THE FAR-INFRARED SPECTRA OF CYCLOPROPYLAMINE

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The infrared spectra of cyclopropylamine (c- $C_3H_5NH_2$) in the region of 35-600 cm⁻¹ have been measured at 298K with a resolution of 0.00096 cm⁻¹ using the far-infrared beamline at the Canadian Light Source synchrotron. We report here the results of the rovibrational analysis of the ν_{27} (253.87 cm⁻¹) –NH₂ torsional fundamental, as well as the pure rotational analysis of transitions associated with the ground state and the first excited state of the –NH₂ torsional mode between 35 and 60 cm⁻¹. The ongoing assignment and analysis of hot bands and overtones involving higher torsional states will also be discussed.