d by Illinois Digital Environment for Ac

TWO-PHOTON IONIZATION STUDY OF THE LOW LYING STATES OF UN+

<u>ROBERT A. VANGUNDY</u>, THOMAS D PERSINGER, MICHAEL HEAVEN, *Department of Chemistry*, *Emory University, Atlanta, GA, USA*.

The electronic structures of UN and UN^+ are of interest for the testing and development of relativistic quantum chemistry methods. The ground state UN was probed by Matthew and Morse¹, who found that the electronic configuration $(5f^27s)$ differed from that of the isoelectronic UO⁺ cation $(5f^3)$. In the present study we examine the ionization energy of UN and the low energy states of UN⁺ by means of pulsed-field ionization zero kinetic energy photoelectron spectroscopy (PFI-ZEKE). Resonantly enhanced two photon ionization (R2PI) coupled with a time of flight mass spectrometer was used to confirm production of the UN molecule and locate suitable electronically excited states for subsequent access to UN⁺ via high-n Rydberg states. The results will be compared to the predictions from ligand field theory and high-level ab ignition calculations.

1. D. J. Matthew and M. D. Morse, J. Chem. Phys. 138, 184303 (2013)