Electronic Supporting Information (ESI)

Chiral 2p-3d heterometallic azido complex with 2,6-pyridinedicarboxylate as co-ligand showing magnetic order

Jiong-Peng Zhao, Qian Yang, Song-De Han, Jie Han, Ran Zhao, Bo-Wen Hu, and Xian-He Bu*

Table S1. Selected bond lengths [Å] for complex 1.

Co1—O1 ⁱ	1.981 (3)	
Co1—N1	2.045 (3)	
Co1—O4 ⁱⁱ	2.095 (3)	
Co1—N2	2.148 (4)	
Co1—O3	2.176 (2)	
Co1—O2	2.221 (3)	
Na1—O2 ⁱ	2.285 (2)	
Na1—O2 ^v	2.285 (2)	
Na1—O5	2.361 (8)	
Na1—O5 ⁱⁱⁱ	2.361 (8)	
Na1—O3 ⁱⁱⁱ	2.526 (3)	
Symmetry codes: (i) $x+1/2$, $-y+3/2$, $-z+7/4$; (ii) $-y+3/2$, $x-1/2$, $z+1/4$; (iii) $-y+2$, $-x+2$, $-z+3/2$; (iv) $x-1/2$, $-y+3/2$, $-z+7/4$; (v) $y+1/2$, $-x+3/2$, $z-1/4$.		

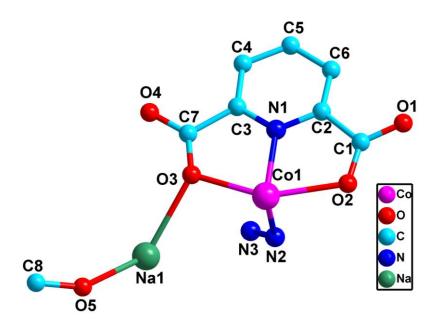


Figure S1. The asymmetry unit of complex **1**.

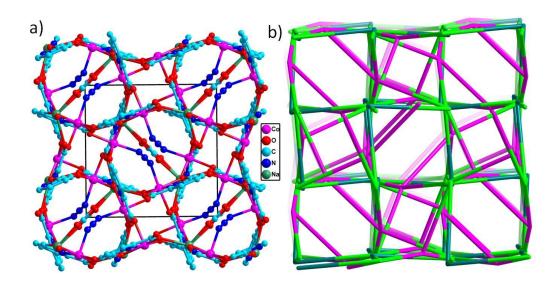


Figure. S2 a) 3D view of the complex along c direction in 1; b) 4,5-connected 3-nodal topology net in 1.

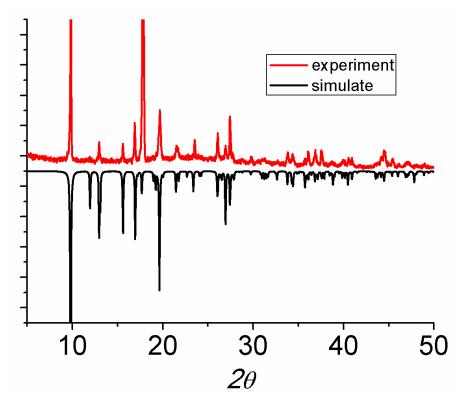


Figure S3. The XRPD diagram for complex **1**.

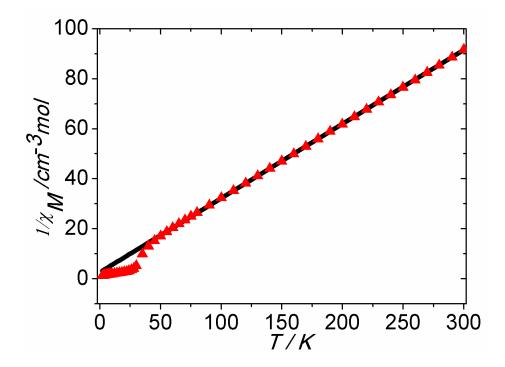


Figure S4. Curie plot for 1. The solid line is the best fit to the Curie-Weiss law.

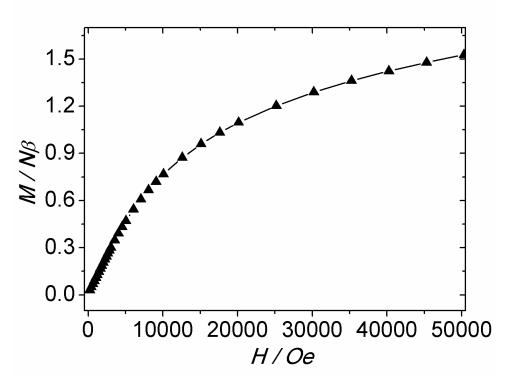


Figure S5. The magnetization curve at 2.0 K of 1.