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Crystallographic and Modeling Methods in Molecular Design Symposium.(1989 : Gulf Shores, Ala.) Title(s): Crystallographic and modeling methods in molecular design. Charles E. Bugg, Steven E. Ealick, editors. Country of Publication: United States Publisher: New York : Springer-Verlag, c1990. Description: xvii, 269 p. : ill. Language: English ISBN: 9780387972107 (alk. paper), 0387972102 (alk

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Abstract. The investigation of low-energy conformations of molecules or substructures is fundamental to the process of molecular modelling. A number of computational procedures exist for this purpose, but suffer from some operational or scientific limitations.

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