

Electronic Supplementary Information (ESI) for

**Syntheses, Structures, and Magnetic Properties of Six Coordination Polymers Based
on 4,5-Di(4'-carboxylphenyl)phthalic Acid and Different Bis(imidazole) Bridging
Linkers**

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ESI

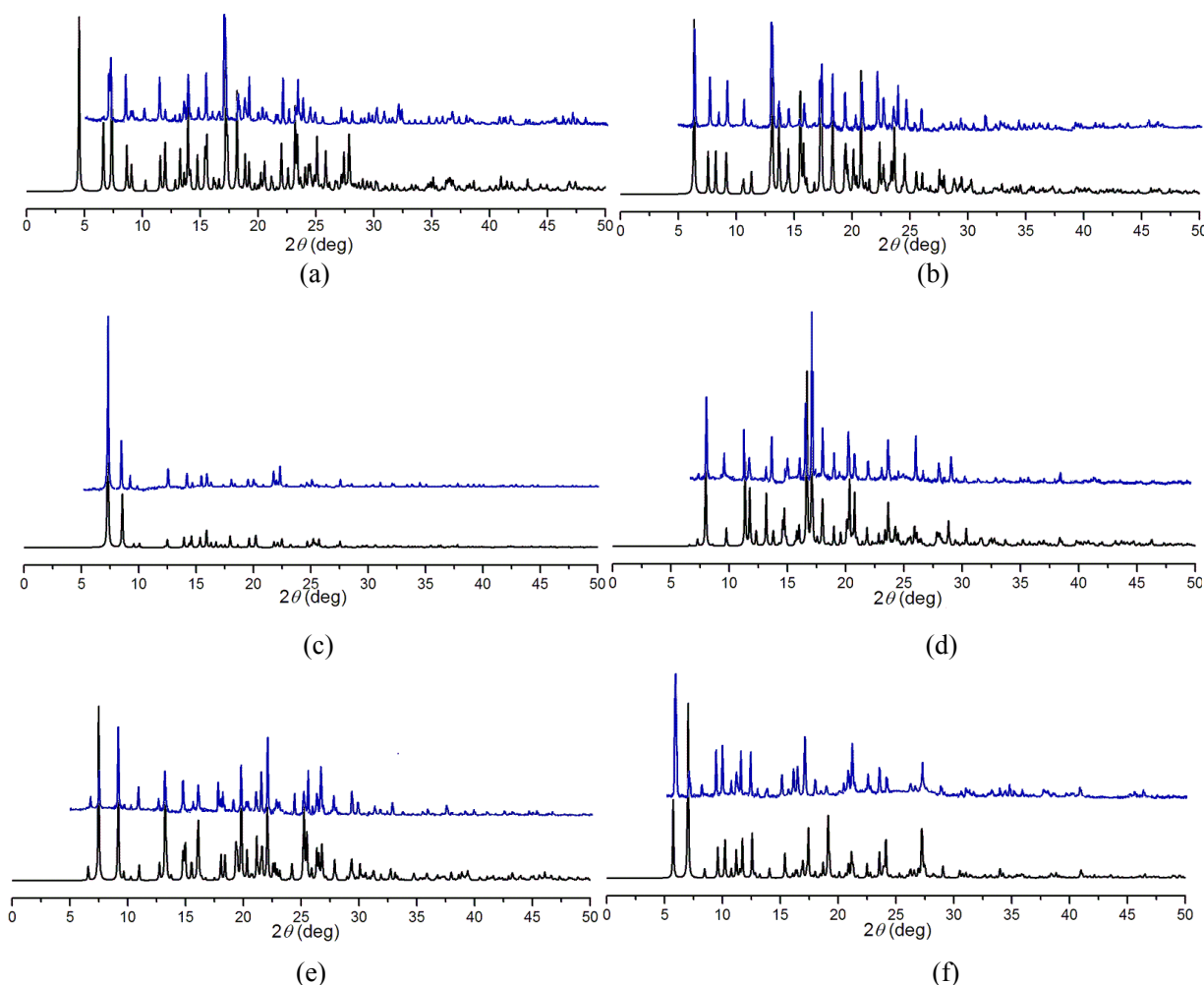
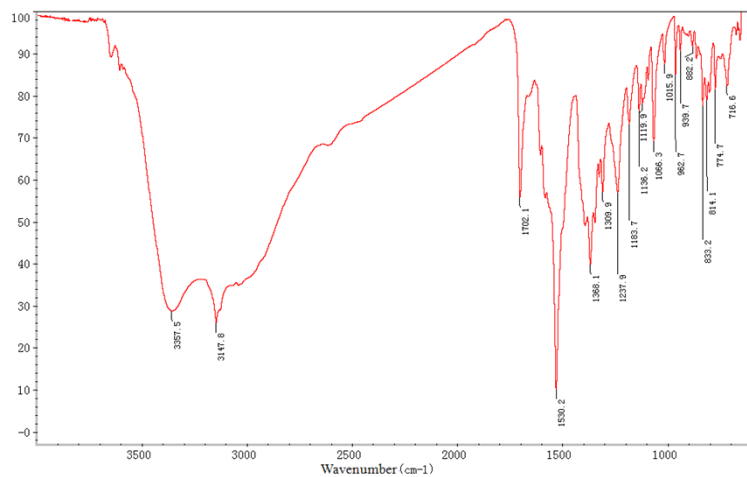
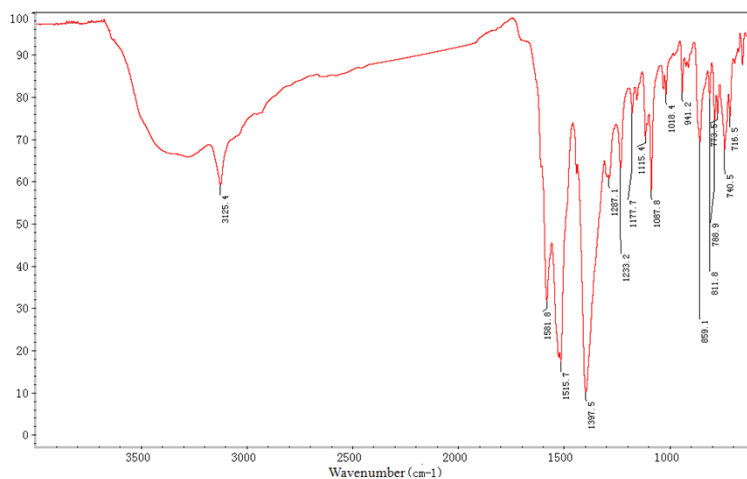


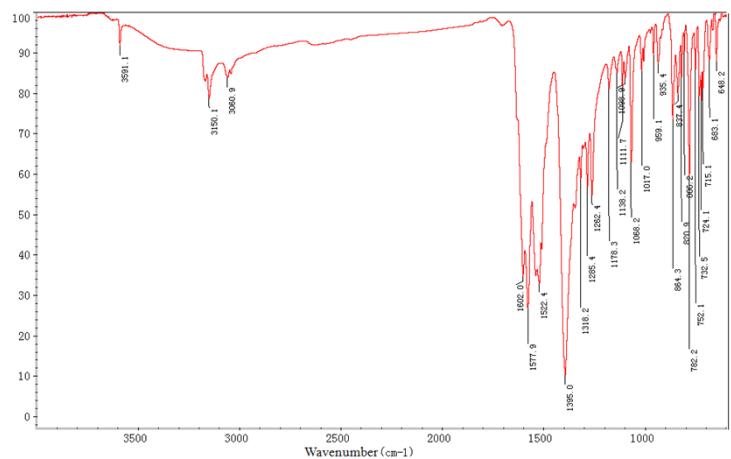
Figure S1. PXRD patterns of **1** (a), **2** (b), **3** (c), **4** (d), **5** (e) and **6** (f). Dark: calculated from the X-ray single-crystal data; Blue: observed for the as-synthesized solids.



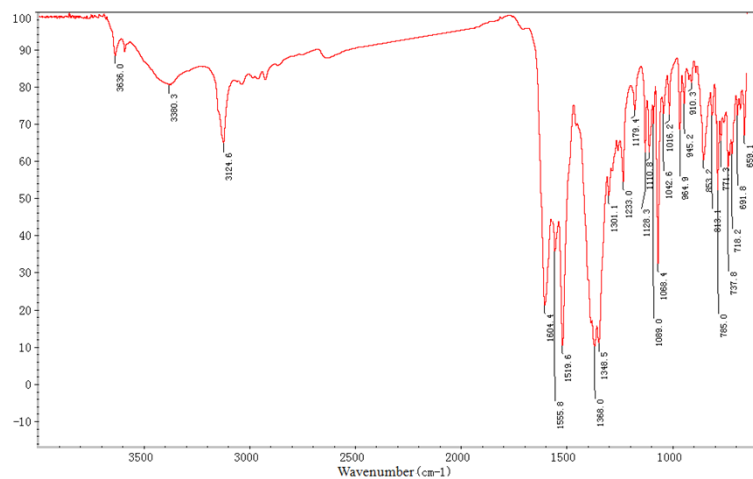
(a)



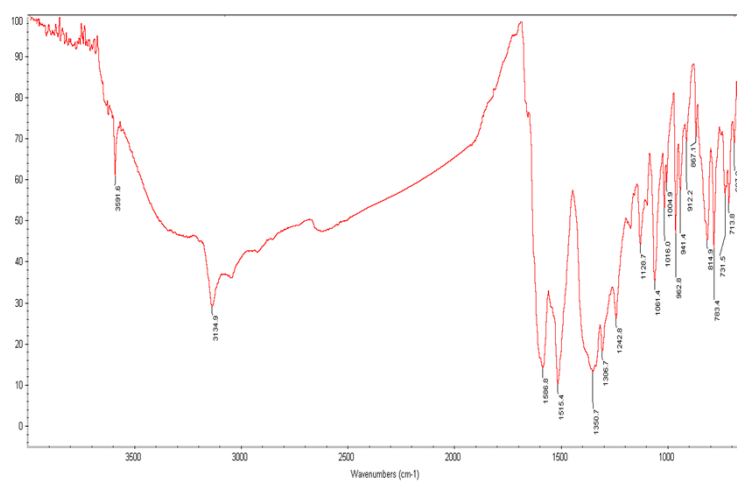
(b)



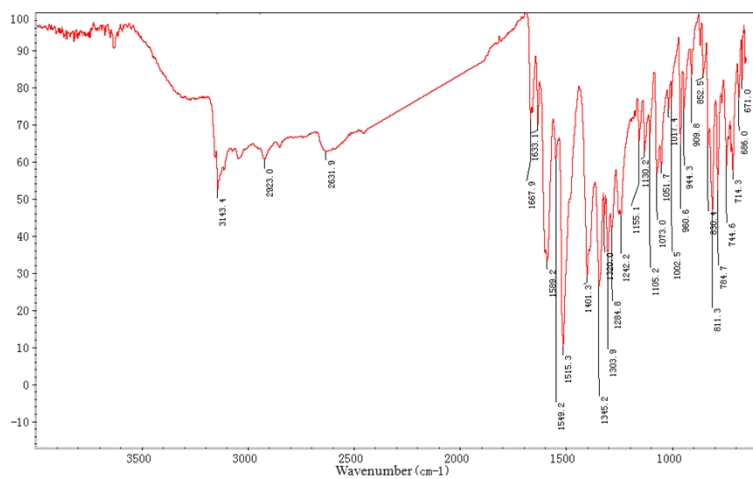
(c)



(d)



(e)



(f)

Figure S2. IR spectra of 1 (a), 2 (b), 3 (c), 4 (d), 5 (e), and 6 (f).

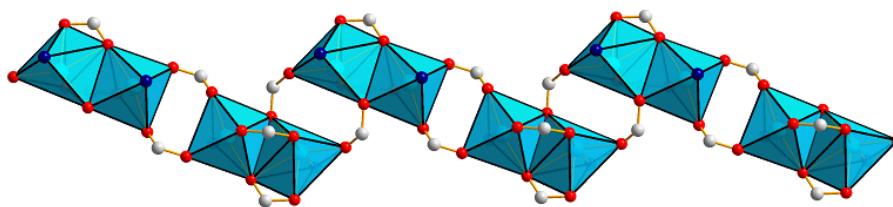


Figure S3. The $[\text{Co}_4(\text{COO})_8]_n$ chain in complex 3.

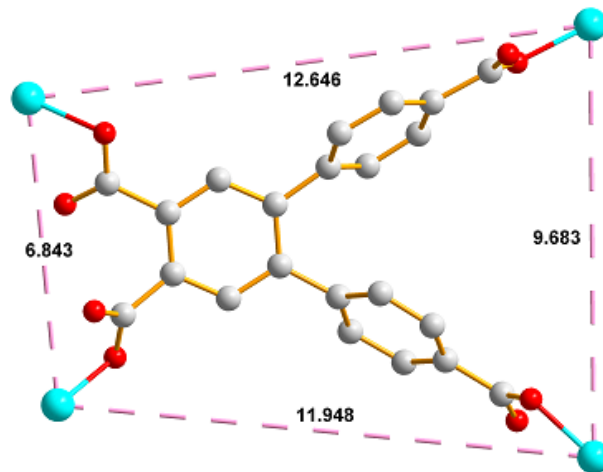


Figure S4. The $\text{Co}\cdots\text{Co}$ distances separated by DCP^{4+} ligand in complex 4.

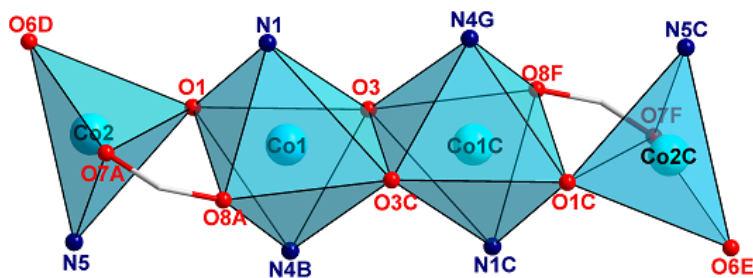


Figure S5. The $\{\text{Co}_4\}$ cluster based SBUs in complex 6.

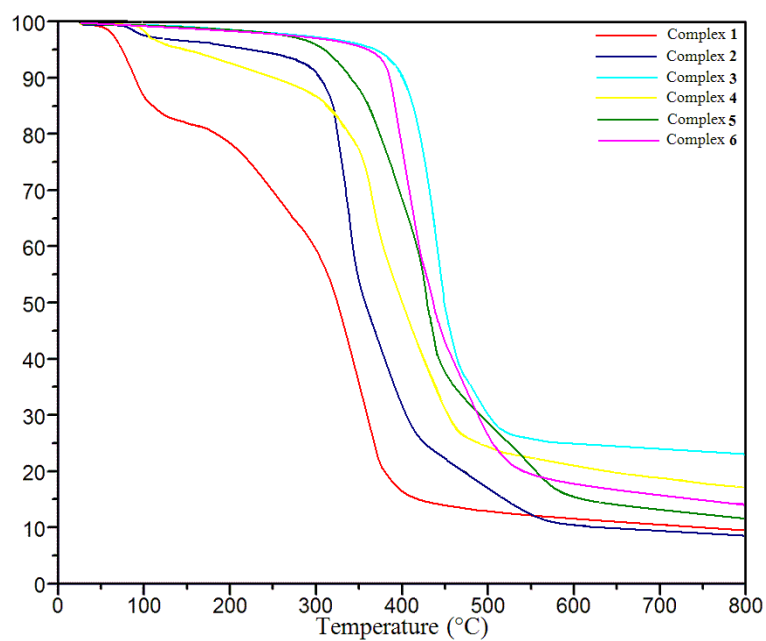


Figure S6. TG curves for complexes 1-6.

Table S1 Selected bond lengths (Å) and angles (°) for **1** – 6.

Complex 1							
N(1)-Ni(1)	2.067(5)	N(3)-Ni(2)	2.080(5)	N(5)-Ni(2)	2.075(6)	Ni(2)-O(13)	2.065(2)
Ni(1)-O(9)	2.071(3)	Ni(2)-O(12)	2.074(2)	Ni(1)-O(1)	2.123(4)	Ni(2)-O(10)	2.083(2)
Ni(2)-O(11)	2.050(2)	N(1)-Ni(1)-O(9)	87.51(17)	N(1) ^{#4} -Ni(1)-O(1)	93.46(19)	O(11)-Ni(1)-O(1)	86.54(19)
N(1)-Ni(1)-O(9) ^{#4}	92.49(17)	O(9)-Ni(1)-O(1)	94.84(14)	O(11)-Ni(2)-O(13)	88.81(8)	O(11)-Ni(2)-O(12)	175.10(12)
O(9) ^{#4} -Ni(1)-O(1)	85.16(14)	O(11)-Ni(2)-N(5)	85.99(17)	O(13)-Ni(2)-N(5)	91.50(18)	O(12)-Ni(2)-N(5)	90.26(18)
O(13)-Ni(2)-O(12)	88.13(8)	O(13)-Ni(2)-N(3)	89.66(17)	O(12)-Ni(2)-N(3)	91.71(17)	N(5)-Ni(2)-N(3)	177.7(2)
O(11)-Ni(2)-N(3)	92.11(17)	O(13)-Ni(2)-O(10)	179.32(12)	O(12)-Ni(2)-O(10)	92.14(8)	N(5)-Ni(2)-O(10)	87.88(19)
O(11)-Ni(2)-O(10)	90.88(8)	N(3)-Ni(2)-O(10)	90.95(17)				
Symmetry codes: #4 -x, -y+1, -z+1.							
Complex 2							
N(4)-Ni(1)	2.081(2)	N(5)-Ni(1)	2.077(2)	N(7)-Ni(1)	2.093(2)	Ni(1)-O(1)	2.0684(17)
Ni(1)-N(1) ^{#4}	2.098(2)	Ni(1)-O(9)	2.1178(19)	O(1)-Ni(1)-N(5)	84.10(8)	O(1)-Ni(1)-N(4)	176.93(8)
N(5)-Ni(1)-N(4)	97.57(8)	O(1)-Ni(1)-N(7)	94.33(8)	N(5)-Ni(1)-N(7)	90.62(8)	N(4)-Ni(1)-N(7)	88.24(8)
O(1)-Ni(1)-N(1) ^{#4}	88.48(8)	N(5)-Ni(1)-N(1) ^{#4}	95.19(8)	N(4)-Ni(1)-N(1) ^{#4}	88.80(8)	N(7)-Ni(1)-N(1) ^{#4}	173.78(8)
O(1)-Ni(1)-O(9)	84.86(7)	N(5)-Ni(1)-O(9)	168.22(8)	N(4)-Ni(1)-O(9)	93.64(8)	N(7)-Ni(1)-O(9)	86.10(9)
N(1) ^{#4} -Ni(1)-O(9)	88.63(9)						
Symmetry codes: #4 -x+3/2, y+1/2, -z+1/2.							
Complex 3							
Co(1)-O(3)	2.012(3)	Co(1)-O(4) ^{#1}	2.048(3)	Co(1)-O(8) ^{#2}	2.079(3)	Co(1)-O(5) ^{#3}	2.111(3)
Co(1)-N(4) ^{#4}	2.138(3)	Co(1)-O(1)	2.281(2)	Co(2)-O(6) ^{#3}	2.023(3)	Co(2)-O(6) ^{#5}	2.023(3)
Co(2)-O(8) ^{#2}	2.338(3)	Co(2)-O(1)	2.084(2)	Co(2)-O(2) ^{#5}	2.027(3)	Co(2)-O(2) ^{#5}	2.027(3)
Co(2)-O(7) ^{#2}	2.162(3)	Co(2)-N(1)	2.067(3)	O(4) ^{#1} -Co(1)-O(8) ^{#2}	96.29(11)	O(4) ^{#1} -Co(1)-O(5) ^{#3}	85.96(11)
O(3)-Co(1)-O(8) ^{#2}	166.19(11)	O(3)-Co(1)-N(4) ^{#4}	93.58(12)	O(3)-Co(1)-O(5) ^{#3}	85.93(11)	O(8) ^{#2} -Co(1)-O(5) ^{#3}	89.70(11)
O(4) ^{#1} -Co(1)-N(4) ^{#4}	98.03(12)	O(5) ^{#3} -Co(1)-N(4) ^{#4}	176.01(12)	O(8) ^{#2} -Co(1)-O(1)	83.22(9)	N(4) ^{#4} -Co(1)-O(1)	83.44(11)
O(8) ^{#2} -Co(1)-N(4) ^{#4}	89.87(12)	O(3)-Co(1)-O(1)	83.89(9)	O(5) ^{#3} -Co(1)-O(1)	92.57(10)	O(3)-Co(1)-O(4) ^{#1}	96.47(11)
O(4) ^{#1} -Co(1)-O(1)	178.45(10)	O(6) ^{#3} -Co(2)-O(2) ^{#5}	82.34(11)	O(2) ^{#5} -Co(2)-O(1)	124.93(10)	O(2) ^{#5} -Co(2)-O(7) ^{#2}	95.94(11)
O(6) ^{#3} -Co(2)-N(1)	168.77(12)	N(1)-Co(2)-O(1)	93.48(11)	N(1)-Co(2)-O(7) ^{#2}	93.54(13)	O(6) ^{#3} -Co(2)-O(8) ^{#2}	93.34(11)
O(2) ^{#5} -Co(2)-N(1)	86.44(12)	O(6) ^{#3} -Co(2)-O(7) ^{#2}	88.04(12)	O(1)-Co(2)-O(7) ^{#2}	138.87(11)	O(2) ^{#5} -Co(2)-O(8) ^{#2}	153.02(10)
O(6) ^{#3} -Co(2)-O(1)	92.61(11)	O(1)-Co(2)-O(8) ^{#2}	81.73(9)	O(7) ^{#2} -Co(2)-O(8) ^{#2}	57.20(10)	N(1)-Co(2)-O(8) ^{#2}	96.86(11)
Symmetry codes: #1 -x+1, -y+1, -z+2. #2 x+1/2, -y+3/2, z+1/2. #3 x, -y+1, z+1/2. #4 x+1/2, y-1/2, z. #5 -x+1/2, -y+3/2, -z+2.							
Complex 4							
Co(1)-O(4) ^{#2}	1.944(5)	Co(1)-N(1)	1.983(5)	Co(2)-O(5) ^{#3}	1.914(4)	Co(2)-O(1)	1.980(4)
Co(1)-N(5)	1.981(5)	Co(1)-O(3) ^{#2}	2.383(6)	Co(2)-N(8) ^{#4}	1.949(5)	Co(2)-N(4) ^{#5}	1.998(5)
O(7)-Co(1)-O(4) ^{#2}	111.0(2)	O(4) ^{#2} -Co(1)-N(5)	123.18(18)	O(4) ^{#2} -Co(1)-N(1)	113.3(2)	O(7)-Co(1)-O(3) ^{#2}	169.41(15)
O(7)-Co(1)-N(5)	104.4(2)	O(7)-Co(1)-N(1)	93.4(2)	N(5)-Co(1)-N(1)	107.2(2)	O(4) ^{#2} -Co(1)-O(3) ^{#2}	59.18(18)
N(5)-Co(1)-O(3) ^{#2}	85.33(19)	O(5) ^{#3} -Co(2)-N(8) ^{#4}	128.19(17)	N(8) ^{#4} -Co(2)-O(1)	110.76(18)	N(8) ^{#4} -Co(2)-N(4) ^{#5}	99.4(2)
N(1)-Co(1)-O(3) ^{#2}	87.59(19)	O(5) ^{#3} -Co(2)-O(1)	102.24(17)	O(5) ^{#3} -Co(2)-N(4) ^{#5}	110.6(2)	O(1)-Co(2)-N(4) ^{#5}	103.44(18)
Symmetry code: #2 x+1, -y+1/2, z+1/2. #3 x, -y+1/2, z-1/2. #4 x-1, y-1, z-1. #5 x, y-1, z.							
Complex 5							
Co(1)-O(2)	1.994(2)	Co(1)-O(8) ^{#2}	2.045(2)	Co(1)-N(1)	2.112(3)	Co(1)-N(4) ^{#3}	2.106(3)
Co(1)-O(7) ^{#1}	2.031(2)	O(2)-Co(1)-O(7) ^{#1}	103.03(11)	O(2)-Co(1)-N(4) ^{#3}	89.15(10)	O(2)-Co(1)-N(1)	84.20(10)
O(8) ^{#2} -Co(1)-N(1)	86.67(9)	O(2)-Co(1)-O(8) ^{#2}	129.58(11)	O(7) ^{#1} -Co(1)-N(4) ^{#3}	94.38(10)	O(7) ^{#1} -Co(1)-N(1)	88.37(10)
N(4) ^{#3} -Co(1)-N(1)	173.23(10)	O(7) ^{#1} -Co(1)-O(8) ^{#2}	126.17(10)	O(8) ^{#2} -Co(1)-N(4) ^{#3}	96.64(10)		
Symmetry code: #1 -x, -y, -z+2. #2 x+1, y, z-1. #3 x-1, y-1, z.							
Complex 6							
Co(1)-O(8) ^{#1}	2.088(2)	Co(1)-N(4) ^{#2}	2.118(3)	Co(1)-O(3) ^{#3}	2.192(2)	Co(1)-O(1)	2.197(2)
Co(1)-N(1)	2.111(3)	Co(1)-O(3)	2.166(2)	Co(2)-O(7) ^{#1}	1.976(2)	Co(2)-O(1)	2.016(2)
Co(2)-N(5)	2.021(3)	Co(2)-O(6) ^{#4}	1.982(3)	O(8) ^{#1} -Co(1)-N(1)	88.11(10)	N(1)-Co(1)-N(4) ^{#2}	172.23(10)
N(1)-Co(1)-O(3)	101.38(9)	N(4) ^{#2} -Co(1)-O(3)	85.37(9)	O(8) ^{#1} -Co(1)-N(4) ^{#2}	85.82(9)	O(8) ^{#1} -Co(1)-O(3)	167.12(8)
N(4) ^{#2} -Co(1)-O(3) ^{#3}	85.76(9)	O(3)-Co(1)-O(3) ^{#3}	79.08(8)	N(4) ^{#2} -Co(1)-O(1)	91.45(9)	O(3) ^{#3} -Co(1)-O(1)	170.49(8)
O(8) ^{#1} -Co(1)-O(3) ^{#3}	90.90(8)	O(8) ^{#1} -Co(1)-O(1)	97.97(8)	O(3)-Co(1)-O(1)	91.64(8)	N(1)-Co(1)-O(1)	84.58(9)
N(1)-Co(1)-O(3) ^{#3}	99.20(9)	O(7) ^{#1} -Co(2)-O(6) ^{#4}	97.37(11)	O(6) ^{#4} -Co(2)-O(1)	109.43(10)	O(6) ^{#4} -Co(2)-N(5)	136.63(13)
O(7) ^{#1} -Co(2)-O(1)	106.13(10)	O(7) ^{#1} -Co(2)-N(5)	98.76(11)	O(1)-Co(2)-N(5)	104.21(10)		
Symmetry code: #1 x, -y+1, z+1/2. #2 x-1/2, y+1/2, z. #3 -x+1, -y+1, -z+1. #4 -x+3/2, -y+3/2, -z+1.							