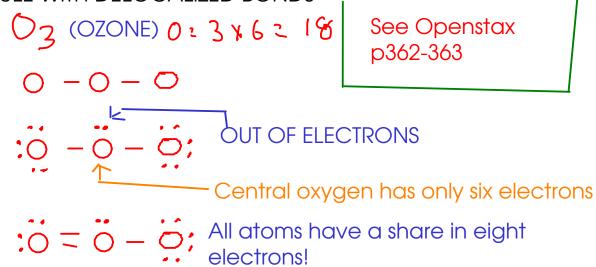
- (1) Count valence electrons
- Pick central atom and draw skeletal structure
  - central atom is usually the one that needs to gain the most electrons!
  - skeletal structure
     has all atoms connected
     to center with single
     bonds
- Distribute remaining valence electrons around structure, outer atoms first. Follow octet rule until you run out of electrons.
- Check octet rule each atom should have a share in 8 electrons (H gets 2). if not, make double or triple bonds.



The structure we drew implies that one of the outer oxygen atoms is closer to the central oxygen atom than the other one.

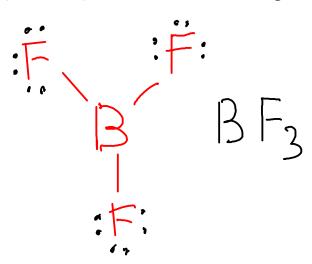
Experimentally, though, we find the two oxygen atoms to be the SAME distance from the center.

In the ozone molecule, electrons are actually being shared between ALL THREE oxygen atoms at the same time. This is called a DELOCALIZED BOND.



The structures in the green box are called RESONANCE STRUCTURES. The "real" structure of ozone is an "average" of the two resonsnce structures. The "double bond" electrons in these structures are actually shared between all three oxygen atoms

- Some atoms do not always obey the octet rule. A few, like BORON, will bond in such a way that they end up with LESS than eight electrons.



... but many more bond in such a way that they end up with a share in MORE THAN EIGHT electrons!

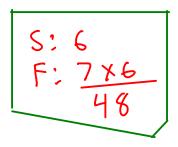
- Any atom in period three or greater can do this. SULFUR and PHOSPHORUS compounds commonly do this!

... these atoms have unfilled "d" orbitals that may participate in bonding!

- All noble gas compounds (example: XENON compounds with oxygen and fluorine) exhibit this behavior!

### **EXAMPLES:**





- The central SULFUR atom has a share in TWELVE total electrons, not eight!
- The SHAPE of the sulfur hexafluoride molecule in three dimensions agrees with the picture of six fluorine atoms each sharing a pair of electrons with a sulfur center.

This structure obeys the octet rule.

This molecule does NOT obey the octet rule. Phosphorus ends up with ten electrons instead of eight.

## FORMAL CHARGE

- You can sometimes draw more than one structure for a molecule that appears correct. How can you determine which one is more likely?

#### - USE FORMAL CHARGE!

- Formal charge is a hypothetical charge on each atom in a structure. It assumes:
  - (1) All bonding electrons are shared EQUALLY between atoms
  - (2) Lone pairs are NOT shared.

FORMAL CHARGE = ORIGINAL # OF VALENCE ELECTRONS — NUMBER OF UNSHARED ELECTRONS

\* The sum of the formal charges of all atoms in a structure should equal to the charge of the molecule (0 for neutral molecules)

The "better" Lewis structure will have:

- Lower magnitudes of formal charge (00 is better than +2-2)
- Negative formal charges on ELECTRONEGATIVE atoms, or positive formal charges on atoms that are less electronegative.

EXAMPLE: LOC/2

... calculate formal charges to tell which structure is more likely!

$$0:6-2-4=0$$
 $0:6-1-6=-1$ 
 $0:6-1-6=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4-4-0=0$ 
 $0:4$ 

Based on formal charge, the structure on the left is preferred. (Lower magnitudes of charge on the left structure!)

VS

... we can determine which of these structures is more likely by calculating \
formal charges!

$$N:S-3-2=6$$

## Which structure is more likely?

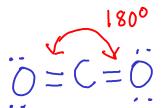
Based on formal charge, the structure on the right is preferred. (lower magnitudes of formal charge)

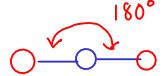
#### PREDICTING MOLECULAR SHAPE

The shape of simple molecules (and parts of larger molecules) can be easily predicted using the VSEPR model

# VSEPR = Valence Shell Electron Pair Repulsion Model

- Each BOND or LONE PAIR OF ELECTRONS around an atom will try to move itself as far away from other bonds or lone pairs as possible!

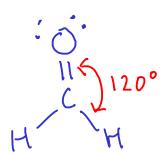


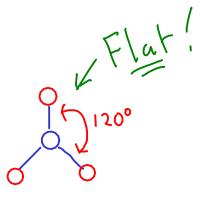


For the two red circles to be farthest apart, they must be 180 degrees apart

LINEAR MOLECULES

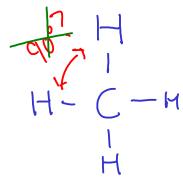
ANY diatomic (two-atom) molecule is linear, but only some three-atom molecules are!





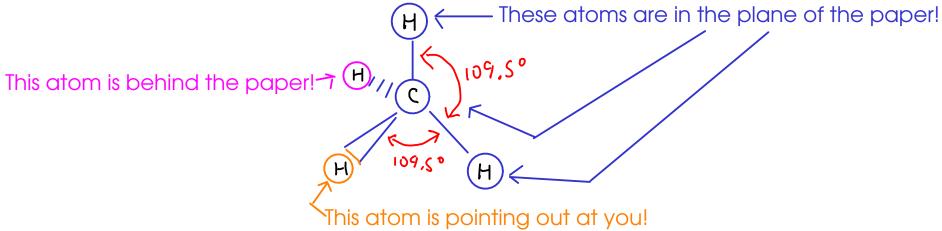
For the three red circles to be farthest apart, they spread out so that each is 120 degrees from the others!

TRIGONAL PLANAR MOLECULES



These hydrogen atoms might appear at first glance to be 90 degrees apart, but remember that molecules exist in THREE DIMENSIONS, not two!

Each hydrogen atom is actually 109.5 degrees apart, forming a TETRAHEDRON.



To see the tetrahedron in three dimensions WITHOUT buying a molecular model kit, just take four balloons, blow them up, and then tie them together. The knot will be the central atom, and the balloons will line themselves up to be 109.5 degrees apart.