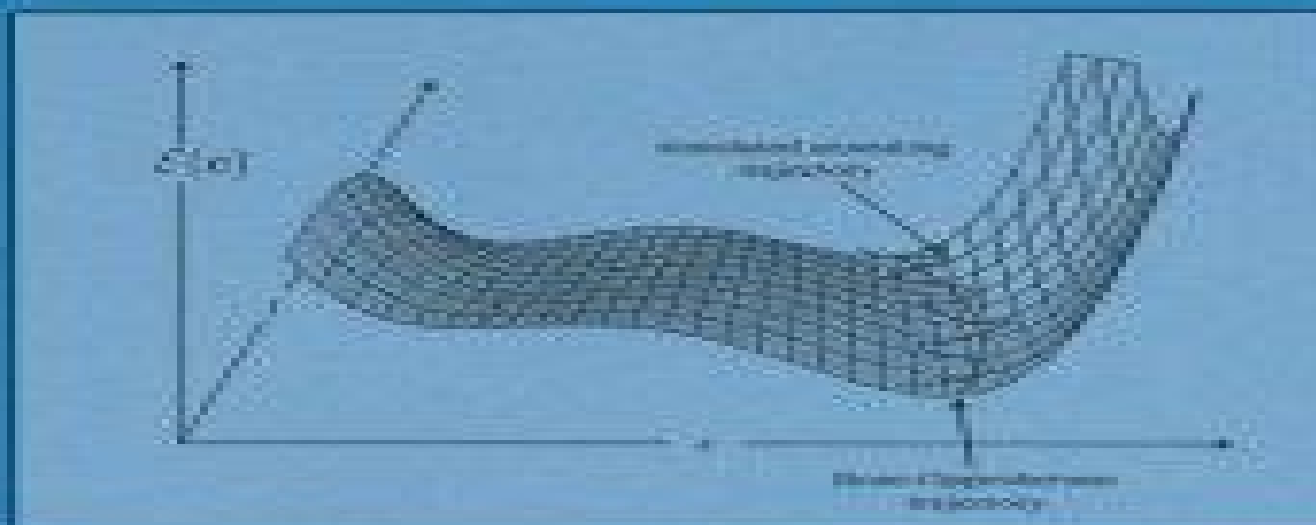


Adaption of simulated annealing to chemical optimization problems

JOHN H. KALVAS
California



Adaption Of Simulated Annealing To Chemical Optimization Problems

J.H. Kalivas

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Adaption of Simulated Annealing to Chemical Optimization Problems J.H. Kalivas, 1995-08-01 Optimization problems occurring regularly in chemistry vary from selecting the best wavelength design for optimal spectroscopic concentration predictions to geometry optimization of atomic clusters and protein folding Numerous optimization tactics have been explored to solve these problems While most optimizers maintain the ability to locate global optima for simple problems few are robust against local optima convergence with regard to difficult or large scale optimization problems Simulated annealing SA has shown a great tolerance to local optima convergence and is often called a global optimizer The optimization algorithm has found wide use in numerous areas such as engineering computer science communication image recognition operation research physics and biology Recently SA and variations thereof have shown considerable success in solving numerous chemical optimization problems The main thrust of this book is to demonstrate the use of SA in a wide range of chemical problems The potentiality of SA GSA and other modifications of SA to serve specific needs in a variety of chemical disciplines are covered A detailed discussion on SA and GSA is given in Chapter 1 presenting the theoretical framework from which a computer program can be written by the reader The remainder of the book describes applications of SA type algorithms to a diverse set of chemical problems The final chapter contains an algorithm for GSA written in the MatLab programming environment This program can be easily adapted to any optimization problem and with only slight modifications can be altered to perform SA A general flowchart is also given

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Design and Optimization in Organic Synthesis Rolf

Carlson,Johan E. Carlson,2005-04-08 Revised and updated Design and Optimization in Organic Synthesis presents strategies to explore experimental conditions and methodologies for systematic studies of entire reaction systems substrates reagent s catalyst s and solvents Chemical phenomena are not usually the result of a single factor and this book describes how statistically designed methods can be used to analyse and evaluate synthetic procedures The methodology is based on multivariate statistical techniques The accompanying CD contains data tables and programmes This book is essential reading for anyone working in process design and development in fine chemicals or the pharmaceutical industry and is suitable for those with no experience in the field Contains recalculated models and redrawn figures as well as new chapters on for example the design of combinatorial libraries Presents strategies to explore experimental conditions and methodologies Enables the analysis and prediction of the best synthetic procedures

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