

## LASERSPECTROSCOPY OF THE B<sup>2</sup>Σ STATE OF LiAr

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The rovibrational structure of the transition B<sup>2</sup>Σ ← X<sup>2</sup>Σ of LiAr has been investigated by means of high-resolution laser spectroscopy using a supersonic beam for production of the molecules<sup>a</sup>. About 100 absorption lines could successfully be assigned for each isotopomer <sup>7</sup>LiAr and <sup>6</sup>LiAr providing in addition the vibrational numbering of the observed levels  $v = 0 \dots 3$  of the B<sup>2</sup>Σ state. The analysis of the experimental observations was complicated by the occurrence of perturbations between rovibrational levels of the B<sup>2</sup>Σ and of the A<sup>2</sup>Π state which have not been taken into account, up to now.

For the B<sup>2</sup>Σ interaction potential we used a Morse–van der Waals potential as an analytical expression reproducing the observed energy levels with an error margin of 0.015cm<sup>-1</sup>. The van der Waals parameter C<sub>6</sub> was fixed to its theoretical value. Our preliminary results for the B<sup>2</sup>Σ state of <sup>7</sup>LiAr are  $R_e = 6.24(5)$  Å and  $D_e = 33.8(2.0)$ cm<sup>-1</sup>.

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<sup>a</sup>R. Brühl and D. Zimmermann, Chem.Phys.Lett. 233 (1995) p.455