Electronic Supplementary Information (ESI) for CrystEngComm

Structural diversity of Ag/3-nitrophthalate coordination polymers controlled by solvent and induction agent

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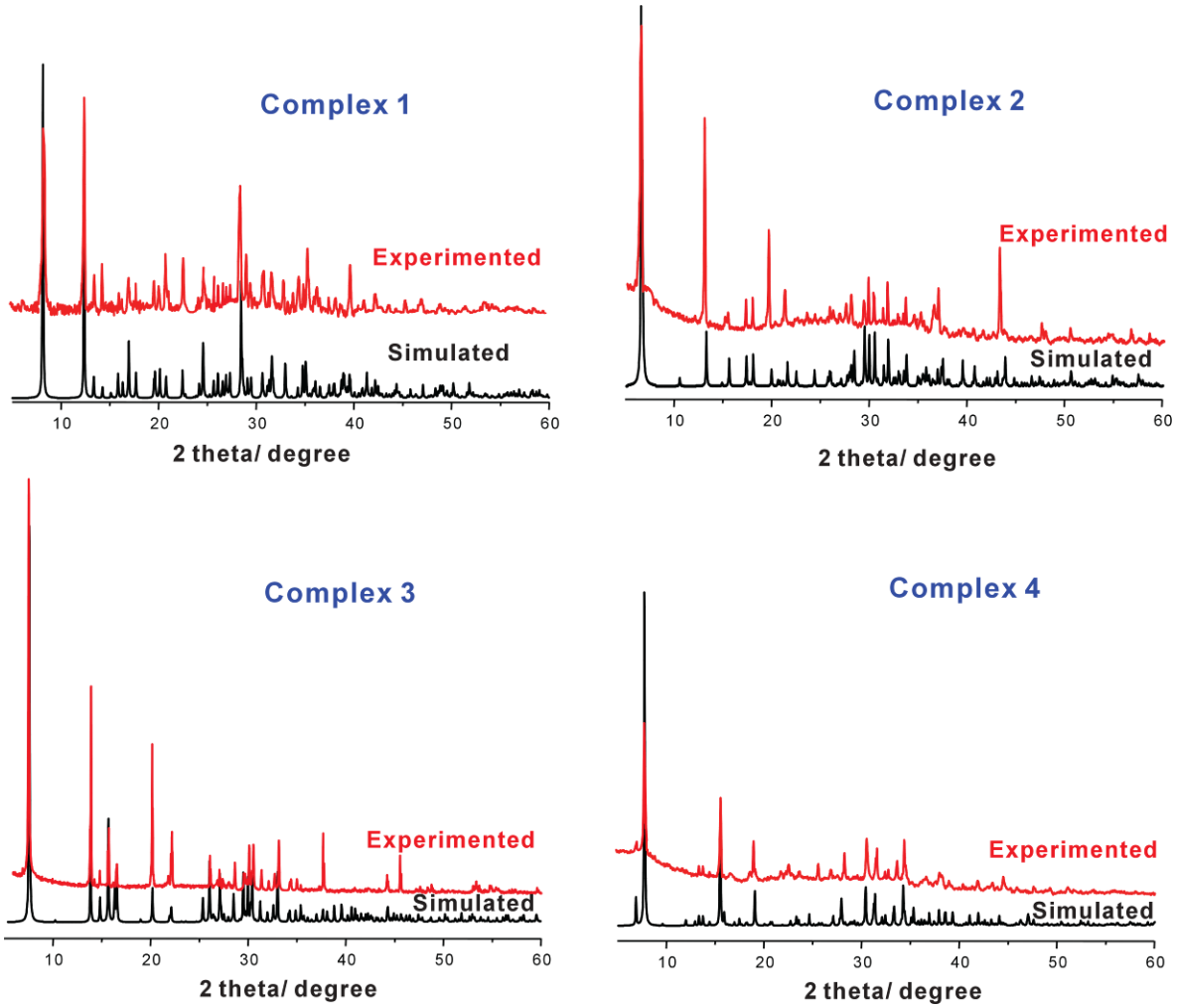
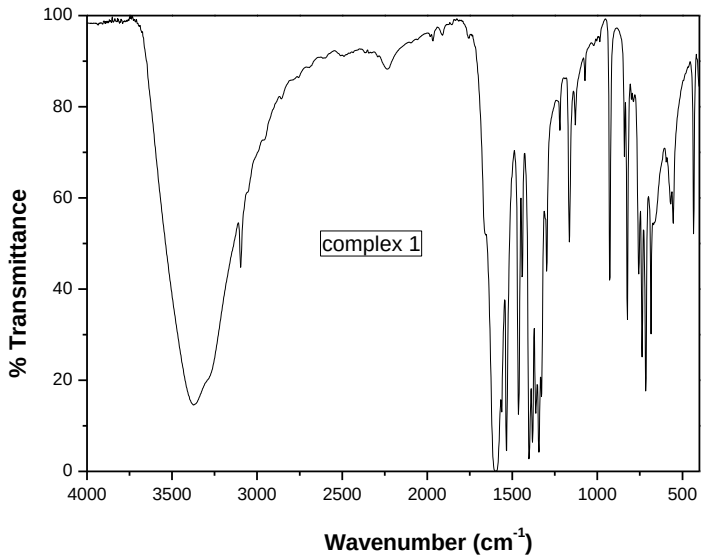
(1) Table S1 The hydrogen bond geometries for **1-4**.

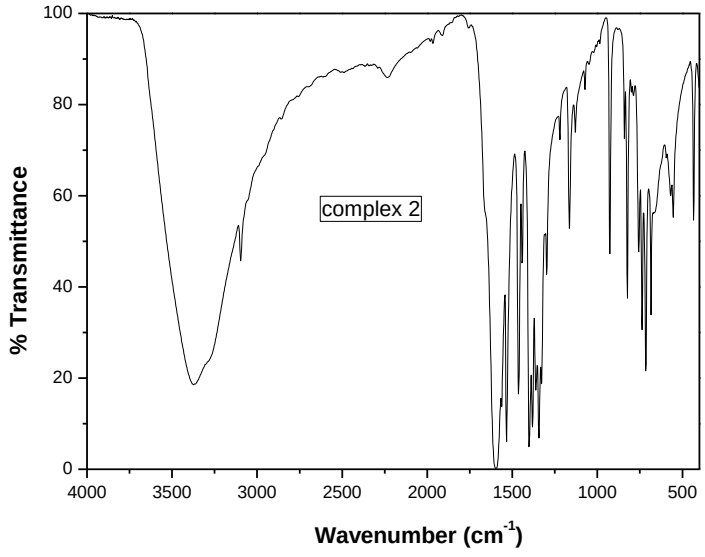
Complex **1**					
D—H···A	D—H	H···*A*	*D*···*A*		*D*—H···*A*
N2—H2B···O4^{iv}	0.91	2.20	3.077	(5)	160.9
N2—H2C···O3	0.91	2.30	3.013	(5)	135.5
N2—H2D···O4^{i}	0.91	2.14	2.977	(5)	152.3
N3—H3C···O4^{v}	0.91	2.17	3.007	(5)	152.2
N3—H3A···O1^{ii}	0.91	2.04	2.947	(5)	173.7
N3—H3B···O6^{vi}	0.91	2.39	3.079	(5)	132.6
Symmetry codes: (i) −*x*+1, −*y*+1, −*z*+2; (ii) *x*, *y*, *z*−1; (iv) *x*−1, *y*, *z*; (v) *x*−1, *y*, *z*−1; (vi) *x*−1/2, −*y*+3/2, *z*−1/2.					
Complex **2**					
N2—H2A···O6^{ii}	0.90	2.06	2.928	(4)	161.2
N2—H2B···O1^{vi}	0.90	2.42	3.314	(5)	174.2
N2—H2C···O2^{iii}	0.90	2.44	3.301	(5)	160.4
Symmetry codes: (ii) *x*, *y*−1, *z*; (iii) *x*, −*y*+3/2, *z*+1/2; (vi) *x*, −*y*+3/2, *z*−1/2.					
Complex **3**					
N2—H2C···O3	0.85	2.09	2.875	(5)	153.2
N2—H2E···O4^{iv}	0.85	1.92	2.764	(5)	173.3
N2—H2D···O3^{ii}	0.85	2.27	3.037	(5)	149.9
N2—H2B···O1W	0.85	2.01	2.851	(5)	173.1
O1W—H1WB···O2^{v}	0.85	2.05	2.827	(5)	151.3
O1W—H1WA···O2	0.85	2.16	2.774	(5)	128.8
Symmetry codes: (ii) −*x*+1, *y*+1/2, −*z*+1/2; (iv) *x*, *y*+1, *z*; (v) −*x*+1/2, *y*+1/2, *z*.					
Complex **4**					
N3—H3B···O3^{iv}	0.91	2.10	3.005	(7)	172.1
N3—H3C···O7^{i}	0.91	2.22	3.085	(6)	158.0
N3—H3D···O11^{vi}	0.91	2.48	3.017	(7)	117.7
N4—H4B···O5^{ii}	0.91	2.26	3.166	(7)	177.9
N4—H4D···O10	0.91	2.47	3.314	(7)	153.5
N4—H4C···O1W^{vii}	0.91	2.14	2.993	(7)	155.3
N5—H5A···O9^{iv}	0.91	2.04	2.879	(6)	152.1
N5—H5B···O1W^{ii}	0.91	2.13	2.973	(7)	153.7
N5—H5C···O10	0.91	2.06	2.848	(6)	143.6
N6—H6A···O2^{i}	0.91	2.21	3.037	(7)	151.2
N6—H6B···O9^{iv}	0.91	2.43	3.072	(7)	127.7
N6—H6C···O8	0.91	2.40	3.054	(7)	128.7
O1W—H1WA···O3	0.85	1.93	2.760	(6)	164.9
O1W—H1WB···O1^{vii}	0.85	2.02	2.848	(6)	164.5
Symmetry codes: (i) −*x*+1, −*y*+1, −*z*+1; (ii) −*x*+1, −*y*+1, −*z*+2; (iv) *x*+1, *y*, *z*; (vi) −*x*+1, *y*−1/2, −*z*+3/2; (vii) −*x*, −*y*+1, −*z*+2.					

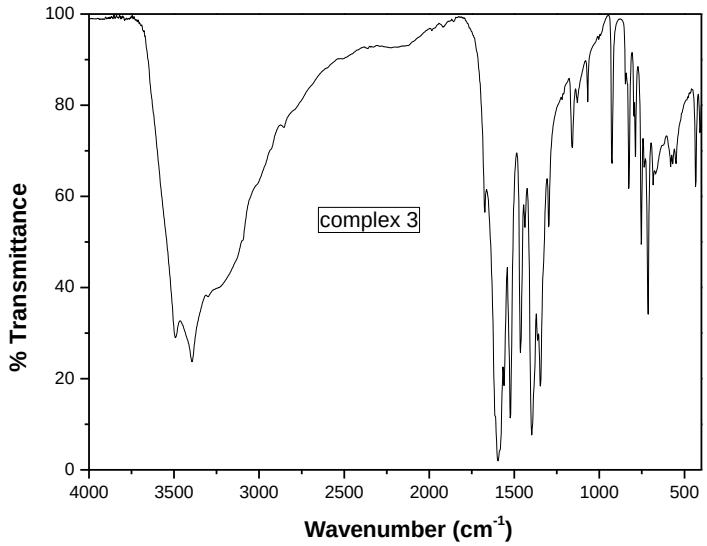
(2) Fig. S1 The powder XRD patterns and the simulated one from the single-crystal diffraction data for complex 1-4

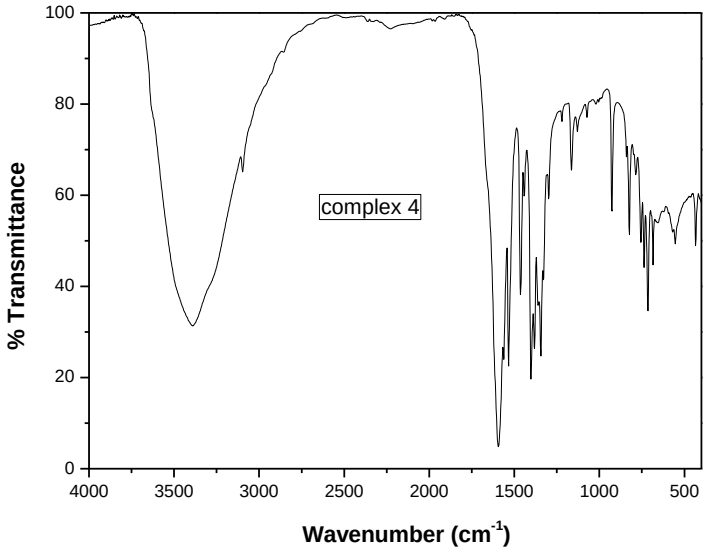


(3) Fig. S2 IR of complexes 1-4

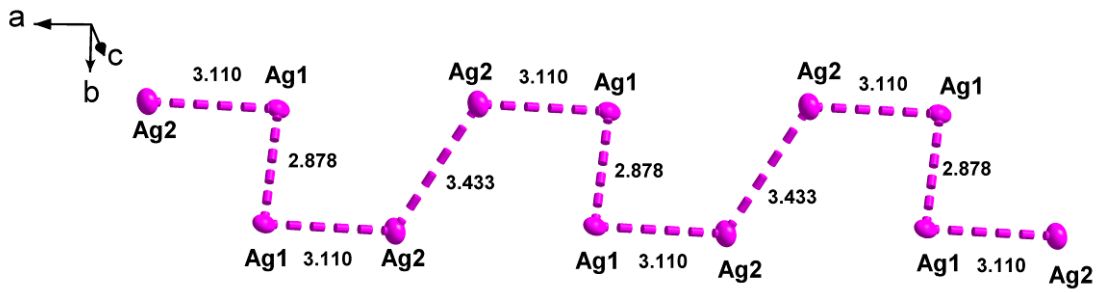




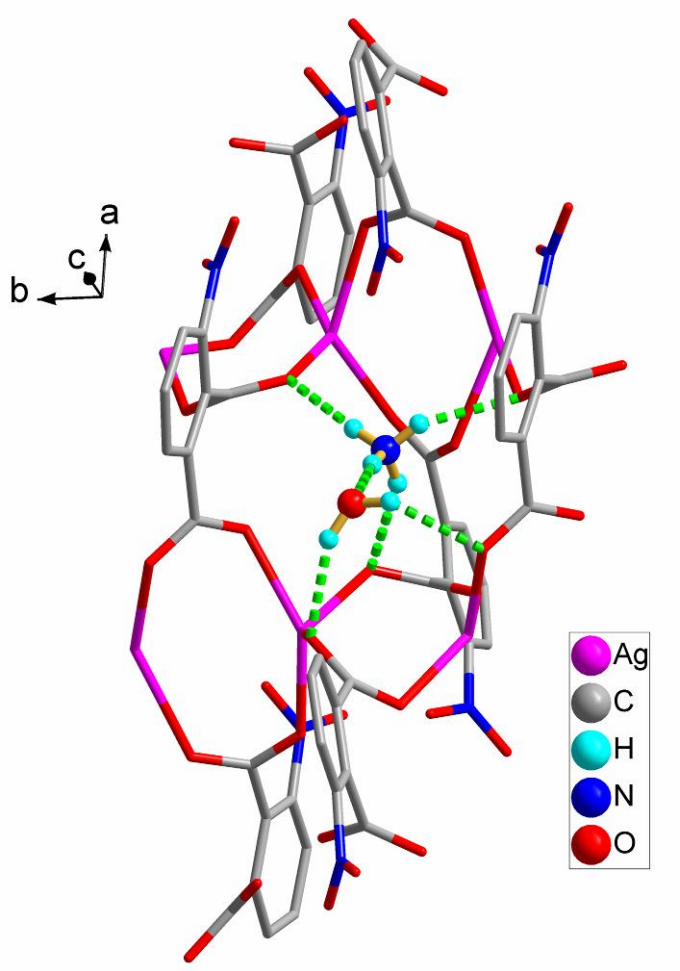




(4) Fig. S3: The silver wire in 1.



(5) Fig. S4: The $[(NH_4)(H_2O)]^+$ cluster in the lacuna of the 2D sheet



(6) Fig. S5: TGA curves for CPs 1-4.