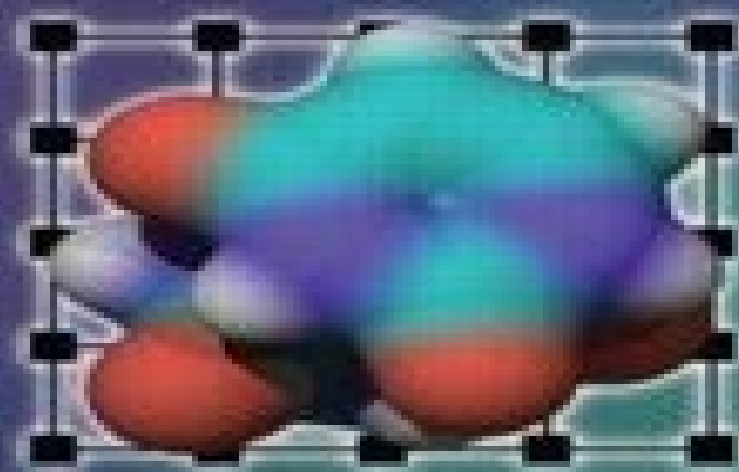


Parallel Computing *in* Quantum Chemistry



Curtis L. Janssen

Ida M. B. Nielsen



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Parallel Computing in Computational Chemistry Timothy G. Mattson, American Chemical Society. Meeting, 1995-07-27 Developed from a symposium sponsored by the Division of Computers in Chemistry at the 207th National Meeting of the American Chemical Society San Diego California March 13 17 1994
Parallel Computing in Quantum Chemistry -- Message Passing and Beyond for a General Ab Initio Program System, 1994 One of the most prominent aims in Computational Chemistry is the modeling of chemical reactions and the prediction of molecular properties Quantum chemical methods are used for the calculation of molecular structures spectra reaction energy profiles and many other interesting quantities Nowadays the accuracy of the theoretical calculations can compete to an increasing extent with the experimental one A great variety of quantum chemical methods exist ranging from the standard Hartree Fock theory to sophisticated electron correlation approaches From a computational point of view all these methods require rather lengthy and complicated program codes and have to handle a large amount of data to be stored on external devices In the simplest case the Hartree Fock SCF method direct algorithms have eliminated the I O and storage bottleneck and have opened the way to parallel implementations For post Hartree Fock methods the situation is much more complicated as will be demonstrated below Therefore most of the previous attempts in parallelizing quantum chemical ab initio programs concentrated on SCF methods The authors investigations presented here are a continuation of their previous work on the parallelization of the COLUMBUS program system The COLUMBUS program is based on the multireference single and double excitation configuration interaction MRSDCI approach It is very portable and runs on a large variety of computers including numerous Unix based workstations VAX VMS minicomputers IBM mainframes and Cray supercomputers

Parallel Computing in Computational Chemistry Timothy G. Mattson, 1995 [Methods in Computational Chemistry](#) Stephen Wilson, 2013-12-01 Recent years have seen the proliferation of new computer designs that employ parallel processing in one form or another in order to achieve maximum performance Although the idea of improving the performance of computing machines by carrying out parts of the computation concurrently is not new indeed the concept was known to Babbage such machines have until fairly recently been confined to a few specialist research laboratories

Nowadays parallel computers are commercially available and they are finding a wide range of applications in chemical calculations. The purpose of this volume is to review the impact that the advent of concurrent computation is already having and is likely to have in the future on chemical calculations. Although the potential of concurrent computation is still far from its full realization it is already clear that it may turn out to be second in importance only to the introduction of the electronic digital computer itself. *Parallel Computing in Computational Chemistry*, 1995

Reviews in Computational Chemistry, Volume 6 Kenny B. Lipkowitz, Donald B. Boyd, 2009-09-22. Volume 6 of the successful series *Reviews in Computational Chemistry* contains articles of interest to pharmaceutical chemists, biological chemists, chemical engineers, inorganic and organometallic chemists, synthetic organic chemists, polymer chemists, and theoretical chemists. The series is designed to help the chemistry community keep current with the many new developments in computational techniques. The writing style is refreshingly pedagogical and non-mathematical, allowing students and researchers access to computational methods outside their immediate area of expertise. **Computational Chemistry**, 2000

Scientific Computing on Supercomputers II J.T. Devreese, 2012-12-06. The International Workshop on The Use of Supercomputers in Theoretical Science took place on November 29 and 30 1989 at the University of Antwerp, UIA, Antwerpen, Belgium. It was the fifth in a series of workshops, the first of which took place in 1984. The principal aim of these workshops is to present the state of the art in scientific large scale and high speed computation. Computational science has developed into a third methodology, equally important now as its theoretical and experimental companions. Gradually, academic researchers acquired access to a variety of supercomputers and as a consequence, computational science has become a major tool for their work. It is a pleasure to thank the Belgian National Science Foundation (NFWO-FNRS) and the Ministry of Scientific Affairs for sponsoring the workshop. It was organized both in the framework of the Third Cycle Vectorization, Parallel Processing, and Supercomputers and the Governmental Program in Information Technology. We also very much would like to thank the University of Antwerp, Universitaire Instelling Antwerpen, UIA, for financial and material support. Special thanks are due to Mrs H. Evans for the typing and editing of the manuscripts and for the preparation of the author and subject index.

Theory and Applications of Computational Chemistry - 2008 Dong-Qing Wei, Xi-Jun Wang, 2009-04-03. The TACC conference deals with the theoretical methods and application of computational chemistry. The theoretical aspects include development of computational and statistical techniques, quantum and statistical mechanical methods that could be used to solve complicated scientific problems which will be applied to various fields of chemistry, physics, and biology. Since the first applications of computer in chemistry in the early 1960s, computational chemistry has evolved today to be at the vanguard of most advances in chemistry, material science, physics, geology, biology, and medicine. Its impact on chemistry can be noted from the fact that within four decades of its existence, five of the top ten most highly cited publications in 125 years of JACS are related to computational chemistry. Apart from technical advances in computers and algorithms, the phenomenal impact

of computational chemistry is also due to the dedicated and sustained efforts of a large number of individuals who participated in this most important conference in this field **Computational Chemistry F.E.C.S. Conference** Francesco Bernardi, Jean-Louis Rivail, 1998-04-15 Parallel Computing 85 Manfred Feilmeier, Gerhard Robert Joubert, Udo Schendel, 1986 Proceedings Parallel Computing Practical Applications of Parallel Computing Laurence Tianruo Yang, 2003 The continuous progress in scientific research is one of the important factors explaining the constantly increasing demand for computational power On the other hand one of the results of such progress is the availability of more powerful computer platforms To that end this volume reviews a broad array of subjects based on the solutions to the daily problems in industrial production research and development *Edmonton Symposium on Computational Quantum Chemistry and Parallel Processing* Edmonton Symposium on Computational Quantum Chemistry and Parallel Processing, 1987 *Vector and Parallel Processors in Computational Science* Iain S. Duff, John Ker Reid, 1985 *Chemical Modelling* Alan Hinchliffe, 2008 Chemical Modelling Applications and Theory comprises critical literature reviews of molecular modelling both theoretical and applied Molecular modelling in this context refers to modelling the structure properties and reactions of atoms molecules materials Each chapter is compiled by experts in their fields and provides a selective review of recent literature With chemical modelling covering such a wide range of subjects this Specialist Periodical Report serves as the first port of call to any chemist biochemist materials scientist or molecular physicist needing to acquaint themselves of major developments in the area Volume 5 covers literature published from June 2005 to May 2007 **Highly Parallel Computing** George S. Almasi, Allan Gottlieb, 1994 This second edition includes new exercises for each chapter a quantitative treatment of speedup seismic migration using a workstation network as a parallel computer recent changes in technology more languages fat trees wormhole switching new SIMD hardware an expanded section on CM 2 new MIMD hardware using workstation clusters as a MIMD system and directory based caches Annotation copyright by Book News Inc Portland OR **Modern Techniques in Computational Chemistry: MOTECCTM-90** E. Clementi, 1990-08-31 SIAM Journal on Scientific Computing , 2006

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