

Intuitively, we know symmetry when we see it.

But how do we put in quantitative terms that allows us to compare, assign, classify?

## Symmetry Operations and Symmetry Elements

#### **Definitions:**

- A symmetry operation is an operation on a body such that, after the operation has been carried out, the result is indistinguishable from the original body (every point of the body is coincident with an equivalent point or the same point of the body in its original orientation).
- A symmetry element is a geometrical entity such as a line, a plane, or a point, with respect to which one or more symmetry operations may be carried out

Symmetry Operation	Symmetry Element	Notation
Identity	-	E
Reflection in a plane	Plane of symmetry	$\sigma_{v}, \sigma_{d}, \sigma_{h}$
Proper rotation	Rotation axis (line)	$C_n$ ; where = 360/angle
Rotation followed by reflection in the plane perpendicular to the rotation axis	Improper rotation axis (line)	S <sub>n</sub>
Inversion	Center of inversion	I

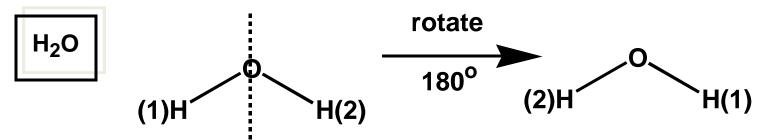
Notes

(i) symmetry operations more fundamental, but elements often easier to spot.

(ii) some symmetry elements give rise to more than one operation - especially rotation - as above.



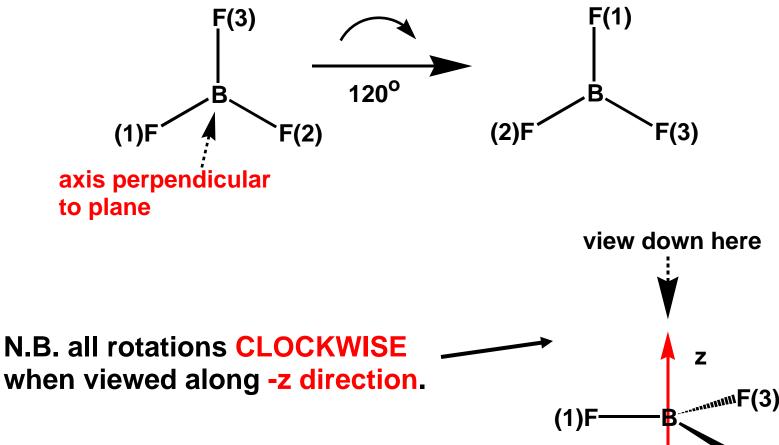
Some examples for different types of molecule: e.g.



Line in molecular plane, bisecting HOH angle is a rotation axis, giving indistinguishable configuration on rotation by 180°.



By VSEPR - trigonal, planar, all bonds equal, all angles 120°. Take as axis a line perpendicular to molecular plane, passing through B atom.



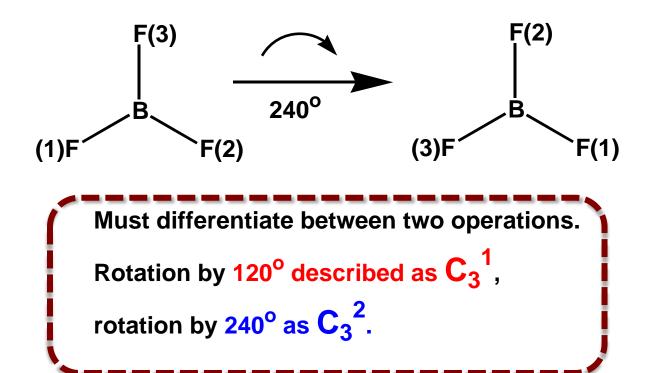
F(2)

### Symbol for axes of symmetry



where rotation about axis gives indistinguishable configuration every (360/n)<sup>o</sup> (i.e. an n-fold axis)

Thus  $H_2O$  has a  $C_2$  (two-fold) axis,  $BF_3$  a  $C_3$  (three-fold) axis. One axis can give rise to >1 rotation, e.g. for  $BF_3$ , what if we rotate by 240°?



In general C<sub>n</sub> axis (minimum angle of rotation (360/n)<sup>o</sup>) gives operations C<sub>n</sub><sup>m</sup>, where both m and n are integers.

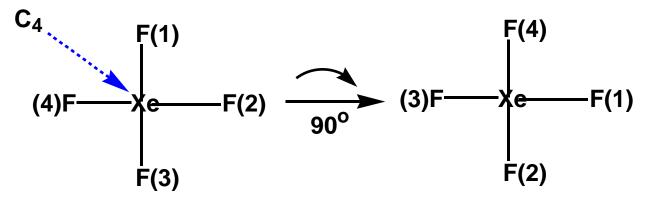
When m = n we have a special case, which introduces a new type of symmetry operation.....

## **IDENTITY OPERATION**

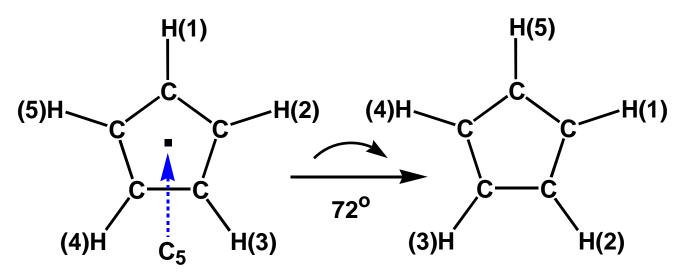
For  $H_2O$ ,  $C_2^2$  and for  $BF_3 C_3^3$  both bring the molecule to an IDENTICAL arrangement to initial one.

Rotation by 360<sup>o</sup> is exactly equivalent to rotation by 0<sup>o</sup>, i.e. the operation of doing NOTHING to the molecule.

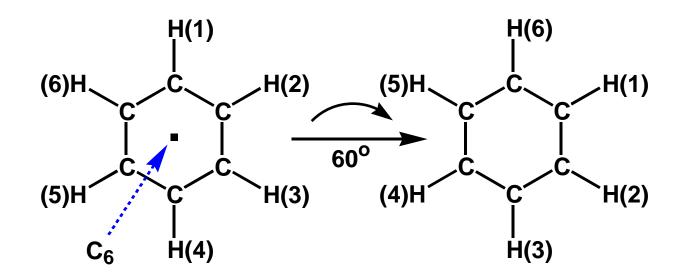
### xenon tetrafluoride, XeF<sub>4</sub>



## cyclopentadienide ion, C<sub>5</sub>H<sub>5</sub><sup>-</sup>

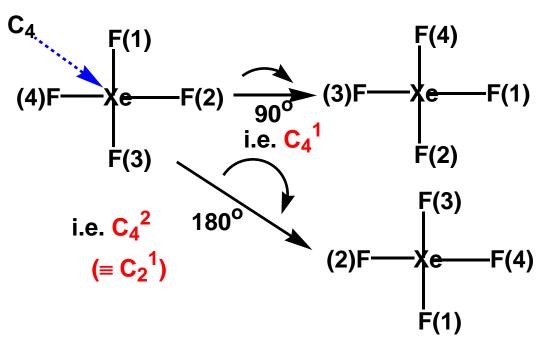


## benzene, C<sub>6</sub>H<sub>6</sub>



Examples also known of  $C_7$  and  $C_8$  axes.

If a C<sub>2n</sub> axis (i.e. even order) present, then C<sub>n</sub> must also be present:



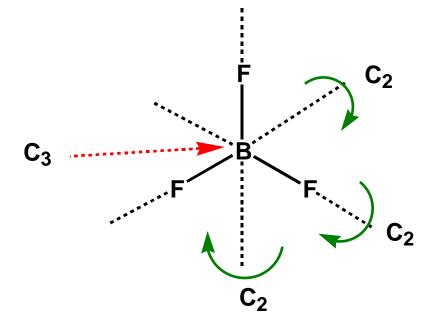
Therefore there must be a  $C_2$  axis coincident with  $C_4$ , and the operations generated by  $C_4$  can be written:

$$C_4^{1}, C_4^{2} (C_2^{1}), C_4^{3}, C_4^{4} (E)$$

Similarly, a  $C_6$  axis is accompanied by  $C_3$  and  $C_2$ , and the operations generated by  $C_6$  are:

 $C_6^{1}, C_6^{2} (C_3^{1}), C_6^{3} (C_2^{1}), C_6^{4} (C_3^{2}), C_6^{5}, C_6^{6} (E)$ 

## Molecules can possess several distinct axes, e.g. BF<sub>3</sub>:



Three  $C_2$  axes, one along each B-F bond, perpendicular to  $C_3$ 

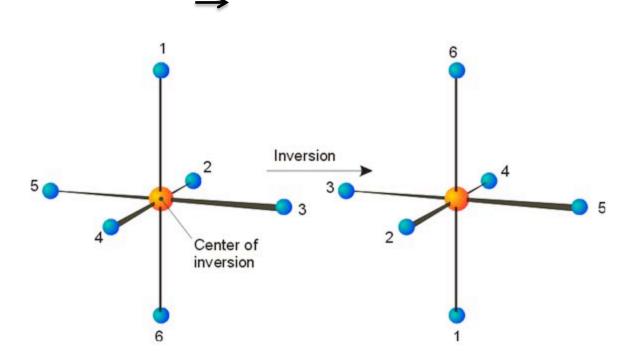
## Inversion (i)

Each atom in the molecule is moved along a straight line through the inversion center to a point an equal distance from the

X,Y,Z

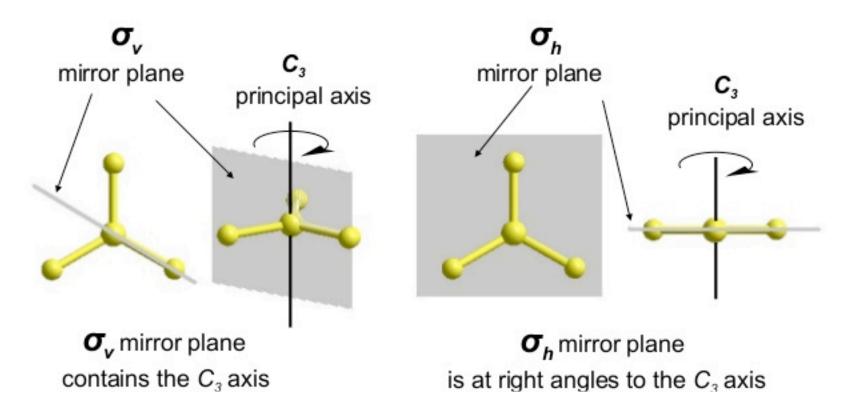
-X, -Y, -Z

inversion center.



## Mirror planes ( $\sigma$ ) of BF<sub>3</sub>:

Mirror planes can contain the principal axis ( $\sigma_v$ ) or be at right angles to it ( $\sigma_h$ ). BF<sub>3</sub> has one  $\sigma_h$  and three  $\sigma_v$  planes: (v = vertical, h = horizontal)



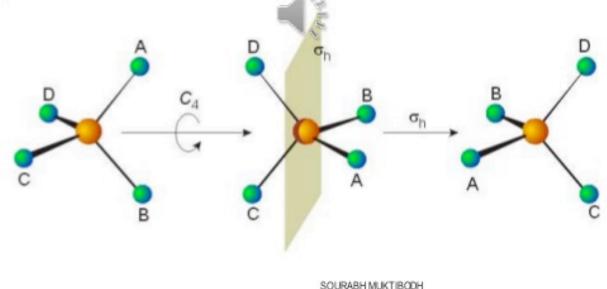
## IMPROPER ROTATION

An improper rotation is rotation, followed by reflection in the plane perpendicular to the axis of rotation. Thus

 $S_n = C_n * i = i * C_n$ 

both independent symmetry operations commute. Essentially





Symmetry elements/operations can be manipulated by Group Theory, Representations and Character Tables



So, What IS a group?





And, What is a Character???



A GROUP is a collection of entities or elements which satisfy the following four conditions:

1) The product of any two elements (including the square of each element) must be an element of the group. For symmetry operations, the multiplication rule is to successively perform operations.

2) One element in the group must commute with all others and leave them unchanged. Therefore the "E",

$$EX = XE = X$$

3) The associative law of multiplication must hold

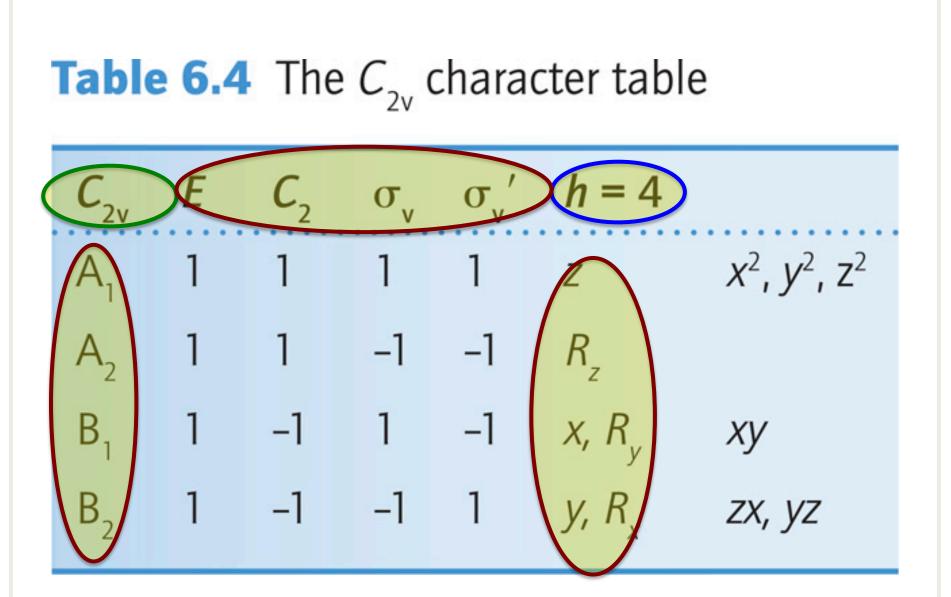
$$A(BC) = (AB)C$$

4) Every element must have a reciprocal which is also an element of the group. i.e.,

$$X(X^{-1}) = (X^{-1}) X = E$$

Note: An element may be its own reciprocal.

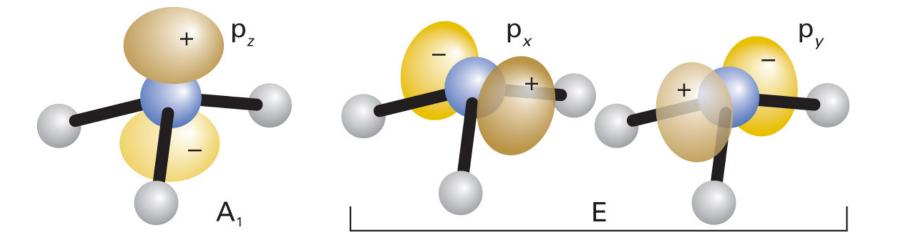
Groups may be composed of anything: symmetry operations, nuclear particles, etc. Simplest is +1, -1.



W. Н.

### Character table for point group C<sub>3v</sub>

C <sub>3v</sub>	E	2C <sub>3</sub> (z)	3σ <sub>v</sub>	linear functions, rotations	quadratic functions	cubic functions
A <sub>1</sub>	+1	+1	+1	Z	x <sup>2</sup> +y <sup>2</sup> , z <sup>2</sup>	$z^3$ , x(x <sup>2</sup> -3y <sup>2</sup> ), z(x <sup>2</sup> +y <sup>2</sup> )
A <sub>2</sub>	+1	+1	-1	Rz	-	y(3x <sup>2</sup> -y <sup>2</sup> )
E	+2	-1	0	(x, y) (R <sub>x</sub> , R <sub>y</sub> )	(x <sup>2</sup> -y <sup>2</sup> , xy) (xz, yz)	$(xz^2, yz^2) [xyz, z(x^2-y^2)]$ [x(x <sup>2</sup> +y <sup>2</sup> ), y(x <sup>2</sup> +y <sup>2</sup> )]



#### Table 6.3 The components of a character table

Name of point group*	Symmetry operations <i>R</i> arranged by class ( <i>E</i> , <i>C<sub>n</sub></i> , etc.)	Functions	Further functions	Order of group, <i>h</i>
Symmetry species (Γ)	Characters ( $\chi$ )	Translations and components of dipole moments ( <i>x</i> , <i>y</i> , <i>z</i> ), of relevance to IR activity; rotations	Quadratic functions such as <i>z</i> <sup>2</sup> , <i>xy</i> , etc., of relevance to Raman activity	

\* Schoenflies symbol.

W. H.

# Consequences of Symmetry

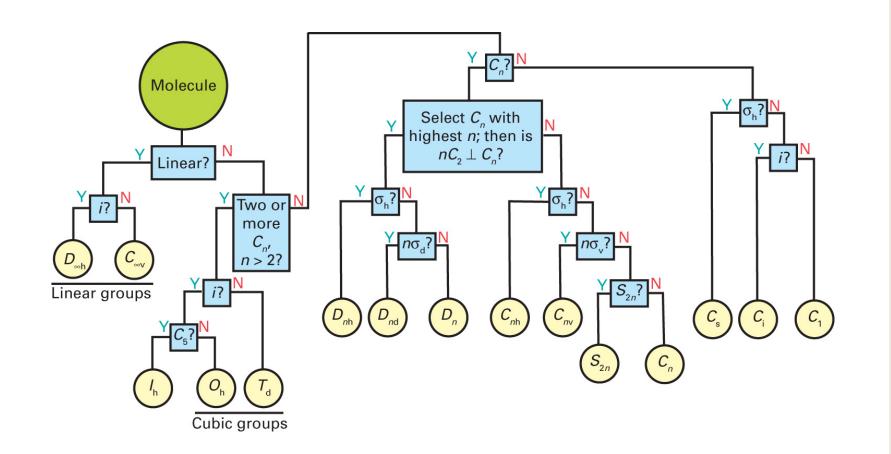
- Only the molecules which belong to the C<sub>n</sub>, C<sub>nv</sub>, or C<sub>s</sub> point group can have a permanent dipole moment.
- A molecule may be chiral only if it does not have an axis of improper rotation **Sn**.
- IR Allowed transitions may be predicted by symmetry operations
- Orbital overlap may be predicted and described by symmetry

# Point Group Assignments and Character Tables



A collection of symmetry operations all of which pass through a single point A point group for a molecule is a quantitative measure of the symmetry of that molecule

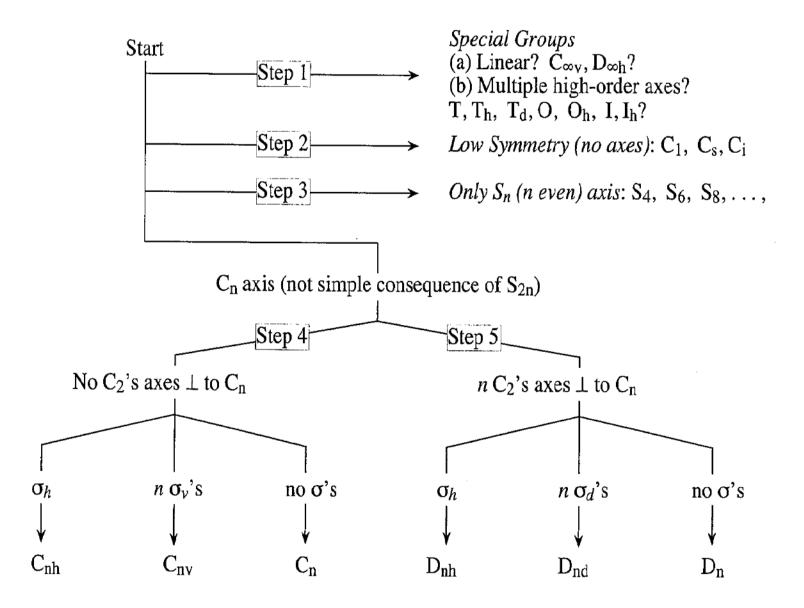
Assignment of Symmetry Elements to Point Group: At first Looks Daunting.



Daunting? However almost all we will be concerned with belong to just a few symmetry point groups

W. H.

## A Simpler Approach



# **POINT GROUPS**

A collection of symmetry operations all of which pass through a single point

A point group for a molecule is a quantitative measure of the symmetry of that molecule

**ASSIGNMENT OF MOLECULES TO POINT GROUPS** 

**STEP 1 : LOOK FOR AN AXIS OF SYMMETRY** 

If one is found - go to STEP 2

If not: look for

(a) plane of symmetry - if one is found, molecule belongs to point group  $C_s$ 

#### Point Group Assignments: MFT Ch. 4

TABLE 4.3	Groups of High Symmetry	
Group	Description	Examples
$C_{\infty \nu}$	These molecules are linear, with an infinite number of rotations and an infinite number of reflection planes containing the rotation axis. They do not have a center of inversion.	C <sub>∞</sub> H—Cl
$D_{\infty h}$	These molecules are linear, with an infinite number of rotations and an infinite number of reflection planes containing the rotation axis. They also have perpendicular $C_2$ axes, a perpendicular reflection plane, and an inversion center.	$C_{\infty} \rightarrow 0 = C_{2}$
T <sub>d</sub>	Most (but not all) molecules in this point group have the familiar tetrahedral geometry. They have four $C_3$ axes, three $C_2$ axes, three $S_4$ axes, and six $\sigma_d$ planes. They have no $C_4$ axes.	
O <sub>h</sub>	These molecules include those of octahedral struc- ture, although some other geometrical forms, such as the cube, share the same set of symmetry opera- tions. Among their 48 symmetry operations are four $C_3$ rotations, three $C_4$ rotations, and an inversion.	F-S-F FF
I <sub>h</sub>	Icosahedral structures are best recognized by their six $C_5$ axes, as well as many other symmetry operations—120 in all.	

#### TABLE 4.3 Groups of High Symmetry

In addition, there are four other groups, T,  $T_h$ , O, and I, which are rarely seen in nature. These groups are discussed at the end of this section.

B<sub>12</sub>H<sub>12</sub><sup>2-</sup> with BH at each vertex of an icosahedron

## LINEAR MOLECULES

Do in fact fit into scheme - but they have an infinite number of symmetry operations.

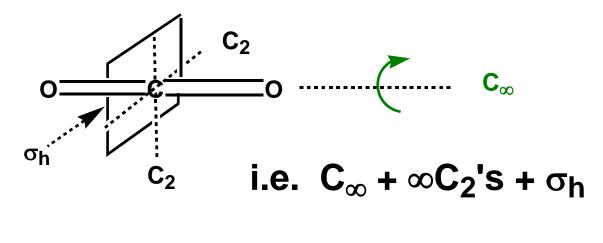
Molecular axis is  $C_{\infty}$  - rotation by any arbitrary angle  $(360/\infty)^{\circ}$ , so infinite number of rotations. Also any plane containing axis is symmetry plane, so infinite number of planes of symmetry.

Divide linear molecules into two groups:

(i) No centre of symmetry, e.g.:  $H \longrightarrow C \longrightarrow N$  ......  $C_{\infty}$ 

No C<sub>2</sub>'s perp. to main axis, but  $\infty \sigma_v$ 's containing main axis: point group  $C_{\infty v}$ 



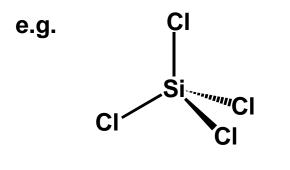


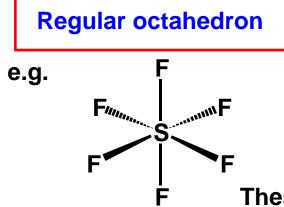
Point group  $D_{\infty h}$ 

Highly symmetrical molecules

A few geometries have several, equivalent, highest order axes. Two geometries most important:

### **Regular tetrahedron**





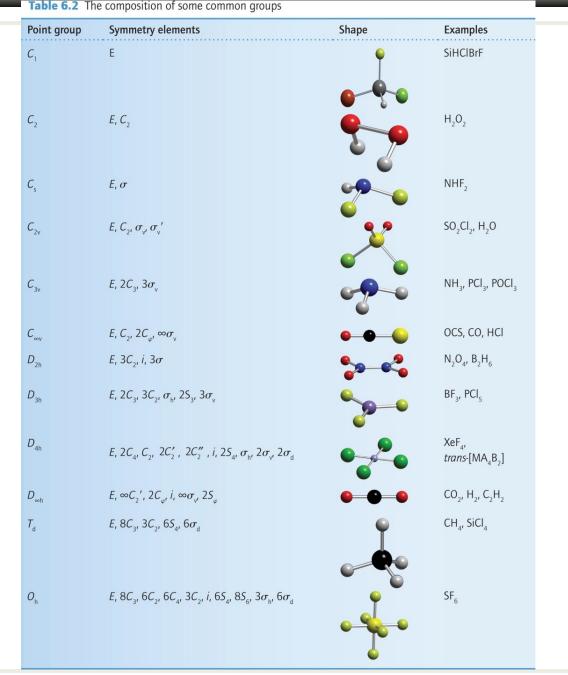
 $3C_4$ 's (along F-S-F axes) also  $4C_3$ 's.  $6C_2$ 's, several planes,  $S_4$ ,  $S_6$  axes, and a centre of symmetry (at S atom) Point group  $O_h$ 

These molecules can be identified without going through the usual steps.

Note: many of the more symmetrical molecules possess many more symmetry operations than are needed to assign the point group.

#### Inorganic Chemistry Chapter 1: Table 6.2 Table 6.2 The composition of some common groups

#### W. H. Freeman



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4. The  $C_{nv}$  Groups

$ \frac{C_{2v}}{A_1} $ $ \frac{A_2}{B_1} $ $ B_2 $	E 1 1 1 1	$ \begin{array}{cccc} C_2 & \sigma_1(xz) \\ \hline 1 & 1 \\ 1 & -1 \\ -1 & 1 \\ -1 & -1 \\ \hline -1 & -1 \end{array} $		$ \frac{z}{R_z} \\ x, R_y \\ y, R_x $	x <sup>2</sup> , y <sup>2</sup> xy xz yz	<sup>2</sup> , z <sup>2</sup>	
$ \begin{array}{c} C_{3v} \\ \hline A_1 \\ A_2 \\ E \end{array} $	E 1 1 2	$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$	$z R_z (x, y)(z)$	R <sub>x</sub> , R <sub>y</sub> )	$x^{2} + \frac{1}{(x^{2} - x^{2})^{2}}$	$y^2, z^2$ $y^2, xy)($	xz, yz)
$ \begin{array}{c} C_{4v} \\ \hline \\ A_1 \\ A_2 \\ B_1 \\ B_2 \\ E \\ \end{array} $	E 1 1 1 1 2	$ \begin{array}{cccccccccccccccccccccccccccccccccccc$	1 1 1 1		x, Ry)	$x^{2} + y$ $x^{2} - y$ $xy$ $(xz, yz$	)
$\frac{C_{50}}{A_1}$ $\frac{A_2}{E_1}$ $E_2$	E 1 1 2 2	1 1 2 cos 72° 2 co	2 <i>C</i> ₅ <sup>2</sup> 1 1 0s 144° 0s 72°	5σ <sub>υ</sub> 1 -1 0 0	z R <sub>z</sub> (x, y)(R	R <sub>x</sub> , R <sub>y</sub> )	$     x^{2} + y^{2}, z^{2} $ (xz, yz) (x <sup>2</sup> - y <sup>2</sup> , xy)
$ \begin{array}{c} C_{6\nu} \\ A_1 \\ A_2 \\ B_1 \\ B_2 \\ E_1 \\ E_2 \end{array} $	E 1 1 1 2 2	$\begin{array}{cccccccccccccccccccccccccccccccccccc$	-1 1	$\begin{array}{c c} 3\sigma_{d} \\ \hline 1 & z \\ -1 & R_{z} \\ -1 & 1 \\ 1 & 0 \\ 0 & (x) \end{array}$	, <i>y</i> )( <i>R</i> <sub>x</sub> ,		$x^{2} + y^{2}, z^{2}$ (xz, yz) (x <sup>2</sup> - y <sup>2</sup> , xy)

4. The  $C_{nv}$  Groups

$ \frac{C_{2v}}{A_1} $ $ \frac{A_2}{B_1} $ $ B_2 $	E 1 1 1 1	$ \begin{array}{cccc} C_2 & \sigma_1(xz) \\ \hline 1 & 1 \\ 1 & -1 \\ -1 & 1 \\ -1 & -1 \\ \hline -1 & -1 \end{array} $		$ \frac{z}{R_z} \\ x, R_y \\ y, R_x $	x <sup>2</sup> , y <sup>2</sup> xy xz yz	<sup>2</sup> , z <sup>2</sup>	
$ \begin{array}{c} C_{3v} \\ \hline A_1 \\ A_2 \\ E \end{array} $	E 1 1 2	$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$	$z R_z (x, y)(z)$	R <sub>x</sub> , R <sub>y</sub> )	$x^{2} + \frac{1}{(x^{2} - x^{2})^{2}}$	$y^2, z^2$ $y^2, xy)($	xz, yz)
$ \begin{array}{c} C_{4v} \\ \hline \\ A_1 \\ A_2 \\ B_1 \\ B_2 \\ E \\ \end{array} $	E 1 1 1 1 2	$ \begin{array}{cccccccccccccccccccccccccccccccccccc$	1 1 1 1		x, Ry)	$x^{2} + y$ $x^{2} - y$ $xy$ $(xz, yz$	)
$\frac{C_{50}}{A_1}$ $\frac{A_2}{E_1}$ $E_2$	E 1 1 2 2	1 1 2 cos 72° 2 co	2 <i>C</i> ₅ <sup>2</sup> 1 1 0s 144° 0s 72°	5σ <sub>υ</sub> 1 -1 0 0	z R <sub>z</sub> (x, y)(R	R <sub>x</sub> , R <sub>y</sub> )	$     x^{2} + y^{2}, z^{2} $ (xz, yz) (x <sup>2</sup> - y <sup>2</sup> , xy)
$ \begin{array}{c} C_{6\nu} \\ A_1 \\ A_2 \\ B_1 \\ B_2 \\ E_1 \\ E_2 \end{array} $	E 1 1 1 2 2	$\begin{array}{cccccccccccccccccccccccccccccccccccc$	-1 1	$\begin{array}{c c} 3\sigma_{d} \\ \hline 1 & z \\ -1 & R_{z} \\ -1 & 1 \\ 1 & 0 \\ 0 & (x) \end{array}$	, <i>y</i> )( <i>R</i> <sub>x</sub> ,		$x^{2} + y^{2}, z^{2}$ (xz, yz) (x <sup>2</sup> - y <sup>2</sup> , xy)

6. The  $D_{nh}$  Groups

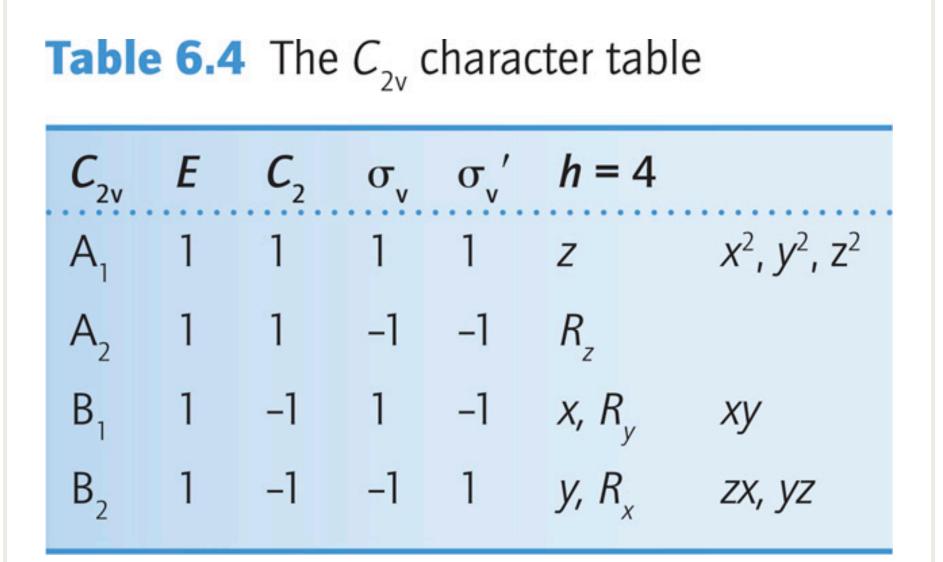
D 2 k	E	$C_2(z)$	C₂(y)	$C_2(x)$	i	σ(xy)	σ(xz)	σ(yz)			
A <sub>0</sub> B <sub>10</sub> B <sub>20</sub> B <sub>30</sub> A <sub>u</sub> B <sub>1u</sub> B <sub>2u</sub> B <sub>3u</sub>	1 1 1 1 1 1 1				1 1 1 -1 -1 -1				Rz Ry Rz z y z	x <sup>2</sup> , j xy xz yz	y <sup>2</sup> , z <sup>2</sup>
D3M	E	2C <sub>3</sub> 2	$C_2 \sigma_h$	253	300	_		-			
A <sub>1</sub> ', A <sub>2</sub> ' A <sub>1</sub> " A <sub>2</sub> " E"	1 1 2 1 1 2	1 1 1 1 1		$ \begin{array}{c} 1 \\ 1 \\ 2 \\ -1 \\ 1 \\ -1 \\ 1 \\ -1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ $		R <sub>z</sub> (x, ) z (R <sub>x</sub>	v) , R <sub>y</sub> )	$x^2 + y$ $(x^2 - y)$ $(xz, yz)$	y², xy)		
DAN	E	2 <i>C</i> _	C <sub>2</sub> 2C <sub>2</sub>	′ 2 <i>C</i> 2 <sup>*</sup>	i	25.	σ <sub>R</sub> 2σ	-, 2σ,		ł	
A 1.9 A 2.9 B 1.9 B 2.9 E 9 A 1.4 A 2.4 B 2.4 B 2.4 E 4 E 4 E 4 E 4 E 4 E 4 E 4 E 4 E 4 E	I I I I I I I I I I 2	$ \begin{array}{c} 1 \\ -1 \\ -1 \\ 0 \\ 1 \\ -1 \\ -1 \\ 0 \\ \end{array} $	1 - 1 1 - 2 1 - 2 1 - 2 1 - 1 1 - 2 1	$ \begin{array}{cccccccccccccccccccccccccccccccccccc$	$ \begin{array}{c} 1 \\ 1 \\ 2 \\ -1 \\ -1 \\ -1 \\ -2 \\ -2 \\ \end{array} $	i 1 1 1 1 1 1 1 1 1 	$ \begin{array}{cccccccccccccccccccccccccccccccccccc$	$ \begin{array}{cccccccccccccccccccccccccccccccccccc$	R <sub>z</sub> (R <sub>x</sub> , z (x, y		$     x^{2} + y^{2}, z^{2}      x^{2} - y^{2}      xy      (xz, yz) $

5. The D<sub>nh</sub> Groups

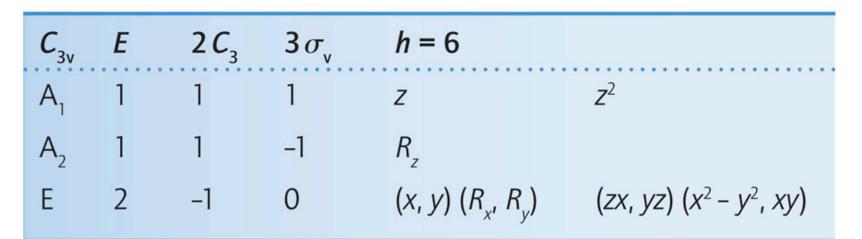
D 28	E	$C_2(z)$	C2())	$C_1(x)$	1	$\sigma(xy)$	σ(xz)	o()/2)			
A. B B.2. B.3. A. B.1. B.1. B.2. B.3. B.3.			-1 -1 -1 -1 -1	1 -1 -1 1 -1 -1 1 -1			-1 -1 -1 -1 -1		R, R, R, z y x	x <sup>2</sup> , xy xz yz	y <sup>2</sup> , z <sup>2</sup>
Dam	E	2C, 2	BC <sub>2</sub> σ <sub>h</sub>	2.53	300		1				
1.1.1 1.1.1.1 1.1.	1 2 1 1 2			-1	$-1 \\ -1 \\ -1 \\ 1 \\ 0$	R. (x, ) (R,	y) . R,)	$x^2 + y$ $(x^2 - z)$ $(xz, yz)$	y², xy)		
D	E	264	C1 2C1	2C2*	i	25. 0	σ <sub>h</sub> 2σ	. 20.	1	I	
A 1.0 A 200 B 200 E 0 A 100 B 200 E 0 A 100 B 200 E 0 E 0 E 0	1 1 2 1 1 1 2		$   \begin{bmatrix}     1 \\     1 \\     -2 \\     -2 \\     1 \\     1 \\     -1 \\     -2 \\     1 \\     -2 \\     1 \\     -2 \\     1 \\     -2 \\    $		1 1 2 1 1 2 1 1 2 1 1 2 1 1 2 1 1 2 1 1 2 1 1 2 1 1 2 1 1 2 1 1 2 1 1 2 1 1 2 1 1 2 1 1 1 2 1 1 1 2 1 1 1 2 1 1 1 1 2 1 1 1 1 1 2 1 1 1 1 2 1	-1 -1 -1 -1 -1 -1 -1 -1 -1 -1 -1 -1 -1 -	-1 -1 -1 -1 -1 -1 -1 -1 -1 -1 -1 -1 -1 -	$ \begin{array}{cccccccccccccccccccccccccccccccccccc$	R. (R z (x. y		$x^{2} + y^{2}, z^{2}$ $x^{2} - y^{2}$ xy (xz, yz)

9. The Cubic Groups (Continued).

T <sub>h</sub>	$E 4C_3 4C_3^2 3C_2 i 4S_6 4S_6^3 3\sigma_h$		$\epsilon = \exp\left(2\pi i/3\right)$
Ag	1 1 1 1 1 1 1		$\frac{1}{x^2+y^2+z^2}$
Au Ee	$ \begin{array}{cccccccccccccccccccccccccccccccccccc$		$\begin{array}{l} (2z^2 - x^2 - y^2, \\ x^2 - y^2) \end{array}$
E,	$\int 1 \varepsilon \varepsilon^* = 1 - 1 - \varepsilon - \varepsilon^* - 1$		$x^2 - y^2$
T <sub>e</sub> T <sub>u</sub>	$\begin{array}{cccccccccccccccccccccccccccccccccccc$	$(R_x, R_y, R_z) (x, y, z)$	(xz, yz, xy)
$T_d$	$E  8C_3  3C_2  6S_4  6\sigma_d$		
$A_1$ $A_2$		$x^2 + y^2 + z^2$	
E	$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$	$(2z^2 - x^2 - y^2),$ $x^2 - y^2)$	
$T_1 \\ T_2$	$\begin{vmatrix} 3 & 0 & -1 & 1 & -1 \\ 3 & 0 & -1 & -1 & 1 \\ \end{vmatrix} \begin{pmatrix} (R_x, R_y, R_z) \\ (x, y, z) \\ \ddots \\ \end{vmatrix}$		
0	$\begin{vmatrix} E & 6C_4 & 3C_2(=C_4^2) & 8C_3 & 6C_2 \end{vmatrix}$	(xy, xz, yz)	
A 1			$x^2 + y^2 + z^2$
A 2 E	$ \begin{array}{cccccccccccccccccccccccccccccccccccc$		$\begin{array}{l} (2z^2 - x^2 - y^2, \\ x^2 - y^2) \end{array}$
$T_1 \\ T_2$	$\begin{vmatrix} 3 & 1 & -1 & 0 & -1 \\ 3 & -1 & -1 & 0 & 1 \end{vmatrix} (R_x, R_y)$	$(R_{z}); (x, y, z)$	• •
0,	$\begin{bmatrix} E & 8C_3 & 6C_2 & 6C_4 & 3C_2(=C_4^2) & i & 6S_4 \end{bmatrix}$	85 <sub>6</sub> 3an 6an	(xy, xz, yz)
A 10	1 1 1 1 1 1	1 1 1	$x^2 + y^2 + z^2$
A 20 Eg	$\begin{vmatrix} 1 & 1 & -1 & -1 & 1 & 1 & -1 \\ 2 & -1 & 0 & 0 & 2 & 2 & 0 \end{vmatrix}$	$\begin{vmatrix} 1 & 1 & -1 \\ -1 & 2 & 0 \end{vmatrix}$	$(2z^2 - x^2 - y^2, x^2 - y^2).$
$T_{1g}$	3 0 -1 1 -1 3 1	0 -1 -1	$(R_x, R_y, R_z)$
T <sub>29</sub> A <sub>11</sub>		$ \begin{array}{cccccccccccccccccccccccccccccccccccc$	(xz, yz, xy)
A <sub>24</sub> E <sub>4</sub>	2 -1 0 0 2 -2 0	$   \begin{bmatrix}     -1 & -1 & 1 \\     1 & -2 & 0   \end{bmatrix} $	
T <sub>1#</sub> T <sub>2#</sub>	$\begin{vmatrix} 3 & 0 & -1 & 1 & -1 & -3 & -1 \\ 3 & 0 & 1 & -1 & -1 & -3 & 1 \end{vmatrix}$	$\begin{array}{cccc} 0 & I & 1 \\ 0 & 1 & -1 \end{array}$	(x, y, z)



### **Table 6.5** The $C_{3v}$ character table



W. H.

### Character table for $C_{\scriptscriptstyle \! \infty \nu}$ point group

	E	2C <sub></sub>	••••	∞ σ <sub>v</sub>	linear, rotations	quadratic
A <sub>1</sub> =Σ <sup>+</sup>	1	1		1	Z	x <sup>2</sup> +y <sup>2</sup> , z <sup>2</sup>
A <sub>1</sub> =Σ <sup>+</sup> A <sub>2</sub> =Σ <sup>-</sup>	1	1		-1	R <sub>z</sub>	
Е <sub>1</sub> =П	2	2cos(Φ)		0	(x, y) (R <sub>x</sub> , R <sub>y</sub> )	(xz, yz)
E <sub>2</sub> =Δ	2	2cos(2φ)		0		(x <sup>2</sup> -y <sup>2</sup> , xy)
Е <sub>3</sub> =Ф	2	2cos(3φ)		0		
•••	•••	•••	•••	•••		

### Character table for $D_{{\scriptscriptstyle \infty}{\scriptscriptstyle h}}$ point group

	E	2C <sub></sub>	 ∞0 <sup>v</sup>	i	2S <sub>∞</sub>	 ∞C'2	linear functions, rotations	quadratic
Α <sub>1g</sub> =Σ+ <sub>g</sub>	1	1	 1	1	1	 1		x <sup>2</sup> +y <sup>2</sup> , z <sup>2</sup>
A <sub>2g</sub> =Σ <sup>-</sup> g	1	1	 -1	1	1	 -1	R <sub>z</sub>	
E <sub>1g</sub> =Π <sub>g</sub>	2	2cos(φ)	 0	2	-2cos(φ)	 0	(R <sub>x</sub> , R <sub>y</sub> )	(xz, yz)
E <sub>2g</sub> =Δ <sub>g</sub>	2	2cos(2φ)	 0	2	2cos(2φ)	 0		(x²-y², xy)
E <sub>3g</sub> =Φ <sub>g</sub>	2	2cos(3ф)	 0	2	-2cos(3ф)	 0		
•••			 			 		
Α <sub>1u</sub> =Σ+ <sub>u</sub>	1	1	 1	-1	-1	 -1	z	
A <sub>2u</sub> =Σ <sup>-</sup> <sub>u</sub>	1	1	 -1	-1	-1	 1		
Е <sub>1u</sub> =П <sub>u</sub>	2	2cos(φ)	 0	-2	2cos(φ)	 0	(x, y)	
E <sub>2u</sub> =Δ <sub>u</sub>	2	2cos(2ф)	 0	-2	-2cos(2ф)	 0		
Е <sub>зи</sub> =Ф <sub>и</sub>	2	2cos(3φ)	 0	-2	2cos(3φ)	 0		
•••			 			 		