

Fourier Transform Spectroscopy Study of Fully Mixed $A^1\Sigma^+$ and $b^3\Pi$ States of KCs Molecule

A. Kruzinsh¹, I. Klincare¹, O. Nikolayeva¹, M. Tamanis¹, R. Ferber¹,
E.A. Pazyuk², A.V. Stolyarov²

¹ *Laser Centre, University of Latvia, 19 Rainis Boulevard, LV-1586 Riga, Latvia*

² *Department of Chemistry, Moscow State University, GSP-2 Leninskie gory 1/3,
Moscow 119992, Russia*

The fully mixed $A^1\Sigma^+ \sim b^3\Pi$ complex of the KCs molecule was studied with laser induced fluorescence $A-b \rightarrow X^1\Sigma^+$ spectra Fourier-transform spectrometry (FTS). Present research is a continuation of the recently performed studies of the $A^1\Sigma^+ \sim b^3\Pi$ complex in the NaRb [1] and NaCs [2] molecules, where high accuracy FTS data have been reproduced with an experimental accuracy by a direct deperturbation analysis by means of the inverted channel-coupling approach with Hund's a -coupling case basic functions. The $A^1\Sigma^+ \sim b^3\Pi$ complex in KCs can be considered as fully mixed system due to very strong spin-orbit interaction.

Direct optical excitation of the KCs $A^1\Sigma^+ \sim b^3\Pi$ complex in a heat-pipe was performed by the diode lasers. Three laser diodes 850 nm, 980 nm and 1020 nm were used with the respective laser frequency tuning ranges 11943 -11550 cm^{-1} , 10515 – 10209 cm^{-1} , and 9940-9700 cm^{-1} . Spectra analysis and determination of the upper state energies was based on the accurate experimental ground state potential energy curve [3]. The laser induced fluorescence spectra typically contained rich collisionally induced rotational relaxation line structure. The analysis of rotational satellites has enlarged the data set of $A-b$ complex rovibronic term values significantly. Currently more than 2200 rovibronic term values have been obtained, including the data on the predominantly $A^1\Sigma^+$ state covering vibrational level v values from $v = 0$ to about 50 and rotational levels J from $J = 9$ to 210, as well as term values data of the predominantly $b^3\Pi_0$ state. We estimate experimental term values uncertainty as 0.01 cm^{-1} .

The experimental term values will be processed by a direct channel-coupling deperturbation approach in order to refine the potential electronic curves of the $A^1\Sigma^+$ and $b^3\Pi$ states of KCs and the spin-orbit coupling functions on the internuclear distance which would reproduce the experimental termvalues and fluorescence intensity distributions within an accuracy of the measurements. Work is in progress.

Moscow team thanks for support the Russian Foundation for Basic Research (grant N06-03-32330a). The Riga team acknowledges Latvian Science Council grant No. 04.1308. , O.N. acknowledges ESF grant.

[1] O. Docenko, M. Tamanis, R. Ferber, E. A. Pazyuk, A. Zaitsevskii, A. Stolyarov, A. Pashov, H. Knöckel, E. Tiemann, *Phys. Rev. A* **75**, 042503 (2007).

[2] J. Zaharova, M. Tamanis, R. Ferber, E. A. Pazyuk, and A.V. Stolyarov, to be published.

[3] R. Ferber, I. Klincare, O. Nikolayeva, M. Tamanis, A. Pashov, H. Knöckel, E. Tiemann, *J. Chem. Phys.*, (2008), submitted.