

OBSERVATION OF A MODERATE STRENGTH INTERACTION OF HYDROGEN WITH A COINAGE METAL HALIDE: THE ROTATIONAL SPECTRUM AND STRUCTURE OF THE *p*-H₂-CuCl AND *o*-H₂-CuCl COMPLEXES

HERBERT M. PICKETT, DANIEL A. OBENCHAIN, G. S. GRUBBS II, and STEWART E. NOVICK, *Department of Chemistry, Wesleyan University, 52 Lawn Avenue, Middletown, CT, 06459-0180, USA.*

Rotational transitions of the *p*-H₂-CuCl and *o*-H₂-CuCl have been observed on a laser ablation equipped FTMW cavity instrument. Computational studies performed using the APFD^a density functional and MP2 level of theory were used to predict the structure of the *p*-H₂-CuCl. Measurements from the J=1–0 to the J=3–2 transitions were used to determine the rotational constants, centrifugal distortion constants, and quadrupole coupling constants for multiple isotopologues of the *p*-H₂-CuCl species. Similar constants, including spin-spin coupling constants, have also been determined for the *o*-H₂-CuCl species for the J=2–1 and the J=3–2 transitions. The *eQq* of the copper in *p*-H₂-⁶³Cu³⁵Cl was found to be 52.058(2) MHz, a change from the monomer ⁶³Cu³⁵Cl value of 16.1712(24) MHz^b.

^aA. Austin, G. A. Petersson, M. J. Frisch, F. J. Dobek, G. Scalmani, and K. J. Throssell. *Chem. Theor. Comp.* **8** (2012) 4989.

^bK. D. Hensel, C. Styger, W. Jager, A. J. Merer, and M. C. L. Gerry, *J. Chem. Phys.* **99**(1993) 3320.