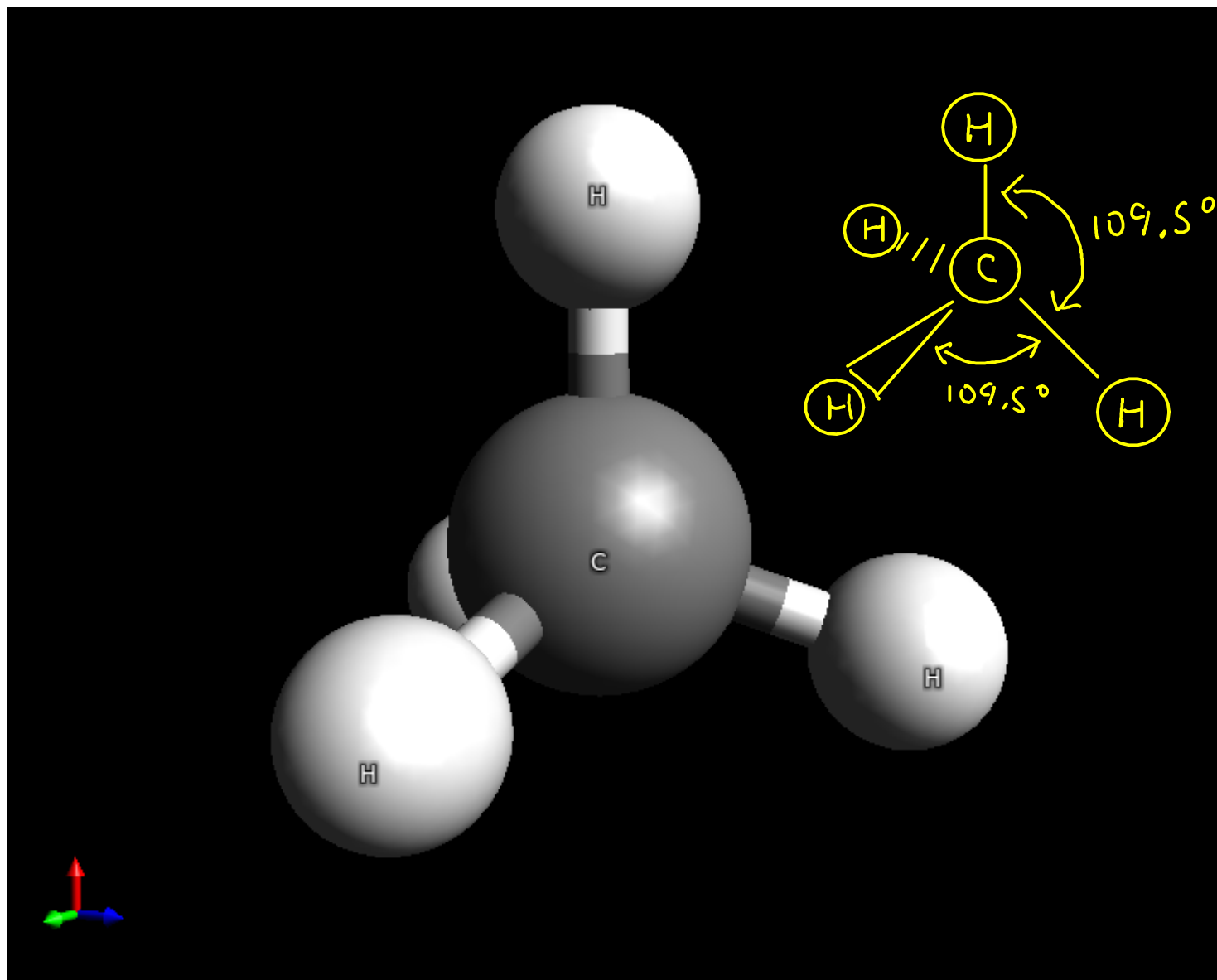
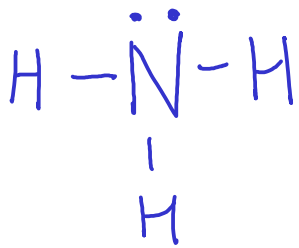


Here's a computer ball-and-stick rendering of the methane molecule.



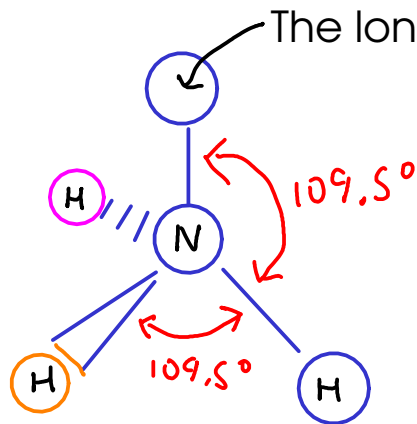
DERIVATIVES OF THE TETRAHEDRON

- What if there are lone pairs? The way the shape of a molecule is described depends on the ATOMS in the molecule, even though lone pairs play a role in the positions of the atoms.

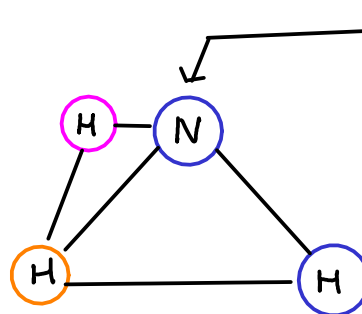


Since there are four "things" around the nitrogen atom, we would expect them to be approximately 109.5 degrees apart (in other words, TETRAHEDRAL). BUT ... only three of these things are atoms.

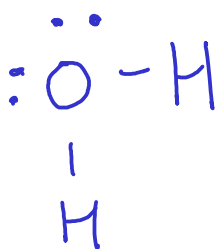
The atoms are arranged in a PYRAMID shape, so we call this molecule PYRAMIDAL!



The lone pair takes one position in the tetrahedron



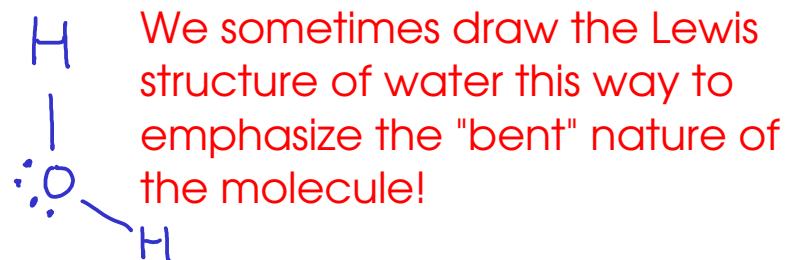
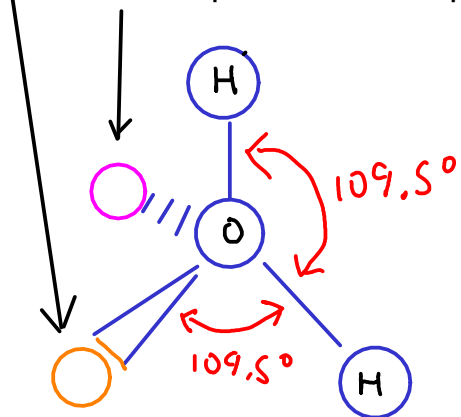
By just looking at the atoms, you can see the pyramid - with the central nitrogen atom as the top and the hydrogen atoms forming the base of the pyramid.



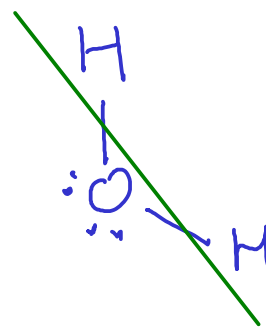
Since there are four "things" around the oxygen atom, we would expect them to be approximately 109.5 degrees apart (in other words, TETRAHEDRAL). BUT... only two of these things are atoms.

The atoms are all in a single plane, but they are not lined up in a straight line. We call this shape "BENT".

Lone pairs take up two positions in the tetrahedron

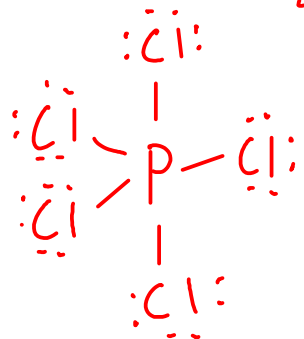


Notice that this molecule has two "sides", one with the oxygen atom and one with hydrogen atoms.

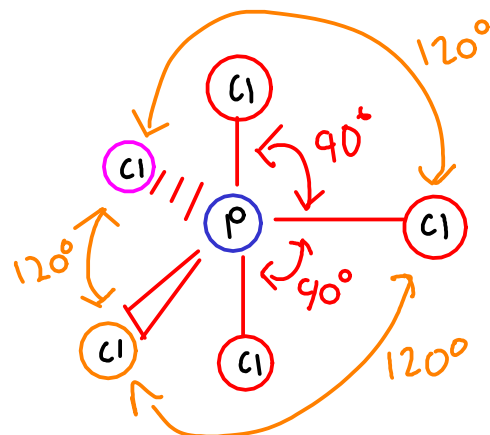


* These atoms are in the same plane, like carbon dioxide. But they are not arranged linearly!

SHAPES OF EXPANDED VALENCE MOLECULES

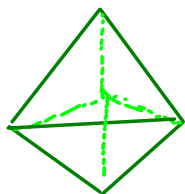


There are five atoms bonded to the central phosphorus atom, and they will attempt to get as far apart as possible from one another!



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.

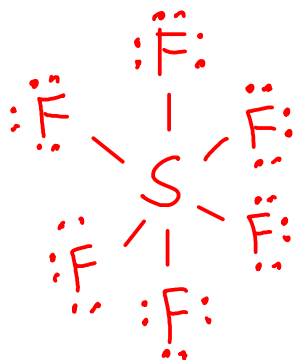


There are actually two DIFFERENT bond angles in this structure. It's called TRIGONAL BIPYRAMIDAL.

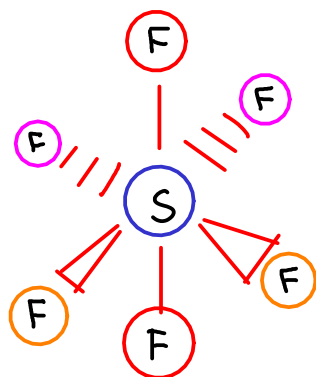
There are several derivatives of the trigonal bipyramidal shape (like the tetrahedral shape) - depending on how many things around the central atom are atoms!



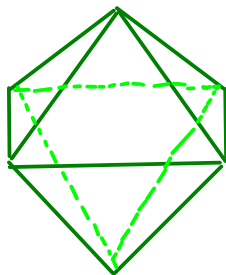
$$\begin{array}{r} \text{S: } 6 \\ \text{F: } 7 \times 6 \\ \hline 48 \end{array}$$



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



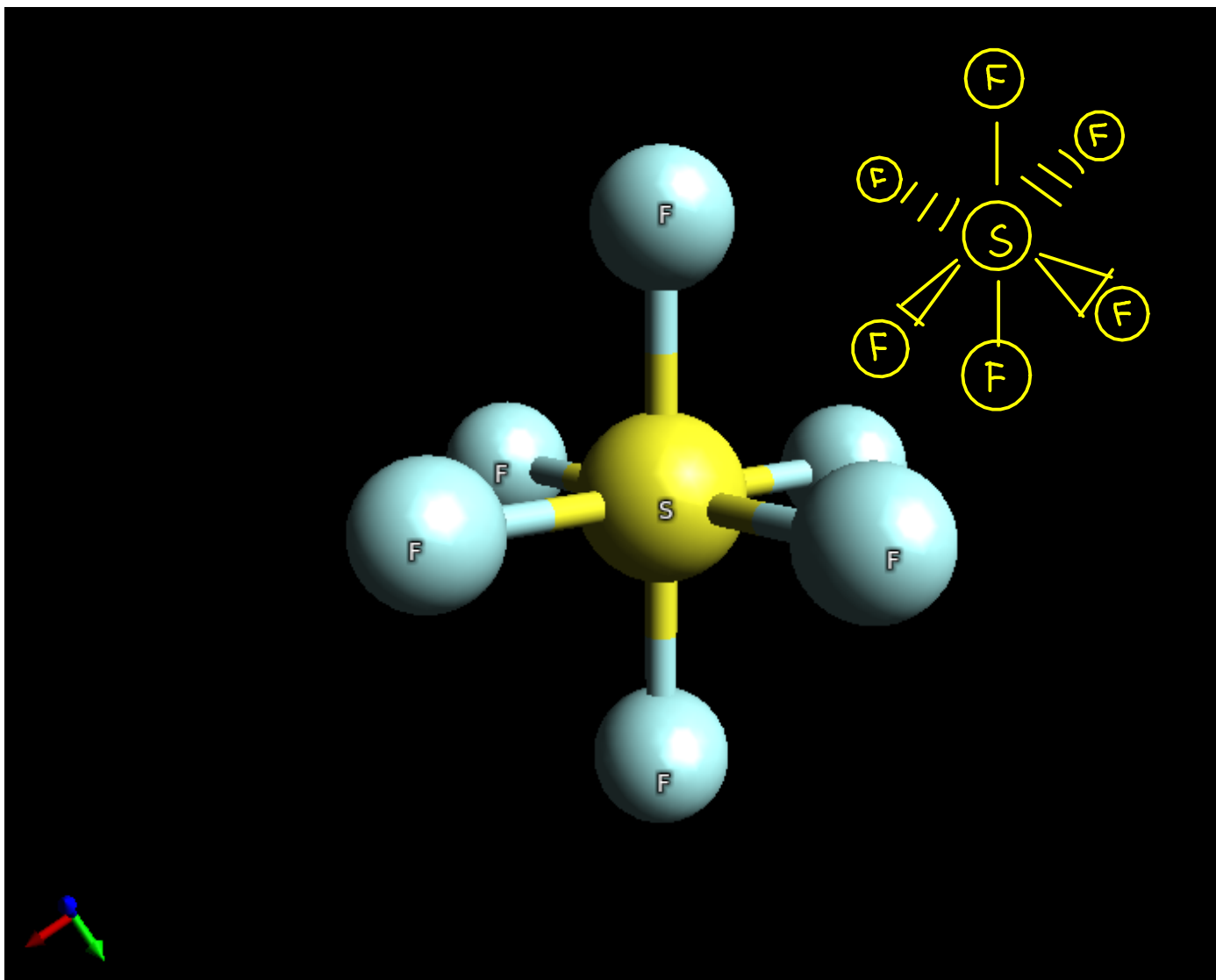
All bond angles in this arrangement are 90 degrees!



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!

Here's a ball-and-stick rendering of the sulfur hexafluoride molecule:



POLARITY

- When atoms share electrons, the electrons might not be EVENLY shared. Shared electrons may spend more time around one atomic nucleus than the other.
- When electrons are shared UNEVENLY, this results in a POLAR BOND.

... 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 (AGAIN!)

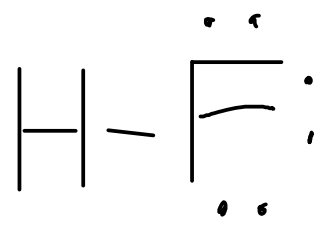
- You may look up electronegativity data in tables, but it helps to know trends!

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



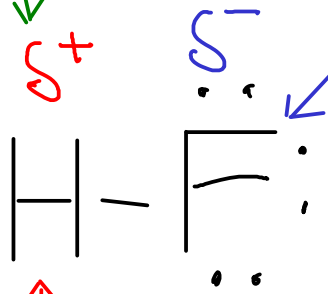
2.1

4.0

Electronegativity values
Difference = 1.9

Fluorine is much more electronegative than hydrogen. This is a POLAR BOND, and the shared electrons will be held more closely to FLUORINE!

δ means "small" here



This end of the molecule will have a slight NEGATIVE charge, since the shared electrons are closer to FLUORINE!

This end of the molecule will have a slight POSITIVE charge, since the shared electrons are pulled away from HYDROGEN!

POLARITY OF MOLECULES

So what can a molecule's LEWIS STRUCTURE, SHAPE, and the POLARITY of its bonds tell us?

... the POLARITY of the overall molecule, which will tell us (among other things) what a given molecule will mix with or dissolve in!

POLAR MOLECULES

- Will dissolve in or dissolve other polar molecules
- Will dissolve some ionic compounds
- Will NOT easily dissolve nonpolar molecules

Example:
WATER

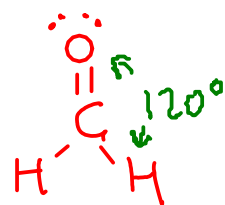
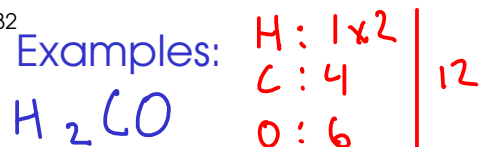
NONPOLAR MOLECULES

- Will dissolve in or dissolve other nonpolar molecules
- Will NOT easily dissolve polar molecules or ionic compounds

Example:
OILS

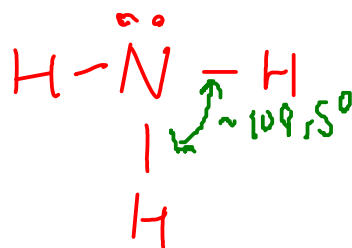
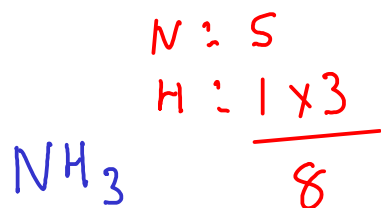
For a molecule to be polar, it must ...

- ① Have polar bonds! (Any molecule that contains no polar bonds must be nonpolar!)
- ② Have polar bonds arranged in such a way that they don't balance each other out! (This is why you need to know the structure and shape of the molecule)



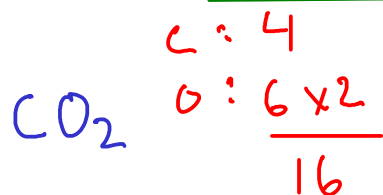
Shape? This is a TRIGONAL PLANAR molecule. There are THREE THINGS around the central atom: =O, -H, -H

Polar? C-H bonds are nonpolar, but the C=O bond should be polar. Electrons are pulled towards the oxygen and this pull is not "cancelled" by the opposing hydrogen atoms (they're not very electronegative). POLAR MOLECULE.



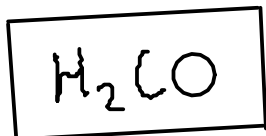
Shape? PYRAMIDAL. There are four things around the central nitrogen, but only three are atoms. Tetrahedral bond angles (109.5), but pyramidal shape.

Polar? N-H bonds are polar, and since the molecule has nitrogen "side" (the top of the pyramid), the molecule is POLAR. The nitrogen end is slightly negative, while the hydrogen end is slightly positive.

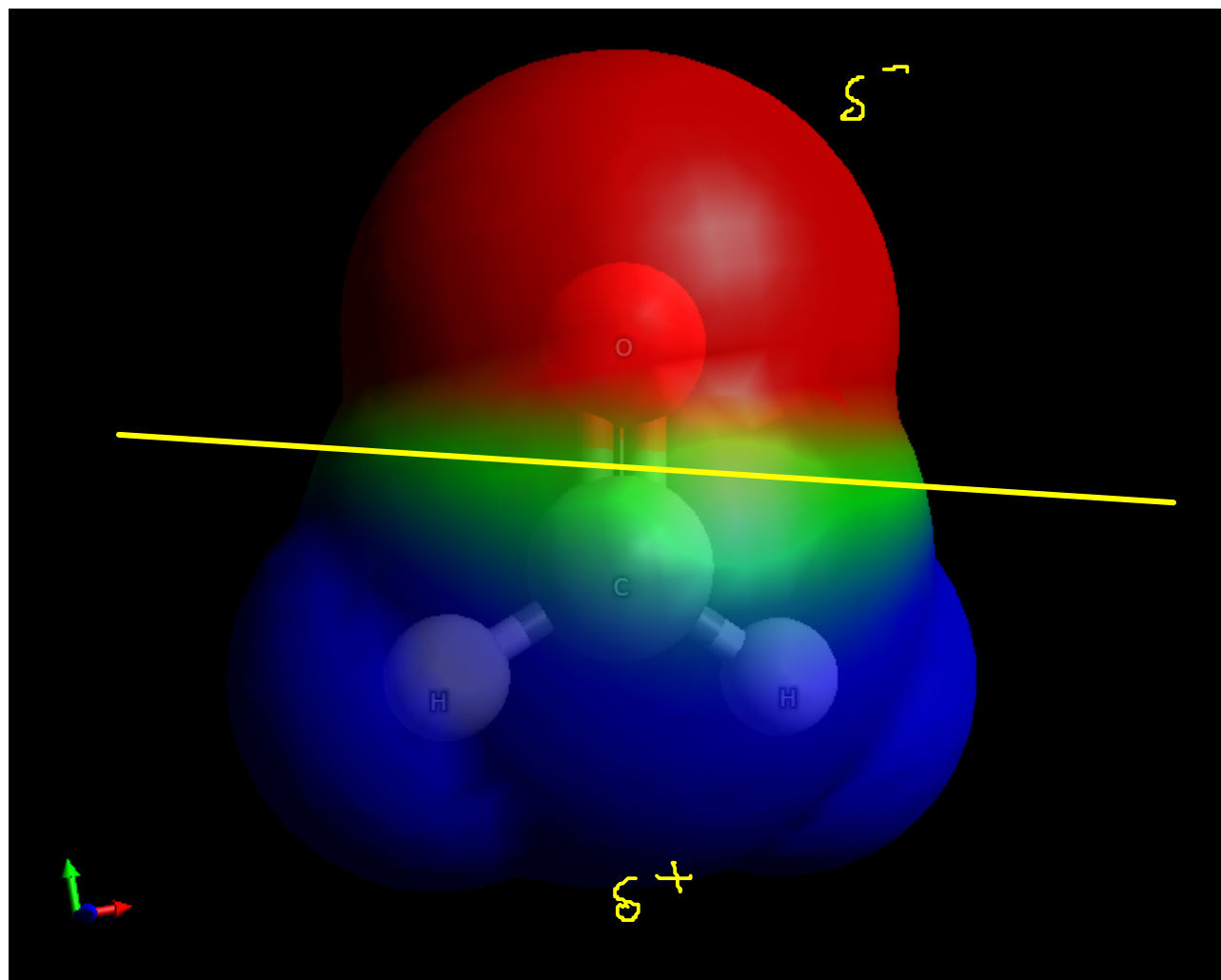


Shape: LINEAR. Only two groups around the central atom

Polarity: We expect C=O bonds to be polar, BUT they are directly opposite one another in this structure, so the overall molecule is NONPOLAR.

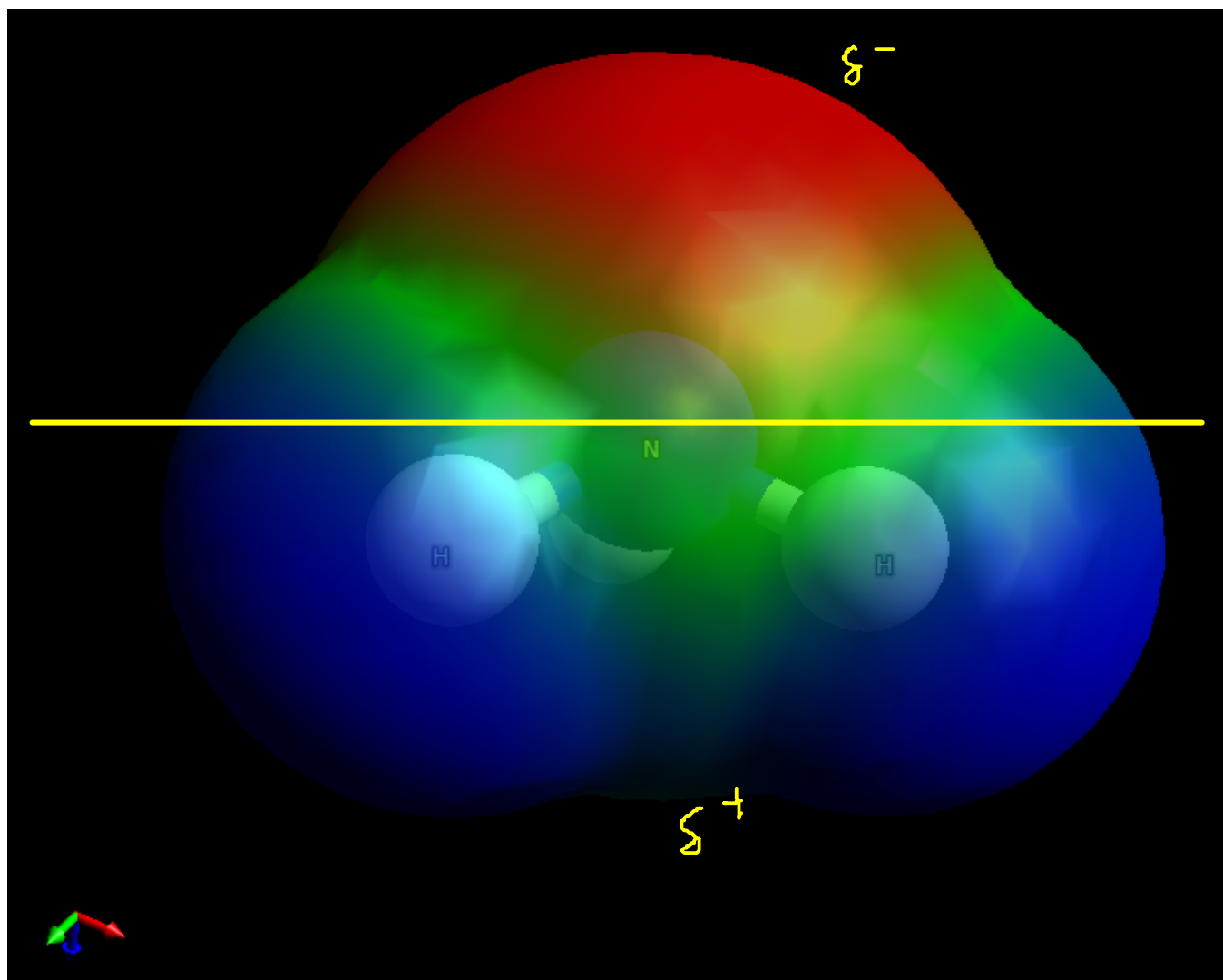
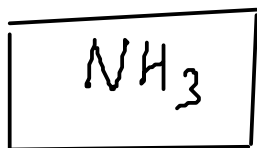


This ball-and-stick model shows electrostatic potential - red for more negative and blue for more positive



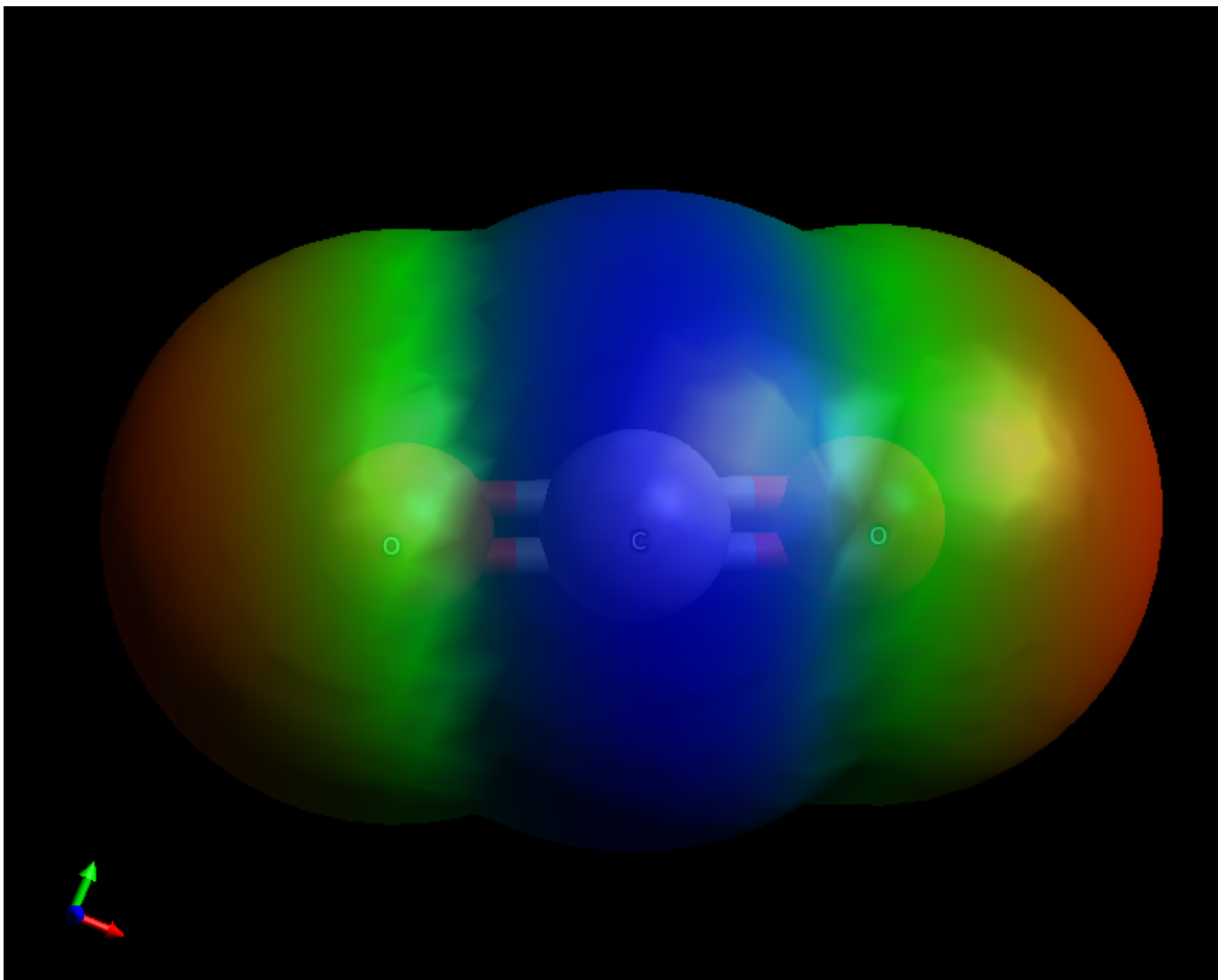
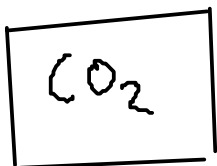
oxygen "side",
slightly negative

hydrogen "side",
slightly positive



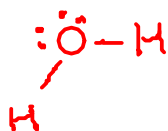
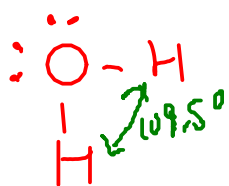
nitrogen "side"
slightly negative

hydrogen "side"
slightly positive



This molecule is NONPOLAR. No positive "side" or negative "side"

EXAMPLES

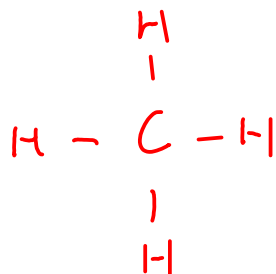
Water, H_2O H: 2×1 O: $\frac{6}{8}$ 

Shape: BENT. Four groups around the central oxygen, but only two are atoms.

Polarity? POLAR. O-H bonds are polar, and oxygen is able to pull electrons through to its "side" of the molecule

methane, CH_4

C: 4

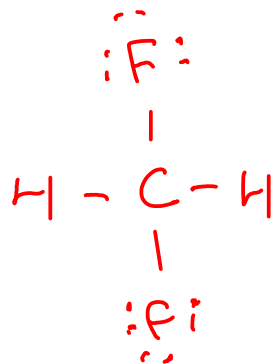
H: 4×1 $\frac{8}{8}$ 

Shape? TETRAHEDRAL. Four groups (all atoms) around central carbon

Polar? NONPOLAR, since we have only nonpolar bonds.

 CH_2F_2

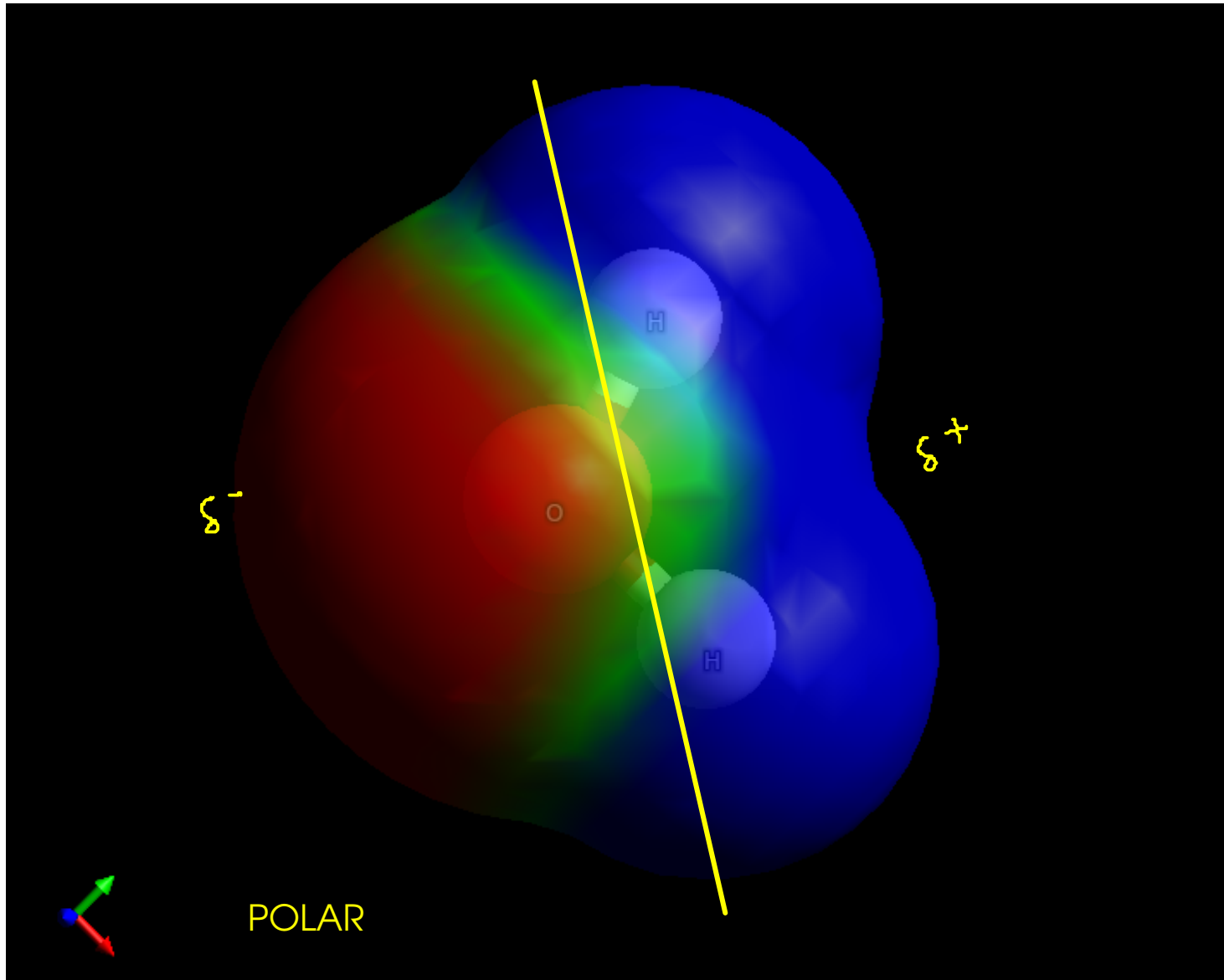
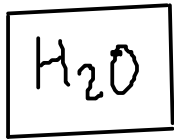
C: 4

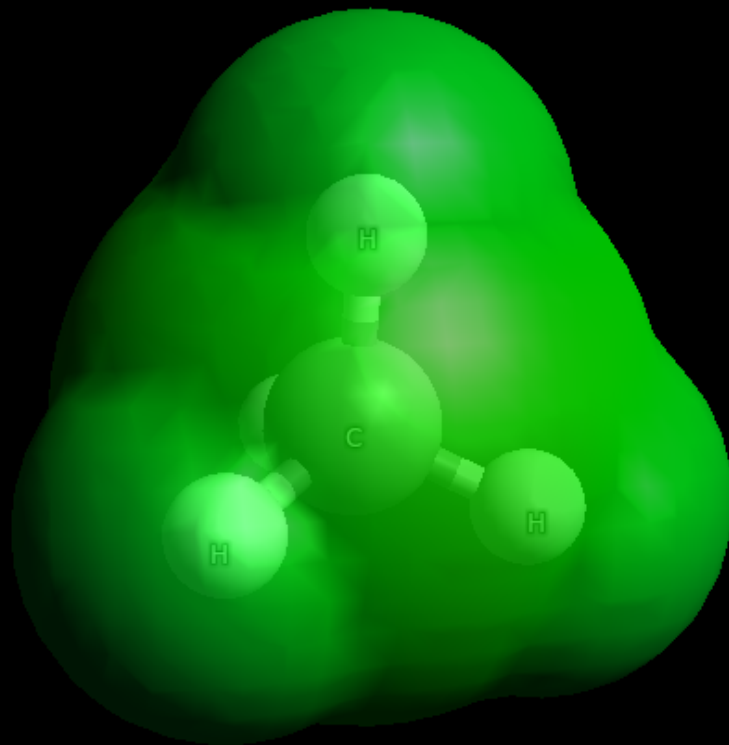
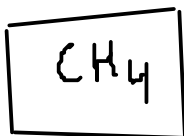
H: $2 \times 1 = 2$ F: $2 \times 7 = 14$ $\frac{20}{20}$ 

Shape? TETRAHEDRAL. Four groups (all atoms) around central carbon

Polar? C-H are nonpolar, but C-F bonds are polar. The molecule is polar, since electrons are pulled towards the fluorine side.

This Lewis structure is CORRECT, but deceptive, since it APPEARS that the fluorines are on opposite sides of the molecule.





NONPOLAR (all bonds are nonpolar)

