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Most identities in chemical thermodynaimcs arise from application of the first and second laws of thermodynamics, particularly the law of conservation of energy, to these state functions. It is the ...

Chemical thermodynamics
Under the single-step nucleation mechanism, the assembly structure is inherited from direct nucleation events in solution. Although the kinetically selected nuclei could give rise to structurally ...

Biomolecular Assemblies: Moving from Observation to Predictive Design
The author's generalized procedure for putting a theory into Hamiltonian form is applied to Einstein's theory of gravitation. It is shown that one can make a change in the action density, not ...

A completely updated, expanded edition of a longstanding and influential text on chemical thermodynamics Covers the logical foundations and interrelationships of thermodynamics and their application to problems that are commonly encountered by the chemist. Explanations of abstract concepts in a clear and simple, yet still rigorous fashion Logical arrangement of the material to facilitate learning, including worked out examples. Computational techniques, graphical, numerical, and analytical, are described fully and are used frequently, both in illustrative and in assigned problems.

Reflecting the growing volume of published work in this field, researchers will find this book an invaluable source of information on current methods and applications.

Annotation This textbook and reference outlines the principles and applications of thermodynamics in geochemistry.

A new, millennium edition of the classic treatment of chemical thermodynamics Widely recognized for half a century for its first-rate, logical introduction to phenomenological thermodynamics, this classic work is now thoroughly revised for the new millennium. The Sixth Edition continues to cover the fundamentals and methods of thermodynamics with exceptional vigor and clarity, while incorporating many new developments. Up-to-date examples are carefully gleaned from the literature for their practical interest to chemists, biochemists, geologists, chemical engineers, and materials scientists. Chemical Thermodynamics: Basic Theory and Methods, Sixth Edition provides readers with clear explanations of essential chemistry, mathematics, and the latest computational tools. Additional new features include: * Liberal reference to Web-based resources and databases * Extensive tables of thermodynamic data organized by source * High-quality exercises with a separate student manual available for solutions to alternate problems * Simple methods for the calculation of partial molar functions from experimental data * Expanded and revised chapters containing discussion of excess thermodynamic functions, a treatment of the Second Law and Equilibrium on the basis of the Planck function as well as the Gibbs function, and treatment of real gases in terms of the Redlich-Kwong equation

Specialist Periodical Reports provide systematic and detailed review coverage of progress in the major areas of chemical research. Written by experts in their specialist fields the series creates a unique service for the active research chemist, supplying regular critical in-depth accounts of progress in particular areas of chemistry. For over 80 years the Royal Society of Chemistry and its predecessor, the Chemical Society, have been publishing reports charting developments in chemistry, which originally took the form of Annual Reports. However, by 1967 the whole spectrum of chemistry could no longer be contained within one volume and the series Specialist Periodical Reports was born. The Annual Reports themselves still existed but were divided into two, and subsequently three, volumes covering Inorganic, Organic and Physical Chemistry. For more general coverage of the highlights in chemistry they remain a 'must'. Since that time the SPR series has altered according to the fluctuating degree of activity in various fields of chemistry. Some titles have remained unchanged, while others have altered their emphasis along with their titles; some have been combined under a new name whereas others have had to be discontinued.

Originally published in 1923, this classic was revised and updated in the early 60s, adding material on solution thermodynamics, results in statistical mechanics, surfaces, gravitational and electromagnetic fields, more. 1961 second edition.

Created for engineers and students working with pure polymers and polymer solutions, this handbook provides up-to-date, easy to use methods to obtain specific volumes and phase equilibrium data. A comprehensive database for the phase equilibria of a wide range of polymer-solvent systems, and PVT behavior of pure polymers are given, as are accurate predictive techniques using group contributions and readily available pure component data. Two computer programs on diskettes are included. POLYPROG implements procedures given for prediction and correlation for specific volume of pure polymer liquids and calculation of vapor-liquid equilibria (VLE) of polymer solutions. POLYDATA provides an easy method of accessing the data contained in the many databases in the book. Both disks require a computer with a math coprocessor. This handbook is a valuable resource in the design and operation of many polymer processes, such as polymerization, devolatilization, drying, extrusion, and heat exchange. Special Details: Hardcover with Disks. Special offer: Purchase this book along with X-131, Handbook of Diffusion and Thermal Properties of Polymers and Polymer Solutions and receive a 20 percent discount off the list or member price.

Solution Thermodynamics and its Application to Aqueous Solutions: A Differential Approach, Second Edition introduces a differential approach to solution thermodynamics, applying it to the study of aqueous solutions. This valuable approach reveals the molecular processes in solutions in greater depth than that gained by spectroscopic and other methods. The book clarifies what a hydrophobe, or a hydrophile, and in turn, an amphiphile, does to H2O. By applying the same methodology to ions that have been ranked by the Hofmeister series, the author shows that the kosmotropes are either hydrophobes or hydration centers, and that chaotropes are hydrophiles. This unique approach and important updates make the new edition a must-have reference for those active in solution chemistry. Unique differential approach to solution thermodynamics allows for experimental evaluation of the intermolecular interaction Incorporates research findings from over 40 articles published since the previous edition Numerical or graphical evaluation and direct experimental determination of third derivatives, enthalpic and volumetric AL-AL interactions and amphiphiles are new to this edition Features new chapters on spectroscopic study in aqueous solutions as well as environmentally friendly and hostile water aqueous solutions

The book is concerned with the application of physical techniques to the study of the structure and interactions of biopolymers. The treatment is confined to those procedures applicable to solutions. The material has been tested on students in actual classes, thereby permitting the elimination of ambiguities and potential points of difficulty. Stress has been placed upon lucidity of treatment, and difficult steps in derivations have been explained. The mathematical exposition has been made as clear and simple as feasible. Examples of actual data are given.

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