Supporting information

Several (4,4)- and (5,6,8)-Connected Lanthanide-Organic Frameworks: Structures, Luminescence and Magnetic Properties

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	2	3	4	5	6
Ln1-O _c	2.350	2.352	2.336	2.317	2.309
Ln1-O _m	2.693	2.671	2.667	2.680	2.668
Ln1-O _w	2.436	2.439	2.430	2.431	2.414
Ln2-O _c	2.377	2.379	2.368	2.348	2.345
Ln2-O _m	2.678	2.665	2.647	2.643	2.636
Ln3-O _c	2.362	2.365	2.352	2.322	2.326
Ln3-O _m	2.518	2.517	2.509	2.508	2.500
Ln3-O _w	2.444	2.445	2.415	2.396	2.389

Table S1. Average bond lengths associated with Ln atoms (Å) of 2-6.

 O_c , oxygen atom of the carboxyl group; O_m , oxygen atom of methyl group; O_w , oxygen atom of the coordinated water molecule.



Fig. S1 The 3D framework of 2 highlighting benzene rings.



Fig. S2 Theoretical and experimental powder X-ray diffraction patterns for 2-7.



Fig. S3 The TGA diagrams of 1, 5, and 7.



Fig. S4 The plot of $1/\chi_M$ versus *T* of compound **4**, and the red line represents the best fitting for Curie-Weiss law $\chi = C/(T-\theta)$