

PHYSICAL METHODS IN CHEMISTRY

DRAGO



Physical Methods For Chemistry Drago

K Morrison



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Technical Library, Pittsburgh, 1968 *Physical Methods for Chemists* Drago, Russell S. Drago, 1992-01-01

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Pharmaceutical and Medical Applications of Near-Infrared Spectroscopy Emil W. Ciurczak, Benoit Igne, 2014-12-15 Since the completion of the first edition of this book major developments have occurred in the pharmaceutical industry that have shaped the field of near infrared NIR spectroscopy A new initiative from the U S Food and Drug Administration FDA to modernize regulations of pharmaceutical manufacturing and drug quality has helped position NIR sp ***Science and Technology of Liquid Metal Coolants in Nuclear Engineering*** Thiagarajan Gnanasekaran, 2024-09-17 Science and Technology of Liquid Metal Coolants in Nuclear Engineering is a comprehensive consolidation of the latest research and knowledge on liquid metal coolants Over the last decades various new technologies have been developed for the liquid metal coolants of fast breeder and fusion reactors and accelerator driven systems Details of pumps and instrumentation used in these coolants and their operating principles are included to provide the reader with a well rounded understanding of the topic and to guide on the operation of different liquid metal coolant systems Methods for the safe handling and control of impurity levels in these coolants are clearly discussed along with alkali metal fires and their management including methods for safe disposal of sodium waste Discusses the thermophysical and chemical properties of liquid metals described with their microscopic origin Includes methods for the safe handling of liquid metal coolants and their purification and management Discusses pumps and instrumentation principles and design **Computational Inorganic and Bioinorganic Chemistry** Edward I. Solomon, Robert A. Scott, R. Bruce King, 2013-02-19 Over the past several decades there have been major advances in our ability to computationally evaluate the electronic structure of inorganic molecules particularly transition metal systems This advancement is due to the Moore s Law increase in computing power as well as the impact of density functional theory DFT and its implementation in commercial and freeware programs for quantum chemical calculations Improved pure and hybrid density functionals are allowing DFT calculations with accuracy comparable to high level Hartree Fock treatments and the results of these calculations can now be evaluated by experiment When calculations are correlated to and supported by experimental data they can provide fundamental insight into electronic structure and its contributions to physical properties and chemical reactivity This interplay continues to expand and contributes to both improved value of experimental results and improved accuracy of computational predictions The purpose of this EIC Book is to provide state of the art presentations of quantum mechanical and related methods and their applications written by many of the leaders in the field Part 1 of this volume focuses on methods their background and implementation and their use in describing bonding properties energies

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