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Nota di contenuto	Calculation of NMR and EPR Parameters; Contents; Foreword; List of Contributors; Part A Introductory Chapters; 1 Introduction: The Quantum Chemical Calculation of NMR and EPR Parameters; 2 Theory of NMR parameters. From Ramsey to Relativity, 1953 to 1983; 2.1 Introduction; 2.2 Spin-Spin Coupling; 2.3 Chemical Shifts; 2.4 General Aspects; 2.5 From 1983 to 2003; 3 Historical Aspects of EPR Parameter Calculations; 4 The Effective Spin Hamiltonian Concept from a Quantum Chemical Perspective; 5 Fundamentals of Nonrelativistic and Relativistic Theory of NMR and EPR Parameters; 5.1 Introduction 5.2 Classical Theory of the Interaction of a Charged Particle with an Electromagnetic Field 5.3 Quantum Mechanical Hamiltonians in a Time-Independent Electromagnetic Field; 5.4 Perturbation Theory of Magnetic

Effects; 5.5 Non-Relativistic Theory of EPR and NMR Parameters; 5.6 Relativistic Theory of Magnetic Properties; 5.7 The Leading Relativistic Corrections; 5.8 Concluding Remarks; Part B NMR Parameters, Methodological Aspects; 6 Chemical Shifts with Hartree-Fock and Density Functional Methods; 6.1 Introduction; 6.2 Linear Response and the Gauge Origin Problem  
6.3 Determination of the First-Order Orbitals  
6.4 Distributed Gauge Origins, IGLO and GIAO Approaches; 6.5 Distributed Gauge Origins in Real Space, a "Continuous Set of Gauge Transformations"; 6.6 Beyond Pure Density Functional Theory; 6.7 Conclusions; 7 Spin-Spin Coupling Constants with HF and DFT Methods; 7.1 Introduction; 7.2 The Calculation of Indirect Nuclear Spin-Spin Coupling Constants; 7.3 Examples of Applications; 7.4 Conclusions; 8 Electron-Correlated Methods for the Calculation of NMR Chemical Shifts; 8.1 Introduction; 8.2 Theoretical Background  
8.3 Electron-Correlated Treatment of NMR Chemical Shifts  
8.4 Special developments; 8.5 Numerical Results; 8.6 Summary and Outlook; 9 Semiempirical Methods for the Calculation of NMR Chemical Shifts; 9.1 Introduction; 9.2 Methods; 9.3 Representative Applications; 9.4 Concluding Remarks: Limitations of Semiempirical Methods for the Calculation of NMR Parameters; 10 Ro-Vibrational Corrections to NMR Parameters; 10.1 Introduction; 10.2 Perturbation Theory; 10.3 Other Approaches for Calculating Vibrationally Averaged NMR Properties; 10.4 Examples of Vibrational Contributions to NMR Properties  
10.5 Summary  
11 Molecular Dynamics and NMR Parameter Calculations; 11.1 Introduction; 11.2 Methods; 11.3 Examples; 11.4 Summary and Conclusions; 12 Use of Continuum Solvent Models in Magnetic Resonance Parameter Calculations; 12.1 Introduction; 12.2 General Features of Continuum Models; 12.3 Applications of Continuum Models to the Prediction of NMR Parameters; 12.4 Applications of Continuum Models to the Prediction of EPR Parameters; 12.5 Conclusions; 13 Perturbational and ECP Calculation of Relativistic Effects in NMR Shielding and Spin-Spin Coupling; 13.1 Introduction  
13.2 Nuclear Shielding and Spin-Spin Coupling

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## Sommario/riassunto

This is the first book to present the necessary quantum chemical methods for both resonance types in one handy volume, emphasizing the crucial interrelation between NMR and EPR parameters from a computational and theoretical point of view. Here, readers are given a broad overview of all the pertinent topics, such as basic theory, methodic considerations, benchmark results and applications for both spectroscopy methods in such fields as biochemistry, bioinorganic chemistry as well as with different substance classes, including fullerenes, zeolites and transition metal compounds. The chapters

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