Syntheses, crystal structures and photoluminescent properties of two novel Ag(I) coordination polymers with benzoguanamine and pyrazine-carboxylate ligands: From 1D helix to 1D→2D interdigitation

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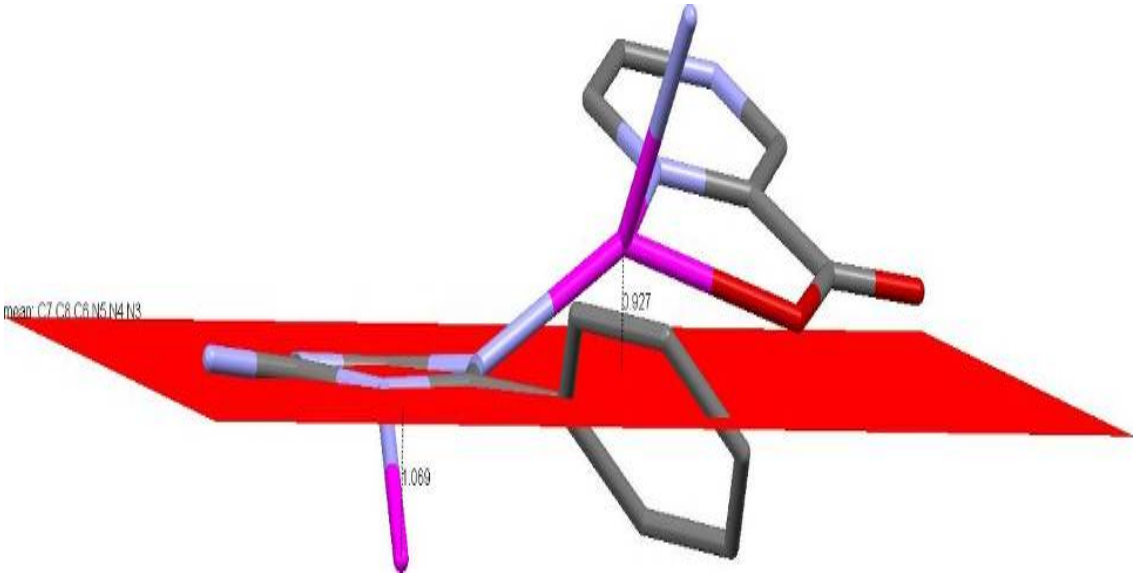
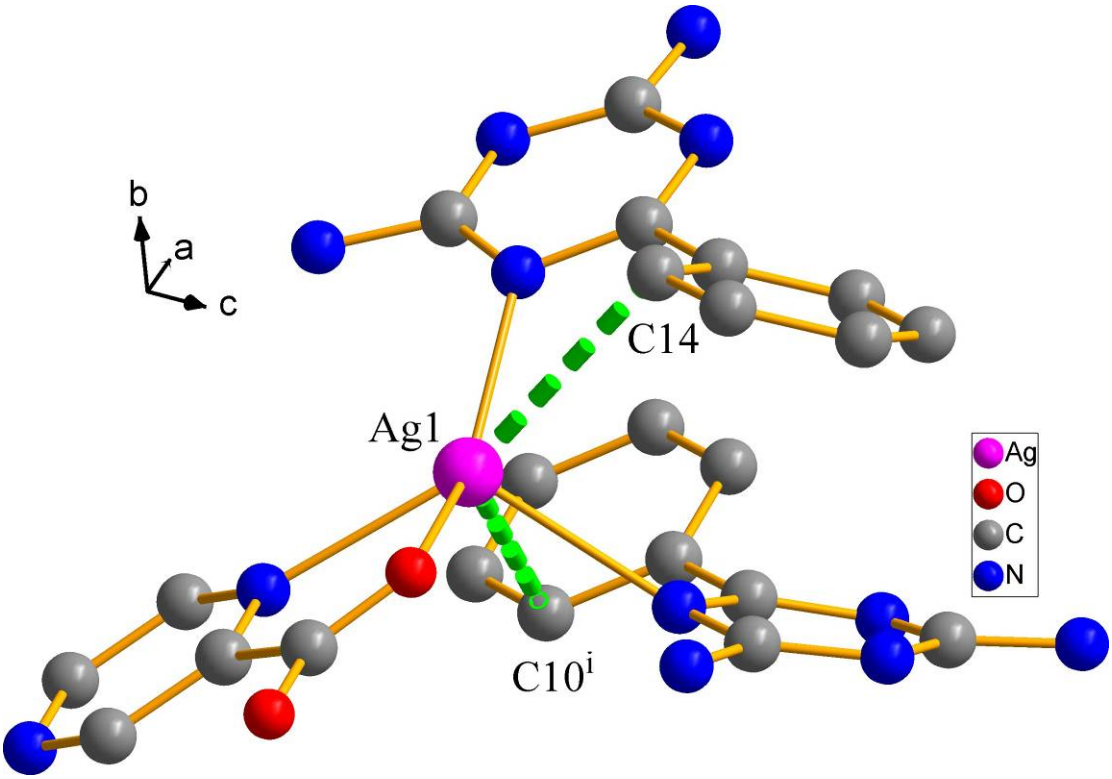
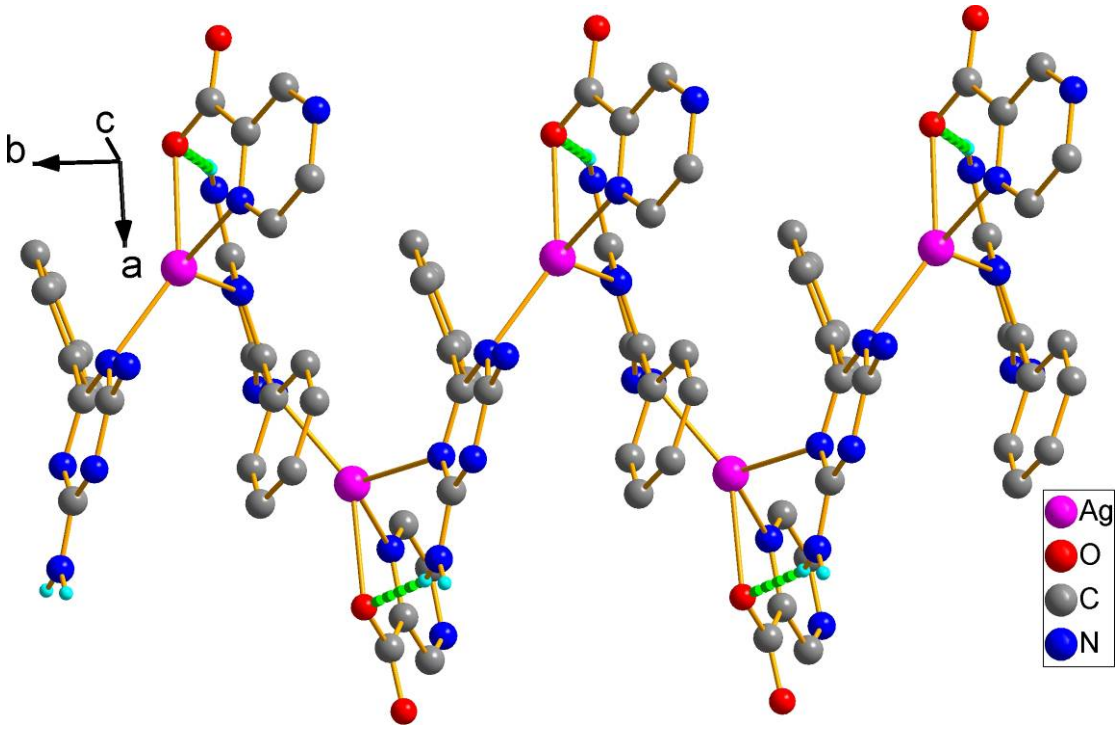
(1) Fig. S1: Relative positions of Ag1 respect to the triazinyl ring in 1.



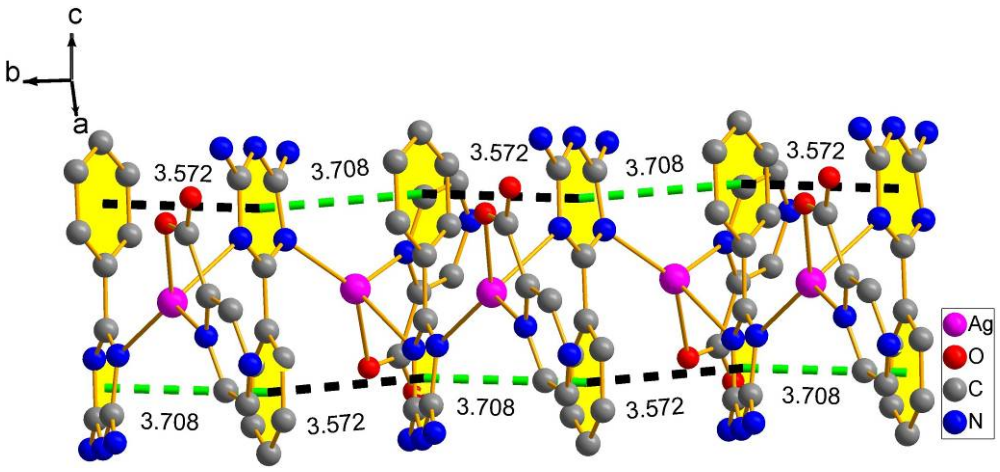
(2) Fig. S2: The Ag···C interactions in 1.



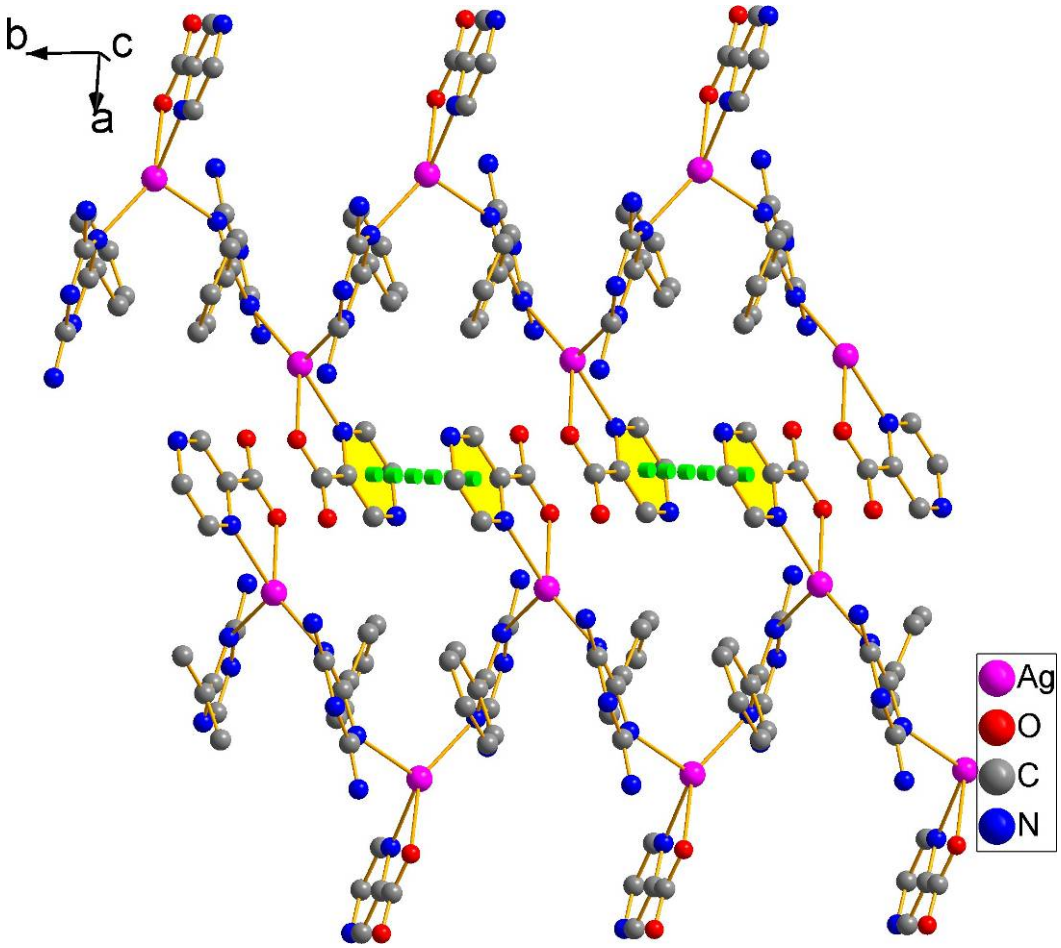
(3) Fig. S3: Intrachain N-H···O hydrogen bonds.



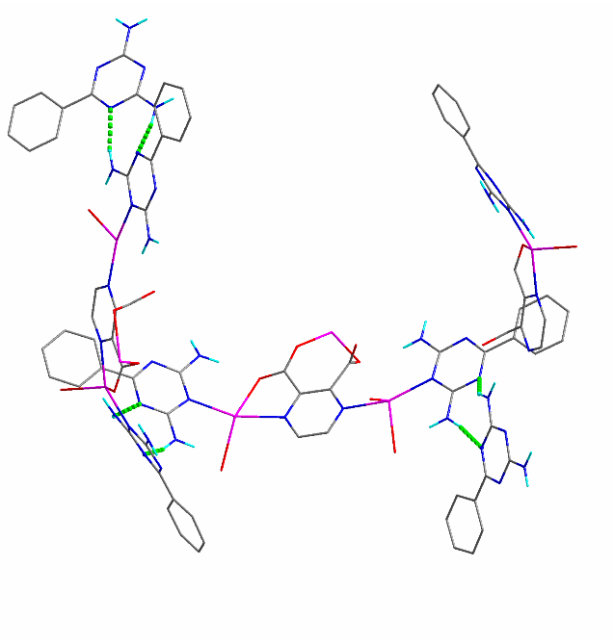
(4) Fig. S4: The face-to-face π···π stacking between adjacent bga ligand in 1.



(5) Fig. S5: The π···π stacking between adjacent pzc ligand in 1.



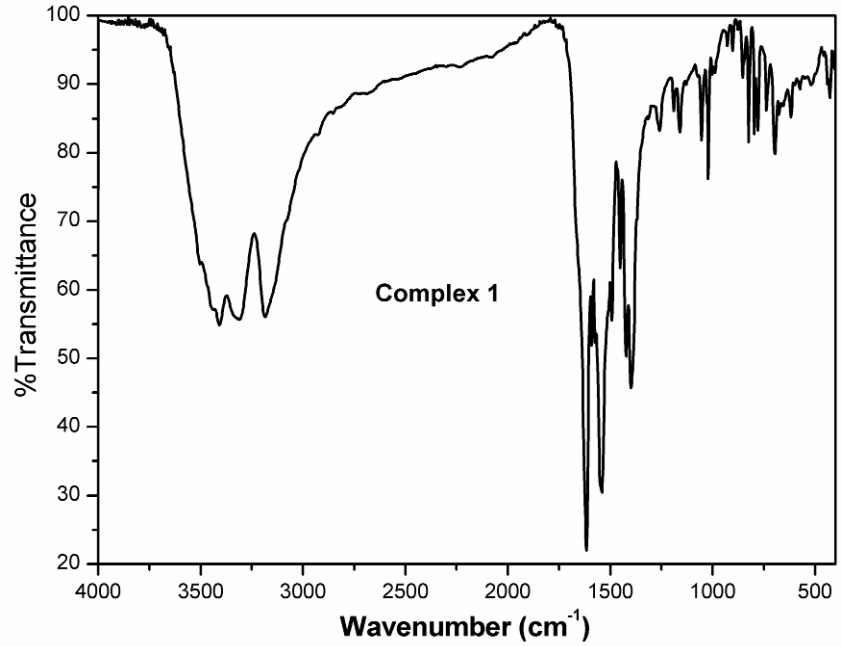
(6) Fig. S6: The complementary N-H···N hydrogen bonds in 2.

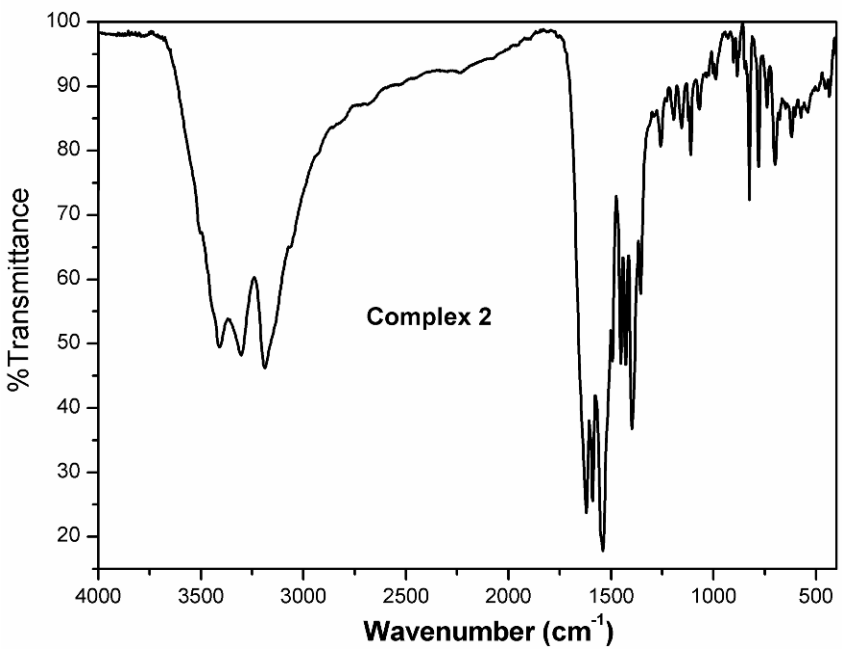


(7) Table S1: The hydrogen bond geometries for 1 and 2

Complex **1**				
D—H···*A*	*D*—H	H···*A*	*D*···*A*	*D*—H···*A*
O1W—H1WA···O2	0.85	2.18	3.031(3)	175
N6—H6A···N3^{ii}	0.88	2.24	3.046(3)	153
N6—H6B···O1^{iii}	0.88	2.17	2.812(3)	129
N7—H7B···O1^{iv}	0.88	2.07	2.940(3)	172
N7—H7C···O2^{v}	0.88	2.05	2.898(3)	160
Symmetry codes: (ii) −*x*+3/2, −*y*+3/2, −*z*+1; (iii) −*x*+1, −*y*+1, −*z*+1; (iv) *x*+1/2, *y*+1/2, *z*; (v) −*x*+3/2, *y*+1/2, −*z*+3/2.				
Complex **2**				
D—H···*A*	*D*—H	H···*A*	*D*···*A*	*D*—H···*A*
N4—H4B···O1W	0.88	2.21	2.943(10)	141
N4—H4A···N2^{ii}	0.88	2.34	3.086(10)	143
N5—H5B···O1	0.88	2.03	2.909(10)	176
N9—H9B···O3^{i}	0.88	1.94	2.817(9)	171
N9—H9C···N6^{iii}	0.88	2.19	3.034(10)	161
N10—H10B···O4	0.88	2.15	2.924(10)	146
N10—H10C···N7^{iv}	0.88	2.23	3.033(10)	153
O1W—H1WC···O2^{v}	0.85	2.07	2.767(8)	139
O1W—H1WA···O4^{vi}	0.85	1.95	2.753(8)	157
Symmetry codes: (i) *x*, −*y*+2, *z*+1/2; (ii) −*x*+1, *y*, −*z*+3/2; (iii) *x*, −*y*+3, *z*+1/2; (iv) *x*, −*y*+3, *z*−1/2; (v) *x*, −*y*+1, *z*+1/2; (vi) *x*, *y*−1, *z*.				

(8) Fig. S7: IR spectra of 1 and 2





(9) Fig. S8: Emission spectrum of free bga ligand