

## OBSERVATION OF HIGH LYING LEVELS OF THE $1_g - 1^1\Pi_g$ STATE OF $K_2$

A. J. ROSS, F. MARTIN and I. RUSSIER-ANTOINE, *Laboratoire de Spectrométrie Ionique et Moléculaire (UMR 5579 CNRS), Bâtiment 205, Université Lyon I, Campus la Doua, 69622 Villeurbanne Cedex, France*;  
H. M. CHEN, M. PICHLER, H. WANG and W. C. STWALLEY, *Department of Physics, University of Connecticut, Storrs, CT 06269-3046, U.S.A.*

Very high-lying vibrational levels  $87 \leq v'' \leq 138$  of the  $1_g - 1^1\Pi_g$  electronic state of  $K_2$  have been observed in a photoassociation experiment performed in Connecticut, using ionization detection. Photoassociation of ultracold potassium atoms furnishes accurate binding energies for  $J \leq 4$  in these levels. These binding energies have been treated with laser induced fluorescence data covering the levels  $0 \leq v'' \leq 105$  at high  $J$  (infrared  $C^1\Pi_u \rightarrow 1^1\Pi_g$  transitions recorded by Fourier transform spectrometry) to construct an accurate potential curve for this electronic state.

Because several vibrational levels were observed both in photoassociation and in fluorescence, an experimental dissociation energy can be deduced from the sum of binding energies (measured in photoassociation) and vibrational energies established with respect to the potential minimum (from fluorescence data), giving  $D_e = 1230.297 \pm 0.002 \text{ cm}^{-1}$ . Applying asymptotic models for the interactions between  $K(4s) + K(4p)$  atoms to the molecular potential energy curve, a slightly lower value,  $1230.292 \pm 0.002 \text{ cm}^{-1}$  is obtained. This is considered to be the best available value for the dissociation energy of the  $1^1\Pi_g$  electronic state into  $K(4s) + K(4p)^2P_{3/2}$  atoms.