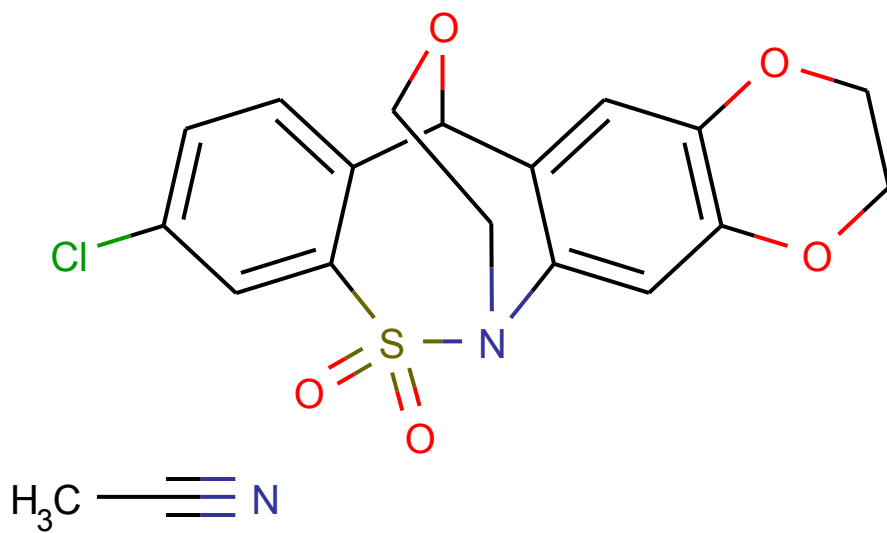


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2182-BGE

Submitted by: Berecz Gabor
Operator: Dancso Andras

X-ray Structure Report



July 14, 2025

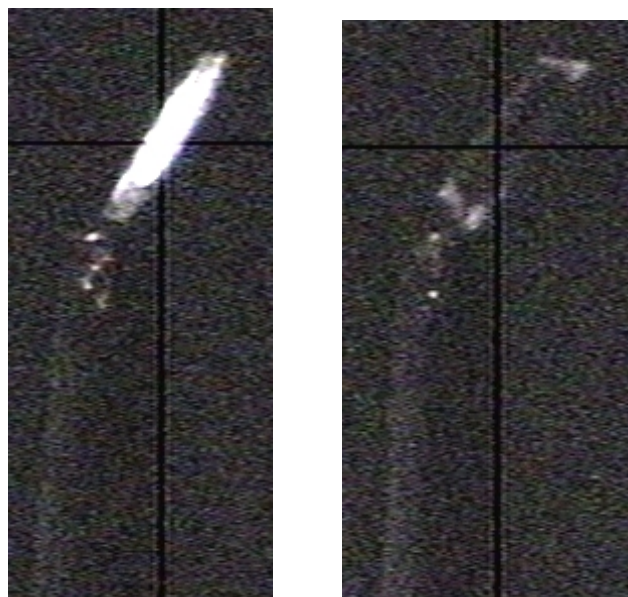


Fig. 1. The crystal

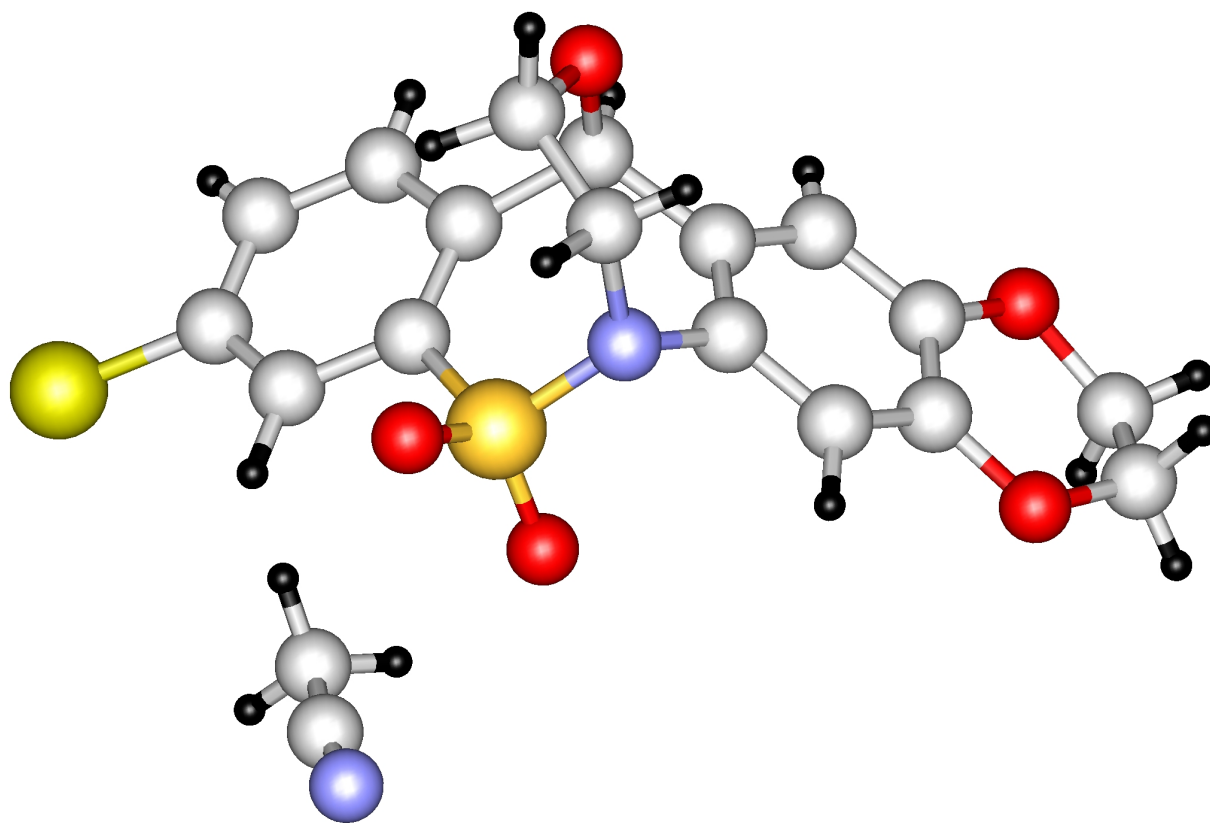


Fig. 2. The molecule (some hydrogens were generated by the software)

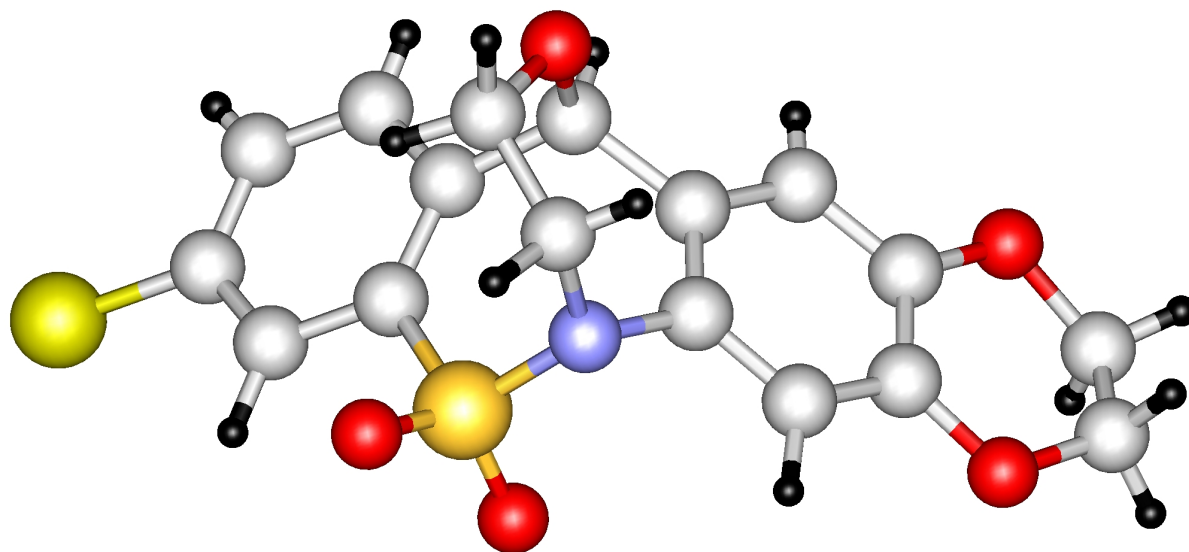


Fig. 3. Fragment 1

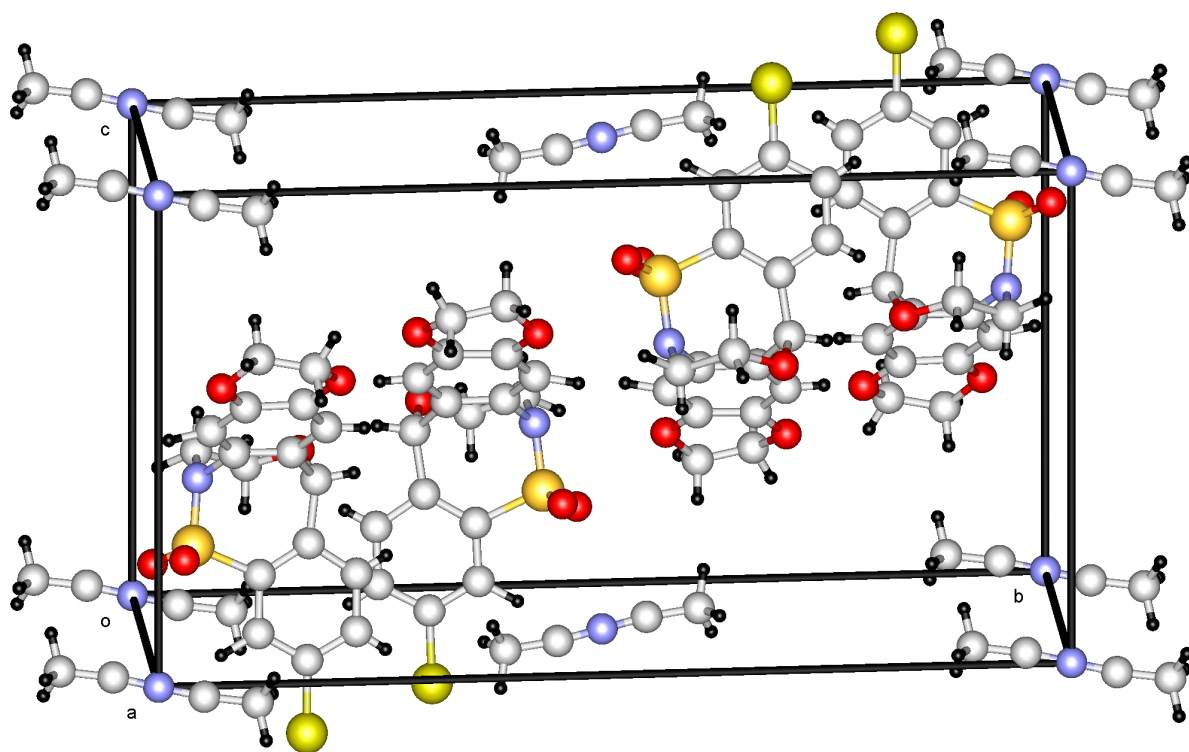


Fig. 4. Packing

Experimental

Data Collection

A colorless prism crystal of $C_{19}H_{17}ClN_2O_5S$ having approximate dimensions of 0.29 x 0.09 x 0.05 mm was mounted on a cactus needle. All measurements were made on a Rigaku RAXIS RAPID imaging plate area detector with graphite monochromated Cu-K α radiation.

Indexing was performed from 4 oscillations that were exposed for 900 seconds. The crystal-to-detector distance was 127.40 mm.

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

$$\begin{aligned}a &= 6.6806(5) \text{ \AA} \\b &= 21.7251(14) \text{ \AA} \quad \beta = 92.707(4)^\circ \\c &= 12.1261(8) \text{ \AA} \\V &= 1758.0(2) \text{ \AA}^3\end{aligned}$$

For $Z = 4$ and F.W. = 420.87, the calculated density is 1.590 g/cm³. The systematic absences of:

$$\begin{aligned}h0l: h \pm 2n \\0k0: k \pm 2n\end{aligned}$$

uniquely determine the space group to be:

$$P2_1/a \text{ (\#14)}$$

The data were collected at a temperature of $20 \pm 1^\circ\text{C}$ to a maximum 2θ value of 143.4° . 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 180.0 [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 180.0 [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 180.0 [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 180.0 [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 180.0 [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 20287 reflections that were collected, 3324 were unique ($R_{\text{int}} = 0.055$).

The linear absorption coefficient, μ , for Cu-K α radiation is 33.675 cm⁻¹. An empirical absorption correction was applied which resulted in transmission factors ranging from 0.549 to 0.845. 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. Some non-hydrogen atoms were refined anisotropically, while the rest 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 13358 observed reflections ($I > 2.00\sigma(I)$) and 279 variable parameters and converged (largest parameter shift was 0.00 times its esd) with unweighted and weighted agreement factors of:

$$R = \sum ||F_o| - |F_c|| / \sum |F_o| = 0.0742$$

$$R_w = [\sum w (|F_o| - |F_c|)^2 / \sum w F_o^2]^{1/2} = 0.0835$$

The standard deviation of an observation of unit weight⁴ was 4.76. Unit weights were used. Plots of $\sum w (|F_o| - |F_c|)^2$ 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 6.13 and -7.57 e⁻/Å³, respectively.

Neutral atom scattering factors were taken from Cromer and Waber⁵. Anomalous dispersion effects were included in F_{calc} ⁶; 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

(5) Cromer, D. T. & Waber, J. T.; "International Tables for X-ray Crystallography", Vol. IV, The Kynoch Press, Birmingham, England, Table 2.2 A (1974).

(6) Ibers, J. A. & Hamilton, W. C.; Acta Crystallogr., 17, 781 (1964).

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(8) Creagh, D. C. & Hubbell, J.H.; "International Tables for Crystallography", Vol C, (A.J.C. Wilson, ed.), Kluwer Academic Publishers, Boston, Table 4.2.4.3, pages 200-206 (1992).

(9) CrystalStructure 3.7.0: Crystal Structure Analysis Package, Rigaku and Rigaku/MSK (2000-2005). 9009 New Trails Dr. The Woodlands TX 77381 USA.

(10) CRYSTALS Issue 10: Watkin, D.J., Prout, C.K. Carruthers, J.R. & Betteridge, P.W. Chemical Crystallography Laboratory, Oxford, UK. (1996)

EXPERIMENTAL DETAILS

A. Crystal Data

Empirical Formula	$\text{C}_{19}\text{H}_{17}\text{ClN}_2\text{O}_5\text{S}$
Formula Weight	420.87
Crystal Color, Habit	colorless, prism
Crystal Dimensions	0.29 X 0.09 X 0.05 mm
Crystal System	monoclinic
Lattice Type	Primitive
Indexing Images	4 oscillations @ 900.0 seconds
Detector Position	127.40 mm
Pixel Size	0.100 mm
Lattice Parameters	$a = 6.6806(5) \text{ \AA}$ $b = 21.7251(14) \text{ \AA}$ $c = 12.1261(8) \text{ \AA}$ $\beta = 92.707(4)^\circ$ $V = 1758.0(2) \text{ \AA}^3$
Space Group	$P2_1/a$ (#14)
Z value	4
D _{calc}	1.590 g/cm ³
F ₀₀₀	872.00
$\mu(\text{CuK}\alpha)$	33.675 cm ⁻¹

B. Intensity Measurements

Diffractometer	Rigaku RAXIS-RAPID
Radiation	CuK α (λ = 1.54187 Å) graphite monochromated
Detector Aperture	280 mm x 256 mm
Data Images	180 exposures
ω oscillation Range (χ =0.0, ϕ =0.0)	20.0 - 200.0°
Exposure Rate	180.0 sec./°
ω oscillation Range (χ =54.0, ϕ =0.0)	20.0 - 200.0°
Exposure Rate	180.0 sec./°
ω oscillation Range (χ =54.0, ϕ =90.0)	20.0 - 200.0°
Exposure Rate	180.0 sec./°
ω oscillation Range (χ =54.0, ϕ =180.0)	20.0 - 200.0°
Exposure Rate	180.0 sec./°
ω oscillation Range (χ =54.0, ϕ =270.0)	20.0 - 200.0°
Exposure Rate	180.0 sec./°
Detector Position	127.40 mm
Pixel Size	0.100 mm
$2\theta_{\text{max}}$	143.4°
No. of Reflections Measured	Total: 20287 Unique: 3324 (R_{int} = 0.055)
Corrections	Lorentz-polarization Absorption (trans. factors: 0.549 - 0.845)

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	143.4 $^{\circ}$
Anomalous Dispersion	All non-hydrogen atoms
No. Observations ($I > 2.00\sigma(I)$)	13358
No. Variables	279
Reflection/Parameter Ratio	47.88
Residuals: R ($I > 2.00\sigma(I)$)	0.0742
Residuals: Rw ($I > 2.00\sigma(I)$)	0.0835
Goodness of Fit Indicator	4.762
Max Shift/Error in Final Cycle	0.000
Maximum peak in Final Diff. Map	6.13 e $^{-}/\text{\AA}^3$
Minimum peak in Final Diff. Map	-7.57 e $^{-}/\text{\AA}^3$

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

atom	x	y	z	B _{eq}	occ
Cl(2)	0.58395(18)	0.17430(6)	-0.17994(9)	6.25(3)	1/2
S(1)	0.53619(16)	0.05103(4)	0.19781(8)	3.03(2)	
O(3)	0.2212(3)	0.17672(10)	0.32766(18)	2.87(6)	
O(4)	0.9481(3)	0.20065(11)	0.60382(19)	3.82(6)	
O(5)	0.4027(3)	0.01039(10)	0.13787(19)	4.20(7)	
O(6)	0.7437(3)	0.03576(10)	0.20798(19)	3.62(6)	
O(7)	0.9868(3)	0.06944(10)	0.6014(2)	4.24(7)	
N(1)	0.4562(4)	0.05808(11)	0.3205(2)	2.64(7)	
N(2)	1.0000	0.0000	0.0000	8.9(3)	
C(1)	0.6941(5)	0.19254(16)	0.4621(2)	2.75(9)	
C(2)	1.0444(11)	0.1070(3)	-0.0289(6)	17.080	
C(9)	0.4761(4)	0.17966(16)	0.1860(2)	2.35(8)	
C(11)	0.5858(5)	0.09478(16)	0.3948(2)	2.60(8)	
C(12)	0.5730(4)	0.15806(14)	0.3902(2)	2.31(8)	
C(13)	0.8308(5)	0.16370(16)	0.5354(2)	2.88(9)	
C(14)	0.8462(5)	0.09983(16)	0.5355(2)	3.00(9)	
C(15)	0.5179(4)	0.12456(14)	0.1320(2)	2.52(8)	
C(16)	0.7224(5)	0.06522(17)	0.4656(3)	3.06(9)	
C(17)	0.5491(5)	0.1222(2)	0.0193(3)	3.33(10)	
C(18)	0.4292(5)	0.18907(14)	0.3066(3)	2.66(9)	
C(19)	0.1402(5)	0.12031(18)	0.2843(3)	3.00(10)	
C(20)	0.4683(5)	0.23208(19)	0.1200(3)	3.02(10)	
C(21)	0.2364(6)	0.06293(19)	0.3343(3)	3.36(11)	
C(22)	0.5015(5)	0.2309(2)	0.0093(3)	3.65(11)	
C(23)	0.5415(5)	0.1759(2)	-0.0394(2)	3.73(10)	
C(25)	1.1064(7)	0.1684(2)	0.6594(4)	8.32(16)	
C(26)	1.0898(8)	0.1076(2)	0.6786(4)	10.08(19)	
C(28)	1.0203(12)	0.0493(3)	-0.0183(6)	16.867	
H(1)	0.437(4)	0.2654(12)	0.150(2)	1.8(8)	
H(2)	0.498(4)	0.2628(12)	-0.030(2)	1.8(7)	
H(3)	0.582(3)	0.0804(11)	-0.018(2)	1.7(6)	
H(4)	0.439(3)	0.2300(10)	0.3199(19)	0.6(5)	
H(5)	0.687(3)	0.2343(11)	0.457(2)	1.7(6)	
H(6)	0.751(4)	0.0185(13)	0.454(2)	3.8(8)	
H(7)	-0.004(4)	0.1256(11)	0.307(2)	1.8(6)	
H(8)	0.152(3)	0.1186(10)	0.200(2)	1.2(6)	
H(9)	0.172(4)	0.0230(13)	0.291(2)	4.0(8)	

Table 1. Atomic coordinates and $B_{\text{iso}}/B_{\text{eq}}$ and occupancy (continued)

atom	x	y	z	B_{eq}	occ
H(10)	0.215(4)	0.0654(12)	0.409(2)	2.3(8)	
H(11)	1.2184	0.1719	0.6143	9.70	
H(12)	1.1345	0.1890	0.7274	9.68	
H(13)	1.0178	0.1051	0.7441	11.73	
H(14)	1.2199	0.0907	0.6917	11.71	
H(15)	0.9174	0.1267	-0.0358	20.56	
H(16)	1.1161	0.1145	-0.0934	20.55	
H(17)	1.1180	0.1230	0.0337	20.56	

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

Table 2. Anisotropic displacement parameters

atom	U ₁₁	U ₂₂	U ₃₃	U ₁₂	U ₁₃	U ₂₃
Cl(2)	0.0789(9)	0.1193(10)	0.0395(6)	0.0085(7)	0.0072(6)	0.0142(6)
S(1)	0.0428(6)	0.0306(5)	0.0415(5)	0.0031(5)	-0.0007(5)	-0.0046(4)
O(3)	0.0300(15)	0.0319(14)	0.0473(16)	0.0024(11)	0.0038(12)	-0.0039(11)
O(4)	0.0522(19)	0.0420(15)	0.0489(17)	-0.0075(13)	-0.0173(15)	-0.0038(13)
O(5)	0.067(2)	0.0372(16)	0.0543(17)	-0.0072(13)	-0.0037(15)	-0.0108(12)
O(6)	0.0370(16)	0.0477(16)	0.0529(17)	0.0165(12)	0.0036(13)	-0.0012(12)
O(7)	0.0594(19)	0.0406(16)	0.0583(18)	0.0031(14)	-0.0248(16)	0.0085(13)
N(1)	0.0339(19)	0.0256(16)	0.0408(19)	-0.0009(14)	0.0024(16)	-0.0027(13)
C(1)	0.042(2)	0.023(2)	0.041(2)	-0.0027(18)	0.007(2)	-0.0017(18)
C(9)	0.017(2)	0.040(2)	0.032(2)	-0.0024(16)	-0.0016(17)	0.0045(17)
C(11)	0.032(2)	0.029(2)	0.038(2)	-0.0017(17)	-0.0008(19)	-0.0008(16)
C(12)	0.027(2)	0.027(2)	0.034(2)	0.0010(16)	0.0087(18)	0.0011(16)
C(13)	0.038(2)	0.037(2)	0.034(2)	-0.0016(19)	-0.0006(19)	-0.0030(17)
C(14)	0.043(2)	0.035(2)	0.035(2)	0.0003(19)	-0.001(2)	0.0083(18)
C(15)	0.022(2)	0.035(2)	0.039(2)	0.0005(17)	-0.0035(18)	0.0022(17)
C(16)	0.041(2)	0.028(2)	0.047(2)	0.0011(18)	-0.000(2)	0.0027(18)
C(17)	0.038(2)	0.054(2)	0.034(2)	0.008(2)	0.004(2)	-0.003(2)
C(18)	0.033(2)	0.022(2)	0.046(2)	-0.0029(17)	-0.000(2)	-0.0015(18)
C(19)	0.023(2)	0.045(2)	0.046(2)	-0.0065(19)	0.007(2)	-0.000(2)
C(20)	0.028(2)	0.038(2)	0.048(2)	0.003(2)	0.002(2)	0.001(2)
C(21)	0.044(2)	0.039(2)	0.044(2)	-0.009(2)	0.006(2)	-0.000(2)
C(22)	0.033(2)	0.057(3)	0.048(3)	-0.001(2)	-0.005(2)	0.026(2)
C(23)	0.033(2)	0.077(3)	0.032(2)	-0.003(2)	0.003(2)	0.002(2)
C(25)	0.102(4)	0.056(3)	0.148(5)	-0.004(3)	-0.093(3)	0.008(3)
C(26)	0.172(5)	0.061(3)	0.138(5)	0.008(3)	-0.123(4)	0.002(3)

The general temperature factor expression: $\exp(-2\pi^2(a^2U_{11}h^2 + b^2U_{22}k^2 + c^2U_{33}l^2 + 2a*b*U_{12}hk + 2a*c*U_{13}hl + 2b*c*U_{23}kl))$

Table 3. Bond lengths (Å)

atom	atom	distance	atom	atom	distance
Cl(2)	C(23)	1.741(3)	S(1)	O(5)	1.429(2)
S(1)	O(6)	1.425(2)	S(1)	N(1)	1.612(2)
S(1)	C(15)	1.787(3)	O(3)	C(18)	1.450(4)
O(3)	C(19)	1.430(4)	O(4)	C(13)	1.373(4)
O(4)	C(25)	1.413(5)	O(7)	C(14)	1.373(4)
O(7)	C(26)	1.405(5)	N(1)	C(11)	1.457(4)
N(1)	C(21)	1.489(5)	N(2)	C(28)	1.104(8)
N(2)	C(28)	1.104(8)	C(1)	C(12)	1.382(4)
C(1)	C(13)	1.393(4)	C(1)	H(5)	0.91(2)
C(2)	C(28)	1.271(10)	C(2)	H(15)	0.950
C(2)	H(16)	0.950	C(2)	H(17)	0.950
C(9)	C(15)	1.399(4)	C(9)	C(18)	1.523(5)
C(9)	C(20)	1.391(5)	C(11)	C(12)	1.378(4)
C(11)	C(16)	1.381(4)	C(12)	C(18)	1.521(4)
C(13)	C(14)	1.391(4)	C(14)	C(16)	1.379(5)
C(15)	C(17)	1.394(5)	C(16)	H(6)	1.04(2)
C(17)	C(23)	1.367(6)	C(17)	H(3)	1.04(2)
C(18)	H(4)	0.91(2)	C(19)	C(21)	1.516(5)
C(19)	H(7)	1.02(2)	C(19)	H(8)	1.04(2)
C(20)	C(22)	1.372(6)	C(20)	H(1)	0.84(2)
C(21)	H(9)	1.09(2)	C(21)	H(10)	0.92(2)
C(22)	C(23)	1.364(6)	C(22)	H(2)	0.84(2)
C(25)	C(26)	1.347(6)	C(25)	H(11)	0.950
C(25)	H(12)	0.950	C(26)	H(13)	0.950
C(26)	H(14)	0.950			

Symmetry Operators:

(1) -X+2,-Y,-Z

Table 4. Bond angles (°)

atom	atom	atom	angle	atom	atom	atom	angle
O(5)	S(1)	O(6)	118.78(14)	O(5)	S(1)	N(1)	107.57(15)
O(5)	S(1)	C(15)	107.30(14)	O(6)	S(1)	N(1)	107.81(14)
O(6)	S(1)	C(15)	107.01(14)	N(1)	S(1)	C(15)	107.96(14)
C(18)	O(3)	C(19)	116.3(2)	C(13)	O(4)	C(25)	113.0(2)
C(14)	O(7)	C(26)	113.8(2)	S(1)	N(1)	C(11)	114.1(2)
S(1)	N(1)	C(21)	118.8(2)	C(11)	N(1)	C(21)	116.9(2)
C(28)	N(2)	C(28) ¹⁾	0(440)	C(12)	C(1)	C(13)	120.4(3)
C(12)	C(1)	H(5)	118.0(16)	C(13)	C(1)	H(5)	121.5(16)
C(28)	C(2)	H(15)	109.5	C(28)	C(2)	H(16)	108.8
C(28)	C(2)	H(17)	110.1	H(15)	C(2)	H(16)	109.5
H(15)	C(2)	H(17)	109.5	H(16)	C(2)	H(17)	109.5
C(15)	C(9)	C(18)	128.2(3)	C(15)	C(9)	C(20)	115.8(3)
C(18)	C(9)	C(20)	116.0(3)	N(1)	C(11)	C(12)	119.1(2)
N(1)	C(11)	C(16)	119.1(3)	C(12)	C(11)	C(16)	121.8(3)
C(1)	C(12)	C(11)	118.8(3)	C(1)	C(12)	C(18)	120.9(2)
C(11)	C(12)	C(18)	120.4(2)	O(4)	C(13)	C(1)	117.5(3)
O(4)	C(13)	C(14)	122.8(3)	C(1)	C(13)	C(14)	119.7(3)
O(7)	C(14)	C(13)	121.9(3)	O(7)	C(14)	C(16)	118.0(3)
C(13)	C(14)	C(16)	120.0(3)	S(1)	C(15)	C(9)	124.6(2)
S(1)	C(15)	C(17)	113.2(2)	C(9)	C(15)	C(17)	122.2(3)
C(11)	C(16)	C(14)	119.2(3)	C(11)	C(16)	H(6)	119.2(15)
C(14)	C(16)	H(6)	120.3(15)	C(15)	C(17)	C(23)	118.4(3)
C(15)	C(17)	H(3)	120.0(14)	C(23)	C(17)	H(3)	121.6(14)
O(3)	C(18)	C(9)	112.8(2)	O(3)	C(18)	C(12)	112.4(2)
O(3)	C(18)	H(4)	102.2(15)	C(9)	C(18)	C(12)	115.4(2)
C(9)	C(18)	H(4)	106.6(15)	C(12)	C(18)	H(4)	106.2(14)
O(3)	C(19)	C(21)	114.3(3)	O(3)	C(19)	H(7)	98.5(14)
O(3)	C(19)	H(8)	110.4(13)	C(21)	C(19)	H(7)	111.8(14)
C(21)	C(19)	H(8)	108.4(13)	H(7)	C(19)	H(8)	113(2)
C(9)	C(20)	C(22)	122.9(3)	C(9)	C(20)	H(1)	117.6(19)
C(22)	C(20)	H(1)	119.4(19)	N(1)	C(21)	C(19)	114.5(3)
N(1)	C(21)	H(9)	105.0(16)	N(1)	C(21)	H(10)	108.4(18)
C(19)	C(21)	H(9)	108.1(15)	C(19)	C(21)	H(10)	105.1(17)
H(9)	C(21)	H(10)	116(2)	C(20)	C(22)	C(23)	119.0(4)
C(20)	C(22)	H(2)	122.5(19)	C(23)	C(22)	H(2)	118.5(19)
Cl(2)	C(23)	C(17)	119.2(3)	Cl(2)	C(23)	C(22)	119.0(3)
C(17)	C(23)	C(22)	121.7(3)	O(4)	C(25)	C(26)	120.2(4)

Table 4. Bond angles ($^{\circ}$) (continued)

atom	atom	atom	angle	atom	atom	atom	angle
O(4)	C(25)	H(11)	106.2	O(4)	C(25)	H(12)	107.0
C(26)	C(25)	H(11)	104.6	C(26)	C(25)	H(12)	109.1
H(11)	C(25)	H(12)	109.5	O(7)	C(26)	C(25)	120.3(4)
O(7)	C(26)	H(13)	105.8	O(7)	C(26)	H(14)	107.5
C(25)	C(26)	H(13)	104.4	C(25)	C(26)	H(14)	109.0
H(13)	C(26)	H(14)	109.5	N(2)	C(28)	C(2)	174.1(8)

Symmetry Operators:

(1) $-X+2, -Y, -Z$

Table 5. Torsion Angles($^{\circ}$)

atom1	atom2	atom3	atom4	angle	atom1	atom2	atom3	atom4	angle
O(5)	S(1)	N(1)	C(11)	175.6(2)	O(5)	S(1)	N(1)	C(21)	-40.0(2)
O(5)	S(1)	C(15)	C(9)	126.2(2)	O(5)	S(1)	C(15)	C(17)	-53.6(2)
O(6)	S(1)	N(1)	C(11)	46.4(2)	O(6)	S(1)	N(1)	C(21)	-169.2(2)
O(6)	S(1)	C(15)	C(9)	-105.3(2)	O(6)	S(1)	C(15)	C(17)	74.9(2)
N(1)	S(1)	C(15)	C(9)	10.5(3)	N(1)	S(1)	C(15)	C(17)	-169.3(2)
C(15)	S(1)	N(1)	C(11)	-68.9(2)	C(15)	S(1)	N(1)	C(21)	75.5(2)
C(18)	O(3)	C(19)	C(21)	64.5(4)	C(19)	O(3)	C(18)	C(9)	48.5(3)
C(19)	O(3)	C(18)	C(12)	-84.0(3)	C(13)	O(4)	C(25)	C(26)	-26.0(5)
C(25)	O(4)	C(13)	C(1)	-168.6(3)	C(25)	O(4)	C(13)	C(14)	10.3(4)
C(14)	O(7)	C(26)	C(25)	-24.9(6)	C(26)	O(7)	C(14)	C(13)	9.0(5)
C(26)	O(7)	C(14)	C(16)	-173.0(3)	S(1)	N(1)	C(11)	C(12)	81.4(3)
S(1)	N(1)	C(11)	C(16)	-96.7(3)	S(1)	N(1)	C(21)	C(19)	-64.9(3)
C(11)	N(1)	C(21)	C(19)	78.5(4)	C(21)	N(1)	C(11)	C(12)	-63.6(4)
C(21)	N(1)	C(11)	C(16)	118.2(3)	C(12)	C(1)	C(13)	O(4)	179.3(3)
C(12)	C(1)	C(13)	C(14)	0.5(5)	C(13)	C(1)	C(12)	C(11)	2.0(5)
C(13)	C(1)	C(12)	C(18)	-177.5(3)	C(15)	C(9)	C(18)	O(3)	-82.0(4)
C(15)	C(9)	C(18)	C(12)	49.0(4)	C(18)	C(9)	C(15)	S(1)	-2.9(4)
C(18)	C(9)	C(15)	C(17)	176.9(3)	C(15)	C(9)	C(20)	C(22)	-0.8(5)
C(20)	C(9)	C(15)	S(1)	180.0(2)	C(20)	C(9)	C(15)	C(17)	-0.2(4)
C(18)	C(9)	C(20)	C(22)	-178.2(3)	C(20)	C(9)	C(18)	O(3)	95.1(3)
C(20)	C(9)	C(18)	C(12)	-133.9(3)	N(1)	C(11)	C(12)	C(1)	179.0(3)
N(1)	C(11)	C(12)	C(18)	-1.6(4)	N(1)	C(11)	C(16)	C(14)	179.6(3)
C(12)	C(11)	C(16)	C(14)	1.5(5)	C(16)	C(11)	C(12)	C(1)	-3.0(5)
C(16)	C(11)	C(12)	C(18)	176.5(3)	C(1)	C(12)	C(18)	O(3)	-114.7(3)
C(1)	C(12)	C(18)	C(9)	114.2(3)	C(11)	C(12)	C(18)	O(3)	65.9(4)
C(11)	C(12)	C(18)	C(9)	-65.3(4)	O(4)	C(13)	C(14)	O(7)	-2.8(5)
O(4)	C(13)	C(14)	C(16)	179.2(3)	C(1)	C(13)	C(14)	O(7)	176.0(3)
C(1)	C(13)	C(14)	C(16)	-2.0(5)	O(7)	C(14)	C(16)	C(11)	-177.0(3)
C(13)	C(14)	C(16)	C(11)	1.0(5)	S(1)	C(15)	C(17)	C(23)	-179.1(2)
C(9)	C(15)	C(17)	C(23)	1.1(5)	C(15)	C(17)	C(23)	Cl(2)	179.0(2)
C(15)	C(17)	C(23)	C(22)	-1.0(5)	O(3)	C(19)	C(21)	N(1)	-59.0(4)
C(9)	C(20)	C(22)	C(23)	0.9(5)	C(20)	C(22)	C(23)	Cl(2)	-179.93(19)
C(20)	C(22)	C(23)	C(17)	0.1(4)	O(4)	C(25)	C(26)	O(7)	35.0(7)

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 6. Distances beyond the asymmetric unit out to 3.60 Å

atom	atom	distance	atom	atom	distance
Cl(2)	H(2) ¹⁾	3.52(2)	Cl(2)	H(11) ²⁾	3.408
Cl(2)	H(12) ²⁾	3.172	Cl(2)	H(12) ³⁾	3.198
Cl(2)	H(13) ⁴⁾	3.429	Cl(2)	H(14) ²⁾	3.358
Cl(2)	H(15)	2.953	Cl(2)	H(16) ⁵⁾	3.588
O(3)	C(1) ⁶⁾	3.284(4)	O(3)	C(18) ⁶⁾	3.511(4)
O(3)	C(20) ⁶⁾	3.566(4)	O(3)	H(1) ⁶⁾	3.08(2)
O(3)	H(4) ⁶⁾	2.77(2)	O(3)	H(5) ⁶⁾	2.51(2)
O(3)	H(11) ⁵⁾	3.479	O(4)	C(1) ¹⁾	3.363(4)
O(4)	H(5) ¹⁾	2.82(2)	O(4)	H(11) ⁶⁾	3.171
O(4)	H(12) ⁶⁾	3.560	O(5)	N(2) ⁵⁾	3.108(2)
O(5)	C(17) ⁷⁾	3.476(4)	O(5)	C(28) ⁵⁾	3.221(8)
O(5)	C(28) ⁷⁾	3.374(8)	O(5)	H(3) ⁷⁾	2.46(2)
O(5)	H(17) ⁵⁾	3.312	O(6)	N(2)	3.210(2)
O(6)	C(19) ⁸⁾	3.319(4)	O(6)	C(26) ⁹⁾	3.561(5)
O(6)	C(28)	3.393(8)	O(6)	C(28) ¹⁰⁾	3.397(8)
O(6)	H(7) ⁸⁾	2.81(2)	O(6)	H(8) ⁸⁾	3.28(2)
O(6)	H(9) ⁸⁾	3.00(2)	O(6)	H(13) ⁹⁾	3.487
O(6)	H(14) ⁹⁾	3.010	O(7)	C(21) ¹¹⁾	3.349(4)
O(7)	H(6) ⁹⁾	2.70(2)	O(7)	H(9) ¹¹⁾	2.64(3)
O(7)	H(10) ⁸⁾	2.85(2)	O(7)	H(10) ¹¹⁾	3.22(2)
N(1)	H(6) ¹¹⁾	3.53(2)	N(2)	O(5) ⁸⁾	3.108(2)
N(2)	O(5) ⁷⁾	3.108(2)	N(2)	O(6)	3.210(2)
N(2)	O(6) ¹⁰⁾	3.210(2)	N(2)	H(3)	3.29(2)
N(2)	H(3) ¹⁰⁾	3.29(2)	C(1)	O(3) ¹⁾	3.284(4)
C(1)	O(4) ⁶⁾	3.363(4)	C(1)	C(18) ¹⁾	3.594(5)
C(1)	H(4) ¹⁾	2.96(2)	C(1)	H(7) ⁸⁾	3.17(2)
C(1)	H(11) ⁶⁾	3.476	C(2)	C(17)	3.403(8)
C(2)	C(17) ⁸⁾	3.411(8)	C(2)	C(22) ¹⁾	3.566(8)
C(2)	C(26) ⁴⁾	3.575(9)	C(2)	H(2) ¹⁾	2.85(2)
C(2)	H(3)	3.15(2)	C(2)	H(8) ⁸⁾	2.84(2)
C(2)	H(12) ⁴⁾	3.527	C(2)	H(13) ⁴⁾	2.751
C(9)	H(1) ¹⁾	3.35(2)	C(9)	H(17) ⁵⁾	3.198
C(11)	H(7) ⁸⁾	3.06(2)	C(12)	H(4) ¹⁾	3.58(2)
C(12)	H(7) ⁸⁾	3.12(2)	C(13)	H(4) ¹⁾	3.59(2)
C(13)	H(5) ¹⁾	3.42(2)	C(13)	H(7) ⁸⁾	3.14(2)
C(14)	H(7) ⁸⁾	3.04(2)	C(14)	H(9) ¹¹⁾	3.40(3)
C(14)	H(10) ⁸⁾	3.06(2)	C(15)	H(15)	3.432

Table 6. Distances beyond the asymmetric unit out to 3.60 Å (continued)

atom	atom	distance	atom	atom	distance
C(15)	H(17) ^{5j}	2.874	C(16)	H(7) ^{8j}	3.01(2)
C(16)	H(9) ^{11j}	3.56(2)	C(16)	H(10) ^{8j}	3.39(2)
C(16)	H(10) ^{11j}	3.24(2)	C(17)	O(5) ^{7j}	3.476(4)
C(17)	C(2) ^{5j}	3.411(8)	C(17)	C(2)	3.403(8)
C(17)	C(28)	3.571(9)	C(17)	H(15)	2.581
C(17)	H(16) ^{5j}	3.145	C(17)	H(17) ^{5j}	2.894
C(18)	O(3) ^{1j}	3.511(4)	C(18)	C(1) ^{6j}	3.594(5)
C(18)	H(5) ^{6j}	3.00(2)	C(19)	O(6) ^{5j}	3.319(4)
C(19)	H(1) ^{6j}	3.23(2)	C(19)	H(4) ^{6j}	3.55(2)
C(19)	H(17) ^{5j}	3.035	C(20)	O(3) ^{1j}	3.566(4)
C(20)	C(20) ^{6j}	3.430(5)	C(20)	C(20) ^{1j}	3.430(5)
C(20)	C(22) ^{6j}	3.432(5)	C(20)	H(1) ^{6j}	3.58(2)
C(20)	H(1) ^{1j}	3.14(2)	C(20)	H(2) ^{6j}	3.55(2)
C(20)	H(8) ^{1j}	3.58(2)	C(20)	H(17) ^{5j}	3.458
C(20)	H(17) ^{6j}	3.479	C(21)	O(7) ^{11j}	3.349(4)
C(21)	H(6) ^{11j}	3.11(2)	C(22)	C(2) ^{6j}	3.566(8)
C(22)	C(20) ^{1j}	3.432(5)	C(22)	C(22) ^{6j}	3.442(5)
C(22)	C(22) ^{1j}	3.442(5)	C(22)	H(1) ^{1j}	3.30(2)
C(22)	H(2) ^{6j}	3.37(2)	C(22)	H(2) ^{1j}	3.38(2)
C(22)	H(15) ^{6j}	3.189	C(22)	H(17) ^{5j}	3.495
C(22)	H(17) ^{6j}	3.279	C(23)	H(2) ^{1j}	3.33(2)
C(23)	H(15)	2.728	C(23)	H(16) ^{5j}	3.180
C(23)	H(17) ^{5j}	3.217	C(25)	H(5) ^{1j}	3.30(2)
C(25)	H(16) ^{12j}	3.217	C(26)	O(6) ^{9j}	3.561(5)
C(26)	C(2) ^{12j}	3.575(9)	C(26)	H(6) ^{9j}	3.37(3)
C(26)	H(9) ^{11j}	3.36(3)	C(26)	H(10) ^{8j}	3.54(2)
C(26)	H(16) ^{12j}	2.767	C(28)	O(5) ^{8j}	3.221(8)
C(28)	O(5) ^{7j}	3.374(8)	C(28)	O(6)	3.393(8)
C(28)	O(6) ^{10j}	3.397(8)	C(28)	C(17)	3.571(9)
C(28)	H(3)	3.00(2)	C(28)	H(8) ^{8j}	3.13(2)
C(28)	H(13) ^{4j}	3.124	H(1)	O(3) ^{1j}	3.08(2)
H(1)	C(9) ^{6j}	3.35(2)	H(1)	C(19) ^{1j}	3.23(2)
H(1)	C(20) ^{6j}	3.14(2)	H(1)	C(20) ^{1j}	3.58(2)
H(1)	C(22) ^{6j}	3.30(2)	H(1)	H(1) ^{6j}	3.41(4)
H(1)	H(1) ^{1j}	3.41(4)	H(1)	H(7) ^{1j}	3.06(3)
H(1)	H(8) ^{1j}	2.95(3)	H(1)	H(15) ^{6j}	3.248
H(1)	H(17) ^{6j}	3.078	H(2)	Cl(2) ^{6j}	3.52(2)

Table 6. Distances beyond the asymmetric unit out to 3.60 Å (continued)

atom	atom	distance	atom	atom	distance
H(2)	C(2) ^{6j}	2.85(2)	H(2)	C(20) ^{1j}	3.55(2)
H(2)	C(22) ^{6j}	3.38(2)	H(2)	C(22) ^{1j}	3.37(2)
H(2)	C(23) ^{6j}	3.33(2)	H(2)	H(2) ^{6j}	3.39(4)
H(2)	H(2) ^{1j}	3.39(4)	H(2)	H(12) ^{3j}	3.291
H(2)	H(15) ^{6j}	2.463	H(2)	H(16) ^{6j}	2.895
H(2)	H(17) ^{6j}	2.708	H(3)	O(5) ^{7j}	2.46(2)
H(3)	N(2)	3.29(2)	H(3)	C(2)	3.15(2)
H(3)	C(28)	3.00(2)	H(3)	H(15)	2.473
H(3)	H(16) ^{5j}	3.289	H(3)	H(17) ^{5j}	3.324
H(4)	O(3) ^{1j}	2.77(2)	H(4)	C(1) ^{6j}	2.96(2)
H(4)	C(12) ^{6j}	3.58(2)	H(4)	C(13) ^{6j}	3.59(2)
H(4)	C(19) ^{1j}	3.55(2)	H(4)	H(4) ^{6j}	3.45(3)
H(4)	H(4) ^{1j}	3.45(3)	H(4)	H(5) ^{6j}	2.54(3)
H(4)	H(7) ^{1j}	3.16(3)	H(5)	O(3) ^{1j}	2.51(2)
H(5)	O(4) ^{6j}	2.82(2)	H(5)	C(13) ^{6j}	3.42(2)
H(5)	C(18) ^{1j}	3.00(2)	H(5)	C(25) ^{6j}	3.30(2)
H(5)	H(4) ^{1j}	2.54(3)	H(5)	H(5) ^{6j}	3.41(3)
H(5)	H(5) ^{1j}	3.41(3)	H(5)	H(11) ^{6j}	2.793
H(6)	O(7) ^{9j}	2.70(2)	H(6)	N(1) ^{11j}	3.53(2)
H(6)	C(21) ^{11j}	3.11(2)	H(6)	C(26) ^{9j}	3.37(3)
H(6)	H(6) ^{9j}	3.55(4)	H(6)	H(7) ^{8j}	3.40(3)
H(6)	H(9) ^{8j}	3.52(4)	H(6)	H(9) ^{11j}	3.24(4)
H(6)	H(10) ^{8j}	3.33(4)	H(6)	H(10) ^{11j}	2.47(4)
H(6)	H(14) ^{9j}	2.972	H(7)	O(6) ^{5j}	2.81(2)
H(7)	C(1) ^{5j}	3.17(2)	H(7)	C(11) ^{5j}	3.06(2)
H(7)	C(12) ^{5j}	3.12(2)	H(7)	C(13) ^{5j}	3.14(2)
H(7)	C(14) ^{5j}	3.04(2)	H(7)	C(16) ^{5j}	3.01(2)
H(7)	H(1) ^{6j}	3.06(3)	H(7)	H(4) ^{6j}	3.16(3)
H(7)	H(6) ^{5j}	3.40(3)	H(7)	H(17) ^{5j}	3.454
H(8)	O(6) ^{5j}	3.28(2)	H(8)	C(2) ^{5j}	2.84(2)
H(8)	C(20) ^{6j}	3.58(2)	H(8)	C(28) ^{5j}	3.13(2)
H(8)	H(1) ^{6j}	2.95(3)	H(8)	H(15) ^{5j}	3.196
H(8)	H(16) ^{5j}	3.550	H(8)	H(17) ^{5j}	2.015
H(9)	O(6) ^{5j}	3.00(2)	H(9)	O(7) ^{11j}	2.64(3)
H(9)	C(14) ^{11j}	3.40(3)	H(9)	C(16) ^{11j}	3.56(2)
H(9)	C(26) ^{11j}	3.36(3)	H(9)	H(6) ^{5j}	3.52(4)
H(9)	H(6) ^{11j}	3.24(4)	H(9)	H(13) ^{11j}	3.079

Table 6. Distances beyond the asymmetric unit out to 3.60 Å (continued)

atom	atom	distance	atom	atom	distance
H(10)	O(7) ⁵⁾	2.85(2)	H(10)	O(7) ¹¹⁾	3.22(2)
H(10)	C(14) ⁵⁾	3.06(2)	H(10)	C(16) ⁵⁾	3.39(2)
H(10)	C(16) ¹¹⁾	3.24(2)	H(10)	C(26) ⁵⁾	3.54(2)
H(10)	H(6) ⁵⁾	3.33(4)	H(10)	H(6) ¹¹⁾	2.47(4)
H(10)	H(11) ⁵⁾	3.402	H(10)	H(14) ⁵⁾	3.475
H(11)	Cl(2) ¹³⁾	3.408	H(11)	O(3) ⁸⁾	3.479
H(11)	O(4) ¹⁾	3.171	H(11)	C(1) ¹⁾	3.476
H(11)	H(5) ¹⁾	2.793	H(11)	H(10) ⁸⁾	3.402
H(12)	Cl(2) ¹³⁾	3.172	H(12)	Cl(2) ¹⁴⁾	3.198
H(12)	O(4) ¹⁾	3.560	H(12)	C(2) ¹²⁾	3.527
H(12)	H(2) ¹⁴⁾	3.291	H(12)	H(15) ¹²⁾	3.546
H(12)	H(16) ¹²⁾	2.718	H(13)	Cl(2) ¹²⁾	3.429
H(13)	O(6) ⁹⁾	3.487	H(13)	C(2) ¹²⁾	2.751
H(13)	C(28) ¹²⁾	3.124	H(13)	H(9) ¹¹⁾	3.079
H(13)	H(15) ¹²⁾	2.821	H(13)	H(16) ¹²⁾	2.058
H(13)	H(17) ¹²⁾	3.565	H(14)	Cl(2) ¹³⁾	3.358
H(14)	O(6) ⁹⁾	3.010	H(14)	H(6) ⁹⁾	2.972
H(14)	H(10) ⁸⁾	3.475	H(14)	H(16) ¹²⁾	2.777
H(15)	Cl(2)	2.953	H(15)	C(15)	3.432
H(15)	C(17)	2.581	H(15)	C(22) ¹⁾	3.189
H(15)	C(23)	2.728	H(15)	H(1) ¹⁾	3.248
H(15)	H(2) ¹⁾	2.463	H(15)	H(3)	2.473
H(15)	H(8) ⁸⁾	3.196	H(15)	H(12) ⁴⁾	3.546
H(15)	H(13) ⁴⁾	2.821	H(16)	Cl(2) ⁸⁾	3.588
H(16)	C(17) ⁸⁾	3.145	H(16)	C(23) ⁸⁾	3.180
H(16)	C(25) ⁴⁾	3.217	H(16)	C(26) ⁴⁾	2.767
H(16)	H(2) ¹⁾	2.895	H(16)	H(3) ⁸⁾	3.289
H(16)	H(8) ⁸⁾	3.550	H(16)	H(12) ⁴⁾	2.718
H(16)	H(13) ⁴⁾	2.058	H(16)	H(14) ⁴⁾	2.777
H(17)	O(5) ⁸⁾	3.312	H(17)	C(9) ⁸⁾	3.198
H(17)	C(15) ⁸⁾	2.874	H(17)	C(17) ⁸⁾	2.894
H(17)	C(19) ⁸⁾	3.035	H(17)	C(20) ⁸⁾	3.458
H(17)	C(20) ¹⁾	3.479	H(17)	C(22) ⁸⁾	3.495
H(17)	C(22) ¹⁾	3.279	H(17)	C(23) ⁸⁾	3.217
H(17)	H(1) ¹⁾	3.078	H(17)	H(2) ¹⁾	2.708
H(17)	H(3) ⁸⁾	3.324	H(17)	H(7) ⁸⁾	3.454
H(17)	H(8) ⁸⁾	2.015	H(17)	H(13) ⁴⁾	3.565

Symmetry Operators:

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