

LECTURE 29/36

NOV/15/02

SQUARE-PLANAR  
TETRAHEDRAL

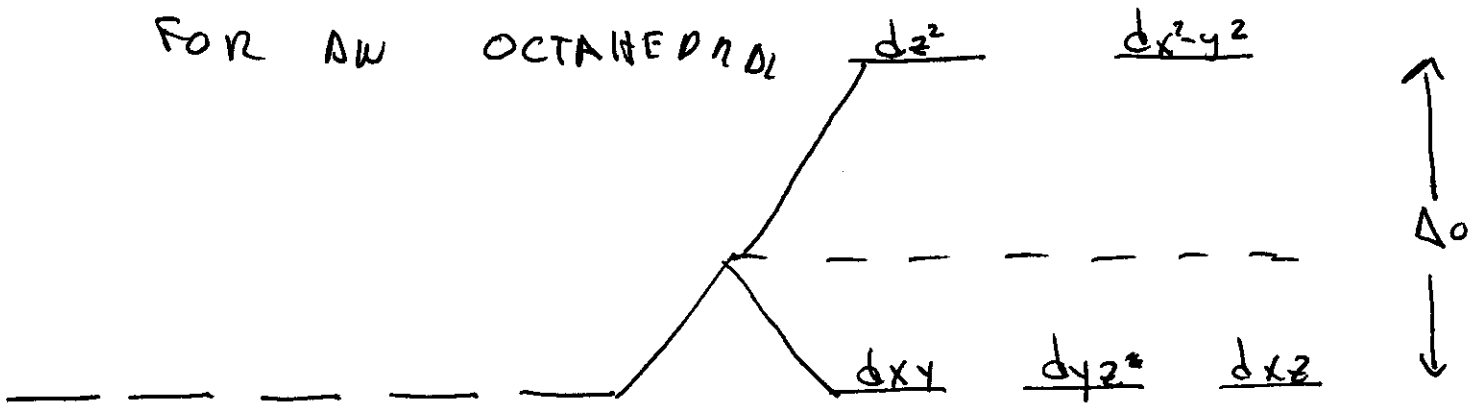
LIGAND FIELD THEORY

READ: TODAY'S LEC GRAY 159-183

NEXT LEC VOL II CTC - 445-468

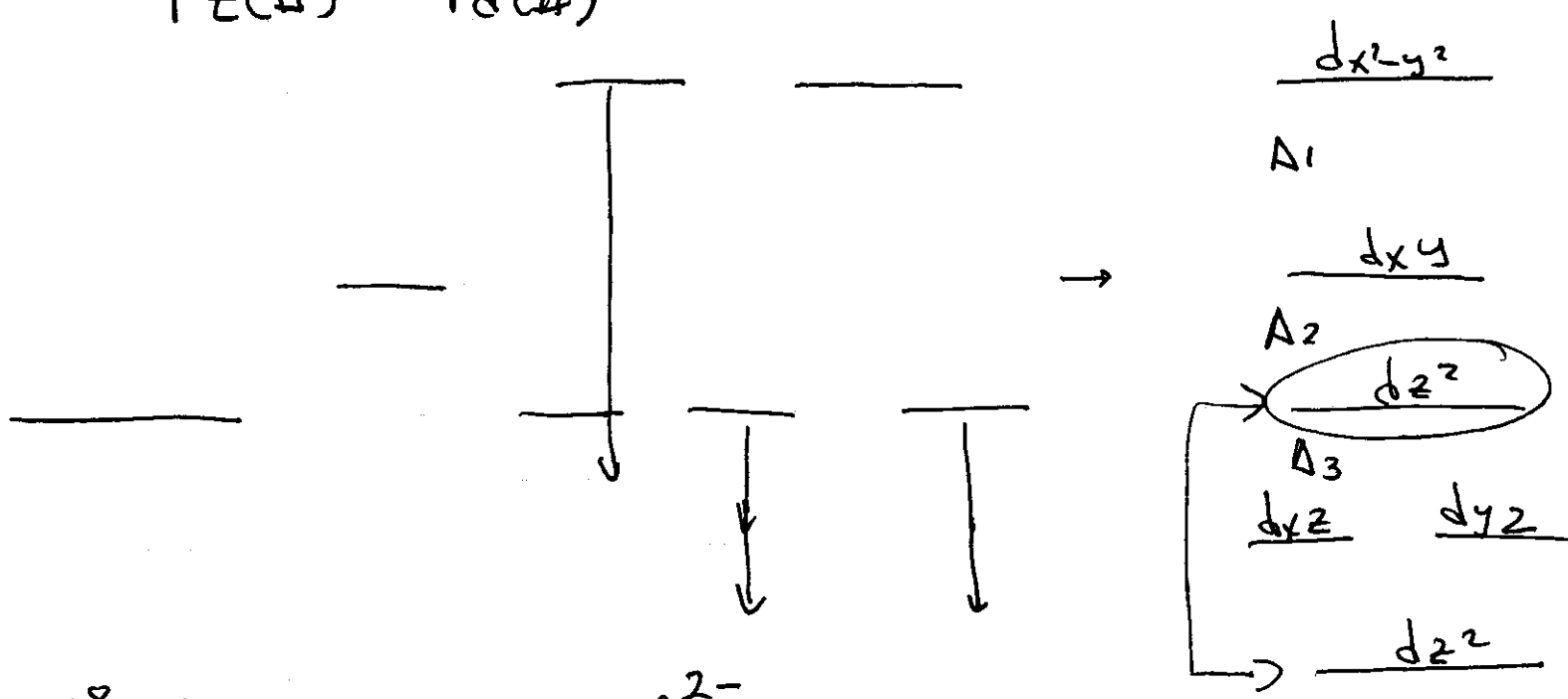
DUE	DATES:	NOV - 18	LAB REPORT
		NOV - 25	PS 7
		DEC - 2	REPORT
		DEC - 6	PS 8

FOR AN OCTAHEDRAL  $d_1$

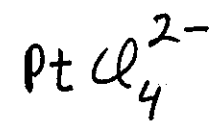


FOR SQUARE-PLANAR COMPLEXES

Pt(II) Pd(II)



$d^8$  low-spin

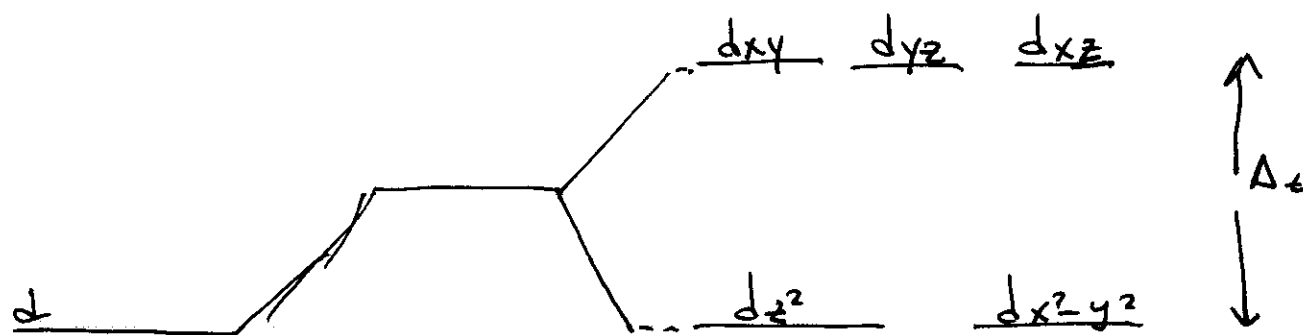


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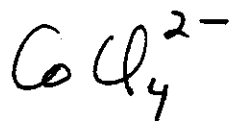
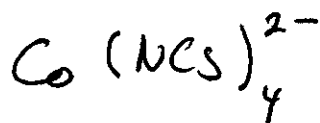
OCTAHEDRAL COMPOUNDS ARE MORE COMMON AND MORE STABLE (6 LIGANDS)

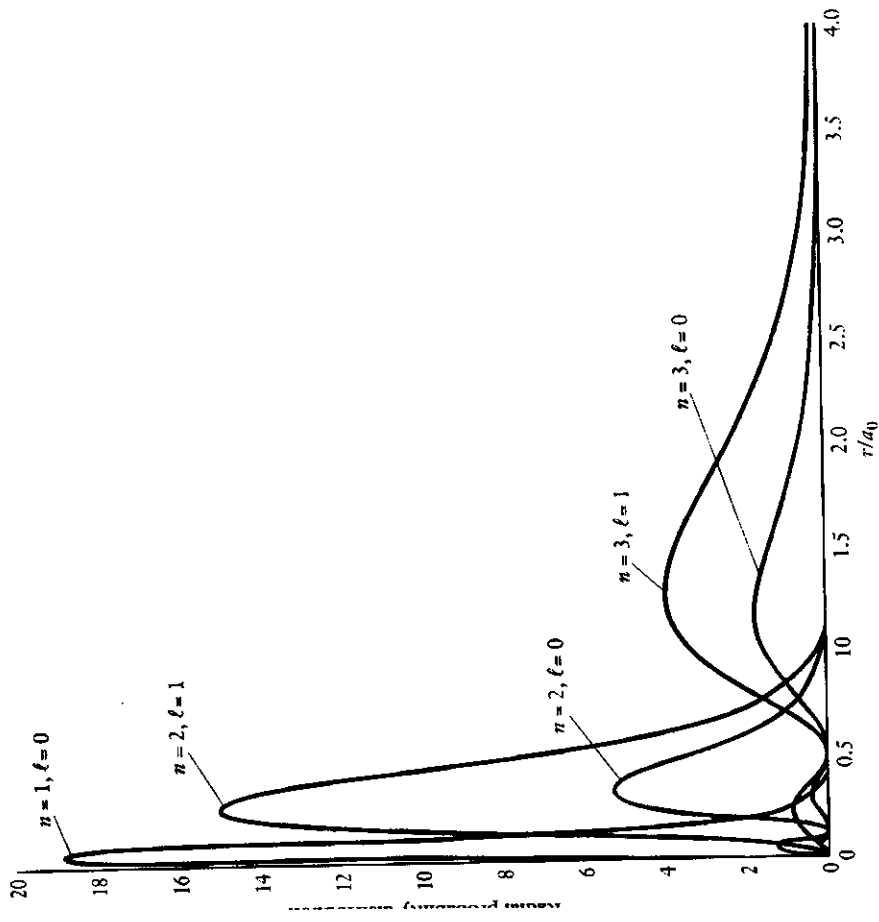
TETRAHEDRAL

POSITION OF THE LIGANDS OPPOSITE TO OCTAHEDRAL



LESS COMMON





**FIGURE 15.34** The shell structure of an argon atom calculated by Hartree's method. The fact that the radial probability for all orbitals with the same value of  $n$  have maxima very near one another suggests that the electrons can be viewed in "shells" made up of these orbitals.

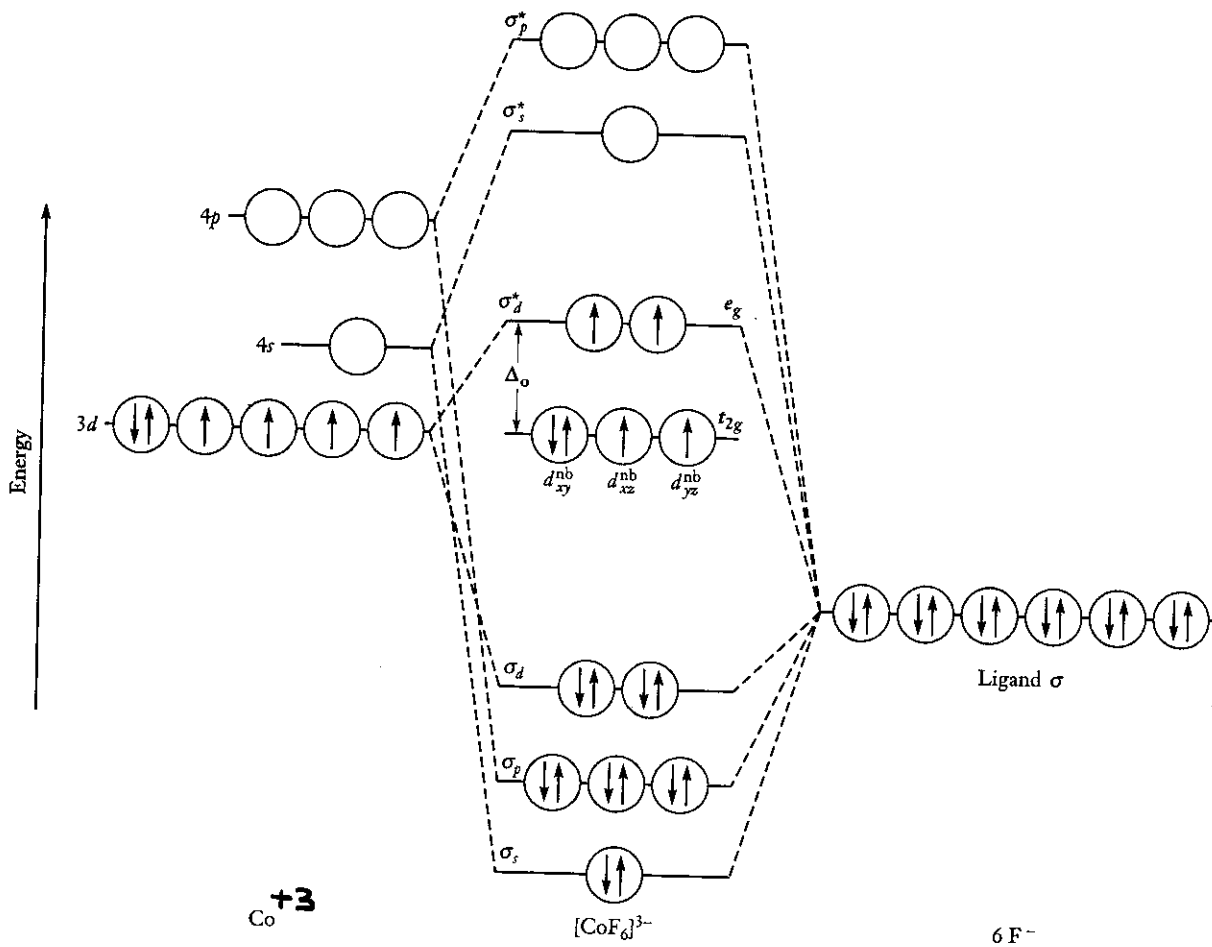


FIGURE 18.22 Orbital correlation diagram for an octahedral ligand field, showing the energy-level filling for a [CoF<sub>6</sub>]<sup>3-</sup> ion.

FOR  $I^-$   $Br^-$   $Cl^-$   $F^-$   $OH^-$

THE NON BONDING d ORBITALS INTERACT w/  $\sigma^-$  IN p ORBITALS  $\perp$  TO THE METAL - LIGAND AXIS

$d_{xy}$  AND  $p_y \Rightarrow d_{xy} d_{xz} d_{yz} \uparrow$

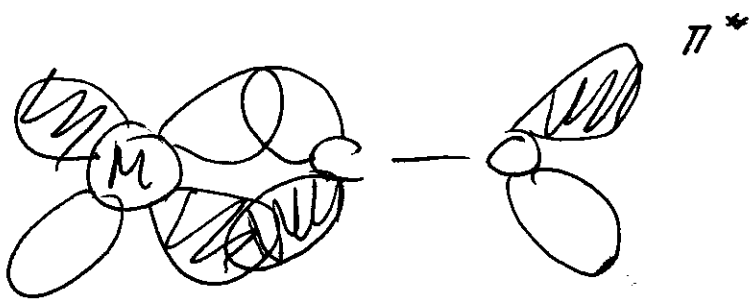
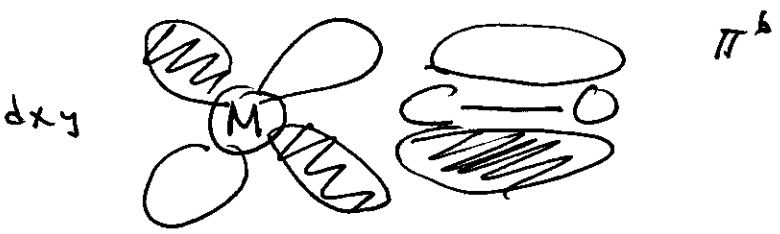
$\Rightarrow$  WEAK-FIELD LIGANDS

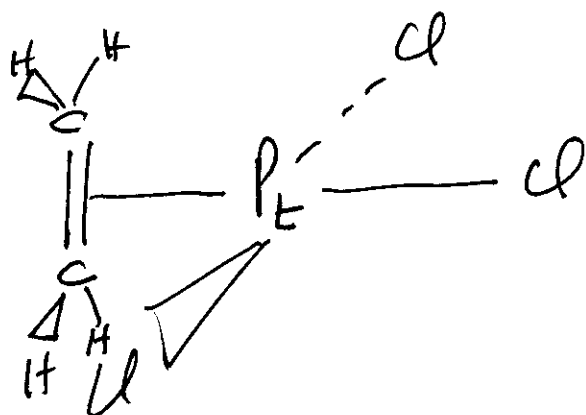


FOR  $CN^-$  AND  $CO$

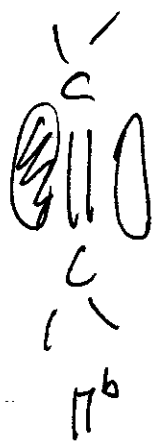
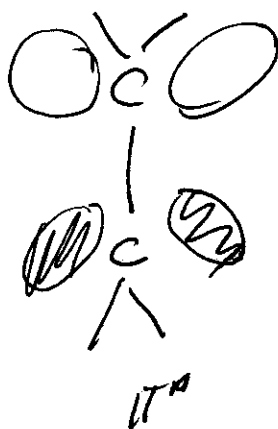
THE ANTI BONDING LIGAND OVERLAPS w/  $e_{2g}$

$\Rightarrow$   $\downarrow$  ENERGY HIGH STRONG-FIELD LIGANDS



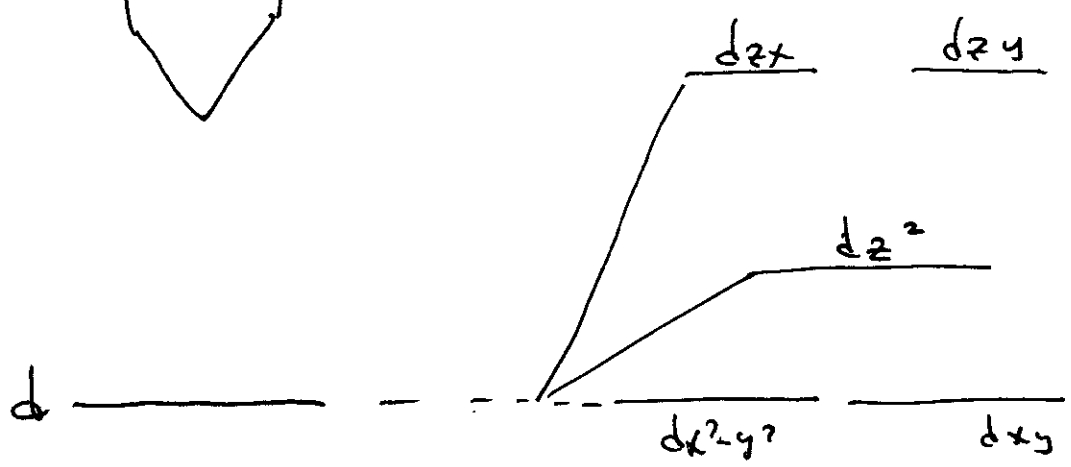
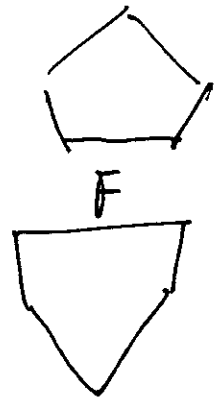
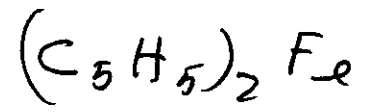
ORGANO METALLIC  $\pi$  COMPLEXES

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 $\sigma$  bond w Ptmetal  $d$   $\rightarrow$   $\pi^*$

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# FERROCENE



PORPHYRINS TETRA DENTATE  
 $Mg^{2+}$      $Fe^{2+}$      $Co^{3+}$      $Zn^{2+}$      $Ni^{2+}$

$Fe^{2+}$ ,  $Fe^{3+}$  → HEME

$Mg^{2+}$  → CHLOROPHYLL