

HIGH RESOLUTION ELECTRONIC SPECTRA OF 1,2-DIMETHOXYBENZENE AND THE 1,2-DIMETHOXYBENZENE/WATER COMPLEX<sup>a</sup>

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High resolution  $S_1 \leftarrow S_0$  fluorescence excitation spectra of 1,2-dimethoxybenzene (DMB) and its water complex have been obtained and analyzed. The derived values of the rotational constants show that DMB itself has a planar heavy-atom structure with trans-disposed methoxy groups in both electronic states. No evidence for any other conformational isomers was found. The water complex origin appears  $127 \text{ cm}^{-1}$  to the blue of the bare molecule origin and is split into two spectra separated by  $0.04 \text{ cm}^{-1}$ . Analyses of these data show that the water molecule is attached via two  $\text{O-H}\cdots\text{O}$  hydrogen bonds to the methoxy groups, and undergoes a hindered internal rotation. Differences in the nature of this motion in the two electronic states will be discussed.

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<sup>a</sup>Work supported by NSF.