Acta Crystallographica Section E
Structure Reports Online
ISSN 1600-5368

Tetra-μ_2-acetato-diaquabis(μ_2-2-{[1,3-dihydroxy-2-(oxidomethyl)-propan-2-yl]iminomethyl}phenolato)-trimanganese(II,III) acetonitrile disolvate dihydrate

Yuhua Guo, Jianping Huang, Yong Huang, Junyue Wang and Youzhu Yu*

Department of Chemistry and Environmental Engineering, Anyang Institute of Technology, Henan 455000, People's Republic of China
Correspondence e-mail: 119yyz@163.com

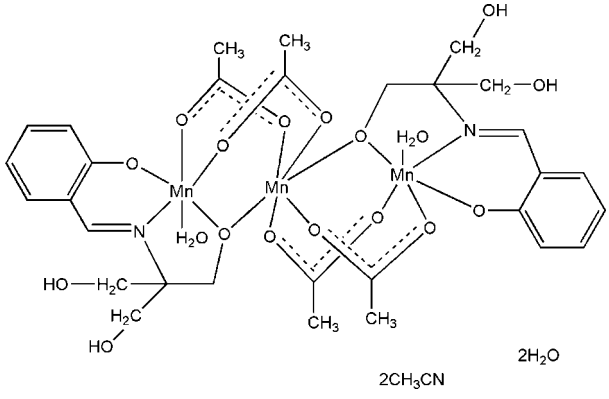
Received 26 June 2011; accepted 12 July 2011

Key indicators: single-crystal X-ray study; T = 293 K; mean σ(C–C) = 0.006 Å; R factor = 0.047; wR factor = 0.133; data-to-parameter ratio = 18.7.

In the title complex, $[Mn^{II}Mn^{III}_2(C_{11}H_{13}NO_4)_2(CH_3CO_2)_4(H_2O)_2]\cdot 2CH_3CN\cdot 2H_2O$, there are two Mn^{III} and one Mn^{II} atoms. The Mn^{II} atom lies on an inversion center and the $Mn^{III}\cdots Mn^{II}\cdots Mn^{III}$ angle is therefore 180°, as required by crystallographic symmetry. The Mn^{III} and Mn^{II} atoms are six-coordinated in a distorted octahedral geometry. In the crystal, complex molecules and solvent molecules are linked into a three-dimensional network by O—H···O and O—H···N hydrogen-bonding interactions.

Related literature

For the importance of Mn complexes in magnetism and biomimetics, see: Stamatatos & Christou (2009); Ferreira *et al.* (2004). For properties and structures of related compounds, see: Kessissoglou *et al.* (1992); Liu *et al.* (2010).



Experimental

Crystal data

$[Mn_3(C_{11}H_{13}NO_4)_2(C_2H_3O_2)_4(H_2O)_2]\cdot 2C_2H_3N\cdot 2H_2O$
M_r = 1001.62
Monoclinic, $P2_1/c$
a = 10.6032 (5) Å
b = 12.2114 (6) Å
c = 19.1608 (9) Å
β = 118.856 (3)°
V = 2172.89 (18) Å^3
Z = 2
Mo $K\alpha$ radiation
μ = 0.94 mm^{-1}
T = 293 K
0.20 × 0.20 × 0.20 mm

Data collection

Bruker APEXII CCD diffractometer
53356 measured reflections
5507 independent reflections
3798 reflections with $I > 2\sigma(I)$
R_{int} = 0.086

Refinement

$R[F^2 > 2\sigma(F^2)]$ = 0.047
$wR(F^2)$ = 0.133
S = 1.07
5507 reflections
294 parameters
7 restraints
H atoms treated by a mixture of independent and constrained refinement
$\Delta\rho_{max}$ = 0.72 e Å^{-3}
$\Delta\rho_{min}$ = −0.61 e Å^{-3}

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

D—H···A	D—H	H···A	D···A	D—H···A
O9—H9A···O10^{i}	0.85 (1)	1.97 (1)	2.806 (4)	167 (3)
O9—H9B···O8^{ii}	0.85 (1)	2.23 (3)	3.008 (3)	153 (5)
O9—H9B···O1^{ii}	0.85 (1)	2.61 (3)	3.322 (3)	142 (5)
O10—H10C···O5^{iii}	0.85 (1)	2.06 (1)	2.907 (4)	176 (5)
O10—H10D···N2^{iv}	0.85 (1)	2.07 (1)	2.914 (6)	174 (6)
O2—H2···O3^{v}	0.82	2.55	3.362 (5)	172
O3—H3···O6^{vi}	0.82	2.00	2.777 (3)	159

Symmetry codes: (i) $x, y+1, z$; (ii) $-x, -y+2, -z+2$; (iii) $-x+1, -y+1, -z+2$; (iv) $x, -y+\frac{3}{2}, z+\frac{1}{2}$; (v) $-x, y-\frac{1}{2}, -z+\frac{3}{2}$; (vi) $-x, y+\frac{1}{2}, -z+\frac{3}{2}$.

Data collection: *APEX2* (Bruker, 1996); cell refinement: *SAINT* (Bruker, 1996); 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*.

This work was supported by the Second Young-Aged Backbone Teachers Foundation of Anyang Institute of Technology.

Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: PV2425).

References

Bruker (1996). *APEX2* and *SAINT*. Bruker AXS Inc., Madison, Wisconsin, USA.
Ferreira, K. N., Iverson, T. M., Maghlaoui, K., Barber, J. & Iwata, S. (2004). *Science*, **303**, 1831–1838.
Kessissoglou, D. P., Kirk, M. L., Lah, M. S., Li, X., Raptopoulou, C., Hatfield, W. E. & Pecoraro, V. L. (1992). *Inorg. Chem.* **31**, 5424–5432.
Liu, D., Zhou, Q., Chen, Y., Yang, F., Yu, Y., Shi, Z. & Feng, S. (2010). *Dalton Trans.* **39**, 5504–5508.
Sheldrick, G. M. (2008). *Acta Cryst.* A**64**, 112–122.
Stamatatos, T. C. & Christou, G. (2009). *Inorg. Chem.* pp. 3308–3322.

supporting information

Acta Cryst. (2011). E**67**, m1098 [doi:10.1107/S1600536811027899]

Tetra-μ_2-acetato-diaquabis(μ_2-2-{[1,3-dihydroxy-2-(oxidomethyl)propan-2-yl]iminomethyl}phenolato)trimanganese(II,III) acetonitrile disolvate dihydrate

Yuhua Guo, Jianping Huang, Yong Huang, Junyue Wang and Youzhu Yu

S1. Comment

The fascination of inorganic chemists with Mn coordination chemistry over the last two decades or so has been primarily driven by the relevance of Mn-complexes to magnetic and biomimetic fields (Stamatatos & Christou, 2009; Ferreira *et al.*2004). As a contribution to these fields, we report here the synthesis and crystal structure of the title compound.

In the title complex (Fig. 1), the Mn^{III} and Mn^{II} atoms are six-coordinated in a distorted octahedral geometry and the two Mn^{III} are in the same coordination environment. The Mn(II) lies on an inversion center, therefore, the angle Mn(III)-Mn(II)—Mn(III) is 180° as required by crystallographic symmetry. The bond lengths and bond angles in the title complex are comparable with those observed in the related complexes (Kessissoglou *et al.*, 1992). In the crystal structure, the complex molecules and the solvent molecules are linked through intermolecular O—H···O and O—H···N hydrogen bonds (Table 1) into a three-dimensional network.

S2. Experimental

To a stirred acetonitrile (20 ml) solution of H_2SALATHM (1 mmol, 225 mg) was added $Mn(OAc)_2.4H_2O$ (1 mmol, 245 mg). The resulting dark-red solution was stirred for 1 h and the filtrate was allowed to stand at room temperature for about three days, whereupon dark block crystal suitable for X-ray diffraction analysis was obtained.

S3. Refinement

H atoms were placed at calculated positions with O–H = 0.82 Å (hydroxyl), and C—H = 0.93 Å (aryl), 0.97 (methylene) and 0.96 Å (methyl) and were refined in the riding-model approximation with U_{iso} = 1.2–1.5 times U_{eq} of the parent atoms. The H-atoms of water of solvation were located from a difference map and were included at distances 0.85 (1) using DFIX commands and were allowed U_{iso} = 1.5 times U_{eq}(O).

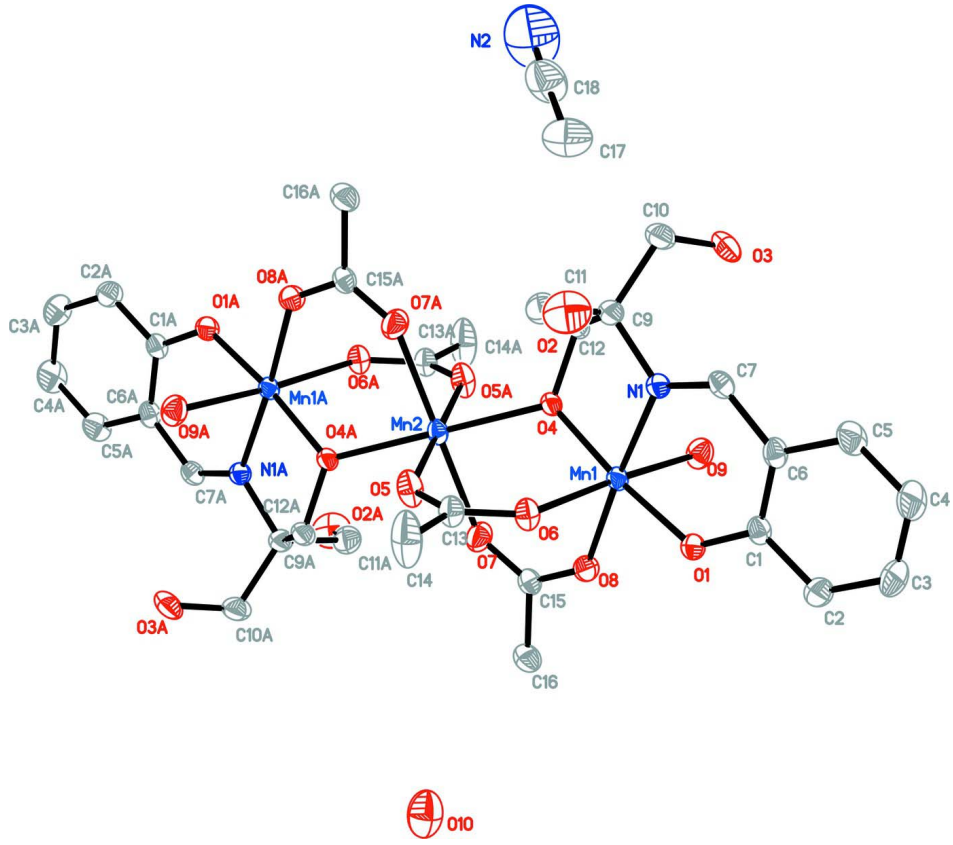


Figure 1

The molecular structure of the title compound, with atom labels and 30% probability displacement ellipsoids.

Tetra-μ_2-acetato-diaquabis(μ_2-2-{[1,3-dihydroxy-2- (oxidomethyl)propan-2-yl]iminomethyl}phenolato)trimanganese(II,III) acetonitrile disolvate dihydrate

Crystal data

$[Mn_3(C_{11}H_{13}NO_4)_2(C_2H_3O_2)_4(H_2O)_2]\cdot 2C_2H_3N\cdot 2H_2O$
$M_r = 1001.62$
Monoclinic, $P2_1/c$
Hall symbol: -P 2ybc
$a = 10.6032\ (5)$ Å
$b = 12.2114\ (6)$ Å
$c = 19.1608\ (9)$ Å
$\beta = 118.856\ (3)°$
$V = 2172.89\ (18)$ Å^3
$Z = 2$
$F(000) = 1038$
$D_x = 1.531$ Mg m^{-3}
Mo $K\alpha$ radiation, $\lambda = 0.71073$ Å
Cell parameters from 9948 reflections
$\theta = 2.4–28.4°$
$\mu = 0.94$ mm^{-1}
$T = 293$ K
Block, black
0.20 × 0.20 × 0.20 mm

Data collection

Bruker APEXII CCD diffractometer
Radiation source: fine-focus sealed tube
Graphite monochromator
φ and ω scans
53356 measured reflections
5507 independent reflections
3798 reflections with $I > 2\sigma(I)$
$R_{int} = 0.086$
$\theta_{max} = 28.6°$, $\theta_{min} = 2.1°$
$h = -14\rightarrow14$
$k = -16\rightarrow16$
$l = -25\rightarrow25$

Refinement

Refinement on F^2
Least-squares matrix: full
$R[F^2 > 2\sigma(F^2)] = 0.047$
$wR(F^2) = 0.133$
$S = 1.07$
5507 reflections
294 parameters
7 restraints
Primary atom site location: structure-invariant direct methods
Secondary atom site location: difference Fourier map
Hydrogen site location: inferred from neighbouring sites
H atoms treated by a mixture of independent and constrained refinement
$w = 1/[\sigma^2(F_o^2) + (0.0525P)^2 + 2.7933P]$ where $P = (F_o^2 + 2F_c^2)/3$
$(\Delta/\sigma)_{max} = 0.024$
$\Delta\rho_{max} = 0.72$ e Å^{-3}
$\Delta\rho_{min} = -0.61$ e Å^{-3}

Fractional atomic coordinates and isotropic or equivalent isotropic displacement parameters (Å²)

	x	*y*	*z*	U_{iso}*/U_{eq}
C1	−0.1569 (3)	0.8511 (2)	0.84008 (18)	0.0317 (6)
C2	−0.2687 (4)	0.8090 (3)	0.8512 (2)	0.0419 (8)
H2A	−0.2506	0.7888	0.9020	0.050*
C3	−0.4053 (4)	0.7971 (3)	0.7877 (2)	0.0548 (10)
H3A	−0.4788	0.7710	0.7966	0.066*
C4	−0.4352 (4)	0.8235 (4)	0.7105 (3)	0.0604 (11)
H4	−0.5274	0.8140	0.6680	0.072*
C5	−0.3275 (4)	0.8635 (3)	0.6980 (2)	0.0475 (9)
H5	−0.3469	0.8808	0.6465	0.057*
C6	−0.1874 (3)	0.8789 (2)	0.76201 (18)	0.0345 (7)
C7	−0.0836 (3)	0.9291 (2)	0.74451 (18)	0.0334 (6)
H7	−0.1103	0.9400	0.6911	0.040*
C9	0.1393 (3)	1.0157 (3)	0.77120 (18)	0.0352 (7)
C10	0.0585 (4)	1.0849 (3)	0.6962 (2)	0.0464 (9)
H10A	0.1265	1.1269	0.6868	0.056*
H10B	0.0049	1.0376	0.6506	0.056*
C11	0.2297 (4)	0.9275 (3)	0.7590 (2)	0.0486 (9)
H11A	0.2880	0.9621	0.7388	0.058*
H11B	0.2946	0.8953	0.8102	0.058*
C12	0.2357 (3)	1.0870 (3)	0.84214 (18)	0.0342 (7)
H12A	0.3199	1.1091	0.8382	0.041*
H12B	0.1840	1.1525	0.8422	0.041*
C13	0.3707 (3)	0.7594 (2)	0.94069 (19)	0.0357 (7)
C14	0.3997 (5)	0.6436 (3)	0.9255 (4)	0.0821 (17)
H14A	0.4021	0.6404	0.8761	0.123*
H14B	0.3248	0.5965	0.9226	0.123*
H14C	0.4907	0.6200	0.9682	0.123*
C15	0.3292 (3)	0.9424 (2)	1.09183 (17)	0.0315 (6)
C16	0.3403 (4)	0.9114 (3)	1.1703 (2)	0.0440 (8)
H16A	0.4374	0.9225	1.2120	0.066*
H16B	0.3147	0.8358	1.1690	0.066*
H16C	0.2761	0.9561	1.1801	0.066*
C17	0.1100 (6)	0.8401 (4)	0.5308 (3)	0.0706 (13)

H17B	0.0142	0.8391	0.4864	0.106*
H17A	0.1057	0.8323	0.5794	0.106*
H17C	0.1644	0.7806	0.5258	0.106*
C18	0.1775 (5)	0.9407 (4)	0.5317 (3)	0.0742 (14)
H9A	0.082 (4)	1.158 (2)	0.938 (3)	0.111*
H9B	−0.014 (5)	1.099 (3)	0.953 (3)	0.111*
H10C	0.320 (2)	0.242 (4)	0.984 (3)	0.111*
H10D	0.235 (5)	0.327 (3)	0.987 (3)	0.111*
N1	0.0433 (3)	0.9604 (2)	0.79651 (14)	0.0299 (5)
N2	0.2282 (7)	1.0199 (4)	0.5303 (5)	0.146 (3)
O1	−0.0288 (2)	0.86304 (18)	0.90298 (12)	0.0356 (5)
O2	0.1473 (4)	0.8430 (3)	0.7059 (2)	0.0775 (9)
H2	0.1298	0.7957	0.7305	0.116*
O3	−0.0374 (3)	1.1568 (2)	0.70565 (16)	0.0569 (7)
H3	−0.0849	1.1909	0.6642	0.085*
O4	0.2792 (2)	1.02733 (16)	0.91436 (11)	0.0290 (4)
O5	0.4759 (2)	0.82092 (18)	0.97622 (15)	0.0466 (6)
O6	0.2412 (2)	0.78562 (16)	0.91590 (13)	0.0359 (5)
O7	0.4363 (2)	0.9780 (2)	1.08976 (14)	0.0480 (6)
O8	0.2062 (2)	0.92758 (18)	1.03081 (12)	0.0358 (5)
O9	0.0281 (3)	1.10148 (19)	0.92495 (16)	0.0451 (6)
O10	0.2353 (3)	0.2668 (3)	0.9640 (2)	0.0640 (8)
Mn1	0.13180 (4)	0.94343 (3)	0.91490 (2)	0.02691 (13)
Mn2	0.5000	1.0000	1.0000	0.02878 (16)

Atomic displacement parameters (Å²)

	U^{11}	U^{22}	U^{33}	U^{12}	U^{13}	U^{23}
C1	0.0261 (14)	0.0260 (14)	0.0358 (16)	−0.0014 (11)	0.0091 (12)	−0.0011 (11)
C2	0.0379 (17)	0.0424 (18)	0.0430 (19)	−0.0061 (14)	0.0176 (15)	0.0013 (14)
C3	0.0326 (17)	0.061 (2)	0.064 (3)	−0.0180 (17)	0.0181 (17)	−0.0055 (19)
C4	0.0317 (18)	0.069 (3)	0.057 (2)	−0.0176 (18)	0.0032 (17)	−0.003 (2)
C5	0.0363 (18)	0.050 (2)	0.0377 (19)	−0.0101 (15)	0.0030 (14)	−0.0016 (15)
C6	0.0258 (14)	0.0306 (15)	0.0382 (17)	−0.0024 (12)	0.0085 (12)	−0.0031 (12)
C7	0.0313 (15)	0.0351 (16)	0.0260 (14)	0.0012 (12)	0.0075 (12)	0.0006 (12)
C9	0.0302 (15)	0.0398 (16)	0.0328 (16)	−0.0026 (13)	0.0129 (13)	0.0056 (13)
C10	0.0397 (18)	0.055 (2)	0.0352 (18)	−0.0057 (16)	0.0104 (15)	0.0132 (15)
C11	0.049 (2)	0.059 (2)	0.042 (2)	0.0034 (17)	0.0253 (17)	−0.0003 (16)
C12	0.0273 (14)	0.0342 (15)	0.0340 (16)	−0.0017 (12)	0.0092 (12)	0.0082 (12)
C13	0.0308 (15)	0.0274 (14)	0.0443 (18)	−0.0002 (12)	0.0145 (14)	−0.0049 (13)
C14	0.045 (2)	0.040 (2)	0.139 (5)	0.0000 (17)	0.026 (3)	−0.030 (3)
C15	0.0321 (15)	0.0281 (14)	0.0313 (15)	0.0043 (12)	0.0130 (12)	−0.0022 (12)
C16	0.0445 (19)	0.0471 (19)	0.0348 (18)	0.0055 (15)	0.0148 (15)	0.0028 (14)
C17	0.102 (4)	0.049 (2)	0.069 (3)	−0.006 (2)	0.047 (3)	−0.004 (2)
C18	0.074 (3)	0.047 (2)	0.087 (3)	0.001 (2)	0.027 (3)	0.011 (2)
N1	0.0262 (12)	0.0323 (13)	0.0284 (12)	−0.0012 (10)	0.0110 (10)	0.0024 (10)
N2	0.124 (5)	0.065 (3)	0.225 (8)	−0.021 (3)	0.065 (5)	0.031 (4)
O1	0.0247 (10)	0.0422 (12)	0.0327 (11)	−0.0046 (9)	0.0080 (9)	0.0057 (9)

O2	0.101 (3)	0.072 (2)	0.068 (2)	−0.0063 (19)	0.047 (2)	−0.0187 (17)
O3	0.0442 (14)	0.0612 (17)	0.0493 (15)	0.0126 (12)	0.0099 (12)	0.0258 (13)
O4	0.0218 (9)	0.0311 (10)	0.0286 (10)	0.0003 (8)	0.0078 (8)	0.0061 (8)
O5	0.0310 (11)	0.0290 (11)	0.0633 (16)	−0.0016 (9)	0.0096 (11)	−0.0065 (11)
O6	0.0285 (10)	0.0283 (10)	0.0440 (13)	−0.0011 (8)	0.0120 (9)	−0.0063 (9)
O7	0.0313 (12)	0.0740 (17)	0.0372 (13)	−0.0035 (12)	0.0152 (10)	−0.0003 (12)
O8	0.0306 (11)	0.0434 (12)	0.0291 (11)	−0.0049 (9)	0.0109 (9)	−0.0004 (9)
O9	0.0436 (14)	0.0384 (13)	0.0596 (16)	0.0039 (10)	0.0299 (12)	−0.0016 (11)
O10	0.0495 (16)	0.0581 (18)	0.081 (2)	0.0004 (13)	0.0286 (16)	−0.0148 (15)
Mn1	0.0215 (2)	0.0281 (2)	0.0258 (2)	−0.00170 (16)	0.00718 (17)	0.00117 (17)
Mn2	0.0204 (3)	0.0280 (3)	0.0314 (3)	−0.0005 (2)	0.0073 (2)	−0.0008 (2)

Geometric parameters (Å, º)

C1—O1	1.319 (3)	C14—H14B	0.9600
C1—C2	1.400 (4)	C14—H14C	0.9600
C1—C6	1.411 (4)	C15—O7	1.234 (4)
C2—C3	1.378 (5)	C15—O8	1.276 (3)
C2—H2A	0.9300	C15—C16	1.498 (4)
C3—C4	1.393 (6)	C16—H16A	0.9600
C3—H3A	0.9300	C16—H16B	0.9600
C4—C5	1.365 (5)	C16—H16C	0.9600
C4—H4	0.9300	C17—C18	1.418 (6)
C5—C6	1.410 (4)	C17—H17B	0.9600
C5—H5	0.9300	C17—H17A	0.9600
C6—C7	1.434 (4)	C17—H17C	0.9600
C7—N1	1.287 (4)	C18—N2	1.113 (6)
C7—H7	0.9300	N1—Mn1	2.004 (2)
C9—N1	1.484 (4)	O1—Mn1	1.882 (2)
C9—C12	1.520 (4)	O2—H2	0.8200
C9—C10	1.525 (4)	O3—H3	0.8200
C9—C11	1.533 (5)	O4—Mn1	1.8729 (19)
C10—O3	1.420 (5)	O4—Mn2	2.1395 (18)
C10—H10A	0.9700	O5—Mn2	2.223 (2)
C10—H10B	0.9700	O6—Mn1	2.244 (2)
C11—O2	1.416 (5)	O7—Mn2	2.147 (2)
C11—H11A	0.9700	O8—Mn1	1.975 (2)
C11—H11B	0.9700	O9—Mn1	2.275 (2)
C12—O4	1.429 (3)	O9—H9A	0.851 (10)
C12—H12A	0.9700	O9—H9B	0.851 (10)
C12—H12B	0.9700	O10—H10C	0.851 (10)
C13—O5	1.241 (4)	O10—H10D	0.849 (10)
C13—O6	1.258 (4)	Mn2—O4^{i}	2.1395 (18)
C13—C14	1.505 (5)	Mn2—O7^{i}	2.147 (2)
C14—H14A	0.9600	Mn2—O5^{i}	2.223 (2)
O1—C1—C2	118.4 (3)	H16A—C16—H16B	109.5
O1—C1—C6	123.4 (3)	C15—C16—H16C	109.5

C2—C1—C6	118.2 (3)	H16A—C16—H16C	109.5
C3—C2—C1	120.7 (3)	H16B—C16—H16C	109.5
C3—C2—H2A	119.7	C18—C17—H17B	109.5
C1—C2—H2A	119.7	C18—C17—H17A	109.5
C2—C3—C4	121.1 (3)	H17B—C17—H17A	109.5
C2—C3—H3A	119.4	C18—C17—H17C	109.5
C4—C3—H3A	119.4	H17B—C17—H17C	109.5
C5—C4—C3	119.3 (3)	H17A—C17—H17C	109.5
C5—C4—H4	120.4	N2—C18—C17	178.1 (7)
C3—C4—H4	120.4	C7—N1—C9	120.5 (3)
C4—C5—C6	120.9 (3)	C7—N1—Mn1	126.1 (2)
C4—C5—H5	119.6	C9—N1—Mn1	113.37 (18)
C6—C5—H5	119.6	C1—O1—Mn1	129.90 (19)
C5—C6—C1	119.8 (3)	C11—O2—H2	109.5
C5—C6—C7	117.3 (3)	C10—O3—H3	109.5
C1—C6—C7	122.8 (3)	C12—O4—Mn1	113.83 (16)
N1—C7—C6	125.4 (3)	C12—O4—Mn2	122.99 (17)
N1—C7—H7	117.3	Mn1—O4—Mn2	121.17 (9)
C6—C7—H7	117.3	C13—O5—Mn2	133.8 (2)
N1—C9—C12	103.9 (2)	C13—O6—Mn1	133.38 (19)
N1—C9—C10	113.5 (3)	C15—O7—Mn2	136.1 (2)
C12—C9—C10	110.8 (3)	C15—O8—Mn1	133.9 (2)
N1—C9—C11	108.0 (3)	Mn1—O9—H9A	116 (2)
C12—C9—C11	109.8 (3)	Mn1—O9—H9B	116 (3)
C10—C9—C11	110.7 (3)	H9A—O9—H9B	109 (2)
O3—C10—C9	109.5 (3)	H10C—O10—H10D	110 (3)
O3—C10—H10A	109.8	O4—Mn1—O1	173.45 (9)
C9—C10—H10A	109.8	O4—Mn1—O8	100.27 (9)
O3—C10—H10B	109.8	O1—Mn1—O8	86.11 (9)
C9—C10—H10B	109.8	O4—Mn1—N1	82.72 (9)
H10A—C10—H10B	108.2	O1—Mn1—N1	90.84 (9)
O2—C11—C9	114.0 (3)	O8—Mn1—N1	176.23 (10)
O2—C11—H11A	108.8	O4—Mn1—O6	92.33 (8)
C9—C11—H11A	108.8	O1—Mn1—O6	89.24 (9)
O2—C11—H11B	108.8	O8—Mn1—O6	88.97 (9)
C9—C11—H11B	108.8	N1—Mn1—O6	93.22 (9)
H11A—C11—H11B	107.6	O4—Mn1—O9	88.58 (9)
O4—C12—C9	109.7 (2)	O1—Mn1—O9	90.39 (9)
O4—C12—H12A	109.7	O8—Mn1—O9	86.27 (9)
C9—C12—H12A	109.7	N1—Mn1—O9	91.54 (10)
O4—C12—H12B	109.7	O6—Mn1—O9	175.23 (9)
C9—C12—H12B	109.7	O4^{i}—Mn2—O4	180.000 (1)
H12A—C12—H12B	108.2	O4^{i}—Mn2—O7^{i}	89.03 (8)
O5—C13—O6	125.6 (3)	O4—Mn2—O7^{i}	90.98 (8)
O5—C13—C14	117.5 (3)	O4^{i}—Mn2—O7	90.97 (8)
O6—C13—C14	117.0 (3)	O4—Mn2—O7	89.02 (8)
C13—C14—H14A	109.5	O7^{i}—Mn2—O7	180.000 (1)
C13—C14—H14B	109.5	O4^{i}—Mn2—O5	88.78 (8)

H14A—C14—H14B	109.5	O4—Mn2—O5	91.22 (8)
C13—C14—H14C	109.5	O7^{i}—Mn2—O5	90.48 (10)
H14A—C14—H14C	109.5	O7—Mn2—O5	89.52 (10)
H14B—C14—H14C	109.5	O4^{i}—Mn2—O5^{i}	91.22 (8)
O7—C15—O8	124.7 (3)	O4—Mn2—O5^{i}	88.78 (8)
O7—C15—C16	119.5 (3)	O7^{i}—Mn2—O5^{i}	89.52 (10)
O8—C15—C16	115.7 (3)	O7—Mn2—O5^{i}	90.48 (10)
C15—C16—H16A	109.5	O5—Mn2—O5^{i}	180.000 (1)
C15—C16—H16B	109.5		

O1—C1—C2—C3	178.6 (3)	Mn2—O4—Mn1—O6	−51.58 (12)
C6—C1—C2—C3	−1.2 (5)	C12—O4—Mn1—O9	−71.9 (2)
C1—C2—C3—C4	1.9 (6)	Mn2—O4—Mn1—O9	123.73 (12)
C2—C3—C4—C5	−1.1 (7)	C1—O1—Mn1—O4	−2.9 (10)
C3—C4—C5—C6	−0.4 (6)	C1—O1—Mn1—O8	164.2 (3)
C4—C5—C6—C1	1.0 (5)	C1—O1—Mn1—N1	−13.6 (3)
C4—C5—C6—C7	−175.3 (4)	C1—O1—Mn1—O6	−106.8 (3)
O1—C1—C6—C5	180.0 (3)	C1—O1—Mn1—O9	78.0 (3)
C2—C1—C6—C5	−0.2 (5)	C15—O8—Mn1—O4	−15.7 (3)
O1—C1—C6—C7	−3.9 (5)	C15—O8—Mn1—O1	165.8 (3)
C2—C1—C6—C7	175.9 (3)	C15—O8—Mn1—N1	−158.1 (13)
C5—C6—C7—N1	171.7 (3)	C15—O8—Mn1—O6	76.5 (3)
C1—C6—C7—N1	−4.5 (5)	C15—O8—Mn1—O9	−103.6 (3)
N1—C9—C10—O3	51.4 (4)	C7—N1—Mn1—O4	−173.7 (3)
C12—C9—C10—O3	−65.0 (3)	C9—N1—Mn1—O4	5.8 (2)
C11—C9—C10—O3	173.0 (3)	C7—N1—Mn1—O1	5.1 (3)
N1—C9—C11—O2	54.4 (4)	C9—N1—Mn1—O1	−175.4 (2)
C12—C9—C11—O2	167.1 (3)	C7—N1—Mn1—O8	−30.9 (16)
C10—C9—C11—O2	−70.3 (4)	C9—N1—Mn1—O8	148.6 (14)
N1—C9—C12—O4	42.2 (3)	C7—N1—Mn1—O6	94.4 (3)
C10—C9—C12—O4	164.4 (3)	C9—N1—Mn1—O6	−86.1 (2)
C11—C9—C12—O4	−73.1 (3)	C7—N1—Mn1—O9	−85.3 (3)
C6—C7—N1—C9	−177.5 (3)	C9—N1—Mn1—O9	94.2 (2)
C6—C7—N1—Mn1	1.9 (4)	C13—O6—Mn1—O4	26.9 (3)
C12—C9—N1—C7	152.4 (3)	C13—O6—Mn1—O1	−159.5 (3)
C10—C9—N1—C7	32.1 (4)	C13—O6—Mn1—O8	−73.4 (3)
C11—C9—N1—C7	−91.0 (3)	C13—O6—Mn1—N1	109.7 (3)
C12—C9—N1—Mn1	−27.1 (3)	C13—O6—Mn1—O9	−74.0 (11)
C10—C9—N1—Mn1	−147.4 (2)	C12—O4—Mn2—O4^{i}	150 (48)
C11—C9—N1—Mn1	89.5 (3)	Mn1—O4—Mn2—O4^{i}	−47 (48)
C2—C1—O1—Mn1	−164.9 (2)	C12—O4—Mn2—O7^{i}	−23.4 (2)
C6—C1—O1—Mn1	14.9 (4)	Mn1—O4—Mn2—O7^{i}	139.47 (13)
C9—C12—O4—Mn1	−41.2 (3)	C12—O4—Mn2—O7	156.6 (2)
C9—C12—O4—Mn2	122.8 (2)	Mn1—O4—Mn2—O7	−40.53 (13)
O6—C13—O5—Mn2	−9.9 (6)	C12—O4—Mn2—O5	−113.9 (2)
C14—C13—O5—Mn2	169.9 (3)	Mn1—O4—Mn2—O5	48.97 (13)
O5—C13—O6—Mn1	3.4 (5)	C12—O4—Mn2—O5^{i}	66.1 (2)
C14—C13—O6—Mn1	−176.4 (3)	Mn1—O4—Mn2—O5^{i}	−131.03 (13)

O8—C15—O7—Mn2	−15.7 (5)	C15—O7—Mn2—O4^{i}	−148.5 (3)
C16—C15—O7—Mn2	163.0 (2)	C15—O7—Mn2—O4	31.5 (3)
O7—C15—O8—Mn1	3.5 (5)	C15—O7—Mn2—O7^{i}	112 (100)
C16—C15—O8—Mn1	−175.3 (2)	C15—O7—Mn2—O5	−59.7 (3)
C12—O4—Mn1—O1	9.0 (9)	C15—O7—Mn2—O5^{i}	120.3 (3)
Mn2—O4—Mn1—O1	−155.3 (8)	C13—O5—Mn2—O4^{i}	166.7 (3)
C12—O4—Mn1—O8	−157.9 (2)	C13—O5—Mn2—O4	−13.3 (3)
Mn2—O4—Mn1—O8	37.79 (13)	C13—O5—Mn2—O7^{i}	−104.3 (3)
C12—O4—Mn1—N1	19.8 (2)	C13—O5—Mn2—O7	75.7 (3)
Mn2—O4—Mn1—N1	−144.53 (13)	C13—O5—Mn2—O5^{i}	30 (100)
C12—O4—Mn1—O6	112.8 (2)		

Symmetry code: (i) −*x*+1, −*y*+2, −*z*+2.

Hydrogen-bond geometry (Å, º)

D—H···*A*	*D*—H	H···*A*	*D*···*A*	*D*—H···*A*
O9—H9*A*···O10^{ii}	0.85 (1)	1.97 (1)	2.806 (4)	167 (3)
O9—H9*B*···O8^{iii}	0.85 (1)	2.23 (3)	3.008 (3)	153 (5)
O9—H9*B*···O1^{iii}	0.85 (1)	2.61 (3)	3.322 (3)	142 (5)
O10—H10*C*···O5^{iv}	0.85 (1)	2.06 (1)	2.907 (4)	176 (5)
O10—H10*D*···N2^{v}	0.85 (1)	2.07 (1)	2.914 (6)	174 (6)
O2—H2···O3^{vi}	0.82	2.55	3.362 (5)	172
O3—H3···O6^{vii}	0.82	2.00	2.777 (3)	159

Symmetry codes: (ii) *x*, *y*+1, *z*; (iii) −*x*, −*y*+2, −*z*+2; (iv) −*x*+1, −*y*+1, −*z*+2; (v) *x*, −*y*+3/2, *z*+1/2; (vi) −*x*, *y*−1/2, −*z*+3/2; (vii) −*x*, *y*+1/2, −*z*+3/2.