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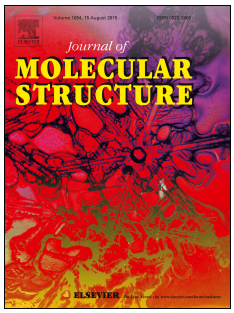

Vibrational spectroscopic analysis, molecular dynamics simulations and molecular docking study of 5-nitro-2-phenoxymethyl benzimidazole

Vidya V. Menon, Egemen Foto, Y. Sheena Mary, Esin Haratas, C. Yohannan Panicker, Gözde Yalcin, Stevan Armaković, Sanja J. Armaković, C. Van Alsenoy, Ilkay Yildiz

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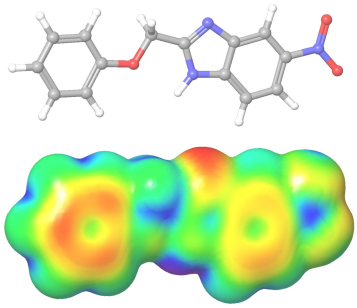

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