

Electronic Supplementary Information for Dalton Transactions
This journal is © The Royal Society of Chemistry 2008

Supporting Information

Contents	Page
Figure S1	3
Figure S2	4
Figure S3	5
Figure S4	6
Figure S5	7
Figure S6	8
Figure S7	9
Figure S8	10
Figure S9	11

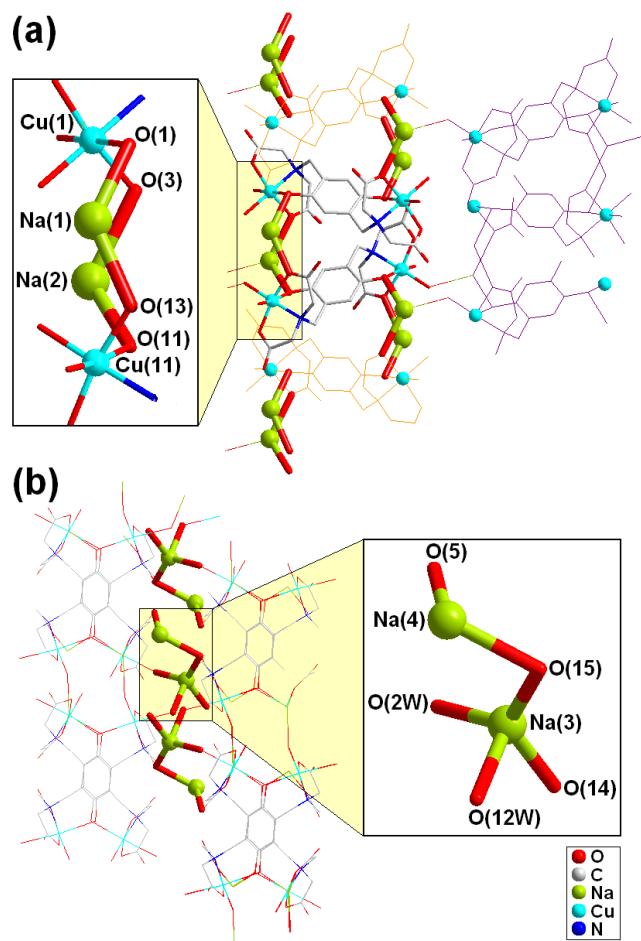


Figure S1. Representative structural drawings of **6** showing the formation of the 3D self-assembly polymeric structure via the interactions of Na^+ with the oxygen atoms of the ligand. (a) The $\text{Na}(1)$ and $\text{Na}(2)$ atoms form bridges [$\text{O}(11)\text{-Na}(2)\text{-O}(3)$ and $\text{O}(1)\text{-Na}(1)\text{-O}(13)$] that connect two complexes in a zig-zag 1D polymer. In addition, $\text{Na}(2)$ forms bonds with the $\text{O}(11)$ of the neighbouring 1D polymers creating 2D sheets. (b) The $\text{Na}(3)$ and $\text{Na}(4)$ atoms operate as bringing atoms [$\text{O}(15)\text{-Na}(3)\text{-O}(2w)\text{O}(12w)\text{O}(14)$, $\text{O}(15)\text{-Na}(4)\text{-O}(3)$] between the 2D polymers resulting in the 3D polymer.

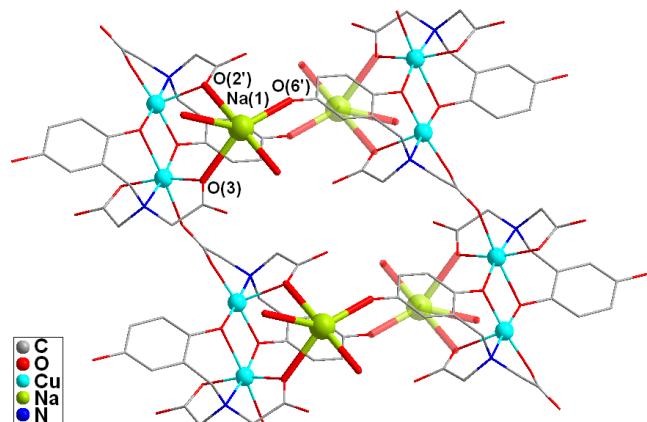


Figure S2. A representative structural drawing of **9** where the 1D polymers are coupled to each other through interaction of the sodium counter ions with the second free phenolic oxygen [$O(6')$], and the carboxylate oxygen atoms [$O(2')$] and [$O(3)$].

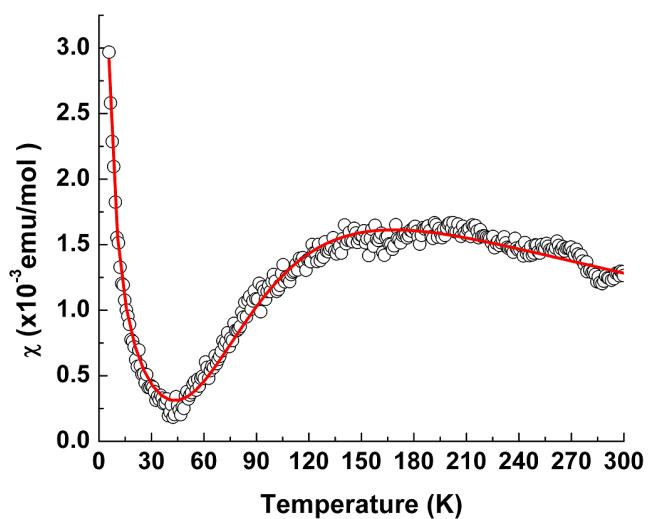


Figure S3. Molar magnetic susceptibility of **7** as a function of temperature. The red line is a fit according to the modified Bleany–Bowers equation (equation 1).

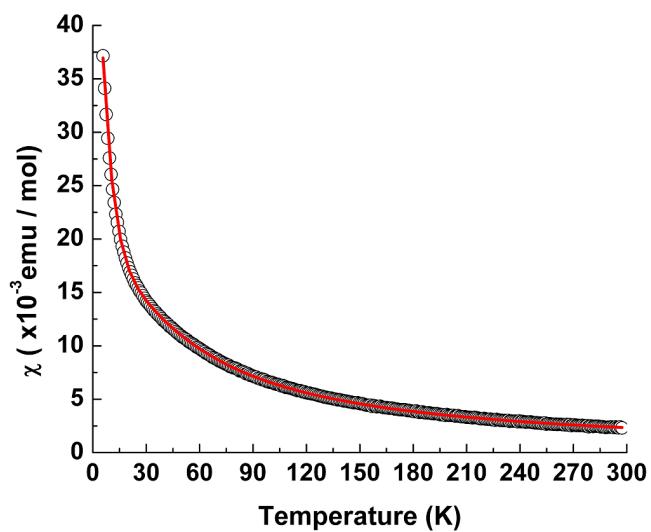


Figure S4. Molar magnetic susceptibility of **9** as a function of temperature. The red line is a fit according to the modified Bleany–Bowers equation (equation 1).

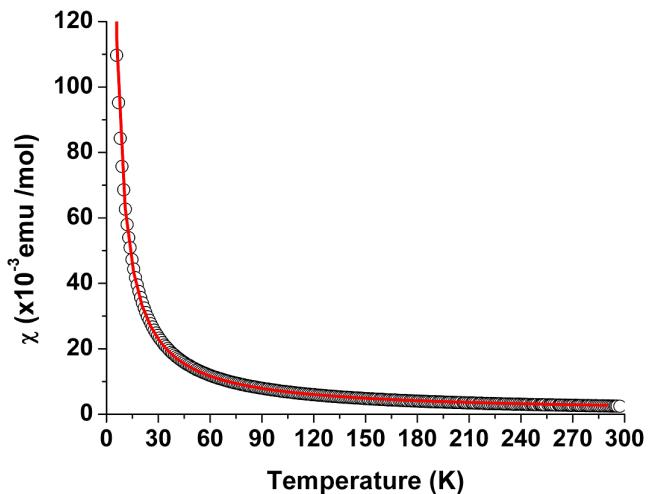


Figure S5. Molar magnetic susceptibility of **3** as a function of temperature. The red line is a fit according to the Curie–Weiss model.

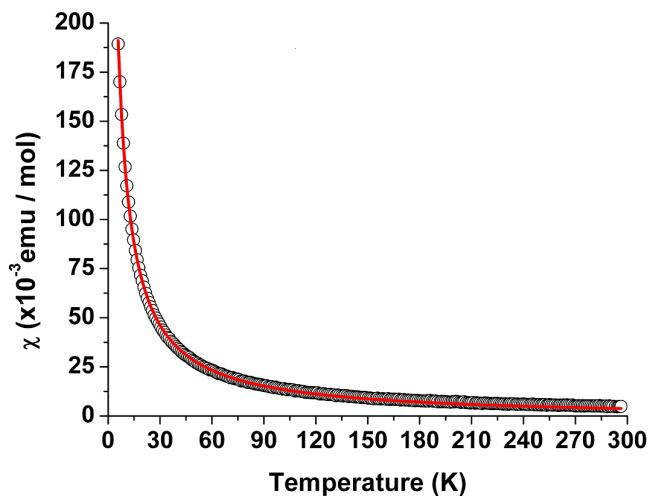


Figure S6. Molar magnetic susceptibility of **6** as a function of temperature. The red line is a fit according to the Curie–Weiss model.

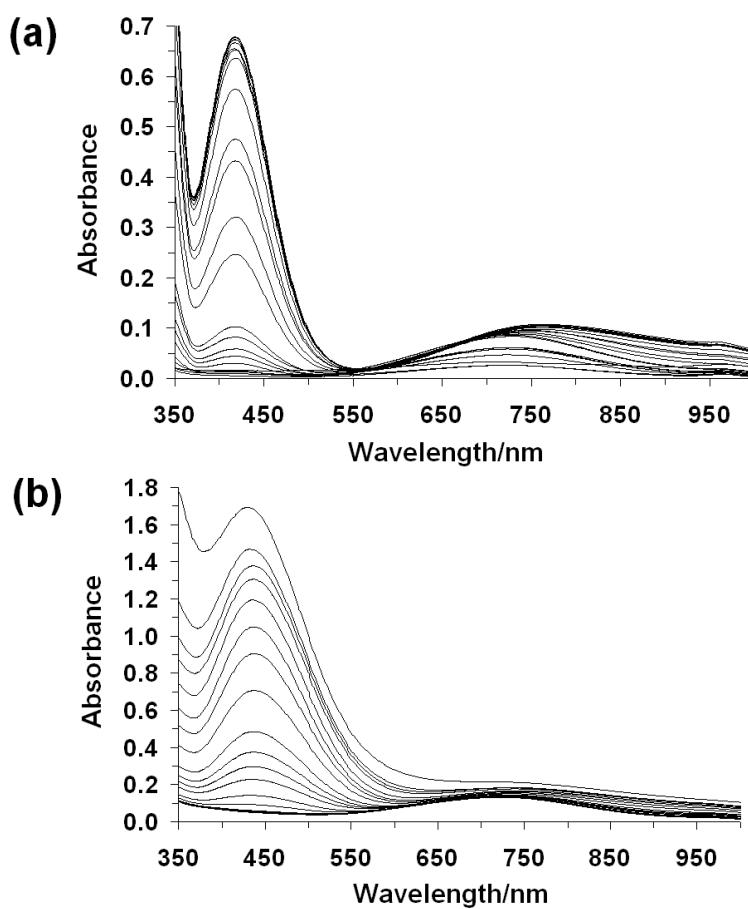


Figure S7. UV-vis absorption spectra of **1** (1.7 mM, a) and **2** (2.5 mM, b) observed over the pH range 2.0 – 10.0 in 0.1M KCl at 25 °C.

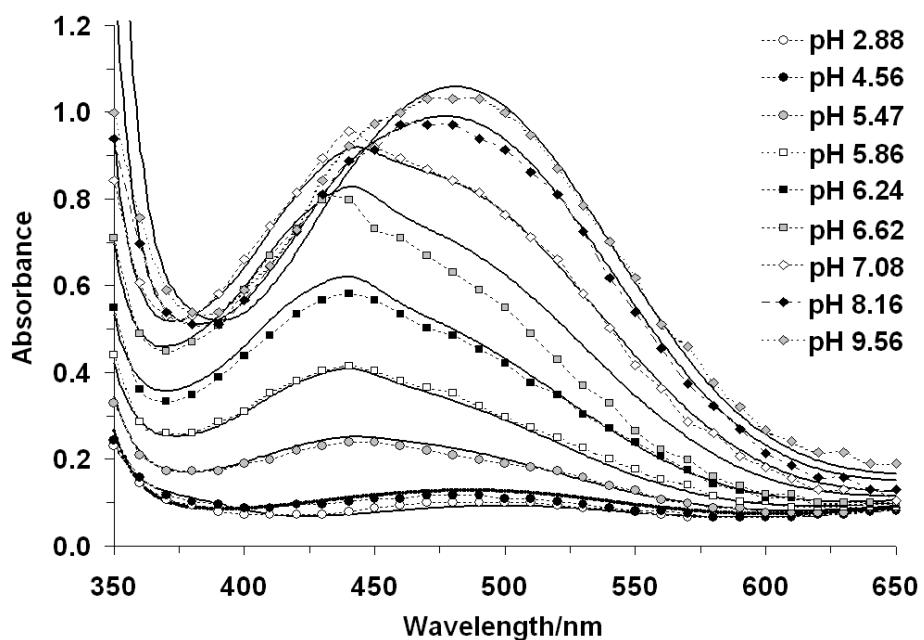


Figure S8. Absorbance spectra of **3** (1.0 mM) over the pH range 3.0 – 10.0. Straight lines refer to the experimental data obtained while points denoted represent the simulated spectra produced from program SQUAD at each pH value studied.

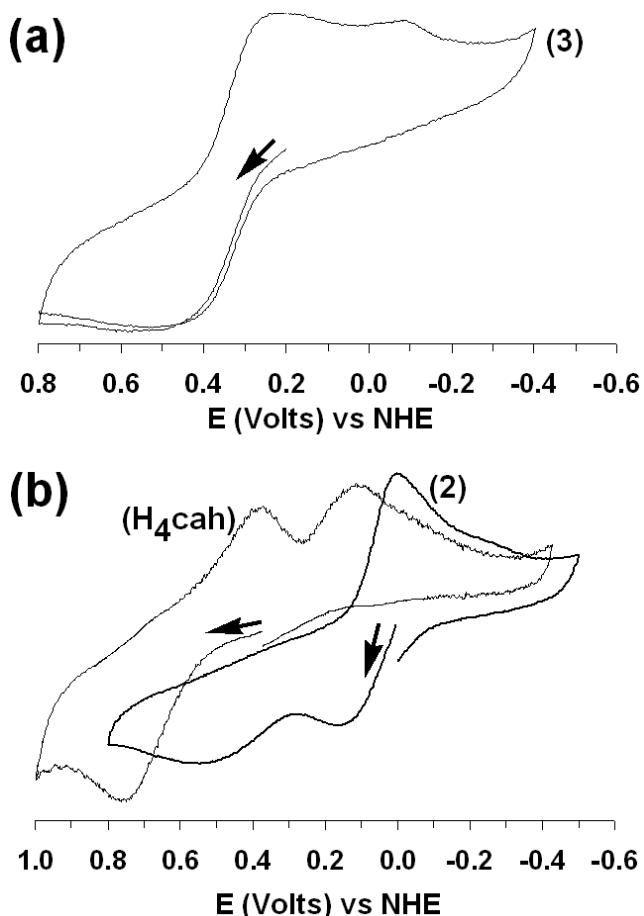


Figure S9. Cyclic voltamograms of (a) **3** and (b) **2** with H_4cah at pH 8.0 (0.1M KNO_3 was used as supporting electrolyte, N_2 , 25 °C). The scan rate was 100 mV s⁻¹. A platinum disk electrode was used as the working electrode, a platinum wire as the auxiliary electrode, and an Hg/HgSO_4 electrode as reference. Arrows indicate the direction of initial potential. The scale of the potential axis has been shifted to match the NHE reference.