

PROBLEM 11 a-f)

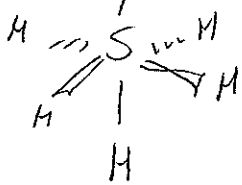
Qh

BASIS SET:

6M : 6S

FROM TABLE

$A_{1g} + T_{1u}$



ρ_h	E	$8C_3$	$6C_2$	$6C_4$	$3C_2$	i	$6S_4$	$8S_6$	$3\sigma_h$	$6\sigma_d$
A_{1g}	1	1	1	-1	1	1	1	1	1	1
A_{2g}	1	1	-1	-1	1	1	-1	1	1	-1
E_g	2	-1	0	0	2	2	0	-1	2	0
T_{1g}	3	0	-1	1	-1	3	1	0	-1	-1
T_{2g}	3	0	1	-1	-1	3	-1	0	-1	1
A_{1u}	1	1	1	1	1	-1	-1	-1	-1	-1
A_{2u}	1	1	-1	-1	1	-1	1	-1	-1	1
E_u	2	-1	0	0	2	-2	0	1	-2	0
T_{1u}	3	0	-1	1	-1	-3	-1	0	1	1
T_{2u}	3	0	1	-1	-1	-3	1	0	1	-1
$\Gamma_{H_{1s}A}$	6	0	0	2	2	0	0	0	4	2

$$\alpha_{A_{1g}} = \frac{1}{48} (6 \cdot 1 \cdot 1 + 0 + 0 + 2 \cdot 1 \cdot 6 + 2 \cdot 1 \cdot 3 + 0 + 0 + 0 + 4 \cdot 1 \cdot 3 + 2 \cdot 1 \cdot 6)$$

$$= \frac{1}{48} (6 + 12 + 12 + 12)$$

$$= 1$$

$$\alpha_{E_g} = \frac{1}{48} (6 \cdot 2 \cdot 2 + 0 + 0 + 0 + 2 \cdot 2 \cdot 3 + 0 + 0 + 0 + 4 \cdot 2 \cdot 3 + 0)$$

$$= \frac{1}{48} (12 + 12 + 24)$$

$$= 1$$

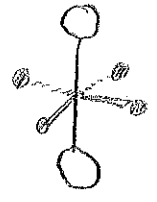
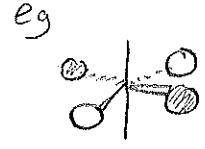
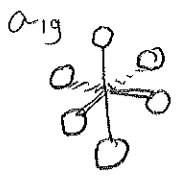
$$\alpha_{T_{1u}} = \frac{1}{48} (6 \cdot 3 \cdot 1 + 0 + 0 + 2 \cdot 1 \cdot 6 + 2 \cdot -1 \cdot 3 + 0 + 0 + 0 + 4 \cdot 1 \cdot 3 + 2 \cdot 1 \cdot 6)$$

$$= \frac{1}{48} (18 + 12 - 6 + 12 + 12)$$

$$= 1$$

$$\Gamma_{H_{1s}A} = A_{1g} + E_g + T_{1u}$$

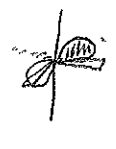
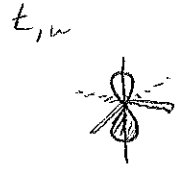
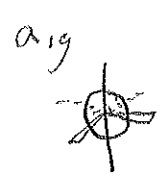
SH₆ (A) SALCS



t_{1u}



5 VALENCE ORBITALS



s

p_z

p_x

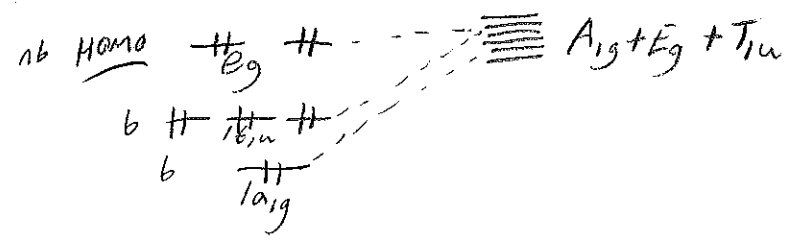
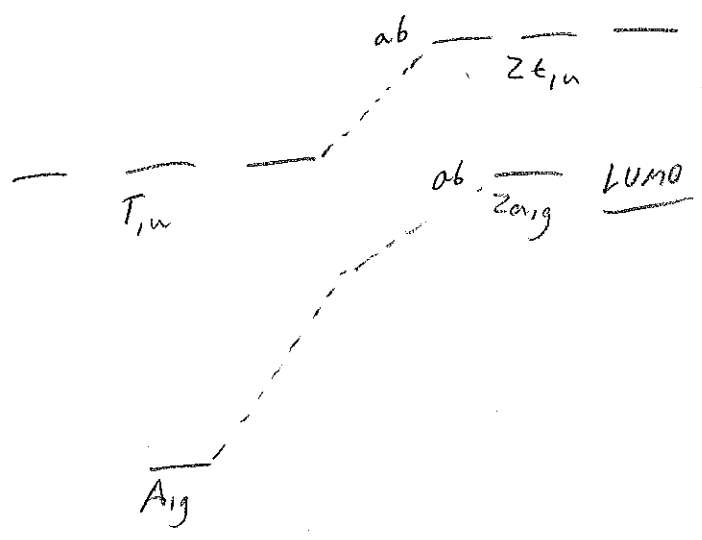
p_y

A MO:

S

SH₆

6H



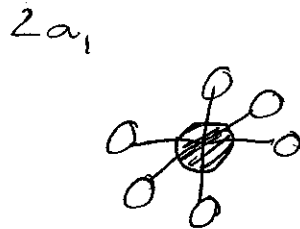
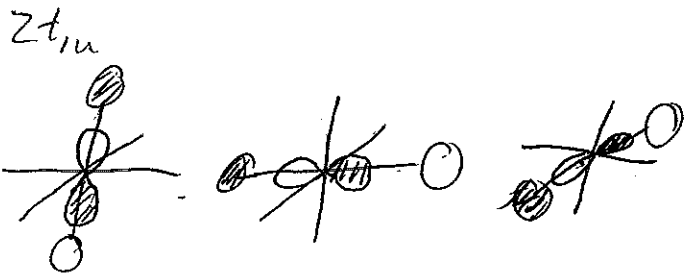
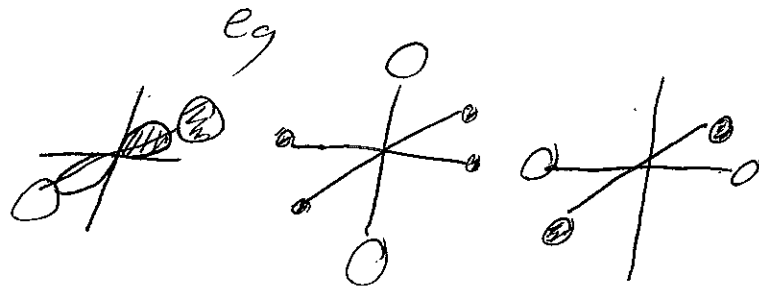
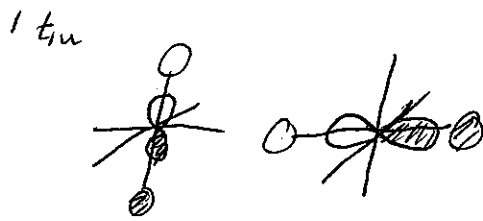
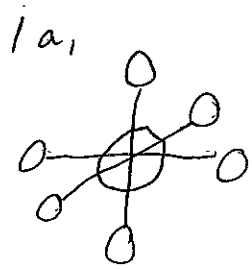
b = BONDING
nb = NON BONDING
ab = ANTIBONDING

PES: 3 PEAKS

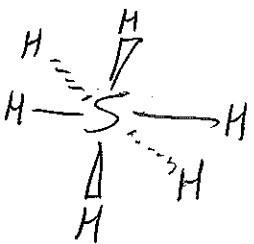
BC: $\frac{8}{2} = 4$

DIAMAGNETIC

SH₆ (Td) MOs



B



D_{3h}

BASIS SET S: S, P ⇒ E' + A₂'' + A₁'
 6M: 6S

D _{3h}	E	2C ₃	3C ₂	σ _h	2S ₃	3σ _v
A ₁ '	1	1	1	1	1	1
A ₂ '	1	1	-1	1	1	-1
E'	2	-1	0	2	-1	0
A ₁ ''	1	1	1	-1	-1	1
A ₂ ''	1	1	-1	-1	-1	-1
E''	2	-1	0	-2	1	0
∑ 6M vs B	6	0	0	0	0	2

$$\alpha_{A_1'} = \frac{1}{12} (6 \cdot 1 \cdot 1 + 0 + 0 + 0 + 0 + 2 \cdot 1 \cdot 3)$$

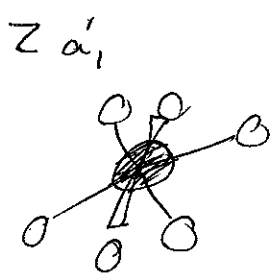
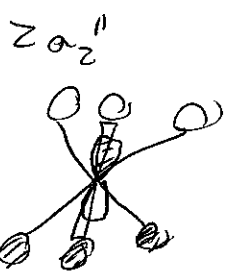
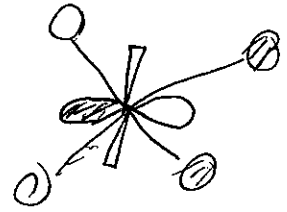
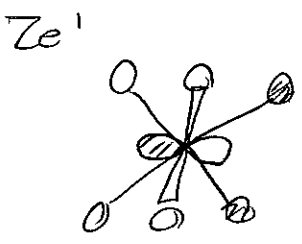
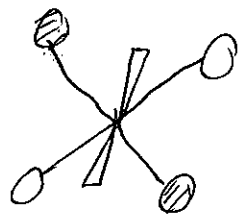
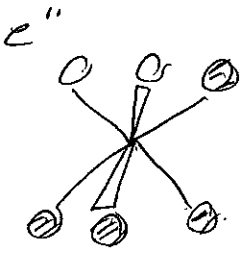
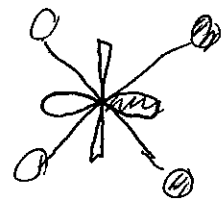
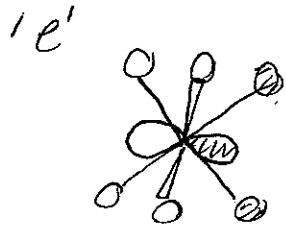
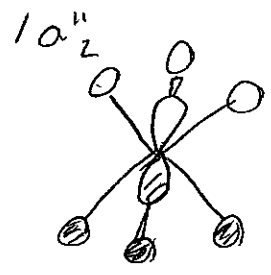
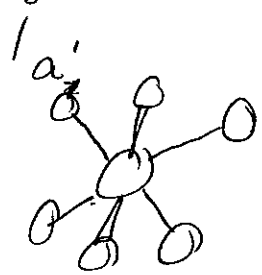
$$= \frac{1}{12} (6 + 6)$$

$$= 1$$

$$\alpha_{E'} = \frac{1}{12} (6 \cdot 2 \cdot 1 + 0 + 0 + 0 + 0 + 2 \cdot 0 \cdot 3)$$

$$= 1$$

SH_6 (B) Mos



g) $A_{1g} + E_g + T_{1g} + T_{2g} + 3T_{1u} + T_{2u}$

TRANSLATION: T_{1u}

ROTATION: T_{1g}

NORMAL VIBRATION: $A_{1g} + E_g + T_{2g} + 2T_{1u} + T_{2u}$

h) IR SPECTRUM: 2 PEAKS ($2 T_{1u}$)

RAMAN SPECTRUM: 3 PEAKS ($A_{1g} + E_g + T_{2g}$)

i)

O_h	E	$8C_3$	$6C_2$	$4C_4$	$3C_2$	i	$6S_4$	$8S_6$	$3\sigma_h$	$6\sigma_d$	
Γ_{S-H}	6	0	0	2	2	0	0	0	4	2	$= A_{1g} + E_g + T_{1u}$

D_{3h}	E	$2C_3$	$3C_2$	σ_h	$2S_6$	$3\sigma_v$	
Γ_{S-H}	6	0	0	0	0	2	$= A_1' + E' + A_2'' + E''$

j) ISOMER A: 1 IR S-H STRETCHES (T_{1u})

2 RAMAN S-H STRETCHES ($A_{1g} + E_g$)

ISOMER B: 2 IR S-H STRETCHES ($E' + A_2''$)

3 RAMAN S-H STRETCHES ($A_1' + E'' + E'$)

2 PART 1

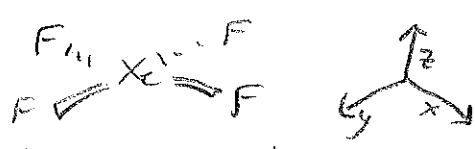
a)

O_h	E	$8C_3$	$6C_2$	$6C_4$	$3C_2$	i	$6S_4$	$8S_6$	$3\sigma_h$	$6\sigma_d$	
$\Gamma_{B_{2z}}$	6	0	0	2	2	0	0	0	4	2	$= A_{1g} + E_g + T_{1u}$

b)

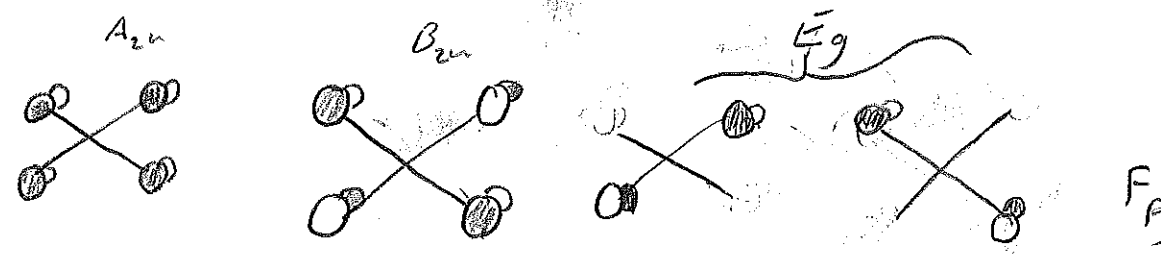
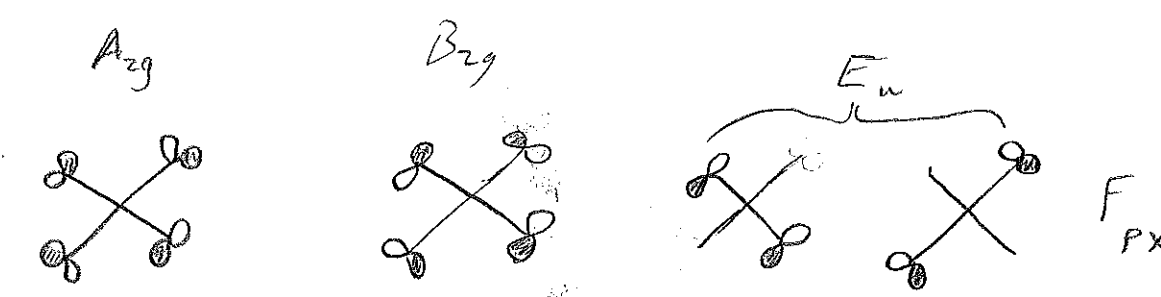
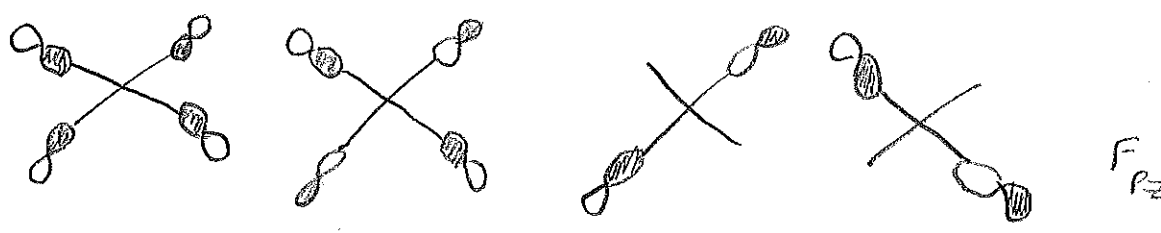
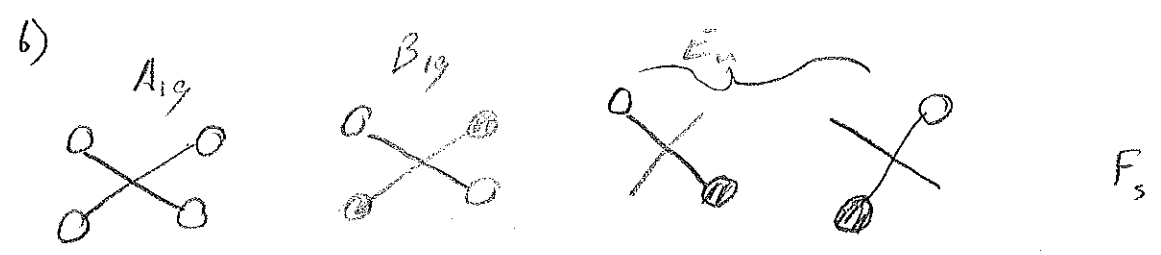
O_h	E	$8C_3$	$6C_2$	$6C_4$	$3C_2$	i	$6S_4$	$8S_6$	$3\sigma_h$	$6\sigma_d$	
$\Gamma_{B_{xy}}$	12	0	0	0	-4	0	0	0	0	0	$= T_{1g} + T_{2g} + T_{1u} + T_{2u}$

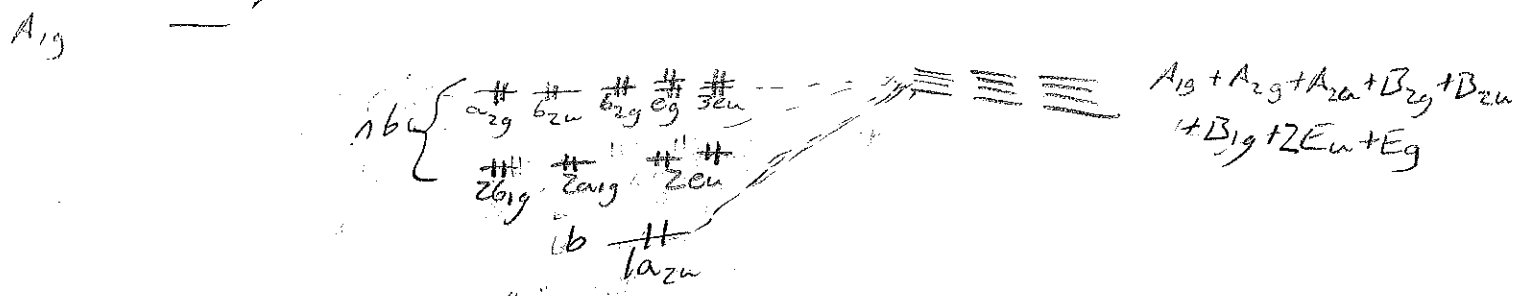
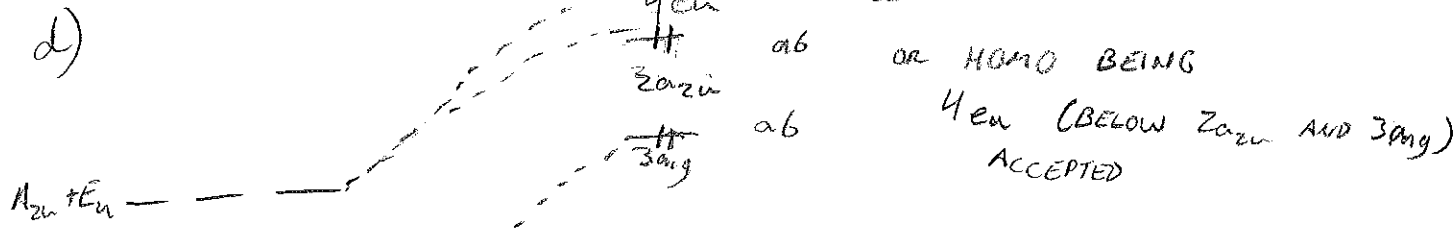
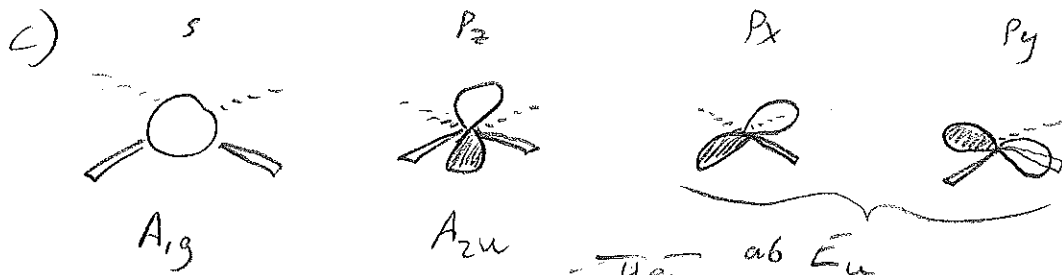
2 PART II



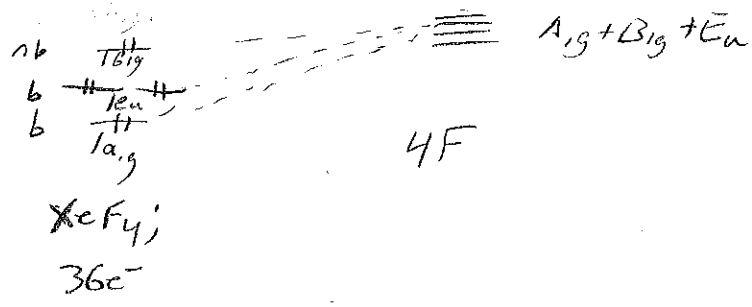
a)

D_{4h}	E	$2C_4$	C_2	$2C_2'$	$2C_2''$	i	$2S_4$	σ_h	$2\sigma_v$	$2\sigma_d$	
Γ_{F_s}	4	0	0	2	0	0	0	4	2	0	$= A_{1g} + B_{1g} + E_u$
$\Gamma_{F_{Pz}}$	4	0	0	2	0	0	0	4	2	0	$= A_{1g} + B_{1g} + E_u$
$\Gamma_{F_{Px}}$	4	0	0	-2	0	0	0	4	-2	0	$= E_u + B_{2g} + A_{2g}$
$\Gamma_{F_{Py}}$	4	0	0	-2	0	0	0	-4	2	0	$= A_{2u} + B_{2u} + E_g$

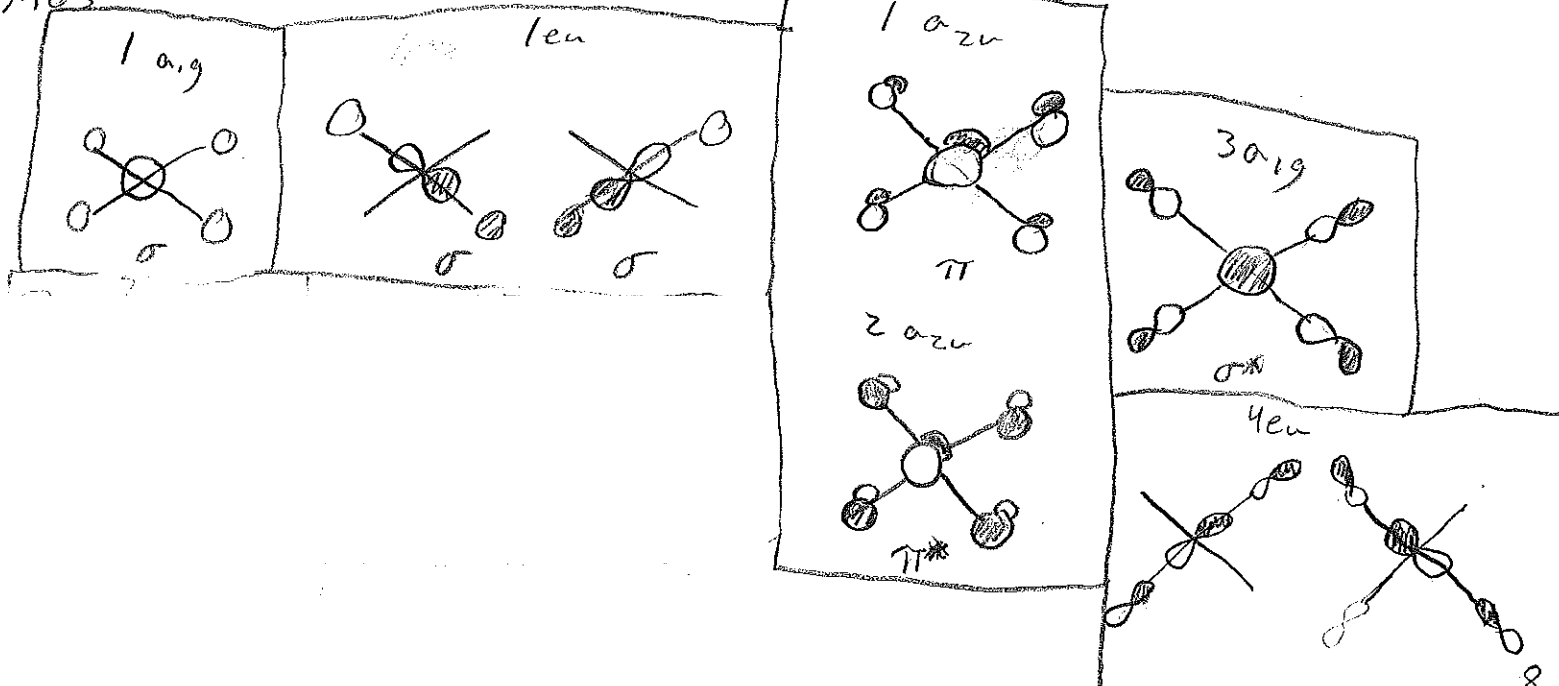




Xe



Mos



#2 f)

IR PEAKS : 141.8 cm^{-1}
 240.0 cm^{-1}
 522.1 cm^{-1}

VALUES DEPEND
ON BASIS SET / FUNCTIONAL

RAMAN PEAKS: 187.4 cm^{-1}
 426.9 cm^{-1}
 446.0 cm^{-1}

STRETCHING MODES FOR RAMAN PEAK AT 446 cm^{-1} , 426.9 cm^{-1}
IR PEAK AT 522.1 cm^{-1}

#3

a)

20 - A_{1g}	21 - E_u	22 - E_u	23 - B_{1g}	24 - A_{1g}
25 - A_{2u}	26 - A_{1g}	27 - E_u	28 - E_u	29 - A_{2u}
30 - B_{2g}	31 - E_g	32 - E_g	33 - B_{1g}	34 - E_u
35 - E_u	36 - B_{2u}	37 - A_{2g}	38 - A_{2u}	39 - A_{1g}
40 - E_u	41 - E_u	42 - E_g	43 - E_g	44 - A_{1g}

b) FLUORINE IS MORE ELECTRONEGATIVE THAN CHLORINE, SO ITS ORBITALS ARE LOWER IN ENERGY. F-BASED MOs ARE THEREFORE LOWER IN ENERGY THAN Cl BASED MOs

c) 33 - B_1
40 - E