

## PROTON TRANSFER MECHANISMS OF o-NITROPHENOL OBSERVED BY MeV ULTRAFAST ELECTRON DIFFRACTION

JOAO P.F. NUNEZ, LAUREN F HEALD, *Department of Physics and Astronomy, University of Nebraska - Lincoln, Lincoln, NE, USA*; MONIKA WILLIAMS, *Department of Chemistry, Stanford University, Stanford, CA, USA*; JIE YANG, *ARD FEL and Beam Physics, SLAC National Accelerator Laboratory, Menlo Park, CA, USA*; THOMAS JA WOLF, *Stanford PULSE Institute, SLAC National Accelerator Laboratory, Menlo Park, CA, USA*; CONOR RANKINE, *Department of Chemistry, University of York, York, United Kingdom*; BRYAN MOORE, *Department of Physics and Astronomy, University of Nebraska - Lincoln, Lincoln, NE, USA*; XIAOZHE SHEN, MING-FU LIN, *Linac Coherent Light Source, SLAC National Accelerator Laboratory, Menlo Park, CA, USA*; TODD MARTINEZ, *Stanford PULSE Institute, SLAC National Accelerator Laboratory, Menlo Park, CA, USA*; XIJIE WANG, *Acceleratory Directory, SLAC National Accelerator Laboratory, Menlo Park, CA, USA*; MARTIN CENTURION, *Department of Physics and Astronomy, University of Nebraska - Lincoln, Lincoln, NE, USA*.

Excited state intramolecular proton transfer (ESIPT) is key to many important biological mechanisms. However, direct observation of the structural dynamics of ESIPT has been limited given both the required spatial and temporal resolution. The combination of femtosecond temporal resolution and sub-Angstrom spatial resolution possible from mega-electronvolt ultrafast electron diffraction (MeV-UED) make it an ideal method for observing ESIPT. Furthermore, the neighboring -OH and -NO<sub>2</sub> groups on o-nitrophenol are known to undergo proton transfer upon excitation to the lowest singlet state (S<sub>1</sub>). Using MeV-UED, the structural dynamics of proton transfer in o-nitrophenol have been resolved following excitation to the S<sub>1</sub> state. In contrast to the S<sub>1</sub> state, higher lying excited states are suspected to follow different relaxation pathways and their structural evolution could provide further insight into the dynamics of ESIPT in o-nitrophenol. This presentation will discuss previous findings of ESIPT following excitation of o-nitrophenol to the S<sub>1</sub> state and will present new findings related to the relaxation dynamics of the S<sub>4</sub> state.