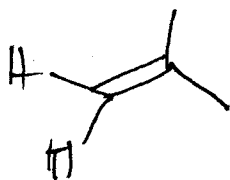
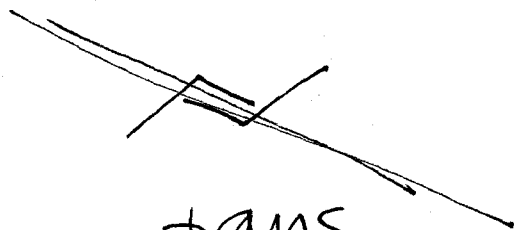


more alkene nomenclature

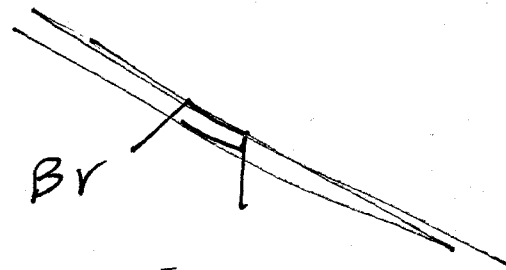
1. cis vs. trans - worry about this when both ends of the double bond have two different substs.



no cis/trans



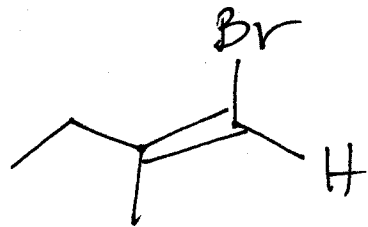
trans
= opposite



cis
= same

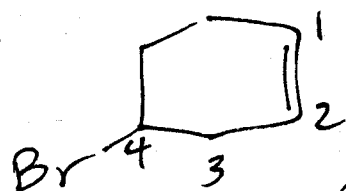
consider the double bond as a fence -
where are the two big groups? (one
from each end of the double bond)

What about:



Hard to say.
(Wait for ch. 7)

2. cyclic alkenes -

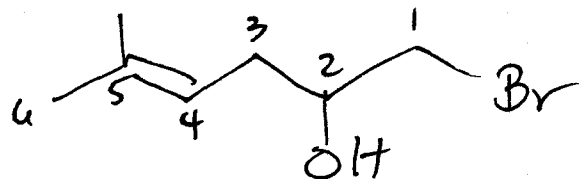


4-bromocyclohexene

(double bond defines #1)

Always is if < 10 C's in ring,
so don't need to specify.

3. What if there's both an OH and an alkene?
OH wins.

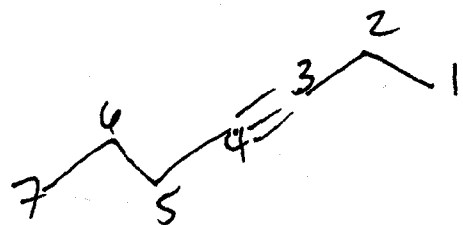


1-bromo-5-methyl-4-hexen-2-ol

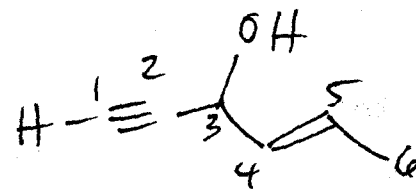
... hex-4-en-2-ol

Alkynes -

longest chain
and becomes yne

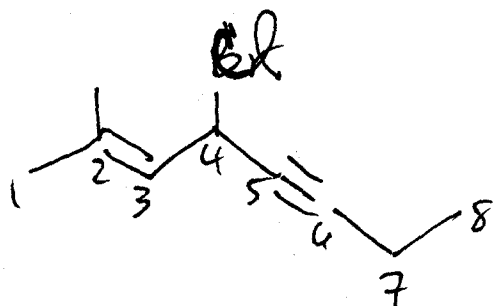


3-heptyne



trans -
hex-4-ene-1-yne-
3-ol

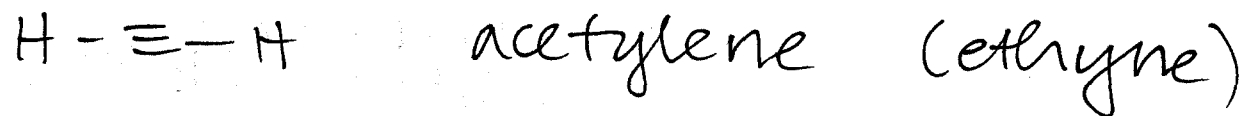
Double and triple - enyne
↑
priority



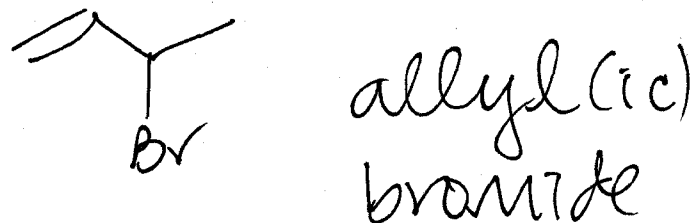
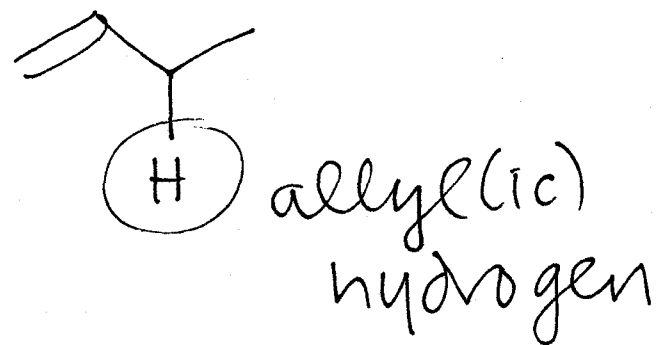
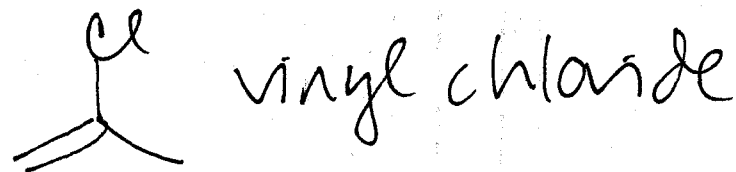
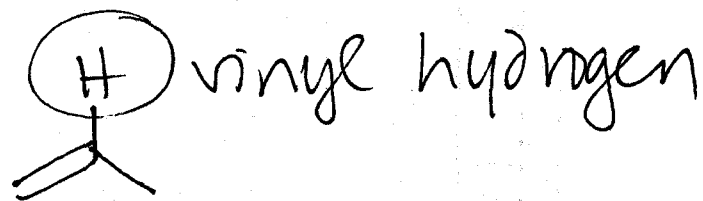
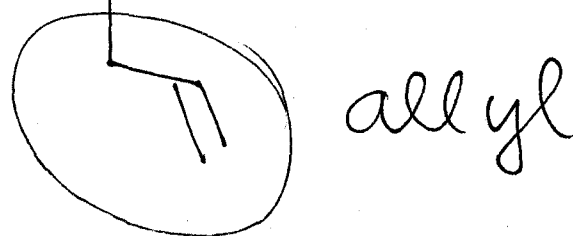
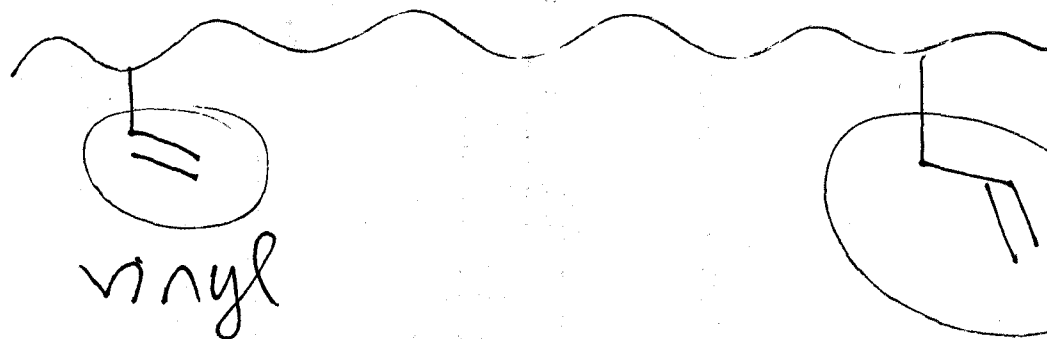
4-chloro-2-methyl-2-octen-5-yne

Add in an OH - now the OH is highest priority
enynol

common names:



common fragments:



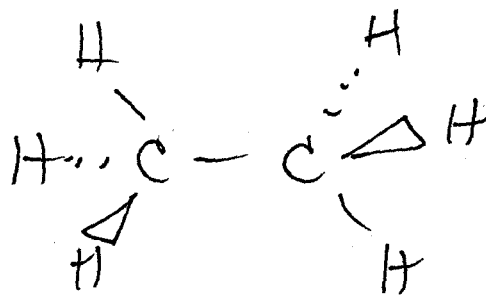
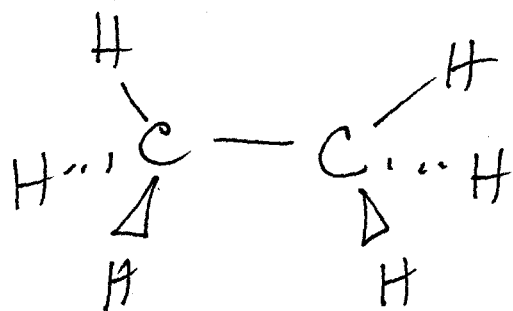
Back when we were discussing bonding models
of ethane, ethene, ethyne

* sigma bonds (axially symmetrical)

* pi bonds (not symmetrical)

- bond line - added in wedges/hatches to
simulate 3D.

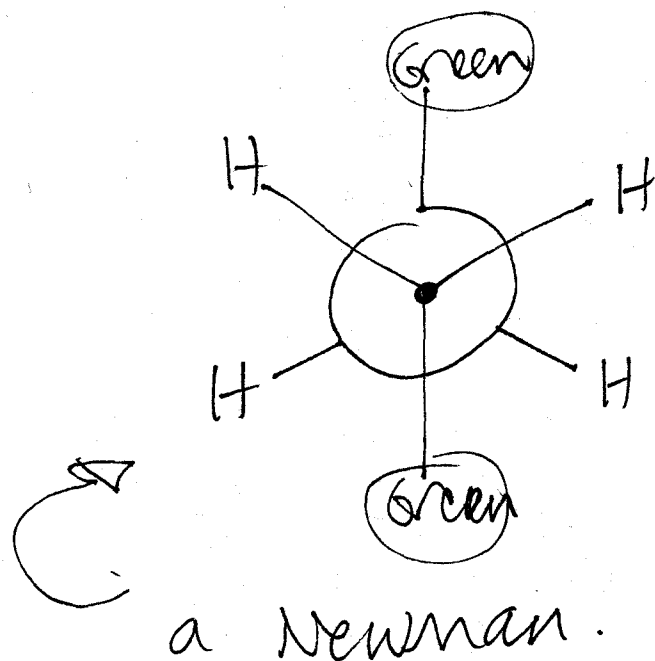
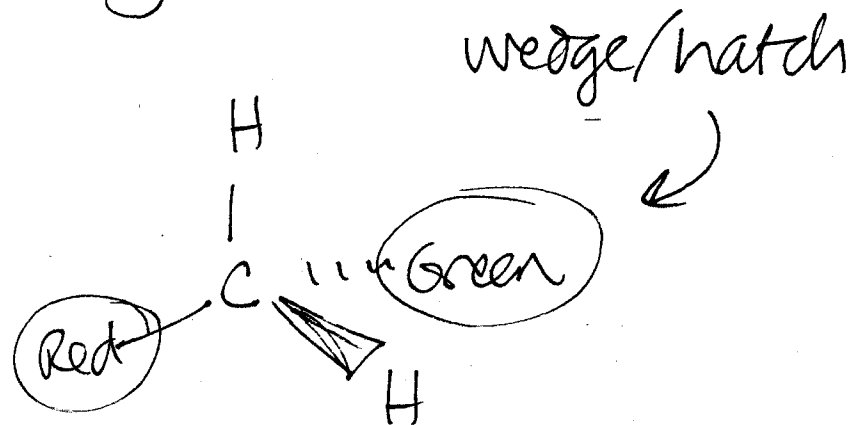
Consider ethane.



two different conformations
of the same molecule.

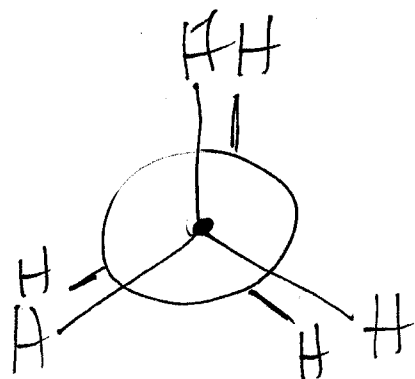
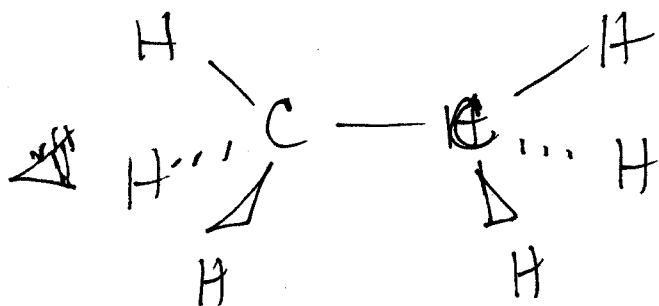
Learn to draw Newman projections.

Rather than a side view of a molecule, we want to sight along a bond.



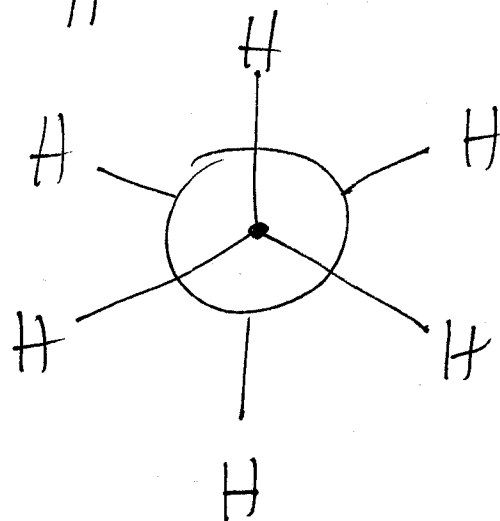
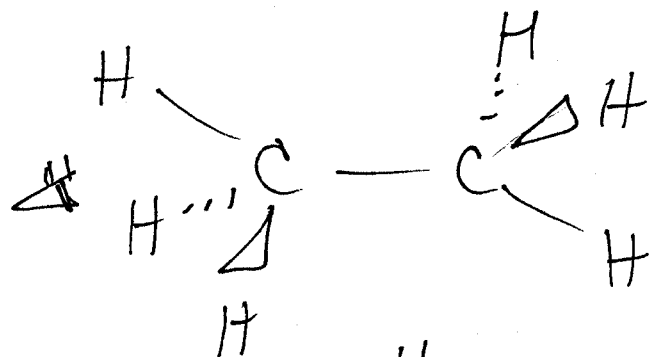
large circle is the atom at the far end of the bond in question.

dot is the atom at the closer end of the bond in question



eclipsed
conformation

high energy



staggered
conformation

low energy