

SLOW ELECTRON VELOCITY-MAP IMAGING OF $\text{La}_2(\text{C}_6\text{H}_6)$ AND $\text{La}(\text{C}_6\text{H}_6)_2$

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Slow electron velocity-map imaging (SEVI) was used to study the structures and electronic states of $\text{La}_2(\text{C}_6\text{H}_6)$ and $\text{La}(\text{C}_6\text{H}_6)_2$ complexes formed in a metal cluster beam source. Electron spectra obtained from SEVI have the energy resolution of $\Delta E/eKE \approx 2\%$ at $eKE=400\text{cm}^{-1}$. The SEVI technique offers much higher data collection efficiency than pulsed-field ionization zero electron kinetic energy spectroscopy. From the SEVI spectra, adiabatic ionization energies were measured to be $32141(5)\text{ cm}^{-1}$ for $\text{La}_2(\text{C}_6\text{H}_6)$ and $39033(5)\text{ cm}^{-1}$ for $\text{La}(\text{C}_6\text{H}_6)_2$. The most active vibrational transition for both complexes was identified to be a metal-ligand stretching mode with a frequency of 180 cm^{-1} in the ion state. In addition, a benzene ring out-of-plane deformation mode was measured to be 303 cm^{-1} for $[\text{La}_2(\text{C}_6\text{H}_6)]^+$ and 408 cm^{-1} for $[\text{La}(\text{C}_6\text{H}_6)_2]^+$. By combining the spectra with theoretical calculations, we identified the ${}^2A_g \leftarrow {}^1A_g$ transition of $\text{La}_2(\text{C}_6\text{H}_6)$ (C_{2h}) and ${}^3A \leftarrow {}^2A$ of $\text{La}(\text{C}_6\text{H}_6)_2$ (C_1). By measuring the anisotropy parameter (β) from photoelectron angular distribution, we found that the outgoing electron was from a largely La $6s$ -based molecular orbital in both complexes.