

# Bis[ $\mu$ -2,2'-[(3-azapentane-1,5-diyl)-bis(nitrilomethylidene)]diphenolato}dicopper(II) dimethyl sulfoxide disolvate

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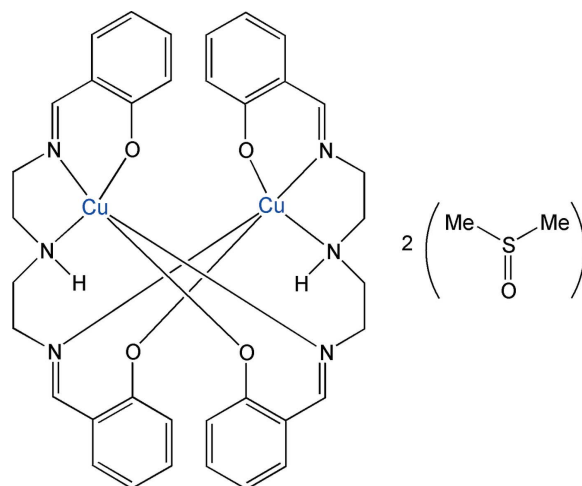
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Key indicators: single-crystal X-ray study;  $T = 298$  K; mean  $\sigma(\text{C}-\text{C}) = 0.009$  Å; disorder in solvent or counterion;  $R$  factor = 0.062;  $wR$  factor = 0.167; data-to-parameter ratio = 12.2.

The title compound,  $[\text{Cu}_2(\text{C}_{18}\text{H}_{19}\text{N}_3\text{O}_2)_2] \cdot 2\text{C}_2\text{H}_6\text{OS}$  or  $[\text{Cu}_2(\text{SalenN}_3\text{H})_2] \cdot 2\text{DMSO}$ , where  $\text{SalenN}_3\text{H}$  is the multi-dentate Schiff base 2,2'-[(3-azapentane-1,5-diyl)bis(nitrilomethylidene)]diphenolate dianion and DMSO is dimethyl sulfoxide, is a solvated dinuclear  $\text{Cu}^{\text{II}}$  complex. The neutral complex is built from two  $\text{Cu}(\text{SalenN}_3\text{H})$  units related by an inversion center. All heteroatoms in the Schiff bases coordinate the  $\text{Cu}^{\text{II}}$  ions, which display highly distorted trigonal bipyramidal geometries. The solvent molecules are located in the structural voids of the complex and are disordered over two positions with occupancies of 0.642 (15) and 0.358 (15). The previously characterized acetone disolvate of the same complex presents identical molecular and crystal structures, and crystallizes with cell parameters very close to those of the DMSO disolvate reported here.

## Related literature

The title compound was synthesized by direct synthesis, using metallic copper as starting material (Gutiérrez *et al.*, 2001; Reyes-Ortega *et al.*, 2005). The same dinuclear  $\text{Cu}^{\text{II}}$  complex was previously characterized with acetone solvent in place of DMSO (McKenzie & Selvey, 1985).



## Experimental

### Crystal data

$[\text{Cu}_2(\text{C}_{18}\text{H}_{19}\text{N}_3\text{O}_2)_2] \cdot 2\text{C}_2\text{H}_6\text{OS}$   
 $M_r = 902.06$   
 Monoclinic,  $P2_1/c$   
 $a = 12.817$  (3) Å  
 $b = 16.783$  (4) Å  
 $c = 9.827$  (3) Å  
 $\beta = 106.732$  (18)°

$V = 2024.5$  (8) Å<sup>3</sup>  
 $Z = 2$   
 Mo  $K\alpha$  radiation  
 $\mu = 1.21$  mm<sup>-1</sup>  
 $T = 298$  (1) K  
 $0.16 \times 0.16 \times 0.12$  mm

### Data collection

Bruker P4 diffractometer  
 Absorption correction:  $\psi$  scan  
 (XSCANS; Siemens, 1996)  
 $T_{\text{min}} = 0.803$ ,  $T_{\text{max}} = 0.866$   
 7380 measured reflections  
 3599 independent reflections

2191 reflections with  $I > 2\sigma(I)$   
 $R_{\text{int}} = 0.076$   
 1 standard reflections  
 every 48 reflections  
 intensity decay: 1%

### Refinement

$R[F^2 > 2\sigma(F^2)] = 0.062$   
 $wR(F^2) = 0.167$   
 $S = 1.05$   
 3599 reflections  
 294 parameters

30 restraints  
 H-atom parameters constrained  
 $\Delta\rho_{\text{max}} = 0.47$  e Å<sup>-3</sup>  
 $\Delta\rho_{\text{min}} = -0.82$  e Å<sup>-3</sup>

**Table 1**

Selected geometric parameters (Å, °).

Cu1—O1	1.932 (4)	Cu1—N3	1.969 (5)
Cu1—N1	1.948 (5)	Cu1—O2	1.978 (4)
Cu1—N2	2.319 (4)		
O1—Cu1—O2	137.2 (2)	N2—Cu1—O1	131.28 (19)
O2—Cu1—N2	90.92 (17)	N1—Cu1—N3	176.4 (2)

Data collection: XSCANS (Siemens, 1996); cell refinement: XSCANS; data reduction: XSCANS; program(s) used to solve structure: SHELXTL-Plus (Sheldrick, 2008); program(s) used to refine structure: SHELXTL-Plus; molecular graphics: Mercury (Macrae *et al.*, 2006); software used to prepare material for publication: SHELXTL-Plus.

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Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: SI2079).

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## References

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## supporting information

*Acta Cryst.* (2008). E64, m631–m632 [doi:10.1107/S1600536808008544]

## Bis{ $\mu$ -2,2'-[(3-azapentane-1,5-diyl)bis(nitrilomethylidene)]diphenolato}dicopper(II) dimethyl sulfoxide disolvate

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### S1. Comment

The title compound was obtained during a general study about direct synthesis in coordination chemistry, *i.e.* effective synthesis of complexes using zero-valent metals, neutral ligands, and polar solvents as starting materials (Gutiérrez *et al.*, 2001). We expect that new molecular arrangements, as well as new compounds including solvent molecules as ligands, may be achieved using such reactions (Reyes-Ortega *et al.*, 2005). We are also interested in the influence of lattice solvents on magnetic properties of paramagnetic coordination compounds.

The title compound is a dinuclear centrosymmetric Cu<sup>II</sup> complex, formed through coordination of two Schiff base dianions SalenN<sub>3</sub>H to two Cu<sup>II</sup> ions; the complex is solvated by two DMSO molecules (Fig. 1). In the complex, all heteroatoms of the Schiff base ligands are coordinated to the metallic centers, which present a highly distorted trigonal-bipyramidal geometry. Atoms O1, O2 and N2 form the equatorial plane, while atoms N1 and N3 occupy apical positions, with an angle N1—Cu1—N3 = 176.4 (2)°. DMSO molecules are located in the structural voids of the complex (Fig. 2), and poorly interact with the Schiff bases, as reflected in the disorder found for this molecule (Fig. 1, inset).

The whole complex presents a rigid conformation, as ten coordination bonds are formed. It may thus be expected to be a good candidate for hosting small solvent molecules with a steric volume similar to that of DMSO. This hypothesis is, at least partially, confirmed by the previous X-ray characterization of the acetone disolvate of the same complex (McKenzie & Selvey, 1985). This compound crystallizes in the same space group, with cell parameters very close to those of the title disolvate. The complex conformation is identical, regardless of the solvent inserted in voids. For example, the non-bonding Cu...Cu separation is virtually not modified: 5.7716 (18) Å in the DMSO disolvate, *vs.* 5.809 Å in the acetone disolvate.

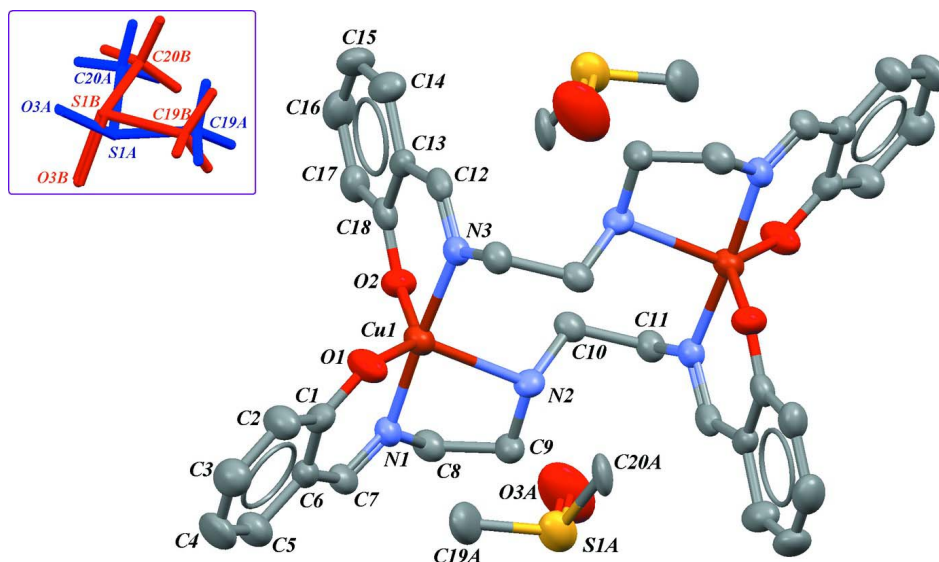
### S2. Experimental

The title compound was prepared by direct synthesis, mixing equimolecular amounts (0.8 mmol) of elemental copper and neutral Schiff base SalenN<sub>3</sub>H<sub>3</sub> in DMSO (2.4 ml). The mixture was heated at 353 K with magnetic stirring for 5.5 hrs, and then filtered. A crystalline compound, was collected after six days. Yield: 30%.

### S3. Refinement

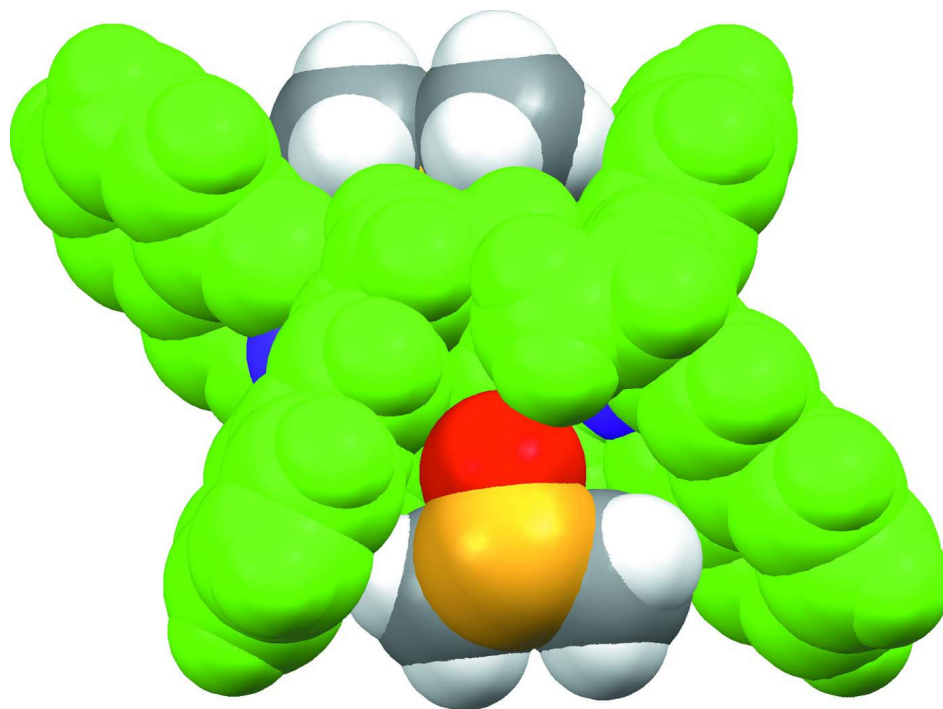
The asymmetric unit contains one DMSO molecule which is clearly disordered over two positions (Fig. 1, inset). S, O and C atoms were splitted over two sites. Refined occupancies converged to 0.642 (15) and 0.358 (15) for sites A and B, respectively. Geometry was regularized through restraints applied to bond lengths: S—O = 1.475 (20) and S—C = 1.750 (20) Å. Finally, sites in each pair of disordered atoms were restrained, with a standard deviation of 0.04 Å<sup>2</sup>, to have the same  $U_{ij}$  components. All H atoms were placed in idealized positions, and were allowed to ride on their carrier atoms,

with C—H bond lengths fixed to 0.93 (aromatic CH), 0.97 (methylene CH<sub>2</sub>) or 0.96 Å (methyl CH<sub>3</sub>), and N—H bond length fixed to 0.91 Å. Isotropic displacement parameters for H atoms were calculated as  $U_{\text{iso}}(\text{H}) = xU_{\text{eq}}(\text{carrier atom})$  where  $x = 1.5$  for methyl groups and  $x = 1.2$  otherwise.



**Figure 1**

The molecular structure of the title compound. Displacement ellipsoids are given at the 40% probability level and all H atoms have been omitted. A single position for DMSO molecules has been retained, which corresponds to the main site occupancy factors, 0.642 (15). The inset represents the model refined for the disordered DMSO molecule. Site occupancy factors are 0.642 (15) and 0.358 (15), for the blue and red molecule, respectively. In the main figure, non labelled atoms are generated with symmetry code  $1 - x, 1 - y, 1 - z$ .

**Figure 2**

A spacefill model for the title compound. All atoms are represented, excepted less occupied disordered sites for DMSO molecules. Colours code: purple: Cu; green: Schiff bases; other: DMSO.

**Bis{ $\mu$ -2,2'-[(3-azapentane-1,5-diy)]bis(nitrilomethylidene)]diphenolato}dicopper(II) dimethyl sulfoxide disolvate**

*Crystal data*

$[\text{Cu}_2(\text{C}_{18}\text{H}_{19}\text{N}_3\text{O}_2)_2] \cdot 2\text{C}_2\text{H}_6\text{OS}$

$M_r = 902.06$

Monoclinic,  $P2_1/c$

Hall symbol: -P 2ybc

$a = 12.817(3) \text{ \AA}$

$b = 16.783(4) \text{ \AA}$

$c = 9.827(3) \text{ \AA}$

$\beta = 106.732(18)^\circ$

$V = 2024.5(8) \text{ \AA}^3$

$Z = 2$

$F(000) = 940$

$D_x = 1.480 \text{ Mg m}^{-3}$

Melting point: 410 K

Mo  $K\alpha$  radiation,  $\lambda = 0.71073 \text{ \AA}$

Cell parameters from 72 reflections

$\theta = 4.8\text{--}10.8^\circ$

$\mu = 1.21 \text{ mm}^{-1}$

$T = 298 \text{ K}$

Irregular, green

$0.16 \times 0.16 \times 0.12 \text{ mm}$

*Data collection*

Bruker P4

diffractometer

Radiation source: fine-focus sealed tube

Graphite monochromator

$2\theta/\omega$  scans

Absorption correction:  $\psi$  scan

(*XSCANS*; Siemens, 1996)

$T_{\min} = 0.803$ ,  $T_{\max} = 0.866$

7380 measured reflections

3599 independent reflections

2191 reflections with  $I > 2\sigma(I)$

$R_{\text{int}} = 0.076$

$\theta_{\max} = 25.1^\circ$ ,  $\theta_{\min} = 2.1^\circ$

$h = -14 \rightarrow 15$

$k = -20 \rightarrow 20$

$l = -11 \rightarrow 7$

1 standard reflections every 48 reflections

intensity decay: 1%

Refinement

Refinement on  $F^2$

Least-squares matrix: full

$R[F^2 > 2\sigma(F^2)] = 0.062$

$wR(F^2) = 0.167$

$S = 1.05$

3599 reflections

294 parameters

30 restraints

0 constraints

Primary atom site location: structure-invariant  
direct methods

Secondary atom site location: difference Fourier  
map

Hydrogen site location: inferred from  
neighbouring sites

H-atom parameters constrained

$w = 1/[\sigma^2(F_o^2) + (0.0556P)^2 + 5.336P]$

where  $P = (F_o^2 + 2F_c^2)/3$

$(\Delta/\sigma)_{\max} < 0.001$

$\Delta\rho_{\max} = 0.47 \text{ e } \text{\AA}^{-3}$

$\Delta\rho_{\min} = -0.82 \text{ e } \text{\AA}^{-3}$

Fractional atomic coordinates and isotropic or equivalent isotropic displacement parameters ( $\text{\AA}^2$ )

	<i>x</i>	<i>y</i>	<i>z</i>	$U_{\text{iso}}^*/U_{\text{eq}}$	Occ. (<1)
Cu1	0.59573 (6)	0.38379 (4)	0.71636 (8)	0.0438 (3)	
O1	0.7294 (3)	0.4094 (3)	0.8606 (5)	0.0597 (12)	
O2	0.4577 (3)	0.3281 (2)	0.7063 (4)	0.0487 (11)	
N1	0.6595 (4)	0.2864 (3)	0.6677 (5)	0.0394 (11)	
N2	0.5491 (4)	0.3917 (3)	0.4705 (5)	0.0391 (11)	
H2B	0.5898	0.4311	0.4478	0.047*	
N3	0.5343 (4)	0.4857 (3)	0.7562 (5)	0.0386 (11)	
C1	0.8131 (5)	0.3647 (4)	0.9119 (6)	0.0460 (15)	
C2	0.8969 (5)	0.3907 (4)	1.0277 (7)	0.0609 (18)	
H2A	0.8908	0.4401	1.0677	0.073*	
C3	0.9878 (6)	0.3459 (5)	1.0844 (8)	0.067 (2)	
H3A	1.0419	0.3655	1.1620	0.080*	
C4	1.0014 (5)	0.2734 (5)	1.0301 (8)	0.073 (2)	
H4A	1.0636	0.2432	1.0704	0.088*	
C5	0.9219 (5)	0.2457 (4)	0.9153 (8)	0.0605 (18)	
H5A	0.9312	0.1966	0.8763	0.073*	
C6	0.8269 (5)	0.2894 (4)	0.8548 (6)	0.0427 (14)	
C7	0.7516 (5)	0.2566 (3)	0.7316 (6)	0.0433 (14)	
H7A	0.7715	0.2097	0.6951	0.052*	
C8	0.5930 (5)	0.2526 (4)	0.5346 (6)	0.0476 (15)	
H8A	0.5215	0.2383	0.5420	0.057*	
H8B	0.6272	0.2052	0.5104	0.057*	
C9	0.5838 (5)	0.3163 (3)	0.4234 (6)	0.0448 (15)	
H9A	0.6538	0.3233	0.4056	0.054*	
H9B	0.5315	0.2998	0.3353	0.054*	
C10	0.4349 (4)	0.4115 (3)	0.4060 (6)	0.0408 (14)	
H10A	0.4141	0.4527	0.4624	0.049*	
H10B	0.3910	0.3647	0.4090	0.049*	
C11	0.4088 (5)	0.4401 (3)	0.2531 (6)	0.0423 (14)	
H11A	0.4302	0.3996	0.1959	0.051*	
H11B	0.3309	0.4483	0.2153	0.051*	
C12	0.4532 (5)	0.4911 (4)	0.8083 (6)	0.0432 (15)	
H12A	0.4395	0.5411	0.8401	0.052*	

C13	0.3816 (4)	0.4273 (4)	0.8226 (6)	0.0401 (14)	
C14	0.2997 (5)	0.4461 (5)	0.8867 (7)	0.0600 (19)	
H14A	0.2965	0.4973	0.9215	0.072*	
C15	0.2255 (5)	0.3910 (5)	0.8987 (8)	0.0622 (19)	
H15A	0.1722	0.4041	0.9422	0.075*	
C16	0.2297 (6)	0.3161 (5)	0.8466 (8)	0.065 (2)	
H16A	0.1796	0.2782	0.8569	0.078*	
C17	0.3054 (5)	0.2953 (4)	0.7795 (7)	0.0532 (17)	
H17A	0.3031	0.2447	0.7403	0.064*	
C18	0.3875 (5)	0.3496 (3)	0.7685 (6)	0.0415 (14)	
S1A	0.8609 (5)	0.4752 (5)	0.4851 (8)	0.068 (2)	0.642 (15)
O3A	0.7492 (8)	0.4440 (7)	0.4334 (14)	0.126 (5)	0.642 (15)
C19A	0.921 (3)	0.449 (2)	0.665 (2)	0.084 (8)	0.642 (15)
H19A	0.9266	0.3926	0.6738	0.126*	0.642 (15)
H19B	0.8767	0.4693	0.7213	0.126*	0.642 (15)
H19C	0.9924	0.4727	0.6974	0.126*	0.642 (15)
C20A	0.845 (2)	0.5781 (8)	0.505 (2)	0.054 (5)	0.642 (15)
H20A	0.8109	0.6014	0.4137	0.081*	0.642 (15)
H20B	0.9154	0.6021	0.5447	0.081*	0.642 (15)
H20C	0.8007	0.5871	0.5670	0.081*	0.642 (15)
S1B	0.8199 (9)	0.4739 (10)	0.4918 (19)	0.082 (4)	0.358 (15)
O3B	0.8755 (13)	0.4476 (9)	0.386 (2)	0.089 (7)	0.358 (15)
C19B	0.901 (5)	0.431 (4)	0.650 (6)	0.12 (2)	0.358 (15)
H19D	0.8815	0.3758	0.6532	0.182*	0.358 (15)
H19E	0.8884	0.4584	0.7299	0.182*	0.358 (15)
H19F	0.9761	0.4351	0.6538	0.182*	0.358 (15)
C20B	0.852 (5)	0.5716 (17)	0.550 (5)	0.107 (18)	0.358 (15)
H20D	0.8245	0.6078	0.4726	0.161*	0.358 (15)
H20E	0.9299	0.5772	0.5855	0.161*	0.358 (15)
H20F	0.8200	0.5835	0.6249	0.161*	0.358 (15)

Atomic displacement parameters ( $\text{\AA}^2$ )

	$U^{11}$	$U^{22}$	$U^{33}$	$U^{12}$	$U^{13}$	$U^{23}$
Cu1	0.0420 (4)	0.0405 (4)	0.0467 (5)	0.0053 (3)	0.0092 (3)	-0.0004 (4)
O1	0.050 (3)	0.062 (3)	0.056 (3)	0.015 (2)	-0.003 (2)	-0.018 (2)
O2	0.047 (2)	0.046 (2)	0.058 (3)	0.003 (2)	0.023 (2)	-0.002 (2)
N1	0.042 (3)	0.036 (3)	0.039 (3)	0.005 (2)	0.010 (2)	0.005 (2)
N2	0.046 (3)	0.036 (3)	0.037 (3)	0.000 (2)	0.014 (2)	-0.008 (2)
N3	0.038 (3)	0.040 (3)	0.035 (3)	0.004 (2)	0.007 (2)	0.003 (2)
C1	0.045 (3)	0.063 (4)	0.031 (3)	0.005 (3)	0.012 (3)	-0.001 (3)
C2	0.056 (4)	0.071 (5)	0.048 (4)	0.013 (4)	0.004 (3)	-0.010 (4)
C3	0.058 (4)	0.084 (6)	0.054 (4)	0.004 (4)	0.007 (4)	-0.004 (4)
C4	0.048 (4)	0.094 (6)	0.066 (5)	0.029 (4)	-0.001 (4)	0.010 (5)
C5	0.050 (4)	0.062 (4)	0.066 (5)	0.012 (3)	0.012 (4)	0.006 (4)
C6	0.041 (3)	0.051 (4)	0.038 (3)	0.009 (3)	0.015 (3)	0.009 (3)
C7	0.047 (3)	0.031 (3)	0.053 (4)	0.005 (3)	0.017 (3)	0.006 (3)
C8	0.045 (3)	0.045 (4)	0.049 (4)	0.005 (3)	0.007 (3)	-0.004 (3)

C9	0.048 (3)	0.049 (4)	0.036 (3)	0.009 (3)	0.010 (3)	-0.001 (3)
C10	0.040 (3)	0.033 (3)	0.049 (4)	0.000 (2)	0.011 (3)	-0.003 (3)
C11	0.044 (3)	0.039 (3)	0.041 (4)	0.001 (3)	0.007 (3)	0.000 (3)
C12	0.053 (4)	0.042 (3)	0.027 (3)	0.009 (3)	-0.001 (3)	-0.001 (3)
C13	0.034 (3)	0.057 (4)	0.029 (3)	0.008 (3)	0.008 (2)	0.007 (3)
C14	0.051 (4)	0.084 (5)	0.045 (4)	0.008 (4)	0.014 (3)	-0.010 (4)
C15	0.050 (4)	0.079 (5)	0.069 (5)	0.010 (4)	0.035 (4)	0.007 (4)
C16	0.055 (4)	0.077 (5)	0.064 (5)	-0.006 (4)	0.020 (4)	0.021 (4)
C17	0.050 (4)	0.055 (4)	0.051 (4)	0.000 (3)	0.010 (3)	0.014 (3)
C18	0.046 (3)	0.037 (3)	0.034 (3)	0.007 (3)	0.001 (3)	0.006 (3)
S1A	0.046 (3)	0.070 (3)	0.085 (4)	-0.009 (3)	0.014 (3)	0.012 (2)
O3A	0.057 (6)	0.100 (8)	0.201 (13)	-0.036 (6)	0.005 (7)	0.026 (8)
C19A	0.098 (16)	0.063 (14)	0.105 (14)	0.017 (15)	0.050 (11)	0.046 (11)
C20A	0.056 (8)	0.052 (7)	0.044 (12)	0.003 (6)	-0.002 (8)	0.028 (6)
S1B	0.052 (9)	0.069 (5)	0.128 (7)	-0.018 (7)	0.032 (8)	0.008 (4)
O3B	0.113 (14)	0.050 (9)	0.103 (15)	0.007 (9)	0.029 (11)	-0.011 (10)
C19B	0.12 (3)	0.11 (4)	0.16 (4)	-0.05 (2)	0.07 (3)	0.04 (3)
C20B	0.11 (3)	0.15 (3)	0.06 (3)	0.00 (2)	0.02 (2)	0.04 (2)

*Geometric parameters (Å, °)*

Cu1—O1	1.932 (4)	C10—H10B	0.9700
Cu1—N1	1.948 (5)	C11—N3 <sup>i</sup>	1.458 (7)
Cu1—N2	2.319 (4)	C11—H11A	0.9700
Cu1—N3	1.969 (5)	C11—H11B	0.9700
Cu1—O2	1.978 (4)	C12—C13	1.444 (8)
Cu1—Cu1 <sup>i</sup>	5.7716 (18)	C12—H12A	0.9300
O1—C1	1.286 (7)	C13—C14	1.407 (8)
O2—C18	1.277 (7)	C13—C18	1.418 (8)
N1—C7	1.270 (7)	C14—C15	1.357 (9)
N1—C8	1.455 (7)	C14—H14A	0.9300
N2—C10	1.456 (7)	C15—C16	1.363 (10)
N2—C9	1.460 (7)	C15—H15A	0.9300
N2—H2B	0.9100	C16—C17	1.367 (9)
N3—C12	1.288 (7)	C16—H16A	0.9300
N3—C11 <sup>i</sup>	1.458 (7)	C17—C18	1.418 (8)
C1—C2	1.392 (8)	C17—H17A	0.9300
C1—C6	1.413 (8)	S1A—O3A	1.471 (10)
C2—C3	1.363 (9)	S1A—C20A	1.756 (13)
C2—H2A	0.9300	S1A—C19A	1.766 (15)
C3—C4	1.361 (10)	C19A—H19A	0.9600
C3—H3A	0.9300	C19A—H19B	0.9600
C4—C5	1.366 (9)	C19A—H19C	0.9600
C4—H4A	0.9300	C20A—H20A	0.9600
C5—C6	1.400 (8)	C20A—H20B	0.9600
C5—H5A	0.9300	C20A—H20C	0.9600
C6—C7	1.424 (8)	S1B—O3B	1.483 (16)
C7—H7A	0.9300	S1B—C20B	1.75 (2)



C8—C9	1.509 (8)	S1B—C19B	1.76 (2)
C8—H8A	0.9700	C19B—H19D	0.9600
C8—H8B	0.9700	C19B—H19E	0.9600
C9—H9A	0.9700	C19B—H19F	0.9600
C9—H9B	0.9700	C20B—H20D	0.9600
C10—C11	1.521 (8)	C20B—H20E	0.9600
C10—H10A	0.9700	C20B—H20F	0.9600
O1—Cu1—O2	137.2 (2)	N2—C9—H9A	109.5
O2—Cu1—N2	90.92 (17)	C8—C9—H9A	109.5
N2—Cu1—O1	131.28 (19)	N2—C9—H9B	109.5
N1—Cu1—N3	176.4 (2)	C8—C9—H9B	109.5
O1—Cu1—N1	91.16 (18)	H9A—C9—H9B	108.1
O1—Cu1—N3	88.98 (18)	N2—C10—C11	114.2 (5)
N1—Cu1—O2	91.38 (18)	N2—C10—H10A	108.7
N3—Cu1—O2	91.00 (18)	C11—C10—H10A	108.7
N1—Cu1—N2	78.03 (17)	N2—C10—H10B	108.7
N3—Cu1—N2	99.17 (17)	C11—C10—H10B	108.7
O1—Cu1—Cu1 <sup>i</sup>	119.32 (15)	H10A—C10—H10B	107.6
N1—Cu1—Cu1 <sup>i</sup>	120.37 (14)	N3 <sup>i</sup> —C11—C10	111.1 (5)
N3—Cu1—Cu1 <sup>i</sup>	56.61 (13)	N3 <sup>i</sup> —C11—H11A	109.4
O2—Cu1—Cu1 <sup>i</sup>	95.61 (12)	C10—C11—H11A	109.4
N2—Cu1—Cu1 <sup>i</sup>	42.83 (11)	N3 <sup>i</sup> —C11—H11B	109.4
C1—O1—Cu1	128.4 (4)	C10—C11—H11B	109.4
C18—O2—Cu1	125.9 (4)	H11A—C11—H11B	108.0
C7—N1—C8	120.8 (5)	N3—C12—C13	126.6 (5)
C7—N1—Cu1	127.3 (4)	N3—C12—H12A	116.7
C8—N1—Cu1	111.7 (3)	C13—C12—H12A	116.7
C10—N2—C9	114.8 (4)	C14—C13—C18	120.1 (6)
C10—N2—Cu1	113.1 (3)	C14—C13—C12	117.0 (6)
C9—N2—Cu1	105.7 (3)	C18—C13—C12	122.8 (5)
C10—N2—H2B	107.7	C15—C14—C13	121.2 (7)
C9—N2—H2B	107.7	C15—C14—H14A	119.4
Cu1—N2—H2B	107.7	C13—C14—H14A	119.4
C12—N3—C11 <sup>i</sup>	116.1 (5)	C14—C15—C16	119.4 (6)
C12—N3—Cu1	123.7 (4)	C14—C15—H15A	120.3
C11 <sup>i</sup> —N3—Cu1	119.6 (4)	C16—C15—H15A	120.3
O1—C1—C2	119.9 (6)	C15—C16—C17	121.8 (7)
O1—C1—C6	123.4 (5)	C15—C16—H16A	119.1
C2—C1—C6	116.7 (6)	C17—C16—H16A	119.1
C3—C2—C1	121.9 (7)	C16—C17—C18	121.2 (7)
C3—C2—H2A	119.0	C16—C17—H17A	119.4
C1—C2—H2A	119.0	C18—C17—H17A	119.4
C4—C3—C2	121.6 (7)	O2—C18—C17	119.7 (6)
C4—C3—H3A	119.2	O2—C18—C13	124.0 (6)
C2—C3—H3A	119.2	C17—C18—C13	116.2 (6)
C3—C4—C5	118.6 (6)	O3A—S1A—C20A	104.8 (10)
C3—C4—H4A	120.7	O3A—S1A—C19A	111.1 (14)

C5—C4—H4A	120.7	C20A—S1A—C19A	99.2 (14)
C4—C5—C6	121.6 (7)	O3B—S1B—C20B	113 (2)
C4—C5—H5A	119.2	O3B—S1B—C19B	103 (3)
C6—C5—H5A	119.2	C20B—S1B—C19B	94 (3)
C5—C6—C1	119.5 (6)	S1B—C19B—H19D	109.5
C5—C6—C7	116.8 (6)	S1B—C19B—H19E	109.5
C1—C6—C7	123.5 (5)	H19D—C19B—H19E	109.5
N1—C7—C6	125.0 (5)	S1B—C19B—H19F	109.5
N1—C7—H7A	117.5	H19D—C19B—H19F	109.5
C6—C7—H7A	117.5	H19E—C19B—H19F	109.5
N1—C8—C9	106.1 (5)	S1B—C20B—H20D	109.5
N1—C8—H8A	110.5	S1B—C20B—H20E	109.5
C9—C8—H8A	110.5	H20D—C20B—H20E	109.5
N1—C8—H8B	110.5	S1B—C20B—H20F	109.5
C9—C8—H8B	110.5	H20D—C20B—H20F	109.5
H8A—C8—H8B	108.7	H20E—C20B—H20F	109.5
N2—C9—C8	110.6 (5)		
N1—Cu1—O1—C1	-11.8 (5)	C6—C1—C2—C3	-0.4 (10)
N3—Cu1—O1—C1	171.8 (5)	C1—C2—C3—C4	0.2 (12)
O2—Cu1—O1—C1	81.5 (6)	C2—C3—C4—C5	0.6 (12)
N2—Cu1—O1—C1	-86.8 (6)	C3—C4—C5—C6	-1.2 (11)
Cu1 <sup>i</sup> —Cu1—O1—C1	-138.1 (5)	C4—C5—C6—C1	1.1 (10)
O1—Cu1—O2—C18	63.5 (5)	C4—C5—C6—C7	177.3 (6)
N1—Cu1—O2—C18	156.7 (5)	O1—C1—C6—C5	178.3 (6)
N3—Cu1—O2—C18	-26.1 (5)	C2—C1—C6—C5	-0.3 (9)
N2—Cu1—O2—C18	-125.3 (4)	O1—C1—C6—C7	2.4 (9)
Cu1 <sup>i</sup> —Cu1—O2—C18	-82.6 (4)	C2—C1—C6—C7	-176.2 (6)
O1—Cu1—N1—C7	6.9 (5)	C8—N1—C7—C6	174.0 (5)
O2—Cu1—N1—C7	-130.3 (5)	Cu1—N1—C7—C6	0.6 (9)
N2—Cu1—N1—C7	139.0 (5)	C5—C6—C7—N1	176.5 (6)
Cu1 <sup>i</sup> —Cu1—N1—C7	132.3 (5)	C1—C6—C7—N1	-7.4 (9)
O1—Cu1—N1—C8	-167.0 (4)	C7—N1—C8—C9	-116.1 (6)
O2—Cu1—N1—C8	55.8 (4)	Cu1—N1—C8—C9	58.2 (5)
N2—Cu1—N1—C8	-34.9 (4)	C10—N2—C9—C8	-102.8 (6)
Cu1 <sup>i</sup> —Cu1—N1—C8	-41.6 (4)	Cu1—N2—C9—C8	22.5 (5)
O1—Cu1—N2—C10	-146.8 (3)	N1—C8—C9—N2	-50.9 (6)
N1—Cu1—N2—C10	132.3 (4)	C9—N2—C10—C11	-75.2 (6)
N3—Cu1—N2—C10	-50.0 (4)	Cu1—N2—C10—C11	163.4 (4)
O2—Cu1—N2—C10	41.1 (4)	N2—C10—C11—N3 <sup>i</sup>	-63.2 (6)
Cu1 <sup>i</sup> —Cu1—N2—C10	-56.2 (3)	C11 <sup>i</sup> —N3—C12—C13	177.1 (5)
O1—Cu1—N2—C9	86.8 (4)	Cu1—N3—C12—C13	-12.0 (8)
N1—Cu1—N2—C9	6.0 (4)	N3—C12—C13—C14	178.2 (5)
N3—Cu1—N2—C9	-176.4 (3)	N3—C12—C13—C18	-5.2 (9)
O2—Cu1—N2—C9	-85.3 (4)	C18—C13—C14—C15	0.2 (9)
Cu1 <sup>i</sup> —Cu1—N2—C9	177.5 (4)	C12—C13—C14—C15	176.9 (6)
O1—Cu1—N3—C12	-114.8 (5)	C13—C14—C15—C16	-0.6 (10)
O2—Cu1—N3—C12	22.3 (5)	C14—C15—C16—C17	-1.4 (11)

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N2—Cu1—N3—C12	113.4 (4)	C15—C16—C17—C18	3.8 (10)
Cu1 <sup>i</sup> —Cu1—N3—C12	118.4 (5)	Cu1—O2—C18—C17	-165.4 (4)
O1—Cu1—N3—C11 <sup>i</sup>	55.8 (4)	Cu1—O2—C18—C13	18.2 (8)
O2—Cu1—N3—C11 <sup>i</sup>	-167.0 (4)	C16—C17—C18—O2	179.4 (6)
N2—Cu1—N3—C11 <sup>i</sup>	-75.9 (4)	C16—C17—C18—C13	-3.9 (9)
Cu1 <sup>i</sup> —Cu1—N3—C11 <sup>i</sup>	-71.0 (4)	C14—C13—C18—O2	178.5 (5)
Cu1—O1—C1—C2	-172.4 (5)	C12—C13—C18—O2	2.0 (8)
Cu1—O1—C1—C6	9.1 (9)	C14—C13—C18—C17	2.0 (8)
O1—C1—C2—C3	-179.0 (6)	C12—C13—C18—C17	-174.6 (5)

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Symmetry code: (i)  $-x+1, -y+1, -z+1$ .