- A CHEMICAL BOND is a strong attractive force between the atoms in a compound.

3 TYPES OF CHEMICAL BOND

| Type | Held together by..n | Example |
| :--- | :--- | :--- |
| lonic bonds | attractive forces between oppositely <br> charged ions | sodium chloride |
| $\underline{\text { Covalent bonds }}$ | sharing of valence electrons between two <br> atoms (sometimes more - "delocalized <br> bonds") | water |
| Metallic bonds | sharing of valence electrons with all atoms <br> in the metal's structure - make the metal <br> conduct electricity | any metal |

*For CHM 110, you don't need to know anything more about metallic bonds than what's in this table. If you take physics, you may learn more about the characteristics of the metallic bond.

197 ... so how can you tell what kind of bond you have? You can use the traditional rules of thumb:

- Metal-Nonmetal bonds will be ionic
- Nonmetal-nonmetal bonds are usually covalent

Metalloids act like NONMETALS, here.
... but for better information about bonding, you can use ELECTRONEGATIVITY.

```
ELECTRONEGATIVITY:
-A measure of how closely to itself an atom will
hold shared electrons
\[
\begin{aligned}
& \text { p346: } \\
& \text { chart of } \\
& \text { electroneg. } \\
& \text { values }
\end{aligned}
\]
```

$p 3 S 2$,
10 th
... in other words, how ELECTRON-GREEDY an atom is!

| Bunds with.w | are... | Examples |
| :--- | :---: | :---: |
| Little or no difference in <br> electronegativity between <br> atoms | NONPOLAR COVALENT | C-C, C-H, etc. |
| Larger differences in <br> electronegativity between <br> atoms | POLAR COVALENT | $\mathrm{H}-\mathrm{F}, \mathrm{C}-\mathrm{F}, \mathrm{C}-\mathrm{Cl}$, etc. |
| Very large differences in <br> electronegativity between <br> atoms | IONIC | $\mathrm{NaCl}, \mathrm{KBr}$, etc. |

* A POLAR bond is a bond where electrons are shared unevenly - electrons spend more time around one atom than another, resulting in a bond with slightly charged ends

198 ELECTRONEGATIVITY TRENDS

- You may look up elecronegativity data in tables, but it helps to know trends!

|  |  | IIA |  |  |  |  |  |  |  |  |  |  | IIIA |  | VA | IA |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 2 | Li | Be |  |  |  |  |  |  |  |  |  |  | B | C | N | $\bigcirc$ | F |
| 3 | Na | Mg |  |  | VB |  | VIIB | - V | VIIIB |  | IB | IIB | Al | Si | P | S | Cl |
| 4 | K | Ca | Sc | Ti | V | Cr | Mn | Fe | Co | Ni | Cu | Zn | Ga | Ge | As | Se | Br |
| 5 | Rb | Sr | Y | Zr | Nb | Mo | Tc | Ru | Rh | Po | Ag | Cd | In | Sn | Sb | Te | 1 |
| 6 | Cs | Ba | * * | Hf | Ta | W | Re | Os | Ir | Pt | Au | Hg | TI | Pb | Bi | Po | At |
| 7 | Fr | Ra | Ac\| | Rf | Db | Sg | Bh | Hs | M $\dagger$ | *"inner" transition metals go here |  |  |  |  |  |  |  |

Notes:

(1) - FLUORINE is the most elecronegative element, while FRANCIUM is the least!
(2) - All the METALS have low electronegativity
(3) - HYDROGEN is similar in electronegativity to CARBON

DESCRIBING CHEMICAL BONDING
"octet rule"

- a "rule of thumb" (NOT a scienfitic law) predicting how atoms will exchange or share electrons to form chemical compounds
- atoms will gain, lose, or share enough electrons so that they end up with full "s" and " $p$ " subshells in their outermost shell.
- Why "octet"? An "s" subshell can hold two electrons, while a " p " subshell can hold six. 2+6 = 8
IONIC COMPOUNDS
- When atoms react to form IONS, they GAIN or LOSE enough electrons to end up with full "s" and "p" subshells.
example:

Aluminum loses its outer $[A r] 3 d^{10} 4 s^{24} 4 p^{5}$ three electrons, and each bromine gains one!

$$
\mathrm{Al}+3 \mathrm{Br} \rightarrow \mathrm{AlBr} 3 / 1 s^{2} 2_{r}^{2} 2 p^{6}
$$

$$
[\mathrm{Ne}] 3 s^{2} 3 p^{\prime}>[\mathrm{Ar}] 3 d^{10} 4 s^{2} 4 p^{5} \quad \mathrm{Al}^{3+}:[\mathrm{Ne}]
$$

$$
\underset{\text { purer }}{\longrightarrow}[A r] 3 d^{10} 4 s^{24} 4 p^{5}
$$

$$
\mathrm{Br}_{r}^{-}:[\mathrm{Ar}] 3 d^{10} 4 s^{24} p^{6}
$$

$$
\mathrm{Br}_{r}^{-}:[\mathrm{Ar}] 3 d^{10} 4 s^{24} p^{6}
$$

$$
\mathrm{Br}^{-}:[\mathrm{Ar}] 3 d^{10} 4 s^{2} 4 p^{6}
$$

... but using electron configurations to describe how aluminum bromide forms is a bit cumbersome! Can we simplify the picture a bit?

## LEWIS NOTATION / ELECTRON-DOT NOTATION

- Lewis notation represents each VALENCE electron with a DOT drawn around the atomic symbol. Since the valence shell of an atom contains only "s" and "p" electrons, the maximum number of dots drawn will be EIGHT.
- To use electron-dot notation, put a dot for each valence electron around the atomic symbol. Put one dot on each "side" of the symbol ( 4 sides), then pair the dots for atoms that have more than four valence electrons.
examples:





Which "side" you draw the dots on isn't important, as long as you have the right number of electrons and the right number of "pairs"


To draw a dot structure for an atom, you need to know HOW MANY valence electrons it has! You can determine this simply from the periodic table, WITHOUT writing the whole electron configuration!

... but how do we use this to describe a reaction that produces ions? Let's look at our previous example!


Aluminum is oxidized!

... this is a bit easier to follow than looking at all those letters and numbers in the electron configurations for these elements!

## MOLECULAR COMPOUNDS

- Form when atoms SHARE electrons instead of transferring them. This results in the formation of MOLECULES ... groups of atoms held together by electron-sharing.

How might atoms SHARE electrons? By coming together close enough so that their atomic ORBITALS overlap each other:


Each hydrogen atom has a single electron in a 1 s orbital.

... so how would this look using dot notation?

a single shared pair of electrons.
This is called a SINGLE BOND

In dot structures, SHARED PAIRS of electrons are often written as DASHES to make the structures look neater.


* Why doesn't hydrogen end up with eight electrons? Because hydrogen has only the first shell, which contains only a single "s" subshell (NO "p"
subshell). This "s" subshell is full with two electrons, and that's all hydrogen needs to get.

Let's look at OXYGEN ...


The oxygen atoms share TWO pairs of electrons. This is called a DOUBLE BOND

OR
$\because \because O O$ On: Each oxygen atom has a share in eight electrons!

A few notes on the double bond:

- For atoms to share more than one pair of electrons, they have to move closer to one another than they would if they were only sharing one pair of electrons. This BOND DISTANCE is measurable!
- It takes more energy to break a double bond between two atoms than it
(2) would to break a single bond between the same two atoms. This BOND ENERGY is also measurable!

: N••••••
- .-..... a

We know that nitrogen exists in air as the diatomic molecule $\mathrm{N}_{2}$
$\because N: M: N$ The nitrogen atoms share THREE pairs of electrons. This is called a TRIPLE BOND

OR
$: N E M: \quad$ bond in nitrogen gas apart!

A few notes on the triple bond:

- For atoms to share three pairs of electrons, they have to move closer to one another than they would if they were sharing one or two pairs of electrons. Triple bonds have the shortest BOND DISTANCE of all covalent bonds.
(2)
- It takes more energy to break a triple bond between two atoms than it would to break either a single or double bond between the same two atoms. The triple bond has the largest BOND ENERGY of all three kinds of covalent bonds.

SO FAR, we've seen that ...
(1) Atoms may share one, two, or three pairs of electrons with each other.
(2) Atoms will usually share enough electrons so that each atom ends up with a share in EIGHT electrons - the "octet rule"

- HYDROGEN will only end up with two electrons!
- Some other atoms may end up with more or less than eight electrons. Exceptions to the octet rule are covered in Chapter 9.

NOW, how could we come up with dot structures for some more complicated (and therefore, more interesting) molecules?

Examples:



$$
\mathrm{H}_{2} \mathrm{CO}_{3} \quad \mathrm{H}-\ddot{O}-{\underset{O}{11}}^{\mathrm{C}}-\underset{O}{-}-\mathrm{H}
$$

$$
\mathrm{CO}_{2} \because \mathrm{O}=\mathrm{C}=\ddot{O} ;
$$

(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.

Check octet rule - each atom should have a share in 8 electrons (H gets 2). if not, make double or triple bonds.
$c: 1 \times 4$
$0: 1 \times 6$
ci: $2 \times 7$
$24 e^{-}$
Choose CARBON as the
$C 1-C-C 1$
central atom since it needs to gain more electrons than the others...

$: \ddot{C}-\ddot{C} \quad$| $\ddot{O}$ |
| :--- |
| Weill pick OXYGEN to form the |
| double bond. It needed two |
| electrons initially, so it's likely |
| to share two of its own |
| electrons to get them. (chlorine |
| only needed one more...) |

(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.

$$
\begin{aligned}
& \text { NOCI } \quad \begin{array}{l}
\text { N: } \\
0: \\
0
\end{array} \\
& \text { 4: } 7 \\
& \overline{18 e^{-}}
\end{aligned}
$$

O-N-Cl We pick NITROGEN as our central atom, since it needs to gain more electrons than the others.
 so we put the last pair of electrons on the central $N$ atom.
Nitrogen has a share in only six electrons, so we'll need to consider double or triple bonds.

We make the double bond with oxygen here for the same reason we used oxygen in the previous (phosgene) structure.
(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.
$\mathrm{CO}_{2} \quad \begin{aligned} & \mathrm{C}: 4 \\ & 0: 6 \mathrm{Cl}^{-}\end{aligned}$
$0-C-O$
$\because \ddot{O}-C-\ddot{O}:$... but the carbon atom has a share in only
- $\because-$-. 4 electrons!
$\because \ddot{O}-C=\ddot{O}: \quad$... now six
$\because \because=C=\ddot{O}$ : Adding a second double bond gives
- $\mathrm{O}=\mathrm{C}=0$ : carbon a share in eight electrons.
$\because-C \equiv 0:$
The two oxygen atoms here are in identical chemical environments (both bonded to C and nothing else), so they should bond in the same way..

EXPERIMENTALLY, we find that the BOND DISTANCES of the two oxygen atoms in carbon dioxide are the same ... which does not agree with the structure in green - and supports the double-bond structure we drew first.
(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.

## $\mathrm{HNO}_{2}$ "nitrous acid"

- In oxyacids, the acidic hydrogen atoms are attached to OXYGEN atoms in the structure!
$H:|x|$
N: 1 $\times 5$
$0: 2 \times 6$
$182^{-}$
$O-N-O-H \quad$ one hydrogen atom bonded directly to oxygen.
$\because \ddot{O}-\ddot{N}-\ddot{O}-H \quad$... but that nitrogen atom has only six electrons.
$\because \ddot{O}=\ddot{N}-\ddot{O}-H$ Here, the two oxygen atoms are in different chemical environments, so we're not surprised to see them bonded differently to the central nitrogen atom.

A DOT STRUCTURE FOR A LARGER MOLECULE

$$
\left.\begin{aligned}
& C: 4 \times 2=8 \\
& H: 1 \times 6=6 \\
& 0: 6 \times 1=6
\end{aligned} \right\rvert\, 20
$$

(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.

$$
\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{OH} \text { ETHANOL! } \quad \begin{aligned}
& \mathrm{H}: 1 \times 6=6 \\
& 0: 6 \times 1=6
\end{aligned}
$$

$L$ The formula gives us a clue on how this molecule is put together. There are THREE PARTS to the molecule ... three "central" atoms!

$$
\begin{aligned}
& \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{OH} \\
& \begin{array}{cc}
H & H \\
1 & 1 \\
H-C & C-O-H \\
1 & 1 \\
H & H
\end{array} \\
& \begin{array}{ccc}
H & H \\
1 & \dot{1} \\
H & C & -\ddot{O} \\
1 & 1 & - \\
H & H & \\
H
\end{array} \\
& \square \\
& \text { Notice that the structures of ethanol and } \\
& \text { water have a similarity. Because of this } \\
& \text { structural similarity, water and ethanol } \\
& \text { dissolve in one another very well. }
\end{aligned}
$$

$0: 3 \times 6=18$
A DOT STRUCTURE FOR A MOLECULE WITH DELOCALIZED BONDS
3 (OZONE)
(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.

$$
\begin{gathered}
\text { See text, 4,7 } \\
\text { p } 356-357
\end{gathered}
$$

$\therefore O-\ddot{O}-\ddot{O}: \quad$ OUT OF ELECTRONS Central oxygen has only six electrons
$\therefore \ddot{O}=\ddot{O}-\ddot{O}:$ All atoms have a share in eight
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.

$$
\because \ddot{O}=\ddot{O}-\ddot{O}: \longleftrightarrow: \ddot{O}-\ddot{O}=\ddot{O}:
$$



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

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.

## $\mathrm{NH}_{4}{ }^{+}$

$$
N: 1 \times S
$$

$$
H: 4 \times 1 \quad . .9 \text { electrons? But all the strucutres we've }
$$

$$
9 e^{-}
$$ seen so far for molecules have an

EVEN number of electrons (it's all about PAIRS...)
$-\downarrow$ Since the ion has a charge of +1 , we $^{-1}$
$8 e^{-}$must account for that. Subtract one electron to get an ion with
$H \quad$ a charge of +1
H -NaH
1
H


Draw brackets around the structure and indicate charge in the upper right-hand corner, much like you normally do with ions.

