

Supporting Information for

Metal-Organic Coordination Architectures of Bis(1,2,4-triazole) Ligands Bearing Different Spacers: Syntheses, Structures and Luminescent Properties

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Table S1 Selected Bond Distances (Å) and Angles (deg) for Complexes **1~6**

1			
Cd(1) - N(1)	2.345(5)	Cd(1) - N(4)	2.317(5)
Cd(1)-O(1)	2.388(6)		
N(4)-Cd(1)-O(1)	82.06(19)	N(4)-Cd(1)-N(1)	90.12(18)
N(4) ^{#1} -Cd(1)-O(1)	97.94(19)	N(4) ^{#1} -Cd(1)-N(1)	89.88(18)
N(1) ^{#1} -Cd(1)-O(1)	96.5(2)	N(1)-Cd(1)-O(1)	83.5(2)
2			
Hg(1)-N(3)	2.455(12)	Hg(1)-Br(3)	2.4610(17)
Hg(1)-Br(2)	2.4712(16)		
N(3)-Hg(1)-Br(3)	95.5(3)	N(3)-Hg(1)-Br(2)	96.5(3)
Br(3)-Hg(1)-Br(2)	167.96(7)		
3			
Mn(1)-N(3)	2.300(4)	Mn(1)-Cl(1)	2.525(2)
N(3) ^{#1} -Mn(1)-N(3) ^{#2}	95.3(2)	N(3)-Mn(1)-N(3) ^{#2}	84.7(2)
N(3) ^{#2} -Mn(1)-N(3) ^{#3}	180.00	N(3) ^{#1} -Mn(1)-Cl(1)	91.78(12)
N(3)-Mn(1)-Cl(1)	88.22(12)		
4			
Cu-N(1)	2.035(4)	Cu-Cl(1)	2.8104(18)
N(1) ^{#1} -Cu-N(1)	87.3(2)	N(1) ^{#1} -Cu-N(1) ^{#2}	92.7(2)
N(1)-Cu-N(1) ^{#2}	87.3(2)	N(1) ^{#1} -Cu-Cl(1)	89.83(11)
N(1)-Cu-Cl(1)	90.17(11)		
5			
Cd(1)-N(1)	2.344(4)	Cd(1)-N(1) ^{#2}	2.345(4)
Cd(1)-O(2) ^{#2}	2.432(15)	Cd(1)-O(2)	2.441(15)
N(1)-Cd(1)-N(1) ^{#2}	83.0(2)	N(1) ^{#1} -Cd(1)-N(1) ^{#2}	97.0(2)
N(1)-Cd(1)-O(2) ^{#3}	99.7(3)	N(1) ^{#1} -Cd(1)-O(2) ^{#3}	80.3(3)
N(1) ^{#2} -Cd(1)-O(2) ^{#3}	79.4(3)	N(1) ^{#3} -Cd(1)-O(2) ^{#3}	100.6(3)
N(1)-Cd(1)-O(2)	101.1(3)	N(1) ^{#1} -Cd(1)-O(2)	78.9(3)
N(1) ^{#2} -Cd(1)-O(2)	79.9(3)	N(1) ^{#3} -Cd(1)-O(2)	100.1(3)
O(2) ^{#3} -Cd(1)-O(2)	148.5(6)	O(2) ^{#2} -Cd(1)-O(2)	31.5(6)
6			
Cd(1)-N(1)	2.307(5)	Cd(1)-N(7)	2.350(5)
Cd(1)-N(4)	2.405(5)		
N(1) ^{#1} -Cd(1)-N(7) ^{#1}	86.69(17)	N(1) ^{#1} -Cd(1)-N(7)	93.31(17)
N(1) ^{#1} -Cd(1)-N(4)	91.38(17)	N(1)-Cd(1)-N(4)	88.62(17)
N(7) ^{#1} -Cd(1)-N(4)	86.18(17)	N(7)-Cd(1)-N(4)	93.82(17)

Symmetry code for **1**: #1 $-x, -y+2, -z+1$. for **3**: #1 $-x+1, -y+1, -z$; #2 $-x+1, y, -z$; #3 $x, -y+1, z$. for **4**: #1 $-x+1, -y, -z$; #2 $-x+1, y, -z$. for **5**: #1 $-x, -y, -z$; #2 $x, -y, z$; #