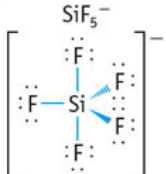
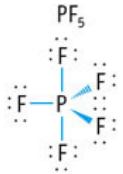
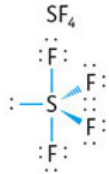
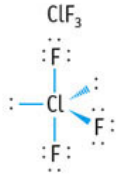
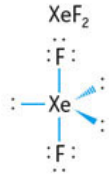
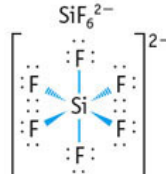
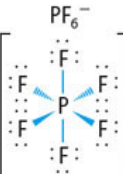
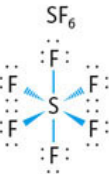
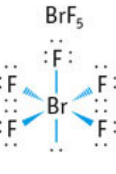
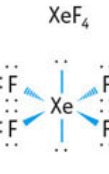


TABLE 8.6 Lewis Structures in Which the Central Atom Exceeds an Octet

Group 4A	Group 5A	Group 6A	Group 7A	Group 8
SiF_5^- 	PF_5 	SF_4 	ClF_3 	XeF_2 
SiF_6^{2-} 	PF_6^- 	SF_6 	BrF_5 	XeF_4 

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Linear

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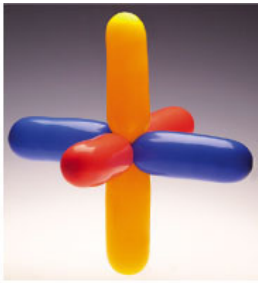
Trigonal planar



Tetrahedral



Trigonal bipyramidal



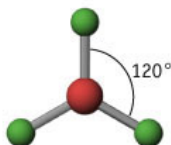
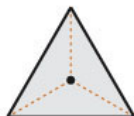
Octahedral

Linear



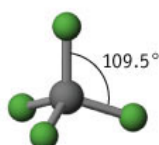
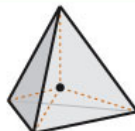
Example: BeF_2

Trigonal-planar



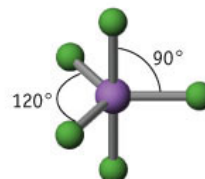
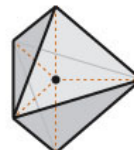
Example: BF_3

Tetrahedral



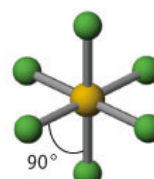
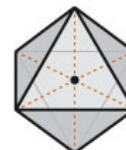
Example: CF_4

Trigonal bipyramidal



Example: PF_5

Octahedral



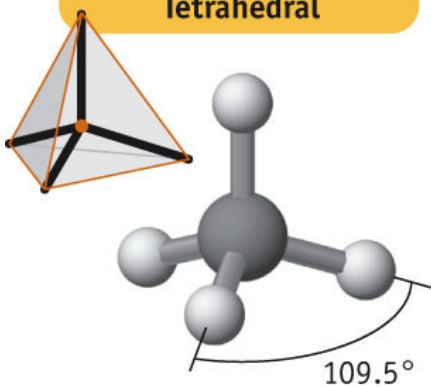
Example: SF_6

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Fig. 8-5, p. 369

FOUR ELECTRON PAIRS
Electron Pair Geometry = tetrahedral

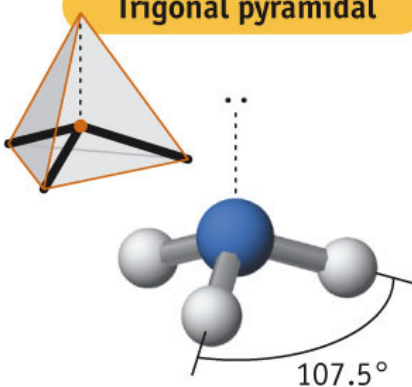
Tetrahedral



Methane, CH₄
4 bond pairs
no lone pairs

(a)

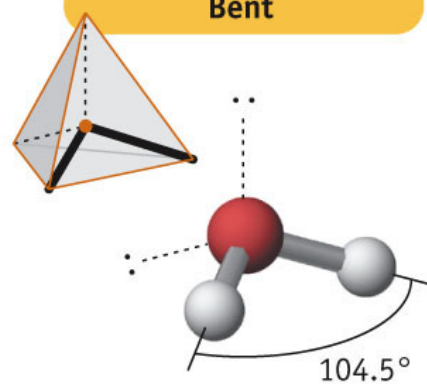
Trigonal pyramidal



Ammonia, NH₃
3 bond pairs
1 lone pair

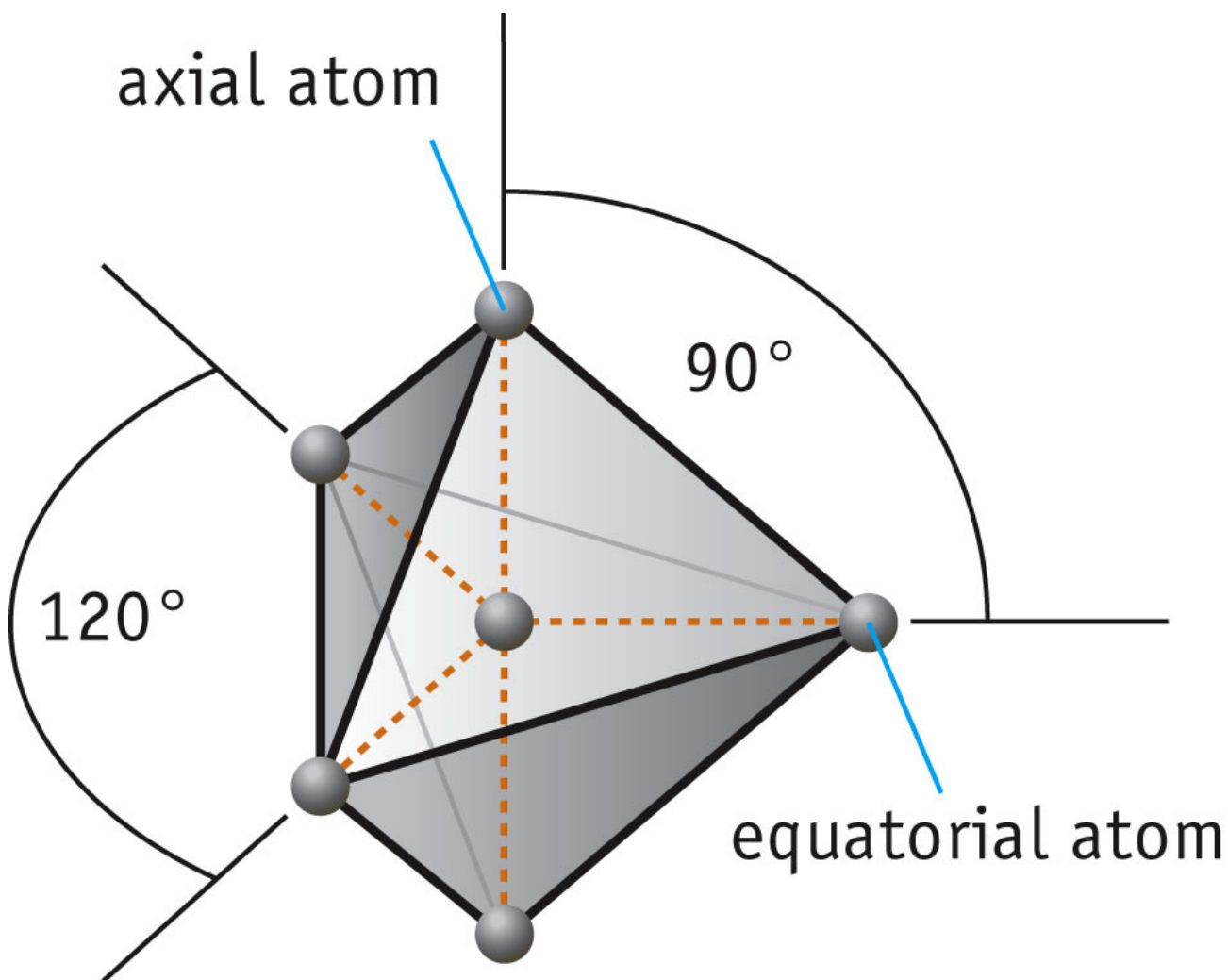
(b)

Bent



Water, H₂O
2 bond pairs
2 lone pairs

(c)



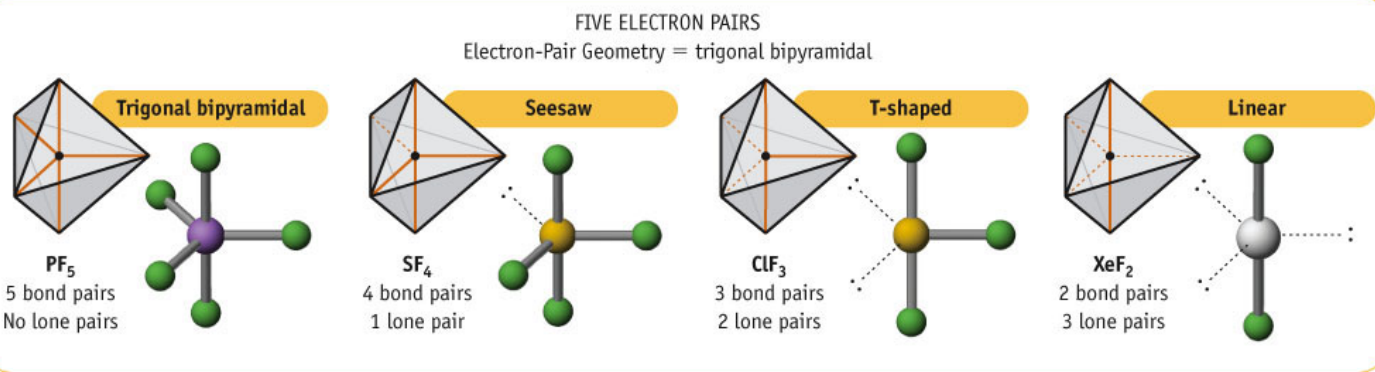
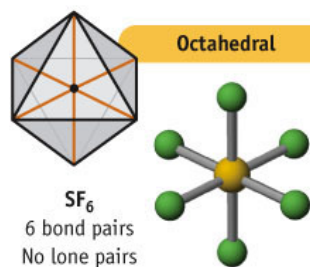


Fig. 8-8, p. 372



SIX ELECTRON PAIRS
Electron-Pair Geometry = octahedral

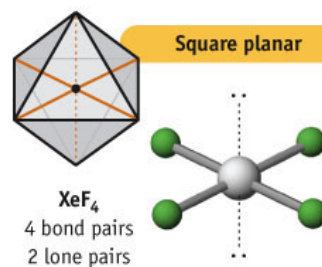
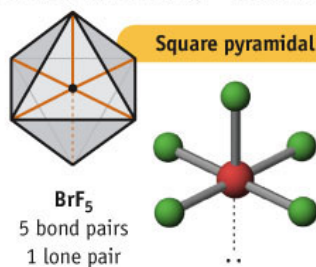
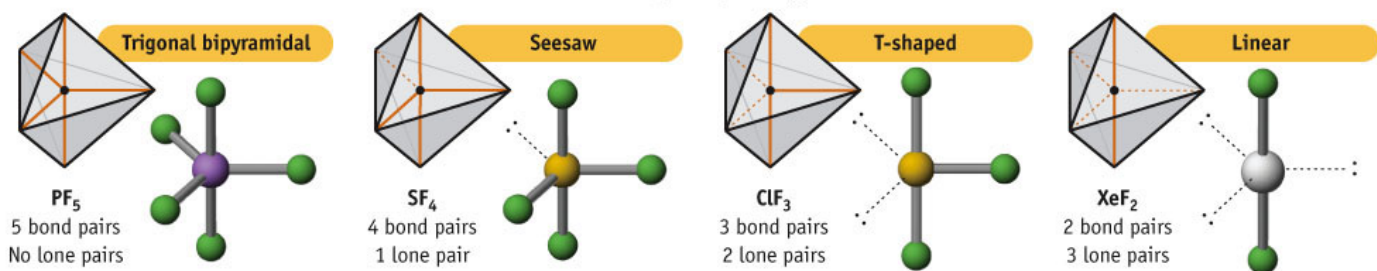


Fig. 8-8, p. 372

FIVE ELECTRON PAIRS
Electron-Pair Geometry = trigonal bipyramidal



SIX ELECTRON PAIRS
Electron-Pair Geometry = octahedral

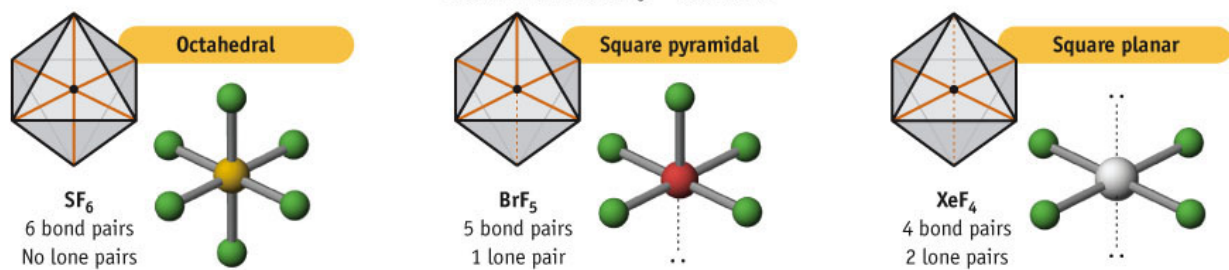
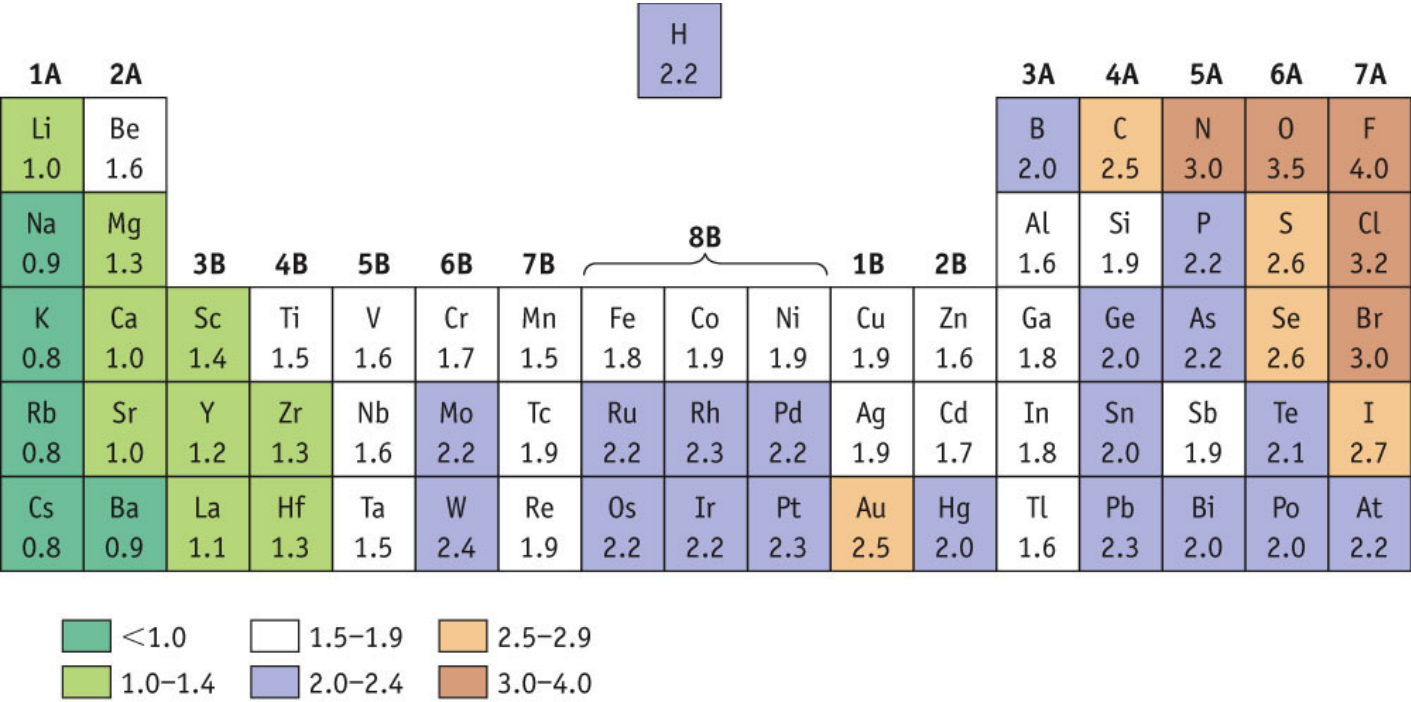


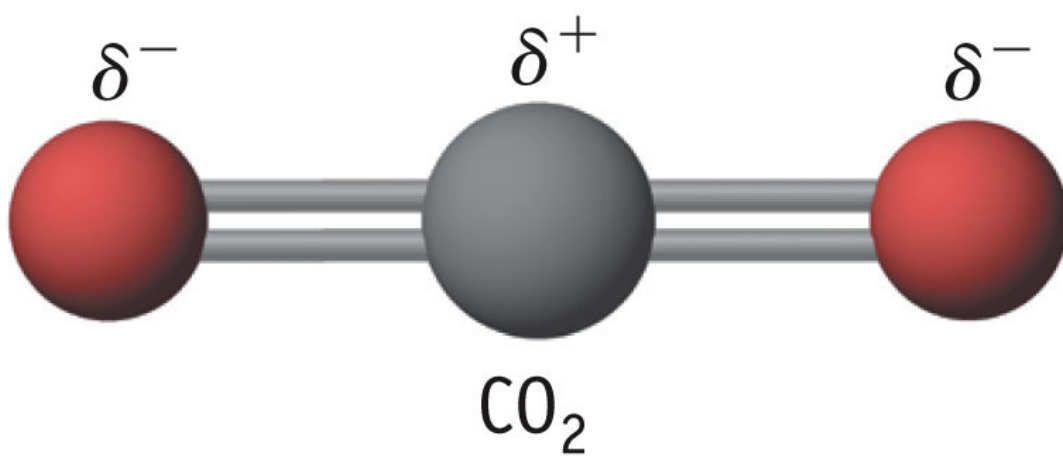
Fig. 8-8, p. 372



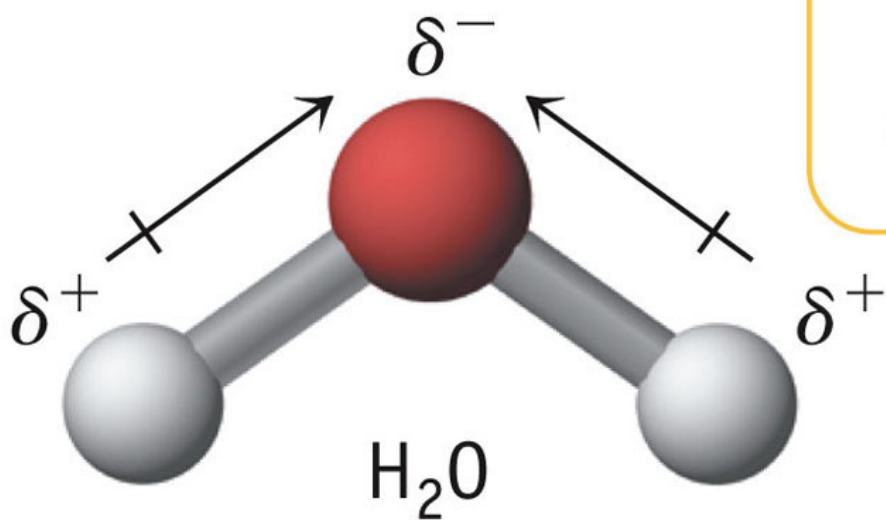
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Fig. 8-11, p. 376

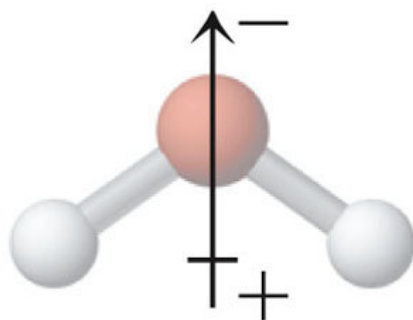
No net dipole
moment



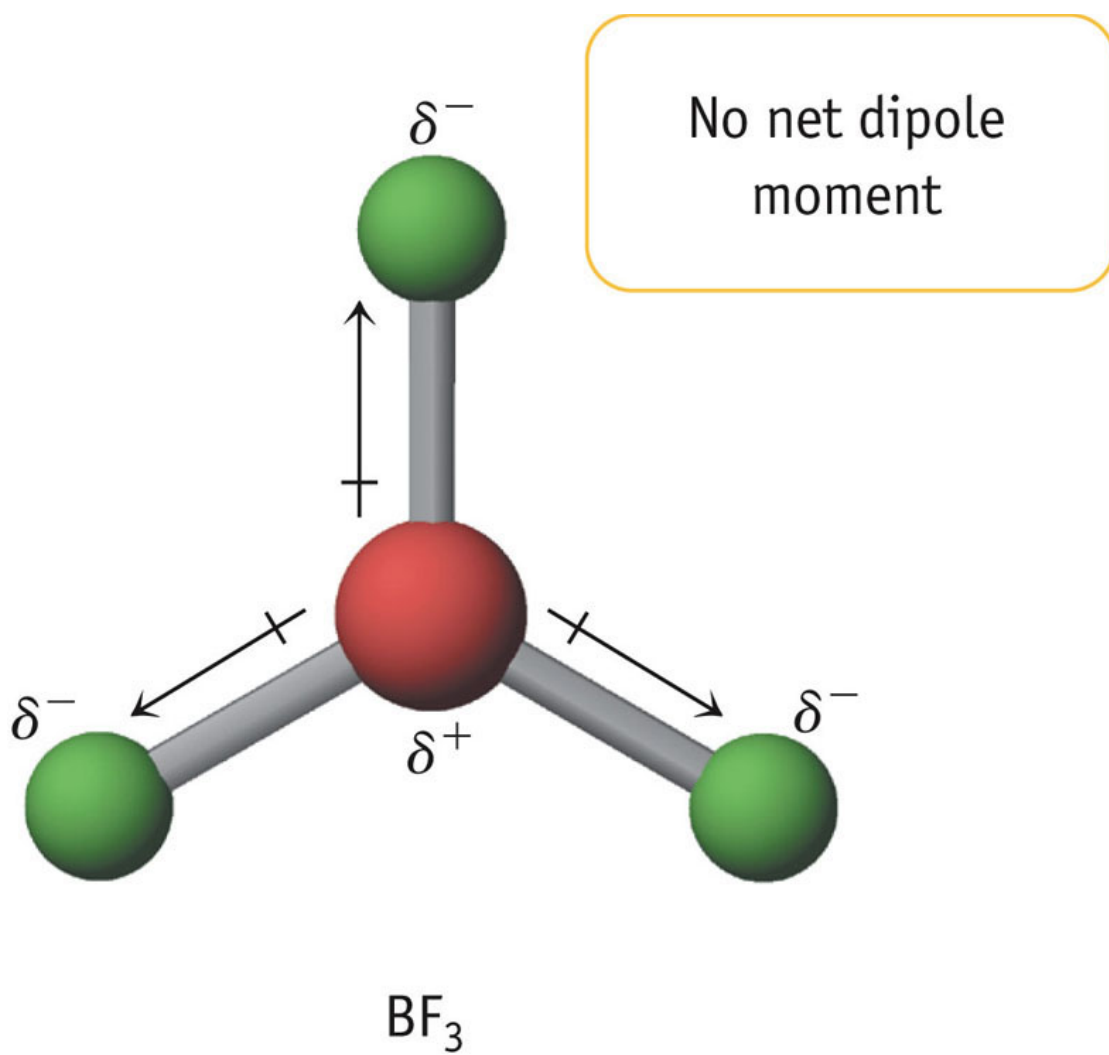
(a)

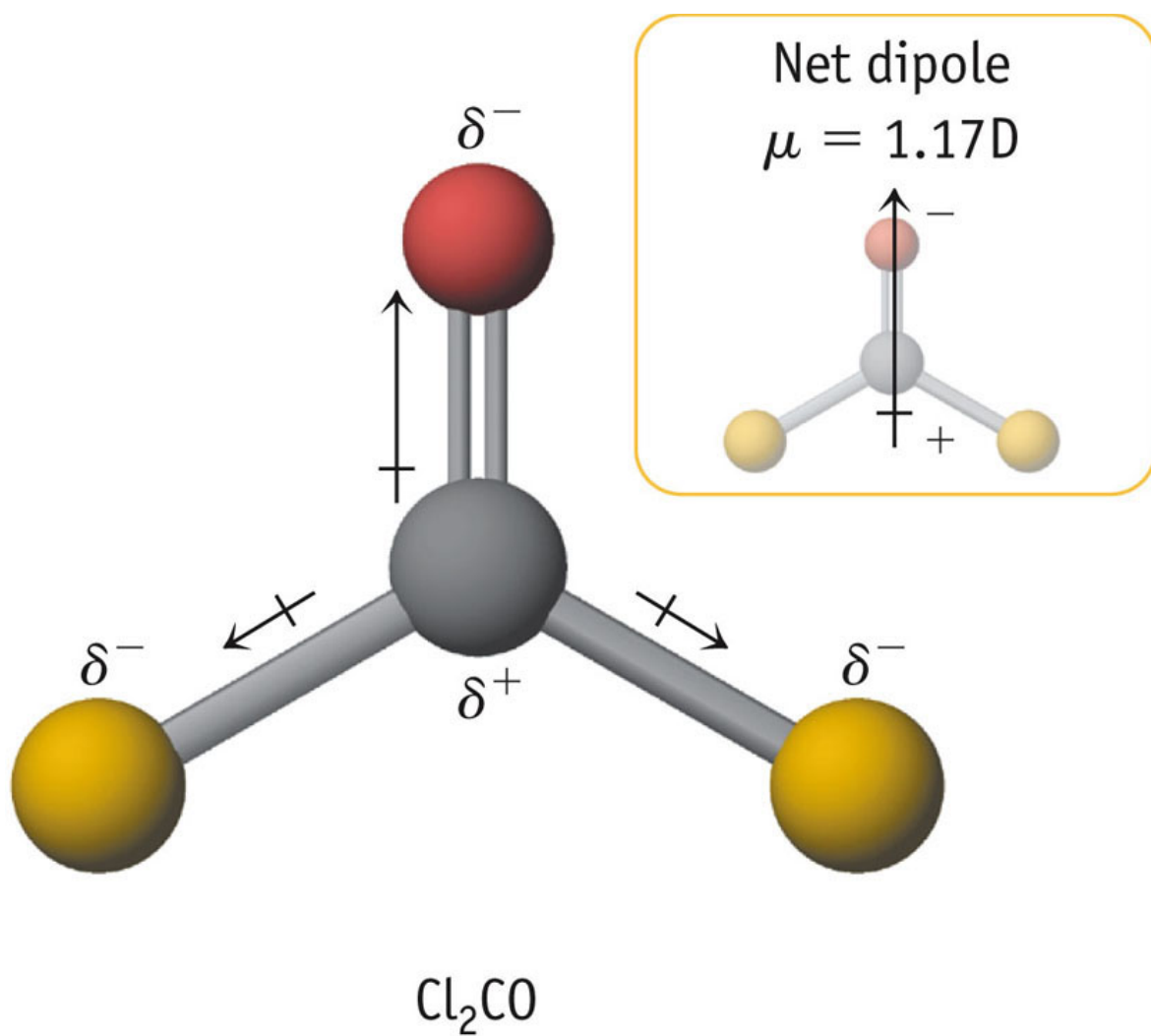


Net dipole
 $\mu = 1.85\text{D}$



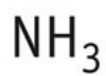
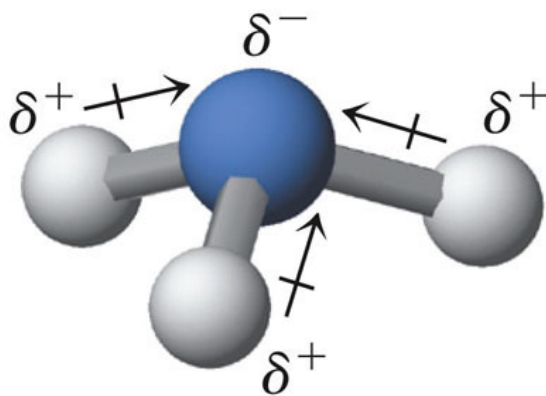
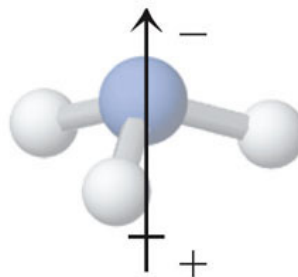
(b)

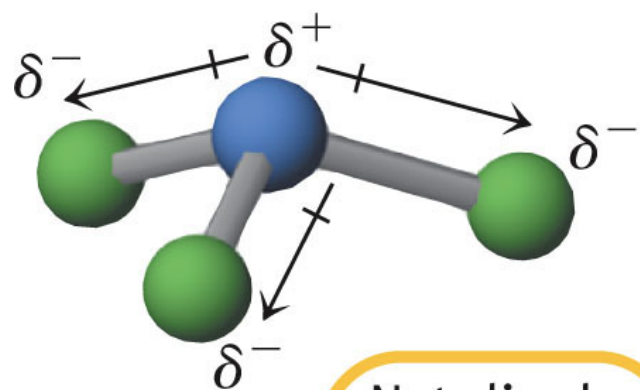




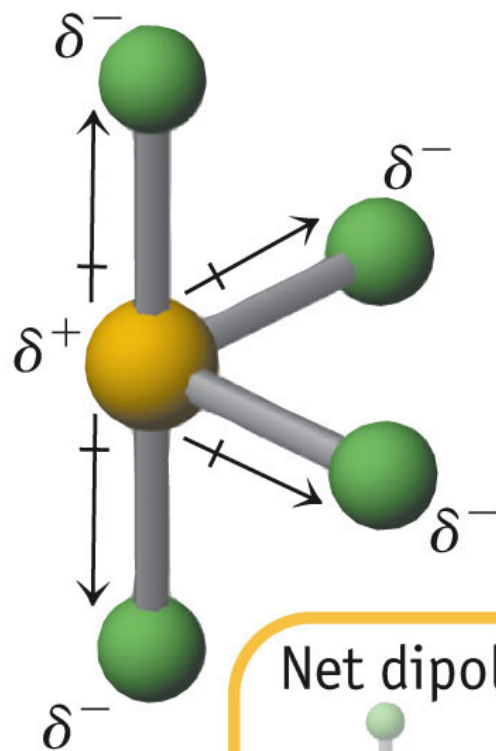
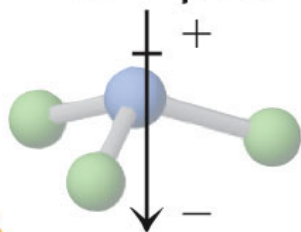
Net dipole

$$\mu = 1.47\text{D}$$





Net dipole



Net dipole

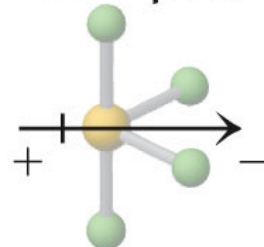


TABLE 8.8 Some Average Single- and Multiple-Bond Lengths in Picometers (pm)*

Single Bond Lengths											
Group											
	1A	4A	5A	6A	7A	4A	5A	6A	7A	7A	7A
	H	C	N	O	F	Si	P	S	Cl	Br	I
H	74	110	98	94	92	145	138	132	127	142	161
C		154	147	143	141	194	187	181	176	191	210
N			140	136	134	187	180	174	169	184	203
O				132	130	183	176	170	165	180	199
F					128	181	174	168	163	178	197
Si						234	227	221	216	231	250
P							220	214	209	224	243
S								208	203	218	237
Cl									200	213	232
Br										228	247
I											266

Multiple Bond Lengths			
C=C	134	C≡C	121
C=N	127	C≡N	115
C=O	122	C≡O	113
N=O	115	N≡O	108

*1 pm = 10⁻¹² m.

TABLE 8.9 Some Average Bond Dissociation Enthalpies (kJ/mol)*

Single Bonds											
	H	C	N	O	F	Si	P	S	Cl	Br	I
H	436	413	391	463	565	328	322	347	432	366	299
C		346	305	358	485	—	—	272	339	285	213
N			163	201	283	—	—	—	192	—	—
O				146	—	452	335	—	218	201	201
F					155	565	490	284	253	249	278
Si						222	—	293	381	310	234
P							201	—	326	—	184
S								226	255	—	—
Cl									242	216	208
Br										193	175
I											151
Multiple Bonds											
N=N				418	C=C				610		
N≡N				945	C≡C				835		
C=N				615	C=O				745		
C≡N				887	C≡O				1046		
O=O (in O ₂)				498							

*Sources of dissociation enthalpies: I. Klotz and R. M. Rosenberg: *Chemical Thermodynamics*, 4th Ed., p. 55, New York, John Wiley, 1994; and J. E. Huheey, E. A. Keiter, and R. L. Keiter: *Inorganic Chemistry* 4th Ed., Table E. 1, New York, Harper-Collins, 1993. See also Lange's *Handbook of Chemistry*, J. A. Dean (ed.), McGraw-Hill Inc., New York.