

# P4987: BOUND STATE CALCULATIONS OF THE VAN DER WAALS $\text{NH}_3\text{--Ne}$ COMPLEX AND FIRST MICROWAVE DETECTION OF THE MISSING (*para*)- $\text{NH}_3\text{--Ne}$ NUCLEAR SPIN ISOMER

L. Surin<sup>1</sup>, I. Tarabukin<sup>1</sup>, C. Pérez<sup>2</sup>, M. Schnell<sup>2</sup>, J. Loreau<sup>3</sup>, and A. van der Avoird<sup>4</sup>

<sup>1</sup> Institute of Spectroscopy of RAS, <sup>2</sup> FS-SMP, Deutsches Elektronen-Synchrotron (DESY), <sup>3</sup> Chemistry, KU Leuven, <sup>4</sup> Institute for Molecules and Materials (IMM), Radboud University Nijmegen

The bound rovibrational energy levels were calculated for total angular momentum  $J = 0 - 10$  for both (*ortho* and *para*) nuclear spin modifications of the complex.

Pure rotational  $J = 1 - 0$  transition of the missing *para* $\text{NH}_3\text{--Ne}$  ( $^{20}\text{Ne}$  /  $^{22}\text{Ne}$ ) has been measured using CP-FTMW spectroscopy.

Nuclear quadrupole structure (from  $^{14}\text{N}$  nucleus) was resolved and analyzed, providing information about angular orientation of  $\text{NH}_3$  within complex.

