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Recall:

- Bonding in organic molecules
- Polar bonds & polarity of molecules
- Drawing conventions

Chapter 2 - Alkanes

*Rec. Problems: 1, 3-12, 14, 15, 23, 24, 26-28, 30, 31, 33, 41, 44-48

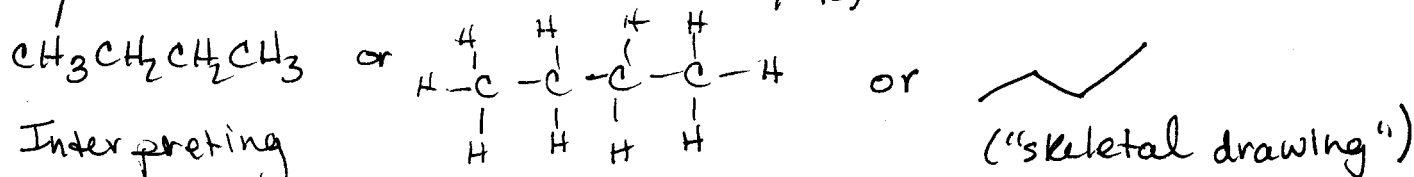
- alkane: molecule comprised of C+H, contains only single bonds

MEMORIZE Names of linear alkanes, $C_1 - C_{10}$ (Table 2.1)

CH_4 = Methane

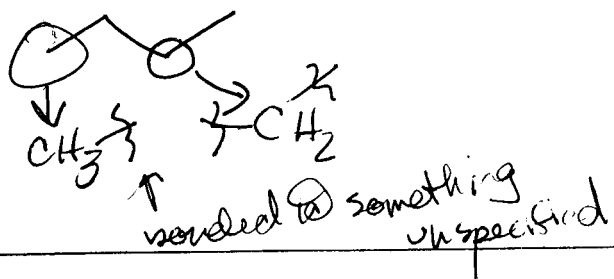
CH_3CH_3 = Ethane ... etc

Graphical Conventions: Butane (C_4H_{10})



Interpreting
skeletal drawings:

- 1) End of every line (including where 2 lines meet) = C
- 2) Every unspecified C valence is H



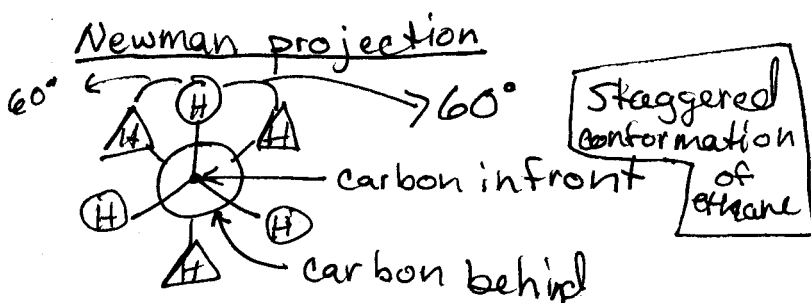
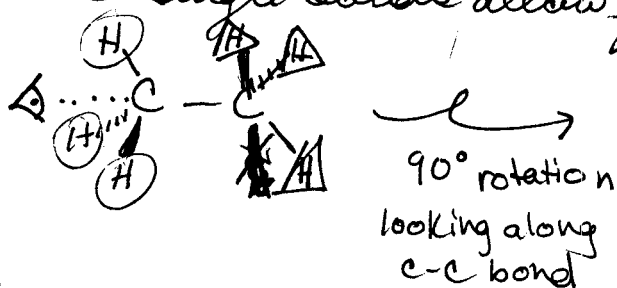
• Don't feel like you always need to use skeletal structures yet

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• These drawings only indicate bond connectivity, not 3D shape!

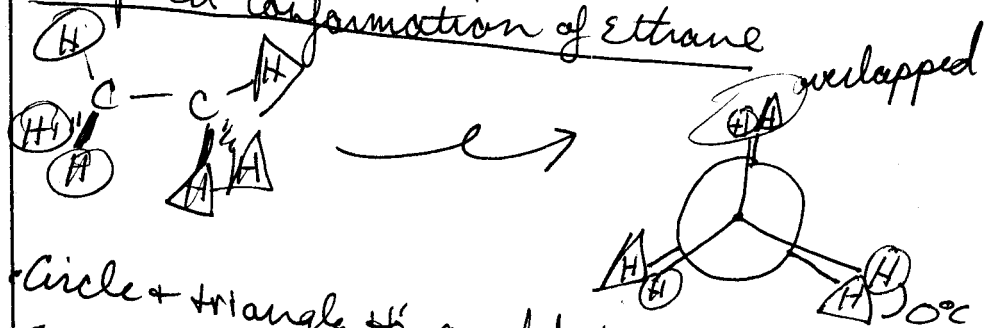
Alkane Shapes or Conformations

• C-C single bonds allow free rotation



• Any 2 H's connected to the same C are 120° apart in the Newman projection
 — not the actual bond angle, just because of perspective

Eclipsed Conformation of Ethane

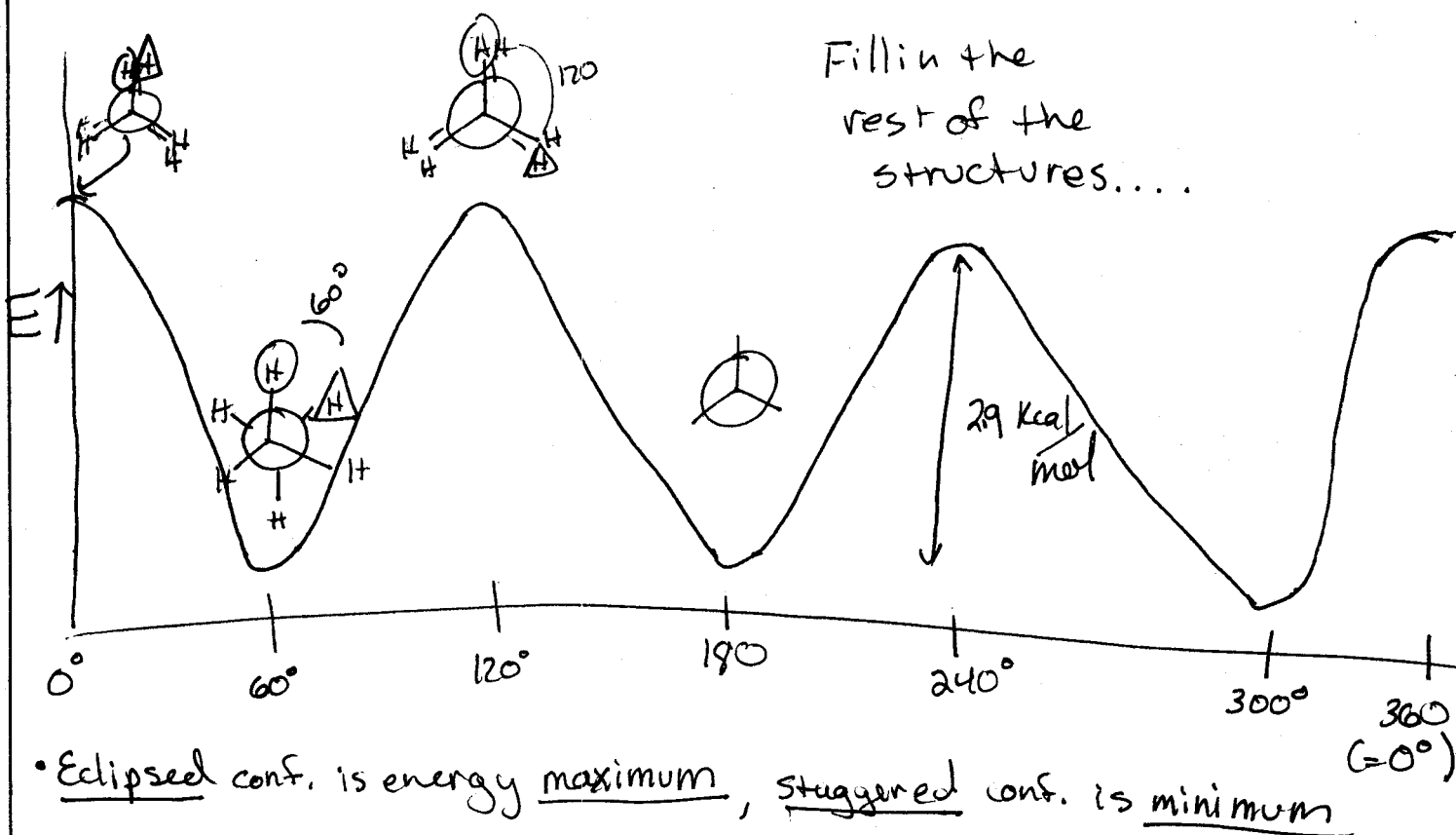


• Circle + triangle H's could have any degree of separation 0° → 60°
 • staggered + eclipsed are max/min cases
 • These names apply to any alkane

Energy Diagram for Rotation about the C-C bond

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Energy Diagram



Note:

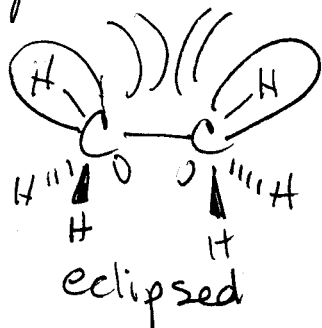
- 1) Eclipsed energy maximum (pl=maximum) ...?
 - 2) staggered E minimum (pl=minimum) ...?
- Tells us eclipsed is least stable conf, staggered is most stable
 - Difference between eclipsed + staggered is $\sim 2.9 \text{ kcal/mol}$ ($\sim 12 \text{ kJ/mol}$)
 → Very small difference in energy

Course CH 343 Sec 003 Lecturer Sam Hellman
Day Monday 1:20 Date 9/12/11
Notes Taken By Ellicia Phelps Total # of Pages 4

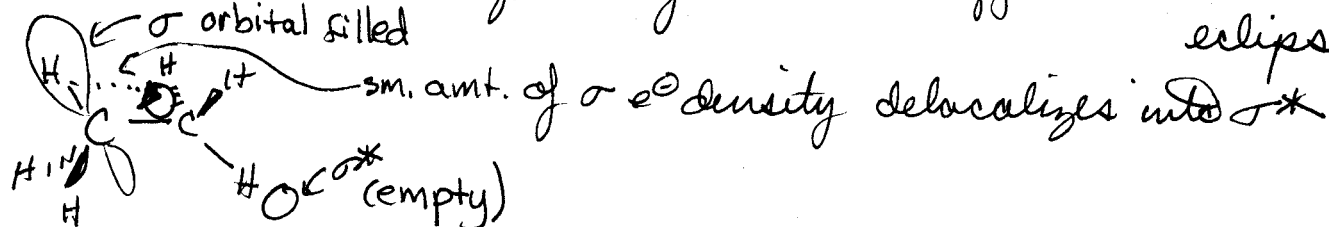
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why is staggered more stable? 2 factors

1) Repulsion between pairs of C-H σ bond e^- s in eclipsed



2) $\sigma \rightarrow \sigma^* e^-$ delocalization (favored in staggered but not in eclipsed)



- Some very sm amt of e^- repulsion in the bond $\rightarrow$ moving a little into σ^* lowers energy by relieving this.