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Nota di contenuto	Computer-Aided Modeling of Reactive Systems; Contents; Chapter 1. Overview; REFERENCES and FURTHER READING; Chapter 2. Chemical Reaction Models; 2.1 STOICHIOMETRY OF REACTION SCHEMES; 2.2 COMPUTABILITY OF REACTION RATES FROM DATA; 2.3 EQUILIBRIA OF CHEMICAL REACTIONS; 2.4 KINETICS OF ELEMENTARY STEPS; 2.5 PROPERTIES OF REACTION NETWORKS; 2.6 EVIDENCE FOR REACTION STEPS; PROBLEMS; REFERENCES and FURTHER READING; Chapter 3. Chemical Reactor Models; 3.1 MACROSCOPIC CONSERVATION EQUATIONS; 3.1.1 Material Balances; 3.1.2 Total Energy Balance; 3.1.3 Momentum Balance; 3.1.4 Mechanical Energy Balance 3.2 HEAT AND MASS TRANSFER IN FIXED BEDS3.3 INTERFACIAL STATES IN FIXED-BED REACTORS; 3.4 MATERIAL TRANSPORT IN POROUS CATALYSTS; 3.4.1 Material Transport in a Cylindrical Pore Segment; 3.4.2 Material Transport in a Pore Network; 3.4.3 Working Models of Flow and Diffusion in Isotropic Media; 3.4.4 Discussion; 3.4.5 Transport and Reaction in Porous Catalysts; 3.5 GAS PROPERTIES AT LOW PRESSURES; 3.6 NOTATION; REFERENCES and FURTHER READING; Chapter 4. Introduction to Probability and Statistics; 4.1 STRATEGY OF DATA-BASED INVESTIGATION; 4.2 BASIC CONCEPTS IN PROBABILITY

## THEORY

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### Sommario/riassunto

Learn to apply modeling and parameter estimation tools and strategies to chemical processes using your personal computer This book introduces readers to powerful parameter estimation and computational methods for modeling complex chemical reactions and reaction processes. It presents useful mathematical models, numerical methods for solving them, and statistical methods for testing and discriminating candidate models with experimental data. Topics covered include: Chemical reaction models Chemical reactor models Probability and statistics Bayesian estimation Process modeling with si

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