Electronic supplementary information (ESI)

Structural Modulation of Silver Complexes and Their Distinctive

Catalytic Properties

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D-H ·· A ^b	Distance $(D \cdot \cdot A^b)$	D-H-A ^b	Angle (D-H-A $^{\rm b}$)
1			
C5-H5B ···F4#1	3.444	C5-H5B-F4	152
C6-H6B ····O7#2	3.260	C6-H6B-O7	140
C8-H8B · · O5#3	3.031	C17-H17A-O8	100
C23-H23B ··· O6#4	2.980	C23-H23B-O6	98
C25-H25A ··· O8#5	3.142	C25-H25A-O8	139
	2		
C2-H2A ···F2#6	2.614(4)	C2-H2A-F2	124
	3		
C19-H7 ··· O4#7	3.170(12)	C1-H1-F6	131
С19-Н7 ·· О3#8	3.411(10)	C4-H4-F1	148
C29-H15 ·· O2#9	2.668(3)	C29-H15-O2	151
	1 0 11 1/0	1/0 //0 1 0	10 1/0 1/0

Table S1 Distances [Å] and angles [] of hydrogen bonding for complexes 1 - 3^a

^a Symmetry codes for **1-3**: #1: -x, 1/2+y, -1/2-z; #2: -1+x, 3/2-y, 1/2+z; #3: x, 3/2-y, -1/2+z; #4: -1+x, 3/2-y, 1/2+z; #5: -x, 2-y, -z; #6: 2-x, -y, 1-z; #7: -1/2+x, 1/2+y, z; #8: 1/2-x, 3/2-y, 1-z; #9: 1/2-x, 1/2-y, 1-z. ^b D: donor; A: acceptor

NMR spectra



Figure S2



Figure S4







Figure S6







Figure S8







Figure S10



Figure S11



Figure S12



Figure S13



Figure S14





Figure S15 (a) 1D double chain of 1 with hydrogen bonds indicated by dashed lines.



Figure S15 (b) 2D network of 1 with hydrogen bonds indicated by dashed lines.



Figure S15 (c) 3D structure of 1 with hydrogen bonds indicated by dashed lines.



Figure S16 3D structure of 2 with hydrogen bonds indicated by dashed lines.



Figure S17 3D structure of 3 with hydrogen bonds indicated by dashed lines, CF_3COO^- are omitted for clarity.