Supporting information

Two water-bridged cobalt(II) chains with isomeric naphthoate spacers: from metamagnetic to single-chain magnetic behaviour

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Fig. S1 FT-IR spectra of 1 and 2.

D–Н…А	<i>d</i> (D–H)	d (H···A)	d (D···A)	∠(D–H…A)
O5−H5A…O4	0.850	1.802	2.637	166.95
O5–H5B…O1 ^{#1}	0.850	1.799	2.635	167.71
O6−H6B…O2 ^{#1}	0.850	1.847	2.696	176.60
O7–H7B…O3 ^{#1}	0.850	1.849	2.697	175.57
O8–H8B…O4	0.850	1.898	2.707	158.56
O7−H7A…O8 ^{#1}	0.850	1.865	2.707	170.65
O9–H9A…O1	0.850	1.880	2.694	159.92
O6−H6A…O9 ^{#1}	0.850	1.857	2.704	173.83
^{<i>a</i>} Symmetry codes: ^{#1} $3/2 - x$, $y - 1/2$, $1/2 - z$.				

Table S1 Hydrogen-bonding parameters (Å, °) for 2^{a}



Fig. S2 TGA of 1.



Fig. S3 Simulated (purple) and experimental (blue) PXRD patterns for 1 and 2.



Fig. S4 2D network of 2 generated by hydrogen-bonding (green broken lines) and C-H $\cdots \pi$ interactions (white broken lines).



Fig. S5 Temperature dependence of χ_M^{-1} for 1 and 2 (Solid lines correspond to the Curie-Weiss fit).



Fig. S6 Temperature dependence of ac susceptibilities for 1 at different frequencies (10 and 1000 Hz) under $H_{ac} = 3.5$ Oe and $H_{dc} = 0$ Oe.



Fig. S7 (a) Field-dependent magnetizations measured at different temperatures for 1 (insert: *d*M/*d*H curves).(b) Temperature-dependent molar susceptibility for 1 collected at different fields.



Fig. S8 Temperature-dependent heat capacity for 1 and 2 measured at different dc fields.



Fig. S9 Field-dependent magnetizations of 2 measured at different temperatures.



Fig. S10 Cole–Cole diagram of 2 at 1.4 K (The solid line represents the least-squares fit obtained with a Debye model).