THE OPEN-SHELL Xe-O₂ VAN DER WAALS COMPLEX: MICROWAVE SPECTRA AND AB INITIO POTENTIAL ENERGY SURFACE

QING WEN AND WOLFGANG JÄGER, Department of Chemistry, University of Alberta, Edmonton, AB T6G 2G2, Canada.

An *ab initio* potential energy surface of the open shell Xe-O₂ complex was calculated at the RCCSD(T) level of theory. The global minimum was found at a T-shape structure with a well depth of 137.0 cm⁻¹. Five rotational transitions of Xe-O₂ were measured using a pulsed-nozzle Fourier transform microwave spectrometer and assigned to the ${}^{3}\Sigma^{-}$ electronic ground state of the complex. Each transition shows magnetic hyperfine structure arising from interactions of the electronic magnetic moment with the earth's magnetic field and these structures were reduced by the use of helmholtz coils. Hyperfine structures due to Fermi contact coupling between the electron spin of O₂ (*S* = 1) and the ¹²⁹Xe (¹³¹Xe) nuclear spin [*I* = 1/2 (3/2)] were also detected. The observed spectra were used to extract information about the geometric and electronic structure of the complex. The results provide detailed insight into the intermolecular interaction between Xe and paramagnetic oxygen. This may help to understand the relaxation mechanism of hyperpolarized ¹²⁹Xe in human blood in *in vivo* magnetic resonance imaging experiments.