Supporting Information.

Table S1. Selected hydrogen bond lengths (Å) and bond angles (°) for **1-7**

D-H···A	d(D-H)	d(H···A)	d(D···A)	∠DHA	Symm code
Complex 1					
O5-H5···N2	0.820	1.835	2.650	172.33	[x-1/2, y-1/2, z]
Complex 2					
O6-H6C···O2	0.847	1.890	2.643	147.35	[-x+5/4, y, -z+5/4]
O7-H7A···O7	0.851	1.492	2.263	148.59	[x, -y+5/4, -z+5/4]
O7-H7A···O2	0.851	2.095	2.643	121.68	[x, -y+5/4, -z+5/4]
O7-H7B···O2	0.851	1.891	2.564	135.00	
O7-H7B···O3	0.851	1.931	2.690	147.94	[-x+2, -y+1, -z+1]
Complex 3					
O5-H5A···N2	0.874	2.091	2.907	155.05	[-x+1, -y, -z]
Complex 4					
O6-H6A···O9	0.843	1.954	2.755	158.23	[-x+1, -y+1, -z+1]
O6-H6B···O4	0.848	1.859	2.697	168.96	[x-1, y-1, z]
O6-H6B···O5	0.848	2.568	2.984	111.42	[x-1, y-1, z]
O7-H7A···O2	0.846	1.956	2.724	150.39	[-x+1, -y+1, -z]
O7-H7B···O2	0.843	1.868	2.669	158.07	
O8-H8A···O4	0.845	1.804	2.640	169.96	[-x+1, -y+2, -z+1]
O8-H8B···O1	0.846	1.904	2.745	172.86	
O9-H9A···O5	0.844	2.131	2.927	157.33	[x-1, y-1, z]
O9-H9B···O7	0.845	2.140	2.953	161.09	
Complex 5					
O6-H6A···O3	0.859	1.854	2.667	157.48	[x, y, z-1]
O6-H6B···O2	0.848	2.056	2.818	149.22	[-x+1, -x+y+1, -z+1/3]
Complex 6					
O6-H6A···O1	0.841	1.868	2.600	144.57	
O7-H7B···O7	0.935	1.309	2.214	160.88	[x, -y+5/4, -z+5/4]
Complex 7					
O11-H11C···O1	0.845	1.871	2.702	167.68	[-x+1, -y+1, -z+2]
O11-H11D···O7	0.953	1.704	2.655	175.14	[-x+2, -y+1, -z+2]