Ammonium Ions Dynamics in NH_4Br – Proton NMR T_1 Measurements and Simulations

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Monte Carlo simulation of correlation functions for ammonium bromide have been carried out and the NMR relaxation times T_1 resulting from these functions were calculated. Results of simulations are compared to experimental values of T_1 measured in the temperature range 92 K to 292 K.

Key words: NMR; NH₄Br; Relaxation; Monte Carlo; Simulation.