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this volume covers the formation of carbon carbon single double and triple bonds by substitution and addition reactions as well as by various rearrangements the formation of carbon carbon multiple bonds by elimination and condensation procedures is fully documented in addition the synthesis of carbon hydrogen bonds principally by substitution and addition reactions is featured as is the preparation of a wide variety of carbon centred anions cations and radicals

most chemists today have either taken part in or been affected by the chemical revolution that has taken place over the course of the last century developments in instrumentation have changed not just what chemists do but also how they think about chemistry new and exciting areas of previously inaccessible research have been opened up as a direct result of this revolution this is the first book to examine this instrumental revolution and goes on to assess the impact on chemical practice in areas ranging from organic chemistry and biochemistry to environmental analysis and process control thus demonstrating how fundamental and extensive are the changes that have occurred with contributions from internationally recognised specialists this lavishly illustrated book provides a focal point for any historian of chemistry or chemist with an interest in this fascinating topic this book is published in association with the science museum london uk and the chemical heritage foundation philadelphia

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computational chemistry is increasingly used in most areas of molecular science including organic inorganic medicinal biological physical and analytical chemistry researchers in these fields who do molecular modelling need to understand and stay current with recent developments this volume

like those prior to it features chapters by experts in various fields of computational chemistry two chapters focus on molecular docking one of which relates to drug discovery and cheminformatics and the other to proteomics in addition this volume contains tutorials on spin orbit coupling and cellular automata modeling as well as an extensive bibliography of computational chemistry books from reviews of the series reviews in computational chemistry remains the most valuable reference to methods and techniques in computational chemistry journal of molecular graphics and modelling one cannot generally do better than to try to find an appropriate article in the highly successful reviews in computational chemistry the basic philosophy of the editors seems to be to help the authors produce chapters that are complete accurate clear and accessible to experimentalists in particular and other nonspecialists in general journal of the american chemical society

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