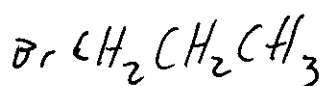
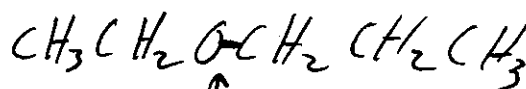
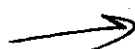
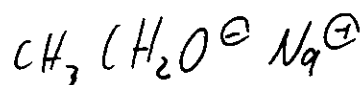


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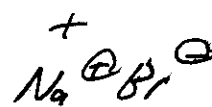
Ch 9 Substitution & Elimination Rxns of alkyl halide

Subst. Rxn Ex:



• Also, called nucleophilic displacement

new bond

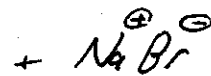
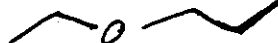


Mechanism

nucleophile



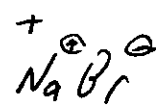
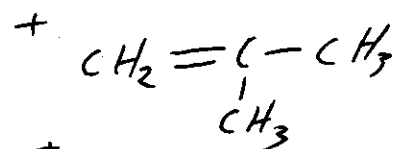
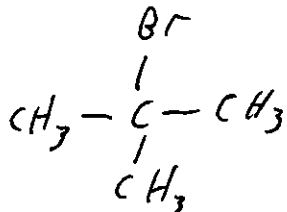
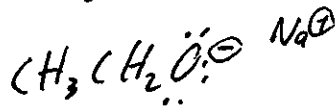
electrophile



leaving group

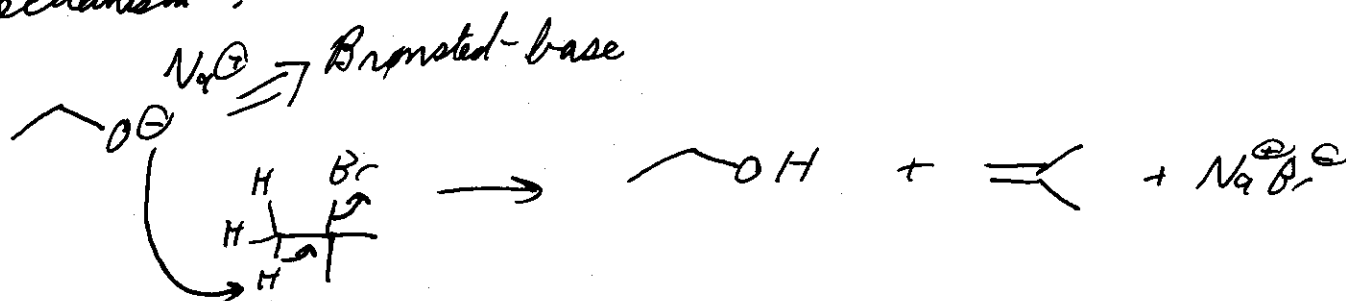
Term: see Ch 3

Ex of elimination



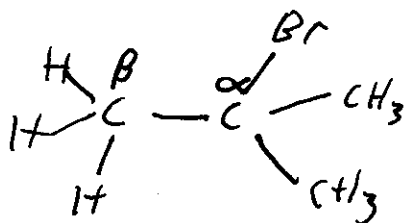
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Mechanism:



Different roles possible (Bronsted base in elim or nucleophile in Nu^+ displacement)

Note: "B-elimination" vs "elimination" pp 378-380
to indicate relative position within molecules



key idea: Greek letters

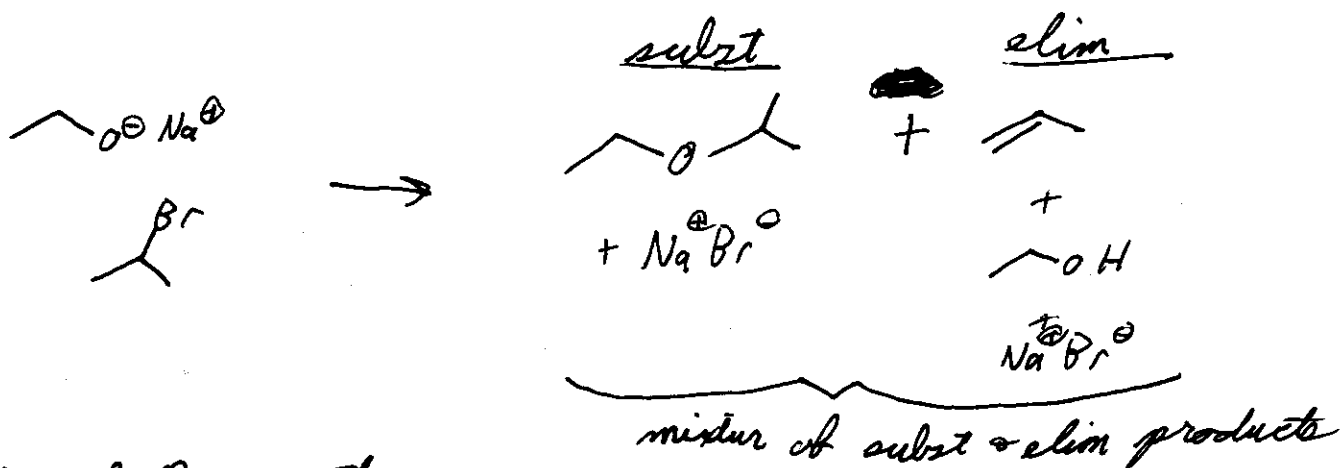
α, β, γ

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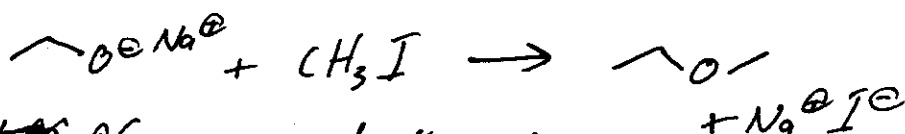
Compare / contrast

• 1° vs 3° → elimination
↓
substitution

Both processes can occur in some cases (like 2°)



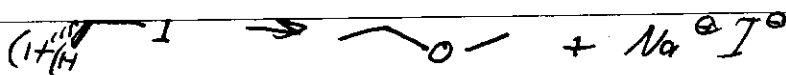
Use of Rxn rates to learn about mechanism
⇒ closer look at substitution



~~Var~~ Vary concentrations of reactants → rate law

rxn rate = $k [\text{CH}_3\text{CH}_2\text{ONa}] [\text{CH}_3\text{I}]$; k = proportionality constant
 based on this rate law... rxn ~~rate~~ overall is "second order"
 (2 molecules used)

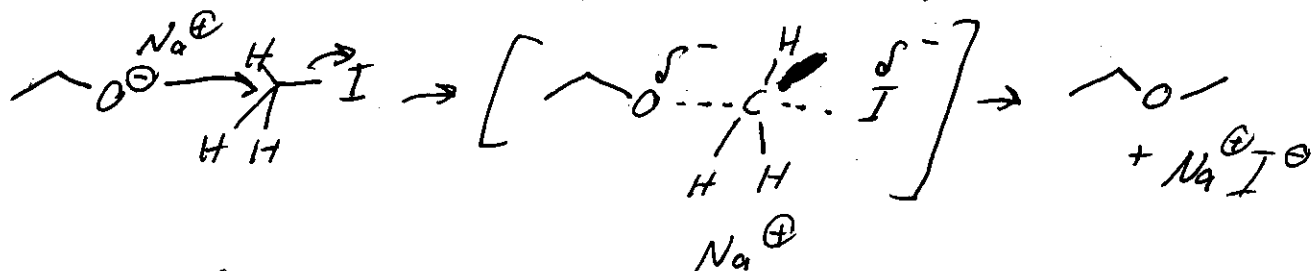
- rxn is first order in $\text{CH}_3\text{CH}_2\text{ONa}$ & first order in CH_3I (2nd order)
- This type of rxn refers to S_N^2 (Substitution Nucleophilic)
- From the rate law, we can deduce that both $\text{CH}_3\text{CH}_2\text{ONa}$ & CH_3I are involved in the rate determining transition state



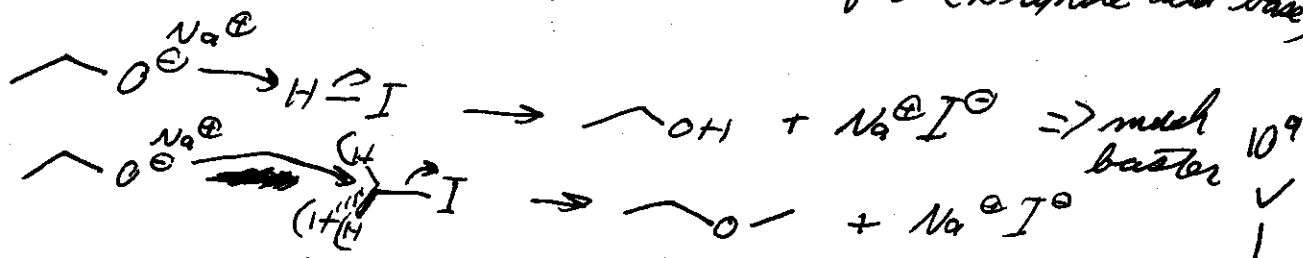
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This rate law is consistent w/ the previous mechanism



Recall (Ch 3): analogy ~~between~~ between displacement at C & proton transfer (Bronsted acid-base)



steric hindrance ("CH" signifies steric interaction)