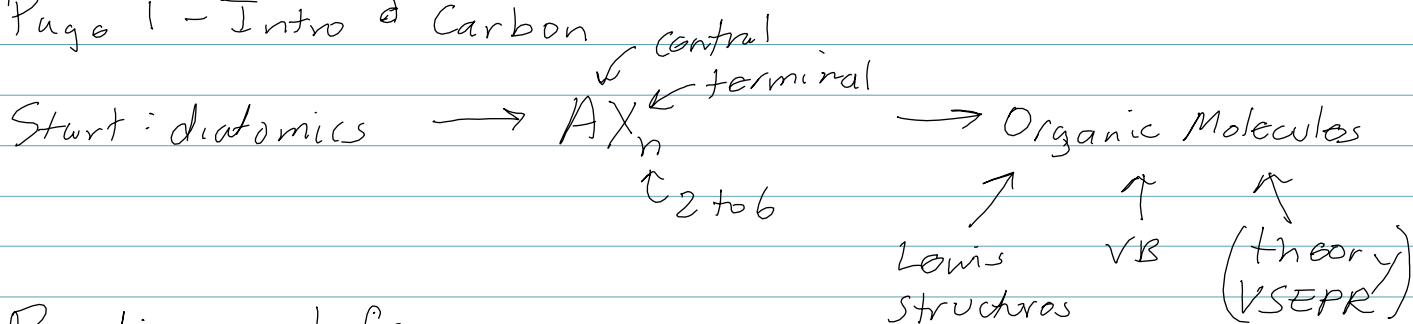


Bonding motifs for common atoms in organic compounds

Page 1 - Intro @ Carbon



Bonding motifs

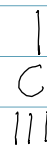
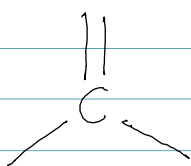
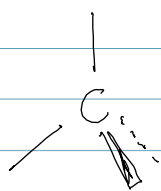
- # of bonds & bond orders
- # of l.p.
- bonding geometry \Rightarrow angles
- formal charge
- hybridization (VB theory)
- octet rule ✓

Carbon - central atom \leftarrow bonded to ≥ 2 atoms

- no formal charge
- no lone pairs
- total 4 bonds ($\sigma + \pi$)

Most common

Rare



tetrahedral
 $\sim 109^\circ$

trigonal planar
 120°

linear
 180°

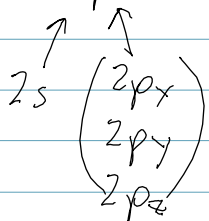
linear
 180°

σ from sp^3

σ from sp^2

σ from sp

σ from sp



p orbital \Rightarrow π bond

2 p orbitals
 make
 π bonds

2 p orbitals
 make
 π bonds
 (one on each end)

Bond motifs in organic compounds

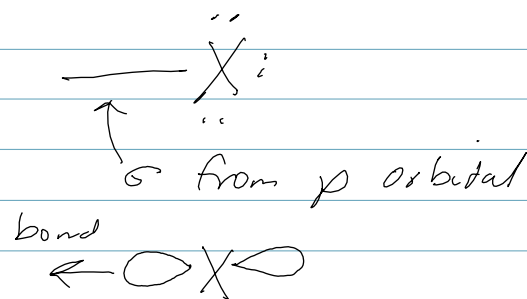
page 2: Hydrogen & Halogens

Inorganic carbon: CO_2 , CN^- , CCl_4

Organic: at least one C-H bond

Hydrogen: always terminal
one bond (σ - 1s orbital from H)
formal charge zero

Halogens: always terminal
F, Cl, Br, I single bond



lone pairs in other
two p orbitals and s

hybridization energy is +

" needed for central atoms \Rightarrow correct angles

" not needed for terminal atoms

Bottom line $X = \text{H, F, Cl, Br, I}$
terminal \swarrow rare exceptions
no formal charges

Bonding motifs in organic compounds

Page 3 - Nitrogen

Mostly formal charge = 0 : single lone pair
three bonds



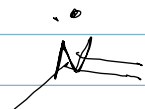
trigonal pyramidal
 $< 109.5^\circ$ (105-108°)

$\approx sp^3$ (25% s / 75% p)

lone pair (34% s / 66% p)

bonds (22% s / 78% p)

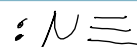
smaller bond angles \leftarrow lessened s character



bent
 $\approx 120^\circ$ ($\sim 118^\circ$)

$\approx sp^2$

p orbital
form π bond



terminal

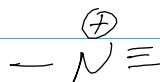
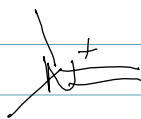
lone pair : 2s

bonds : 2p

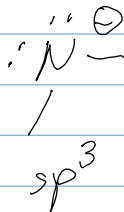
\uparrow
6 + 2 π
bonds

N most likely element to have + formal charge

4 bonds, No lone pairs



Very Rare N^{\ominus}



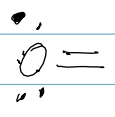
Bonding Motifs in Organic Compounds

Page 4 - Oxygen and Sulfur

Oxygen most common: no f.c.
two bonds
two l.p.

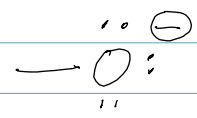


sp³
≈ 105°



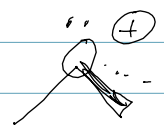
terminal
VB: no hybridization
σ & π bonds from p orbitals
lone pair: remaining p & s

Most likely to be ⊖



σ bond of p orbital on O

Can be ⊕



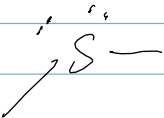
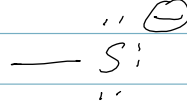
trigonal pyramidal

eg H₃O⁺



bent

Sulfur

most common  or 

Rare: EC exceeds 8 in organics

Later: functional groups