

CHARACTERISATION OF THE GROUND AND LOW-LYING EXCITED STATES OF MgO^+ BY PFI-ZEKE PHOTOELECTRON SPECTROSCOPY

JOEL R. SCHMITZ, CARLA KREIS, FRÉDÉRIC MERKT, *Department of Chemistry and Applied Biosciences, ETH Zürich, Zürich, Switzerland*

We report on the characterisation of the rovibrational structure of the ground and first excited electronic states of MgO^+ by high-resolution pulsed-field ionization zero-kinetic-energy (PFI-ZEKE) photoelectron spectroscopy. Rotationally cold ($T_{\text{rot}} = 5\text{K}$) MgO molecules in the $X^1\Sigma^+$ ($v = 0 - 2$) levels were generated in a supersonic expansion of a 0.1% $\text{N}_2\text{O}:\text{He}$ gas mixture following laser ablation off a magnesium (Mg) rod [1]. The rovibrational ionization thresholds corresponding to both spin-orbit components ($\Omega = \frac{1}{2}, \frac{3}{2}$) of the $X^+ 2\Pi_\Omega$ ($v^+ = 0 - 10$) states as well as to the first excited $A^+ 2\Sigma_{\frac{1}{2}}^+$ ($v^+ = 0 - 10$) state were reached in a resonant $1+1'$ two-photon excitation sequence via the rovibrational levels of the $F^1\Pi$, $E^1\Sigma^+$, $G^1\Pi$ and $3^3\Pi_2$ intermediate levels of MgO studied previously by Breckenridge and coworkers [2,3]. Our new results include accurate values for the adiabatic ionization energy of MgO and for the dissociation energies of the $\text{MgO } X^1\Sigma^+$ and $\text{MgO}^+ X^+ 2\Pi_\Omega$ and $A^+ 2\Sigma_{\frac{1}{2}}^+$ states. This work is carried out in the context of our studies of the rovibrational structure of doubly charged dications by high-resolution PFI-ZEKE spectroscopy of singly-charged cations following the approach recently taken to characterise the ground state of the thermodynamically stable dication MgAr^{2+} [4]. The talk will present a roadmap towards characterising the ground state of MgO^{2+} by resonant multiphoton excitation via electronically excited states of MgO^+ . The experiments will reveal whether MgO^{2+} is thermodynamically stable as predicted in Ref. [5] or metastable as predicted in Ref. [6].

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