Chem 103, Section F0F Unit VI - Compounds Part II: Covalent Compounds Lecture 17

- Using the Valence-Shell Electron-Pair Repulsion (VSEPR) Theory to predict molecular shapes
- Molecular shape and polarity

Lecture 17 - Covalent Bonding

Reading in Silberberg

- Chapter 10, Section 2
 - Valance-Shell Electron-Pair Repulsion (VSEPR) Theory and Molecular Shape
- Chapter 10, Section 3
 - Molecular Shape and Molecular Polarity

Lecture 17 - Introduction

In this lecture we will look into predicting molecular shapes and polarity

- Using a simply theory called the Valence-Shell Electron-Pair Repulsion (VSEPR) Theory it is possible to predict the shape of a molecule starting with its Lewis structure.
- The molecular structure can then be coupled with electronegativity to predict molecular polarity.
- Molecular shape and polarity have a marked influence on a molecule's physical properties.
 - This is particularly the case for biological molecules.

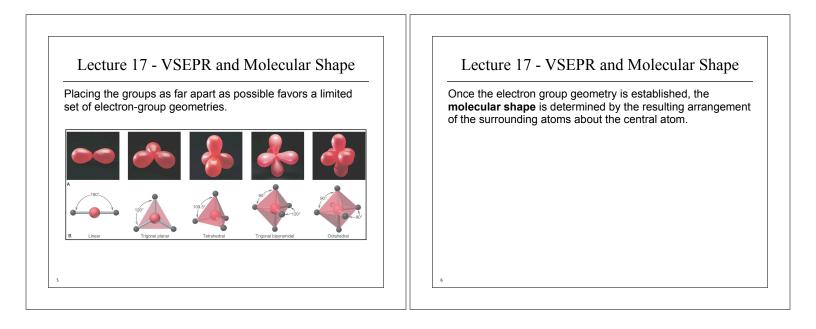
Lecture 17 - VSEPR and Molecular Shape

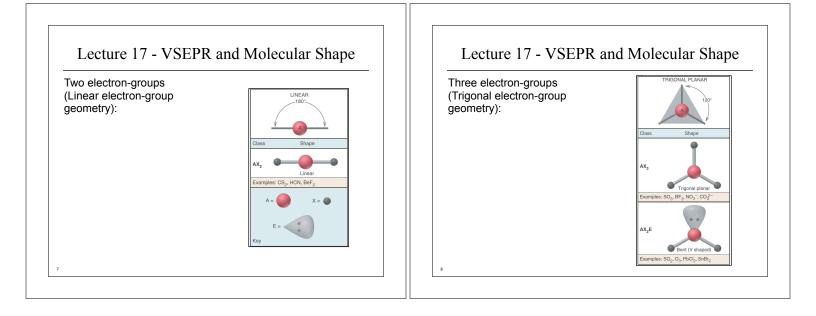
The main principle of VSEPR is

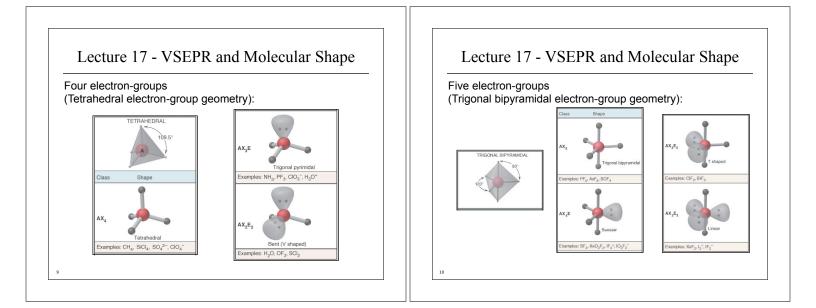
• Each group of valence electrons (VSEP) around a central atoms is located as far away as possible from the other groups (R).

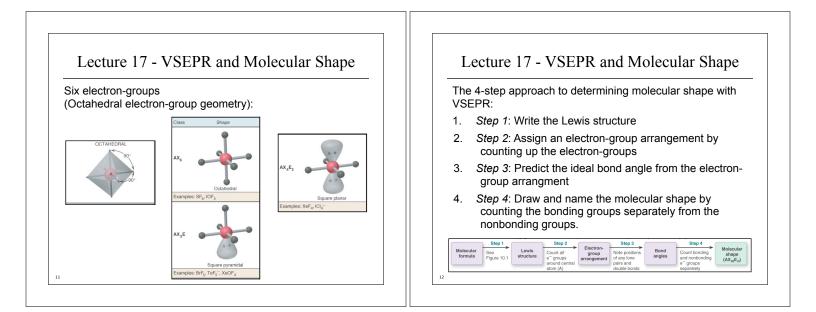
A "group" is any number of electrons that are located near to one another:

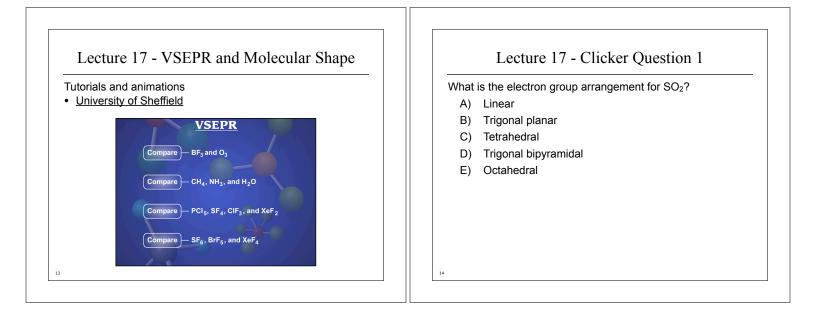
- A non bonded pair of electrons (2 e-'s)
- A single bond (2 e⁻'s)
- A double bond (4 e⁻'s)
- A triple bond (6 e⁻'s)
- A lone free radical electron (1 e^{-})

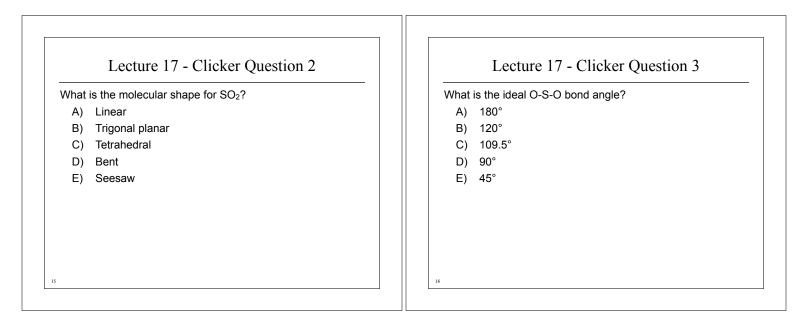






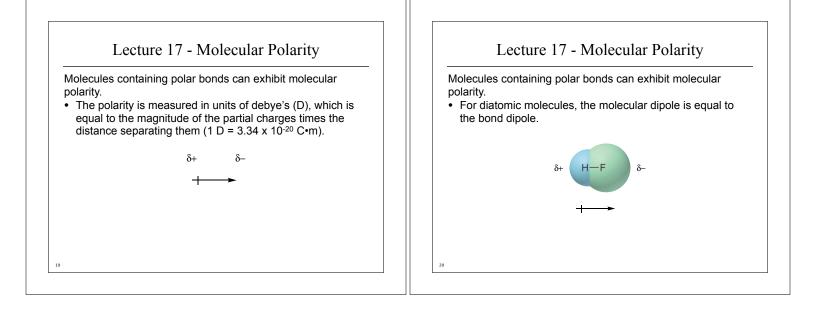


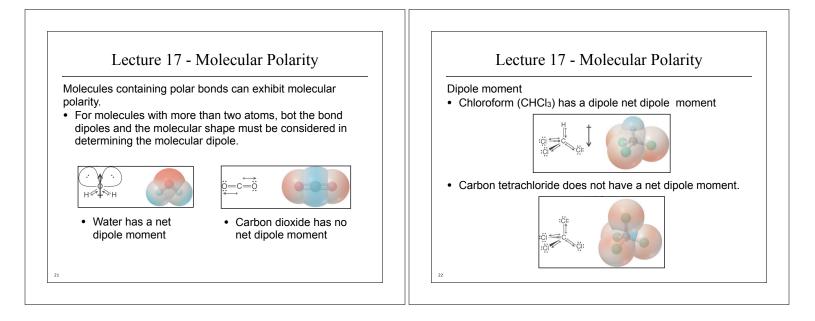


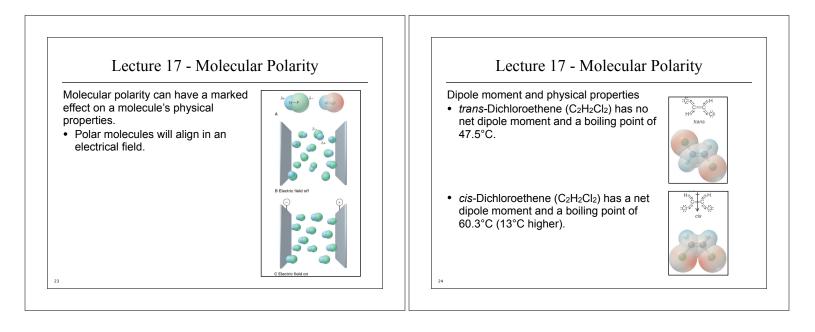


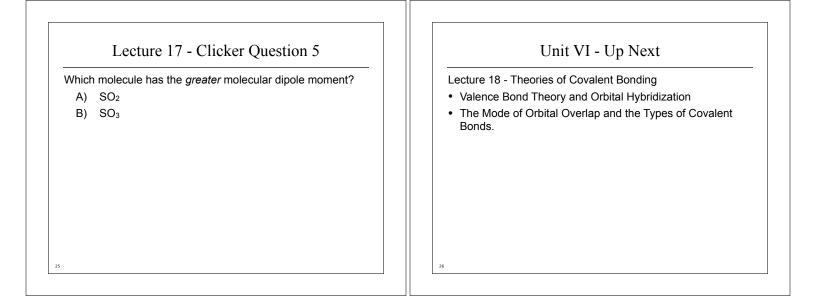
Lecture 17 - Question 4	Lecture 17 - VSEPR and Molecular Shape
Determine the electron-group arrangement, molecular shape, and ideal bond angle for N ₂ O (N is central)? Using formal charges, which resonance structure for N ₂ O is the most probable?	When molecules have more than one central atom, each is treated separately.
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