

**Versatility in the binding modes of 4,6-dimethyl-1,2,3-triazolo[4,5-*d*]-pyrimidin-5,7-dionato in presence of bipyrimidine. Analyses of their supramolecular architectures through H-bonds.**

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**Electronic supplementary information**

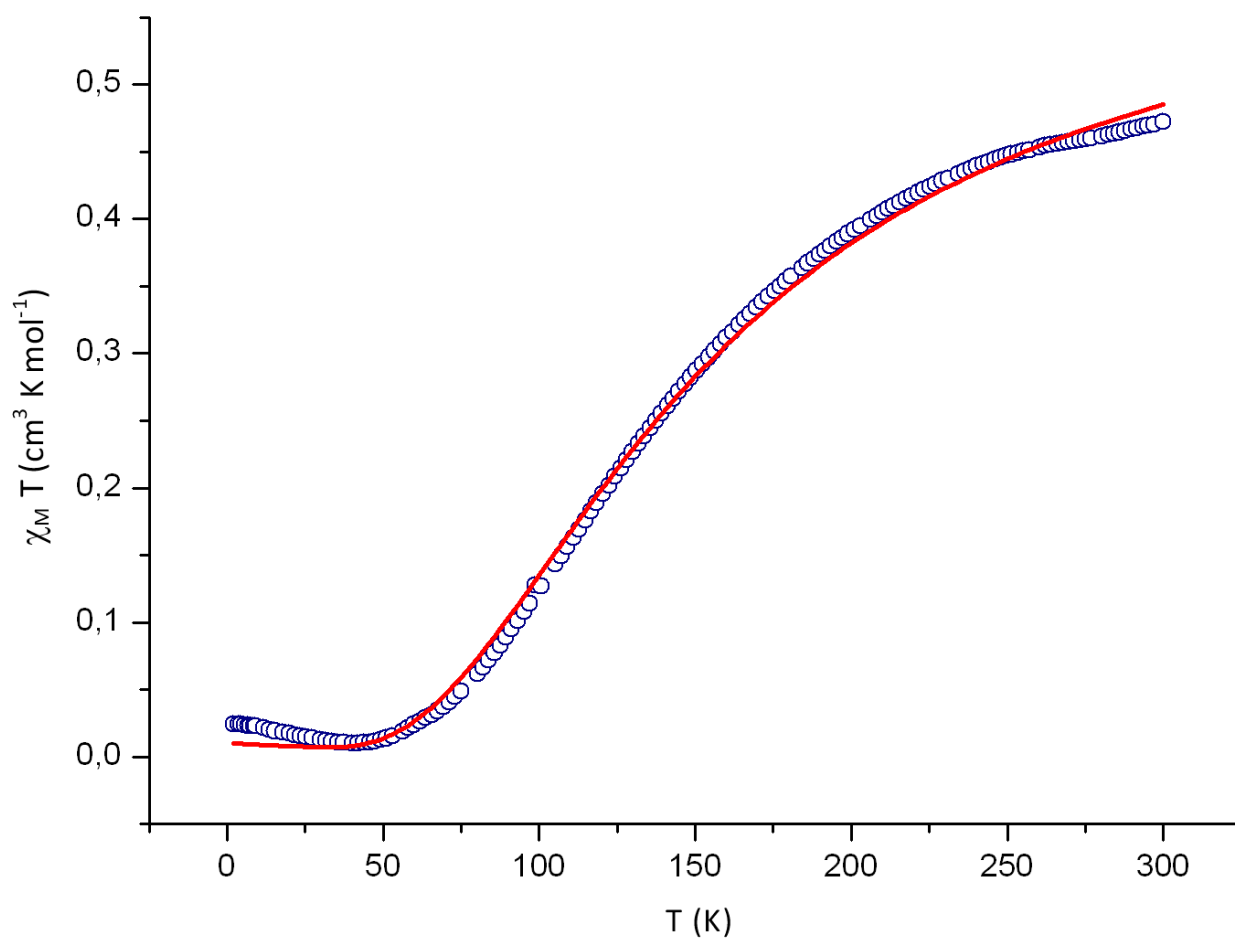


Figure S-1: Experimental and calculated (from Bleany-Bowers equation, see text) molar magnetic susceptibility of compound **3**.

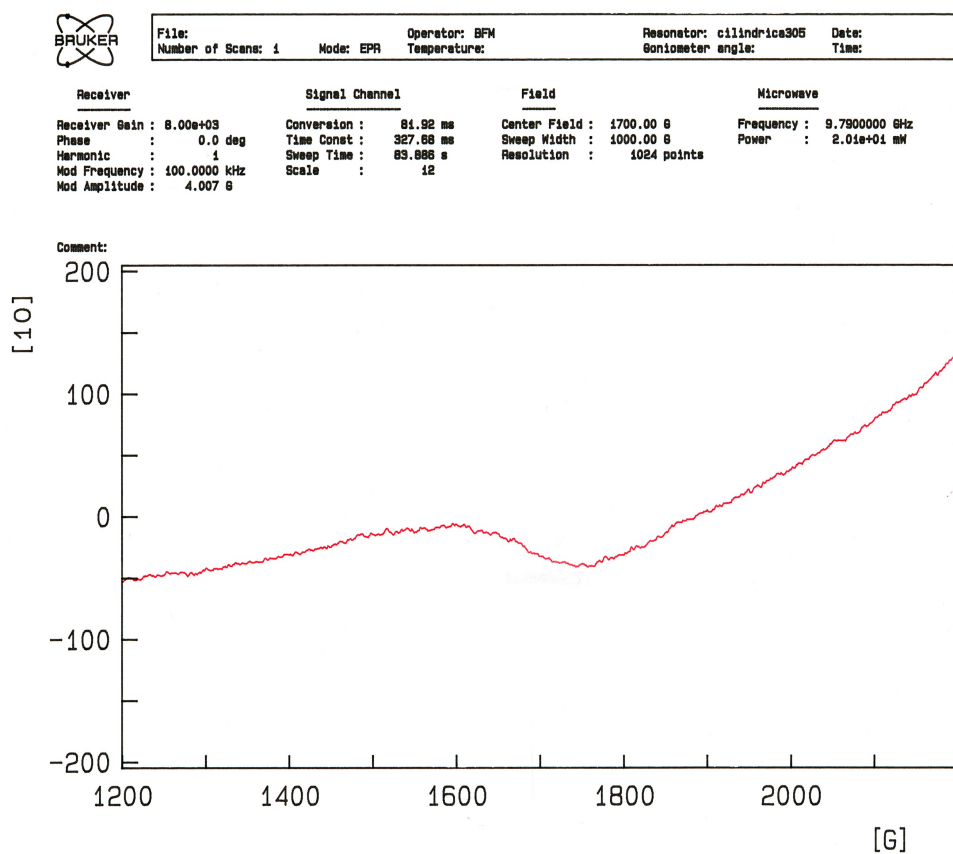
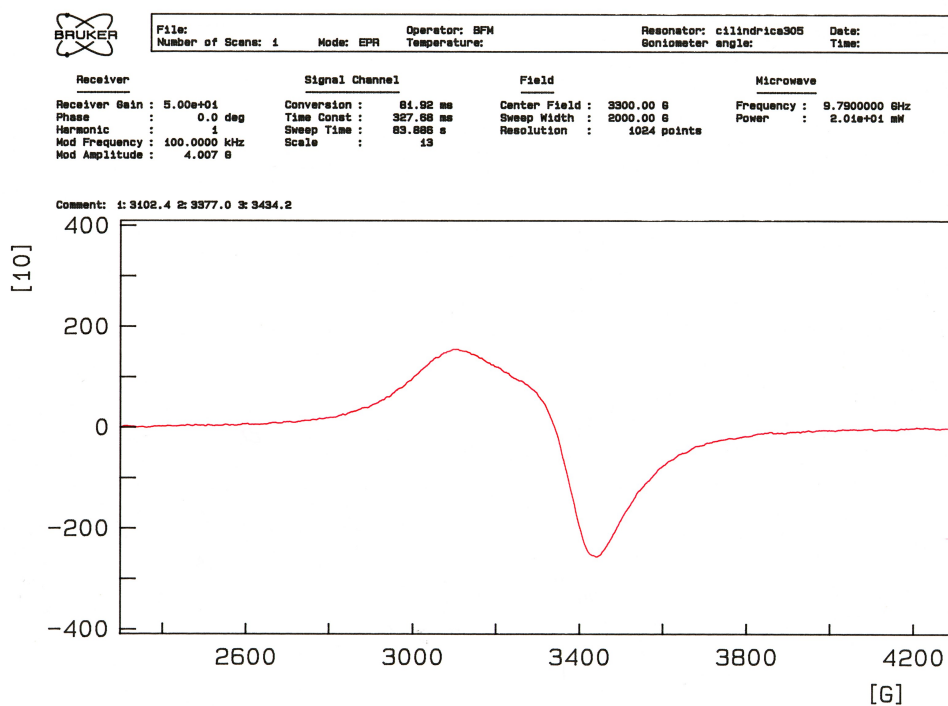


Figure S-2: X-band EPR spectrum of compound **3**: g=2 region (top) and half-field region (bottom).