

2 ve⁻

Ba

Ba⁺²

cation
(+) ion

give away
2 ve⁻ becoming
+2 cation

7 ve⁻



I⁻¹

b/c easier to take
1 e⁻ = anion

Covalent

Sharing of e⁻ gets
no charge designation

1 ve⁻

K

+1

b/c easier
to give 1 e⁻ =
cation

transferring (give or take) of
e⁻ gets a charge & is
called a cation or anion

Ionic

H₂

each stick/line
= 2e⁻



single
(sharing 2 e⁻... 1e⁻
from each H)

Linear Diatomic

180°

F₂



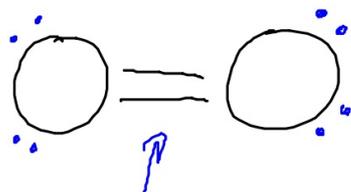
single bond

each F has 3 non-bonding
sites & 1 bonding site

Must always Draw the
NBPs

Linear Diatomic

180°



Double Bond

* 2 NBP on each

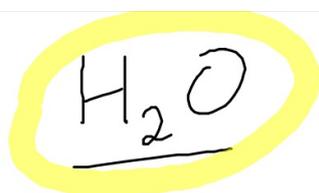
Linear Diatomic
180°



Triple Bond

* 1 NBP on each N

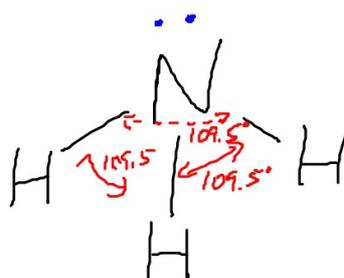
Linear Diatomic
180°



Central Atom = O

2 NBP on O
2 Single Bonds

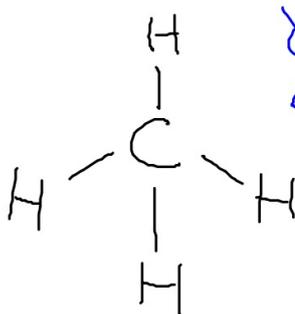
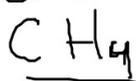
Bent
120°



Central atom = N

1 NBP on N
3 Single Bonds

Pyramidal
109.5°



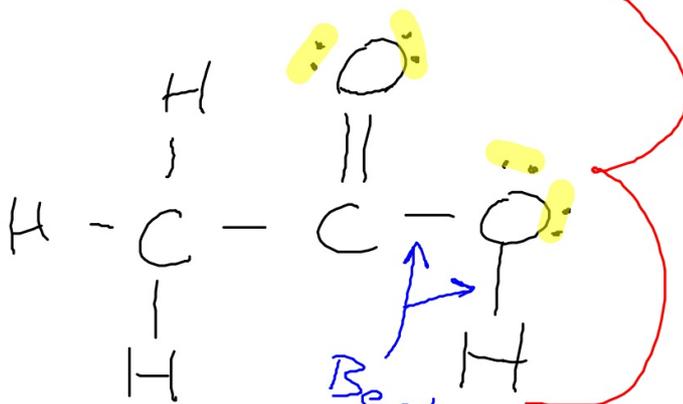
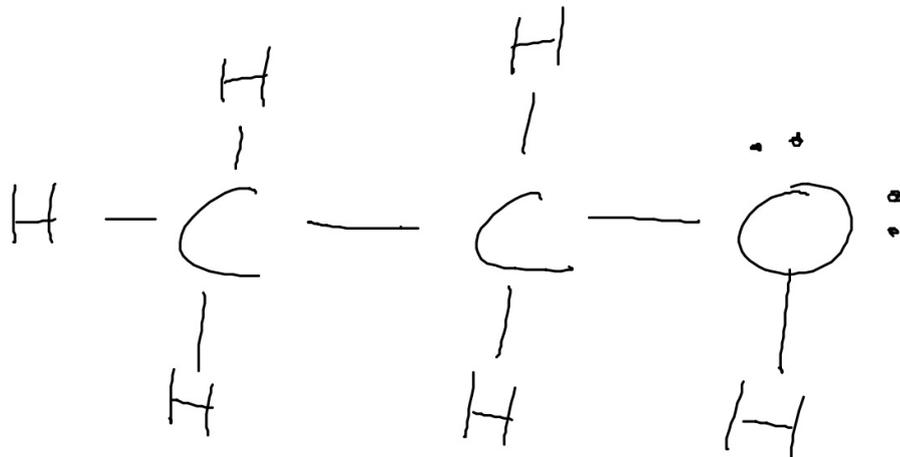
Central atom = C
4 NBP

4 single bonds

Tetrahedral
109.5°



* 2 + Carbons = C-backbone structure



Bent structure when O has 2 single bonds



C is central atom b/c it has the most available bonding sites

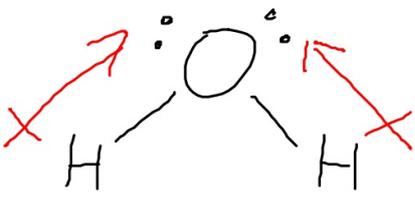


arrow points to atom w/ higher e⁻N

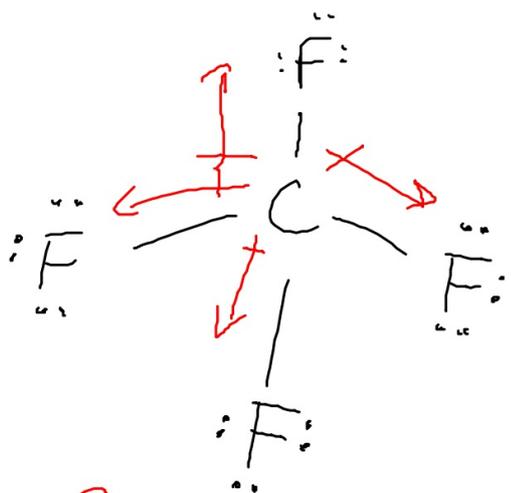
Linear Triatomic
180°

Each C=O bond is polar (unequal sharing ... 1 atom pulls with a higher e⁻N than the other) ... BUT overall CO₂ is nonpolar (equal sharing) b/c each C=O bond is pulling in opposite directions

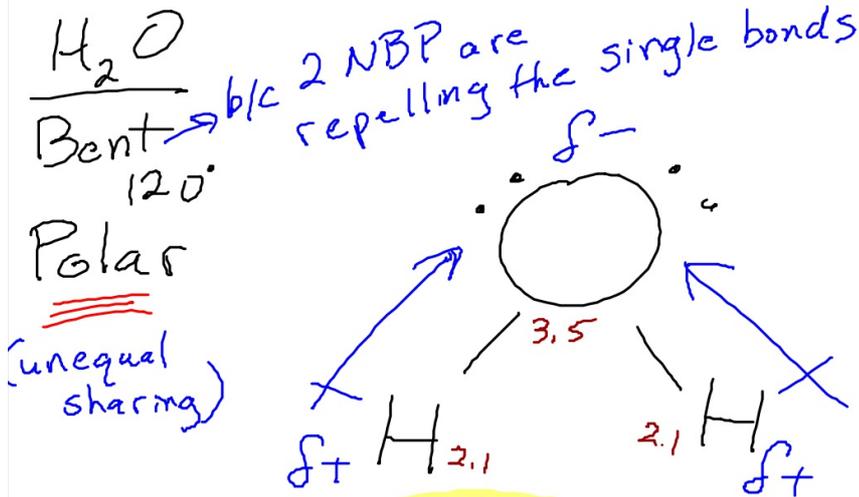
Polar



Nonpolar



Equal pulling



Do for each bond

δ^+
 partial (+) charge

δ^-
 partial (-) charge

→
 point to the more EN atom in the bond

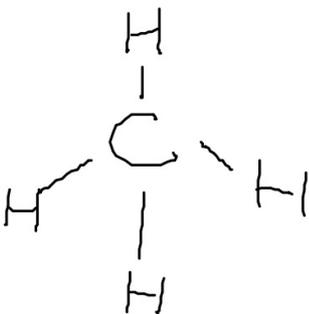
Calculating Individual Bond Polarity:

H-O → 2.1 - 3.5 = 1.4

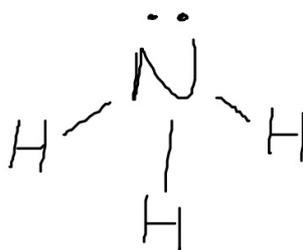
H-O → 2.1 - 3.5 = 1.4

Very Polar Covalent

Tetrahedral



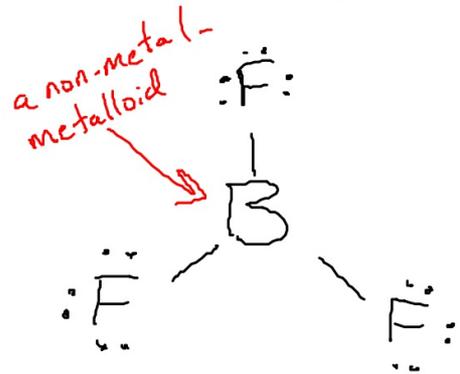
Pyramidal



109.5° → 107°

the 1 NBP has a slightly strong push to the single bonds

Trigonal Planar



120°

b/c B is an exception w/ NO NBPs & will only need 6 total e⁻

Intermolecular Attractions

What type of bond is it?

