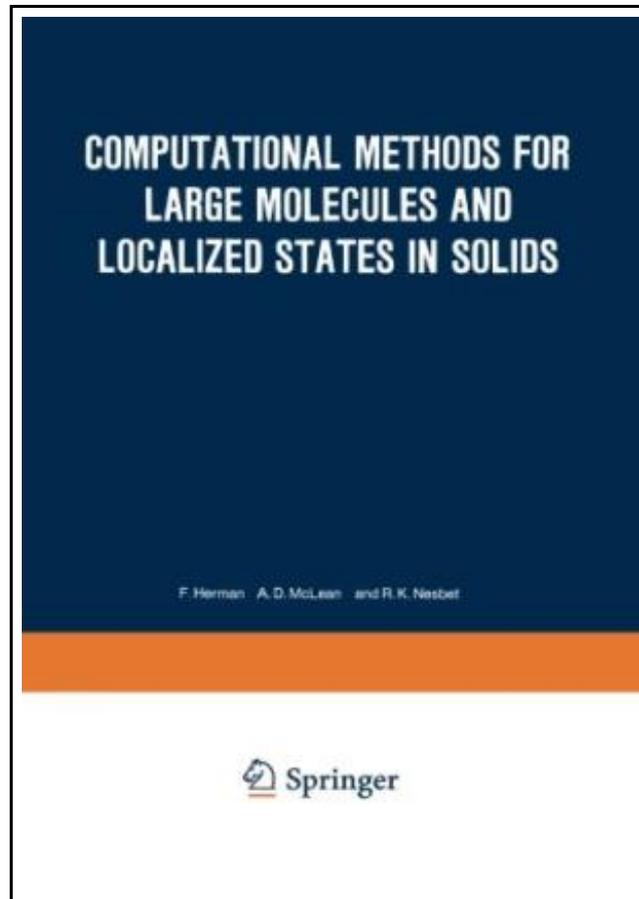


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COMPUTATIONAL METHODS FOR LARGE MOLECULES AND LOCALIZED STATES IN SOLIDS



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