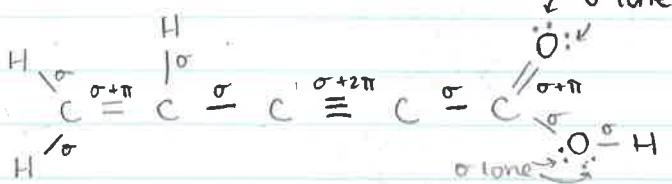


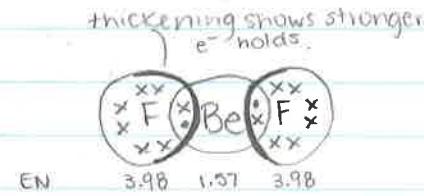
## Lewis Dot Diagrams → Shape

Oct. 21st

$\sigma$  vs.  $\pi$   
sigma pi



e.g.  $\text{BeF}_2$   
covalent!

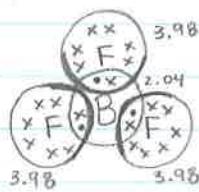


total valence pairs  
↓ σ ↓ lone ↓ π  
2 2 O O

base shape = linear  
actual shape = linear  
bond angle =  $180^\circ$

oxidation states:  $\text{Be}^{2+}$   $\text{F}^{-}$

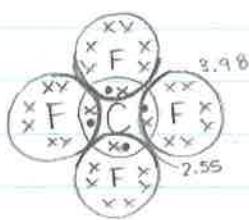
e.g.  $\text{BF}_3$   
covalent!



total  
↓ σ ↓ lone ↓ π  
3 3 O O  
NOTE: when zero, these are the same  
base shape = trigonal planar  
actual shape = trigonal planar  
bond angle =  $120^\circ$

oxidation states:  $\text{B}^{3+}$   $\text{F}^{1-}$

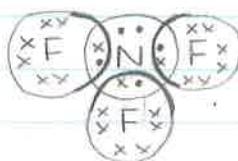
e.g.  $\text{CF}_4$   
covalent!



total  
↓ σ ↓ lone ↓ π  
4 4 O O  
base shape = tetrahedral  
actual shape = tetrahedral  
bond angle =  $109.5^\circ$

oxidation states:  $\text{C}^{4+}$   $\text{F}^{1-}$

e.g.  $\text{NF}_3$



oxidation states:  $\text{N}^{3+}$   $\text{F}^{1-}$

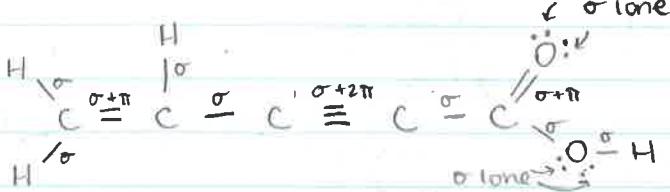
total  
↓ σ ↓ lone ↓ π  
4 3 1 O  
common!  
base shape = tetrahedral (e-shape based on σ pairs)  
actual shape = pyramidal (trigonal pyramidal)  
bond angle =  $< 109.5^\circ$

(lone pair occupy more space around central atom & ∴ repel more effectively  
"weak")

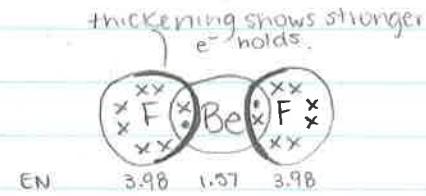
## Lewis Dot Diagrams → Shape

Oct. 21st

$\sigma$  vs.  $\pi$   
sigma pi



e.g.  $\text{BeF}_2$   
covalent!

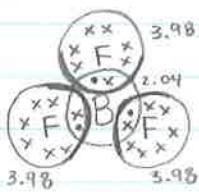


total valence pairs  
↓ σ ↓ lone ↓ π  
2 2 O O

base shape = linear  
actual shape = linear  
bond angle =  $180^\circ$

oxidation states:  $\text{Be}^{2+}$   $\text{F}^{-}$

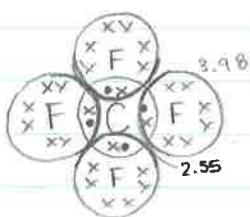
e.g.  $\text{BF}_3$   
covalent!



total  
↓ σ ↓ lone ↓ π  
3 3 O O  
NOTE when zero  
these are the same  
base shape = trigonal planar  
actual shape = trigonal planar  
bond angle =  $120^\circ$

oxidation states:  $\text{B}^{3+}$   $\text{F}^{-}$

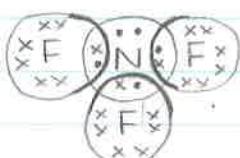
e.g.  $\text{CF}_4$   
covalent!



total  
↓ σ ↓ lone ↓ π  
4 4 O O  
base shape = tetrahedral  
actual shape = tetrahedral  
bond angle =  $109.5^\circ$

oxidation states:  $\text{C}^{4+}$   $\text{F}^{-}$

e.g.  $\text{NF}_3$



oxidation states:  $\text{N}^{3+}$   $\text{F}^{-}$

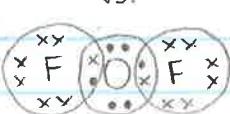
total  
↓ σ ↓ lone ↓ π  
4 3 O  
common!  
base shape = tetrahedral (e-shape based on σ pairs)  
actual shape = pyramidal (trigonal pyramidal)  
bond angle =  $< 109.5^\circ$

(lone pair occupy more space around central atom & ∴ repel more effectively "weak")

e.g.  $\text{OF}_2$



$\text{O}^{2+}$   $\text{F}^-$



↓ total  
↓ σ ↓ lone ↓ π

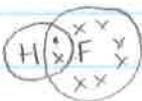
4 | 2 2 | 0

base shape = tetrahedral

actual shape = angular

bond angle =  $< 109.5^\circ$

e.g.  $\text{HF}$



↓ total  
↓ σ ↓ lone ↓ π

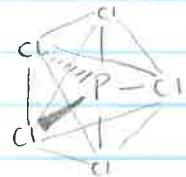
4 | 1 3 | 0

base shape = tetrahedral

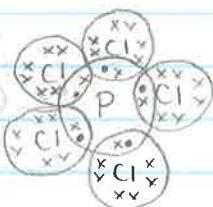
actual shape = linear

bond angle = NA

e.g.  $\text{PCl}_5$



$\text{P}^{5+}$



5 | 5 0 | 0

base shape = trigonal bi-pyramidal

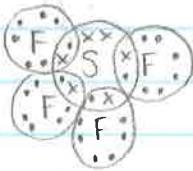
actual shape = trigonal bi-pyramidal

bond angle =  $120^\circ$  and  $90^\circ$

↑  
in the plane

↑  
above & below  
the plane

e.g.  $\text{SF}_4$



5 4 | 1 | 0

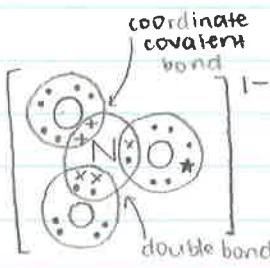
base shape = trigonal bi-pyramidal

actual shape = see-saw

bond angle =  $< 120^\circ$  and  $< 90^\circ$

↑  
in the plane

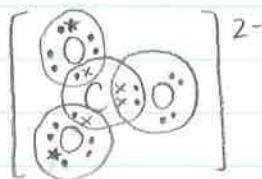
↑  
above & below  
the plane

e.g.  $\text{NO}_3^{1-}$  $\text{N}^{5+}$ 

4  $\boxed{3 \text{ O}} \text{ } 1$  ↗ (but does not determine shape)

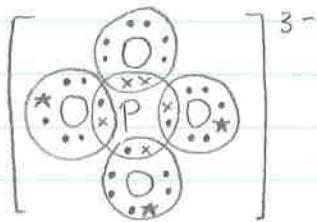
base shape = trigonal planar

actual shape = trigonal planar

bond angle =  $120^\circ$ e.g.  $\text{CO}_3^{2-}$ 4  $\boxed{3 \text{ O}} \text{ } 1$ 

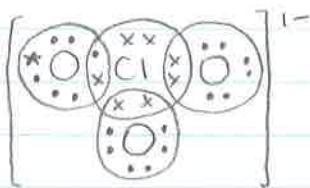
trigonal planar

trigonal planar

 $120^\circ$ e.g.  $\text{PO}_4^{3-}$  $\text{P}^{5+}$ 4  $\boxed{4 \text{ O}} \text{ } 0$ 

tetrahedral

tetrahedral

 $109.5^\circ$ e.g.  $\text{ClO}_3^{1-}$  $\text{Cl}^{5+}$ 4  $\boxed{3 \text{ O}} \text{ } 1$ 

tetrahedral

pyramidal

 $109.5^\circ$ 

presence of lone pairs  
cause  $\text{O}$ s to be <

(square planar  
is an exception)

e.g.  $\text{SF}_6$  $\text{S}^{6+}$ 6  $\boxed{6 \text{ O}} \text{ } 0$ 

octahedral

octahedral

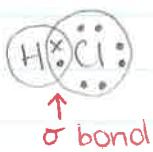
 $90^\circ$

## Covalent Bonding (including polyatomic ions)

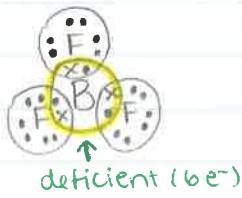
1. All halogens get their way. They form simple single sigma bonds,  $\sigma$ , even if it means breaking the octet rule.

↳ Octet deficient & expanded valence shells can result.

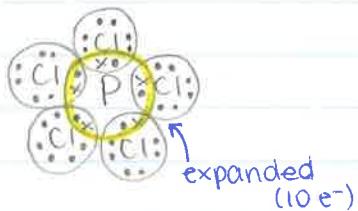
e.g.s: HCl



BF<sub>3</sub>

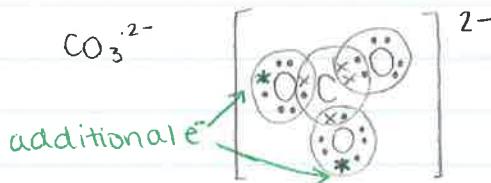


PCl<sub>5</sub>



2. Any extra electrons ( $x^{2-}$ ) will go to the most electronegative element.  
(usually oxygen)

e.g. CO<sub>3</sub><sup>2-</sup>



\* NOTE: added e<sup>-</sup> are drawn farthest from central atom /  $\sigma$  bonds (II)

3. Oxygen always finds a way to satisfy the octet rule for the central atom.

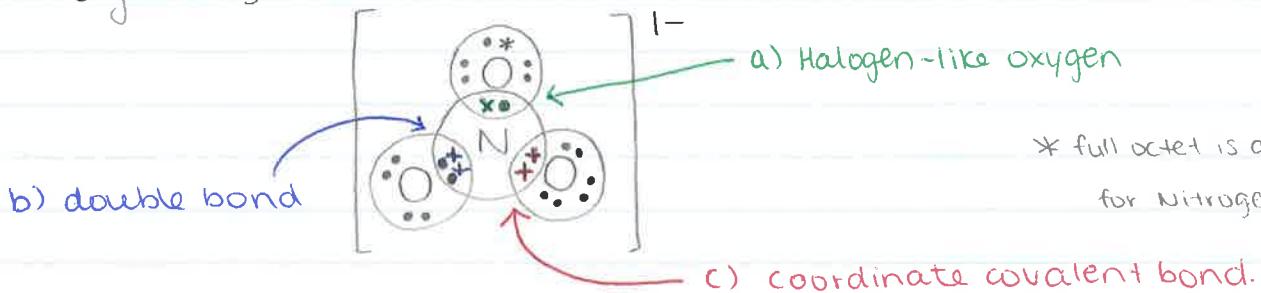
1<sup>st</sup> step → a) Halogen-like oxygens\* (see #2) behave like halogens and always form simple single sigma bonds

{ 2<sup>nd</sup> step → b) Double bonds will add two electrons to the central atom

(use until you have a full octet, 8e<sup>-</sup>)

{ 3<sup>rd</sup> step → c) Coordinate Covalent Bonds\*\* will keep the number of electrons around the central atom the same.

e.g. NO<sub>3</sub><sup>1-</sup>



\* full octet is achieved for Nitrogen!!

\* includes oxygens that already have one single bond, i.e.

\*\* single bonds in which both bonding electrons are contributed by one of the bonded atoms:

↳ indistinguishable from regular covalent bonds

