A DOT STRUCTURE FOR A LARGER MOLECULE

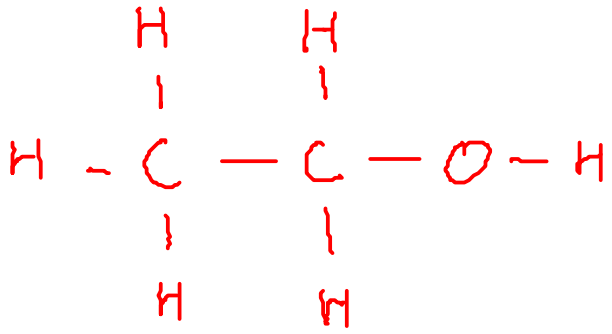
CH_3CH_2OH ETHANOL!

C : 4 x 2 = 8
H : 1 x 6 = 6
O : 6 x 1 = 6
| 20

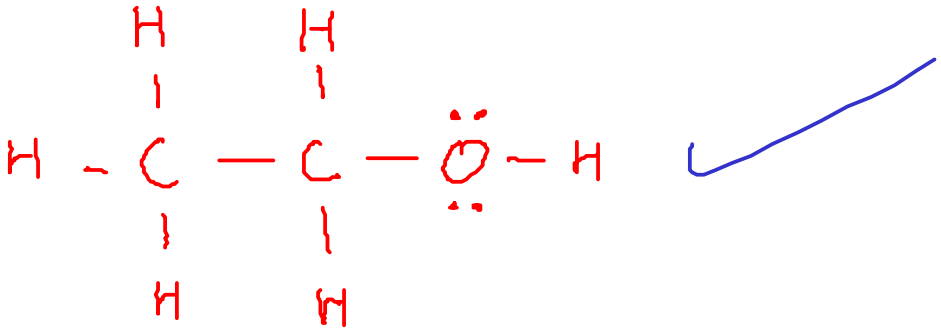
This formula gives us a hint to the structure of the molecule. Ethanol has THREE centers: the two carbon atoms and the oxygen atom.

1. Count valence electrons

2. 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



3. Distribute remaining valence electrons around structure, outer atoms first. Follow octet rule until you run out of electrons.



4. Check octet rule - each atom should have a share in 8 electrons (H gets 2). if not, make double or triple bonds.

A DOT STRUCTURE FOR A MOLECULE WITH DELOCALIZED BONDS

O_3 (OZONE)

$O = 3 \times 6 = 18$

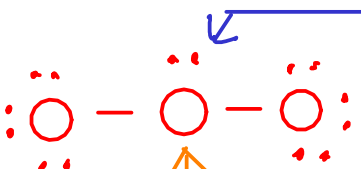
See text, 9.7 p 350-352

1. Count valence electrons

2. 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

3. Distribute remaining valence electrons around structure, outer atoms first. Follow octet rule until you run out of electrons.

4. Check octet rule - each atom should have a share in 8 electrons (H gets 2). if not, make double or triple bonds.

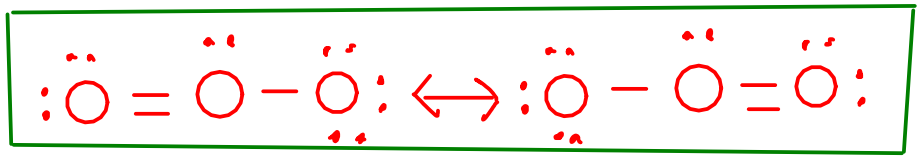


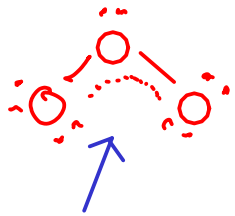
All atoms have a share in eight electrons here.

This structure suggests that one of the outer oxygen atoms is closer to the central oxygen atom than the other one!

Experimentally, we observe that both outer oxygen atoms are the SAME distance from the center.

In the molecule, electrons are actually being shared between ALL THREE oxygen atoms. This is a DELOCALIZED bond!





These are RESONANCE structures. The real structure is an "average" of these two. The "double bond"'s electrons are shared between all three oxygen atoms!

A DOT STRUCTURE FOR A POLYATOMIC ION

1. Count valence electrons

2. 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

3. Distribute remaining valence electrons around structure, outer atoms first. Follow octet rule until you run out of electrons.

4. Check octet rule - each atom should have a share in 8 electrons (H gets 2). if not, make double or triple bonds.

NH_4^+

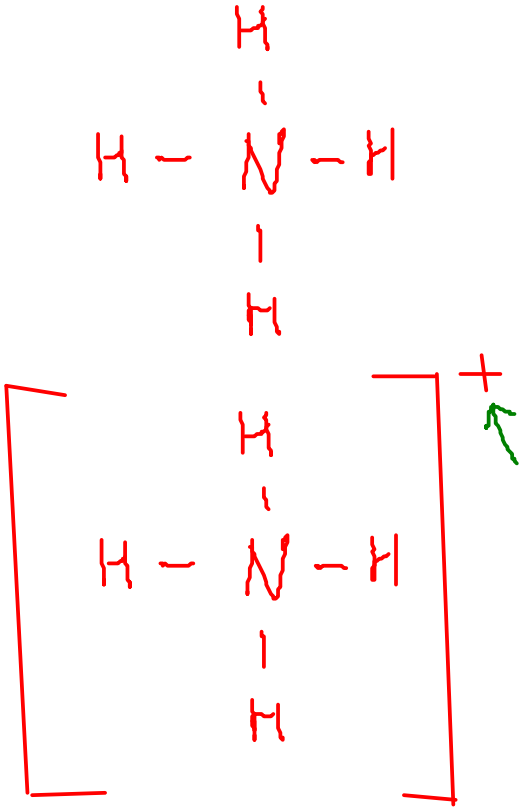
N : 5

H : 1 x H

9 ← An odd number of electrons? But Lewis structures deal in PAIRS of electrons!

−1 ← Subtract an electron to account for ammonium's +1 charge

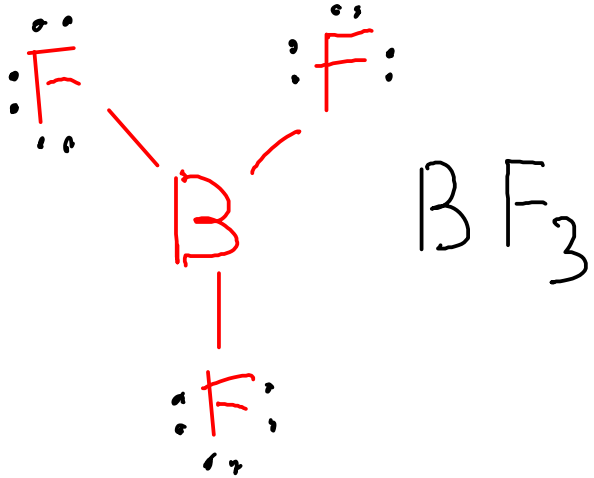
8



Draw brackets around the complete structure for an ion, and then indicate the charge as shown!

EXPANDED VALENCE and other exceptions to the "octet rule"

- 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!

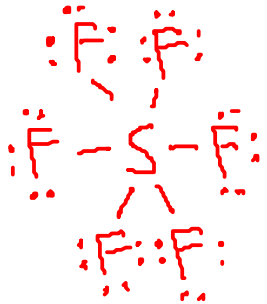
- Any atom in period three or greater can do this. SULFUR and PHOSPHORUS compounds commonly do this!

... 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:

SF_6



S: 6
F: 7 x 6
48

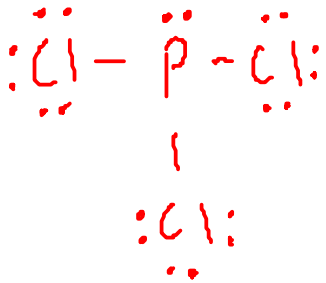
- The central SULFUR atom has a share in TWELVE total electrons, not eight!

- 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.

PCl_3 vs PCl_5

PCl_3

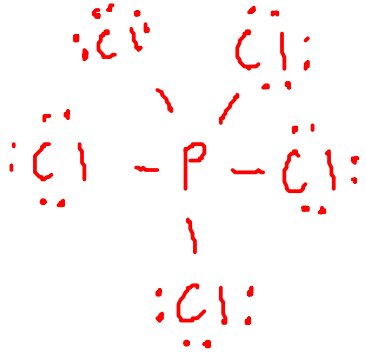
P : 5
Cl : 7 x 3 = 21
26



This structure obeys the octet rule.

PCl_5

P : 5
Cl : 7 x 5 = 35
40



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

FORMAL CHARGE

- You can often 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.

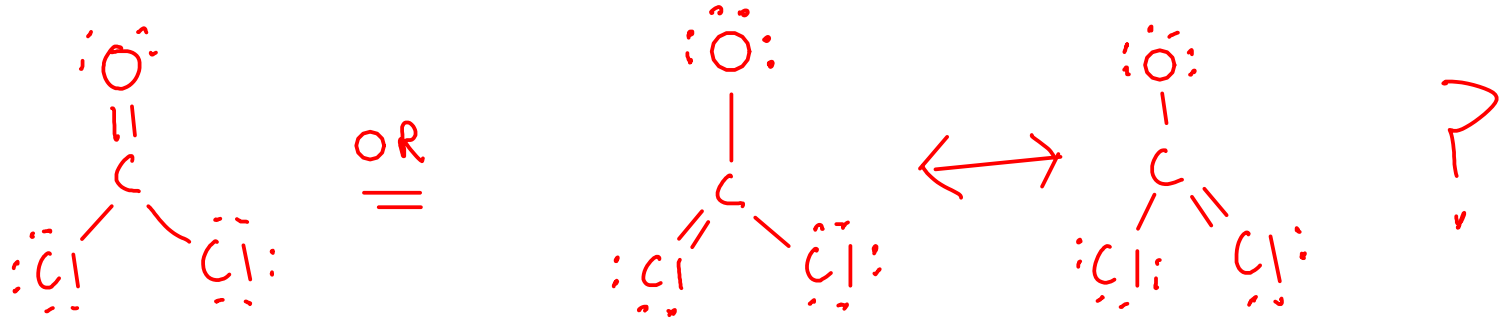
FORMAL CHARGE = ORIGINAL # OF VALENCE ELECTRONS − NUMBER OF BONDS − NUMBER OF UNSHARED ELECTRONS

* 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 (0 0 is better than +2 -2)
- Negative formal charges on ELECTRONEGATIVE atoms, or positive formal charges on atoms that are less electronegative.

EXAMPLE: $COCl_2$



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

O: 6 − 2 − 4 = 0

C: 4 − 4 − 0 = 0

Cl: 7 − 1 − 6 = 0

Cl: 7 − 1 − 6 = 0

O: 6 − 1 − 6 = −1

C: 4 − 4 − 0 = 0

=Cl: 7 − 2 − 4 = +1

−Cl: 7 − 1 − 6 = 0

* The sum of the formal charges on both structures equals zero - so we've drawn them correctly. (This is a neutral molecule, not a polyatomic ion)

* The structure on the left is preferred. It has LOWER formal charges (all zeroes) than the structure on the right (+1/-1/0). The structure on the right also places a positive formal charge on the electronegative element CHLORINE.

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

H: 1 − 1 − 0 = 0

C: 4 − 3 − 2 = −1

N: 5 − 4 − 0 = +1

H: 1 − 1 − 0 = 0

C: 4 − 4 − 0 = 0

N: 5 − 3 − 2 = 0

Which structure is more likely?

* The HCN structure (right) is more likely. It has lower formal charges than the HNC structure. Also, the structure on the left places a positive formal charge on electronegative NITROGEN, while the less electronegative CARBON gets a negative formal charge.

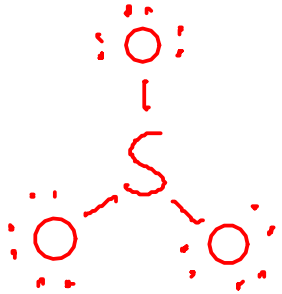
Let's look at sulfur trioxide. SO_3

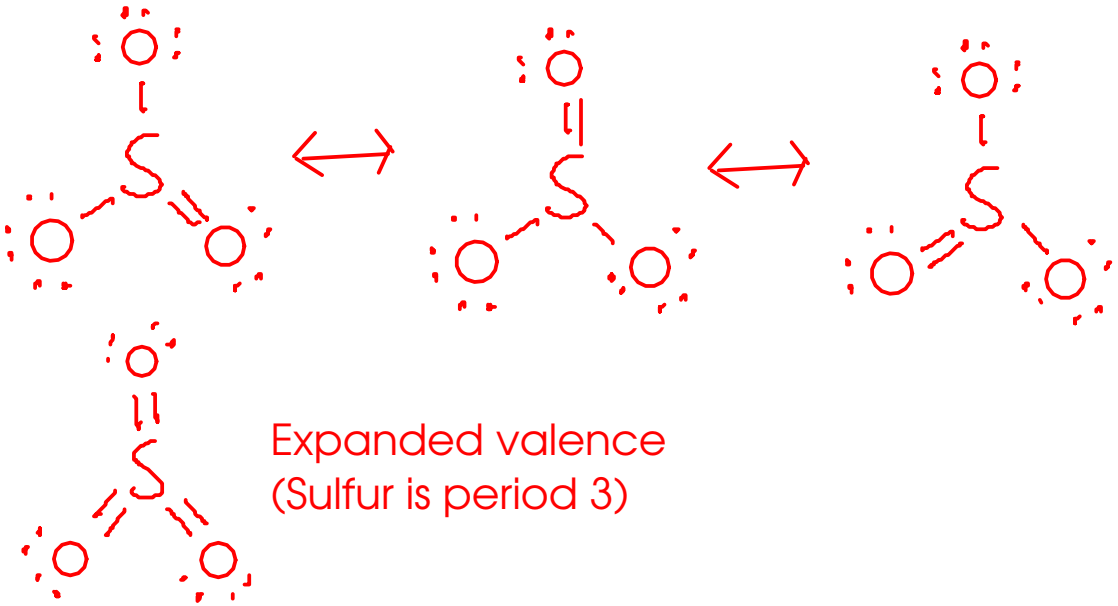
S: 6

O: 6 x 3 = 18

24 e^-

Skeletal structure:

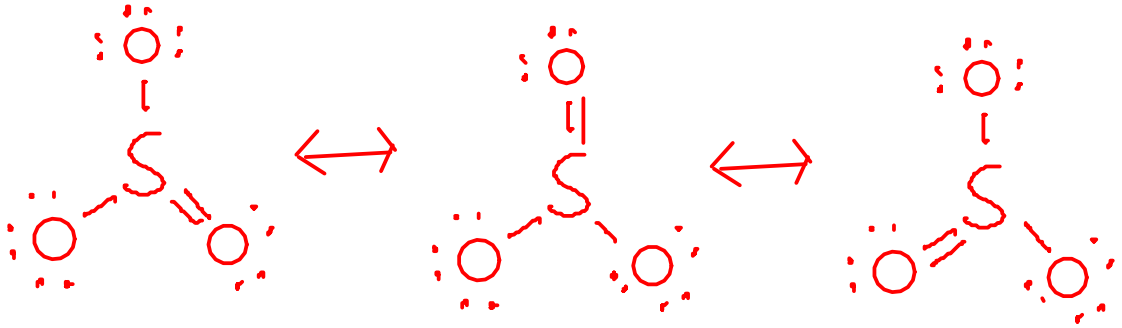




Resonance structures.

Expanded valence
(Sulfur is period 3)

To decide which structure is preferred, let's look at formal charges.



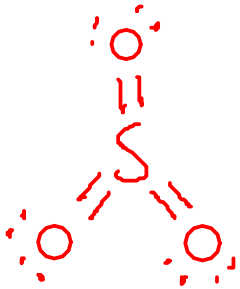
Resonance structures.

S: 6 - 4 - 0 = +2

O-: 6 - 1 - 6 = -1

O-: 6 - 1 - 6 = -1

O=: 6 - 2 - 4 = 0



Expanded valence
(Sulfur is period 3)

S: 6 - 6 - 0 = 0

O=: 6 - 2 - 4 = 0

O=: 6 - 2 - 4 = 0

O=: 6 - 2 - 4 = 0

BASED ON FORMAL CHARGES, the expanded valence structure is preferred.

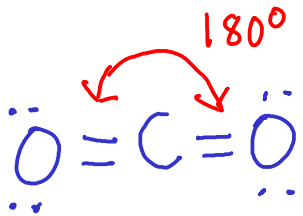
The correct structure is typically the one with minimized formal charges - even if it violates the octet rule. (Just remember - period 2 NEVER violates thie octet rule - no unfilled "d" orbitals)

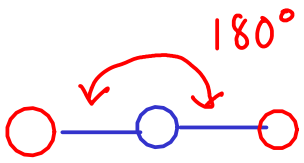
PREDICTING MOLECULAR SHAPE

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!

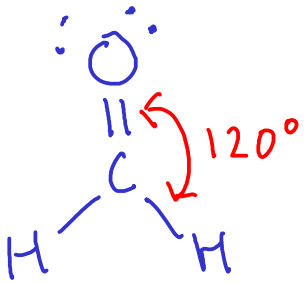


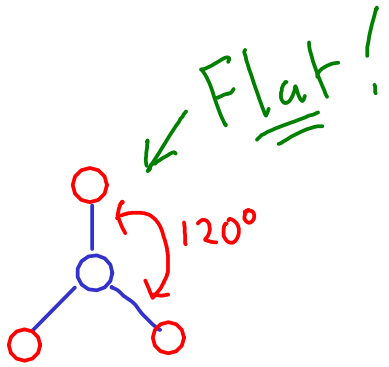


For the two red circles to be farthest apart, they must be 180 degrees apart

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