

$$Cl = Be = Cl$$
 $Be = Cl$
 $Be = Cl$

$$F = F \longrightarrow F$$

$$F = F \longrightarrow F$$

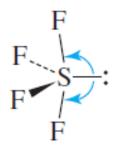
$$F = F$$

$$F =$$

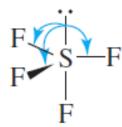
Predicted

The B—F bond length is 131 pm; the calculated single-bond length is 152 pm.

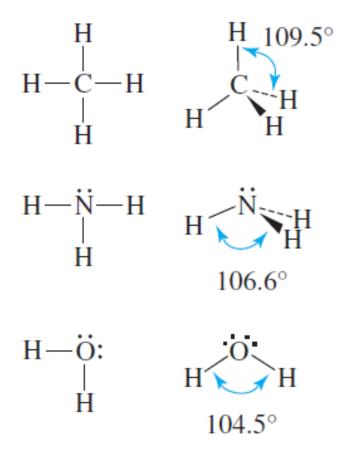
Steric Number	Geometry	Examples	Calculated Bond Angles	
2	Linear	CO_2	180°	0=c=0
3	Trigonal (triangular)	SO ₃	120°	o S
4	Tetrahedral	CH_4	109.5°	HHH
5	Trigonal bipyramidal	PCl ₅	120°, 90°	CI P—CI CI CI
6	Octahedral	SF_6	90°	$F = F \\ F = F$
7	Pentagonal bipyramidal	${\rm IF}_7$	72°, 90°	F F F
8	Square antiprismatic	[TaF ₈] ^{3–}	70.5°, 99.6°, 109.5°	F F F



Equatorial lone pair (observed structure)



Axial lone pair



		Number of Lone Pairs on Central Atom					
Steric Number	None	1	2	3			
2	:Cl=Be=Cl						
3	F F	Sn Cl 95° Cl					
4	H H C H	H N H H 106.6°	H-O:				
5	Cl Cl → P — Cl Cl ✓ Cl	173° F \ 101.6° F \ F F	F 86.2°	Xe—:			
6	F-S-F	F F 81.9°	F—Xe—F				

Molecule	X–P–X Angle (°)	Molecule	X–S–X Angle (°)
PF_3	97.8	OSF ₂	92.3
PCl ₃	100.3	$OSCl_2$	96.2
PBr_3	101.0	$OSBr_2$	98.2

Molecule	Bond Angle (°)	Molecule	Bond Angle (°)
H_2O	104.5	NCl_3	106.8
H_2S	92.1	PCl ₃	100.3
H_2Se	90.6	$AsCl_3$	98.9

TABLE 3.6 Bond Angles and Lengths

Molecule	Bond Angle (°)	Bond Length (pm)	Molecule	Bond Angle (°)	Bond Length (pm)	Molecule	Bond Angle (°)	Bond Length (pm)	Molecule	Bond Angle (°)	Bond Length (pm)
H_2O	104.5	97	OF_2	103.3	96	OCl ₂	110.9	170			
H_2S	92.1	135	SF ₂	98.0	159	SCI ₂	102.7	201			
H ₂ Se	90.6	146				SeCl ₂	99.6	216			
H ₂ Te	90.2	169				TeCl ₂	97.0	233			
NH ₃	106.6	101.5	NF ₃	102.2	137	NCl ₃	106.8	175			
PH ₃	93.2	142	PF_3	97.8	157	PCI ₃	100.3	204	PBr ₃	101.0	220
AsH ₃	92.1	151.9	AsF_3	95.8	170.6	AsCl ₃	98.9	217	$AsBr_3$	99.8	236
SbH ₃	91.6	170.7	SbF ₃	87.3	192	SbCl ₃	97.2	233	SbBr ₃	98.2	249

Source: N. N. Greenwood and A. Earnshaw, Chemistry of the Elements, 2nd ed., Butterworth-Heinemann, Oxford, 1997, pp. 557, 767; A. F. Wells, Structural Inorganic Chemistry, 5th ed., Oxford University Press, Oxford, 1987, pp. 705, 793, 846, and 879; R. J. Gillespie and I. Hargittai, The VSEPR Model of Molecular Geometry, Allyn and Bacon, Needham Heights, MA, 1991.