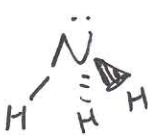
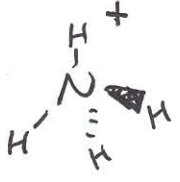
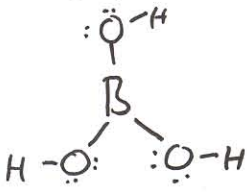
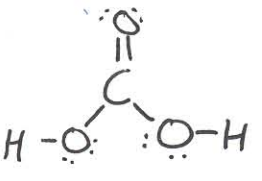

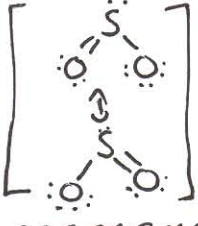
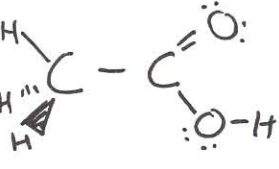
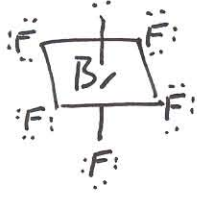
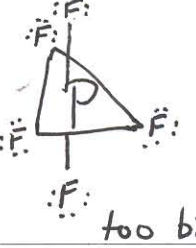
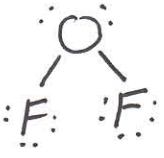
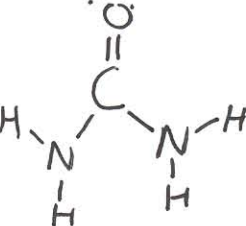
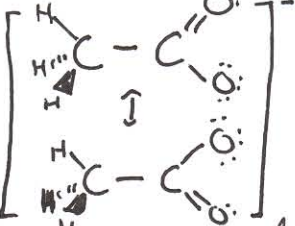
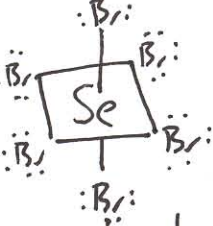


### Fall 2006 CH301 Worksheet 4

In each box, write the total number of electrons, the number of electrons around the perimeter atoms, and the 3D Lewis structure for the given compound. Identify multiple bonds or resonance. Indicate compounds that do not follow octet/duplet structures and the reason (too big, too small, too odd).

<p>CO<sub>2</sub> 16 total, 16 perimeter</p> <p style="text-align: center;"><math>\text{:O}=\text{C}=\text{O:}</math></p> <p>2 double bonds</p>	<p>O<sub>3</sub> 18 total, 16 perimeter</p> <p style="text-align: center;">  resonance         </p>	<p>NO<sub>2</sub><sup>-</sup> 18 total, 16 perimeter</p> <p style="text-align: center;"> </p>	<p>PCl<sub>5</sub> 40 total, 40 perimeter</p> <p style="text-align: center;">  too big         </p>
<p>SF<sub>6</sub> 48 total, 48 perimeter</p> <p style="text-align: center;">  too big         </p>	<p>CH<sub>2</sub>Cl<sub>2</sub> 20 total, 20 perimeter</p> <p style="text-align: center;"> </p>	<p>CH<sub>3</sub>CHO 18 total, 16 perimeter</p> <p style="text-align: center;">  multiple central atoms         </p>	<p>NO 11 total, 11 perimeter</p> <p style="text-align: center;"><math>\text{:N}=\text{O:}</math></p> <p>too odd</p>
<p>CO 10 total, 8 perimeter</p> <p style="text-align: center;"><math>\text{:C}\equiv\text{O:}</math></p> <p>triple bond</p>	<p>BaO  <math>\text{Ba}^{+2}, \text{O}^{2-}</math></p> <p>ionic</p>	<p>H<sub>2</sub>O 8 total, 4 perimeter</p> <p style="text-align: center;"> </p>	<p>H<sub>2</sub>O<sub>2</sub> 14 total, 4 perimeter</p> <p style="text-align: center;">  multiple central atoms         </p>
<p>I<sub>3</sub><sup>-</sup> 22 total, 16 perimeter</p> <p style="text-align: center;">  too big         </p>	<p>BF<sub>3</sub> 24 total, 24 perimeter</p> <p style="text-align: center;">  too small         </p>	<p>HCN 10 total, 10 perimeter</p> <p style="text-align: center;"><math>\text{H}-\text{C}\equiv\text{N:}</math></p> <p>triple bond</p>	<p>SiCl<sub>4</sub> 32 total, 32 perimeter</p> <p style="text-align: center;"> </p>
<p>O<sub>2</sub> 12 total, 8 perimeter</p> <p style="text-align: center;"><math>\text{:O}=\text{O:}</math></p> <p>double bond</p>	<p>CH<sub>2</sub>CHOH 11 total, 8 perimeter</p> <p style="text-align: center;">  multiple central atoms         </p>	<p>NaCl  <math>\text{Na}^+, \text{Cl}^-</math></p> <p>ionic</p>	<p>SOCl<sub>2</sub> (S is the central atom) 26 total, 24 perimeter</p> <p style="text-align: center;"> </p>

<p>NH<sub>3</sub> 8 total, 6 perimeter</p> 	<p>NH<sub>4</sub><sup>+</sup> 8 total, 8 perimeter</p> 	<p>N<sub>2</sub> 14 total, 8 perimeter</p> <p>:N≡N:</p> <p>triple bond</p>	<p>B(OH)<sub>3</sub> 24 total, 6 perimeter</p>  <p>too small</p>
<p>H<sub>2</sub>CO<sub>3</sub> 24 total, 12 perimeter</p>  <p>multiple central atom</p>	<p>Br<sub>2</sub>O 20 total, 16 perimeter</p> 	<p>KCl</p> <p>K<sup>+</sup>, Cl<sup>-</sup></p> <p>ionic</p>	<p>SO<sub>2</sub> 18 total, 16 perimeter</p>  <p>resonance</p>
<p>CH<sub>3</sub>COOH 24 total, 16 perimeter</p>  <p>multiple central atom</p>	<p>BrF<sub>5</sub> 42 total, 40 perimeter</p>  <p>too big</p>	<p>K<sub>2</sub>S</p> <p>2K<sup>+</sup>, S<sup>-2</sup></p> <p>ionic</p>	<p>Mg<sub>3</sub>N<sub>2</sub></p> <p>3Mg<sup>+2</sup>, 2N<sup>-3</sup></p> <p>ionic</p>
<p>PF<sub>5</sub> 40 total, 40 perimeter</p>  <p>too big</p>	<p>OF<sub>2</sub> 20 total, 16 perimeter</p> 	<p>ClF 14 total, 14 perimeter</p> <p>:Cl-F:</p>	<p>HBr 8 total, 8 perimeter</p> <p>H-Br:</p>
<p>(NH<sub>2</sub>)<sub>2</sub>CO 24 total, 16 perimeter</p>  <p>multiple central atom</p>	<p>CNO<sup>-</sup> (C is the central atom) 16 total, 16 perimeter</p> <p>:N=C=O<sup>-</sup></p> <p>2 double bonds</p>	<p>CH<sub>3</sub>COO<sup>-</sup> 24 total, 22 perimeter</p>  <p>multiple central atom</p>	<p>SeBr<sub>6</sub> 48 total, 48 perimeter</p>  <p>too big</p>