

Insights on Uranium-Halogen Bonding Derived from Charge-Density Studies at 20 K

Christopher G. Gianopoulos,[†] Vladimir V. Zhurov,[†] Stefan G. Minasian,[‡] Enrique R. Batista,[¶]
Christian Jelsch,[§] A. Alan Pinkerton[†]

[†]Department of Chemistry, School of Green Chemistry and Engineering, The University of Toledo, Toledo, Ohio 43606

[‡]Lawrence Berkeley National Laboratory (LBNL), Berkeley, California 94720

[¶]Theoretical Division, Los Alamos National Laboratory, Los Alamos, New Mexico 87545

[§]CRM2, CNRS, Institute Jean Barriol, Université de Lorraine, Vandoeuvre les Nancy CEDEX, France

Accurate electron density distributions can be obtained from high-resolution X-ray diffraction data measured at low temperatures, even in the case of systems containing very heavy elements such as uranium. Modeling of the experimental electron density distribution with an augmented Hansen-Coppens multipolar formalism and subsequent topological analysis of the model density provides insight into the nature of U-X bonding in the electronic ground-state. The results of charge density studies measured at 20 K for systems containing uranium-halogen bonds are presented and compared to complementary results obtained from gas-phase theoretical calculations.