

CARBOCYCLIC CAGE COMPOUNDS

CHEMISTRY AND APPLICATIONS

EIJI OSAWA - OSAMU YONEMITSU



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Carbocyclic Cage Compounds Eiji Osawa, Osamu Yonemitsu, 1992
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Organic Conformational Analysis and Stereochemistry from Circular Dichroism Spectroscopy David A. Lightner, Jerome E. Gurst, 2000-03-31
A unique guide to variable temperature CD spectroscopy and its application in organic chemistry This timely original thought provoking work looks at organic stereochemistry from the perspective of circular dichroism CD using variable temperature CD spectroscopy to determine the conformation or absolute configuration of chiral molecules With an emphasis on the analysis of optically active ketones and the carbonyl chromophore the authors demonstrate the advantages of this highly sensitive spectroscopic tool for obtaining stereochemical information in diverse areas of organic chemistry biochemistry and medicinal pharmaceutical chemistry They combine detailed examples of stereochemical analysis with clear thorough presentations correlating chiroptical data with molecular mechanics calculations as well as data from NMR spectroscopy and other spectroscopic techniques In addition they provide a systematic survey of the professional literature featuring an extraordinary collection of original CD spectra run at varying temperatures Coverage includes Chiroptical measurements CD and ORD Optical Rotatory Dispersion Conformational analysis of compounds ranging from simple cyclic ketones to polycyclics Conjugated and homoconjugated systems Stereochemistry of the carbon carbon double bond Stereochemistry from exciton coupling of two or more chromophores An interesting historical account of the development of stereochemical concepts

Crystal Structure Analysis for Chemists and Biologists Jenny P. Glusker, Mitchell Lewis, Miriam Rossi, 1996-12-17
The fields of structural chemistry and biochemistry have blossomed in the last seventy years since X ray diffraction was discovered in 1912 Dorothy Hodgkin who obtained a Nobel Prize in 1965 for her X ray diffraction work wrote a great advantage of X ray analysis as a method of chemical structure analysis is its power to show some totally unexpected and surprising structure with at the same time complete certainty The results of all X ray diffraction studies are used by chemists and biochemists but these scientists need to be able to appreciate the significance and extent to which these results may be used A number of books written for practicing crystallographers cover the theory and applications of X ray diffraction but few are of real practical use to non specialists In Crystal Structure Analysis for Biologists and Chemists the general principles of crystal structure are presented in a highly readable way The book of Glusker who is internationally renowned provides good coverage of theory including data and understanding their significance

The CH/ π Interaction Motohiro Nishio, Minoru Hirota, Yoji Umezawa, 1998-06-10
A systematic exploration of the fascinating area of CH π interaction from the frontier of chemical research The CH π bond may be the weakest of the hydrogen bonds but it has been found to play important roles in the physical chemical and biological properties of a variety of substances This book offers a complete overview of the experimental basis theoretical background and nature of the CH π interaction to facilitate a clearer common understanding of the interaction between CH bonds and π systems Featured topics

include Evidence and methods of detection Quantum chemical treatment of CH π interaction Comparison with other types of weak molecular interactions Conformational consequences Chiroptical properties Selectivity in organic reactions Crystal structures of clathrates Inclusion compounds Interligand interactions in coordination chemistry Specific interactions in protein structures Prospects and possibilities for use in biochemistry and materials science The CH π Interaction makes an important contribution to the understanding of weak intermolecular forces and their potential value to the biochemical and pharmaceutical sciences It will be welcomed by physical and synthetic organic chemists structural chemists and pharmaceutical and biochemists *Perspectives on Structure and Mechanism in Organic Chemistry* Felix A.

Carroll, 2011-09-20 Helps to develop new perspectives and a deeper understanding of organic chemistry Instructors and students alike have praised *Perspectives on Structure and Mechanism in Organic Chemistry* because it motivates readers to think about organic chemistry in new and exciting ways Based on the author's first hand classroom experience the text uses complementary conceptual models to give new perspectives on the structures and reactions of organic compounds The first five chapters of the text discuss the structure and bonding of stable molecules and reactive intermediates These are followed by a chapter exploring the methods that organic chemists use to study reaction mechanisms The remaining chapters examine different types of acid base substitution addition elimination pericyclic and photochemical reactions This Second Edition has been thoroughly updated and revised to reflect the latest findings in physical organic chemistry Moreover this edition features New references to the latest primary and review literature More study questions to help readers better understand and apply new concepts in organic chemistry Coverage of new topics including density functional theory quantum theory of atoms in molecules Marcus theory molecular simulations effect of solvent on organic reactions asymmetric induction in nucleophilic additions to carbonyl compounds and dynamic effects on reaction pathways The nearly 400 problems in the text do more than allow students to test their understanding of the concepts presented in each chapter They also encourage readers to actively review and evaluate the chemical literature and to develop and defend their own ideas With its emphasis on complementary models and independent problem solving this text is ideal for upper level undergraduate and graduate courses in organic chemistry *Conformational Analysis of Molecules in Excited States* Jacek Waluk, 2000-06-13 A unique look at some of the hottest topics in photophysics and photochemistry today The study of molecules in excited states has exploded over the past decade providing new insights into conformational changes in organic molecules and opening up research opportunities for scientists and professionals in chemistry physics biology medicine and materials engineering Using conformational analysis as a unifying concept this important new work provides readers with a cohesive and cutting edge overview of this fascinating and challenging field From conformational changes accompanying photoinduced electron transfer to elementary photophysical and photochemical processes in living systems the most representative and challenging topics are carefully gleaned from the vast literature highlighting major conceptual problems along with the relevant

experimental techniques Authoritative detailed contributions from both experimentalists and theoreticians include coverage of Conformational changes in intramolecular excited state electron transfer Conformational aspects of excited state proton transfer The novel topic of solute solvent friction in chemical reactions Mechanisms and structural aspects of exciplex formations Conformational aspects of organic photochemistry Calculations of excited state conformational properties

Modern Conformational Analysis Helena Dodziuk,1996-12-17 Most books on organic conformational deal with stereochemistry while neglecting the spatial structure of organic molecules especially the unusual ones This ambitious monograph presents a global view of organic structure and conformational analysis in a very readable style The author s fascination with the beauty and variety of shapes of organic molecules comes through in her description of the importance of molecular organization in organic chemical and biological systems the quantum foundations of molecular spatial structure molecular symmetry chirality and physical methods for studying molecular geometry She provides ample bases for the new trends in organic conformational chemistry including the application of Computer Assisted Molecular Design CAMD to organic synthesis Strained Hydrocarbons Helena Dodziuk,2009-03-30 In clearly structured chapters this book covers the fascinating world of hydrocarbons providing an insight into the fundamental principles of chemistry The monograph covers modern aspects of the topic such as carbon nanotubes molecular flask inclusion and fullerenes with new synthetic procedures for the build up of the structural lattice included New Technical Books New York Public Library,1993

McGraw-Hill Encyclopedia of Science & Technology: Bio-cha ,2002 This resource provides in depth coverage of major scientific and technological developments It offers illustrated detailed coverage of the discoveries advances and milestones that continue to shape our lives *Bulletin of the Chemical Society of Japan* Nihon Kagakkai,1996

Aromaticity and Antiaromaticity Vladimir I. Minkin,Mikhail Nikolaevich Glukhovtsev,Mikhail N. Glukhovtsev,Boris I[A]kovlevich Simkin,1994-04-14 Designed to assist chemists in integrating the results of calculations on molecules and ions into their general body of chemical knowledge Contains recent contributions from theoretical and computational chemistry to the development of the concept of aromaticity antiaromaticity and its expansion into new areas such as organometallic and cluster compounds and three dimensional structures Updates the modern status of aromaticity and covers basic principles and experimental applications **Applications of Mass Spectrometry to Organic Stereochemistry** Janet S. Splitter,František Tureček,1994 **Conformational Behavior of Six-membered Rings** Eusebio Juaristi,1995

Carbon-13 NMR Chemical Shifts in Structural and Stereochemical Analysis Kalevi Pihlaja,Erich Kleinpeter,1994 A review of recent research on strategies and applications of the C 13 chemical shift a method for determining configuration of organic compounds Introduces C 13 NMR spectroscopy and describes conditions for collecting the FID for data handling and for obtaining a well resolved C 13 NMR spectrum as well as various substituent effect correlations their derivations and the origin of the effects Also discusses the use of multidimensional NMR methods For organic physical and natural products

chemists Includes bandw diagrams Annotation copyright by Book News Inc Portland OR **The Nuclear Overhauser Effect in Structural and Conformational Analysis** David Neuhaus, Michael P. Williamson, 2000-04-24 An authoritative review of the state of the art in the Nuclear Overhauser Effect essential information for organic chemists biochemists biophysicists and NMR spectroscopists The field of NMR spectroscopy has seen tremendous growth in the last twenty years particularly advances relating to Nuclear Overhauser Effect NOE spectroscopy the most powerful technique for obtaining structural information on molecules in solution Extensive and engaging the Second Edition of the leading reference on the NOE is significantly updated to reflect the latest changes and new approaches in the field Neuhaus and Williamson provide an essential guide to the complexities and use of the NOE in a readily accessible straightforward manner Their practical handbook features a new chapter addressing the use of NOE data to calculate biomolecular structures Chapters dealing with the kinetics of the NOE the effects of exchange and internal motion and applications of the NOE are also extensively revised Cross referenced in remarkable depth The Nuclear Overhauser Effect is organized into three main parts Part I describes the theory of the Nuclear Overhauser Effect in a clear comprehensive fashion Part II discusses the considerations involved in implementing NOE experiments including full coverage of all necessary details for both new and established techniques Part III offers examples of how the NOE is used including applications to defining molecular geometry stereochemistry conformation and biomolecular structure and interactions The Nuclear Overhauser Effect in Structural and Conformational Analysis Second Edition uniquely explains the NOE in detail making it an indispensable resource for the novice as well as the experienced NMR researcher Phosphorus-31 NMR Spectral Properties in Compound Characterization and Structural Analysis Louis D. Quin, John G. Verkade, 1994 Presents advances in ³¹P NMR and coupling effects of the past six years since VCH published its first jointly edited monograph on the subject A wide cross section of experts contribute a concise account of their use of ³¹P NMR in their own area of expertise giving major emphasis to the research results they have obtained since the earlier monograph appeared The volume is divided into nine sections chemical shift theory conformational analysis of phosphorus containing ring systems compounds with three four and five coordinate phosphorus low coordinate phosphorus compounds metal complexes of phosphorus ligands biological phosphorus systems mechanism elucidation in phosphorus chemistry analytical and interpretive aspects and NMR with solid phosphorus containing systems Annotation copyright by Book News Inc Portland OR **Carbon-rich Compounds** A. de Meijeres, 1998 Carbon Rich Compounds are defined here as carbon skeletons with a carbon to hydrogen ratio of 1 centro polyhydrindanes as well as oligoquinanes in particular highly unsaturated ones are discussed **Bibliography of Agriculture** , 1995

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