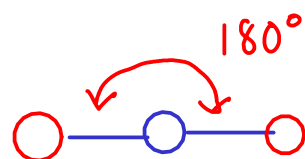
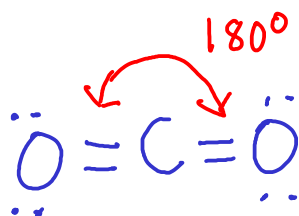


# PREDICTING MOLECULAR SHAPE

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

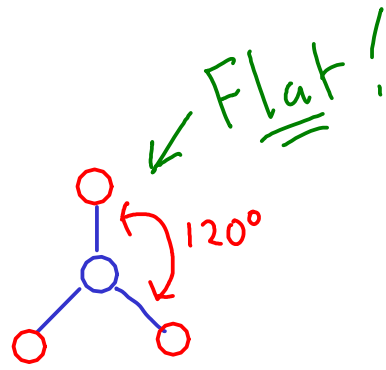
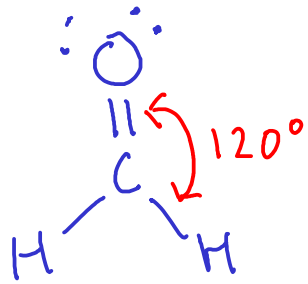
- Each BOND or LONE PAIR OF ELECTRONS around an atom will try to move itself as far away from other bonds or lone pairs as possible!



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

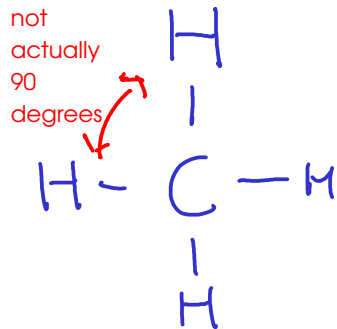
LINEAR  
MOLECULES

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



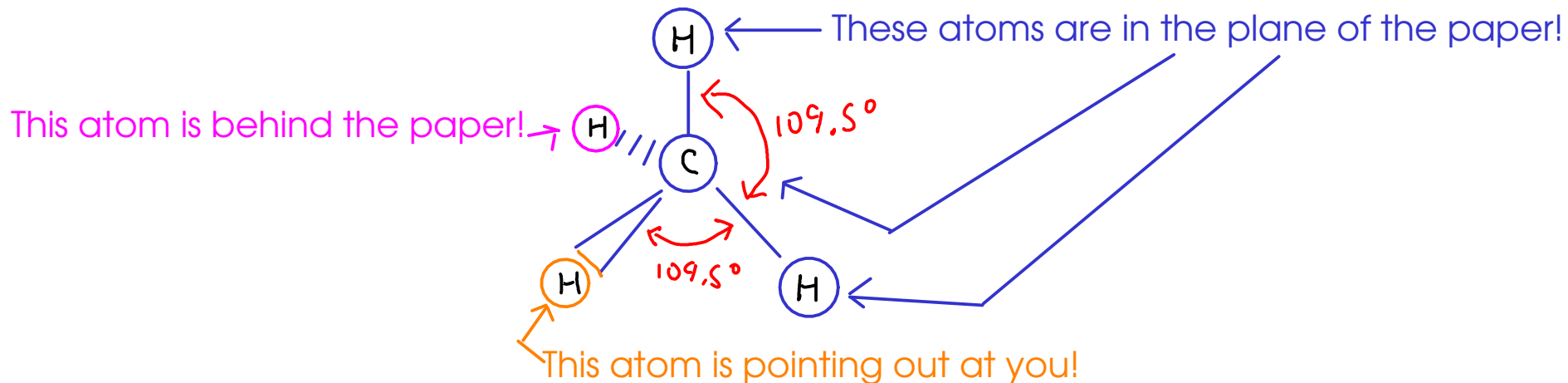
For the three red circles to be farthest apart, they spread out so that each is 120 degrees from the others!

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.



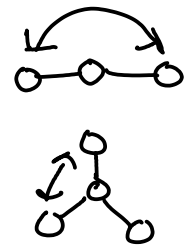
To see the tetrahedron in three dimensions WITHOUT buying a molecular model kit, just take four balloons, blow them up, and then tie them together. The knot will be the central atom, and the balloons will line themselves up to be 109.5 degrees apart.

VSEPR shapes

\* "Groups" can be either BONDS or LONE PAIRS!

VSEPR shapes:

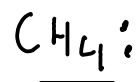
Groups* around central atom	Shape	Bond angle(s) in degrees
2	linear	180
3	trigonal planar	120
<u>4</u>	tetrahedral / pyramidal / bent	109.5
5	trigonal bipyramidal (and derivatives)	90 and 120
6	octahedral (and derivatives)	90



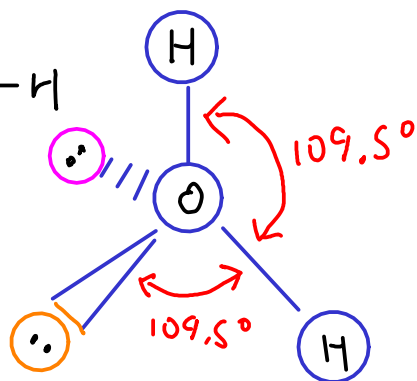
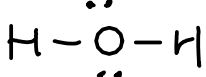
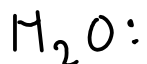
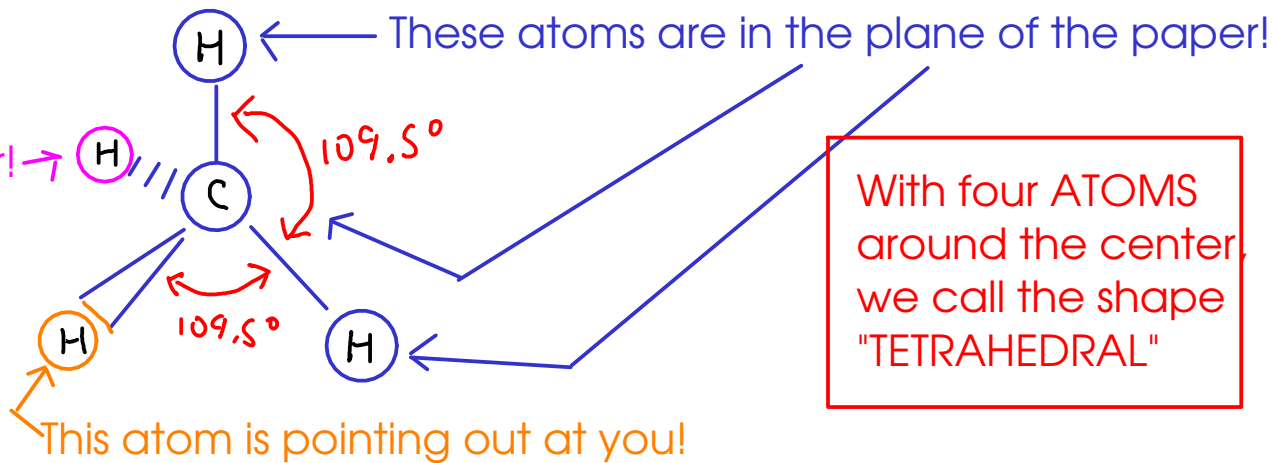
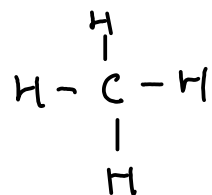
5 and 6 violate "octet rule"

5 More on "4 things around a central atom":

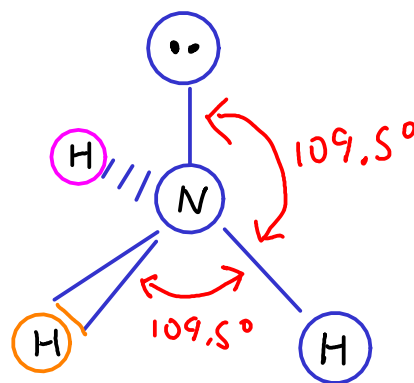
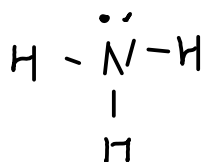
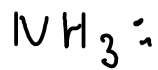
- A compound that obeys the octet rule can have a maximum of four groups around its central atom. But we describe the molecular shape based on how ATOMS are arranged around the center. What if some of those groups aren't atoms, but pairs of UNSHARED electrons?



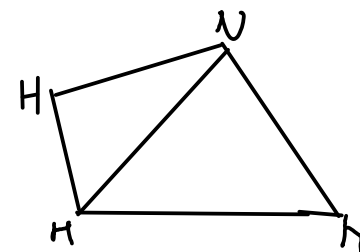
This atom is behind the paper! →



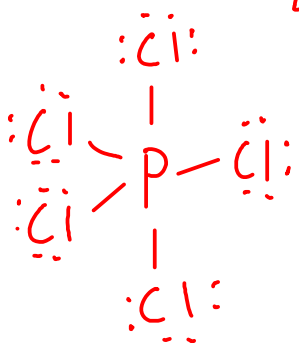
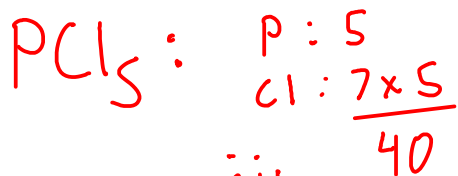
With two ATOMS and two LONE PAIRS, we call the shape "BENT"



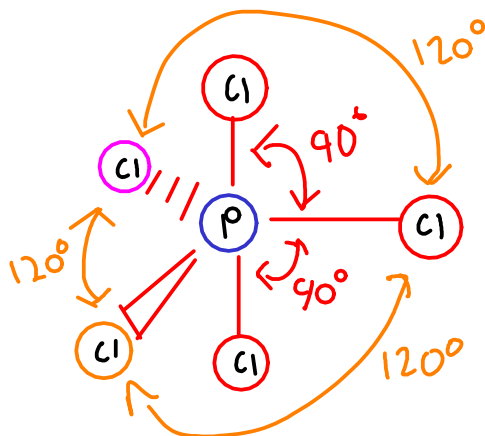
With three ATOMS and one LONE PAIR, we call the shape "PYRAMIDAL"



## 6 SHAPES OF EXPANDED VALENCE MOLECULES

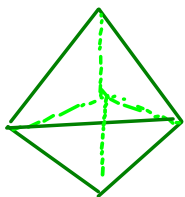


There are five atoms bonded to the central phosphorus atom, and they will attempt to get as far apart as possible from one another!



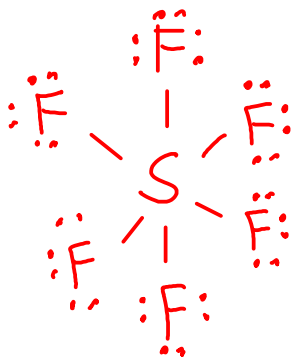
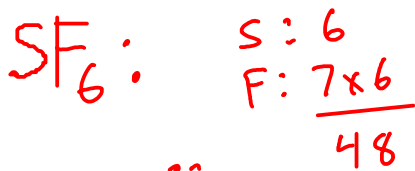
The top and bottom atoms are 90 degrees apart from the atoms around the center.

The atoms around the center are 120 degrees apart from each other.

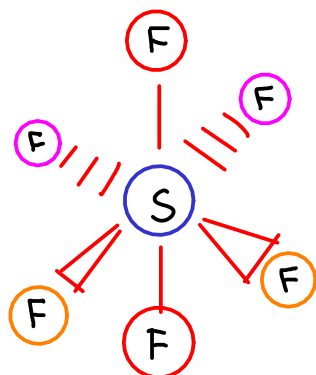


There are actually two DIFFERENT bond angles in this structure. It's called TRIGONAL BIPYRAMIDAL.

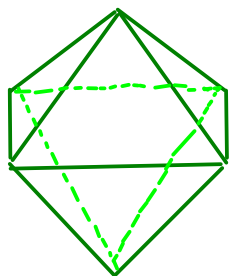
There are several derivatives of the trigonal bipyramidal shape (like the tetrahedral shape) - depending on how many things around the central atom are atoms!



There are six atoms bonded to the central sulfur atom, and they will attempt to get as far apart as possible from one another!



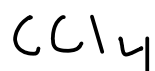
All bond angles in this arrangement are 90 degrees!



This shape is called OCTAHEDRAL, since it has eight sides.

Like the tetrahedral and trigonal bipyramidal arrangements, there are several derivatives of the octahedron - depending on how many of the six things around the center are atoms!

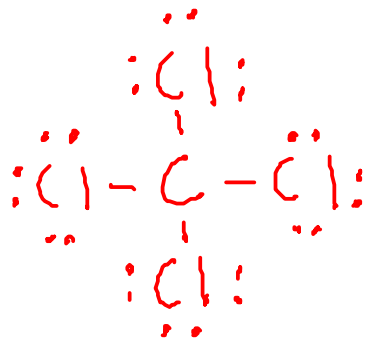
8 Examples:



$C: 4$

$Cl: 7 \times 4 = 28$

32



Molecular shape? There are FOUR atoms bonded to the central carbon atom (no lone pairs on the carbon). This gives a TETRAHEDRAL molecule.



$C: 4$

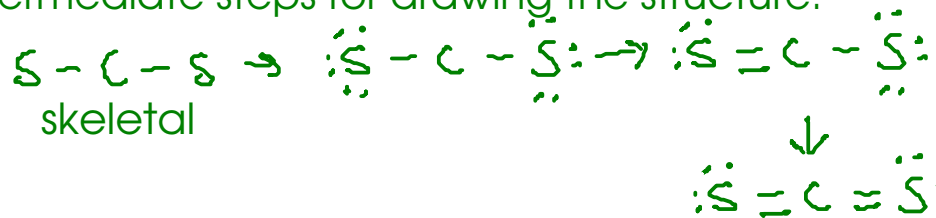
$S: 6 \times 2 = 12$

16



Molecular shape? There are TWO sulfur atoms bonded to the central carbon (no lone pairs on carbon). This is a LINEAR molecule.

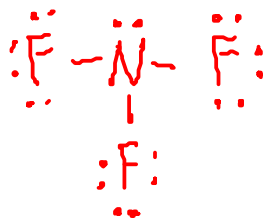
Intermediate steps for drawing the structure:



$N: 5$

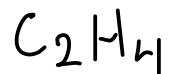
$F: 7 \times 3 = 21$

26

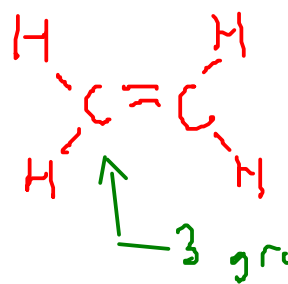
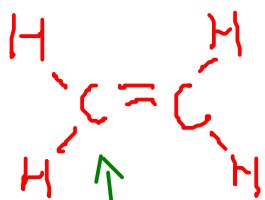


Molecular shape? There are FOUR GROUPS around the central atom, but only THREE atoms. (one lone pair on nitrogen). This is a PYRAMIDAL structure - the lone pair pushes the fluorine atoms out of the plane.



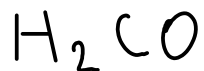
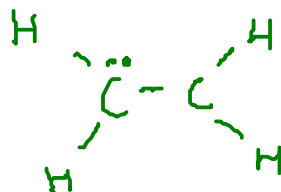


$$\begin{array}{r} \text{C} : 4 \times 2 = 8 \\ \text{H} : 1 \times 4 = 4 \\ \hline 12 \end{array}$$

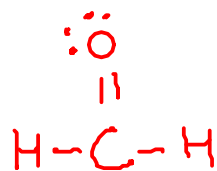


Shape? This molecule has TWO central atoms, so let's describe the shape of the molecule around each one. Each carbon center is TRIGONAL PLANAR.

Intermediate step for this structure:

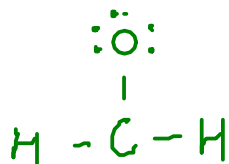


$$\begin{array}{r} \text{H} : 1 \times 2 = 2 \\ \text{C} : 4 \\ \text{O} : 6 \\ \hline 12 \end{array}$$



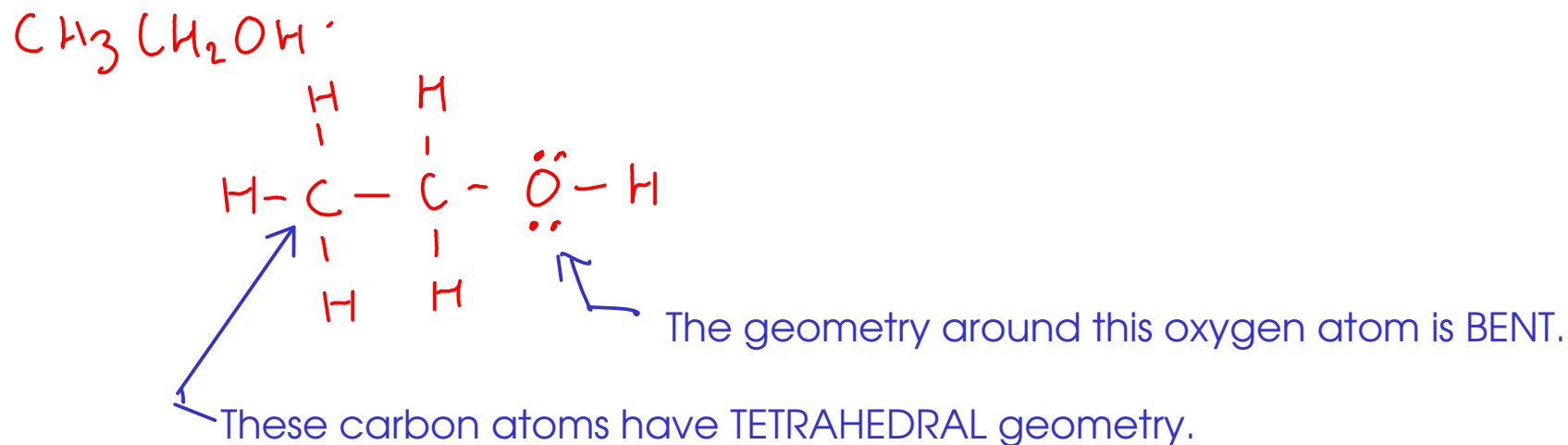
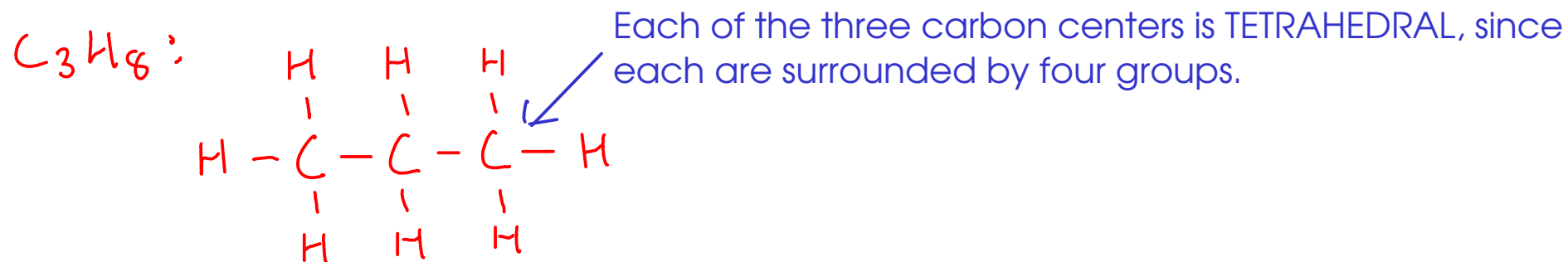
Shape? This molecule has three atoms (and no lone pairs) around the central carbon atoms. It's TRIGONAL PLANAR.

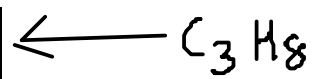
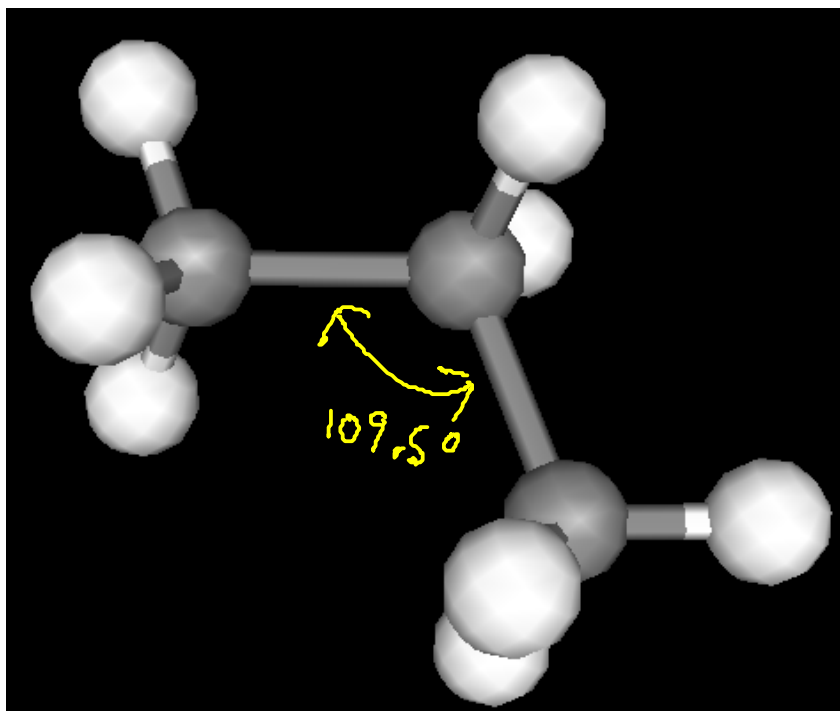
Intermediate step for this structure:



## VSEPR and large molecules

- Large molecules have more than one "center" atom
- Describe the molecule by describing the shape around each "center".

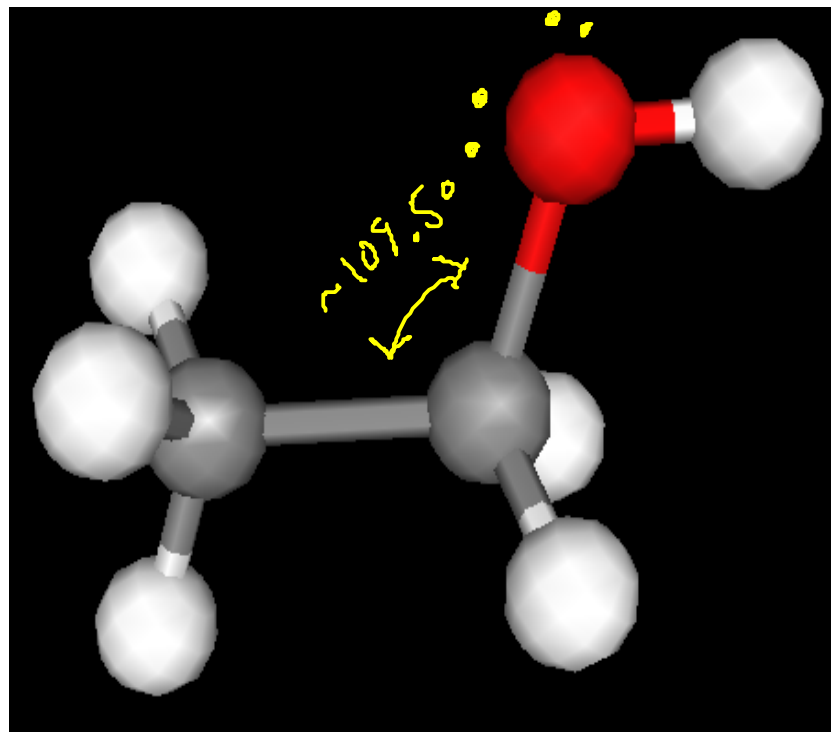




All bond angles in the propane molecule are 109.5 degrees



Like propane, the bond angles in ethanol are also close to 109.5 degrees.



## 12 POLARITY and shape:

- A polar molecule has an uneven distribution of electron density, making it have ends (poles) that are slightly charged.

POLARITY influences several easily observable properties.

- Melting point. (Polar substances have higher melting points than nonpolar substances of similar molecular weight.)

- Boiling point. (Polar substances have higher boiling points than nonpolar substances of similar molecular weight.)

- Solubility. (Polar substances tend to dissolve in other polar substances, while being insoluble in nonpolar substances. Nonpolar substances dissolve other nonpolar substances, and generally have poor solubility in polar solvents.)

- Polar molecules contain POLAR BONDS arranged in such a way that they do not cancel each other out.

... but how can we tell whether or not a bond will be POLAR? Use experimental data on ELECTRONEGATIVITY!

### ELECTRONEGATIVITY:

- A measure of how closely to itself an atom will hold shared electrons

- A bond where there is a LARGE electronegativity difference between atoms will be either POLAR or (for very large differences) IONIC!

- A bond with little or no electronegativity difference between atoms will be NONPOLAR

## ELECTRONEGATIVITY TRENDS

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

INCREASING  
ELECTRO-  
NEGATIVITY

	IA	IIA										IIIA	IVA	VA	VIA	VIIA	
2	Li	Be										B	C	N	O	F	
3	Na	Mg	IIIB	IVB	VB	VIB	VII B	VIII B	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	Pd	Ag	Cd	In	Sn	Sb	Te	I
6	Cs	Ba	La*	Hf	Ta	W	Re	Os	Ir	Pt	Au	Hg	Tl	Pb	Bi	Po	At
7	Fr	Ra	Ac*	Rf	Db	Sg	Bh	Hs	Mt	*"inner" transition metals go here							

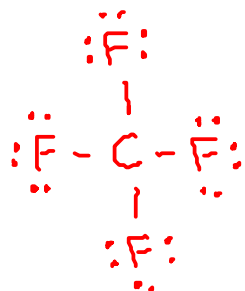
### Notes:

- ① - FLUORINE is the most electronegative element, while FRANCIUM is the least!
- ② - All the METALS have low electronegativity, and metal/nonmetal combinations form IONIC bonds
- ③ - HYDROGEN is similar in electronegativity to CARBON, so C-H bonds are considered NONPOLAR

## Examples:



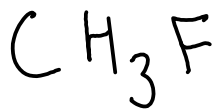
$$\begin{array}{l} \text{C: } 4 \\ \text{F: } 7 \times 4 = 28 \\ \hline 32 \end{array}$$



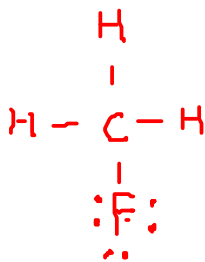
Polar or nonpolar?

\* POLAR BONDS? YES: Big electronegativity difference between C and F

\* GEOMETRY: Tetrahedral. All C-F bonds are arranged symmetrically around the center. So the molecule is NONPOLAR



$$\begin{array}{l} \text{C: } 4 \\ \text{F: } 7 \\ \text{H: } 1 \times 3 \\ \hline 14 \end{array}$$



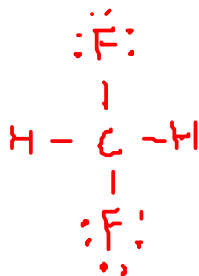
Polar or nonpolar?

\* POLAR BONDS? YES: Big electronegativity difference between C and F. C-H bonds are NONPOLAR.

\* GEOMETRY: Tetrahedral. Fluorine will pull electrons towards itself. The fluorine end of the molecule will have a slight negative charge. POLAR MOLECULE



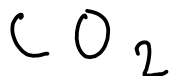
$$\begin{array}{l} \text{C: } 4 \\ \text{F: } 7 \times 2 = 14 \\ \text{H: } 1 \times 2 = 2 \\ \hline 20 \end{array}$$



Polar or nonpolar?

\* POLAR BONDS? YES: Big electronegativity difference between C and F. C-H bonds are NONPOLAR.

\* GEOMETRY: Tetrahedral. Fluorines will pull electrons towards themselves. In three dimensions, the fluorine atoms are on one side of the molecule, while the hydrogens are on the other. POLAR MOLECULE



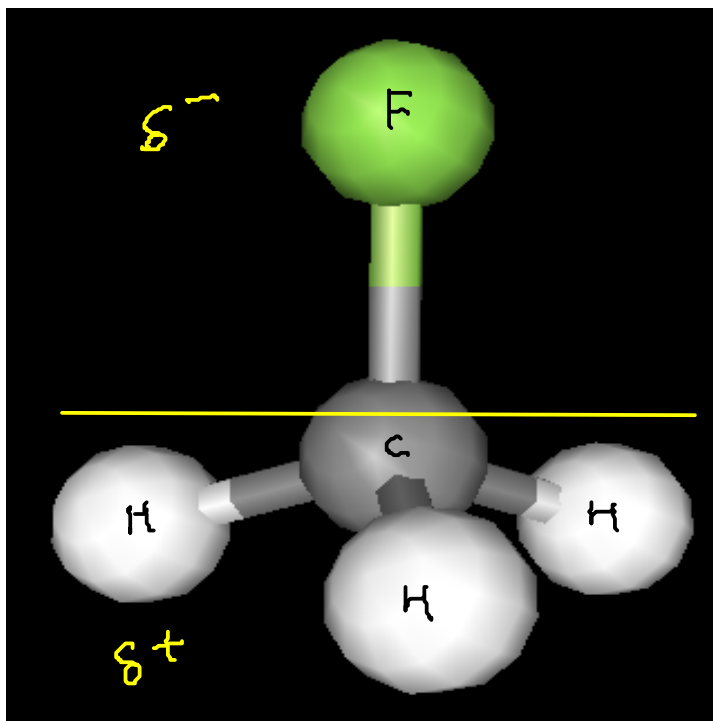
$$\begin{array}{l} \text{C: } 4 \\ \text{O: } 6 \times 2 = 12 \\ \hline 16 \end{array}$$



Polar or nonpolar?

\* POLAR BONDS? YES: Big electronegativity difference between C and O.

\* GEOMETRY: Linear. The two C=O bonds are directly opposite one another, so the overall molecule is nonpolar.



←  $\text{CH}_3\text{F}$  "fluoromethane"

Fluorine is able to pull electron density through the molecule, as it is being opposed by much less electronegative hydrogen atoms.

"difluoromethane"  $\text{CH}_2\text{F}_2$  →

In 2D, the fluorine atoms appear to be on the opposite sides of the molecule, but in 3D they are on the same side.

