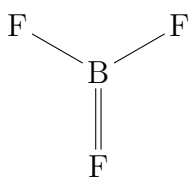


Chapter 7 – Covalent Bonds

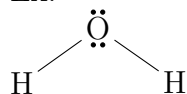
Michael Brodskiy

Instructor: Mr. Morgan

November 19, 2020

- Lewis Structures:
 1. Sum Valence Electrons
 2. Connect atoms with lines
 3. Arrange remaining e^- to satisfy octet rule
- Resonance – Being able to draw more than one Lewis structure
- Halogens will almost never have double bonds
- Molecular Shapes:
 1. Linear
 - Bond angle equals 180°
 - Usually non-polar
 - Ex:
 $\ddot{\text{O}}=\text{C}=\ddot{\text{O}}$
 2. Triangular Planar
 - Bond angle equals 120°
 - Usually non-polar
 - Ex:

 3. Bent
 - Bond angle equals 109.5°
 - Appears as a linear, but bent
 - Always polar

– Ex:



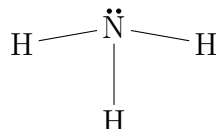
4. Tri-Pyramid

– Bond angle equals 109.5°

– Appears as a three dimensional, triangular pyramid

– Always polar

– Ex:



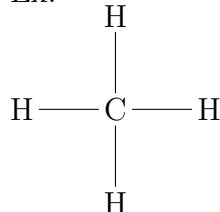
5. Tetrahedral

– Bond angle equals 109.5°

– Appears as a three dimensional, square pyramid

– Usually non-polar

– Ex:



• Expanded Octets: $5e^-$ pairs

1. Tri-Bipyramid (5 atoms)

(a) Usually non-polar

2. Seesaw (4 atoms)

(a) Always polar

3. T Shape (3 atoms)

(a) Always polar

4. Linear (2 atoms)

(a) Usually non-polar

• Expanded Octets: $6e^-$ pairs

– Octahedral (6 atoms)

1. Usually non-polar

– Square Pyramid (5 atoms)

1. Always polar

– Square Planar (4 atoms)

1. Usually non-polar

- VESPR – e^- pairs repel, and double and triple bonds count as one pair.
- Molecular Polarity – The negative side is what repels branching atoms, while the branching atoms are attracted to the positive side.
- Exceptions – NF_3 has a bond angle of 102° , CF_4 has a bond angle of 109.5° . This is because lone pairs repel more than bonding pairs.
- These types of graphs show the appropriate bond distance between two atoms:

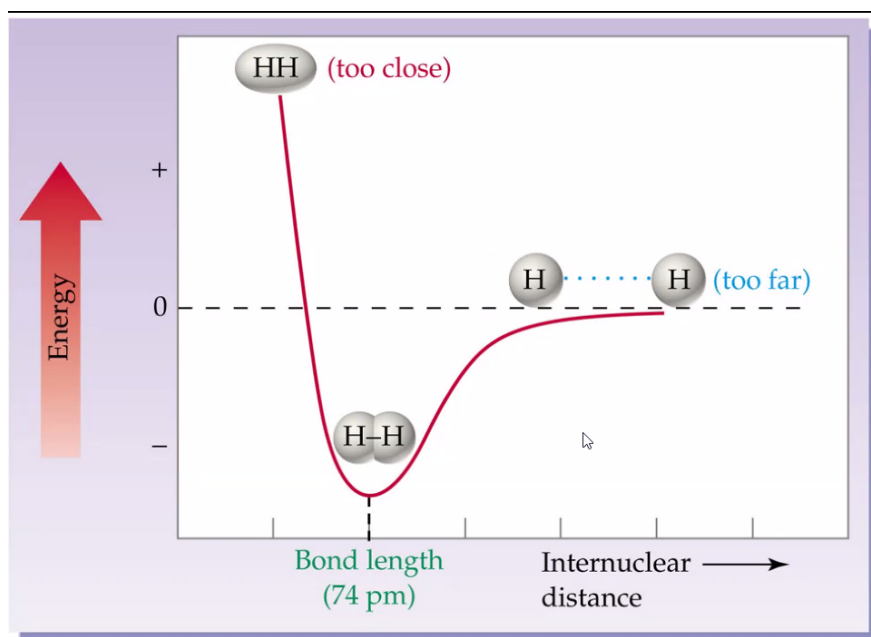


Figure 1: Best Bonding Distance For H_2

- The AP Exam will show a graph like the following and ask what is incorrect. Horizontal is determined by protons, while vertical is determined by size of the atom. In this case, F_2 should be left of I_2

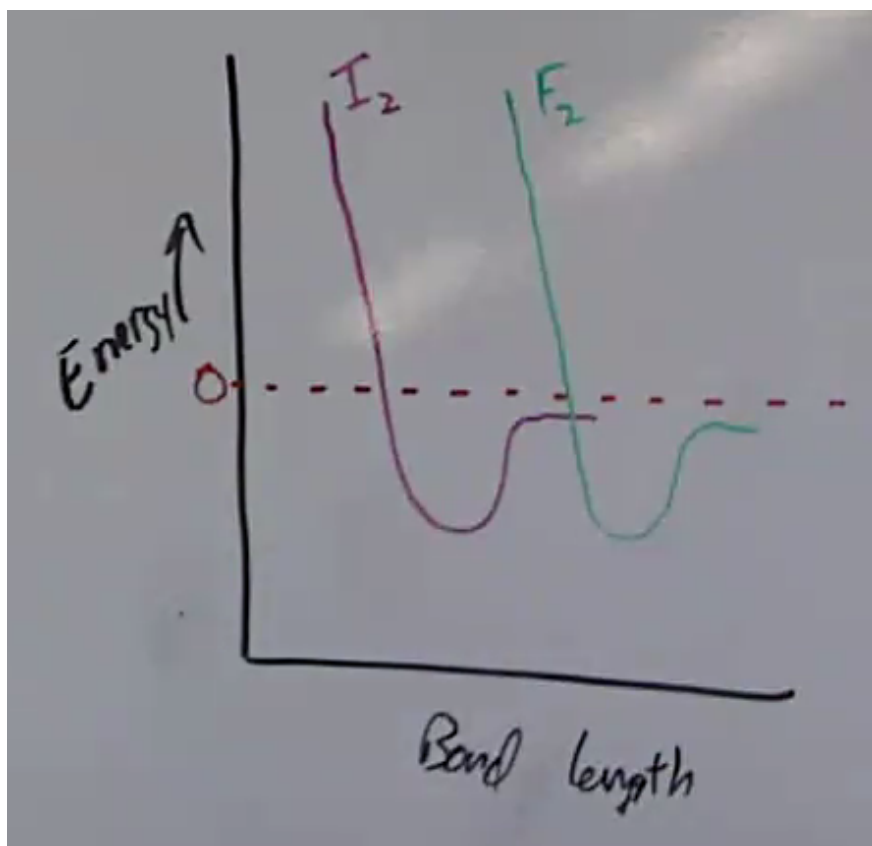


Figure 2: Incorrect Energy Required to Break I_2 and F_2 Bonds