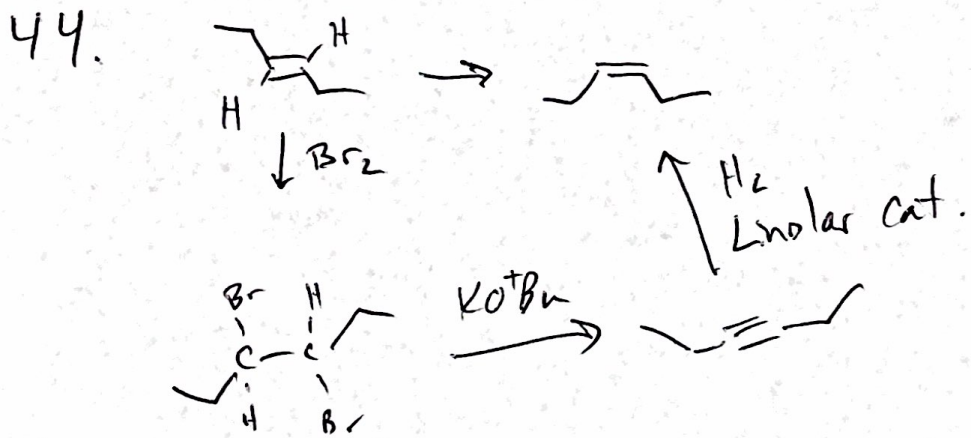
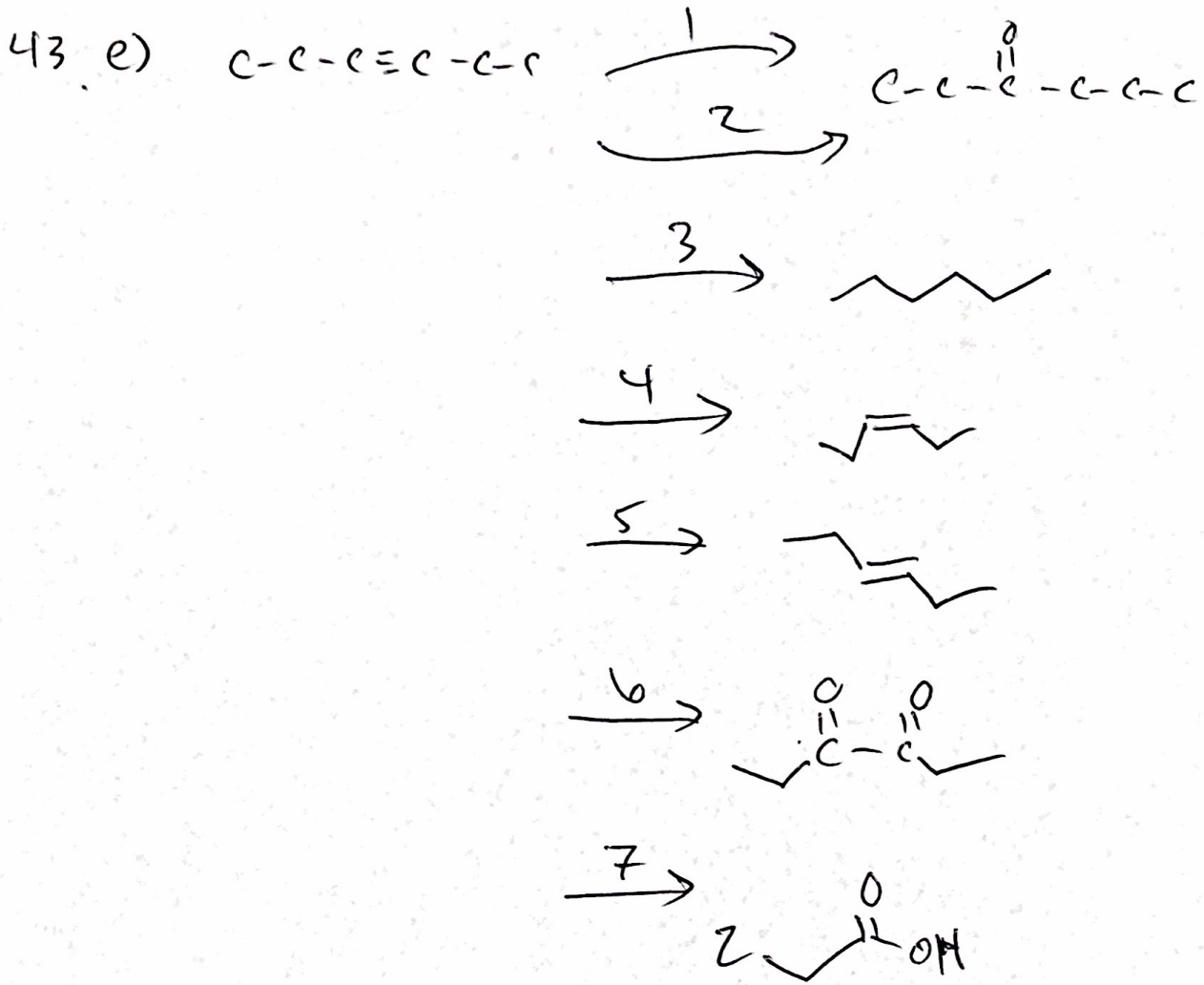
- 40)
- $\text{K}^+\text{O}^-\text{Bu}$ 2. $\text{Br}_2, \text{H}_2\text{O}$
 1. $\text{Br}_2, \text{H}_2\text{O}$ 2. $\text{K}^+\text{O}^-\text{Bu}$ 3. O_3 2) Me_2S
 1. H_2SO_4 2. 1) $\text{H}_2\text{OAc}_2, \text{H}_2\text{O}$
2) NaBH_4
 1. $\text{K}^+\text{O}^-\text{Bu}$ 2. $\text{OsO}_4, \text{H}_2\text{O}_2$
 1. H_2SO_4 2. mCPBA
 1. $\text{K}^+\text{O}^-\text{Bu}$ 2. CH_2N_2

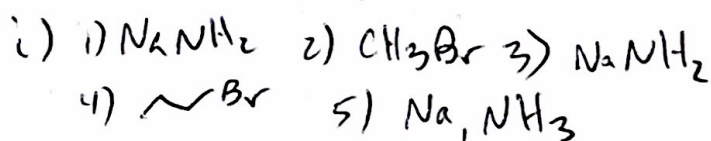
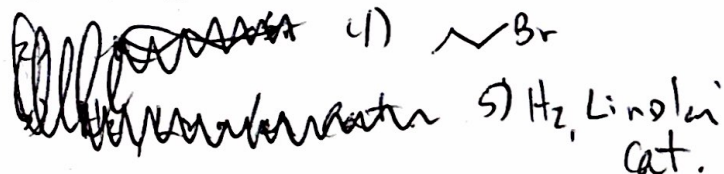
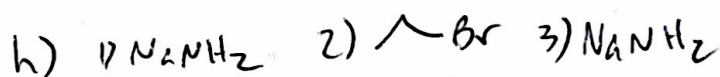
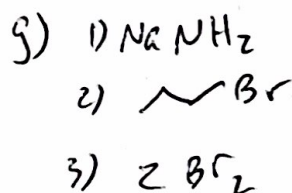
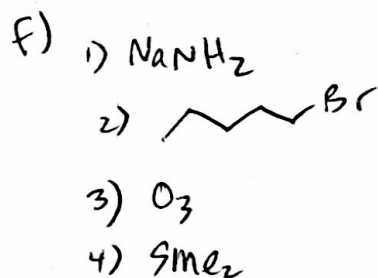
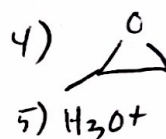
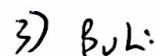
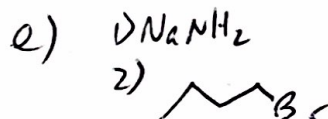
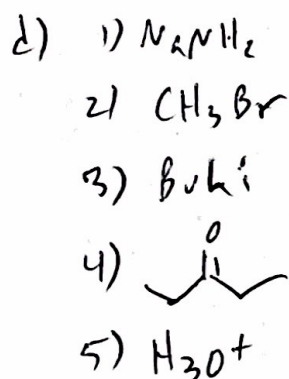
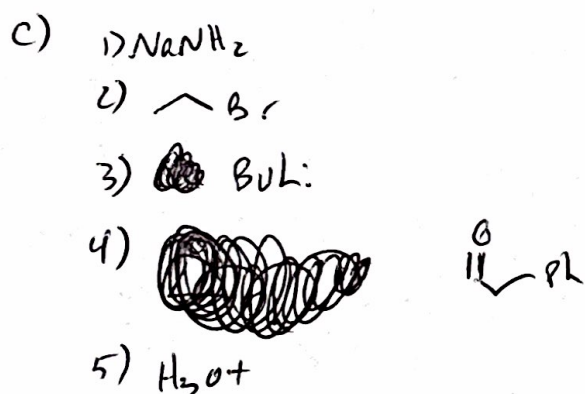
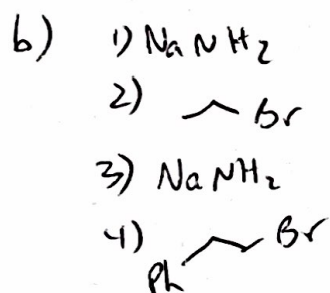
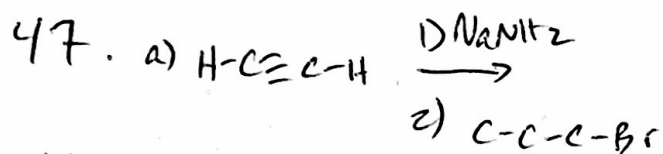
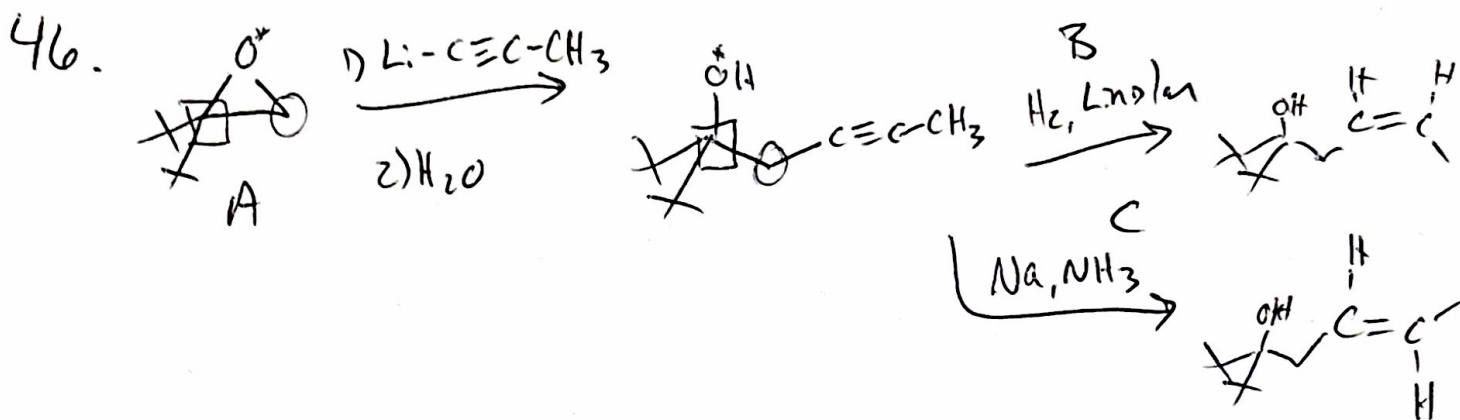


4/3.









47. j) 1) NaNH_2 2) MeBr 3) NaNH_2 4) MeBr ~~5) H_2 , Lindlar cat.~~
 5) H_2 , Lindlar cat. 6) OsO_4 , H_2O_2

k) 1) NaNH_2 2) MeBr 3) NaNH_2 4) MeBr 5) H_2 , Lindlar cat.
 6) CH_2N_2

l) 1) NaNH_2 2) CH_3Br 3) NaNH_2 4) CH_3Br 5) KMnO_4 , cold, dil. aq.

m) 1) NaNH_2 2) CH_3Br 3) HgSO_4 , H_2SO_4

n) 1) NaNH_2 2) CH_3Br 3) BuLi 4) $\text{H}-\text{C}(=\text{O})-\text{H}$ 5) H_3O^+

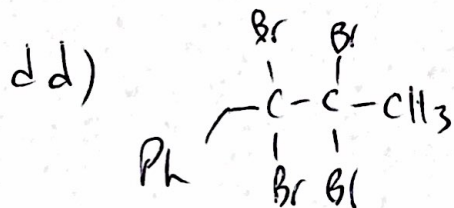
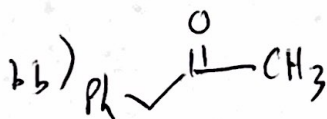
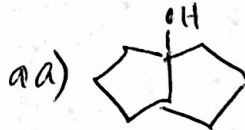
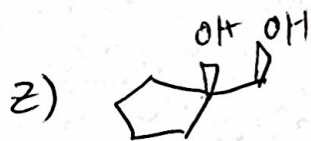
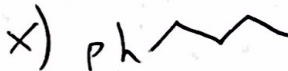
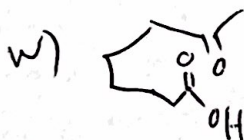
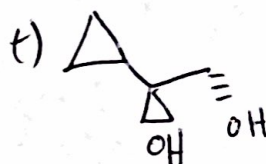
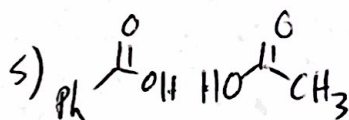
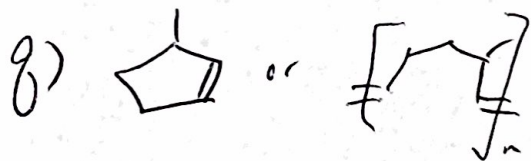
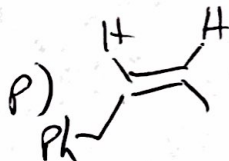
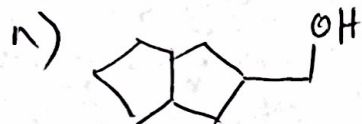
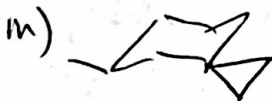
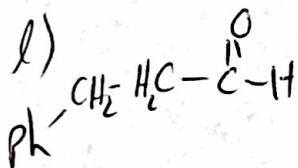
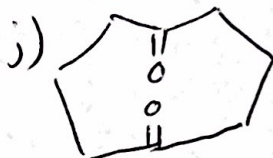
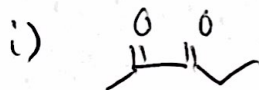
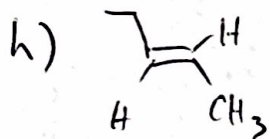
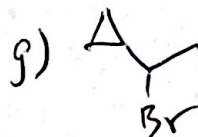
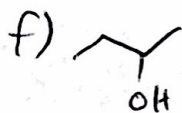
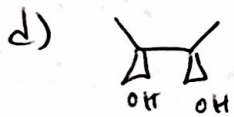
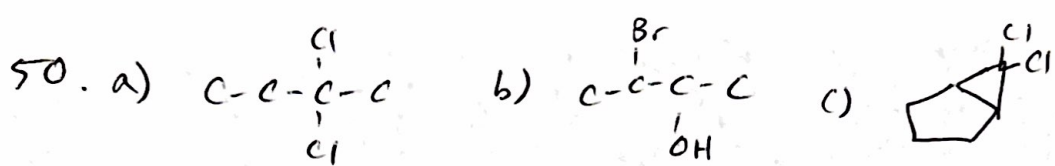
o) 1) NaNH_2 2) CH_3Br 3) 2HBr

48) a) $\text{C}-\text{C}-\text{C}\equiv\text{C}-\text{CH}_3$

b) $\text{CH}_3-\text{C}\equiv\text{C}-\text{C}-\text{C}-\text{C}-\text{C}\equiv\text{C}-\text{CH}_3$

49) a) $\text{C}-\overset{\text{Br}}{\underset{|}{\text{C}}}-\text{C} \xrightarrow{\text{KO}^t\text{Bu}} \text{C}=\text{C}-\text{C} \xrightarrow{\text{Br}_2} \text{C}-\overset{\text{Br}}{\underset{\text{Br}}{|}{\text{C}}}-\text{C}$

b) 
 $\text{1-methylcyclopentanol} \xrightarrow{\text{H}_2\text{SO}_4} \text{1-methylcyclopentene} \xrightarrow[\text{Perox.}]{\text{HBr}} \text{2-bromo-1-methylcyclopentane}$



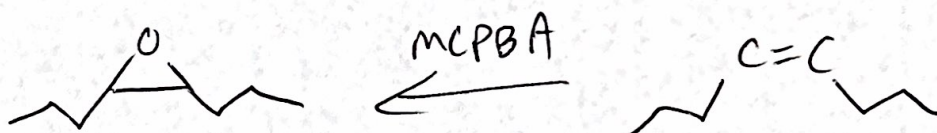
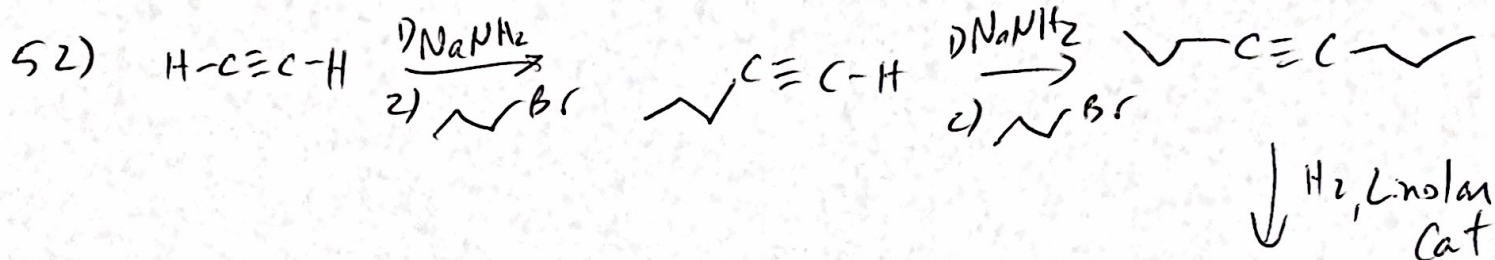
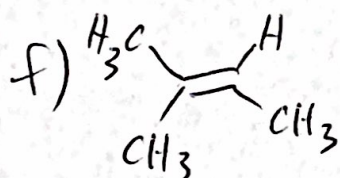
51) a) 1) O_3
2) me_{2s}

b) 1) DSa_2BH
c) H_2O_2, OH^-

c) BF_3, H_2O

d) $HBr, ROOR$

e) 1) $Hg(OAc)_2, EtOH$
2) $NaBH_4$



53) a) 1 degree of unsat. $\frac{2C+2-H}{2} = \frac{2(6)+2-12}{2} = 1$ 11 H, 1 Cl
 b) NO $C=C$. The degree of unsat. must be a ring.

c) $\frac{2(6)+2-10}{2} = 2$ degrees of unsat.

d) Now we have a $C=C$.

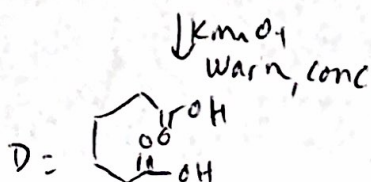


e) 1 type

f) 1 degree of unsat. (The ring)

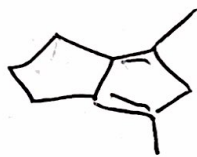


g) A = B = C =



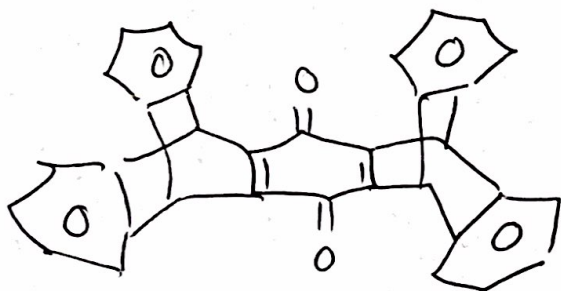
54. a) E contains a π bond

b)

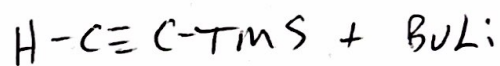


55.

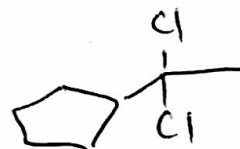
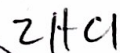
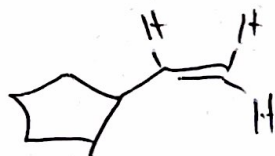
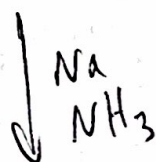
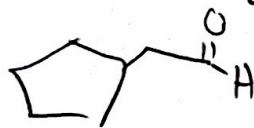
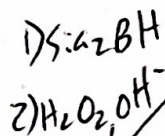
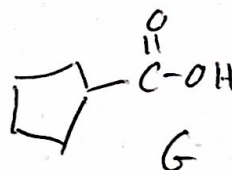
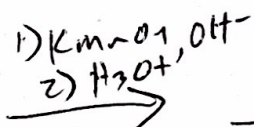
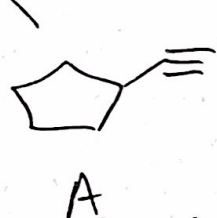
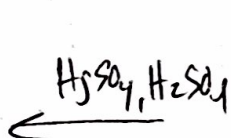
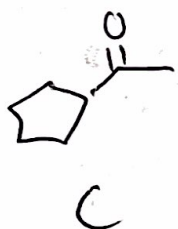
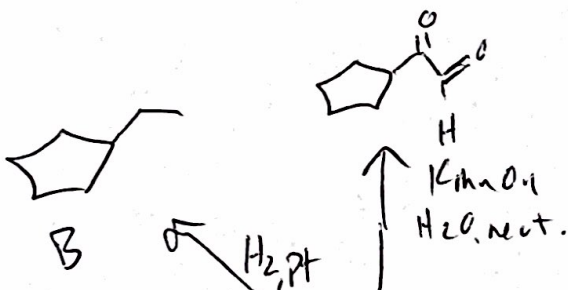
a)



b)



56.



F