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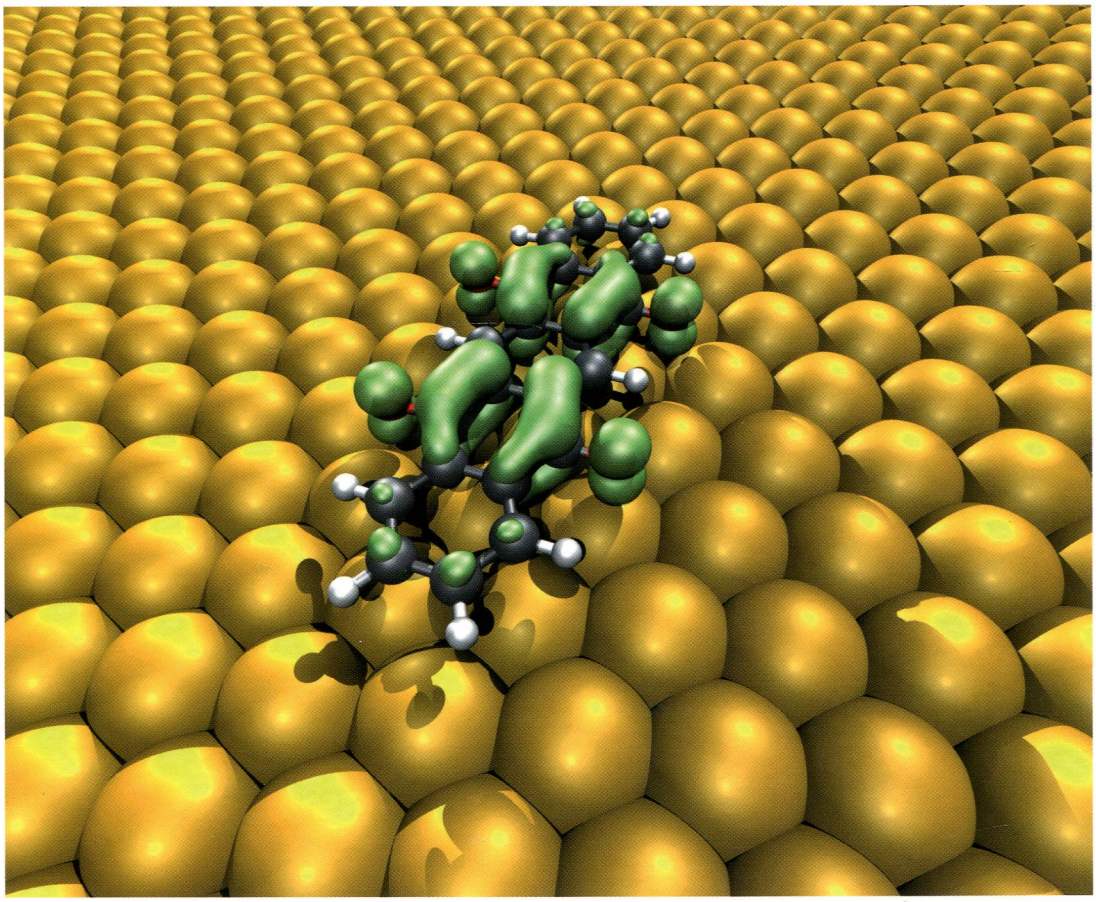
The Molecule-Metal Interface



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Cover Picture
Lowest unoccupied molecular orbital of 5,7,12,14-pentacenetetrone on the Au(111) surface as calculated by density-functional theory. The illustration was created with VMD. VMD was developed by the Theoretical and Computational Biophysics Group in the Beckman Institute for Advanced Science and Technology at the University of Ilinois at Urbana-Champaign. Courtesy of G. Heimel.

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