

Combining oxime-based $[Mn_6]$ clusters with cyanometalates: 1D chains of $[Mn_6]$ SMMs from $[M(CN_2)]^-$ ($M = Au, Ag$)

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Experimental Procedures

Solvents and reagents were used as received from commercial suppliers.

Synthesis of compound 1: $Mn(NO_3)_2 \cdot 4H_2O$ (125 mg, 0.5 mmol), 2-hydroxypropiophenone oxime (80 mg, 0.5 mmol), 3-ethynlpyridine (50 mg, 0.25 mmol) and $KAu(CN)_2$ (144 mg, 0.5 mmol) were dissolved in MeOH (20 ml). After 5 minutes of stirring $LiOMe$ (37 mg, 1 mmol) was added, and the solution stirred for a further 1 h, before being filtered and allowed to stand. Black rod-like X-ray quality crystals were obtained after room temperature evaporation of the mother liquor over 4 days. Elemental analysis (%) calculated (found) for $C_{38}H_{39}AuMn_3N_6O_9$ (1085.23): C 42.03 (41.92), H 3.62 (3.49), N 7.74 (7.58).

Synthesis of compound 2: $Mn(NO_3)_2 \cdot 4H_2O$ (125 mg, 0.5 mmol), 2-hydroxyacetophenone Oxime (75 mg, 0.5 mmol), 3-ethynlpyridine (50 mg, 0.25mmol), $KAu(CN)_2$ (128 mg, 0.5 mmol), were dissolved in MeOH (20 ml). After 5 minutes of stirring $LiOMe$ (37 mg, 1 mmol) was added, and the solution stirred for a further 1 h, before being filtered. The filtrate was collected and allowed to stand for 24 h before being filtered again. Black rod-like X-ray quality crystals were obtained after room temperature evaporation of the mother liquor over 4 days. Elemental analysis (%) calculated (found) for $C_{34}H_{30}AgMn_3N_6O_8$ (923.33): C 44.21 (44.14), H 3.28 (3.24), N 9.10 (9.01).

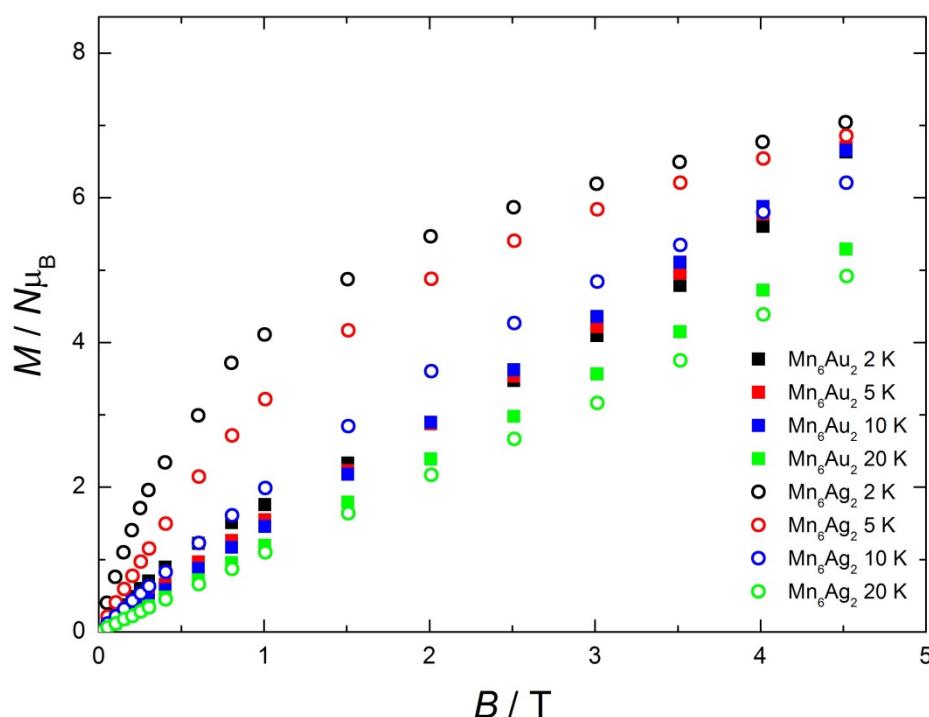


Figure S1. Plot of magnetisation versus field for **1** and **2** in the 2.0 to 20 K temperature range, in fields up to 5 T.

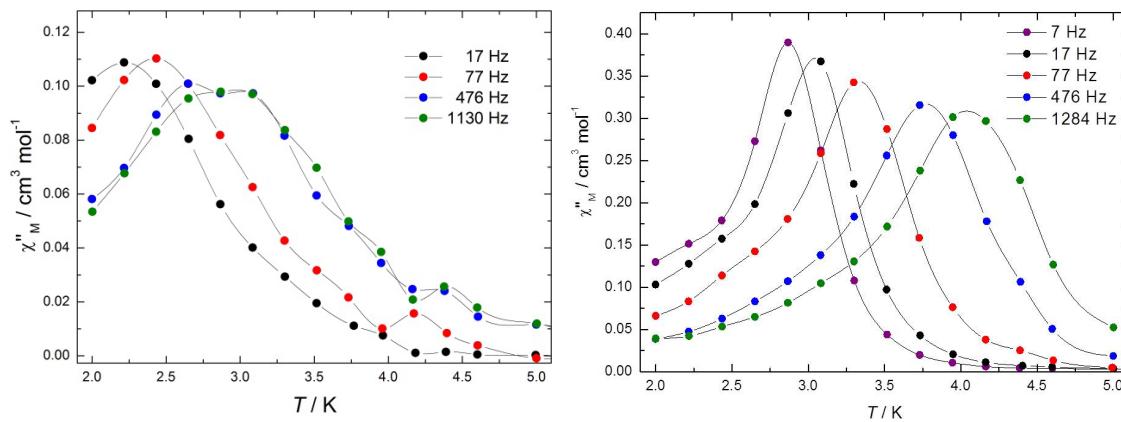


Figure S2. Plot of χ'' vs. T for compound compound **1** (left) and **2** (right) in the indicated temperature and frequency ranges.

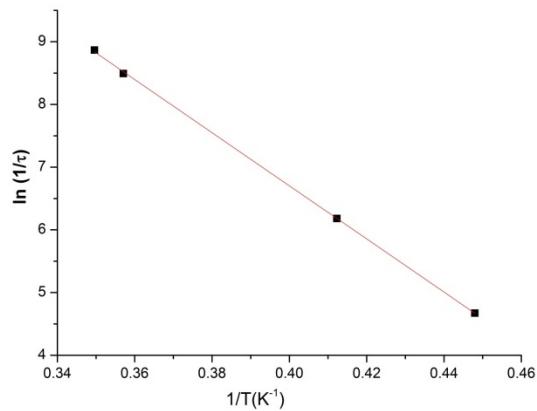


Figure S3. Arrhenius analysis of the ac susceptibility data for **1**; $\tau_0 = 1.5 \times 10^{-10}$ s and $U_{\text{eff}} = 39.9$ K.

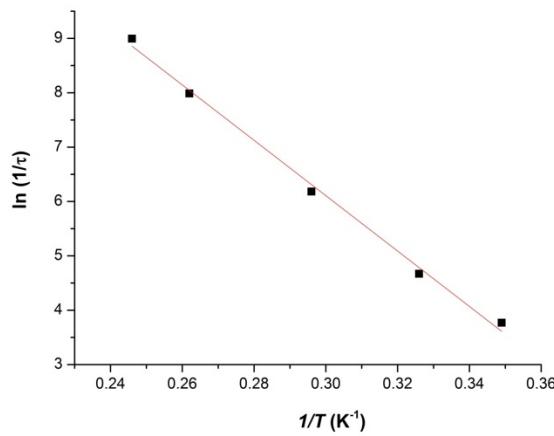


Figure S4. Arrhenius analysis of the ac susceptibility data for **2**; $\tau_0 = 5.4 \times 10^{-10}$ s and $U_{\text{eff}} = 50.7$ K.

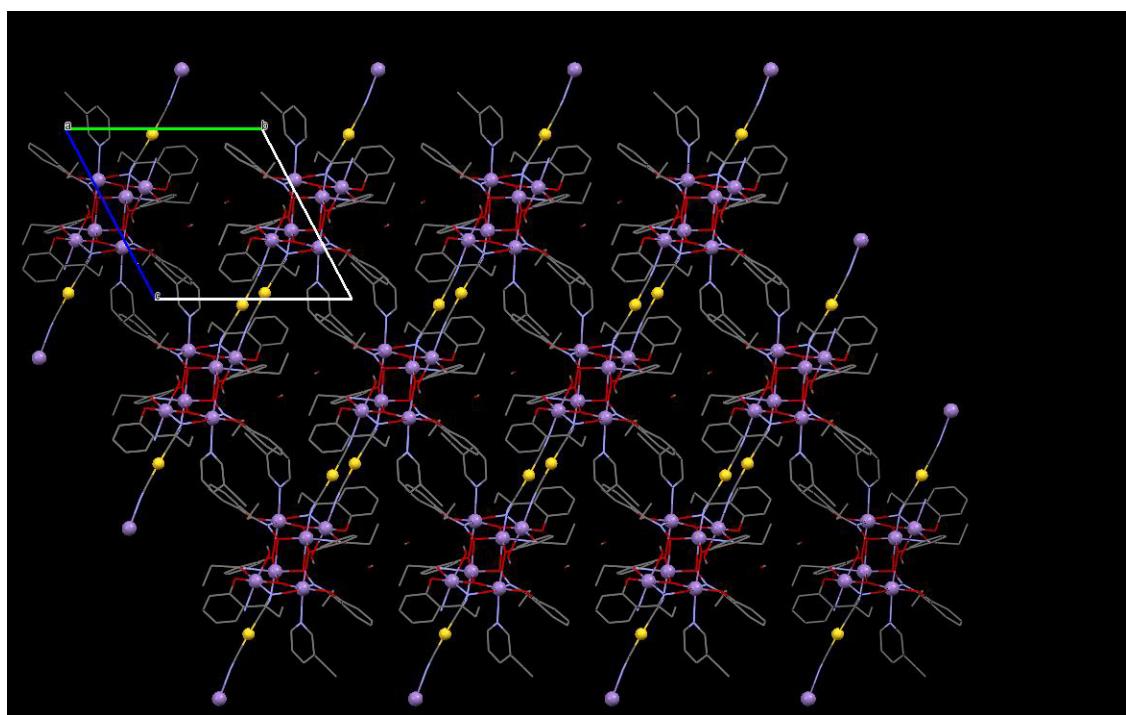


Figure S5. Packing of the chains of **1** in the crystal. H-atoms omitted for clarity.

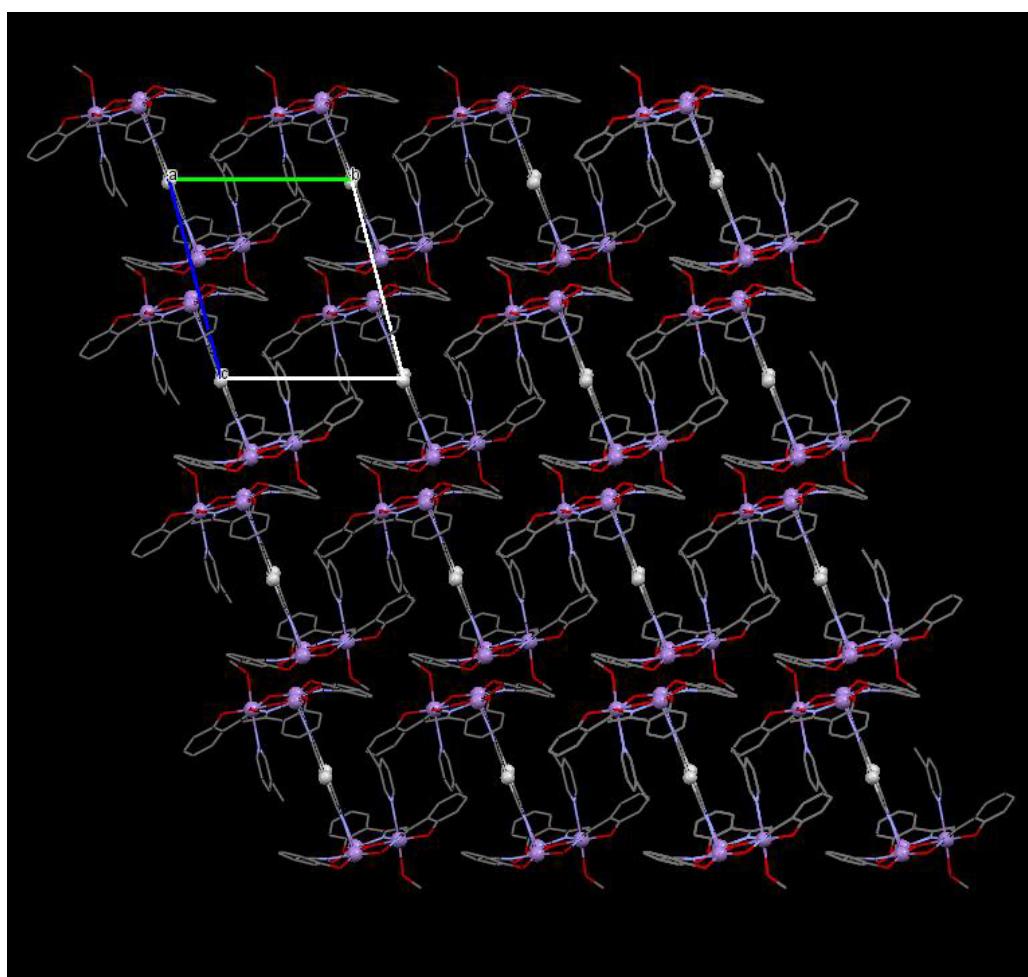


Figure S6. Packing of the chains of **2** in the crystal. H-atoms omitted for clarity.