Supporting information

Self-Assembly of Discrete and Polymeric Metal-Organic Frameworks via a Semirigid Glycoluril-Based Molecular Clip

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Figure S1. Packing diagram of **1** along the crystallographic *bc* axis. The oxygen atoms of water molecules are shown in space-filling model (red).





(b)



(c)









Figure S2. (a) ORTEP view of **3** showing four ligands surrounding one Co(II) octahedra (30% probability ellipsoids). (b) View of **3** showing the 1D wavelike chain (left) and individual rhombuslike unit (right). (c) The 2D sheet driven by hydrogen bond interactions and its 1D channels. (d) The 2D hydrogen bond network in the framework **3**. Metal atoms and ligands **L** have been omitted for clarity. (e) Space-filling model for topview of **3**. Hydrogen atoms and free water molecules have been omitted for clarity.



Figure S4. TGA curve of 2.







Figure S6. TGA curve of 4.

Photoluminescence



Figure S7. Fluorescent emission spectrum of complex L in solid state at room temperature.



Figure S8. Fluorescent emission spectrum of complex 3 in solid state at room temperature.



Figure S9. Fluorescent emission spectrum of complex 4 in solid state at room temperature.



(c)

Figure S10. X-ray power diffraction diagram of (a) compound **1**, (b) **1** after removing the free water molecules and (c) the simulated spectra from single crystal data.



Figure S11. X-ray power diffraction diagram of (a) compound **2**, (b) **2** after removing the free water molecules and (c) the simulated spectra from single crystal data.



Figure S12. X-ray power diffraction diagram of (a) compound **3**, (b) **3** after removing the free water molecules and (c) the simulated spectra from single crystal data.



Figure S13. X-ray power diffraction diagram of (a) compound **4**, (b) **4** after removing the free DMF molecules and (c) the simulated spectra from single crystal data.