Acta Crystallographica Section E
Structure Reports Online
ISSN 1600-5368

Poly[[aqua(μ_7-biphenyl-3,3′,4,4′-tetracarboxylato)(1,10-phenanthroline)-dicobalt(II)] monohydrate]

Hailiang Yin,[a]* Fengjuan Yin[b] and Yukun Lu[a]

[a]College of Chemistry and Chemical Engineering, China University of Petroleum, Qingdao 255666, People's Republic of China, and [b]Environmental Protection Bureau of Dongying, Dongying 257091, People's Republic of China
Correspondence e-mail: yinhl2000@163.com

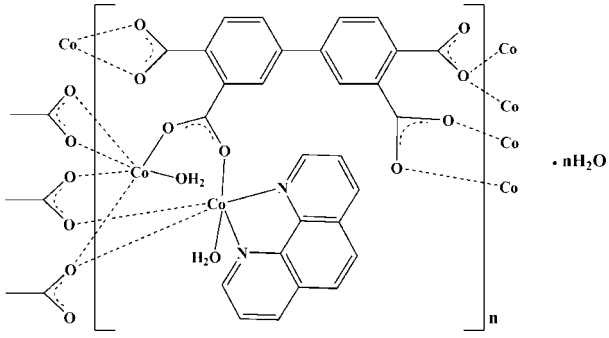
Received 26 June 2011; accepted 14 July 2011

Key indicators: single-crystal X-ray study; T = 298 K; mean σ(C–C) = 0.007 Å; R factor = 0.052; wR factor = 0.113; data-to-parameter ratio = 12.0.

In the title compound, $\{[Co_2(C_{16}H_6O_8)(C_{12}H_8N_2)(H_2O)_2]\cdot H_2O\}_n$, one Co^{II} ion has a $\{CoN_2O_4\}$ distorted octahedral environment defined by two N atoms of one 1,10-phenanthroline (phen) ligand, three O atoms of the carboxylate groups of three biphenyl-3,3′,4,4′-tetracarboxylate (BPTC) ligands, one of which is bidentate, and one O atom from one coordinated water molecule. The other Co^{II} atom is surrounded by six O atoms from four different BPTC ligands and one coordinated water molecule. Each BPTC ligand forms eight coordination bonds with seven Co^{II} atoms, leading to a layer structure along the *ac* plane. Uncoordinated water molecules occupy the space between the layers, and interact *via* interlayer O—H···O hydrogen bonds along the *b* axis, generating a three-dimensional supramolecular network.

Related literature

For applications of compounds with metal-organic framework structures (MOFs), see: Rowsell & Yaghi (2005). For related structures, see: Zhu *et al.* (2008); Konar *et al.* (2004).



Experimental

Crystal data

$[Co_2(C_{16}H_6O_8)(C_{12}H_8N_2)(H_2O)_2]\cdot H_2O$
M_r = 678.32
Triclinic, $P\overline{1}$
a = 9.793 (3) Å
b = 10.885 (3) Å
c = 12.453 (3) Å
α = 97.567 (4)°
β = 102.608 (4)°
γ = 95.653 (4)°
V = 1273.1 (6) Å^3
Z = 2
Mo $K\alpha$ radiation
μ = 1.38 mm^{-1}
T = 298 K
0.40 × 0.17 × 0.16 mm

Data collection

Bruker SMART CCD area-detector diffractometer
Absorption correction: multi-scan (*SADABS*; Bruker, 2000)
T_{min} = 0.609, T_{max} = 0.810
6716 measured reflections
4641 independent reflections
3194 reflections with $I > 2\sigma(I)$
R_{int} = 0.034

Refinement

$R[F^2 > 2\sigma(F^2)]$ = 0.052
$wR(F^2)$ = 0.113
S = 0.98
4641 reflections
388 parameters
H-atom parameters constrained
$\Delta\rho_{max}$ = 0.46 e Å^{-3}
$\Delta\rho_{min}$ = −0.50 e Å^{-3}

Table 1
Hydrogen-bond geometry (Å, °).

D—H···A	D—H	H···A	D···A	D—H···A
O11—H11B···O7^{i}	0.85	2.04	2.886 (6)	173
O11—H11A···O7^{ii}	0.85	2.15	2.931 (5)	152
O10—H10B···O7^{iii}	0.85	2.20	2.668 (4)	115
O10—H10A···O2^{iv}	0.85	1.95	2.756 (4)	159
O9—H9A···O3	0.85	2.09	2.661 (4)	124

Symmetry codes: (i) $-x+1, -y+1, -z+2$; (ii) $x, y-1, z-1$; (iii) $x, y, z-1$; (iv) $x+1, y, z$.

Data collection: *SMART* (Bruker, 2000); cell refinement: *SAINT* (Bruker, 2000); data reduction: *SAINT*; program(s) used to solve structure: *SHELXS97* (Sheldrick, 2008); program(s) used to refine structure: *SHELXL97* (Sheldrick, 2008); molecular graphics: *SHELXTL* (Sheldrick, 2008); software used to prepare material for publication: *SHELXTL*.

The authors thank the National Key Fundamental Research Development Project (973 Project: No. 2010CB226905) and the Fundamental Research Funds for the Central Universities (No. 09CX04045A).

Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: BG2411).

References

Bruker (2000). *SMART*, *SAINT* and *SADABS*. Bruker AXS Inc., Madison, Wisconsin, USA.
Konar, S., Zangrando, E., Drew, M. G. B., Ribas, J. & Chaudhuri, N. R. (2004). *Dalton Trans.* pp. 260–266.
Rowsell, J. L. C. & Yaghi, O. M. (2005). *Angew. Chem. Int. Ed.* **44**, 4670–4679.
Sheldrick, G. M. (2008). *Acta Cryst.* A**64**, 112–122.
Zhu, S. R., Zhang, H., Shao, M., Zhao, Y. M. & Li, M. X. (2008). *Transition Met. Chem.* **33**, 669–680.

supplementary materials

Acta Cryst. (2011). E**67**, m1118 [**doi:10.1107/S1600536811028224**]

Poly[[aqua(μ_7-biphenyl-3,3',4,4'-tetracarboxylato)(1,10-phenanthroline)dicobalt(II)] monohydrate]

H. Yin, F. Yin and Y. Lu

Comment

The assembly of coordination architectures has attracted much attention in recent years due to their potential applications in separation, sorption, hydrogen storage, and catalysis, as well as due to their intriguing topologies such as molecular ladders, grids, rings, boxes, honeycombs, and diamondoids (Rowsell & Yaghi, 2005). Coordination polymers containing biphenylpolycarboxylate and 1,10-phenanthroline as ligands have played a important role in the area of modern coordination chemistry. A few coordination polymers dealing with 3,3',4,4'-biphenyltetracarboxylate (H_4BPTC) and 1,10-phenanthroline (phen) have been reported (Zhu *et al.*, 2008). Herein, we report a new cobalt coordination polymer, $\{[Co_2(C_{16}H_6O_8)(C_{12}H_8N_2)(H_2O)_2]H_2O\}_n$, resulting from reaction of Co^{2+} cations, phen and H_4BPTC under hydrothermal conditions.

As shown in Fig. 1, the asymmetric unit consists of two crystallographically independent Co^{2+} ions, one fully deprotonated $BPTC^{4-}$ anion, a chelating phen ligand, two coordinated water molecules and one lattice water molecule. The Co1 center is in an octahedral environment defined by two N atoms of one phen ligand, three O atoms of carboxylate groups from three BPTC ligands, and one O atom from one coordinated water molecule. The Co1–O bond lengths fall in the range 2.001 (3)–2.149 (3) Å and the two Co1-N distances are 2.108 (4) and 2.143 (4) Å, thus falling in the expected region (Konar, *et al.*, 2004). The Co2 atom is surrounded by six O atoms from four different BPTC ligands and one coordinated water molecule with Co–O distances in the range 2.025 (3)–2.210 (3) Å, and O–Co–O angles varying from 61.42 (11)°–168.23 (11)°. The octahedral coordination around the Co atoms is strongly distorted since the diametrical and non-diametrical bond angles indicate significant deviations from 180° and 90°, respectively. $BPTC^{4-}$ forms eight coordination bonds with seven Co centers. Two carboxylates of $BPTC^{4-}$ act as monodentate bridging and adopt a μ_2-η^2:η^0 coordinated mode, one carboxylate acts as bidentate bridging and adopts a μ_2-η^1:η^1 coordinated mode, while the remaining carboxylate chelates a Co cation. As a result, each $BPTC^{4-}$ forms eight coordination bonds with seven Co centers, leading to a 2D layer structure parallel to the *ac* plane. Lattice water molecules occupy the space between 2D layers, and interact *via* interlayer O–H···O hydrogen bonds along the *b*-axis to generate a 3D supramolecular network (Table 1 and Fig.2).

Experimental

A mixture of $Co(NO_3)_2.6H_2O$ (146 mg, 0.5 mmol),3,3',4,4' -biphenyltetracarboxylate (74 mg, 0.25 mmol), phen (99 mg, 0.5 mmol), NaOH (40 mg, 1.0 mmol) and water (15 ml) were heated at 393 K for 4 days in a sealed 25 ml Teflon-lined stainless steel vessel under autogenous pressure. Slow cooling of the reaction mixture at 2 K/min to room temperature gave salmon pink block crystals.

Refinement

Hydrogen atoms attached to carbon were idealized and included as riding atoms; those attached to oxygen were located in the difference map, idealized and refined as riding. [d(O—H) = 0.85 Å; U_{iso}(H) = 1.2U_{eq}(O)]

Figures

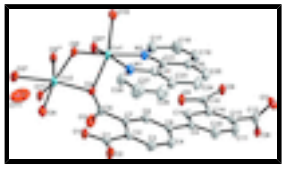


Fig. 1. : View of (I), showing 30% displacement ellipsoids. Symmetry codes: (i) x, y, z - 2; (ii) -x, 1 - y, 1 - z; (iii) 1 - x, 1 - y, 2 - z.

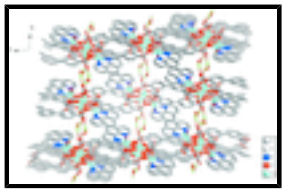


Fig. 2. : Crystal packing of (I) as viewed down the crystallographic a axis.

Poly[[aqua(μ$_7$-biphenyl-3,3',4,4'-tetracarboxylato)(1,10- phenanthroline)dicobalt(II)] monohydrate]

Crystal data

$[Co_2(C_{16}H_6O_8)(C_{12}H_8N_2)(H_2O)_2]\cdot H_2O$	$Z = 2$
$M_r = 678.32$	$F(000) = 688$
Triclinic, $P\overline{1}$	$D_x = 1.770$ Mg m^{-3}
Hall symbol: -P 1	Mo $K\alpha$ radiation, λ = 0.71073 Å
$a = 9.793$ (3) Å	Cell parameters from 1160 reflections
$b = 10.885$ (3) Å	θ = 2.3–22.3°
$c = 12.453$ (3) Å	μ = 1.38 mm^{-1}
α = 97.567 (4)°	$T = 298$ K
β = 102.608 (4)°	Block, red
γ = 95.653 (4)°	0.40 × 0.17 × 0.16 mm
$V = 1273.1$ (6) Å^3	

Data collection

Bruker SMART CCD area-detector diffractometer	4641 independent reflections
Radiation source: fine-focus sealed tube	3194 reflections with $I > 2\sigma(I)$
graphite	$R_{int} = 0.034$
φ and ω scans	$\theta_{max} = 25.5°$, $\theta_{min} = 1.7°$
Absorption correction: multi-scan (*SADABS*; Bruker, 2000)	$h = -11\rightarrow9$
$T_{min} = 0.609$, $T_{max} = 0.810$	$k = -9\rightarrow13$
6716 measured reflections	$l = -14\rightarrow15$

Refinement

Refinement on F^2	Primary atom site location: structure-invariant direct methods
Least-squares matrix: full	Secondary atom site location: difference Fourier map
$R[F^2 > 2\sigma(F^2)] = 0.052$	Hydrogen site location: inferred from neighbouring sites
$wR(F^2) = 0.113$	H-atom parameters constrained
$S = 0.98$	$w = 1/[\sigma^2(F_o^2) + (0.0467P)^2]$ where $P = (F_o^2 + 2F_c^2)/3$
4641 reflections	$(\Delta/\sigma)_{max} = 0.001$
388 parameters	$\Delta\rho_{max} = 0.46$ e Å^{-3}
0 restraints	$\Delta\rho_{min} = -0.50$ e Å^{-3}

Special details

Geometry. All e.s.d.'s (except the e.s.d. in the dihedral angle between two l.s. planes) are estimated using the full covariance matrix. The cell e.s.d.'s are taken into account individually in the estimation of e.s.d.'s in distances, angles and torsion angles; correlations between e.s.d.'s in cell parameters are only used when they are defined by crystal symmetry. An approximate (isotropic) treatment of cell e.s.d.'s is used for estimating e.s.d.'s involving l.s. planes.

Refinement. Refinement of F^2 against ALL reflections. The weighted *R*-factor *wR* and goodness of fit *S* are based on F^2, conventional *R*-factors *R* are based on *F*, with *F* set to zero for negative F^2. The threshold expression of $F^2 > \sigma(F^2)$ is used only for calculating *R*-factors(gt) *etc*. and is not relevant to the choice of reflections for refinement. *R*-factors based on F^2 are statistically about twice as large as those based on *F*, and *R*- factors based on ALL data will be even larger.

Fractional atomic coordinates and isotropic or equivalent isotropic displacement parameters (Å^2)

	x	*y*	*z*	U_{iso}*/U_{eq}
C1	−0.1205 (5)	0.5902 (4)	0.6684 (4)	0.0227 (11)
C2	−0.0391 (5)	0.6133 (4)	0.7872 (3)	0.0209 (10)
C3	−0.0950 (5)	0.6775 (4)	0.8670 (4)	0.0284 (12)
H3	−0.1796	0.7094	0.8450	0.034*
C4	−0.0286 (5)	0.6957 (4)	0.9786 (4)	0.0298 (12)
H4	−0.0699	0.7388	1.0300	0.036*
C5	0.0979 (5)	0.6512 (4)	1.0160 (3)	0.0189 (10)
C6	0.1527 (5)	0.5852 (4)	0.9355 (3)	0.0234 (11)
H6	0.2372	0.5535	0.9583	0.028*
C7	0.0872 (5)	0.5642 (4)	0.8227 (3)	0.0191 (10)
C8	0.1543 (5)	0.4826 (5)	0.7467 (3)	0.0224 (11)
C9	0.1711 (4)	0.6716 (4)	1.1364 (3)	0.0194 (10)
C10	0.1249 (5)	0.7513 (4)	1.2133 (4)	0.0297 (12)
H10	0.0491	0.7939	1.1891	0.036*
C11	0.1904 (5)	0.7684 (5)	1.3263 (4)	0.0317 (12)
H11	0.1585	0.8230	1.3765	0.038*
C12	0.3014 (5)	0.7054 (4)	1.3644 (4)	0.0231 (11)
C13	0.3634 (5)	0.7284 (5)	1.4884 (4)	0.0248 (11)

C14	0.3493 (5)	0.6257 (4)	1.2883 (3)	0.0209 (10)
C15	0.4728 (5)	0.5573 (4)	1.3218 (4)	0.0230 (11)
C16	0.2826 (5)	0.6097 (4)	1.1761 (3)	0.0225 (10)
H16	0.3144	0.5550	1.1260	0.027*
C17	0.5542 (5)	0.7004 (5)	0.9733 (4)	0.0369 (13)
H17	0.6056	0.6388	0.9503	0.044*
C18	0.5861 (6)	0.7519 (5)	1.0863 (4)	0.0459 (15)
H18	0.6582	0.7257	1.1365	0.055*
C19	0.5106 (6)	0.8406 (5)	1.1219 (4)	0.0446 (15)
H19	0.5308	0.8756	1.1968	0.054*
C20	0.4023 (6)	0.8790 (5)	1.0452 (4)	0.0343 (13)
C21	0.3795 (5)	0.8238 (4)	0.9332 (4)	0.0272 (11)
C22	0.2701 (5)	0.8587 (4)	0.8500 (4)	0.0267 (11)
C23	0.1811 (6)	0.9417 (5)	0.8828 (4)	0.0380 (13)
C24	0.2065 (6)	0.9964 (5)	0.9968 (5)	0.0476 (15)
H24	0.1481	1.0524	1.0184	0.057*
C25	0.3128 (7)	0.9689 (5)	1.0736 (5)	0.0475 (15)
H25	0.3290	1.0088	1.1469	0.057*
C26	0.0685 (7)	0.9626 (5)	0.7995 (5)	0.0550 (17)
H26	0.0046	1.0151	0.8173	0.066*
C27	0.0522 (6)	0.9065 (6)	0.6930 (5)	0.0533 (17)
H27	−0.0217	0.9215	0.6373	0.064*
C28	0.1466 (6)	0.8265 (5)	0.6673 (4)	0.0381 (13)
H28	0.1328	0.7869	0.5941	0.046*
Co1	0.39769 (6)	0.67515 (6)	0.72127 (5)	0.02274 (18)
Co2	0.24578 (6)	0.44660 (6)	0.52994 (5)	0.02263 (18)
N1	0.2553 (4)	0.8043 (4)	0.7428 (3)	0.0277 (9)
N2	0.4548 (4)	0.7350 (4)	0.8980 (3)	0.0273 (9)
O1	−0.0796 (3)	0.5208 (3)	0.5955 (2)	0.0285 (8)
O2	−0.2316 (3)	0.6395 (3)	0.6414 (2)	0.0304 (8)
O3	0.1484 (4)	0.3712 (3)	0.7556 (3)	0.0385 (9)
O4	0.2165 (3)	0.5344 (3)	0.6805 (2)	0.0229 (7)
O5	0.4838 (3)	0.4631 (3)	1.2562 (2)	0.0284 (8)
O6	0.5598 (3)	0.6015 (3)	1.4135 (2)	0.0264 (8)
O7	0.4224 (4)	0.8357 (3)	1.5296 (2)	0.0352 (9)
O8	0.3428 (3)	0.6420 (3)	1.5449 (2)	0.0217 (7)
O9	0.1558 (4)	0.2681 (3)	0.5513 (2)	0.0376 (9)
H9A	0.1290	0.2480	0.6078	0.045*
H9B	0.1354	0.2042	0.5007	0.045*
O10	0.5871 (3)	0.7894 (3)	0.7169 (2)	0.0306 (8)
H10A	0.6525	0.7612	0.6901	0.037*
H10B	0.5721	0.8574	0.6925	0.037*
O11	0.3441 (5)	0.0884 (4)	0.5642 (4)	0.0968 (19)
H11A	0.3480	0.0115	0.5697	0.116*
H11B	0.4080	0.1111	0.5312	0.116*

Atomic displacement parameters (Å^2)

	U^{11}	U^{22}	U^{33}	U^{12}	U^{13}	U^{23}
C1	0.015 (2)	0.028 (3)	0.023 (2)	−0.001 (2)	0.001 (2)	0.003 (2)
C2	0.019 (2)	0.027 (3)	0.014 (2)	0.003 (2)	−0.0016 (19)	0.003 (2)
C3	0.021 (3)	0.041 (3)	0.022 (3)	0.013 (2)	0.001 (2)	0.003 (2)
C4	0.033 (3)	0.038 (3)	0.018 (2)	0.011 (2)	0.006 (2)	−0.002 (2)
C5	0.021 (3)	0.021 (3)	0.015 (2)	0.004 (2)	0.0047 (19)	0.0034 (19)
C6	0.021 (3)	0.028 (3)	0.021 (2)	0.007 (2)	0.002 (2)	0.005 (2)
C7	0.023 (3)	0.025 (3)	0.011 (2)	0.004 (2)	0.0062 (18)	0.0035 (19)
C8	0.016 (2)	0.035 (3)	0.013 (2)	0.008 (2)	−0.0029 (19)	0.000 (2)
C9	0.019 (2)	0.024 (3)	0.015 (2)	0.004 (2)	0.0021 (19)	0.003 (2)
C10	0.034 (3)	0.032 (3)	0.023 (3)	0.017 (2)	0.001 (2)	0.005 (2)
C11	0.040 (3)	0.040 (3)	0.016 (2)	0.022 (3)	0.004 (2)	−0.001 (2)
C12	0.024 (3)	0.022 (3)	0.023 (2)	0.001 (2)	0.004 (2)	0.006 (2)
C13	0.025 (3)	0.034 (3)	0.016 (2)	0.016 (2)	0.001 (2)	0.001 (2)
C14	0.023 (3)	0.025 (3)	0.015 (2)	0.007 (2)	0.0029 (19)	0.005 (2)
C15	0.016 (2)	0.034 (3)	0.022 (2)	0.007 (2)	0.008 (2)	0.010 (2)
C16	0.024 (3)	0.024 (3)	0.019 (2)	0.006 (2)	0.002 (2)	0.002 (2)
C17	0.033 (3)	0.050 (4)	0.026 (3)	0.015 (3)	0.002 (2)	0.004 (3)
C18	0.045 (4)	0.062 (4)	0.024 (3)	0.004 (3)	−0.004 (3)	0.005 (3)
C19	0.055 (4)	0.049 (4)	0.022 (3)	−0.014 (3)	0.007 (3)	−0.003 (3)
C20	0.042 (3)	0.036 (3)	0.023 (3)	−0.006 (3)	0.012 (2)	−0.004 (2)
C21	0.032 (3)	0.027 (3)	0.021 (2)	−0.004 (2)	0.009 (2)	−0.002 (2)
C22	0.031 (3)	0.021 (3)	0.028 (3)	0.003 (2)	0.008 (2)	0.001 (2)
C23	0.052 (4)	0.028 (3)	0.039 (3)	0.012 (3)	0.017 (3)	0.004 (2)
C24	0.066 (4)	0.035 (3)	0.050 (4)	0.017 (3)	0.033 (3)	−0.006 (3)
C25	0.065 (4)	0.039 (4)	0.036 (3)	0.001 (3)	0.019 (3)	−0.012 (3)
C26	0.068 (5)	0.045 (4)	0.063 (4)	0.034 (3)	0.027 (4)	0.012 (3)
C27	0.056 (4)	0.064 (4)	0.047 (4)	0.035 (3)	0.011 (3)	0.017 (3)
C28	0.045 (4)	0.043 (4)	0.028 (3)	0.015 (3)	0.010 (3)	0.004 (3)
Co1	0.0226 (4)	0.0293 (4)	0.0144 (3)	0.0090 (3)	0.0007 (3)	−0.0016 (3)
Co2	0.0197 (4)	0.0316 (4)	0.0136 (3)	0.0072 (3)	−0.0016 (3)	−0.0011 (3)
N1	0.032 (2)	0.030 (2)	0.023 (2)	0.0107 (19)	0.0061 (18)	0.0034 (18)
N2	0.024 (2)	0.035 (3)	0.022 (2)	0.0073 (19)	0.0032 (18)	0.0003 (19)
O1	0.0256 (19)	0.041 (2)	0.0142 (16)	0.0109 (16)	−0.0018 (14)	−0.0061 (15)
O2	0.0232 (19)	0.050 (2)	0.0155 (16)	0.0147 (16)	−0.0021 (14)	0.0003 (15)
O3	0.057 (3)	0.033 (2)	0.032 (2)	0.0186 (19)	0.0175 (18)	0.0101 (17)
O4	0.0182 (17)	0.033 (2)	0.0159 (15)	0.0016 (14)	0.0036 (13)	−0.0008 (14)
O5	0.030 (2)	0.033 (2)	0.0183 (16)	0.0156 (16)	−0.0006 (14)	−0.0055 (15)
O6	0.0196 (18)	0.039 (2)	0.0164 (16)	0.0134 (15)	−0.0035 (14)	−0.0042 (15)
O7	0.047 (2)	0.024 (2)	0.0256 (18)	−0.0012 (17)	−0.0046 (16)	−0.0016 (16)
O8	0.0254 (18)	0.0271 (19)	0.0116 (15)	0.0054 (14)	0.0011 (13)	0.0031 (14)
O9	0.052 (2)	0.033 (2)	0.0264 (18)	−0.0014 (17)	0.0126 (17)	−0.0021 (16)
O10	0.031 (2)	0.030 (2)	0.0276 (18)	0.0052 (15)	0.0039 (15)	−0.0027 (15)
O11	0.104 (4)	0.038 (3)	0.184 (5)	0.023 (3)	0.094 (4)	0.031 (3)

Geometric parameters (Å, °)

C1—O2	1.260 (5)	C19—H19	0.9300
C1—O1	1.262 (5)	C20—C21	1.404 (6)
C1—C2	1.495 (6)	C20—C25	1.437 (7)
C1—Co2^{i}	2.467 (4)	C21—N2	1.357 (6)
C2—C3	1.381 (6)	C21—C22	1.437 (6)
C2—C7	1.402 (6)	C22—N1	1.358 (5)
C3—C4	1.379 (6)	C22—C23	1.399 (7)
C3—H3	0.9300	C23—C26	1.398 (7)
C4—C5	1.383 (6)	C23—C24	1.424 (7)
C4—H4	0.9300	C24—C25	1.337 (8)
C5—C6	1.389 (6)	C24—H24	0.9300
C5—C9	1.491 (5)	C25—H25	0.9300
C6—C7	1.390 (6)	C26—C27	1.354 (7)
C6—H6	0.9300	C26—H26	0.9300
C7—C8	1.509 (6)	C27—C28	1.389 (7)
C8—O3	1.229 (5)	C27—H27	0.9300
C8—O4	1.283 (5)	C28—N1	1.322 (6)
C9—C16	1.374 (6)	C28—H28	0.9300
C9—C10	1.386 (6)	Co1—O5^{ii}	2.001 (3)
C10—C11	1.393 (6)	Co1—N1	2.108 (4)
C10—H10	0.9300	Co1—O8^{iii}	2.117 (3)
C11—C12	1.373 (6)	Co1—N2	2.143 (4)
C11—H11	0.9300	Co1—O10	2.144 (3)
C12—C14	1.388 (6)	Co1—O4	2.149 (3)
C12—C13	1.508 (6)	Co2—O6^{ii}	2.025 (3)
C13—O7	1.241 (5)	Co2—O4	2.080 (3)
C13—O8	1.273 (5)	Co2—O1^{i}	2.088 (3)
C14—C16	1.387 (6)	Co2—O9	2.127 (3)
C14—C15	1.493 (6)	Co2—O2^{i}	2.184 (3)
C15—O5	1.253 (5)	Co2—O8^{iii}	2.210 (3)
C15—O6	1.269 (5)	Co2—C1^{i}	2.467 (4)
C16—H16	0.9300	O9—H9A	0.8500
C17—N2	1.318 (6)	O9—H9B	0.8499
C17—C18	1.399 (6)	O10—H10A	0.8499
C17—H17	0.9300	O10—H10B	0.8501
C18—C19	1.359 (8)	O11—H11A	0.8520
C18—H18	0.9300	O11—H11B	0.8548
C19—C20	1.402 (7)		
O2—C1—O1	119.9 (4)	C25—C24—H24	119.3
O2—C1—C2	119.5 (4)	C23—C24—H24	119.3
O1—C1—C2	120.6 (4)	C24—C25—C20	121.3 (5)
O2—C1—Co2^{i}	62.1 (2)	C24—C25—H25	119.3
O1—C1—Co2^{i}	57.8 (2)	C20—C25—H25	119.3
C2—C1—Co2^{i}	177.7 (3)	C27—C26—C23	120.1 (5)

C3—C2—C7	118.1 (4)	C27—C26—H26	119.9
C3—C2—C1	119.1 (4)	C23—C26—H26	119.9
C7—C2—C1	122.6 (4)	C26—C27—C28	119.6 (5)
C4—C3—C2	121.7 (4)	C26—C27—H27	120.2
C4—C3—H3	119.1	C28—C27—H27	120.2
C2—C3—H3	119.1	N1—C28—C27	122.6 (5)
C3—C4—C5	121.6 (4)	N1—C28—H28	118.7
C3—C4—H4	119.2	C27—C28—H28	118.7
C5—C4—H4	119.2	$O5^{ii}$—Co1—N1	161.87 (14)
C4—C5—C6	116.5 (4)	$O5^{ii}$—Co1—$O8^{iii}$	97.71 (11)
C4—C5—C9	121.9 (4)	N1—Co1—$O8^{iii}$	96.93 (13)
C6—C5—C9	121.6 (4)	$O5^{ii}$—Co1—N2	88.47 (13)
C5—C6—C7	123.2 (4)	N1—Co1—N2	77.93 (14)
C5—C6—H6	118.4	$O8^{iii}$—Co1—N2	172.24 (13)
C7—C6—H6	118.4	$O5^{ii}$—Co1—O10	86.77 (13)
C6—C7—C2	118.9 (4)	N1—Co1—O10	104.03 (14)
C6—C7—C8	116.8 (4)	$O8^{iii}$—Co1—O10	89.41 (11)
C2—C7—C8	124.2 (4)	N2—Co1—O10	86.25 (13)
O3—C8—O4	124.7 (4)	$O5^{ii}$—Co1—O4	87.22 (13)
O3—C8—C7	117.0 (4)	N1—Co1—O4	86.00 (14)
O4—C8—C7	118.2 (4)	$O8^{iii}$—Co1—O4	76.00 (11)
C16—C9—C10	117.5 (4)	N2—Co1—O4	109.13 (13)
C16—C9—C5	121.9 (4)	O10—Co1—O4	163.32 (11)
C10—C9—C5	120.6 (4)	$O6^{ii}$—Co2—O4	98.38 (11)
C9—C10—C11	121.0 (4)	$O6^{ii}$—Co2—$O1^{i}$	153.17 (12)
C9—C10—H10	119.5	O4—Co2—$O1^{i}$	107.16 (12)
C11—C10—H10	119.5	$O6^{ii}$—Co2—O9	89.10 (13)
C12—C11—C10	120.7 (4)	O4—Co2—O9	93.31 (12)
C12—C11—H11	119.7	$O1^{i}$—Co2—O9	97.35 (13)
C10—C11—H11	119.7	$O6^{ii}$—Co2—$O2^{i}$	93.38 (11)
C11—C12—C14	119.0 (4)	O4—Co2—$O2^{i}$	168.19 (12)
C11—C12—C13	116.7 (4)	$O1^{i}$—Co2—$O2^{i}$	61.42 (11)
C14—C12—C13	124.3 (4)	O9—Co2—$O2^{i}$	85.65 (12)
O7—C13—O8	124.2 (4)	$O6^{ii}$—Co2—$O8^{iii}$	89.12 (12)
O7—C13—C12	116.5 (4)	O4—Co2—$O8^{iii}$	75.45 (11)
O8—C13—C12	119.0 (4)	$O1^{i}$—Co2—$O8^{iii}$	89.44 (12)
C16—C14—C12	119.4 (4)	O9—Co2—$O8^{iii}$	168.23 (11)
C16—C14—C15	117.9 (4)	$O2^{i}$—Co2—$O8^{iii}$	106.07 (12)
C12—C14—C15	122.6 (4)	$O6^{ii}$—Co2—$C1^{i}$	123.70 (13)
O5—C15—O6	125.4 (4)	O4—Co2—$C1^{i}$	137.77 (14)
O5—C15—C14	117.5 (4)	$O1^{i}$—Co2—$C1^{i}$	30.76 (13)
O6—C15—C14	117.0 (4)	O9—Co2—$C1^{i}$	90.96 (13)
C9—C16—C14	122.4 (4)	$O2^{i}$—Co2—$C1^{i}$	30.68 (13)

C9—C16—H16	118.8	O8^{iii}—Co2—C1^{i}	99.70 (13)
C14—C16—H16	118.8	C28—N1—C22	117.8 (4)
N2—C17—C18	123.2 (5)	C28—N1—Co1	127.4 (3)
N2—C17—H17	118.4	C22—N1—Co1	114.3 (3)
C18—C17—H17	118.4	C17—N2—C21	117.7 (4)
C19—C18—C17	119.3 (5)	C17—N2—Co1	128.8 (3)
C19—C18—H18	120.4	C21—N2—Co1	113.5 (3)
C17—C18—H18	120.4	C1—O1—Co2^{i}	91.4 (3)
C18—C19—C20	119.6 (5)	C1—O2—Co2^{i}	87.2 (3)
C18—C19—H19	120.2	C8—O4—Co2	125.8 (3)
C20—C19—H19	120.2	C8—O4—Co1	128.5 (3)
C19—C20—C21	117.1 (5)	Co2—O4—Co1	98.35 (12)
C19—C20—C25	124.6 (5)	C15—O5—Co1^{ii}	131.1 (3)
C21—C20—C25	118.3 (5)	C15—O6—Co2^{ii}	127.3 (3)
N2—C21—C20	123.1 (5)	C13—O8—Co1^{iv}	121.5 (3)
N2—C21—C22	116.8 (4)	C13—O8—Co2^{iv}	142.9 (3)
C20—C21—C22	120.1 (5)	Co1^{iv}—O8—Co2^{iv}	95.37 (11)
N1—C22—C23	123.3 (5)	Co2—O9—H9A	128.0
N1—C22—C21	117.4 (4)	Co2—O9—H9B	124.2
C23—C22—C21	119.3 (4)	H9A—O9—H9B	107.7
C26—C23—C22	116.5 (5)	Co1—O10—H10A	123.2
C26—C23—C24	124.1 (5)	Co1—O10—H10B	113.4
C22—C23—C24	119.4 (5)	H10A—O10—H10B	107.7
C25—C24—C23	121.4 (5)	H11A—O11—H11B	107.1
O2—C1—C2—C3	6.6 (7)	C26—C27—C28—N1	−1.6 (9)
O1—C1—C2—C3	−172.6 (4)	C27—C28—N1—C22	2.8 (8)
O2—C1—C2—C7	−178.3 (4)	C27—C28—N1—Co1	174.4 (4)
O1—C1—C2—C7	2.5 (7)	C23—C22—N1—C28	−3.7 (7)
C7—C2—C3—C4	1.1 (7)	C21—C22—N1—C28	173.6 (4)
C1—C2—C3—C4	176.4 (4)	C23—C22—N1—Co1	−176.4 (4)
C2—C3—C4—C5	0.5 (8)	C21—C22—N1—Co1	0.9 (5)
C3—C4—C5—C6	−1.4 (7)	O5^{ii}—Co1—N1—C28	−129.0 (5)
C3—C4—C5—C9	179.3 (4)	O8^{iii}—Co1—N1—C28	14.7 (4)
C4—C5—C6—C7	0.6 (7)	N2—Co1—N1—C28	−171.2 (5)
C9—C5—C6—C7	179.9 (4)	O10—Co1—N1—C28	105.8 (4)
C5—C6—C7—C2	1.0 (7)	O4—Co1—N1—C28	−60.7 (4)
C5—C6—C7—C8	−175.6 (4)	O5^{ii}—Co1—N1—C22	42.8 (6)
C3—C2—C7—C6	−1.9 (7)	O8^{iii}—Co1—N1—C22	−173.5 (3)
C1—C2—C7—C6	−176.9 (4)	N2—Co1—N1—C22	0.6 (3)
C3—C2—C7—C8	174.5 (4)	O10—Co1—N1—C22	−82.4 (3)
C1—C2—C7—C8	−0.6 (7)	O4—Co1—N1—C22	111.2 (3)
C6—C7—C8—O3	68.8 (6)	C18—C17—N2—C21	−0.8 (7)
C2—C7—C8—O3	−107.7 (5)	C18—C17—N2—Co1	176.4 (4)
C6—C7—C8—O4	−109.2 (5)	C20—C21—N2—C17	−0.4 (7)
C2—C7—C8—O4	74.4 (6)	C22—C21—N2—C17	−179.0 (4)
C4—C5—C9—C16	168.3 (5)	C20—C21—N2—Co1	−178.1 (4)

C6—C5—C9—C16	−10.9 (7)	C22—C21—N2—Co1	3.3 (5)
C4—C5—C9—C10	−9.3 (7)	O5^{ii}—Co1—N2—C17	12.6 (4)
C6—C5—C9—C10	171.5 (4)	N1—Co1—N2—C17	−179.5 (5)
C16—C9—C10—C11	0.6 (7)	O10—Co1—N2—C17	−74.3 (4)
C5—C9—C10—C11	178.3 (4)	O4—Co1—N2—C17	99.1 (4)
C9—C10—C11—C12	−0.8 (8)	O5^{ii}—Co1—N2—C21	−170.1 (3)
C10—C11—C12—C14	1.1 (7)	N1—Co1—N2—C21	−2.1 (3)
C10—C11—C12—C13	−178.9 (4)	O10—Co1—N2—C21	103.1 (3)
C11—C12—C13—O7	−65.5 (6)	O4—Co1—N2—C21	−83.5 (3)
C14—C12—C13—O7	114.5 (5)	O2—C1—O1—Co2^{i}	2.7 (4)
C11—C12—C13—O8	108.9 (5)	C2—C1—O1—Co2^{i}	−178.1 (4)
C14—C12—C13—O8	−71.1 (6)	O1—C1—O2—Co2^{i}	−2.6 (4)
C11—C12—C14—C16	−1.2 (7)	C2—C1—O2—Co2^{i}	178.2 (4)
C13—C12—C14—C16	178.8 (4)	O3—C8—O4—Co2	28.2 (6)
C11—C12—C14—C15	177.6 (4)	C7—C8—O4—Co2	−154.1 (3)
C13—C12—C14—C15	−2.4 (7)	O3—C8—O4—Co1	−115.2 (4)
C16—C14—C15—O5	−21.1 (6)	C7—C8—O4—Co1	62.6 (5)
C12—C14—C15—O5	160.0 (4)	O6^{ii}—Co2—O4—C8	−93.6 (3)
C16—C14—C15—O6	157.3 (4)	O1^{i}—Co2—O4—C8	94.8 (3)
C12—C14—C15—O6	−21.6 (7)	O9—Co2—O4—C8	−4.0 (3)
C10—C9—C16—C14	−0.8 (7)	O2^{i}—Co2—O4—C8	80.6 (7)
C5—C9—C16—C14	−178.5 (4)	O8^{iii}—Co2—O4—C8	179.6 (3)
C12—C14—C16—C9	1.1 (7)	C1^{i}—Co2—O4—C8	91.1 (4)
C15—C14—C16—C9	−177.8 (4)	O6^{ii}—Co2—O4—Co1	58.29 (14)
N2—C17—C18—C19	1.0 (8)	O1^{i}—Co2—O4—Co1	−113.39 (13)
C17—C18—C19—C20	0.0 (8)	O9—Co2—O4—Co1	147.87 (13)
C18—C19—C20—C21	−1.1 (8)	O2^{i}—Co2—O4—Co1	−127.6 (6)
C18—C19—C20—C25	177.9 (5)	O8^{iii}—Co2—O4—Co1	−28.59 (11)
C19—C20—C21—N2	1.4 (7)	C1^{i}—Co2—O4—Co1	−117.05 (18)
C25—C20—C21—N2	−177.8 (5)	O5^{ii}—Co1—O4—C8	82.0 (4)
C19—C20—C21—C22	179.9 (4)	N1—Co1—O4—C8	−81.2 (4)
C25—C20—C21—C22	0.8 (7)	O8^{iii}—Co1—O4—C8	−179.4 (4)
N2—C21—C22—N1	−2.9 (6)	N2—Co1—O4—C8	−5.5 (4)
C20—C21—C22—N1	178.4 (4)	O10—Co1—O4—C8	150.9 (4)
N2—C21—C22—C23	174.5 (5)	O5^{ii}—Co1—O4—Co2	−68.77 (12)
C20—C21—C22—C23	−4.1 (7)	N1—Co1—O4—Co2	128.06 (14)
N1—C22—C23—C26	3.2 (8)	O8^{iii}—Co1—O4—Co2	29.89 (11)
C21—C22—C23—C26	−174.1 (5)	N2—Co1—O4—Co2	−156.19 (13)
N1—C22—C23—C24	−178.6 (5)	O10—Co1—O4—Co2	0.2 (5)
C21—C22—C23—C24	4.2 (8)	O6—C15—O5—Co1^{ii}	16.8 (7)
C26—C23—C24—C25	177.3 (6)	C14—C15—O5—Co1^{ii}	−165.0 (3)
C22—C23—C24—C25	−0.8 (9)	O5—C15—O6—Co2^{ii}	10.6 (7)
C23—C24—C25—C20	−2.6 (9)	C14—C15—O6—Co2^{ii}	−167.6 (3)
C19—C20—C25—C24	−176.4 (5)	O7—C13—O8—Co1^{iv}	3.7 (6)

C21—C20—C25—C24	2.6 (8)	C12—C13—O8—Co1^{iv}	−170.3 (3)
C22—C23—C26—C27	−1.8 (9)	O7—C13—O8—Co2^{iv}	176.9 (3)
C24—C23—C26—C27	−180.0 (6)	C12—C13—O8—Co2^{iv}	3.0 (7)
C23—C26—C27—C28	1.1 (9)		

Symmetry codes: (i) −*x*, −*y*+1, −*z*+1; (ii) −*x*+1, −*y*+1, −*z*+2; (iii) *x*, *y*, *z*−1; (iv) *x*, *y*, *z*+1.

Hydrogen-bond geometry (Å, °)

D—H···*A*	*D*—H	H···*A*	*D*···*A*	*D*—H···*A*
O11—H11B···O7^{ii}	0.85	2.04	2.886 (6)	173.
O11—H11A···O7^{v}	0.85	2.15	2.931 (5)	152.
O10—H10B···O7^{iii}	0.85	2.20	2.668 (4)	115.
O10—H10A···O2^{vi}	0.85	1.95	2.756 (4)	159.
O9—H9A···O3	0.85	2.09	2.661 (4)	124.

Symmetry codes: (ii) −*x*+1, −*y*+1, −*z*+2; (v) *x*, *y*−1, *z*−1; (iii) *x*, *y*, *z*−1; (vi) *x*+1, *y*, *z*.

Fig. 1

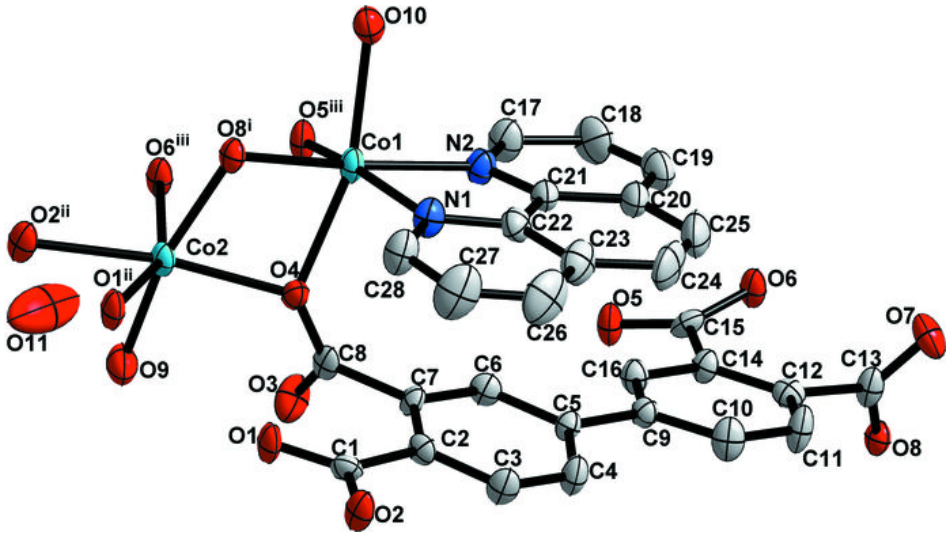

O10
O5iii
C17
C18
O8i
Co1
N2
C19
O6iii
C21
C20
O2ii
N1
C22
C25
Co2
C23
C27
O1ii
O4
C28
C24
O6
O5
O7
O11
C26
C15
C8
C14
C12
C13
O9
C16
C6
C7
O3
C5
C10
O1
C2
C9
C1
C3
C11
O8
C4
O2

Fig. 2