

Diverse hydrogen-bonded structural motifs in 1,4-diazabicyclo[2.2.2]octane N,N'-dioxide salts with oxoanions

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Supporting Information

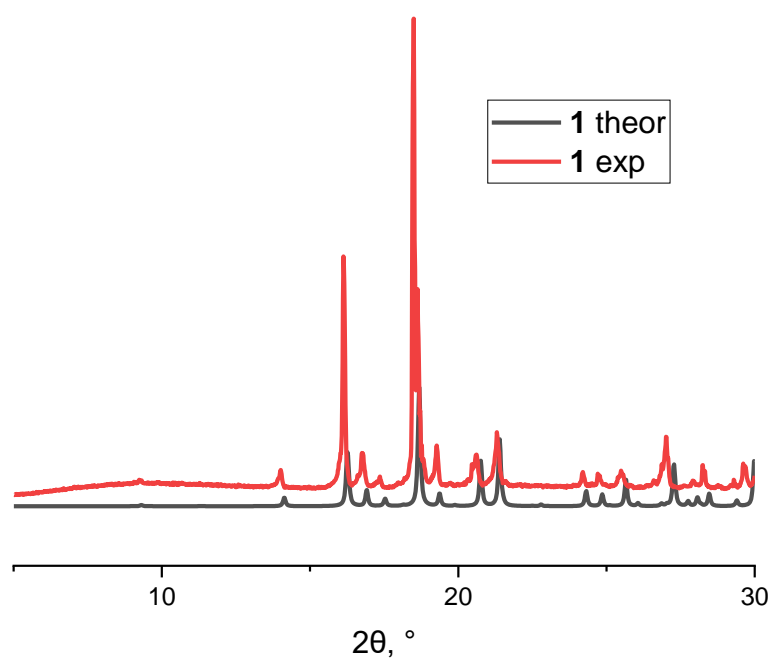


Figure S1. Experimental PXRD pattern of 1 sample compared to the theoretical one.

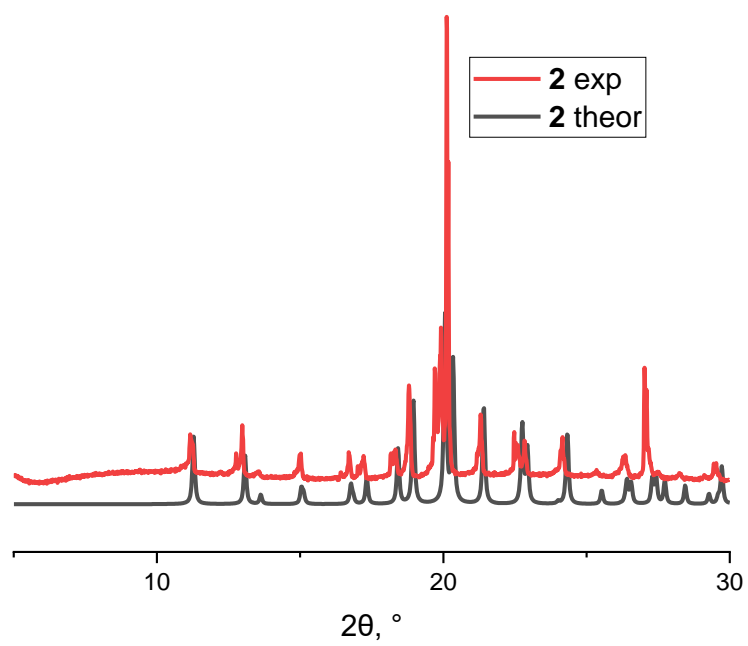


Figure S2. Experimental PXRD pattern of 2 sample compared to the theoretical one.

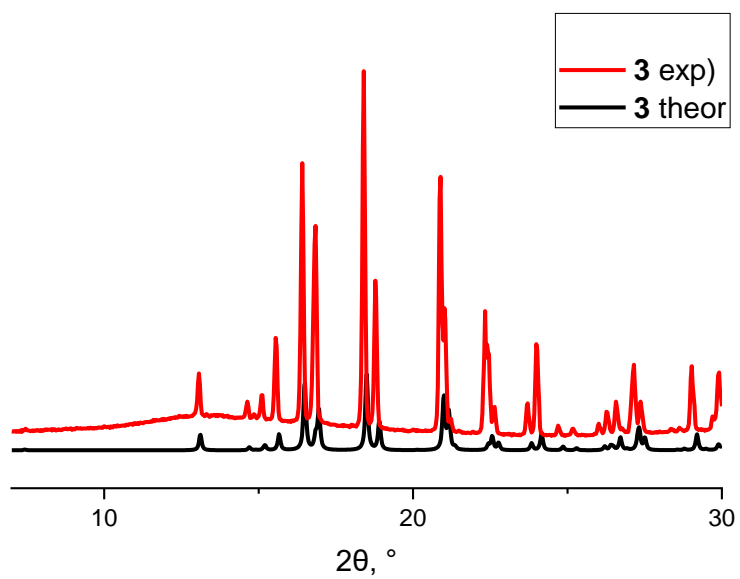


Figure S3. Experimental PXRD pattern of 3 sample compared to the theoretical one.

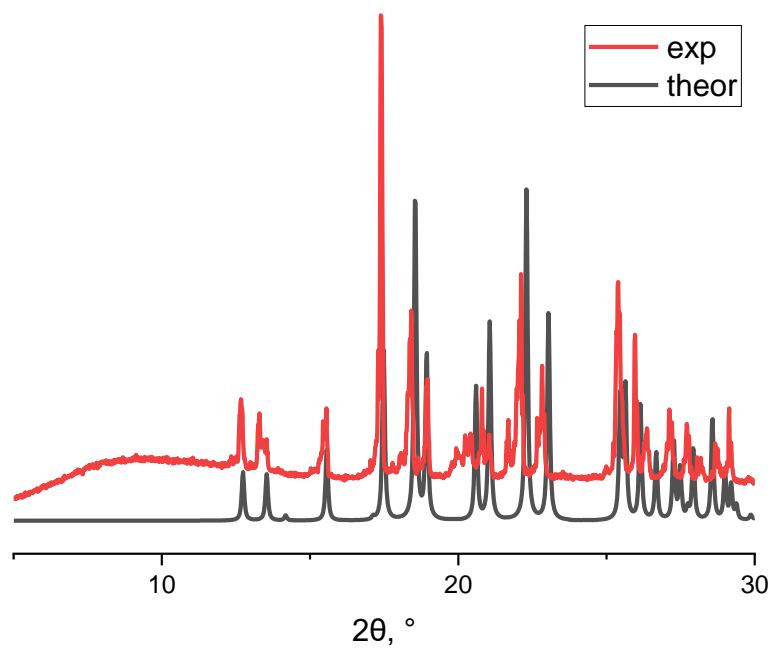


Figure S4. Experimental PXRD pattern of 4 sample compared to the theoretical one.

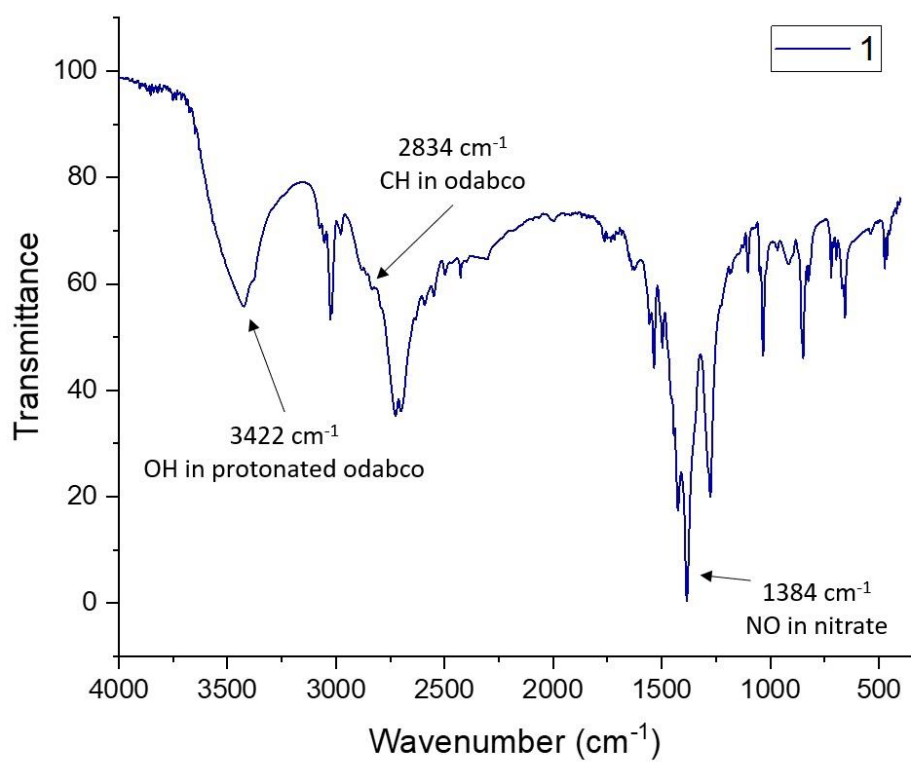


Figure S5. Infrared spectrum of 1.

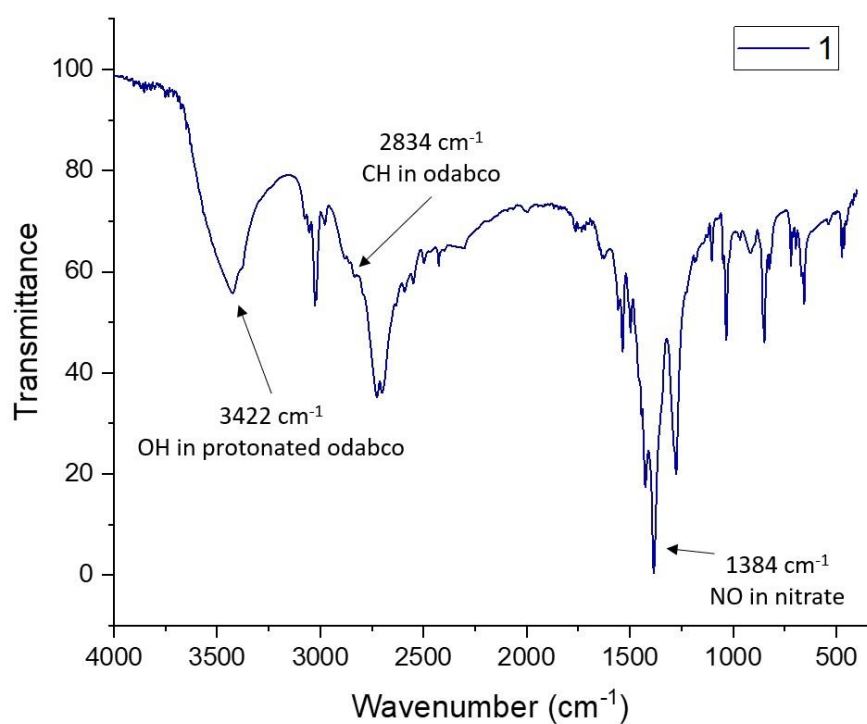


Figure S6. Infrared spectrum of 2.

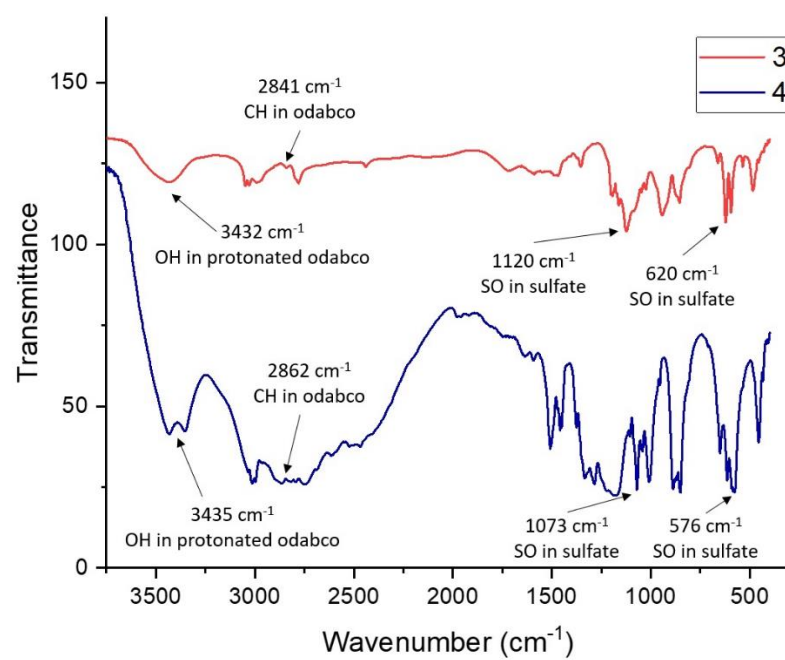


Figure S7. Infrared spectra of 3 and 4.

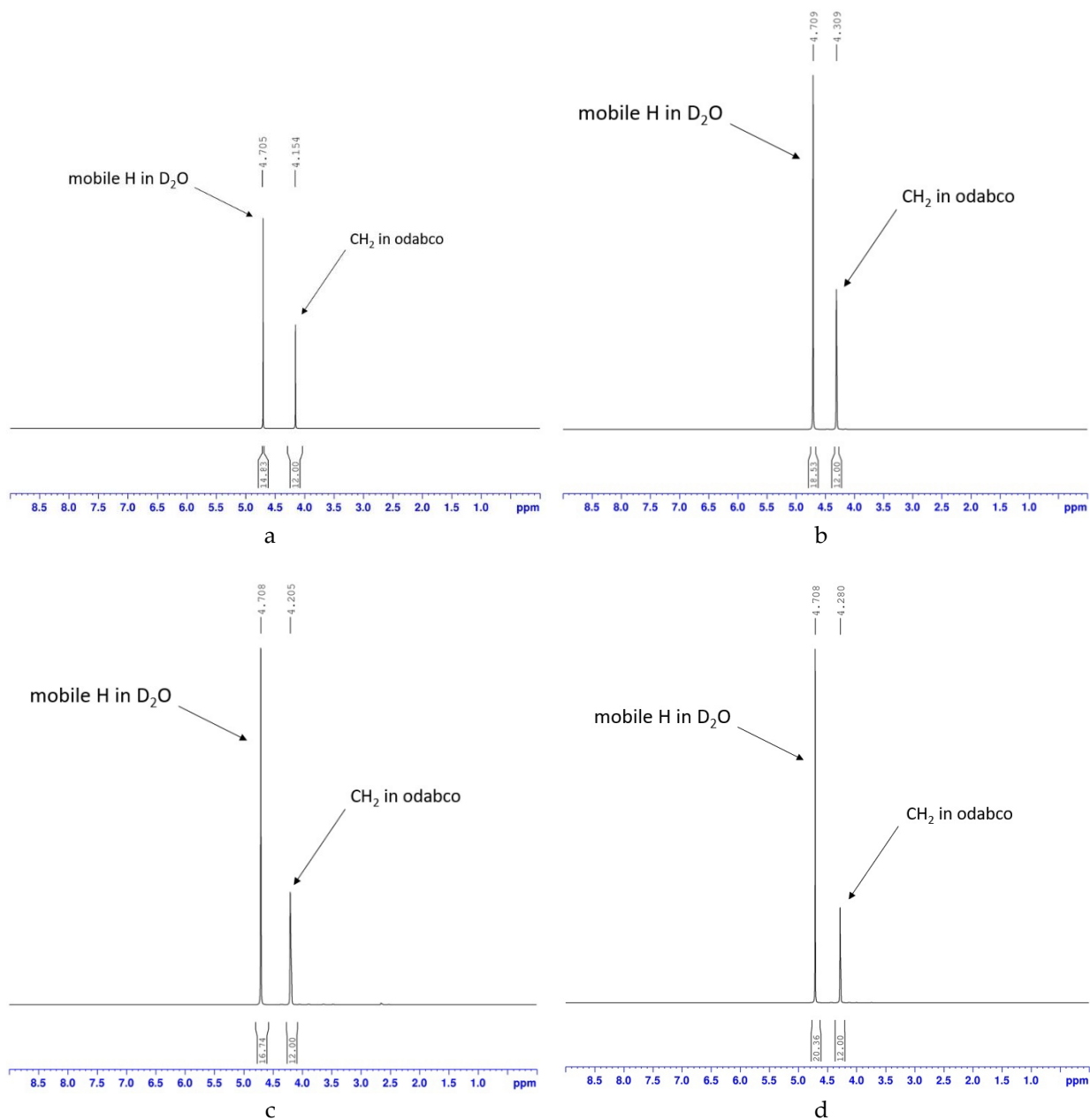


Figure S8. ^1H NMR spectra for **1** (a), **2** (b), **3** (c) and **4** (d) in D_2O . Peaks at ~ 4.71 ppm correspond to the residual mobile ^1H protons in D_2O .