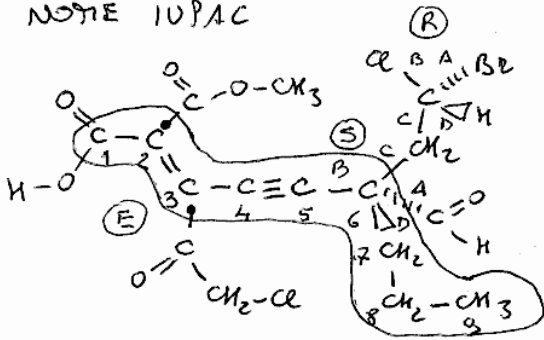


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1) NOME IUPAC

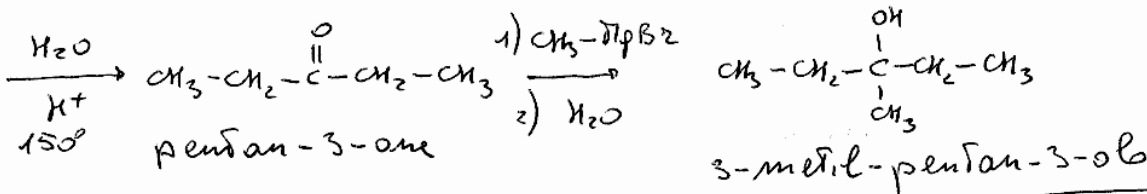
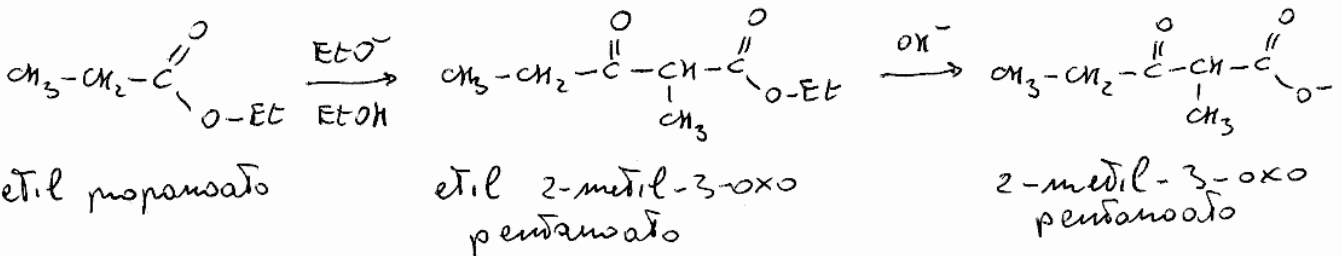
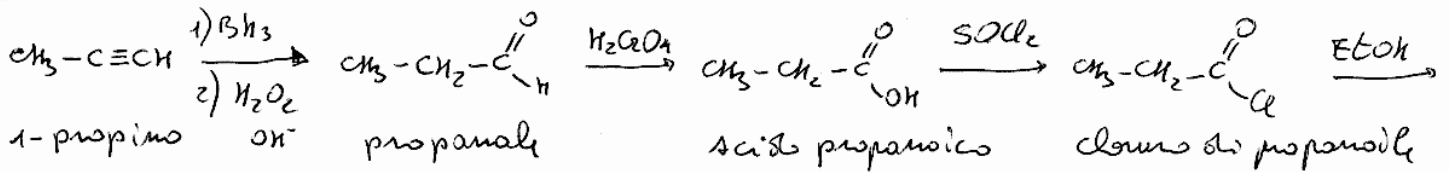
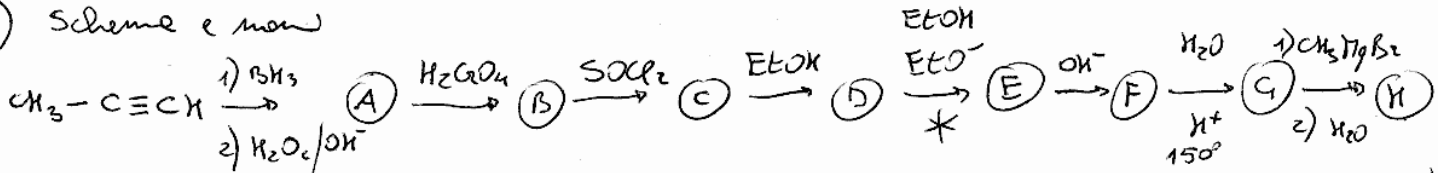


NONANO  
 ACIDO NONANOICO  
 ACIDO NON-2-EN-4-INOICO  
 (2E, 6S)  
 2-(metoxycarbonyl)  
 3-(2-cloroetil)

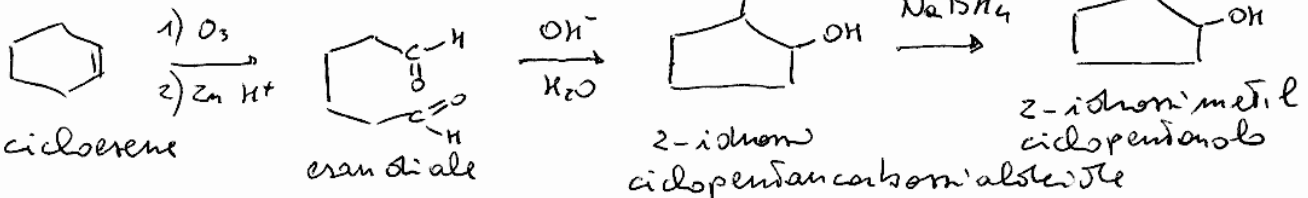
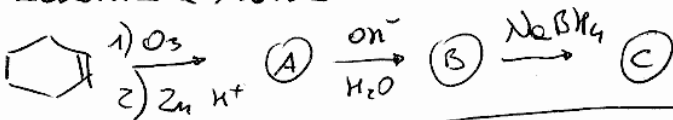
6-formil  $6-[ (2R)-2-bromo-2-cloroetil ]$

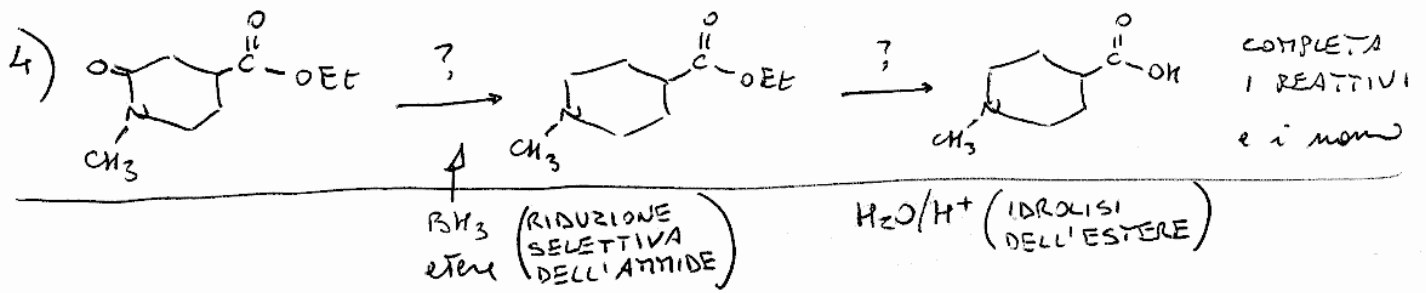
Acido (2E, 6S)-6-[ (2R)-2-bromo-2-cloroetil ]-3-(2-cloroetil)-6-formil  
 2-(metoxycarbonyl)-NON-2-en-4-inoico

2) Scheme e nomi



3) Scheme e nomi



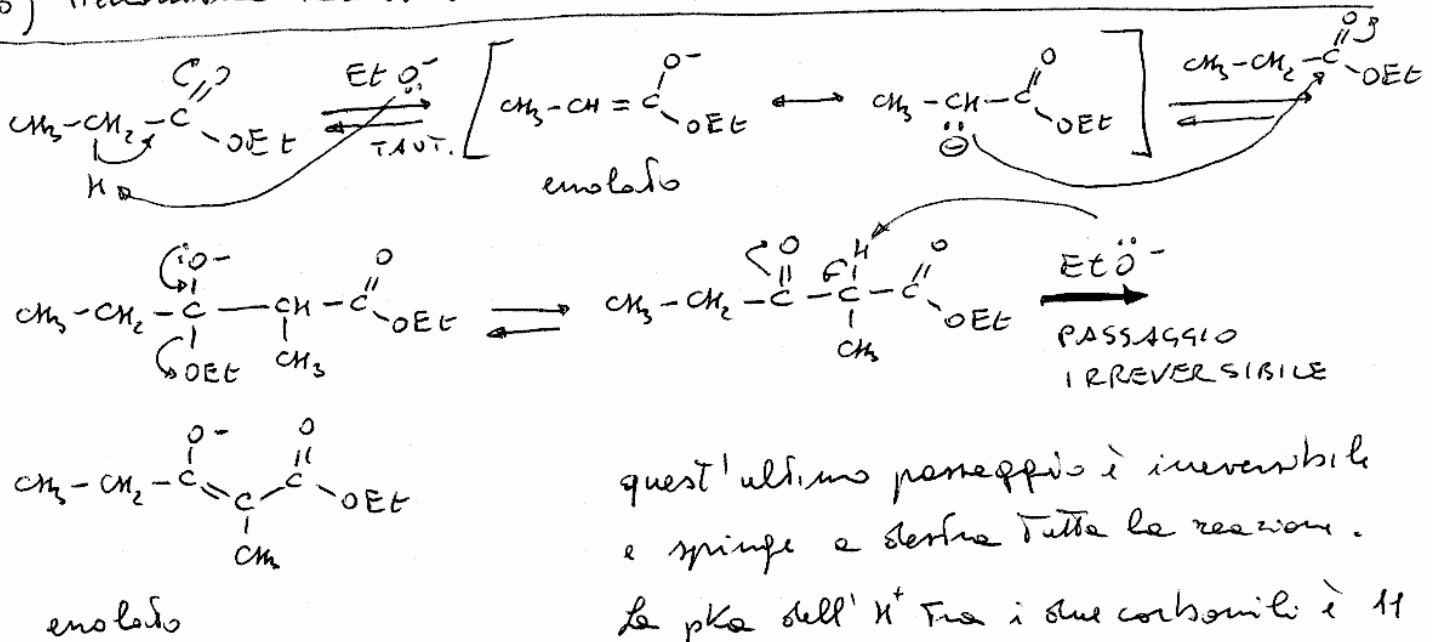


1-metil-2-oxo-piperidin-4-carbomilato di etile

1-metil-piperidin-4-carbomilato di etile

Acido 1-metil-piperidin-4-carbomilico

5) Meccanismo reazione con  $\text{EtO}^-$  \* nell'esercizio 2



quest'ultimo passaggio è irreversibile e spinge a destra tutta la reazione. La pKa dell' $\text{H}^+$  tra i due carbomili è 11 quindi questo  $\text{H}^+$  viene frappato al 100% da  $\text{EtO}^-$  che ha  $\text{pKa} \approx 16$