Abstract Submitted for the DAMOP09 Meeting of The American Physical Society

Global analysis of data on the  $A^{1}\Sigma^{+} \sim b^{3}\Pi$  states of NaK<sup>1</sup> HOUS-SAM SALAMI, THOMAS BERGEMAN, SUNY Stony Brook, AMANDA ROSS, U. Lyon 1 — NaK electronic states have been extensively studied over recent years. Many of these studies have involved the lowest excited states,  $A^{1}\Sigma^{+}$  and  $b^{3}\Pi$ , as they offer pathways to higher states. These states have regained attention as they are used as intermediaries in the production of ultracold molecules. Recently, the  $b^{3}\Pi_{0} \sim A^{1}\Sigma^{+}$  spin-orbit interactions have been examined [1,2]. However analysis was based on band-by-band local deperturbation which is not consistent with experimental errors limits. In this study, we collect existing (published and unpublished) data from various experiments (FT-LIF, PLS and PFOODR) performed at Orsay and Lyon (Ross et al.), Warsaw (Kowalczyk et al.) and Lehigh (Huennekens et al.) universities. Transitions from  $D^{-1}\Pi$ ,  $B^{-1}\Pi$ , and  $C^{-1}\Sigma^+$  states simultaneously to the  $A \sim b$  and X states constitute the bulk of the data as the upper term energy values can be deduced precisely from the well known X state parameters. Combined data will be modeled using a global deperturbation approach employing the discrete variable representation (DVR) so as to fit potential energy and spin-orbit functions. 1. P. Burns et al., J. Chem. Phys. **122**, 074306 (2005).

2. R. Ferber et al., J. Chem . Phys.  ${\bf 112},\,5740$  (2000).

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