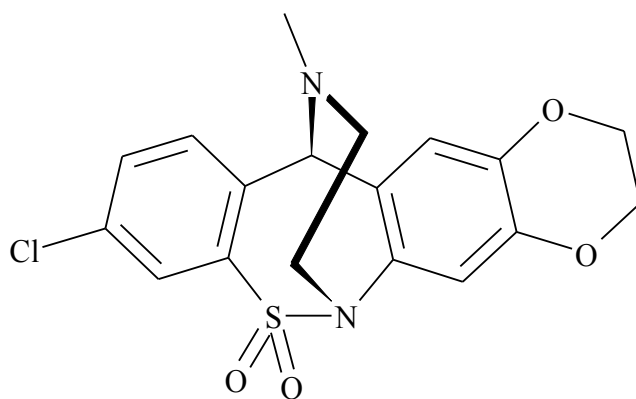


85623

2366-BGE

Submitted by: Berecz Gabor
Operator: Dancso Andras

X-ray Structure Report



April 1, 2010



Fig. 1. The crystal

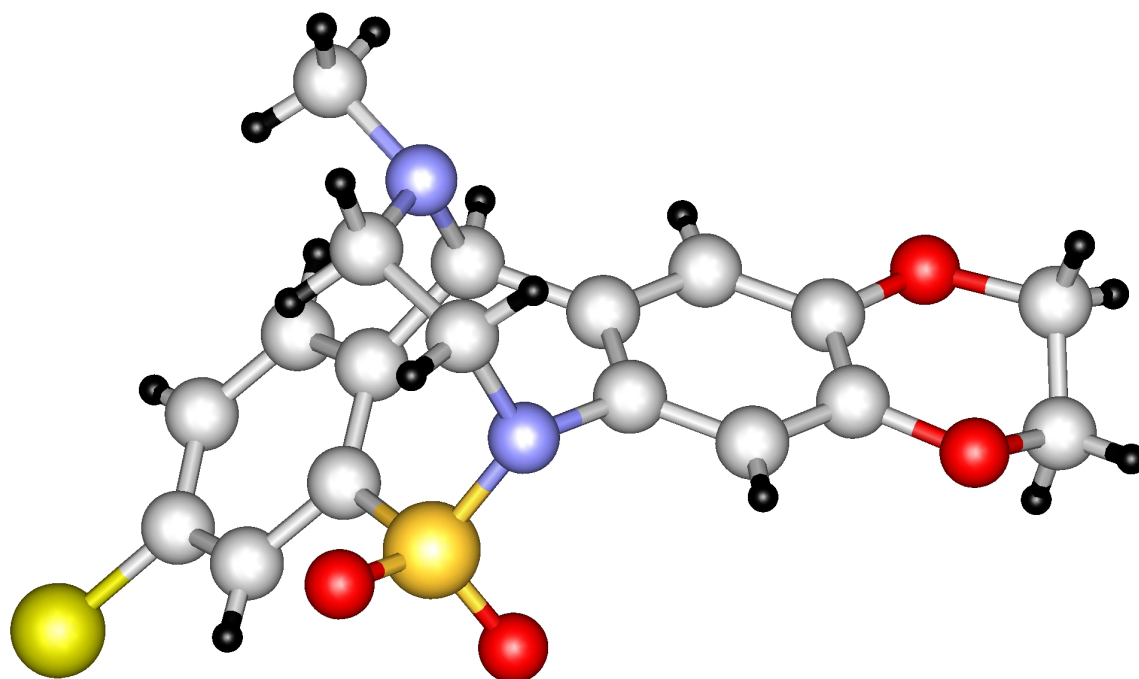


Fig. 2. The molecule

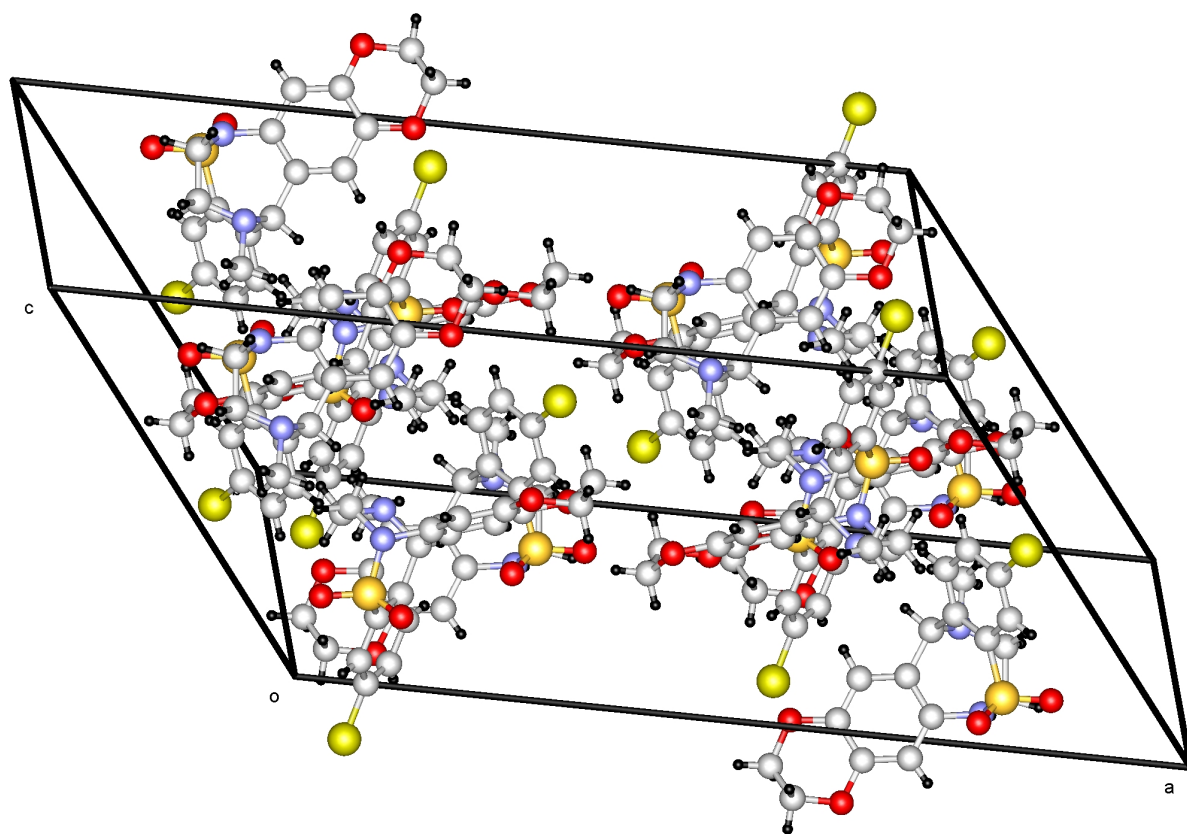


Fig. 3. Packing

Experimental

Data Collection

A colorless prism crystal of $C_{18}H_{17}ClN_2O_4S$ having approximate dimensions of 0.37 x 0.10 x 0.04 mm was mounted on a cactus needle. All measurements were made on a Rigaku RAXIS RAPID imaging plate area detector with graphite monochromated Mo-K α radiation.

Indexing was performed from 4 oscillations that were exposed for 300 seconds. The crystal-to-detector distance was 127.40 mm.

Cell constants and an orientation matrix for data collection corresponded to a C-centered monoclinic cell with dimensions:

$$\begin{aligned}a &= 26.820(2) \text{ \AA} \\b &= 10.0640(10) \text{ \AA} \quad \beta = 127.6638(16)^\circ \\c &= 16.4457(15) \text{ \AA} \\V &= 3514.0(6) \text{ \AA}^3\end{aligned}$$

For $Z = 8$ and F.W. = 392.86, the calculated density is 1.485 g/cm³. Based on the systematic absences of:

$$\begin{aligned}\text{hkl: } &h+k \pm 2n \\ \text{h0l: } &l \pm 2n\end{aligned}$$

packing considerations, a statistical analysis of intensity distribution, and the successful solution and refinement of the structure, the space group was determined to be:

C2/c (#15)

The data were collected at a temperature of $20 \pm 1^\circ\text{C}$ to a maximum 2θ value of 54.9° . A total of 180 oscillation images were collected. A sweep of data was done using ω scans from 20.0 to 200.0° in 5.0° step, at $\chi=0.0^\circ$ and $\phi = 0.0^\circ$. The exposure rate was 120.0 [sec./ $^\circ$]. A second sweep was performed using ω scans from 20.0 to 200.0° in 5.0° step, at $\chi=54.0^\circ$ and $\phi = 0.0^\circ$. The exposure rate was 120.0 [sec./ $^\circ$]. Another sweep was performed using ω scans from 20.0 to 200.0° in 5.0° step, at $\chi=54.0^\circ$ and $\phi = 90.0^\circ$. The exposure rate was 120.0 [sec./ $^\circ$]. Another sweep was performed using ω scans from 20.0 to 200.0° in 5.0° step, at $\chi=54.0^\circ$ and $\phi = 180.0^\circ$. The exposure rate was 120.0 [sec./ $^\circ$]. Another sweep was performed using ω scans from 20.0 to 200.0° in 5.0° step, at $\chi=54.0^\circ$ and $\phi = 270.0^\circ$. The exposure rate was 120.0 [sec./ $^\circ$]. The crystal-to-detector distance was 127.40 mm. Readout was performed in the 0.100 mm pixel mode.

Data Reduction

Of the 67252 reflections that were collected, 4020 were unique ($R_{\text{int}} = 0.098$).

The linear absorption coefficient, μ , for Mo-K α radiation is 3.631 cm⁻¹. The data were corrected for Lorentz and polarization effects.

Structure Solution and Refinement

The structure was solved by direct methods¹ and expanded using Fourier techniques². The non-hydrogen atoms were refined isotropically. Some hydrogen atoms were refined isotropically and the rest were refined using the riding model. The final cycle of full-matrix least-squares refinement³ on F was based on 26624 observed reflections ($I > 2.00\sigma(I)$) and 169 variable parameters and converged (largest parameter shift was 0.00 times its esd) with unweighted and weighted agreement factors of:

$$R = \Sigma ||F_o| - |F_c|| / \Sigma |F_o| = 0.0905$$

$$R_w = [\Sigma w (|F_o| - |F_c|)^2 / \Sigma w F_o^2]^{1/2} = 0.0947$$

The standard deviation of an observation of unit weight⁴ was 3.16. Unit weights were used. Plots of $\Sigma w (|F_o| - |F_c|)^2$ versus $|F_o|$, reflection order in data collection, $\sin \theta/\lambda$ and various classes of indices showed no unusual trends. The maximum and minimum peaks on the final difference Fourier map corresponded to 15.90 and -20.60 e⁻/Å³, respectively.

Neutral atom scattering factors were taken from Cromer and Waber⁵. Anomalous dispersion effects were included in Fcalc⁶; the values for $\Delta f'$ and $\Delta f''$ were those of Creagh and McAuley⁷. The values for the mass attenuation coefficients are those of Creagh and Hubbell⁸. All calculations were performed using the CrystalStructure^{9,10} crystallographic software package.

References

- (1) SIR92: Altomare, A., Cascarano, G., Giacovazzo, C., Guagliardi, A., Burla, M., Polidori, G., and Camalli, M. (1994) J. Appl. Cryst., 27, 435.
- (2) DIRDIF99: Beurskens, P.T., Admiraal, G., Beurskens, G., Bosman, W.P., de Gelder, R., Israel, R. and Smits, J.M.M. (1999). The DIRDIF-99 program system, Technical Report of the Crystallography Laboratory, University of Nijmegen, The Netherlands.

(3) Least Squares function minimized:

$$\sum w(|F_o| - |F_c|)^2 \quad \text{where } w = \text{Least Squares weights.}$$

(4) Standard deviation of an observation of unit weight:

$$[\sum w(|F_o| - |F_c|)^2 / (N_o - N_v)]^{1/2}$$

where: N_o = number of observations

N_v = number of variables

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EXPERIMENTAL DETAILS

A. Crystal Data

Empirical Formula	$\text{C}_{18}\text{H}_{17}\text{ClN}_2\text{O}_4\text{S}$
Formula Weight	392.86
Crystal Color, Habit	colorless, prism
Crystal Dimensions	0.37 X 0.10 X 0.04 mm
Crystal System	monoclinic
Lattice Type	C-centered
Indexing Images	4 oscillations @ 300.0 seconds
Detector Position	127.40 mm
Pixel Size	0.100 mm
Lattice Parameters	$a = 26.820(2) \text{ \AA}$ $b = 10.0640(10) \text{ \AA}$ $c = 16.4457(15) \text{ \AA}$ $\beta = 127.6638(16)^\circ$ $V = 3514.0(6) \text{ \AA}^3$
Space Group	C2/c (#15)
Z value	8
D_{calc}	1.485 g/cm^3
F_{000}	1632.00
$\mu(\text{MoK}\alpha)$	3.631 cm^{-1}

B. Intensity Measurements

Diffractometer	Rigaku RAXIS-RAPID
Radiation	MoK α ($\lambda = 0.71075 \text{ \AA}$) graphite monochromated
Detector Aperture	280 mm x 256 mm
Data Images	180 exposures
ω oscillation Range ($\chi=0.0$, $\phi=0.0$)	20.0 - 200.0 $^{\circ}$
Exposure Rate	120.0 sec./ $^{\circ}$
ω oscillation Range ($\chi=54.0$, $\phi=0.0$)	20.0 - 200.0 $^{\circ}$
Exposure Rate	120.0 sec./ $^{\circ}$
ω oscillation Range ($\chi=54.0$, $\phi=90.0$)	20.0 - 200.0 $^{\circ}$
Exposure Rate	120.0 sec./ $^{\circ}$
ω oscillation Range ($\chi=54.0$, $\phi=180.0$)	20.0 - 200.0 $^{\circ}$
Exposure Rate	120.0 sec./ $^{\circ}$
ω oscillation Range ($\chi=54.0$, $\phi=270.0$)	20.0 - 200.0 $^{\circ}$
Exposure Rate	120.0 sec./ $^{\circ}$
Detector Position	127.40 mm
Pixel Size	0.100 mm
$2\theta_{\text{max}}$	54.9 $^{\circ}$
No. of Reflections Measured	Total: 67252 Unique: 4020 ($R_{\text{int}} = 0.098$)
Corrections	Lorentz-polarization

C. Structure Solution and Refinement

Structure Solution	Direct Methods (SIR92)
Refinement	Full-matrix least-squares on F
Function Minimized	$\Sigma w (Fo - Fc)^2$
Least Squares Weights	1
$2\theta_{\text{max}}$ cutoff	54.9 $^{\circ}$
Anomalous Dispersion	All non-hydrogen atoms
No. Observations ($I > 2.00\sigma(I)$)	26624
No. Variables	169
Reflection/Parameter Ratio	157.54
Residuals: R ($I > 2.00\sigma(I)$)	0.0905
Residuals: Rw ($I > 2.00\sigma(I)$)	0.0947
Goodness of Fit Indicator	3.160
Max Shift/Error in Final Cycle	0.000
Maximum peak in Final Diff. Map	15.90 e $^{-}/\text{\AA}^3$
Minimum peak in Final Diff. Map	-20.60 e $^{-}/\text{\AA}^3$

Table 1. Atomic coordinates and B_{iso}/B_{eq}

atom	x	y	z	B _{eq}
Cl(1)	0.97763(4)	-0.09655(8)	0.60461(6)	5.370(18)
S(1)	0.83848(3)	-0.10063(6)	0.19520(5)	2.957(11)
O(1)	0.57102(8)	-0.12741(18)	-0.12311(13)	4.47(3)
O(3)	0.89632(8)	-0.11378(18)	0.20867(12)	4.09(3)
O(4)	0.80468(8)	-0.21760(17)	0.18502(12)	4.06(3)
O(5)	0.55017(9)	0.06711(18)	-0.01904(13)	4.72(4)
N(1)	0.79148(8)	-0.00902(18)	0.09347(13)	2.78(3)
N(8)	0.79954(9)	0.2728(2)	0.16943(14)	3.24(3)
C(9)	0.72828(10)	0.0078(2)	0.06310(17)	2.70(4)
C(10)	0.67927(12)	-0.0662(2)	-0.01596(19)	3.17(4)
C(11)	0.90348(12)	-0.0721(2)	0.39796(19)	3.19(4)
C(12)	0.82340(10)	0.0986(2)	0.30048(17)	2.91(4)
C(13)	0.85573(10)	-0.0134(2)	0.30505(17)	2.81(4)
C(14)	0.71953(10)	0.0987(2)	0.11707(17)	2.97(4)
C(15)	0.81880(12)	0.1073(2)	0.0781(2)	3.36(4)
C(16)	0.65876(11)	0.1169(2)	0.08688(19)	3.22(4)
C(17)	0.61900(11)	-0.0483(2)	-0.04466(18)	3.10(4)
C(18)	0.77380(11)	0.1797(2)	0.20514(18)	3.04(4)
C(19)	0.84544(12)	0.2168(2)	0.1587(2)	3.39(4)
C(20)	0.91847(11)	-0.0207(2)	0.48804(19)	3.49(4)
C(21)	0.60946(11)	0.0439(2)	0.00676(18)	3.30(4)
C(22)	0.84100(12)	0.1462(2)	0.3954(2)	3.66(5)
C(23)	0.82580(16)	0.3931(3)	0.2340(2)	4.64(6)
C(24)	0.88814(12)	0.0877(2)	0.4878(2)	4.01(5)
C(25)	0.49882(16)	0.0044(3)	-0.1157(2)	5.60(7)
C(26)	0.51629(16)	-0.1277(3)	-0.1234(2)	5.10(6)
H(1)	0.9253(7)	-0.1416(17)	0.3995(12)	0.9(3)
H(2)	0.8176(11)	0.222(2)	0.3941(18)	5.1(6)
H(3)	0.8987(13)	0.123(2)	0.550(2)	7.5(8)
H(4)	0.7528(9)	0.239(2)	0.2247(15)	3.0(4)
H(5)	0.7921(11)	0.432(2)	0.2313(17)	4.8(6)
H(6)	0.8609(11)	0.363(2)	0.3001(19)	4.5(5)
H(7)	0.8337(12)	0.449(2)	0.193(2)	6.5(7)
H(8)	0.8587(10)	0.289(2)	0.1321(18)	4.9(5)
H(9)	0.8841(9)	0.1800(19)	0.2224(16)	2.9(4)
H(10)	0.7856(8)	0.1366(17)	0.0106(14)	1.5(3)
H(11)	0.8553(11)	0.079(2)	0.0786(17)	4.8(5)

Table 1. Atomic coordinates and B_{iso}/B_{eq} (continued)

atom	x	y	z	B _{eq}
H(12)	0.6549(11)	0.177(2)	0.128(2)	6.2(7)
H(13)	0.6881(9)	-0.123(2)	-0.0477(16)	3.7(5)
H(14)	0.5258(13)	-0.189(3)	-0.066(2)	7.9(8)
H(15)	0.4917(14)	-0.170(2)	-0.187(2)	7.2(8)
H(16)	0.4625	-0.0005	-0.1183	6.72
H(17)	0.4897	0.0563	-0.1716	6.72

$$B_{eq} = 8/3 \pi^2 (U_{11}(aa^*)^2 + U_{22}(bb^*)^2 + U_{33}(cc^*)^2 + 2U_{12}(aa^*bb^*)\cos \gamma + 2U_{13}(aa^*cc^*)\cos \beta + 2U_{23}(bb^*cc^*)\cos \alpha)$$

Table 2. Bond lengths (Å)

atom	atom	distance	atom	atom	distance
Cl(1)	C(20)	1.749(2)	S(1)	O(3)	1.433(2)
S(1)	O(4)	1.432(2)	S(1)	N(1)	1.6312(17)
S(1)	C(13)	1.796(2)	O(1)	C(17)	1.387(2)
O(1)	C(26)	1.465(5)	O(5)	C(21)	1.392(3)
O(5)	C(25)	1.466(3)	N(1)	C(9)	1.455(3)
N(1)	C(15)	1.483(4)	N(8)	C(18)	1.482(4)
N(8)	C(19)	1.459(4)	N(8)	C(23)	1.475(3)
C(9)	C(10)	1.373(2)	C(9)	C(14)	1.391(4)
C(10)	C(17)	1.391(4)	C(10)	H(13)	0.90(2)
C(11)	C(13)	1.390(2)	C(11)	C(20)	1.377(4)
C(11)	H(1)	0.90(2)	C(12)	C(13)	1.396(3)
C(12)	C(18)	1.531(2)	C(12)	C(22)	1.410(4)
C(14)	C(16)	1.395(4)	C(14)	C(18)	1.518(2)
C(15)	C(19)	1.525(3)	C(15)	H(10)	0.951(15)
C(15)	H(11)	1.01(3)	C(16)	C(21)	1.377(2)
C(16)	H(12)	0.96(3)	C(17)	C(21)	1.381(4)
C(18)	H(4)	1.00(2)	C(19)	H(8)	1.02(3)
C(19)	H(9)	0.991(17)	C(20)	C(24)	1.359(4)
C(22)	C(24)	1.381(3)	C(22)	H(2)	0.98(2)
C(23)	H(5)	0.96(3)	C(23)	H(6)	0.95(2)
C(23)	H(7)	0.99(3)	C(24)	H(3)	0.94(3)
C(25)	C(26)	1.440(5)	C(25)	H(16)	0.950
C(25)	H(17)	0.950	C(26)	H(14)	1.02(3)
C(26)	H(15)	0.93(2)			

Table 3. Bond angles (°)

atom	atom	atom	angle	atom	atom	atom	angle
O(3)	S(1)	O(4)	119.37(12)	O(3)	S(1)	N(1)	107.21(13)
O(3)	S(1)	C(13)	107.48(11)	O(4)	S(1)	N(1)	107.91(9)
O(4)	S(1)	C(13)	105.96(14)	N(1)	S(1)	C(13)	108.54(11)
C(17)	O(1)	C(26)	111.2(2)	C(21)	O(5)	C(25)	113.5(2)
S(1)	N(1)	C(9)	115.2(2)	S(1)	N(1)	C(15)	117.94(13)
C(9)	N(1)	C(15)	116.31(19)	C(18)	N(8)	C(19)	115.8(2)
C(18)	N(8)	C(23)	111.5(2)	C(19)	N(8)	C(23)	109.7(2)
N(1)	C(9)	C(10)	120.2(2)	N(1)	C(9)	C(14)	118.22(17)
C(10)	C(9)	C(14)	121.6(2)	C(9)	C(10)	C(17)	119.5(2)
C(9)	C(10)	H(13)	117.2(13)	C(17)	C(10)	H(13)	123.3(13)
C(13)	C(11)	C(20)	119.3(2)	C(13)	C(11)	H(1)	120.5(11)
C(20)	C(11)	H(1)	120.2(11)	C(13)	C(12)	C(18)	127.4(2)
C(13)	C(12)	C(22)	115.97(19)	C(18)	C(12)	C(22)	116.5(2)
S(1)	C(13)	C(11)	113.3(2)	S(1)	C(13)	C(12)	124.80(15)
C(11)	C(13)	C(12)	121.8(2)	C(9)	C(14)	C(16)	118.34(19)
C(9)	C(14)	C(18)	121.6(2)	C(16)	C(14)	C(18)	120.0(2)
N(1)	C(15)	C(19)	116.1(3)	N(1)	C(15)	H(10)	103.7(13)
N(1)	C(15)	H(11)	110.7(14)	C(19)	C(15)	H(10)	113.5(11)
C(19)	C(15)	H(11)	104.9(12)	H(10)	C(15)	H(11)	108(2)
C(14)	C(16)	C(21)	120.0(2)	C(14)	C(16)	H(12)	115.9(15)
C(21)	C(16)	H(12)	123.9(16)	O(1)	C(17)	C(10)	117.8(2)
O(1)	C(17)	C(21)	122.8(2)	C(10)	C(17)	C(21)	119.4(2)
N(8)	C(18)	C(12)	114.8(2)	N(8)	C(18)	C(14)	111.6(2)
N(8)	C(18)	H(4)	104.1(14)	C(12)	C(18)	C(14)	114.71(19)
C(12)	C(18)	H(4)	107.2(11)	C(14)	C(18)	H(4)	103.0(10)
N(8)	C(19)	C(15)	113.7(2)	N(8)	C(19)	H(8)	108.7(17)
N(8)	C(19)	H(9)	115.6(19)	C(15)	C(19)	H(8)	104.0(16)
C(15)	C(19)	H(9)	106.3(13)	H(8)	C(19)	H(9)	107.8(19)
Cl(1)	C(20)	C(11)	119.2(2)	Cl(1)	C(20)	C(24)	119.6(2)
C(11)	C(20)	C(24)	121.2(2)	O(5)	C(21)	C(16)	116.9(2)
O(5)	C(21)	C(17)	122.06(19)	C(16)	C(21)	C(17)	121.1(2)
C(12)	C(22)	C(24)	122.4(2)	C(12)	C(22)	H(2)	117.6(14)
C(24)	C(22)	H(2)	120.0(14)	N(8)	C(23)	H(5)	106.3(13)
N(8)	C(23)	H(6)	105.5(14)	N(8)	C(23)	H(7)	100.3(16)
H(5)	C(23)	H(6)	115(2)	H(5)	C(23)	H(7)	109(2)
H(6)	C(23)	H(7)	119(2)	C(20)	C(24)	C(22)	119.3(3)
C(20)	C(24)	H(3)	120.9(16)	C(22)	C(24)	H(3)	119.8(16)

Table 3. Bond angles ($^{\circ}$) (continued)

atom	atom	atom	angle	atom	atom	atom	angle
O(5)	C(25)	C(26)	110.6(2)	O(5)	C(25)	H(16)	109.2
O(5)	C(25)	H(17)	109.2	C(26)	C(25)	H(16)	109.2
C(26)	C(25)	H(17)	109.2	H(16)	C(25)	H(17)	109.5
O(1)	C(26)	C(25)	112.2(3)	O(1)	C(26)	H(14)	108(2)
O(1)	C(26)	H(15)	92(2)	C(25)	C(26)	H(14)	113(2)
C(25)	C(26)	H(15)	118.3(17)	H(14)	C(26)	H(15)	111(2)

Table 4. Torsion Angles($^{\circ}$)

atom1	atom2	atom3	atom4	angle	atom1	atom2	atom3	atom4	angle
O(3)	S(1)	N(1)	C(9)	175.85(15)	O(3)	S(1)	N(1)	C(15)	-40.6(2)
O(3)	S(1)	C(13)	C(11)	-54.3(2)	O(3)	S(1)	C(13)	C(12)	128.7(2)
O(4)	S(1)	N(1)	C(9)	46.1(2)	O(4)	S(1)	N(1)	C(15)	-170.4(2)
O(4)	S(1)	C(13)	C(11)	74.4(2)	O(4)	S(1)	C(13)	C(12)	-102.6(2)
N(1)	S(1)	C(13)	C(11)	-169.9(2)	N(1)	S(1)	C(13)	C(12)	13.0(2)
C(13)	S(1)	N(1)	C(9)	-68.32(18)	C(13)	S(1)	N(1)	C(15)	75.2(2)
C(17)	O(1)	C(26)	C(25)	47.7(3)	C(26)	O(1)	C(17)	C(10)	162.5(2)
C(26)	O(1)	C(17)	C(21)	-16.0(3)	C(21)	O(5)	C(25)	C(26)	41.6(4)
C(25)	O(5)	C(21)	C(16)	169.5(2)	C(25)	O(5)	C(21)	C(17)	-10.8(3)
S(1)	N(1)	C(9)	C(10)	-100.6(2)	S(1)	N(1)	C(9)	C(14)	78.6(2)
S(1)	N(1)	C(15)	C(19)	-64.0(3)	C(9)	N(1)	C(15)	C(19)	79.1(2)
C(15)	N(1)	C(9)	C(10)	115.3(2)	C(15)	N(1)	C(9)	C(14)	-65.6(3)
C(18)	N(8)	C(19)	C(15)	63.2(2)	C(19)	N(8)	C(18)	C(12)	49.2(2)
C(19)	N(8)	C(18)	C(14)	-83.5(2)	C(23)	N(8)	C(18)	C(12)	-77.1(2)
C(23)	N(8)	C(18)	C(14)	150.1(2)	C(23)	N(8)	C(19)	C(15)	-169.5(2)
N(1)	C(9)	C(10)	C(17)	-179.5(2)	N(1)	C(9)	C(14)	C(16)	178.8(2)
N(1)	C(9)	C(14)	C(18)	0.3(3)	C(10)	C(9)	C(14)	C(16)	-2.1(4)
C(10)	C(9)	C(14)	C(18)	179.4(2)	C(14)	C(9)	C(10)	C(17)	1.4(4)
C(9)	C(10)	C(17)	O(1)	-178.4(2)	C(9)	C(10)	C(17)	C(21)	0.2(3)
C(13)	C(11)	C(20)	Cl(1)	178.9(2)	C(13)	C(11)	C(20)	C(24)	-1.1(4)
C(20)	C(11)	C(13)	S(1)	-175.5(2)	C(20)	C(11)	C(13)	C(12)	1.6(4)
C(13)	C(12)	C(18)	N(8)	-78.3(3)	C(13)	C(12)	C(18)	C(14)	53.0(4)
C(18)	C(12)	C(13)	S(1)	-8.4(4)	C(18)	C(12)	C(13)	C(11)	174.8(2)
C(13)	C(12)	C(22)	C(24)	1.2(4)	C(22)	C(12)	C(13)	S(1)	175.2(2)
C(22)	C(12)	C(13)	C(11)	-1.6(4)	C(18)	C(12)	C(22)	C(24)	-175.6(3)
C(22)	C(12)	C(18)	N(8)	98.1(3)	C(22)	C(12)	C(18)	C(14)	-130.6(3)
C(9)	C(14)	C(16)	C(21)	1.2(4)	C(9)	C(14)	C(18)	N(8)	65.3(3)
C(9)	C(14)	C(18)	C(12)	-67.5(3)	C(16)	C(14)	C(18)	N(8)	-113.2(3)
C(16)	C(14)	C(18)	C(12)	114.0(3)	C(18)	C(14)	C(16)	C(21)	179.8(2)
N(1)	C(15)	C(19)	N(8)	-58.1(2)	C(14)	C(16)	C(21)	O(5)	179.9(2)
C(14)	C(16)	C(21)	C(17)	0.3(4)	O(1)	C(17)	C(21)	O(5)	-2.2(4)
O(1)	C(17)	C(21)	C(16)	177.5(2)	C(10)	C(17)	C(21)	O(5)	179.4(2)
C(10)	C(17)	C(21)	C(16)	-1.0(4)	Cl(1)	C(20)	C(24)	C(22)	-179.3(2)
C(11)	C(20)	C(24)	C(22)	0.7(5)	C(12)	C(22)	C(24)	C(20)	-0.8(5)
O(5)	C(25)	C(26)	O(1)	-61.5(4)					

The sign is positive if when looking from atom 2 to atom 3 a clock-wise motion of atom 1 would superimpose it on atom 4.

Table 5. Distances beyond the asymmetric unit out to 3.60 Å

atom	atom	distance	atom	atom	distance
Cl(1)	O(3) ¹⁾	3.5571(16)	Cl(1)	O(5) ²⁾	3.5654(19)
Cl(1)	H(1) ¹⁾	3.57(2)	Cl(1)	H(9) ¹⁾	3.108(18)
Cl(1)	H(11) ³⁾	3.04(3)	S(1)	H(12) ²⁾	3.59(3)
S(1)	H(13) ⁴⁾	3.47(2)	O(1)	O(1) ⁵⁾	3.536(2)
O(1)	O(3) ⁴⁾	3.322(3)	O(1)	C(24) ²⁾	3.375(3)
O(1)	C(26) ⁵⁾	3.306(4)	O(1)	H(2) ²⁾	3.41(2)
O(1)	H(3) ²⁾	2.69(2)	O(1)	H(15) ⁵⁾	2.52(3)
O(1)	H(17) ⁵⁾	3.279	O(3)	Cl(1) ¹⁾	3.5571(16)
O(3)	O(1) ⁴⁾	3.322(3)	O(3)	C(24) ⁶⁾	3.512(4)
O(3)	H(3) ⁶⁾	2.66(4)	O(3)	H(13) ⁴⁾	3.44(2)
O(3)	H(15) ⁷⁾	2.98(2)	O(4)	C(22) ²⁾	3.547(3)
O(4)	H(2) ²⁾	2.76(2)	O(4)	H(4) ²⁾	2.75(3)
O(4)	H(7) ⁸⁾	3.43(2)	O(4)	H(12) ²⁾	2.77(3)
O(4)	H(13) ⁴⁾	2.87(2)	O(5)	Cl(1) ⁹⁾	3.5654(19)
O(5)	O(5) ¹⁰⁾	3.394(3)	O(5)	C(25) ¹⁰⁾	3.274(6)
O(5)	H(1) ⁹⁾	3.363(18)	O(5)	H(6) ²⁾	3.52(2)
O(5)	H(14) ¹⁰⁾	3.33(4)	O(5)	H(16) ¹⁰⁾	2.570
N(1)	H(3) ⁶⁾	3.55(4)	N(8)	C(15) ¹¹⁾	3.508(2)
N(8)	H(10) ¹¹⁾	2.559(16)	C(9)	H(2) ²⁾	3.36(2)
C(9)	H(7) ¹¹⁾	3.49(3)	C(9)	H(8) ¹¹⁾	3.28(2)
C(10)	H(2) ²⁾	2.89(3)	C(10)	H(7) ¹¹⁾	2.96(3)
C(10)	H(8) ¹¹⁾	3.18(2)	C(11)	H(12) ²⁾	2.86(2)
C(12)	H(5) ²⁾	3.27(3)	C(13)	H(12) ²⁾	3.37(2)
C(14)	H(5) ²⁾	3.17(3)	C(14)	H(8) ¹¹⁾	3.46(2)
C(14)	H(10) ¹¹⁾	3.34(2)	C(15)	N(8) ¹¹⁾	3.508(2)
C(15)	C(24) ⁶⁾	3.590(5)	C(15)	H(3) ⁶⁾	3.38(3)
C(15)	H(10) ¹¹⁾	3.409(18)	C(16)	C(23) ²⁾	3.525(5)
C(16)	H(1) ⁹⁾	3.41(2)	C(16)	H(5) ²⁾	3.05(2)
C(16)	H(6) ²⁾	3.38(3)	C(16)	H(8) ¹¹⁾	3.47(3)
C(17)	H(2) ²⁾	3.03(2)	C(17)	H(3) ²⁾	3.34(3)
C(17)	H(7) ¹¹⁾	3.52(4)	C(17)	H(8) ¹¹⁾	3.21(2)
C(18)	H(5) ²⁾	3.56(3)	C(18)	H(10) ¹¹⁾	3.41(2)
C(19)	H(10) ¹¹⁾	3.213(16)	C(20)	H(11) ³⁾	2.92(3)
C(20)	H(12) ²⁾	3.49(2)	C(20)	H(14) ⁹⁾	3.56(2)
C(21)	H(5) ²⁾	3.59(2)	C(21)	H(6) ²⁾	3.31(3)
C(21)	H(8) ¹¹⁾	3.34(3)	C(21)	H(16) ¹⁰⁾	3.407
C(22)	O(4) ⁹⁾	3.547(3)	C(22)	H(5) ²⁾	3.55(2)

Table 5. Distances beyond the asymmetric unit out to 3.60 Å (continued)

atom	atom	distance	atom	atom	distance
C(22)	H(11) ³⁾	3.60(2)	C(22)	H(14) ⁹⁾	3.34(2)
C(23)	C(16) ⁹⁾	3.525(5)	C(23)	H(10) ¹¹⁾	3.242(16)
C(23)	H(12) ⁹⁾	3.49(2)	C(23)	H(16) ¹²⁾	3.103
C(24)	O(1) ⁹⁾	3.375(3)	C(24)	O(3) ³⁾	3.512(4)
C(24)	C(15) ³⁾	3.590(5)	C(24)	C(26) ⁹⁾	3.573(4)
C(24)	H(11) ³⁾	2.73(3)	C(24)	H(14) ⁹⁾	2.90(3)
C(25)	O(5) ¹⁰⁾	3.274(6)	C(25)	H(6) ¹³⁾	3.33(2)
C(25)	H(7) ¹³⁾	3.57(2)	C(25)	H(16) ¹⁰⁾	3.319
C(26)	O(1) ⁵⁾	3.306(4)	C(26)	C(24) ²⁾	3.573(4)
C(26)	H(1) ¹⁴⁾	3.55(2)	C(26)	H(3) ²⁾	3.10(3)
C(26)	H(15) ⁵⁾	3.02(4)	H(1)	Cl(1) ¹⁾	3.57(2)
H(1)	O(5) ²⁾	3.363(18)	H(1)	C(16) ²⁾	3.41(2)
H(1)	C(26) ⁷⁾	3.55(2)	H(1)	H(12) ²⁾	2.64(3)
H(1)	H(14) ⁷⁾	2.94(4)	H(1)	H(15) ⁷⁾	3.45(4)
H(2)	O(1) ⁹⁾	3.41(2)	H(2)	O(4) ⁹⁾	2.76(2)
H(2)	C(9) ⁹⁾	3.36(2)	H(2)	C(10) ⁹⁾	2.89(3)
H(2)	C(17) ⁹⁾	3.03(2)	H(2)	H(13) ⁹⁾	3.05(4)
H(2)	H(14) ⁹⁾	3.45(3)	H(3)	O(1) ⁹⁾	2.69(2)
H(3)	O(3) ³⁾	2.66(4)	H(3)	N(1) ³⁾	3.55(4)
H(3)	C(15) ³⁾	3.38(3)	H(3)	C(17) ⁹⁾	3.34(3)
H(3)	C(26) ⁹⁾	3.10(3)	H(3)	H(11) ³⁾	2.53(4)
H(3)	H(13) ⁹⁾	3.44(4)	H(3)	H(14) ⁹⁾	2.66(5)
H(3)	H(15) ⁹⁾	3.16(3)	H(4)	O(4) ⁹⁾	2.75(3)
H(4)	H(5) ²⁾	3.54(3)	H(4)	H(10) ¹¹⁾	3.57(3)
H(5)	C(12) ⁹⁾	3.27(3)	H(5)	C(14) ⁹⁾	3.17(3)
H(5)	C(16) ⁹⁾	3.05(2)	H(5)	C(18) ⁹⁾	3.56(3)
H(5)	C(21) ⁹⁾	3.59(2)	H(5)	C(22) ⁹⁾	3.55(2)
H(5)	H(4) ⁹⁾	3.54(3)	H(5)	H(10) ¹¹⁾	3.24(3)
H(5)	H(12) ⁹⁾	3.07(3)	H(6)	O(5) ⁹⁾	3.52(2)
H(6)	C(16) ⁹⁾	3.38(3)	H(6)	C(21) ⁹⁾	3.31(3)
H(6)	C(25) ¹²⁾	3.33(2)	H(6)	H(12) ⁹⁾	3.49(3)
H(6)	H(14) ⁹⁾	3.52(3)	H(6)	H(16) ¹²⁾	2.581
H(6)	H(17) ¹²⁾	3.291	H(7)	O(4) ¹⁵⁾	3.43(2)
H(7)	C(9) ¹¹⁾	3.49(3)	H(7)	C(10) ¹¹⁾	2.96(3)
H(7)	C(17) ¹¹⁾	3.52(4)	H(7)	C(25) ¹²⁾	3.57(2)
H(7)	H(10) ¹¹⁾	3.01(2)	H(7)	H(12) ⁹⁾	3.60(4)
H(7)	H(13) ¹¹⁾	2.73(4)	H(7)	H(16) ¹²⁾	2.956

Table 5. Distances beyond the asymmetric unit out to 3.60 Å (continued)

atom	atom	distance	atom	atom	distance
H(7)	H(17) ¹²⁾	3.329	H(8)	C(9) ¹¹⁾	3.28(2)
H(8)	C(10) ¹¹⁾	3.18(2)	H(8)	C(14) ¹¹⁾	3.46(2)
H(8)	C(16) ¹¹⁾	3.47(3)	H(8)	C(17) ¹¹⁾	3.21(2)
H(8)	C(21) ¹¹⁾	3.34(3)	H(8)	H(10) ¹¹⁾	3.16(2)
H(8)	H(13) ¹¹⁾	3.56(3)	H(8)	H(17) ¹²⁾	3.365
H(9)	Cl(1) ¹⁾	3.108(18)	H(9)	H(17) ¹²⁾	3.473
H(10)	N(8) ¹¹⁾	2.559(16)	H(10)	C(14) ¹¹⁾	3.34(2)
H(10)	C(15) ¹¹⁾	3.409(18)	H(10)	C(18) ¹¹⁾	3.41(2)
H(10)	C(19) ¹¹⁾	3.213(16)	H(10)	C(23) ¹¹⁾	3.242(16)
H(10)	H(4) ¹¹⁾	3.57(3)	H(10)	H(5) ¹¹⁾	3.24(3)
H(10)	H(7) ¹¹⁾	3.01(2)	H(10)	H(8) ¹¹⁾	3.16(2)
H(10)	H(10) ¹¹⁾	2.86(2)	H(11)	Cl(1) ⁶⁾	3.04(3)
H(11)	C(20) ⁶⁾	2.92(3)	H(11)	C(22) ⁶⁾	3.60(2)
H(11)	C(24) ⁶⁾	2.73(3)	H(11)	H(3) ⁶⁾	2.53(4)
H(12)	S(1) ⁹⁾	3.59(3)	H(12)	O(4) ⁹⁾	2.77(3)
H(12)	C(11) ⁹⁾	2.86(2)	H(12)	C(13) ⁹⁾	3.37(2)
H(12)	C(20) ⁹⁾	3.49(2)	H(12)	C(23) ²⁾	3.49(2)
H(12)	H(1) ⁹⁾	2.64(3)	H(12)	H(5) ²⁾	3.07(3)
H(12)	H(6) ²⁾	3.49(3)	H(12)	H(7) ²⁾	3.60(4)
H(12)	H(16) ¹⁰⁾	3.534	H(13)	S(1) ⁴⁾	3.47(2)
H(13)	O(3) ⁴⁾	3.44(2)	H(13)	O(4) ⁴⁾	2.87(2)
H(13)	H(2) ²⁾	3.05(4)	H(13)	H(3) ²⁾	3.44(4)
H(13)	H(7) ¹¹⁾	2.73(4)	H(13)	H(8) ¹¹⁾	3.56(3)
H(14)	O(5) ¹⁰⁾	3.33(4)	H(14)	C(20) ²⁾	3.56(2)
H(14)	C(22) ²⁾	3.34(2)	H(14)	C(24) ²⁾	2.90(3)
H(14)	H(1) ¹⁴⁾	2.94(4)	H(14)	H(2) ²⁾	3.45(3)
H(14)	H(3) ²⁾	2.66(5)	H(14)	H(6) ²⁾	3.52(3)
H(14)	H(16) ¹⁰⁾	3.433	H(15)	O(1) ⁵⁾	2.52(3)
H(15)	O(3) ¹⁴⁾	2.98(2)	H(15)	C(26) ⁵⁾	3.02(4)
H(15)	H(1) ¹⁴⁾	3.45(4)	H(15)	H(3) ²⁾	3.16(3)
H(15)	H(15) ⁵⁾	2.37(6)	H(15)	H(17) ⁵⁾	3.501
H(16)	O(5) ¹⁰⁾	2.570	H(16)	C(21) ¹⁰⁾	3.407
H(16)	C(23) ¹³⁾	3.103	H(16)	C(25) ¹⁰⁾	3.319
H(16)	H(6) ¹³⁾	2.581	H(16)	H(7) ¹³⁾	2.956
H(16)	H(12) ¹⁰⁾	3.534	H(16)	H(14) ¹⁰⁾	3.433
H(16)	H(16) ¹⁰⁾	3.102	H(17)	O(1) ⁵⁾	3.279
H(17)	H(6) ¹³⁾	3.291	H(17)	H(7) ¹³⁾	3.329

Table 5. Distances beyond the asymmetric unit out to 3.60 Å (continued)

atom	atom	distance	atom	atom	distance
H(17)	H(8) ¹³⁾	3.365	H(17)	H(9) ¹³⁾	3.473
H(17)	H(15) ⁵⁾	3.501	H(17)	H(17) ⁵⁾	2.950

Symmetry Operators:

- | | |
|-----------------------------|-------------------------------|
| (1) -X+2,-Y,-Z+1 | (2) -X+1/2+1,Y+1/2-1,-Z+1/2 |
| (3) X,-Y,Z+1/2 | (4) -X+1/2+1,-Y+1/2-1,-Z |
| (5) -X+1,Y,-Z+1/2-1 | (6) X,-Y,Z+1/2-1 |
| (7) X+1/2,-Y+1/2-1,Z+1/2 | (8) X,Y-1,Z |
| (9) -X+1/2+1,Y+1/2,-Z+1/2 | (10) -X+1,-Y,-Z |
| (11) -X+1/2+1,-Y+1/2,-Z | (12) X+1/2,-Y+1/2,Z+1/2 |
| (13) X+1/2-1,-Y+1/2,Z+1/2-1 | (14) X+1/2-1,-Y+1/2-1,Z+1/2-1 |
| (15) X,Y+1,Z | |