

ZEKE SPECTROSCOPY OF METAL COMPLEXES WITH A MULTIDENTATE LIGAND:  $M(\text{H}_2\text{NCH}_2\text{CH}_2\text{NH}_2)$  ( $M = \text{Al, Ga, In}$ )

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Metal coordination to an ethylenediamine ligand may form a chain-like or a cyclic structure. The complex is expected to be a chain if the metal atom is bound to a single nitrogen atom or a five-member ring if the metal is attached to both nitrogen atoms. ZEKE (zero-electron-kinetic-energy) spectra show that  $M(\text{H}_2\text{NCH}_2\text{CH}_2\text{NH}_2)$  and  $M^+(\text{H}_2\text{NCH}_2\text{CH}_2\text{NH}_2)$  ( $M = \text{Al, Ga, In}$ ) complexes have a cyclic conformation. These cyclic complexes have significantly lower adiabatic ionization potentials (AIPs) and metal-nitrogen stretching frequencies than simple metal-primary amines. For example, the AIP of  $\text{Ga}(\text{H}_2\text{NCH}_2\text{CH}_2\text{NH}_2)$  was determined to be  $33322 \text{ cm}^{-1}$ ,  $6008 \text{ cm}^{-1}$  less than that of  $\text{Ga}(\text{H}_2\text{NCH}_3)$ . The  $\text{Ga}^+\text{-N}$  stretching frequency was measured to be  $204 \text{ cm}^{-1}$  in gallium-ethylenediamine and  $299 \text{ cm}^{-1}$  in gallium-methylamine.<sup>a</sup>

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<sup>a</sup>S. Li, G. K. Rothschof, D. Pillai, B. R. Sohnlein, B. M. Wilson, D. -S. Yang, *J. Chem. Phys.* **115**, 7968 (2001)