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Nota di contenuto	Frontmatter Preface Contents List of Algorithms Chapter 1. Introduction to computer simulation Chapter 2. Scientific Computing in C Chapter 3. Fundamentals of statistical physics Chapter 4. Inter- and intramolecular potentials Chapter 5. Molecular Dynamics simulations Chapter 6. Monte Carlo simulations Chapter 7. Advanced topics, and applications in soft matter Appendix A. The software development life cycle Appendix B. Installation guide to Cygwin Appendix C. Introduction to the UNIX/Linux programming environment Appendix D. Sample program listings Appendix E. Reserved keywords in C Appendix F. Functions of the standard library <string.h> Appendix G. Elementary combinatorial problems  Appendix H. Some useful constants Appendix I. Installing the GNU Scientific Library, GSL Appendix J. Standard header files of the ANSI-C library Appendix K. The central limit theorem Bibliography Glossary of Acronyms Index Authors</string.h>
Sommario/riassunto	This work is a needed reference for widely used techniques and methods of computer simulation in physics and other disciplines, such as materials science. Molecular dynamics computes a molecule's reactions and dynamics based on physical models; Monte Carlo uses

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random numbers to image a system's behaviour when there are different possible outcomes with related probabilities. The work conveys both the theoretical foundations as well as applications and "tricks of the trade", that often are scattered across various papers. Thus it will meet a need and fill a gap for every scientist who needs computer simulations for his/her task at hand. In addition to being a reference, case studies and exercises for use as course reading are included.