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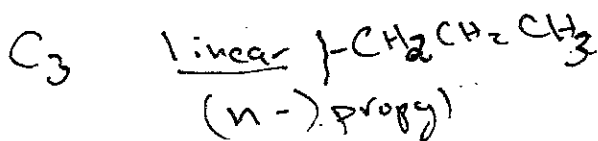
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Last time: - Conformation analysis of alkanes
 - staggered, gauche, and eclipsed conformations (energy values)
 - local and global energy minima + maxima
 - nomenclature

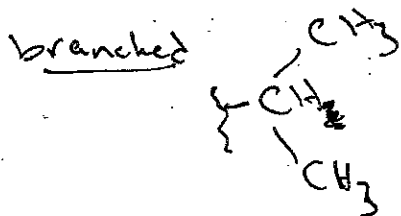
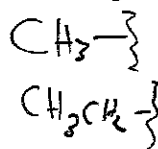
Branched alkyl substituents

→ know the names

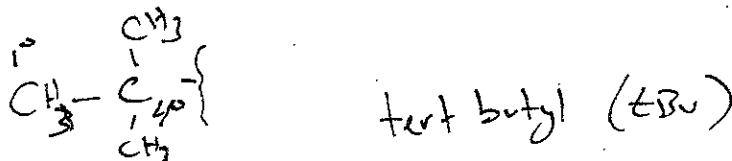
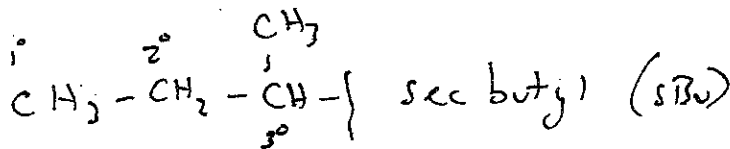
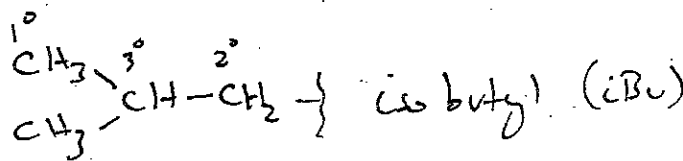
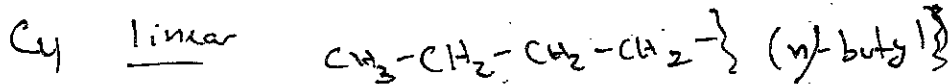
attached to something
↓



C₁ methyl substituent
 C₂ ethyl-substituent



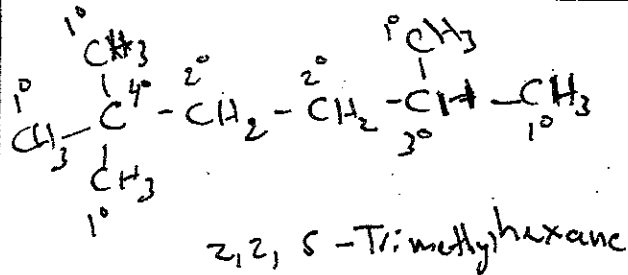
isopropyl (iPr)



Differently substituted carbons → 1°, 2°, 3°, 4° carbons
 attached to ↓
 1 other C ↓ 2 C's ↓ 3 C's ↓ 4 C's

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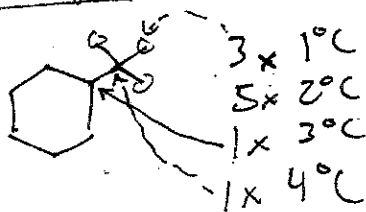


Different substituted sp^3 hybridized

C can be formed, differ by # of C's attached

- Primary	/	Secondary	/	Tertiary	/	Quaternary
↓		↓		↓		↓
1 other carbon attached		2 "		3 "		4 "
1°		2°		3°		4°

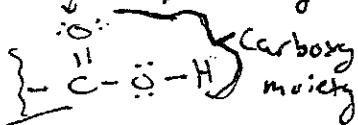
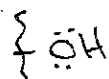
Other example



Functional Groups

- remember names for nomenclature

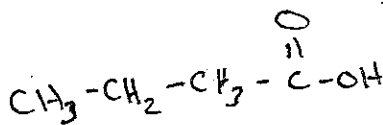
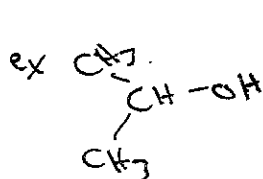
Commonly bonded atoms that have similar chemical properties



Alcohol (R=alkyl chain)

Carboxylic acid
→ sp^2 hybridized

hydroxyl group



2-propanol (isopropanol)

n-butanoic acid

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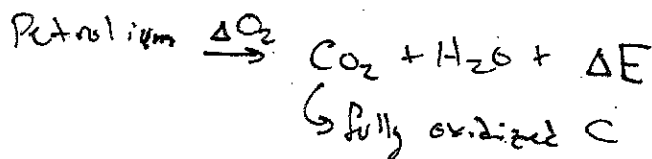
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Oxidation/Reduction

oxidation: ↑ O that are bound to carbon

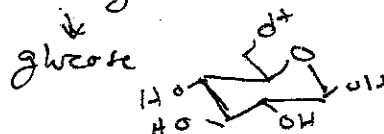
reduction: ↑ H " " "

Petroleum oxidation (alkanes, aromatics)



CH_4 → fully reduced

Reduction: plants $CO_2 \xrightarrow{h\nu}$ Carbohydrates (photosynthesis)

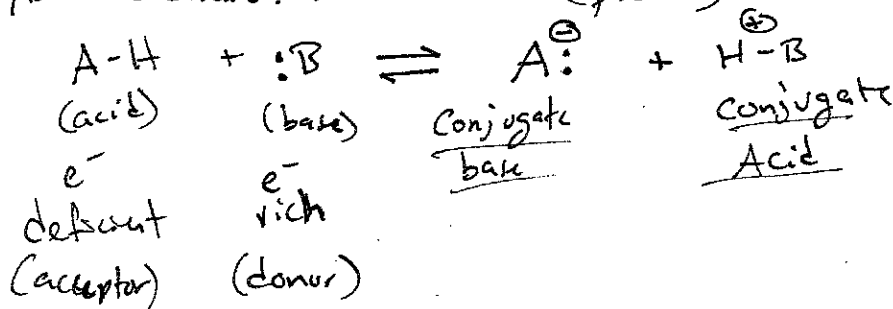


Animals: Glucose $\xrightarrow{O_2}$ $CO_2 + Energy$ (respiration) → oxidative process
 ↳ stored as **Fat** or **ATP**

CH. 3 - Acids + Bases

Problems (3.X) x = 1-10, 15, 21, 24, 27, 39, 42 a-c, 43, 45, 49, 50

A/B reactions: H^+ transfer (proton)



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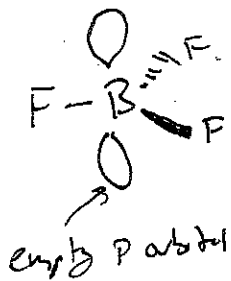
Three definitions:

~~Arrhenius~~ Arrhenius: A forms H^+
B forms OH^- } happens in an aqueous system

Bronsted-Lowry: A proton donor
B proton acceptor

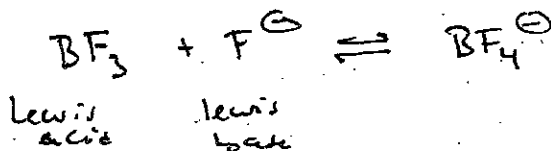
Lewis: A - e^- deficient system (e^- acceptor)
B - e^- rich system (e^- donor)

Example: Boron trifluoride



B - 6 valence electrons (wants 8) → can accept 2 more in the empty p orbital

Lewis acid



Use of arrows to indicate e^- movement

- start @ the e^- donor (or e^- in the bond)
- end @ the e^- acceptor

