

Figure S1: FT-IR spectrum of **1**.

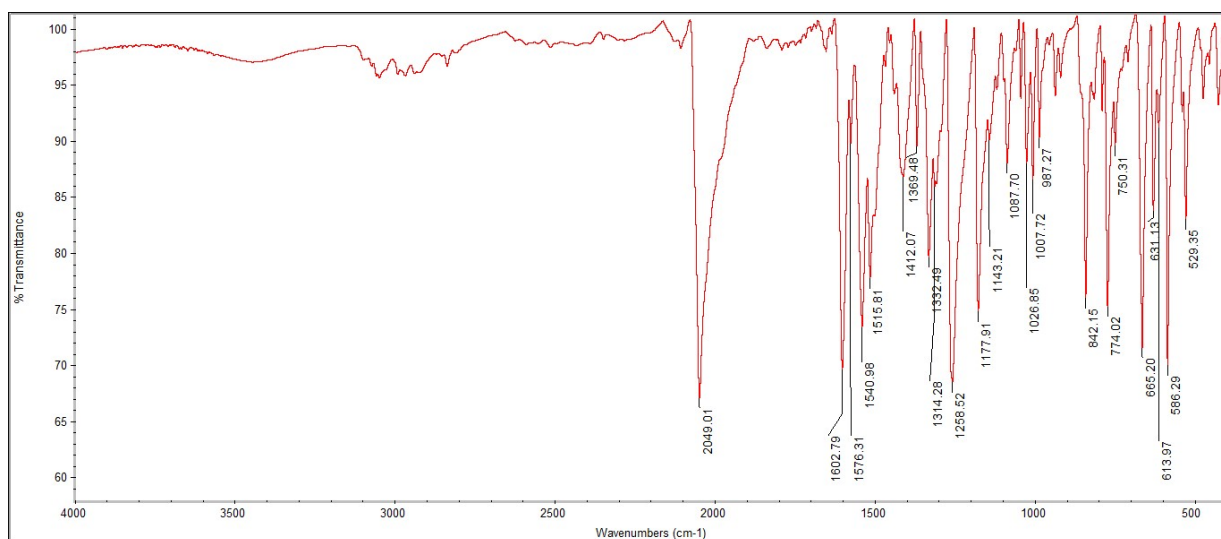


Figure S2: FT-IR spectrum of **2**.

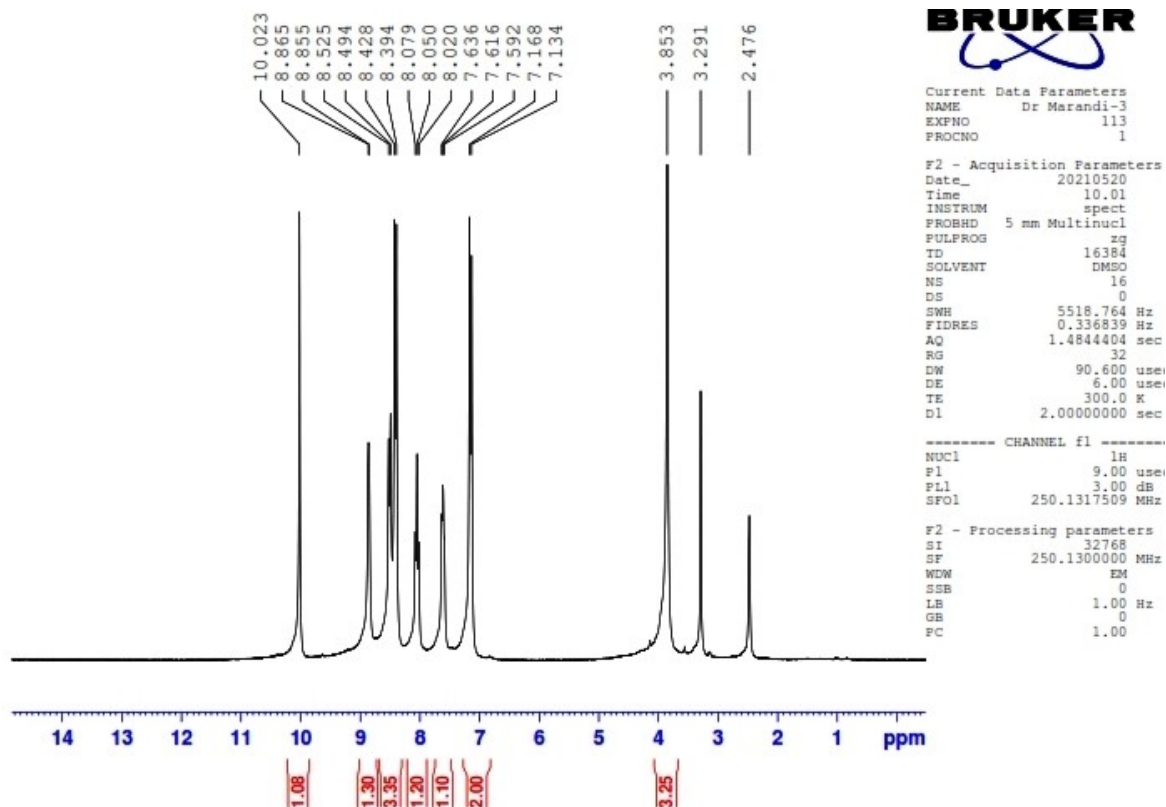


Figure S3: 250 MHz ^1H NMR spectrum of **1** in $\text{DMSO-}d_6$.

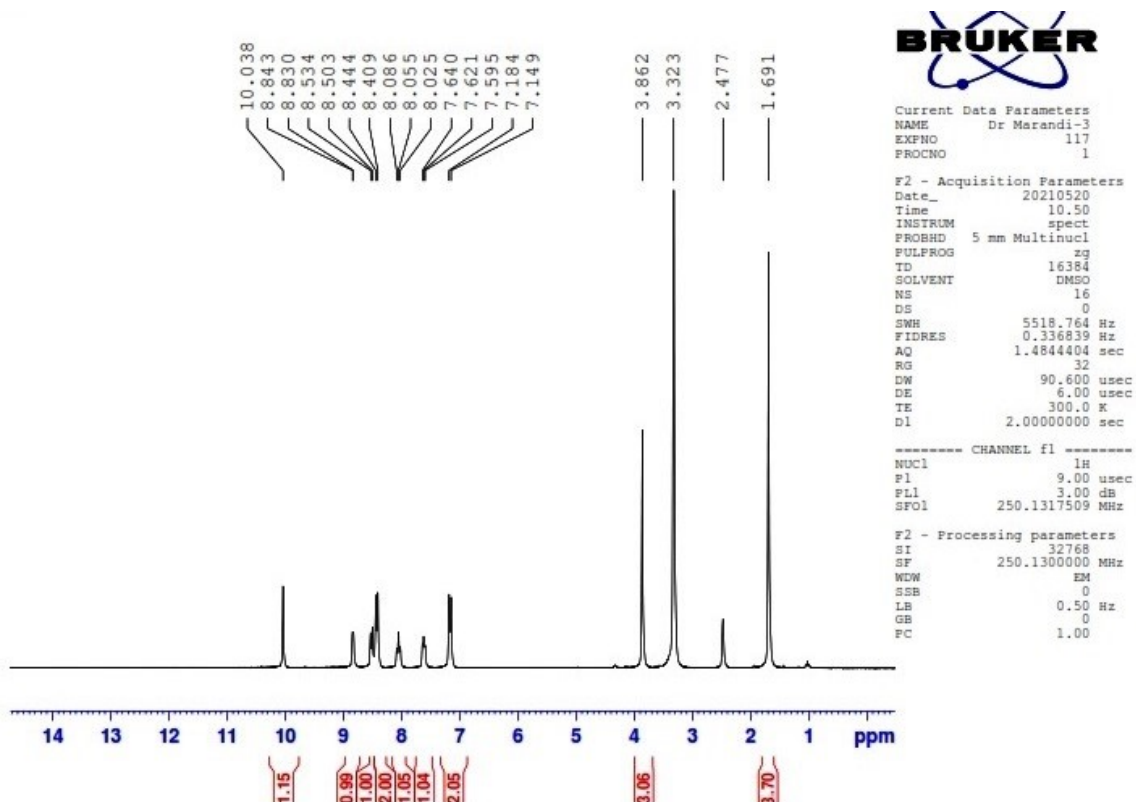


Figure S4: 250 MHz ^1H NMR spectrum of **2** in $\text{DMSO-}d_6$.

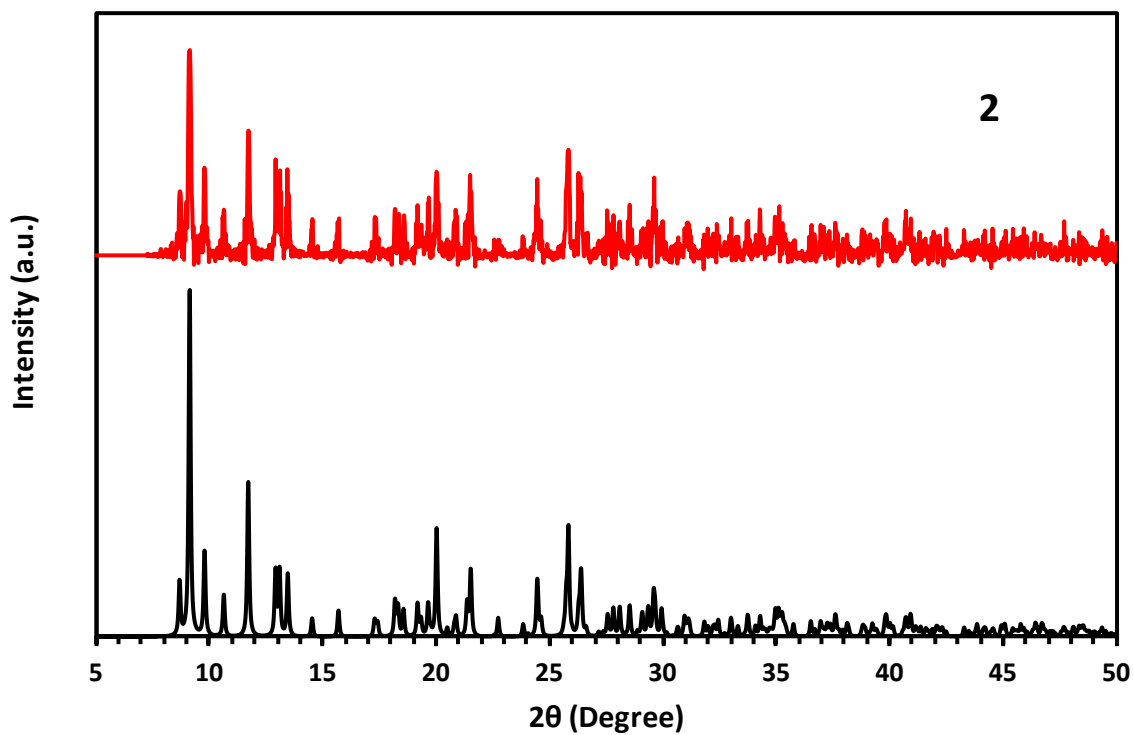
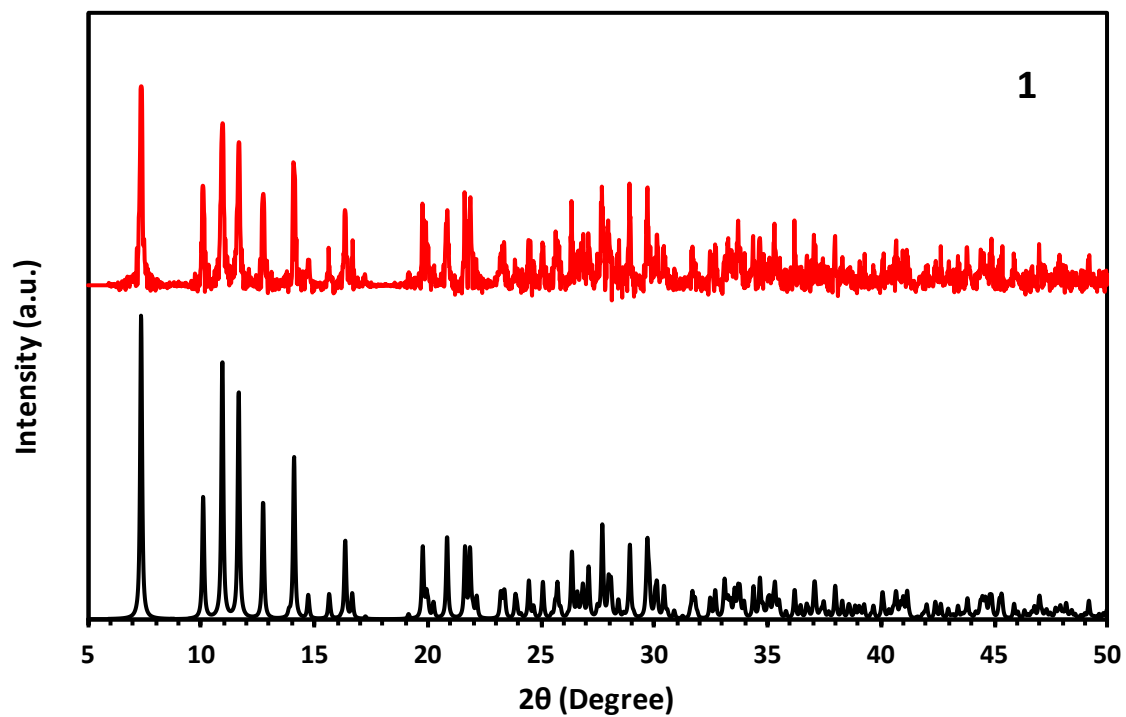
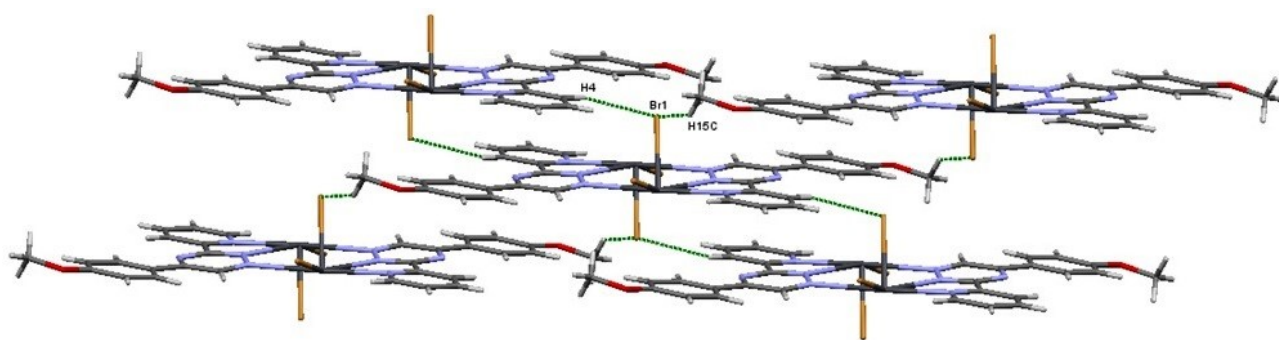
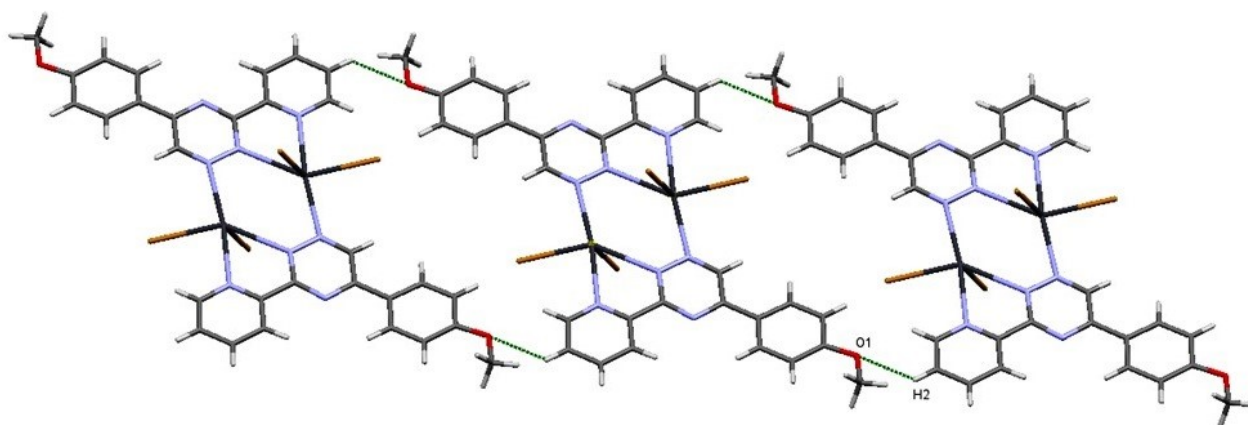


Figure S5. Experimental (in red) and simulated (in black) powder X-ray diffraction patterns of **1** and **2**. The simulated pattern is based on the structure refined by single-crystal X-ray diffraction analysis.



(a) •



(b) •

Figure S6: Crystal packing in **1** by a) C-H...Br b) C-H...O intermolecular interactions.

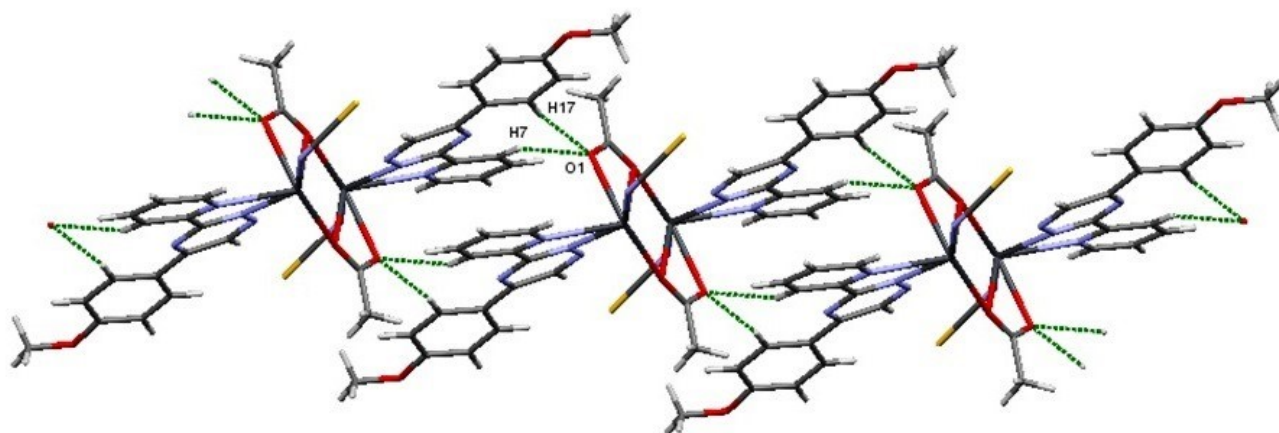


Figure S7: Crystal packing by C-H...O intermolecular interactions and π - π stacking in **2**