

Molecular Geometry

PURPOSE

- A To explore some simple molecular structures
- B To explore the relationship between bond order and bond length
- C To explore resonance structures

GOALS

- To compare Lewis structures to three-dimensional models
- To visualize the three-dimensional structures of some common molecules
- To obtain bond angle, bond length, and hybridization data for molecules
- To rationalize differences in predicted and measured values
- To learn how to use computer modeling software

INTRODUCTION

The chemical and physical properties of covalently bonded materials are related to the spatial arrangement of the atoms and other electrons not involved in the actual bond formation. There are many ways to depict the spatial arrangement in both two and three dimensions. A **Lewis structure** is a two-dimensional representation of the arrangement of the atoms, bonding electrons and non-bonding (lone pair) electrons in a covalent material.

In a Lewis structure, the nucleus is represented by the atomic symbol with a line between the atoms in a bond depicting each pair of shared bonding electrons in the structure. Non-bonding electrons around the atoms are depicted as dots. The steps to building a Lewis structure representation of a molecule are shown below. The Lewis structure of the formate ion, CHO_2^- , will be used as an example.

- 1 Calculate the **electrons required (ER)** = the minimum number of electrons necessary to satisfy the octet rule for the non-hydrogen atoms and the duet rule for hydrogen. For CHO_2^- , this would be (2 electrons \times 1 hydrogen atom) + (8 electrons \times 3 non-hydrogen atoms) = $2 + 24 = 26$ electrons required.
- 2 Calculate the number of available **valence electrons (VE)** = the total number of electrons available for the molecule. For example, in CHO_2^- , this would be (1 C atom \times 4 electrons) + (1 H atom \times 1 electron) + (2 O atoms \times 6 electrons) + (1 electron as the ion has a charge of -1) = $4 + 1 + 12 + 1 = 18$ valence electrons. NOTE: For ions, the charge must be included in this by adding the charge on an anion or subtracting the charge on a cation.
- 3 Calculate the **Shared Pairs (SP)** = the number of electrons to be shared in bonds. The SP is $1/2(\text{ER} - \text{VE})$; for CHO_2^- , this would be $1/2(26 - 18) = 4$ shared pairs or four bonds.

- 4 Calculate the **Lone Pairs (LP)** = the number of electron pairs belonging to only one atom. The LP is $1/2(\text{VE} - (2 \times \text{SP}))$; for CHO_2^- , this would be $1/2(18 - (2 \times 4)) = 5$ lone pairs. Notice that $\text{VE} = 2 \times (\text{SP} + \text{LP})$.
- 5 Place the first atom in the molecular formula as the central atom, surrounded by the other atoms in the compound.

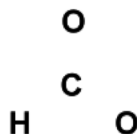


Figure 1

- 6 Draw bonds (shared pairs) from the central atom to each surrounding atom. The bonds are represented as lines; each line represents two electrons. The number of lines should be equal to the number of shared pairs calculated in step 3, which in this case is four. Since hydrogen follows the duet rule, it only prefers one bond. The fourth bond can be drawn to either one of the oxygen atoms.

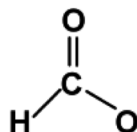


Figure 2

- 7 Draw lone pairs on each of the non-hydrogen atoms. A lone pair is represented as two dots; each dot represents an electron. Each non-hydrogen atom prefers eight electrons in the vicinity of the atom. If an atom has 1 bond, it requires 3 lone pairs. If an atom has 2 bonds, it requires 2 lone pairs. If an atom has 3 bonds, it requires 1 lone pair. If an atom has 4 bonds, do not add any additional lone pairs. In our example, C requires no lone pairs, one oxygen requires 3 lone pairs and one oxygen requires 2 lone pairs.

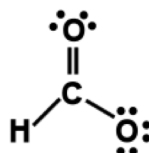


Figure 3

8 Calculate the **formal charge (FC)** on each atom. For an atom:

$$FC = VE - SP - (2 \times LP) \quad (1)$$

where SP is the number of shared pairs on the atom and LP is the number of lone pairs on the atom. For molecules the sum of the formal charges of all the atoms must be zero; for an ion, the sum will be the ionic charge. To choose the more favorable of two Lewis structures, the one with the lowest FC on each individual atom would be the most likely candidate. Remember that the more electronegative atom will generally prefer the negative formal charge. The formal charge for each atom on the formate ion is shown below. Note how the sum of the formal charges ($0 + 0 + 0 + -1$) adds up to the -1 charge on the formate ion.

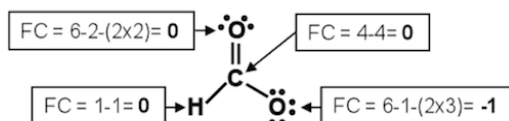


Figure 4

The structure shown above is the Lewis structure representation of the formate ion. Note that the structure could also have been drawn with two bonds directed towards the oxygen on the right. This is known as a resonance structure. Resonance structures have the same number of electrons, but may have a different number of shared pairs or the atoms in the structure may have different formal charges.

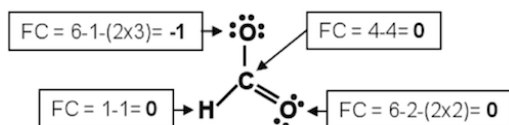


Figure 5

Once a Lewis structure has been drawn, one can evaluate the **bond order** of shared pairs. A shared pair is a single bond with a bond order of 1; two shared pairs in a bond is a double bond (bond order = 2) and three shared pairs in a bond is a triple bond (bond order = 3). In general, the bond strength increases and bond length decreases as the bond order increases. (The bond length is the distance between the centers of the two atoms bonded together). Thus, **a single bond is longer than a triple bond, but the triple bond is stronger**. NOTE: This general statement applies when comparing the bond order of similarly bonded atoms.

However, as noted, the Lewis structure is a two-dimensional representation of the molecule or

ion. The two-dimensional diagram does not show us the geometry or shape of the molecule. Over the years, many theories have attempted to explain the shape of covalently bonded substances; one of the simplest is the **Valence Shell Electron Pair Repulsion** (VSEPR) theory. In most molecules, electrons occur in pairs, either as part of a bond (bonding pair) or as a lone pair; these pairs occupy space around an atom. The theory proposes that these areas or electron regions repel one another. VSEPR theory's main postulate is that the regions around a given atom will arrange themselves to minimize this repulsion by positioning themselves as far apart as possible. This leads to a predictable shape based on simple geometry.

For the purposes of the VSEPR theory, it is more helpful to think about electron groups (rather than just electron pairs) and the regions the groups occupy in space. A lone pair or bonding pair are each considered to be an electron group. However, the four electrons in a double bond or the six electrons in a triple bond are also considered to be **one** electron group. In each case they occupy one region where they are bonding two atoms together.

The position of the electron groups determines the angles at which centers of those electron regions are located. As a result, the angle between atoms that may be at the ends of those electron groups is also determined. These angles for covalent compounds that obey the octet rule are shown in the table below. The values in the table give the angles between the centers of the electron regions. If these are bonding groups (single, double or triple bonds), this number also is the **bond angle**.

# of electron regions (# of electron groups)	angle formed (bond angle)	hybrid orbitals	# of non-central atoms	molecular shape
2	180°	sp	2	linear
3	120°	sp ²	3 2	trigonal planar bent
4	109°	sp ³	4 3 2	tetrahedral trigonal pyramidal bent

Figure 6

Also shown in the above table are the **hybridized orbitals** used to form the σ -bonds (sigma bonds) and to hold the lone pairs. Chemists invoke hybrid orbitals based on the assumption that one atom will point its electron region directly at an atom to which it bonds. s orbitals do not point in any direction, and simple p atomic orbitals lie 90° apart along the x , y and z -axes. Regions in space that point 109°, 120° or 180° apart result by taking mathematical combinations (hybrids) of the simple atomic orbitals. Hybrid orbitals have thus been incorporated into the language that describes covalent bonding.

When we use the term molecular geometry or molecular shape, we are not describing the shape of the electron regions, but rather, the location of the atoms. The words used to describe the shapes are therefore describing the location of the atoms. Where four atoms surround a central atom, the shape would be **tetrahedral**. Where three atoms surround a central atom, the shape would either be **trigonal planar** or **pyramidal**. Where two atoms surround a central atom, the shape would either be **linear** or **bent**.

It is difficult to represent three-dimensional structures, like molecules, on a flat piece of paper.

It is equally as difficult to look at a molecular structure drawn on a piece of paper and imagine what it would look like in three-dimensions. With the help of a molecular model kit and a computer modeling program, you will be able to visualize a molecule in three-dimensions. In this lab, you will use a computer program within WebAssign that allows molecules to be rotated, just like you could manually rotate a model built with a model kit. You will also be able to use the computer program to obtain experimentally determined measurements, such as bond angles and bond lengths, for the molecule under investigation.

In introductory chemistry courses, we often predict bond angles and bond lengths in simple molecules based on VSEPR or hybridization. However, our predictions are simply approximations because other factors influence the structure of real molecules. In this lab, you will compare the bond angles and bond lengths predicted from theory to the experimentally determined values. You will then be asked to explain differences that occur between the experimental values and the theoretical values.

EQUIPMENT

- 1 Chem-Tutor model kit
- 1 computer with internet browser
- 1 Molecular Geometry In-Lab assignment in WebAssign

About the model kit

The Chem-Tutor model kit was designed by Professor Samuel G. Levine for use by his organic chemistry students at North Carolina State University.

Atoms: Atoms are represented by jacks that have one, two, three or four pegs coming out. Each peg represents an electron region (or electron group). The electron regions (groups) can be single bonds, double bonds, triple bonds or lone pairs. In general, black jacks are used for carbon, blue for nitrogen, red for oxygen, white for hydrogen and yellow for sulfur. More important than color is the number of pegs. **The type of atom/jack to use is determined by the Lewis structure of the molecule.**

Bonds: Green bonds can be used for single, double or triple bonds. Do not bend the green bonds to force a molecule into a particular shape. Although you are strong enough to bend these bonds, it takes energy for atoms and molecules to bend away from the predicted angles and shapes. The molecules prefer to exist in the lowest energy configurations (“straight” green bonds).

Black bonds between two carbon atoms can be used for carbon-carbon double bonds. You will notice that the pegs on each carbon define a plane, and that the two carbons are coplanar. The atoms at the ends of the bond are **fixed in place**. **DO NOT** attempt to twist this bond.

White bonds are made out of rubberized plastic, and are the only ones that are meant to bend. They are only used for special “strained” structures that you should not encounter in this lab.

About the computer modeling software

The molecules in your WebAssign In-Lab are viewed with jmol, a javascript plug-in. The computer files viewed with jmol contain real data from measurements on these molecules. Often

we find that molecules follow what is predicted from textbooks. Occasionally, they do not. Do not be surprised if what you observe is sometimes different than what you might predict!

Rotating a molecule: A molecule in the viewing area can be rotated in all directions. This can be achieved by holding down the left mouse button while moving the mouse. After a few tries at this, you should be able to manipulate the molecule any way that you want.

To determine a bond length: Double click on one of the atoms in the bond. The cursor will change to a "+". Then double click on the other atom in the bond. The measurement will be displayed in nanometers (nm). To make the measurement disappear, re-measure the bond in the same way.

To determine a bond angle: Double click on one of the atoms of the angle. The cursor will change to a "+". Then **single** click on the vertex of the angle. And last, double click on the third atom in the angle. The measurement will be displayed in degrees. To make the measurement disappear, re-measure the angle in the same way.

REAGENTS

No chemicals will be used in this experiment.

SAFETY

None of the materials being used in this experiment present a safety hazard. However, the work is being done in a laboratory and the usual rules about eye protection and proper clothing apply.

WASTE DISPOSAL

Since no chemicals are being used in this experiment, there will not be any waste for disposal.

LAB PROCEDURE

You will notice the format of this lab experiment is different than other experiments. You will build a series of models and investigate them on the computer. Questions to help you with your observations are intermingled with the procedure. Please answer the questions in your lab manual along with any other observations you make while you are building the structures.

- Launch Internet Explorer.
- Open one partner's Molecular Geometry In-Lab in WebAssign.

Please print the worksheet for this lab. You will need this sheet to record your data.

Part A: Exploring Simple Structures

For each of the following molecules,

- Review the correct, complete Lewis structure(s), including any resonance structures and any formal charges that you drew in your Prelab assignment.
- Build the molecule with the model set.
- Look at the molecule in your In-Lab assignment on WebAssign.

- Fill in the table and answer the questions below.

1 Two Electron Regions

Dinitrogen Oxide (N₂O) Hint: Nitrogen is the central atom.

2 Three Electron Regions

- a **Sulfur Dioxide (SO₂)**
- b **Formaldehyde (CH₂O)**

3 Four Electron Regions

- a **Water (H₂O)**
- b **Ammonia (NH₃)**
- c **Methane (CH₄)**

Part B: Bond Order vs Bond Length

For each of the following molecules,

- Review the correct, complete Lewis structure(s), including any resonance structures and any formal charges that you drew in your Prelab assignment.
- Build the molecule with the model set.
- Look at the molecule in your In-Lab assignment on WebAssign.
- Fill in the table and answer the questions below.

1 **Ethane (C₂H₆)**

2 **Ethene (C₂H₄)**

3 **Ethyne (C₂H₂)**

Part C: Resonance Structures

Draw the Lewis structures for Part C on your lab worksheet.

For each of the following molecules,

- Review the correct, complete Lewis structure(s), including any resonance structures and any formal charges that you drew in your Prelab assignment.
- Build the molecule with the model set.
- Look at the molecule in your In-Lab assignment on WebAssign.
- Fill in the table and answer the questions below.

1 **Benzene (C₆H₆)** Hint: Place the carbons in a hexagon.

2 Carbonate Ion (CO_3^{2-})

3 Thiocyanate Ion (SCN^-) Hint: Carbon is the central atom.

- When you are done, clean up any model set structures and leave the sets as you found them when you arrived in the lab.
- Before leaving, go to a computer in the laboratory and enter your results in the In-Lab assignment. If all results are scored as correct, log out. If not all results are correct, try to find the error or consult with your lab instructor. When all results are correct, note them and log out of WebAssign. The In-Lab assignment must be completed by the end of the lab period. If additional time is required, please consult with your lab instructor.