# Semiclassical Mechanics with Molecular Applications

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# **Semiclassical Mechanics With Molecular Applications**

Stuart A. Rice

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systems and thus nonlinear mechanics is the proper theory to elucidate molecular dynamics by investigating invariant structures in phase space Intramolecular energy transfer and the breaking and forming of a chemical bond have now found a rigorous explanation by studying phase space structures Unimolecular Reaction Dynamics Tomas Baer, William L. Hase, 1996-06-27 This book provides a penetrating and comprehensive description of energy selected reactions from a theoretical as well as experimental view Three major aspects of unimolecular reactions involving the preparation of the reactants in selected energy states the rate of dissociation of the activated molecule and the partitioning of the excess energy among the final products are fully discussed with the aid of 175 illustrations and over 1 000 references most from the recent literature Examples of both neutral and ionic reactions are presented Many of the difficult topics are discussed at several levels of sophistication to allow access by novices as well as experts Among the topics covered for the first time in monograph form is a discussion of highly excited vibrational rotational states and intramolecular vibrational energy redistribution Problems associated with the application of RRKM theory are discussed with the aid of experimental examples Detailed comparisons are also made between different statistical models of unimolecular decomposition Both quantum and classical models not based on statistical assumptions are described Finally a chapter devoted to the theory of product energy distribution includes the application of phase space theory to the dissociation of small and large clusters The work will be welcomed as a valuable resource by practicing researchers and graduate students in physical chemistry and those involved in the study of chemical reaction dynamics Theory of Chemical Reaction Dynamics Antonio Laganà, 2005 Degenerate States in Chemistry, Volume 124 Michael Baer, Gert Due Billing, 2003-03-31 Edited by Nobel Prize winner Ilya Prigogine and renowned authority Stuart A Rice the Advances in Chemical Physics series provides a forum for critical authoritative evaluations in every area of the discipline In a format that encourages the expression of individual points of view experts in the field present comprehensive analyses of subjects of interest This stand alone special topics volume edited by Gert D Billing of the University of Copenhagen and Michael Baer of the Soreg Nuclear Research Center in Yavne Israel reports recent advances on the role of degenerate states in chemistry Volume 124 collects innovative papers on Complex States of Simple Molecular Systems Electron Nuclear Dynamics Conical Intersections and the Spin Orbit Interaction and many more related topics Advances in Chemical Physics remains the premier venue for presentations of new findings in its field Modern Trends in Chemical Reaction Dynamics Xueming Yang, Kopin Liu, 2004 The field of chemical reaction dynamics has made huge progress during the last decade or so The aim of these volumes is to provide graduate students and experts in the field with a picture of the current status of advanced experimental and theoretical research in chemical reaction dynamics Modern Methods for Multidimensional Dynamics Computations in Chemistry Donald Leo Thompson, 1998 This volume describes many of the key practical theoretical techniques that have been developed to treat chemical dynamics problems in many atom systems It contains thorough treatments of fundamental theory and prescriptions

for performing computations The selection of methods ranging from gas phase bimolecular reactions to complex processes in condensed phases reflects the breadth of the field The book is an excellent reference for proven and accepted methods as well as for theoretical approaches that are still being developed It is appropriate for graduate students and other novices who wish to begin working in chemical dynamics as well as active researchers who wish to acquire a wider knowledge of the field

Vibronic Interactions: Jahn-Teller Effect in Crystals and Molecules Michael D. Kaplan, George O. Zimmerman, 2012-12-06 This book is based mostly on the reports presented at the XVth International lahn Teller Symposium on Vibronic Interactions in Crystals and Molecules and NATO Advanced Research Workshop Colossal Magnetoresistance and Vibronic Interactions that took place at Boston on August 16 22 of the year 2000 This is the first time the Symposium took place in the USA where recently the giant splash of the attention to the 1 ahn Teller effect occurred This tremendous interest to the field all over the world is reflected not only in the numerous publications in many American and European 10urnals but of the leading scientists from additionally in the Symposium's participation the well known Universities National Laboratories and industrial companies which was the largest in the history of the Symposium The renaissance of the 1ahn Teller physics is closely related to the three fundamental discoveries in science The most significant among them is the discovery of high Tc superconductivity by K A Muller and G Bednorz for whom the 1ahn Teller idea was the motivation in their search The result of this search is well known a wide spectrum of the 1ahn Teller ion based materials with Tc between 24K and 135K were found The second discovery is the existence of a new polymorph of carbon the C60 The microscopic analysis of all physical chemical and biological properties of the buckyballs is based on 1ahn Teller type of interactions The third is colossal magnetoresistance Molecular Quantum Mechanics Peter W. Atkins, Ronald S. Friedman, 2011 This text unravels those fundamental physical principles which explain how all matter behaves It takes us from the foundations of quantum mechanics through quantum models of atomic molecular and electronic structure and on to discussions of spectroscopy and the Theory of Molecular Collisions Gabriel G Balint-Kurti, Alexander electronic and magnetic properties of molecules Palov, 2015-07-02 Almost 100 years have passed since Trautz and Lewis put forward their collision theory of molecular processes Today knowledge of molecular collisions forms a key part of predicting and understanding chemical reactions This book begins by setting out the classical and quantum theories of atom atom collisions Experimentally observable aspects of the scattering processes their relationship to reaction rate constants and the experimental methods used to determine them are described The quantum mechanical theory of reactive scattering is presented and related to experimental observables The role of lasers in the measurement and analysis of reactive molecular collisions is also discussed Written with postgraduates and newcomers to the field in mind mathematics is kept to a minimum and readers are guided to appendices and further reading to gain a deeper understanding of the mathematics involved Advances in Chemical Physics, Volume 138 Stuart A. Rice, 2008-04-30 This series provides the chemical physics field with a forum for critical authoritative

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