

HIGH RESOLUTION INFRARED AND MICROWAVE STUDY OF $^{10}\text{BF}_2\text{OH}$ AND $^{11}\text{BF}_2\text{OH}$: the 5^1 , 6^1 , 7^1 , 8^1 , 9^1 AND $8^1 9^1$ VIBRATIONNALLY EXCITED STATES.

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High resolution ($2\text{-}3 \times 10^{-3} \text{ cm}^{-1}$) Fourier transform infrared spectra of gas phase ^{10}B and ^{11}B enriched isotopic and natural samples of BF_2OH (difluoroboric acid) were recorded in the $400\text{-}4000 \text{ cm}^{-1}$ spectral range. Starting from the results of a previous study^a which involved the ν_8 (BF_2 out-of-plane bending) and ν_9 (OH torsion) bands of $^{11}\text{BF}_2\text{OH}$, it has been possible to perform the first rovibrational analysis of the ν_5 (BF_2 bending), ν_8 , ν_9 and $\nu_8+\nu_9$ bands of $^{10}\text{BF}_2\text{OH}$, and of the ν_5 , ν_7 (F_2BO in-plane bending) and $\nu_8+\nu_9$ bands of $^{11}\text{BF}_2\text{OH}$ up to very high rotational quantum numbers. In addition microwave transitions within the 5^1 , 6^1 , 7^1 , 8^1 and 9^1 vibrational states of $^{11}\text{BF}_2\text{OH}$ were measured using predictions performed from ab initio calculations. For the ν_5 and ν_7 bands of $^{11}\text{BF}_2\text{OH}$, C-type Coriolis resonances coupling the 5^1 and 7^1 energy levels with those of the 7^2 and 6^1 dark states respectively were accounted for in the calculations. The other bands appeared to be unperturbed,

^aD.Collet, A.Perrin, H.Bürger, and J.-M.Flaud, *J. Mol. Spectrosc.* 212, 118 (2002)