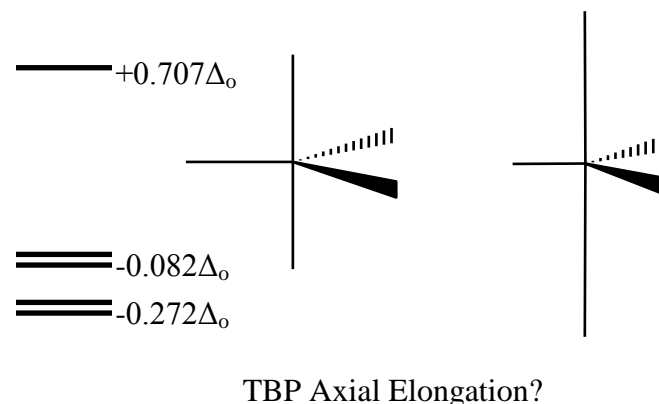
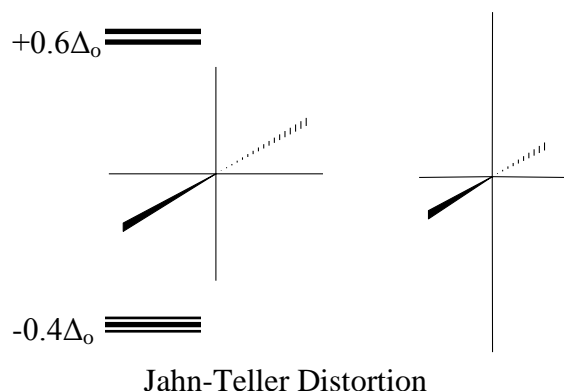


- Calculate the magnetic moment you would expect for each of the following complexes:  $[\text{Fe}(\text{CN})_6]^{3-}$ ,  $[\text{CoF}_6]^{3-}$ ,  $[\text{Cu}(\text{NH}_3)_4]^{2+}$ . Explain all answers and assumptions.
- In mixed-ligand complexes, sterically demanding ligands can distort the ideal geometry one might expect for a complex. Consider the two “octahedral” complexes  $\text{MF}_6$  and  $\text{MF}_3\text{I}_3$ . Describe any differences in the structure and crystal field splitting of these two complexes. Specify any assumptions you make and explain all your answers.
- Some octahedral complexes can undergo a Jahn-Teller distortion (axial elongation) to achieve a lower energy (more stable) state. Are there any d-electron counts that would favor an axial elongation in a trigonal bipyramidal complex? Show how you would expect the crystal field splitting to change when the TBP axial elongation occurs. Explain all answers and assumptions.



4. Describe the crystal field splitting you would expect for the f-orbitals pictured below in an octahedral geometry. Explain your splitting assignments. How would the splitting change if the octahedral complex lost a ligand to become a square pyramid? How would the splitting change if the square pyramid rearranged to form a trigonal bipyramid? Explain all assignments and assumptions. If you have trouble visualizing any of these orbitals, *ask!*  $f_2$ - $f_5$  have the same basic shape with different orientations;  $f_6$  and  $f_7$  have the same shape with different orientations,  $f_1$  is unique. There are a number of different versions of the f-orbitals, your assignments must be for *this* set of f-orbitals and must use *this* naming system.

