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Molecular Orbitals Of Transition Metal Complexes





Synopsis

This book starts with the most elementary ideas of molecular orbital theory and leads the reader progressively to an understanding of the electronic structure, geometry and, in some cases, reactivity of transition metal complexes. The qualitative orbital approach, based on simple notions such as symmetry, overlap and electronegativity, is the focus of the presentation and a substantial part of the book is associated with the mechanics of the assembly of molecular orbital diagrams. The first chapter recalls the basis for electron counting in transition metal complexes. The main ligand fields (octahedral, square planar, tetrahedral, etc.) are studied in the second chapter (sigma interactions) and the structure of the "d block" is used to trace the relationships between the electronic structure and the geometry of the complexes. The third chapter studies the change in analysis when the ligands have pi-type interactions with the metal. All these ideas are then used in the fourth chapter to study a series of selected applications of varying complexity (structure, reactivity). The fifth chapter deals with the "isolobal analogy" which points out the resemblance between the molecular orbitals of inorganic and organic species and provides a bridge between these two subfields of chemistry. The last chapter is devoted to a presentation of basic Group Theory with applications to some of the complexes studied in the earlier chapters.

Book Information

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Customer Reviews

This book was assigned for my introductory inorganic chemistry course at Caltech. Its strong points are that it covers most common inorganic complex shapes and derives their MO diagrams in a

simple and easy-to-follow way. The exercises are also relatively well-written and contain answers. The caveat is that you have to do some work on the side to understand what is going on, and some diagrams could be better listed. For example, the LGOs are shown 150+ pages away (in the appendices) from where their relevant groups are discussed, and a brief introduction to the Huckel method would have been appreciated. The source of character tables could have been explained a bit more clearly as well, and there is little to no underlying derivation of the group theory (the words dihedral group, etc. are never mentioned). But all in all, it's an OK introduction.

I highly recommend this book if you try to have insights into the details of structure and reactivity in organometallic chemistry and catalysis. Very helpful for further understanding of MO theory.

The book arrived in a timely manor and was in excellent condition. The price was good. I would recommend to others using the text for a class, or looking for a good reference book.

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