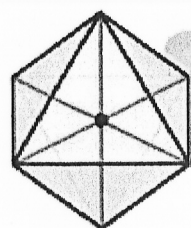
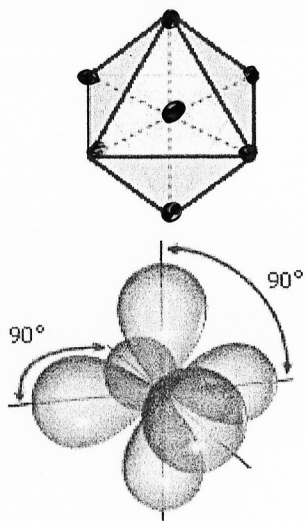


Molecular Geometry = Electron Pair Geometry, *ignoring lone pairs*

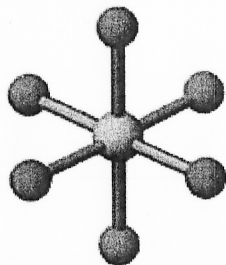
6

Octahedral



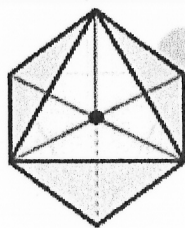
SF_6

6 bond pairs
No lone pairs



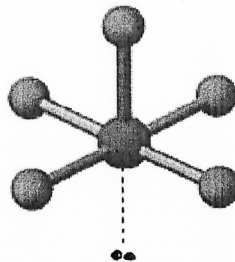
mg
Octahedral

SIX ELECTRON PAIRS
Electron-Pair Geometry = octahedron

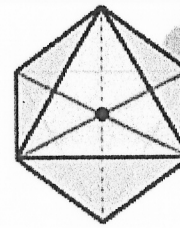


BrF_5

5 bond pairs
1 lone pair

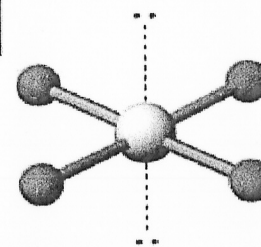


mg
Square-pyramidal



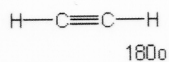
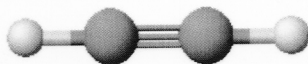
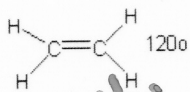
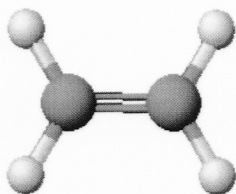
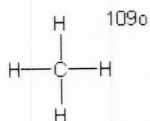
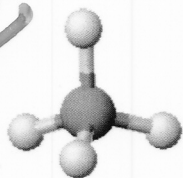
XeF_4

4 bond pairs
2 lone pairs

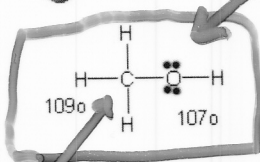
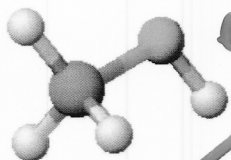


mg
Square-planar

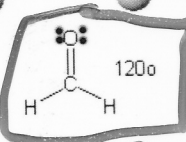
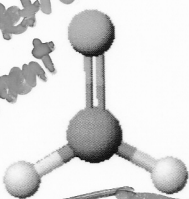
Methane



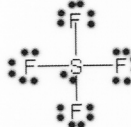
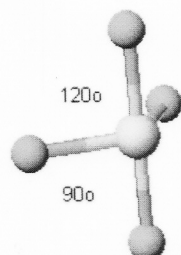
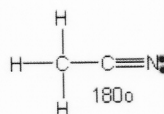
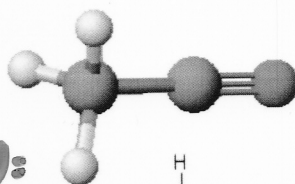
Methanol



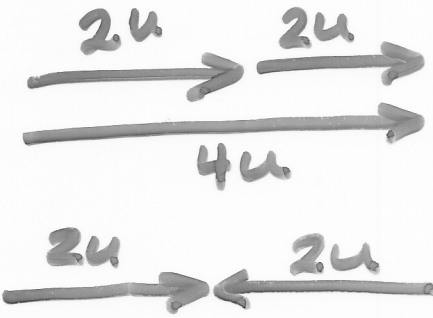
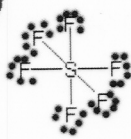
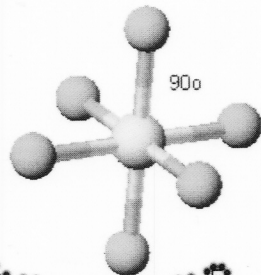
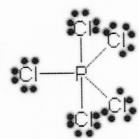
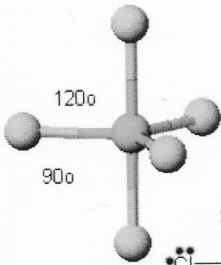
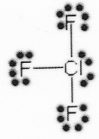
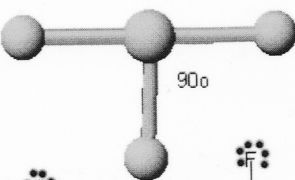
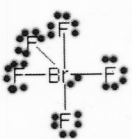
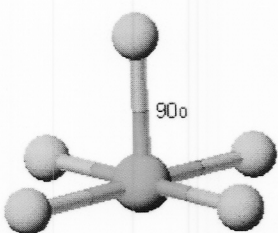
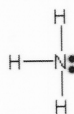
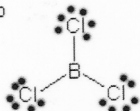
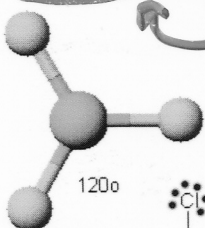
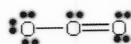
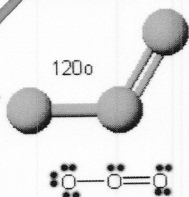
eg: tetrah.
mg: bent



mg: trig. planer



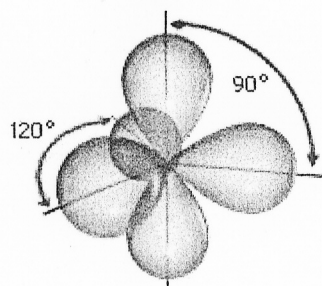
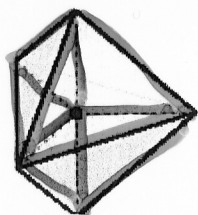
C
eg: tetrah.
mg: tetrah.



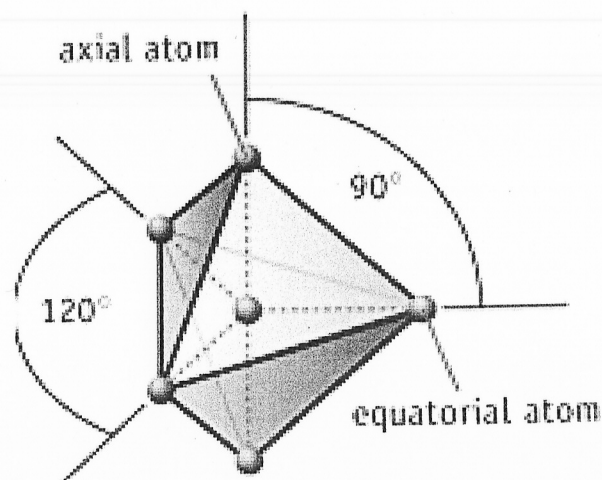
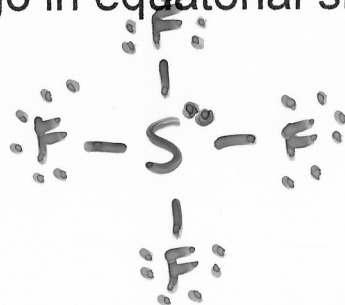
Molecular Geometry = Electron Pair Geometry, *ignoring lone pairs*

5

Trigonal-bipyramidal

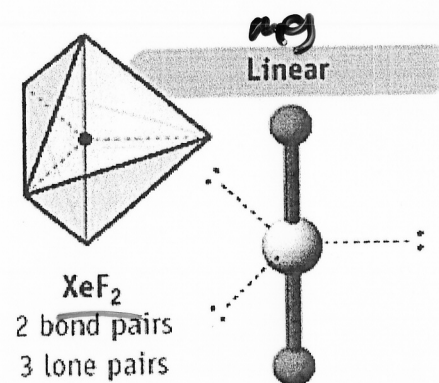
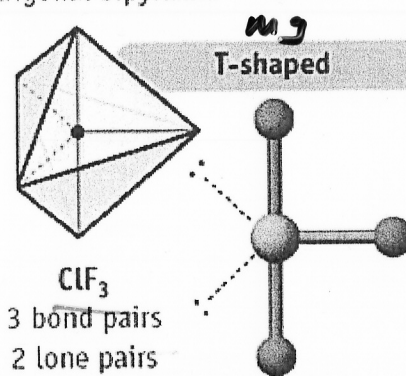
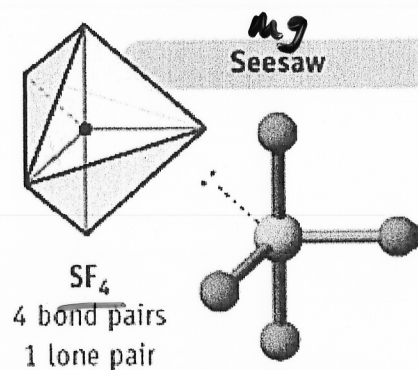
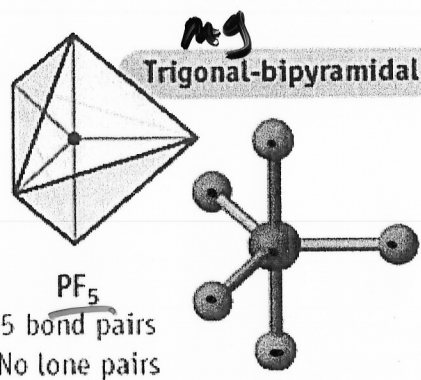


Lone pairs always
go in equatorial sites



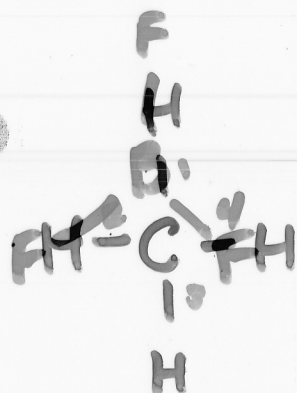
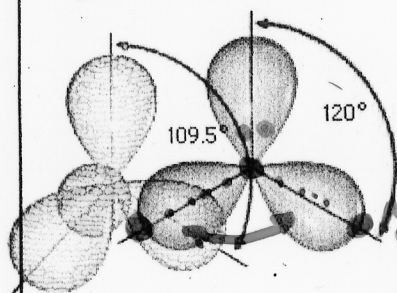
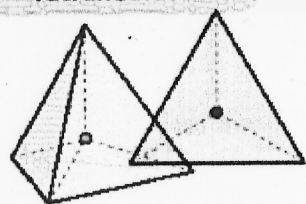
FIVE ELECTRON PAIRS

Electron-Pair Geometry = trigonal bipyramid



Molecular Geometry from Electron Pair Geometry

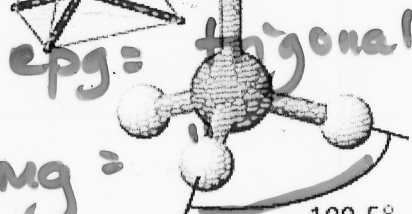
4
3
Trigonal-planar
Tetrahedral



epg: tetrahedral

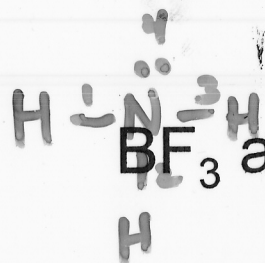
Tetrahedral

$\# \text{ sp} = 3$



mg =
angle = 109.5°
Methane, CH_4
4 bond pairs
no lone pairs

mg =

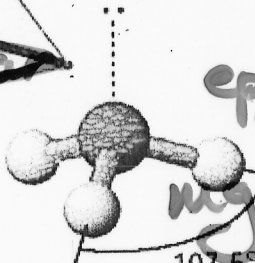


BF_3 and O_3



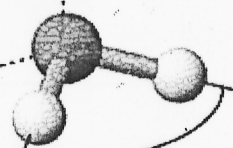
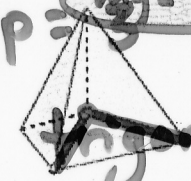
FOUR ELECTRON PAIRS
Electron Pair Geometry = tetrahedral

mg: Trigonal-pyramidal



Ammonia, NH_3
3 bond pairs
1 lone pair

mg: Bent



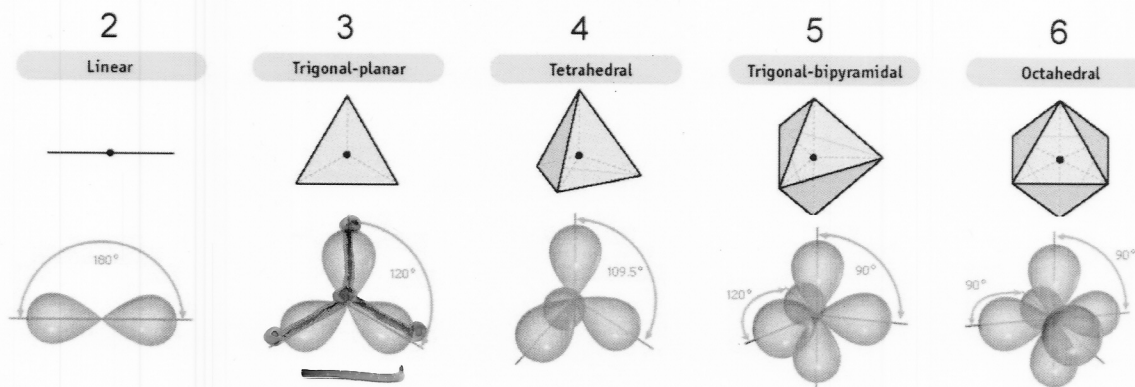
Water, H_2O
2 bond pairs
2 lone pairs
angle = close to 120°
slightly less than 120°

Molecular Structure

Ideal Electron Repulsion Shapes

Number of Pairs*	Electron-Pair Geometry	Bond Angles
2	linear	180°
3	trigonal planar	120°
4	tetrahedral	109.5°
5	trigonal bipyramidal	90° and 120°
6	octahedral	90°

*Add number of lone pairs and bonded atoms



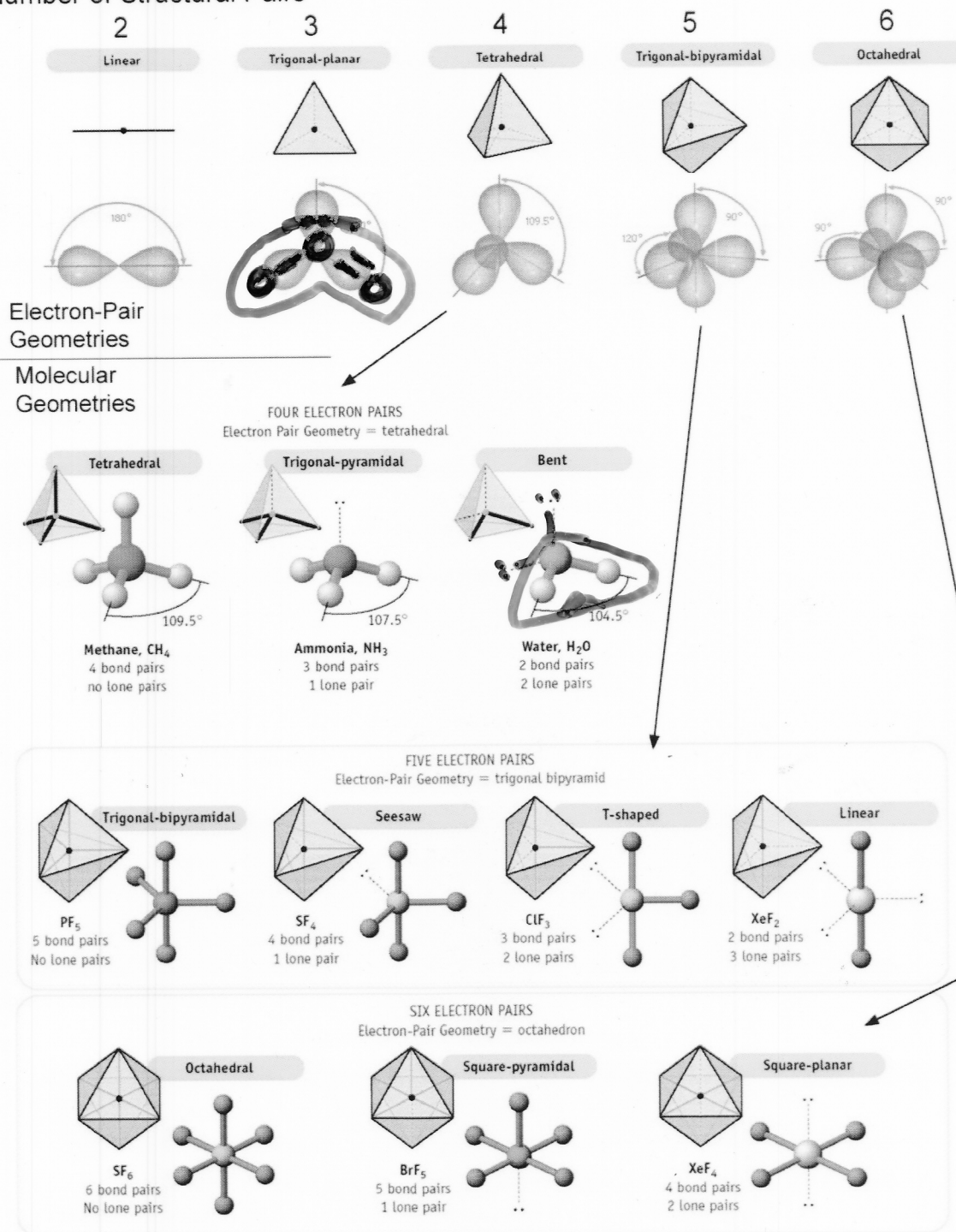
↓
Structural
Pairs

EPG

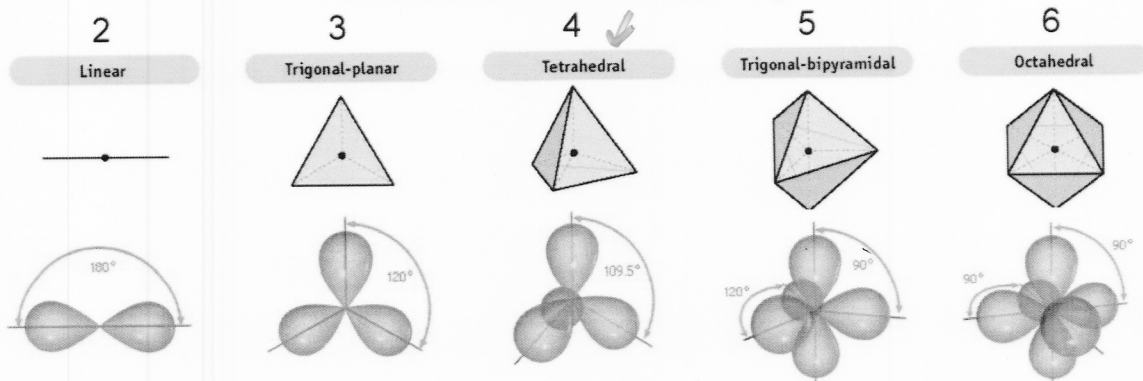
MG

Number of Pairs	Electron-Pair Geometry	Molecular Geometry			
		0	1	2	3
2	Linear	Linear			
3	trigonal planar	trigonal planar			
4	tetrahedral	tetrahedral	trigonal pyramidal	bent	
5	trigonal bipyramidal	trigonal bipyramidal	see-saw	T-shaped	
6	octahedral	octahedral	square pyramidal	square planar	

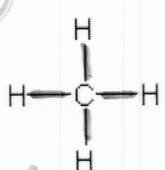
Number of Structural Pairs



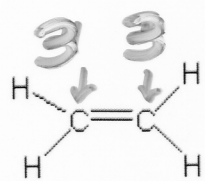
Electron Pair Geometry (EPG)



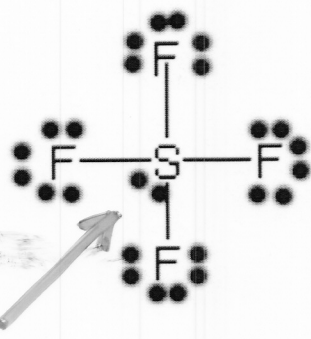
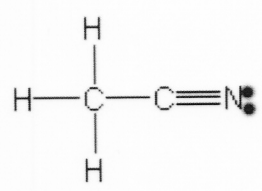
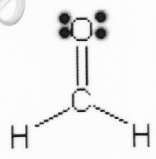
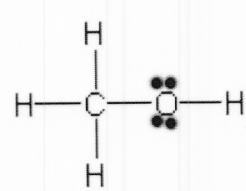
S.P. = 4



EPG = Tetrahedral

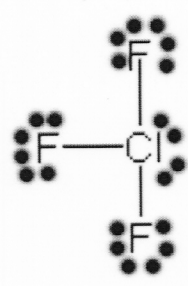
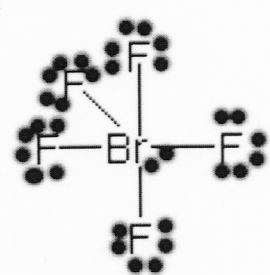


EPG = Trigonal planar



S.P. = 6

EPG = Octahedral



S.P. = 5

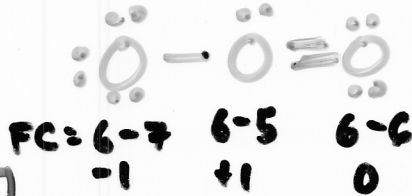
EPG = Trig. bipy.



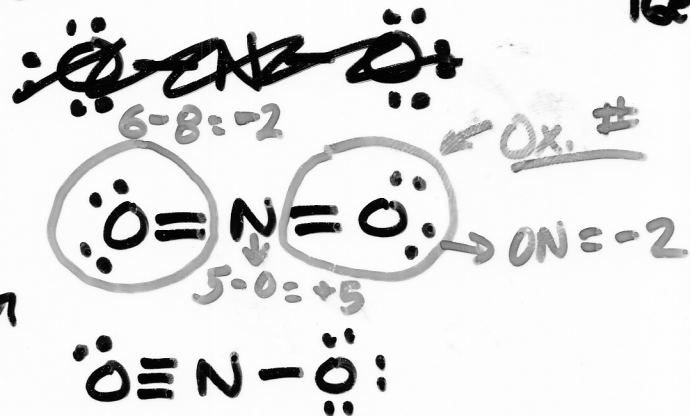
- Formal charge (FC)
 - Electrons are shared evenly throughout molecule
 - Closer to "real" charge than oxidation number
 - Use to determine which nonequivalent resonance structure is more likely
- Oxidation Number
 - Bonding electrons "belong" to more electronegative atom
 - Quick and easy to determine without Lewis structure
 - Doesn't really reflect actual charges on atoms
 - Use to determine whether compounds are oxidized/reduced in chemical reactions
- Partial Charges
 - How electrons are really distributed in a molecule
 - Need computer to calculate
 - Used only when you need very accurate description- molecular modeling

Examples:

Ozone
O₃



NO₂⁺



CH₃COOH



"CORRECT" SCN⁻ Picking "best" structure → FC's close to 0 → O has neg. FC

CO₂

