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{6,6′-Dimethoxy-2,2′-[cyclohexane-1,2-diylbis(nitrilomethylidyne)]diphenolato-$\kappa^4O^1,N,N',O^{1'}$}iron(II) monohydrate

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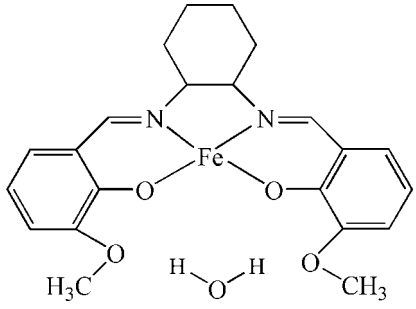
Key indicators: single-crystal X-ray study; T = 291 K; mean σ(C–C) = 0.005 Å; R factor = 0.042; wR factor = 0.107; data-to-parameter ratio = 16.8.

In the title complex, $[Fe(C_{22}H_{24}N_2O_4)]\cdot H_2O$, the Fe^{II} center is four-coordinated by two O and two N atoms from 2,2′-[6,6′-dimethoxycyclohexane-1,2-diylbis(nitrilomethylidyne)]-diphenolate (L) ligands in a distorted square-planar geometry. Uncoordinated water and FeL molecules are paired *via* intermolecular water–methoxy O—H···O hydrogen bonds.

Related literature

For a manganese complex of a similar Schiff base ligand, see: Watkinson *et al.* (1999). For the isotypic Co^{II} compound, see: Bao *et al.* (2009).



Experimental

Crystal data

$[Fe(C_{22}H_{24}N_2O_4)]\cdot H_2O$ M_r = 454.30
Monoclinic, $P2_1/n$
a = 11.243 (5) Å
b = 10.617 (3) Å
c = 17.863 (7) Å
β = 107.042 (14)°
V = 2038.5 (13) Å^3
Z = 4
Mo $K\alpha$ radiation
μ = 0.78 mm^{-1}
T = 291 K
0.22 × 0.21 × 0.18 mm

Data collection

Rigaku R-AXIS RAPID diffractometer
Absorption correction: multi-scan (*ABSCOR*; Higashi, 1995)
T_{min} = 0.846, T_{max} = 0.873
18680 measured reflections
4584 independent reflections
3446 reflections with $I > 2\sigma(I)$
R_{int} = 0.054

Refinement

$R[F^2 > 2\sigma(F^2)]$ = 0.042
$wR(F^2)$ = 0.107
S = 1.03
4584 reflections
273 parameters
H-atom parameters constrained
$\Delta\rho_{max}$ = 0.30 e Å^{-3}
$\Delta\rho_{min}$ = −0.35 e Å^{-3}

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

D—H···A	D—H	H···A	D···A	D—H···A
O5—H2···O2	0.85	2.15	2.915 (4)	149
O5—H1···O4	0.85	2.06	2.898 (4)	169

Data collection: *RAPID-AUTO* (Rigaku, 1998); cell refinement: *RAPID-AUTO*; data reduction: *CrystalStructure* (Rigaku/MSC, 2002); 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: *SHELXL97*.

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Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: CV2567).

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supplementary materials

Acta Cryst. (2009). E**65**, m832 [**doi:10.1107/S1600536809022235**]

{6,6'-Dimethoxy-2,2'-[cyclohexane-1,2-diylbis(nitrilomethylidyne)]diphenolato-$\kappa^4O^1,N,N',O^{1'}$}iron(II) monohydrate

P.-F. Yan, Y. Bao, H.-F. Li and G.-M. Li

Comment

In the title compound (Fig. 1), the Fe^{II} ion is four-coordinated by the tetradentate Schiff base ligand in a square planar environment in a manner observed earlier for a manganese complex (Watkinson *et al.*, 1999). Uncoordinated water molecule is paired with the main molecule by the O—H···O hydrogen bonds (Table 1, Fig. 1).

Experimental

The title complex was obtained by the treatment of anhydrous ferrous chloride with the Schiff base in methanol/acetone (2:3). The yellow clear mixture turned to black precipitation immediately, stirred for 4 h; diethyl ether was allowed to diffuse slowly into the solution of the filtrate. Black single crystals were obtained after several days. Analysis calculated for $C_{22}H_{26}Fe_N2_O5$: C, 58.16; H, 5.77; N, 6.17; Fe, 12.29; found: C, 57.56; H, 5.23; N, 6.77; F, 12.79%.

Refinement

H atoms bound to C atoms were placed in calculated positions and treated as riding on their parent atoms, with C—H = 0.93 Å (aromatic C), C—H = 0.97 Å (methylene C), C—H = 0.98 Å (methine C), and with U_{iso}(H) = 1.2U_{eq}(C) or C—H = 0.96 Å (methly C) and with U_{iso}(H) = 1.5U_{eq}(C). Water H atoms were initially located in a difference Fourier map, but they were treated as riding on their parent atoms with O—H = 0.85 Å and with with U_{iso}(H) = 1.5U_{eq}(O).

Figures

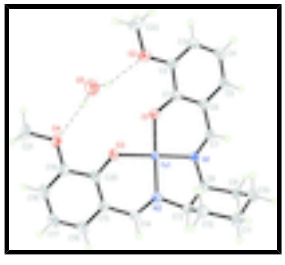


Fig. 1. View of the title compound showing the atomic numbering and 30% probability displacement ellipsoids. Dashed lines indicate the hydrogen-bonding interactions.

{6,6'-Dimethoxy-2,2'-[cyclohexane-1,2-diylbis(nitrilomethylidyne)]diphenolato- $\kappa^4O^1,N,N',O^{1'}$}iron(II) monohydrate

Crystal data

$[Fe(C_{22}H_{24}N_2O_4)]\cdot H_2O$ $F_{000} = 952$

M_r = 454.30	D_x = 1.480 Mg m^{-3}
Monoclinic, $P2_1/n$	Mo $K\alpha$ radiation, λ = 0.71073 Å
Hall symbol: -P 2yn	Cell parameters from 13436 reflections
a = 11.243 (5) Å	θ = 3.1–27.5º
b = 10.617 (3) Å	μ = 0.78 mm^{-1}
c = 17.863 (7) Å	T = 291 K
β = 107.042 (14)º	Block, black
V = 2038.5 (13) Å^3	0.22 × 0.21 × 0.18 mm
Z = 4	

Data collection

Rigaku R-AXIS RAPID diffractometer	4584 independent reflections
Radiation source: fine-focus sealed tube	3446 reflections with $I > 2\sigma(I)$
Monochromator: graphite	R_{int} = 0.054
T = 291 K	θ_{max} = 27.5º
ω scans	θ_{min} = 3.1º
Absorption correction: multi-scan (ABSCOR; Higashi, 1995)	h = −14→14
T_{min} = 0.846, T_{max} = 0.873	k = −13→13
18680 measured reflections	l = −23→23

Refinement

Refinement on F^2	Secondary atom site location: difference Fourier map
Least-squares matrix: full	Hydrogen site location: inferred from neighbouring sites
$R[F^2 > 2\sigma(F^2)]$ = 0.042	H-atom parameters constrained
$wR(F^2)$ = 0.107	$w = 1/[\sigma^2(F_o^2) + (0.0436P)^2 + 1.092P]$ where $P = (F_o^2 + 2F_c^2)/3$
S = 1.03	$(\Delta/\sigma)_{max}$ = 0.001
4584 reflections	$\Delta\rho_{max}$ = 0.30 e Å^{-3}
273 parameters	$\Delta\rho_{min}$ = −0.35 e Å^{-3}
Primary atom site location: structure-invariant direct methods	Extinction correction: none

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}
Fe1	0.65626 (3)	0.46981 (3)	0.162776 (18)	0.02766 (11)
N1	0.7304 (2)	0.52432 (19)	0.08782 (12)	0.0356 (5)
O1	0.55829 (18)	0.34802 (16)	0.09933 (10)	0.0414 (4)
O3	0.57113 (19)	0.43064 (17)	0.23484 (11)	0.0447 (5)
N2	0.7650 (2)	0.5782 (2)	0.23003 (12)	0.0397 (5)
C7	0.6991 (3)	0.4904 (2)	0.01541 (15)	0.0385 (6)
H7	0.7339	0.5350	−0.0178	0.046*
C1	0.5517 (2)	0.3230 (2)	0.02601 (15)	0.0370 (6)
C2	0.4762 (3)	0.2205 (3)	−0.01171 (17)	0.0439 (6)
C20	0.5684 (3)	0.5024 (2)	0.29420 (15)	0.0401 (6)
O4	0.3917 (2)	0.3866 (2)	0.29798 (13)	0.0597 (6)
C19	0.4729 (3)	0.4826 (3)	0.33048 (17)	0.0481 (7)
O2	0.4206 (2)	0.1564 (2)	0.03582 (13)	0.0636 (6)
C8	0.8176 (3)	0.6325 (2)	0.11520 (15)	0.0394 (6)
H8	0.7676	0.7097	0.1075	0.047*
C14	0.7552 (3)	0.6262 (2)	0.29455 (15)	0.0414 (6)
H14	0.8169	0.6814	0.3220	0.050*
C6	0.6156 (2)	0.3903 (2)	−0.01837 (15)	0.0372 (6)
C4	0.5227 (3)	0.2627 (3)	−0.13222 (17)	0.0517 (7)
H4	0.5117	0.2439	−0.1847	0.062*
C15	0.6547 (3)	0.5989 (3)	0.32603 (15)	0.0421 (6)
C5	0.5980 (3)	0.3603 (3)	−0.09773 (16)	0.0464 (7)
H5	0.6381	0.4075	−0.1269	0.056*
C13	0.8726 (3)	0.6129 (2)	0.20281 (16)	0.0411 (6)
H13	0.9091	0.6918	0.2277	0.049*
C12	0.9700 (3)	0.5089 (3)	0.22201 (18)	0.0502 (7)
H12A	0.9306	0.4287	0.2040	0.060*
H12B	1.0067	0.5038	0.2783	0.060*
C3	0.4622 (3)	0.1911 (3)	−0.08882 (17)	0.0495 (7)
H3	0.4125	0.1234	−0.1121	0.059*
C9	0.9182 (3)	0.6509 (3)	0.07480 (17)	0.0486 (7)
H9A	0.8798	0.6520	0.0186	0.058*
H9B	0.9578	0.7319	0.0899	0.058*
C10	1.0165 (3)	0.5485 (3)	0.09490 (19)	0.0594 (8)
H10A	1.0822	0.5690	0.0719	0.071*
H10B	0.9797	0.4692	0.0727	0.071*
C11	1.0719 (3)	0.5339 (4)	0.1833 (2)	0.0648 (9)
H11A	1.1307	0.4646	0.1943	0.078*
H11B	1.1165	0.6101	0.2049	0.078*
C18	0.4656 (3)	0.5554 (3)	0.3928 (2)	0.0633 (9)
H18	0.4018	0.5412	0.4152	0.076*
C21	0.2942 (3)	0.3601 (4)	0.3316 (2)	0.0694 (10)

H21A	0.3291	0.3303	0.3843	0.104*
H21B	0.2405	0.2967	0.3012	0.104*
H21C	0.2472	0.4354	0.3322	0.104*
C17	0.5525 (4)	0.6499 (3)	0.4229 (2)	0.0674 (10)
H17	0.5468	0.6982	0.4651	0.081*
C16	0.6452 (3)	0.6712 (3)	0.39035 (18)	0.0551 (8)
H16	0.7032	0.7342	0.4106	0.066*
O5	0.3096 (3)	0.3010 (3)	0.13694 (18)	0.1061 (11)
H1	0.3424	0.3299	0.1827	0.159*
H2	0.3649	0.2594	0.1238	0.159*
C22	0.3307 (3)	0.0621 (3)	0.0015 (2)	0.0666 (10)
H22A	0.2609	0.1004	−0.0360	0.100*
H22B	0.3032	0.0219	0.0416	0.100*
H22C	0.3674	0.0005	−0.0243	0.100*

Atomic displacement parameters (Å^2)

	U^{11}	U^{22}	U^{33}	U^{12}	U^{13}	U^{23}
Fe1	0.02801 (19)	0.03099 (18)	0.02628 (18)	−0.00228 (14)	0.01153 (13)	−0.00086 (13)
N1	0.0334 (12)	0.0387 (11)	0.0353 (11)	−0.0007 (9)	0.0108 (9)	0.0018 (9)
O1	0.0425 (11)	0.0466 (10)	0.0377 (10)	−0.0046 (8)	0.0157 (8)	−0.0025 (7)
O3	0.0515 (12)	0.0491 (10)	0.0391 (10)	−0.0053 (9)	0.0221 (9)	−0.0030 (8)
N2	0.0406 (13)	0.0432 (12)	0.0382 (12)	−0.0011 (10)	0.0163 (10)	−0.0013 (9)
C7	0.0384 (15)	0.0438 (14)	0.0351 (14)	0.0024 (11)	0.0136 (12)	0.0046 (10)
C1	0.0351 (15)	0.0369 (13)	0.0386 (14)	0.0047 (10)	0.0101 (12)	−0.0025 (10)
C2	0.0398 (16)	0.0464 (15)	0.0470 (16)	0.0015 (12)	0.0148 (13)	−0.0043 (12)
C20	0.0409 (16)	0.0498 (15)	0.0318 (13)	0.0055 (11)	0.0141 (12)	0.0041 (10)
O4	0.0513 (14)	0.0833 (15)	0.0531 (13)	−0.0119 (11)	0.0289 (11)	0.0000 (11)
C19	0.0444 (18)	0.0620 (18)	0.0408 (16)	0.0010 (14)	0.0169 (13)	0.0047 (13)
O2	0.0712 (16)	0.0618 (13)	0.0629 (14)	−0.0281 (11)	0.0277 (12)	−0.0130 (10)
C8	0.0363 (15)	0.0393 (14)	0.0432 (15)	−0.0015 (11)	0.0129 (12)	0.0038 (11)
C14	0.0414 (16)	0.0463 (15)	0.0359 (14)	0.0002 (12)	0.0101 (12)	−0.0035 (11)
C6	0.0342 (14)	0.0418 (14)	0.0349 (13)	0.0054 (11)	0.0086 (11)	−0.0004 (10)
C4	0.0539 (19)	0.0608 (18)	0.0390 (16)	0.0104 (14)	0.0112 (14)	−0.0096 (13)
C15	0.0470 (17)	0.0469 (15)	0.0340 (14)	0.0011 (12)	0.0143 (12)	−0.0002 (11)
C5	0.0456 (18)	0.0559 (17)	0.0383 (15)	0.0046 (13)	0.0134 (13)	−0.0003 (12)
C13	0.0387 (16)	0.0445 (14)	0.0417 (15)	−0.0055 (11)	0.0143 (12)	−0.0025 (11)
C12	0.0397 (17)	0.0653 (19)	0.0439 (16)	0.0075 (13)	0.0098 (13)	0.0074 (13)
C3	0.0473 (18)	0.0480 (16)	0.0498 (17)	0.0019 (13)	0.0086 (14)	−0.0111 (13)
C9	0.0451 (18)	0.0587 (18)	0.0440 (16)	−0.0118 (13)	0.0162 (14)	0.0028 (12)
C10	0.0436 (19)	0.085 (2)	0.056 (2)	0.0016 (16)	0.0244 (16)	0.0000 (16)
C11	0.0369 (18)	0.098 (3)	0.061 (2)	0.0095 (17)	0.0161 (15)	0.0101 (18)
C18	0.057 (2)	0.093 (3)	0.0510 (19)	0.0029 (18)	0.0322 (17)	−0.0024 (16)
C21	0.046 (2)	0.106 (3)	0.065 (2)	−0.0016 (19)	0.0290 (17)	0.0186 (19)
C17	0.077 (3)	0.084 (2)	0.051 (2)	−0.001 (2)	0.0337 (19)	−0.0176 (17)
C16	0.062 (2)	0.0628 (19)	0.0436 (17)	0.0018 (15)	0.0206 (15)	−0.0097 (13)
O5	0.080 (2)	0.167 (3)	0.080 (2)	−0.021 (2)	0.0366 (17)	−0.0236 (19)
C22	0.055 (2)	0.0543 (19)	0.085 (3)	−0.0180 (15)	0.0124 (19)	−0.0009 (16)

Geometric parameters (Å, °)

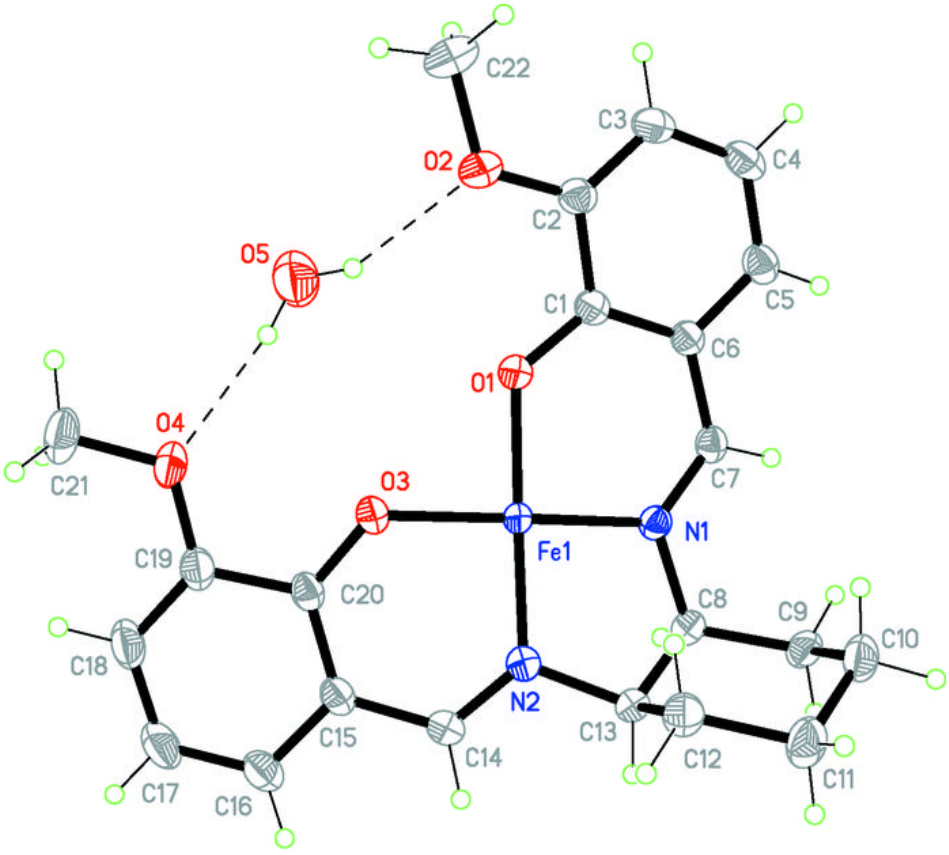
Fe1—N2	1.844 (2)	C4—H4	0.9300
Fe1—O1	1.8541 (18)	C15—C16	1.412 (4)
Fe1—O3	1.8623 (19)	C5—H5	0.9300
Fe1—N1	1.864 (2)	C13—C12	1.523 (4)
N1—C7	1.288 (3)	C13—H13	0.9800
N1—C8	1.496 (3)	C12—C11	1.525 (4)
O1—C1	1.317 (3)	C12—H12A	0.9700
O3—C20	1.313 (3)	C12—H12B	0.9700
N2—C14	1.294 (3)	C3—H3	0.9300
N2—C13	1.476 (3)	C9—C10	1.517 (4)
C7—C6	1.429 (4)	C9—H9A	0.9700
C7—H7	0.9300	C9—H9B	0.9700
C1—C6	1.410 (4)	C10—C11	1.525 (5)
C1—C2	1.423 (4)	C10—H10A	0.9700
C2—O2	1.373 (3)	C10—H10B	0.9700
C2—C3	1.376 (4)	C11—H11A	0.9700
C20—C15	1.411 (4)	C11—H11B	0.9700
C20—C19	1.424 (4)	C18—C17	1.394 (5)
O4—C19	1.378 (4)	C18—H18	0.9300
O4—C21	1.425 (4)	C21—H21A	0.9600
C19—C18	1.377 (4)	C21—H21B	0.9600
O2—C22	1.428 (4)	C21—H21C	0.9600
C8—C13	1.519 (4)	C17—C16	1.352 (5)
C8—C9	1.522 (4)	C17—H17	0.9300
C8—H8	0.9800	C16—H16	0.9300
C14—C15	1.431 (4)	O5—H1	0.8500
C14—H14	0.9300	O5—H2	0.8500
C6—C5	1.409 (4)	C22—H22A	0.9600
C4—C5	1.365 (4)	C22—H22B	0.9600
C4—C3	1.397 (4)	C22—H22C	0.9600
N2—Fe1—O1	174.23 (9)	C8—C13—C12	112.4 (2)
N2—Fe1—O3	93.78 (9)	N2—C13—H13	109.8
O1—Fe1—O3	86.17 (8)	C8—C13—H13	109.8
N2—Fe1—N1	85.62 (10)	C12—C13—H13	109.8
O1—Fe1—N1	95.05 (9)	C13—C12—C11	111.0 (2)
O3—Fe1—N1	173.78 (9)	C13—C12—H12A	109.4
C7—N1—C8	120.2 (2)	C11—C12—H12A	109.4
C7—N1—Fe1	125.91 (19)	C13—C12—H12B	109.4
C8—N1—Fe1	113.05 (16)	C11—C12—H12B	109.4
C1—O1—Fe1	126.75 (17)	H12A—C12—H12B	108.0
C20—O3—Fe1	124.53 (17)	C2—C3—C4	120.0 (3)
C14—N2—C13	119.1 (2)	C2—C3—H3	120.0
C14—N2—Fe1	127.8 (2)	C4—C3—H3	120.0
C13—N2—Fe1	113.07 (16)	C10—C9—C8	112.8 (2)
N1—C7—C6	125.7 (2)	C10—C9—H9A	109.0
N1—C7—H7	117.1	C8—C9—H9A	109.0

C6—C7—H7	117.1	C10—C9—H9B	109.0
O1—C1—C6	124.6 (2)	C8—C9—H9B	109.0
O1—C1—C2	118.5 (2)	H9A—C9—H9B	107.8
C6—C1—C2	116.9 (2)	C9—C10—C11	111.5 (3)
O2—C2—C3	124.5 (3)	C9—C10—H10A	109.3
O2—C2—C1	113.8 (2)	C11—C10—H10A	109.3
C3—C2—C1	121.7 (3)	C9—C10—H10B	109.3
O3—C20—C15	124.8 (3)	C11—C10—H10B	109.3
O3—C20—C19	118.8 (3)	H10A—C10—H10B	108.0
C15—C20—C19	116.4 (2)	C12—C11—C10	110.7 (3)
C19—O4—C21	117.7 (3)	C12—C11—H11A	109.5
C18—C19—O4	124.5 (3)	C10—C11—H11A	109.5
C18—C19—C20	121.2 (3)	C12—C11—H11B	109.5
O4—C19—C20	114.3 (2)	C10—C11—H11B	109.5
C2—O2—C22	118.4 (3)	H11A—C11—H11B	108.1
N1—C8—C13	105.1 (2)	C19—C18—C17	120.9 (3)
N1—C8—C9	116.7 (2)	C19—C18—H18	119.6
C13—C8—C9	111.8 (2)	C17—C18—H18	119.6
N1—C8—H8	107.6	O4—C21—H21A	109.5
C13—C8—H8	107.6	O4—C21—H21B	109.5
C9—C8—H8	107.6	H21A—C21—H21B	109.5
N2—C14—C15	123.6 (3)	O4—C21—H21C	109.5
N2—C14—H14	118.2	H21A—C21—H21C	109.5
C15—C14—H14	118.2	H21B—C21—H21C	109.5
C5—C6—C1	120.4 (2)	C16—C17—C18	119.8 (3)
C5—C6—C7	118.3 (2)	C16—C17—H17	120.1
C1—C6—C7	121.3 (2)	C18—C17—H17	120.1
C5—C4—C3	120.1 (3)	C17—C16—C15	120.7 (3)
C5—C4—H4	120.0	C17—C16—H16	119.6
C3—C4—H4	120.0	C15—C16—H16	119.6
C20—C15—C16	121.0 (3)	H1—O5—H2	107.8
C20—C15—C14	121.1 (2)	O2—C22—H22A	109.5
C16—C15—C14	117.8 (3)	O2—C22—H22B	109.5
C4—C5—C6	120.8 (3)	H22A—C22—H22B	109.5
C4—C5—H5	119.6	O2—C22—H22C	109.5
C6—C5—H5	119.6	H22A—C22—H22C	109.5
N2—C13—C8	104.4 (2)	H22B—C22—H22C	109.5
N2—C13—C12	110.5 (2)		

Hydrogen-bond geometry (Å, °)

D—H···*A*	*D*—H	H···*A*	*D*···*A*	*D*—H···*A*
O5—H2···O1	0.85	2.52	3.102 (4)	126
O5—H2···O2	0.85	2.15	2.915 (4)	149
O5—H1···O3	0.85	2.69	3.254 (4)	125
O5—H1···O4	0.85	2.06	2.898 (4)	169

Fig. 1


C22
C3
C4
O2
C2
O5
C1
C5
C6
O1
O4
C7
C21
O3
Fe1
N1
C19
C20
C8
C9
C10
C18
N2
C13
C15
C14
C12
C11
C17
C16