

Errata

Calculation of Optimum Geometries and Force Fields by the CNDO/Force Method

Meenakshisundaram Kanakavel, Jayaraman Chandrasekhar, Sankaran
Subramanian and Surjit Singh

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Ref. [30] should read: Schachtschneider, J. H., in: Vibrational spectra of polyatomic molecules VI,
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