

C_{2v} -TOP INTERNAL ROTATION IN THE HIGH-RESOLUTION UV SPECTRA OF WEAKLY BOUND COMPLEXES: II. APPLICATIONS.^a

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The model of a C_{2v} top – C_s frame internal rotation discussed in Part I has been used in the analysis of rotationally resolved $S_1 \leftarrow S_0$ fluorescence excitation spectra of three different complexes; aniline–N₂, *para*-difluorobenzene–N₂, and benzonitrile–H₂O. The results of the internal rotation analysis as well as the structure determination will be presented and the limits of the applied model will be discussed.

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