

Analysis of infrared spectra of neat liquid N-methylpyrrole.

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Streszczenie

The vibrational properties of neat liquid N-methylpyrrole were investigated. Quantitative near-infrared and mid-infrared spectra recordings including thin film MIR transmission studies were performed. Based on these data the optical constants of N-methylpyrrole in the 12500-560 cm⁻¹ region were determined. The experimental results were supported by an anharmonic analysis based on Möller-Plesset perturbation theory and density functional theory level of theory, including the hybrid B2PLYP functional. Assignments based on potential energy distribution calculations and a comparison of experimental and simulated spectra were presented. The inaccuracy of static calculations in the reproduction of experimental spectra in the CH stretching region has been observed and explained by the rotation of methyl group. A vibrational spectrum was therefore recalculated by the use of CarParinello molecular dynamics method for ab initio simulation of experimental spectrum.

Adres publiczny

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<http://www.ifpan.edu.pl>