< Electronic Supplementary Information>

Supramolecular isomerism between cyclodimeric and sinusoidal 1D coodination polymers: competition of tunable argentophilic vs electrostatic interactions

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Electronic Supplementary Information (ESI) available: ¹H NMR spectra and IR spectra of ligand, $[AgL]_2(NO_3)_2$, $[AgL]_2(BF_4)_2$, $[AgL]_2(ClO_4)_2$, $[AgL]_2(PF_6)_2 \cdot 5C_4H_8O_2$, $[AgL](BF_4)$, $[AgL](ClO_4)$ and $[AgL](PF_6)$. TGA-DSC curves of $[AgL]_2(NO_3)_2$, $[AgL]_2(BF_4)_2$, $[AgL]_2(ClO_4)_2$, $[AgL]_2(PF_6)_2 \cdot 5C_4H_8O_2$, $[AgL](BF_4)$, $[AgL](ClO_4)$ and $[AgL](PF_6)$. Crystal structures of $[AgL]_2(NO_3)_2$, $[AgL]_2(BF_4)_2$, $[AgL]_2(BF_4)_2$, $[AgL]_2(PF_6)_2 \cdot 5C_4H_8O_2$. Solid-state PL spectra of $[AgL](ClO_4)_2$ and $[AgL](ClO_4)_2$ and $[AgL](ClO_4)_2$.



Fig. S1 ¹H NMR spectra for L (a), $[AgL]_2(NO_3)_2$ (b), $[AgL]_2(BF_4)_2$ (c), $[AgL]_2(ClO_4)_2$ (d), $[AgL]_2(PF_6)_2 \cdot 5C_4H_8O_2$ (e), $[AgL](BF_4)$ (f), $[AgL](ClO_4)$ (g), and $[AgL](PF_6)$ (h) in Me₂SO-*d*₆.



Fig. S2 ¹H NMR spectra for L (a), $[AgL]_2(ClO_4)_2$ (b), $[AgL](ClO_4)$ (c) in acetone- d_6 .



Fig. S3 IR spectra for L (a), $[AgL]_2(NO_3)_2$ (b), $[AgL]_2(BF_4)_2$ (c), $[AgL]_2(CIO_4)_2$ (d), $[AgL]_2(PF_6)_2 \cdot 5C_4H_8O_2$ (e), $[AgL](BF_4)$ (f), $[AgL](CIO_4)$ (g), and $[AgL](PF_6)$ (h).



Fig. S4 TGA and DSC curves for $[AgL]_2(NO_3)_2$ (a), $[AgL]_2(BF_4)_2$ (b), $[AgL]_2(ClO_4)_2$ (c), $[AgL]_2(PF_6)_2 \cdot 5C_4H_8O_2$ (d), $[AgL](BF_4)$ (e), $[AgL](ClO_4)$ (f) and $[AgL](PF_6)$ (g).



Fig. S5 Crystal structures (side views) of [AgL]₂(NO₃)₂ (a), [AgL]₂(BF₄)₂ (b), [AgL]₂(ClO₄)₂ (c), [AgL]₂(PF₆)₂·5C₄H₈O₂ (d).



Fig. S6 Solid-state PL spectra for $[AgL]_2(ClO_4)_2$ (top, a) and $[AgL](ClO_4)$ (bottom, b) upon additions of a drop of solvent at $\lambda_{ex} = 321$ nm.