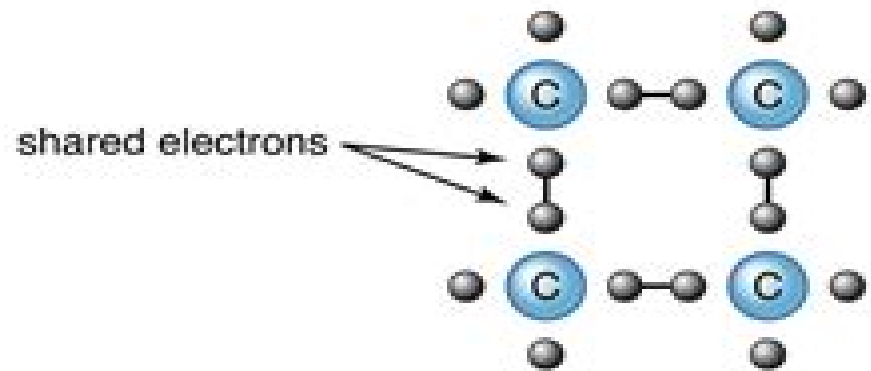
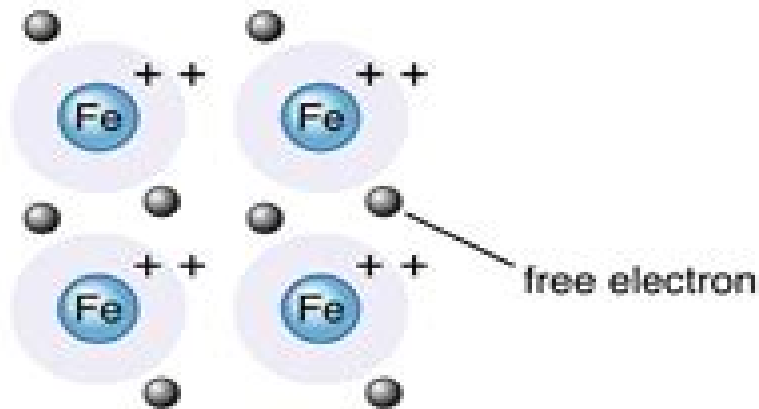


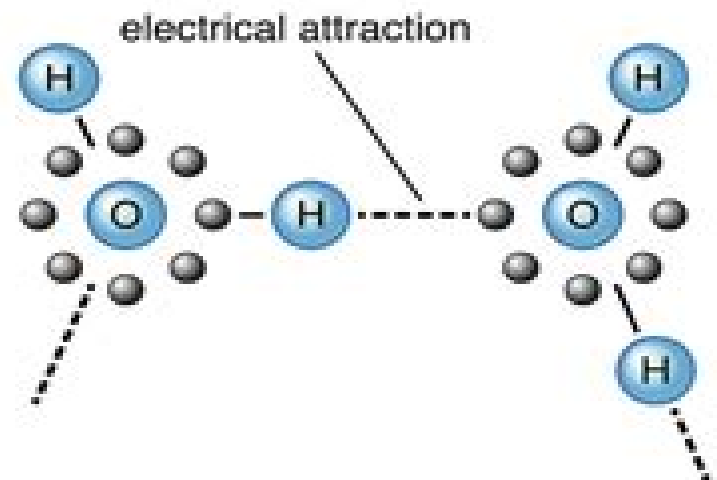
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# Electronic Structure And Chemical Binding

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principles and techniques of the contemporary computational and conceptual quantum chemistry are presented within both the wave function and electron density theories This background material is followed by a discussion of chemical concepts including stages of the bond formation processes chemical valence and bond multiplicity indices the hardness softness descriptors of molecules and reactants and general chemical reactivity stability principles The insights from Information Theory the basic elements of which are briefly introduced including the entropic origins and Orbital Communication Theory of the chemical bond are the subject of Part IV The importance of the non additive interference information tools in exploring patterns of chemical bonds and their covalent and ionic components will be emphasized      *Electronic Structure*

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**Modern Electronic Structure Theory And Applications In Organic Chemistry** Ernest R Davidson, 1997-10-31 This volume focuses on the use of quantum theory to understand and explain experiments in organic chemistry High level ab initio calculations when properly performed are useful in making quantitative distinctions between various possible interpretations of structures reactions and spectra Chemical reasoning based on simpler quantum models is however essential to enumerating the likely possibilities The simpler models also often suggest the type of wave function likely to be involved in ground and excited states at various points along reaction paths This preliminary understanding is needed in order to select the appropriate higher level approach since most higher level models are designed to describe improvements to some reasonable zeroth order wave function Consequently most of the chapters in this volume begin with experimental facts and model functions and then progress to higher level theory only when quantitative results are required In the first chapter Zimmerman discusses a wide variety of thermal and photochemical reactions of organic molecules Gronert discusses the use of ab initio calculations and experimental facts in deciphering the mechanism of elimination reactions in the gas phase Bettinger et al focus on carbene structures and reactions with comparison of the triplet and singlet states Next Hrovat and Borden discuss more general molecules with competitive triplet and singlet contenders for the ground state structure Cave explains the difficulties and considerations involved with many of the methods and illustrates the difficulties by comparing with the UV spectra of short polyenes Jordan et al discuss long range electron transfer using model compounds and model Hamiltonians Finally Hiberty discusses the breathing orbital valence bond model as a different approach to introducing the crucial correlation that is known to be important in organic reactions

**Unified Valence Bond Theory of Electronic Structure** N. D. Epitidis, 2012-12-06

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conceptual and interpretive aspects of the theory in line with the general philosophy adopted in the series The Chemical Bond Gernot Frenking, Sason Shaik, 2014-06-13 A unique overview of the different kinds of chemical bonds that can be found in the periodic table from the main group elements to transition elements lanthanides and actinides It takes into account the many developments that have taken place in the field over the past few decades due to the rapid advances in quantum chemical models and faster computers This is the perfect complement to Chemical Bonding Fundamentals and Models by the same editors who are two of the top scientists working on this topic each with extensive experience and important connections within the community

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