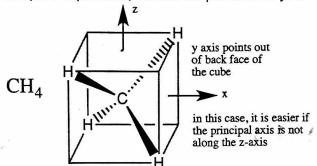
SALC/MO DIAGRAM EXERCISE 2

More SALC exercises! In the interest of time, I have done the first few steps for you so you can focus on figuring out what each SALC looks like and draw the MO diagram. It is up to you to make sure you can get to the same answers that I come to for the first few steps.

Note: for SALC derivation, the rule for character assignments when determining Γ_{orbital} are: +1 if orbital stays in same place, same phase, -1 if orbitalsstays in same place, changes phase, 0 if orbital moves

CH₄: In this example, we define the x,y,and z axes slightly differently-the z axis in this case is NOT along the principle rotation axis. If different axis orientations are required to easily complete a problem, this will be provided to you.



Character table for T _d point group								
	E	8C ₃	3C ₂	6S ₄	6σ _d	linear, rotations	quadratic	
A ₁	1	1	1	1	1		$x^2+y^2+z^2$	
A ₂	1	1	1	-1	-1			
E	2	-1	2	0	0		$(2z^2-x^2-y^2, x^2-y^2)$	
T ₁	3	0	or live	<i>§</i> 1	-1	(R_x, R_y, R_z)		
T ₂	3	0	-1	-1	1	(x, y, z)	(xy, xz, yz)	

- 1) Identify point group: T_d
- 2) Identify relevant orbitals:

Valence orbitals of C? 2s, 2px, 2py, 2pz

Valence orbitals of H? 1s

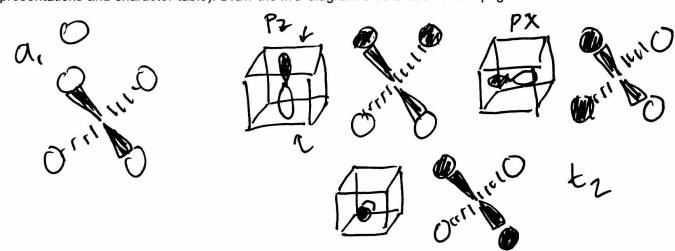
3) Determine Γ_{orbital} for all of the relevant valence orbitals of the H's in CH₄ In CH₄, this means we only have to do the 'gamma' for the 1s orbitals of the H's.

$$T_d$$
 E 8C₃ 3C₂ 6S₄ 6 σ_d
 Γ_{1s} = 4 1 0 0 2

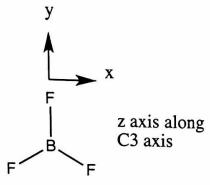
4) Use the reduction formula to find irreducible representations

$$\Gamma_{1s} = A_1 + T_2$$

5) Using the above results, draw the 4 SALCs of the H atoms in CH₄, matching their symmetry to the symmetry of the appropriate orbitals on the central atom (as determined by the irreducible representations and character table). Draw the MO diagram on the back of this page.



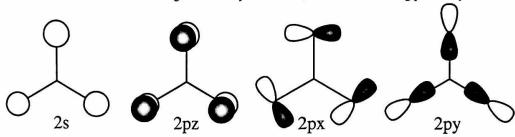
1 2px 2px 2px 2px 2px 1<u>v</u> 25 a, tz



Character table for D_{3h} point group

	E	2C ₃	3C'2	$\sigma_{\mathbf{b}}$	2S ₃	3σ _v	linear, rotations	quadratic
A'1	1	1	1	,1	1	1		x^2+y^2, z^2
A'2	1	1	-1	1	1	-1	R _z	
E'	2	-1	0	2	-1	0	(x, y)	(x^2-y^2, xy)
A'' ₁	1	1	1	-1	-1	-1		
A''2	1	1	-1	-1	-1	1	z	
E"	2	-1	0	-2	'n	0	(R_x, R_y)	(xz, yz)

orbital orientations used for deriving reducible representations, use these as a starting point for your SALCs



BF₃: Note the orientations of the py and px orbitals.

- 1) Identify point group: D_{3h}
- 2) Identify relevant orbitals:

Valence orbitals of F? 2s, 2px, 2py, 2pz

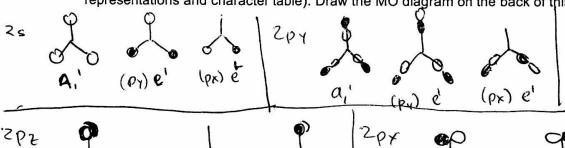
Valence orbitals of B? 2s, 2px, 2py, 2pz

3) Determine Γ_{orbital} for all of the relevant valence orbitals of the H's in CH₄ In CH₄, this means we only have to do the 'gamma' for the 1s orbitals of the H's.

D_{3h}	E	2C ₃	3C'2	σ_{h}	2S ₃	$3\sigma_{\rm v}$
$\Gamma_{2s} =$	3	0	1	3	0	1
$\Gamma_{2py} =$	3	0	1	3	0 ′	1
	3	0	-1	-3	0	1
	3	0	-1	3	0	Water Street

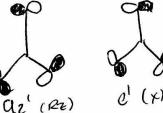
4) Use the reduction formula to find irreducible representations

 $\Gamma_{2s} = A_1' + E'$ $\Gamma_{2py} = A_1' + E'$ $\Gamma_{2pz} = A_2'' + E''$ $\Gamma_{2pz} = A_2'' + E''$ $\Gamma_{2px} = A_2' + E''$ $\Gamma_{2px} = A_2' + E''$ $\Gamma_{2px} = A_2' + E''$ $\Gamma_{2px} = A_2'' + E''$ $\Gamma_{2px} = A_$ to the symmetry of the appropriate orbitals on the central atom (as determined by the irreducible representations and character table). Draw the MO diagram on the back of this page.











see also el mo MFT Rigures energy levels 5.31 and 5.32 are mickey, e' 2 d from B, 4 e' from 3F ai' =) 2 bonding, 2 almost nonbonding (in between energy) 92" 2 antibonding =) sets of 2 because 202 2px 2py 2 P SALCS more strang ai' nan 25 closer in évergy) 11 ei ei 71 (non bonding ZS SALCE 'a,' essentially) 24 electrons to ta)