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## Structure Reports

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# Low-temperature redetermination of 4-chloro-2-[tris(hydroxymethyl)methyl-aminomethyl]phenol as zwitterionic 4-chloro-2-[tris(hydroxymethyl)methyl-iminiomethyl]phenolate

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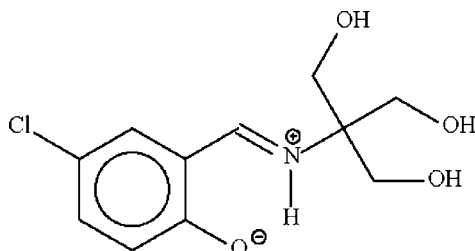
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 Key indicators: single-crystal X-ray study;  $T = 100$  K; mean  $\sigma(\text{C}-\text{C}) = 0.003$  Å; disorder in main residue;  $R$  factor = 0.042;  $wR$  factor = 0.117; data-to-parameter ratio = 14.9.

The title Schiff base,  $\text{C}_{11}\text{H}_{14}\text{ClNO}_4$ , originally refined as a neutral molecule [Chumakov, Antosyak, Mazus, Tsapkov & Samus (2000), *Crystallogr. Rep.* **45**, 945–950], is interpreted as a zwitterionic compound. There are two independent zwitterions in the asymmetric unit. Five of the six hydroxy groups of the  $-\text{CH}_2\text{OH}$  fragments are disordered. In one zwitterion, two are disordered over three sites [0.53 (1):0.33 (1):0.14 (1) and 0.65 (1):0.18 (1):0.17 (1)] and the third over two sites [0.66 (1):0.34 (1)]. In the second zwitterion, two are disordered over two sites [0.84 (1):0.16 (1) and 0.83 (1):0.17 (1)] and the third hydroxy group is ordered.

## Related literature

The room-temperature crystal structure was interpreted as a neutral molecule; the disordered structure (the disorder confined to the hydroxy part of one of the six  $-\text{CH}_2\text{OH}$  groups) refined to 0.072, see: Chumakov *et al.* (2000). The unsubstituted parent Schiff base is also a zwitterion; its structure has been determined several times, see: Asgedom *et al.* (1996); Odabaşoğlu *et al.* (2003); Tatar *et al.* (2005); Zhang *et al.* (2000).



## Experimental

## Crystal data

$\text{C}_{11}\text{H}_{14}\text{ClNO}_4$	$\gamma = 85.258 (1)^\circ$
$M_r = 259.68$	$V = 1172.27 (2) \text{ \AA}^3$
Triclinic, $P\bar{1}$	$Z = 4$
$a = 7.0174 (1) \text{ \AA}$	Mo $K\alpha$ radiation
$b = 10.1935 (1) \text{ \AA}$	$\mu = 0.33 \text{ mm}^{-1}$
$c = 16.7234 (2) \text{ \AA}$	$T = 100 (2) \text{ K}$
$\alpha = 79.520 (1)^\circ$	$0.30 \times 0.20 \times 0.05 \text{ mm}$
$\beta = 89.031 (1)^\circ$	

## Data collection

Bruker SMART APEX diffractometer	11241 measured reflections
Absorption correction: multi-scan (SADABS; Sheldrick, 1996)	5361 independent reflections
$T_{\min} = 0.908$ , $T_{\max} = 0.984$	4771 reflections with $I > 2\sigma(I)$
	$R_{\text{int}} = 0.018$

## Refinement

$R[F^2 > 2\sigma(F^2)] = 0.042$	H atoms treated by a mixture of independent and constrained refinement
$wR(F^2) = 0.117$	$\Delta\rho_{\text{max}} = 0.41 \text{ e \AA}^{-3}$
$S = 1.01$	$\Delta\rho_{\text{min}} = -0.74 \text{ e \AA}^{-3}$
5361 reflections	
360 parameters	
47 restraints	

Data collection: *APEX2* (Bruker, 2007); cell refinement: *SAINT* (Bruker, 2007); data reduction: *SAINT*; program(s) used to solve structure: *SHELXS97* (Sheldrick, 2008); program(s) used to refine structure: *SHELXL97* (Sheldrick, 2008); molecular graphics: *X-SEED* (Barbour, 2001); software used to prepare material for publication: *pubCIF* (Westrip, 2008).

I thank the University of Malaya for supporting this study.

Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: TK2330).

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## supporting information

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## Low-temperature redetermination of 4-chloro-2-[tris(hydroxymethyl)methyl-iminomethyl]phenol as zwitterionic 4-chloro-2-[tris(hydroxymethyl)methyl-iminiomethyl]phenolate

Seik Weng Ng

### S1. Comment

A reevaluation of the title compound, originally reported as a neutral molecule (Chumakov *et al.*, 2000), shows a zwitterionic compound, at low temperature, Fig. 1.

### S2. Experimental

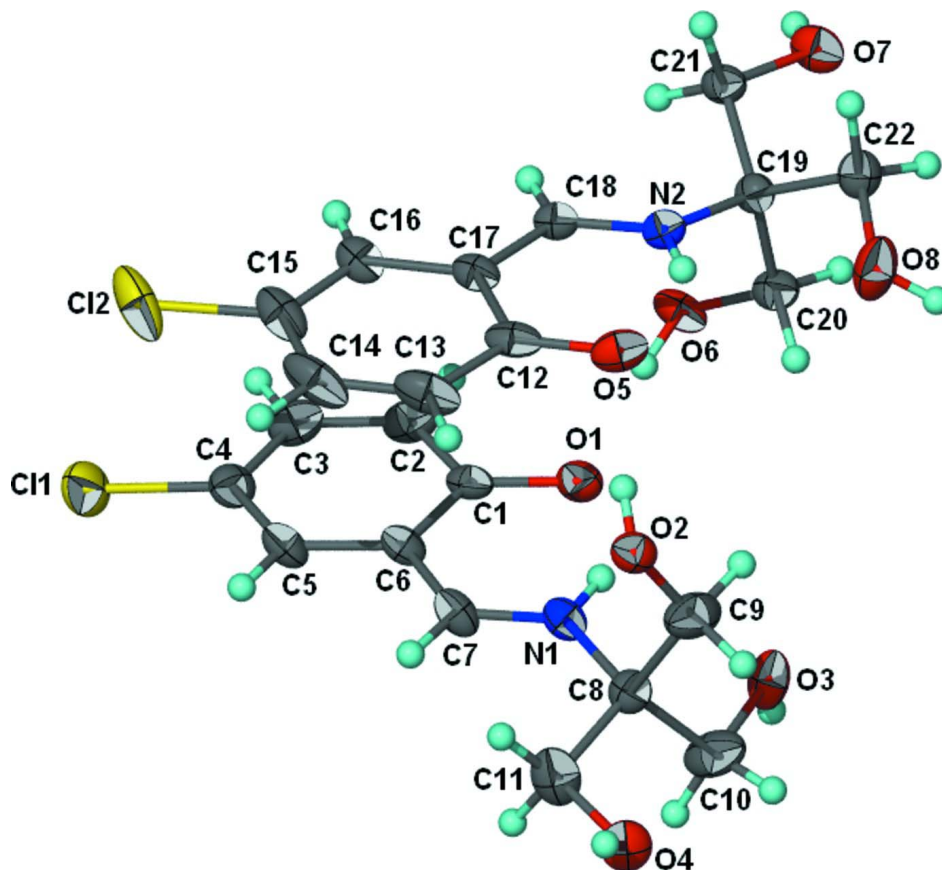
The compound was synthesized from the reaction of 5-chlorosalicylaldehyde and tris(hydroxymethyl)aminomethane (Chumakov *et al.*, 2000), and was recrystallized from ethanol.

### S3. Refinement

Five of the six hydroxy groups of the  $-\text{CH}_2\text{OH}$  fragments are disordered. In one zwitterion, two are disordered over three sites [0.53 (1): 0.33: 0.14] and the third over two sites [0.65 (1): 0.18: 0.17]. In the second zwitterion, two are disordered over two sites [0.84 (1): 0.16 and 0.83 (1): 0.17] and the third is ordered. For each disordered  $-\text{CH}_2\text{OH}$  fragment, the C–O distances (of the unprimed and primed atoms) were restrained to within 0.01 Å. The temperature factors of the primed atom(s) were restrained to those of the unprimed ones.

Carbon- and oxygen-bound hydrogen atoms were placed at calculated positions (C–H 0.95–0.99, O–H 0.84 Å) and were treated as riding on their parent atoms, with  $U(\text{H})$  set to 1.2–1.5 times  $U_{\text{eq}}(\text{C}, \text{O})$ . The iminium hydrogen atoms were located in a difference Fourier map, and were refined with a distance restraint of N–H 0.88±0.01 Å; their temperature factors were freely refined.

Owing to the disorder, hydrogen bonding interactions involving the hydroxyl groups were not computed.



**Figure 1**

Thermal ellipsoid plot (Barbour, 2001) of  $C_{22}H_{28}Cl_2N_2O_8$  at the 70% probability level. Hydrogen atoms are drawn as spheres of arbitrary radius. The minor disorder components are not shown for clarity.

#### 4-chloro-2-[tris(hydroxymethyl)methyliminomethyl]phenolate

##### Crystal data

$C_{11}H_{14}ClNO_4$

$M_r = 259.68$

Triclinic,  $P\bar{1}$

Hall symbol:  $-P\ 1$

$a = 7.0174$  (1) Å

$b = 10.1935$  (1) Å

$c = 16.7234$  (2) Å

$\alpha = 79.520$  (1)°

$\beta = 89.031$  (1)°

$\gamma = 85.258$  (1)°

$V = 1172.27$  (2) Å<sup>3</sup>

$Z = 4$

$F(000) = 544$

$D_x = 1.471$  Mg m<sup>-3</sup>

Mo  $K\alpha$  radiation,  $\lambda = 0.71073$  Å

Cell parameters from 6864 reflections

$\theta = 2.2\text{--}28.3^\circ$

$\mu = 0.33$  mm<sup>-1</sup>

$T = 100$  K

Prism, yellow

$0.30 \times 0.20 \times 0.05$  mm

##### Data collection

Bruker SMART APEX  
diffractometer

Radiation source: fine-focus sealed tube

Graphite monochromator

$\omega$  scans

Absorption correction: multi-scan  
(*SADABS*; Sheldrick, 1996)

$T_{\min} = 0.908$ ,  $T_{\max} = 0.984$

11241 measured reflections

5361 independent reflections

4771 reflections with  $I > 2\sigma(I)$

$R_{\text{int}} = 0.018$   
 $\theta_{\text{max}} = 27.5^\circ$ ,  $\theta_{\text{min}} = 1.2^\circ$   
 $h = -9 \rightarrow 9$

$k = -13 \rightarrow 13$   
 $l = -21 \rightarrow 21$

*Refinement*

Refinement on  $F^2$   
 Least-squares matrix: full  
 $R[F^2 > 2\sigma(F^2)] = 0.042$   
 $wR(F^2) = 0.117$   
 $S = 1.01$   
 5361 reflections  
 360 parameters  
 47 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  
 $w = 1/[\sigma^2(F_o^2) + (0.0573P)^2 + 1.0948P]$   
 where  $P = (F_o^2 + 2F_c^2)/3$   
 $(\Delta/\sigma)_{\text{max}} = 0.001$   
 $\Delta\rho_{\text{max}} = 0.41 \text{ e } \text{\AA}^{-3}$   
 $\Delta\rho_{\text{min}} = -0.74 \text{ e } \text{\AA}^{-3}$

*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)
Cl1	0.74259 (7)	0.67565 (5)	0.56473 (3)	0.03176 (13)	
Cl2	0.23984 (7)	0.77985 (6)	0.56151 (3)	0.03809 (15)	
O1	0.54240 (19)	0.73990 (13)	0.21969 (8)	0.0239 (3)	
O5	0.03446 (18)	0.95420 (12)	0.21807 (8)	0.0246 (3)	
O2	0.1961 (3)	1.1785 (2)	0.17629 (14)	0.0204 (4)	0.535 (2)
H2	0.1471	1.1063	0.1939	0.031*	0.535 (2)
O2'	0.2372 (5)	1.0814 (4)	0.0932 (2)	0.0204 (4)	0.328 (3)
H2'	0.1688	1.0397	0.1292	0.031*	0.328 (3)
O2''	0.3318 (13)	1.2743 (9)	0.0813 (5)	0.0204 (4)	0.137 (3)
H2''	0.4077	1.3189	0.1015	0.031*	0.137 (3)
O3	0.6388 (3)	0.98149 (19)	0.05581 (12)	0.0256 (4)	0.650 (3)
H3	0.7445	0.9411	0.0476	0.038*	0.650 (3)
O3'	0.8488 (9)	1.1295 (10)	0.0883 (6)	0.0256 (4)	0.183 (6)
H3'	0.8968	1.0856	0.1317	0.038*	0.183 (6)
O3''	0.8481 (9)	1.0984 (11)	0.0690 (7)	0.0256 (4)	0.166 (6)
H3''	0.8995	1.0227	0.0894	0.038*	0.166 (6)
O4	0.6047 (3)	1.35087 (18)	0.13329 (12)	0.0254 (5)	0.660 (5)
H4	0.5375	1.4132	0.1492	0.038*	0.660 (5)
O4'	0.5213 (6)	1.2777 (4)	0.2551 (2)	0.0254 (5)	0.34
H4'	0.4631	1.3531	0.2394	0.038*	0.340 (5)
O6	0.3862 (2)	0.54296 (15)	0.17552 (10)	0.0241 (5)	0.835 (6)
H6	0.4403	0.6031	0.1925	0.036*	0.835 (6)
O6'	0.2819 (11)	0.5226 (7)	0.0539 (5)	0.0241 (5)	0.16
H6'	0.1780	0.4913	0.0462	0.036*	0.165 (6)
O7	-0.0249 (2)	0.40172 (15)	0.12281 (10)	0.0251 (5)	0.831 (6)
H7	0.0171	0.3226	0.1412	0.038*	0.831 (6)
O7'	0.1139 (11)	0.3978 (7)	0.2403 (4)	0.0251 (5)	0.17
H7'	0.2264	0.4154	0.2276	0.038*	0.169 (6)
O8	-0.0449 (2)	0.81865 (13)	0.05073 (8)	0.0336 (3)	
H8	-0.0069	0.8306	0.0023	0.050*	

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N1	0.5472 (2)	0.99524 (14)	0.21010 (9)	0.0194 (3)	
H1N	0.537 (4)	0.9209 (17)	0.1907 (15)	0.043 (7)*	
N2	0.0469 (2)	0.70644 (14)	0.20403 (9)	0.0174 (3)	
H2N	0.029 (3)	0.7936 (11)	0.1854 (14)	0.033 (6)*	
C1	0.5840 (2)	0.72652 (17)	0.29646 (11)	0.0192 (3)	
C2	0.6121 (3)	0.59874 (18)	0.34689 (12)	0.0271 (4)	
H2A	0.5990	0.5209	0.3243	0.032*	
C3	0.6579 (3)	0.58514 (18)	0.42746 (12)	0.0268 (4)	
H3A	0.6760	0.4983	0.4598	0.032*	
C4	0.6782 (2)	0.69780 (18)	0.46259 (11)	0.0232 (4)	
C5	0.6502 (2)	0.82333 (18)	0.41675 (11)	0.0224 (3)	
H5	0.6619	0.8996	0.4411	0.027*	
C6	0.6041 (2)	0.84008 (16)	0.33387 (11)	0.0188 (3)	
C7	0.5840 (2)	0.97254 (17)	0.28681 (11)	0.0216 (3)	
H7A	0.5984	1.0463	0.3131	0.026*	
C8	0.5387 (2)	1.12384 (17)	0.15371 (10)	0.0203 (3)	
C9	0.3328 (3)	1.1578 (2)	0.12391 (13)	0.0329 (4)	
H9A	0.3316	1.2392	0.0813	0.039*	0.535 (2)
H9B	0.2978	1.0842	0.0972	0.039*	0.535 (2)
H9C	0.2577	1.1804	0.1709	0.039*	0.328 (3)
H9D	0.3335	1.2417	0.0834	0.039*	0.328 (3)
H9E	0.2957	1.0929	0.0908	0.039*	0.137 (3)
H9F	0.2433	1.1569	0.1704	0.039*	0.137 (3)
C10	0.6699 (3)	1.1019 (2)	0.08285 (13)	0.0338 (5)	
H10A	0.6471	1.1782	0.0373	0.041*	0.650 (3)
H10B	0.8048	1.0987	0.1000	0.041*	0.650 (3)
H10C	0.6697	1.0070	0.0769	0.041*	0.183 (6)
H10D	0.6146	1.1573	0.0325	0.041*	0.183 (6)
H10E	0.6352	1.0150	0.0712	0.041*	0.166 (6)
H10F	0.6163	1.1698	0.0371	0.041*	0.166 (6)
C11	0.6064 (3)	1.23560 (18)	0.19279 (12)	0.0289 (4)	
H11A	0.5204	1.2514	0.2382	0.035*	0.660 (5)
H11B	0.7372	1.2106	0.2146	0.035*	0.660 (5)
H11C	0.7408	1.2083	0.2092	0.035*	0.340 (5)
H11D	0.6095	1.3150	0.1489	0.035*	0.340 (5)
C12	0.0782 (2)	0.91510 (17)	0.29366 (12)	0.0216 (3)	
C13	0.1062 (3)	1.00731 (19)	0.34634 (14)	0.0292 (4)	
H13	0.0916	1.1006	0.3253	0.035*	
C14	0.1531 (3)	0.9650 (2)	0.42592 (15)	0.0335 (5)	
H14	0.1714	1.0291	0.4593	0.040*	
C15	0.1751 (3)	0.8282 (2)	0.45953 (12)	0.0271 (4)	
C16	0.1489 (2)	0.73448 (18)	0.41255 (11)	0.0219 (3)	
H16	0.1619	0.6419	0.4357	0.026*	
C17	0.1027 (2)	0.77585 (16)	0.32963 (11)	0.0187 (3)	
C18	0.0882 (2)	0.67582 (16)	0.28079 (10)	0.0175 (3)	
H18	0.1095	0.5842	0.3053	0.021*	
C19	0.0507 (2)	0.61584 (16)	0.14497 (10)	0.0172 (3)	
C20	0.2570 (2)	0.60202 (17)	0.11302 (11)	0.0213 (3)	

H20A	0.2621	0.5465	0.0701	0.026*	0.835 (6)
H20B	0.2959	0.6917	0.0884	0.026*	0.835 (6)
H20C	0.2963	0.6924	0.0902	0.026*	0.165 (6)
H20D	0.3425	0.5636	0.1593	0.026*	0.165 (6)
C21	-0.0160 (2)	0.47961 (16)	0.18370 (10)	0.0197 (3)	
H21A	0.0747	0.4344	0.2267	0.024*	0.831 (6)
H21B	-0.1436	0.4914	0.2088	0.024*	0.831 (6)
H21C	-0.1389	0.4944	0.2119	0.024*	0.169 (6)
H21D	-0.0407	0.4299	0.1398	0.024*	0.169 (6)
C22	-0.0847 (3)	0.68282 (17)	0.07586 (11)	0.0229 (3)	
H22A	-0.0676	0.6350	0.0294	0.027*	
H22B	-0.2189	0.6784	0.0947	0.027*	

*Atomic displacement parameters (Å<sup>2</sup>)*

	$U^{11}$	$U^{22}$	$U^{33}$	$U^{12}$	$U^{13}$	$U^{23}$
C11	0.0346 (3)	0.0301 (2)	0.0279 (2)	-0.00246 (19)	-0.01100 (18)	0.00271 (18)
C12	0.0343 (3)	0.0474 (3)	0.0384 (3)	0.0139 (2)	-0.0164 (2)	-0.0295 (2)
O1	0.0282 (7)	0.0225 (6)	0.0243 (6)	-0.0092 (5)	0.0037 (5)	-0.0100 (5)
O5	0.0241 (6)	0.0153 (6)	0.0333 (7)	-0.0013 (5)	0.0054 (5)	-0.0018 (5)
O2	0.0200 (9)	0.0169 (9)	0.0237 (9)	-0.0017 (7)	-0.0014 (7)	-0.0023 (7)
O2'	0.0200 (9)	0.0169 (9)	0.0237 (9)	-0.0017 (7)	-0.0014 (7)	-0.0023 (7)
O2''	0.0200 (9)	0.0169 (9)	0.0237 (9)	-0.0017 (7)	-0.0014 (7)	-0.0023 (7)
O3	0.0374 (10)	0.0208 (9)	0.0184 (9)	0.0056 (7)	0.0022 (7)	-0.0069 (7)
O3'	0.0374 (10)	0.0208 (9)	0.0184 (9)	0.0056 (7)	0.0022 (7)	-0.0069 (7)
O3''	0.0374 (10)	0.0208 (9)	0.0184 (9)	0.0056 (7)	0.0022 (7)	-0.0069 (7)
O4	0.0304 (10)	0.0163 (8)	0.0289 (10)	0.0014 (7)	-0.0012 (7)	-0.0035 (7)
O4'	0.0304 (10)	0.0163 (8)	0.0289 (10)	0.0014 (7)	-0.0012 (7)	-0.0035 (7)
O6	0.0191 (8)	0.0217 (8)	0.0350 (9)	-0.0028 (6)	-0.0028 (6)	-0.0135 (6)
O6'	0.0191 (8)	0.0217 (8)	0.0350 (9)	-0.0028 (6)	-0.0028 (6)	-0.0135 (6)
O7	0.0273 (9)	0.0185 (8)	0.0326 (9)	-0.0019 (6)	0.0009 (6)	-0.0125 (6)
O7'	0.0273 (9)	0.0185 (8)	0.0326 (9)	-0.0019 (6)	0.0009 (6)	-0.0125 (6)
O8	0.0575 (10)	0.0221 (7)	0.0194 (6)	-0.0016 (6)	-0.0017 (6)	0.0000 (5)
N1	0.0198 (7)	0.0167 (7)	0.0231 (7)	-0.0006 (5)	-0.0021 (5)	-0.0075 (5)
N2	0.0191 (7)	0.0132 (6)	0.0200 (7)	-0.0008 (5)	0.0024 (5)	-0.0042 (5)
C1	0.0150 (7)	0.0201 (8)	0.0246 (8)	-0.0050 (6)	0.0038 (6)	-0.0082 (6)
C2	0.0326 (10)	0.0189 (8)	0.0318 (10)	-0.0069 (7)	0.0093 (8)	-0.0088 (7)
C3	0.0286 (9)	0.0187 (8)	0.0315 (10)	-0.0017 (7)	0.0052 (7)	-0.0014 (7)
C4	0.0185 (8)	0.0250 (9)	0.0250 (9)	-0.0013 (6)	-0.0032 (6)	-0.0020 (7)
C5	0.0199 (8)	0.0210 (8)	0.0275 (9)	-0.0013 (6)	-0.0078 (7)	-0.0068 (7)
C6	0.0145 (7)	0.0181 (8)	0.0244 (8)	-0.0002 (6)	-0.0039 (6)	-0.0059 (6)
C7	0.0211 (8)	0.0192 (8)	0.0262 (9)	0.0016 (6)	-0.0069 (6)	-0.0094 (7)
C8	0.0208 (8)	0.0198 (8)	0.0203 (8)	-0.0005 (6)	0.0008 (6)	-0.0044 (6)
C9	0.0230 (9)	0.0364 (11)	0.0336 (10)	-0.0047 (8)	-0.0029 (8)	0.0098 (8)
C10	0.0383 (11)	0.0307 (10)	0.0367 (11)	-0.0115 (8)	0.0183 (9)	-0.0150 (8)
C11	0.0396 (11)	0.0174 (8)	0.0301 (10)	0.0014 (7)	-0.0068 (8)	-0.0066 (7)
C12	0.0132 (7)	0.0175 (8)	0.0360 (10)	-0.0019 (6)	0.0056 (6)	-0.0101 (7)
C13	0.0212 (9)	0.0185 (8)	0.0515 (12)	-0.0012 (7)	-0.0006 (8)	-0.0161 (8)

C14	0.0225 (9)	0.0308 (10)	0.0549 (13)	0.0016 (7)	-0.0065 (8)	-0.0292 (9)
C15	0.0175 (8)	0.0338 (10)	0.0344 (10)	0.0042 (7)	-0.0049 (7)	-0.0202 (8)
C16	0.0174 (8)	0.0230 (8)	0.0273 (9)	0.0019 (6)	-0.0009 (6)	-0.0114 (7)
C17	0.0136 (7)	0.0185 (8)	0.0261 (8)	-0.0013 (6)	0.0024 (6)	-0.0096 (6)
C18	0.0161 (7)	0.0150 (7)	0.0219 (8)	-0.0011 (6)	0.0022 (6)	-0.0052 (6)
C19	0.0191 (8)	0.0158 (7)	0.0176 (7)	-0.0014 (6)	0.0016 (6)	-0.0056 (6)
C20	0.0204 (8)	0.0213 (8)	0.0236 (8)	-0.0034 (6)	0.0042 (6)	-0.0076 (6)
C21	0.0211 (8)	0.0167 (7)	0.0225 (8)	-0.0037 (6)	0.0018 (6)	-0.0054 (6)
C22	0.0251 (9)	0.0214 (8)	0.0215 (8)	0.0007 (7)	-0.0012 (7)	-0.0036 (6)

*Geometric parameters (Å, °)*

C11—C4	1.7442 (19)	C6—C7	1.431 (2)
C12—C15	1.744 (2)	C7—H7A	0.9500
O1—C1	1.302 (2)	C8—C11	1.524 (2)
O5—C12	1.289 (2)	C8—C10	1.526 (2)
O2—C9	1.318 (3)	C8—C9	1.528 (3)
O2—H2	0.8400	C9—H9A	0.9900
O2'—C9	1.251 (4)	C9—H9B	0.9900
O2'—H2'	0.8400	C9—H9C	0.9900
O2"—C9	1.268 (9)	C9—H9D	0.9900
O2"—H2"	0.8400	C9—H9E	0.9900
O3—C10	1.417 (3)	C9—H9F	0.9900
O3—H3	0.8400	C10—H10A	0.9900
O3'—C10	1.318 (6)	C10—H10B	0.9900
O3'—H3'	0.8400	C10—H10C	0.9900
O3"—C10	1.266 (6)	C10—H10D	0.9900
O3"—H3"	0.8400	C10—H10E	0.9900
O4—C11	1.394 (3)	C10—H10F	0.9900
O4—H4	0.8400	C11—H11A	0.9900
O4'—C11	1.313 (4)	C11—H11B	0.9900
O4'—H4'	0.8400	C11—H11C	0.9900
O6—C20	1.411 (2)	C11—H11D	0.9900
O6—H6	0.8400	C12—C13	1.427 (2)
O6'—C20	1.389 (6)	C12—C17	1.435 (2)
O6'—H6'	0.8400	C13—C14	1.360 (3)
O7—C21	1.405 (2)	C13—H13	0.9500
O7—H7	0.8400	C14—C15	1.402 (3)
O7'—C21	1.428 (6)	C14—H14	0.9500
O7'—H7'	0.8400	C15—C16	1.368 (2)
O8—C22	1.421 (2)	C16—C17	1.409 (2)
O8—H8	0.8400	C16—H16	0.9500
N1—C7	1.288 (2)	C17—C18	1.428 (2)
N1—C8	1.467 (2)	C18—H18	0.9500
N1—H1N	0.885 (10)	C19—C21	1.529 (2)
N2—C18	1.297 (2)	C19—C22	1.532 (2)
N2—C19	1.468 (2)	C19—C20	1.539 (2)
N2—H2N	0.886 (10)	C20—H20A	0.9900

C1—C2	1.418 (2)	C20—H20B	0.9900
C1—C6	1.430 (2)	C20—H20C	0.9900
C2—C3	1.370 (3)	C20—H20D	0.9900
C2—H2A	0.9500	C21—H21A	0.9900
C3—C4	1.400 (3)	C21—H21B	0.9900
C3—H3A	0.9500	C21—H21C	0.9900
C4—C5	1.368 (2)	C21—H21D	0.9900
C5—C6	1.405 (2)	C22—H22A	0.9900
C5—H5	0.9500	C22—H22B	0.9900
C9—O2—H2	109.5	H10E—C10—H10F	105.1
C9—O2'—H2'	109.5	O4'—C11—O4	103.6 (2)
C9—O2"—H2"	109.5	O4'—C11—C8	123.9 (2)
C10—O3—H3	109.5	O4—C11—C8	107.90 (16)
C10—O3'—H3'	109.5	O4—C11—H11A	110.1
C10—O3"—H3"	109.5	C8—C11—H11A	110.1
C11—O4—H4	109.5	O4'—C11—H11B	100.6
C11—O4'—H4'	109.5	O4—C11—H11B	110.1
C20—O6—H6	109.5	C8—C11—H11B	110.1
C20—O6'—H6'	109.5	H11A—C11—H11B	108.4
C21—O7—H7	109.5	O4'—C11—H11C	106.4
C21—O7'—H7'	109.5	O4—C11—H11C	108.0
C22—O8—H8	109.5	C8—C11—H11C	106.4
C7—N1—C8	127.86 (14)	O4'—C11—H11D	106.4
C7—N1—H1N	112.8 (18)	C8—C11—H11D	106.4
C8—N1—H1N	119.1 (18)	H11B—C11—H11D	108.8
C18—N2—C19	127.48 (14)	H11C—C11—H11D	106.4
C18—N2—H2N	114.0 (16)	O5—C12—C13	122.12 (16)
C19—N2—H2N	118.0 (16)	O5—C12—C17	121.79 (15)
O1—C1—C2	121.63 (15)	C13—C12—C17	116.09 (17)
O1—C1—C6	121.54 (15)	C14—C13—C12	121.68 (18)
C2—C1—C6	116.83 (16)	C14—C13—H13	119.2
C3—C2—C1	121.40 (16)	C12—C13—H13	119.2
C3—C2—H2A	119.3	C13—C14—C15	120.98 (17)
C1—C2—H2A	119.3	C13—C14—H14	119.5
C2—C3—C4	120.73 (17)	C15—C14—H14	119.5
C2—C3—H3A	119.6	C16—C15—C14	120.36 (19)
C4—C3—H3A	119.6	C16—C15—C12	120.69 (16)
C5—C4—C3	120.11 (17)	C14—C15—C12	118.94 (14)
C5—C4—C11	120.71 (14)	C15—C16—C17	119.69 (17)
C3—C4—C11	119.18 (14)	C15—C16—H16	120.2
C4—C5—C6	120.31 (16)	C17—C16—H16	120.2
C4—C5—H5	119.8	C16—C17—C18	118.51 (15)
C6—C5—H5	119.8	C16—C17—C12	121.20 (15)
C5—C6—C1	120.61 (15)	C18—C17—C12	120.24 (16)
C5—C6—C7	118.89 (15)	N2—C18—C17	121.93 (15)
C1—C6—C7	120.46 (15)	N2—C18—H18	119.0
N1—C7—C6	122.31 (15)	C17—C18—H18	119.0



N1—C7—H7A	118.8	N2—C19—C21	111.20 (13)
C6—C7—H7A	118.8	N2—C19—C22	106.69 (13)
N1—C8—C11	112.26 (14)	C21—C19—C22	109.93 (14)
N1—C8—C10	105.88 (15)	N2—C19—C20	107.43 (13)
C11—C8—C10	109.77 (15)	C21—C19—C20	111.18 (14)
N1—C8—C9	108.14 (14)	C22—C19—C20	110.29 (14)
C11—C8—C9	110.66 (16)	O6'—C20—O6	105.2 (4)
C10—C8—C9	110.00 (16)	O6'—C20—C19	114.2 (3)
O2'—C9—C8	124.6 (2)	O6—C20—C19	111.78 (14)
O2"—C9—C8	104.9 (4)	O6—C20—H20A	109.3
O2—C9—C8	119.54 (19)	C19—C20—H20A	109.3
O2'—C9—H9A	103.4	O6'—C20—H20B	106.9
O2—C9—H9A	107.4	O6—C20—H20B	109.3
C8—C9—H9A	107.4	C19—C20—H20B	109.3
O2"—C9—H9B	118.1	H20A—C20—H20B	107.9
O2—C9—H9B	107.4	O6'—C20—H20C	108.7
C8—C9—H9B	107.4	O6—C20—H20C	108.0
H9A—C9—H9B	107.0	C19—C20—H20C	108.7
O2'—C9—H9C	106.2	H20A—C20—H20C	109.8
C8—C9—H9C	106.2	O6'—C20—H20D	108.7
H9A—C9—H9C	108.2	C19—C20—H20D	108.7
H9B—C9—H9C	120.1	H20A—C20—H20D	112.6
O2'—C9—H9D	106.2	H20B—C20—H20D	109.0
O2—C9—H9D	106.2	H20C—C20—H20D	107.6
C8—C9—H9D	106.2	O7—C21—O7'	101.9 (3)
H9B—C9—H9D	109.8	O7—C21—C19	108.48 (14)
O2"—C9—H9E	110.8	O7'—C21—C19	115.5 (3)
O2—C9—H9E	110.1	O7—C21—H21A	110.0
C8—C9—H9E	110.8	C19—C21—H21A	110.0
O2"—C9—H9F	110.8	O7—C21—H21B	110.0
C8—C9—H9F	110.8	O7'—C21—H21B	110.7
H9E—C9—H9F	108.8	C19—C21—H21B	110.0
O3'—C10—C8	116.5 (5)	H21A—C21—H21B	108.4
O3—C10—C8	111.74 (16)	O7—C21—H21C	114.3
O3—C10—H10A	109.3	O7'—C21—H21C	108.4
C8—C10—H10A	109.3	C19—C21—H21C	108.4
O3—C10—H10B	109.3	H21A—C21—H21C	105.6
C8—C10—H10B	109.3	O7'—C21—H21D	108.4
H10A—C10—H10B	107.9	C19—C21—H21D	108.4
O3'—C10—H10C	108.2	H21A—C21—H21D	116.7
C8—C10—H10C	108.2	H21B—C21—H21D	103.1
O3'—C10—H10D	108.2	H21C—C21—H21D	107.5
C8—C10—H10D	108.2	O8—C22—C19	109.62 (14)
H10C—C10—H10D	107.3	O8—C22—H22A	109.7
C8—C10—H10E	103.1	C19—C22—H22A	109.7
O3"—C10—H10F	103.1	O8—C22—H22B	109.7
O3'—C10—H10F	104.6	C19—C22—H22B	109.7
O3—C10—H10F	101.4	H22A—C22—H22B	108.2

C8—C10—H10F	103.1		
O1—C1—C2—C3	179.32 (17)	C9—C8—C11—O4'	58.1 (3)
C6—C1—C2—C3	-0.5 (3)	N1—C8—C11—O4	176.23 (16)
C1—C2—C3—C4	0.0 (3)	C10—C8—C11—O4	58.8 (2)
C2—C3—C4—C5	0.8 (3)	C9—C8—C11—O4	-62.8 (2)
C2—C3—C4—C11	-178.35 (15)	O5—C12—C13—C14	-179.70 (17)
C3—C4—C5—C6	-1.1 (3)	C17—C12—C13—C14	0.0 (3)
C11—C4—C5—C6	178.02 (13)	C12—C13—C14—C15	-0.3 (3)
C4—C5—C6—C1	0.6 (3)	C13—C14—C15—C16	-0.2 (3)
C4—C5—C6—C7	-177.26 (16)	C13—C14—C15—C12	178.80 (15)
O1—C1—C6—C5	-179.64 (15)	C14—C15—C16—C17	1.0 (3)
C2—C1—C6—C5	0.2 (2)	C12—C15—C16—C17	-177.95 (13)
O1—C1—C6—C7	-1.8 (2)	C15—C16—C17—C18	176.05 (16)
C2—C1—C6—C7	178.03 (16)	C15—C16—C17—C12	-1.4 (3)
C8—N1—C7—C6	-174.53 (16)	O5—C12—C17—C16	-179.45 (15)
C5—C6—C7—N1	178.05 (16)	C13—C12—C17—C16	0.9 (2)
C1—C6—C7—N1	0.2 (3)	O5—C12—C17—C18	3.2 (2)
C7—N1—C8—C11	7.8 (2)	C13—C12—C17—C18	-176.51 (15)
C7—N1—C8—C10	127.59 (19)	C19—N2—C18—C17	171.78 (15)
C7—N1—C8—C9	-114.5 (2)	C16—C17—C18—N2	-179.76 (15)
N1—C8—C9—O2'	-52.8 (3)	C12—C17—C18—N2	-2.3 (2)
C11—C8—C9—O2'	-176.2 (3)	C18—N2—C19—C21	36.7 (2)
C10—C8—C9—O2'	62.4 (3)	C18—N2—C19—C22	156.57 (16)
N1—C8—C9—O2"	174.6 (5)	C18—N2—C19—C20	-85.16 (19)
C11—C8—C9—O2"	51.2 (5)	N2—C19—C20—O6'	-177.7 (4)
C10—C8—C9—O2"	-70.2 (5)	C21—C19—C20—O6'	60.4 (4)
N1—C8—C9—O2	63.7 (2)	C22—C19—C20—O6'	-61.8 (4)
C11—C8—C9—O2	-59.6 (2)	N2—C19—C20—O6	62.94 (17)
C10—C8—C9—O2	178.90 (19)	C21—C19—C20—O6	-58.93 (18)
N1—C8—C10—O3"	-80.6 (7)	C22—C19—C20—O6	178.86 (14)
C11—C8—C10—O3"	40.8 (7)	N2—C19—C21—O7	175.00 (14)
C9—C8—C10—O3"	162.8 (7)	C22—C19—C21—O7	57.08 (18)
N1—C8—C10—O3'	-92.1 (5)	C20—C19—C21—O7	-65.33 (18)
C11—C8—C10—O3'	29.3 (5)	N2—C19—C21—O7'	-71.5 (4)
C9—C8—C10—O3'	151.3 (4)	C22—C19—C21—O7'	170.6 (4)
N1—C8—C10—O3	46.0 (2)	C20—C19—C21—O7'	48.2 (4)
C11—C8—C10—O3	167.35 (18)	N2—C19—C22—O8	48.05 (18)
C9—C8—C10—O3	-70.7 (2)	C21—C19—C22—O8	168.73 (14)
N1—C8—C11—O4'	-62.9 (3)	C20—C19—C22—O8	-68.33 (18)
C10—C8—C11—O4'	179.7 (3)		

## Hydrogen-bond geometry (Å, °)

<i>D</i> —H... <i>A</i>	<i>D</i> —H	H... <i>A</i>	<i>D</i> ... <i>A</i>	<i>D</i> —H... <i>A</i>
N1—H1n...O1	0.87 (1)	1.82 (2)	2.582 (2)	143 (2)
N2—H2n...O5	0.87 (1)	1.82 (2)	2.573 (2)	141 (2)