

# SINGLE VIBRONIC LEVEL EMISSION SPECTROSCOPY OF CHCl AND CDCl: VIBRATIONAL STRUCTURE OF THE $X^1A'$ AND $a^3A''$ STATES

C. MUKARAKATE, C. TAO, AND S. A. REID, *Department of Chemistry, Marquette University, P.O. Box 1881, Milwaukee, WI 53201-1881.*

We report on single vibronic level (SVL) emission spectra from bands in the progressions  $2_0^n$ ,  $2_0^n 3_0^m$ ,  $1_0^1 2_0^n$ , and  $1_0^1 2_0^n 3_0^1$  in the  $A^1A''$ - $X^1A'$  system of  $\text{CH}^{35}\text{Cl}$ ,  $\text{CH}^{37}\text{Cl}$ ,  $\text{CD}^{35}\text{Cl}$  and  $\text{CD}^{37}\text{Cl}$ . As in previous studies of CHF/CDF and CHBr/CDBr, the carbenes were generated using a pulsed discharge source, and SVL emission spectra obtained using a 0.3 m spectrograph in combination with a gated, intensified CCD detector. These spectra reveal rich new detail regarding the vibrational structure of the  $X^1A'$  and  $a^3A''$  states, and spin-orbit induced mixing between them, up to  $9000\text{ cm}^{-1}$  above the vibrationless level of the  $X^1A'$  state. For CHCl, we observe around three times the number of  $X^1A'$  levels previously reported,<sup>a</sup> and a number of new  $a^3A''$  state levels. The results of Dunham expansion fits to the vibrational term energies, and comparisons with previous experimental and recent high quality theoretical studies<sup>b,c</sup> will be reported. Overall, the derived vibrational parameters of the  $X^1A'$  and  $a^3A''$  states are in excellent agreement with *ab initio* predictions, including our own DFT calculations.

---

<sup>a</sup>C.-S. Lin et al., J. Chem. Phys. 121, 4164 (2004).

<sup>b</sup>G. Tarczay et al., Phys. Chem. Chem. Phys. 7, 2881 (2005).

<sup>c</sup>H.-G. Yu et al., Mol. Phys. 104, 47 (2006).