Auxiliary ligand-directed and counter anion-templated effects on coordination networks based on semirigid 2-aminodiacetic terephthalic acid ligand

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Fig. S1. Power X-ray diffraction patterns for complexes 1-4.



Fig. S2. TG curves for complexes 1-4.



Fig. S3. Hydrogen bonds motif in **1.** symmetry codes, A: -*x*, -*y*, 2-*z*; B: *x*, *y*, 1+z. O9…O9A distance is 2.8894(61) Å and O8B…O10A is 2.6866(60) Å.



Fig. S4. Luminescent excitation spectra for complexes 1, 3, 4 and H_4 adtp.



Fig. S5. Luminescent excitation spectra for 2a at room temperature (left) and 10 K (right).