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*Simulation of FT-IR and FT-Raman Spectra Based on Scaled DFT Calculations,  
Vibrational Assignments, Hyperpolarizability, NMR Chemical Shifts and Homo-Lumo Analysis of 1-  
Chloro-4-Nitrobenzene*

*by*

*S. Seshadri*

*after review is found suitable and has been published in*

*Volume 4, Issue VIII, August 2016*

*in*

*By*

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