

(2-Methoxybenzyl)(2-methoxybenzylidene)azanium (2-methoxyphenyl)-methanaminium tetrachloridozincate(II) monohydrate

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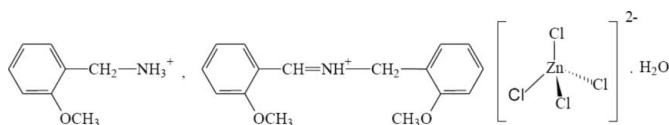
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Key indicators: single-crystal X-ray study; $T = 293$ K; mean $\sigma(\text{C}-\text{C}) = 0.003$ Å; R factor = 0.036; wR factor = 0.087; data-to-parameter ratio = 20.0.

The title compound, $(\text{C}_8\text{H}_{12}\text{NO})(\text{C}_{16}\text{H}_{18}\text{NO}_2)[\text{ZnCl}_4]\cdot\text{H}_2\text{O}$, was obtained as a by-product of the Zn^{2+} and HCl catalyzed condensation of (2-methoxyphenyl)methanamine in water. Both cations feature an intramolecular $\text{N}-\text{H}\cdots\text{O}$ hydrogen bond. In the crystal, the components are linked by an extensive three-dimensional network of $\text{N}-\text{H}\cdots\text{O}$, $\text{O}-\text{H}\cdots\text{Cl}$ and $\text{N}-\text{H}\cdots\text{Cl}$ hydrogen bonds (three of them bifurcated). Weak $\text{C}-\text{H}\cdots\text{O}$ interactions also occur.

Related literature

For related *meta*-chlorido complexes, see: Ben Gharbia *et al.* (2005, 2008). For $\text{Zn}-\text{Cl}$ distances and $\text{Cl}-\text{Zn}-\text{Cl}$ bond angles, see: Gayathri *et al.* (2008); Hosseinian & Mahjoub (2009).



Experimental

Crystal data

$(\text{C}_8\text{H}_{12}\text{NO})(\text{C}_{16}\text{H}_{18}\text{NO}_2)[\text{ZnCl}_4]\cdot\text{H}_2\text{O}$

$M_r = 619.69$
Triclinic, $P\bar{1}$
 $a = 8.0884$ (9) Å
 $b = 12.424$ (2) Å
 $c = 14.678$ (2) Å
 $\alpha = 98.23$ (1)°

$\beta = 97.43$ (1)°
 $\gamma = 90.29$ (1)°
 $V = 1447.1$ (3) Å³
 $Z = 2$
Mo $K\alpha$ radiation
 $\mu = 1.25$ mm⁻¹
 $T = 293$ K
 $0.54 \times 0.47 \times 0.25$ mm

Data collection

Oxford Diffraction Xcalibur diffractometer
Absorption correction: analytical (*CrysAlis PRO*; Oxford Diffraction, 2009; Clark & Reid, 1995)
 $T_{\min} = 0.576$, $T_{\max} = 0.749$
12584 measured reflections
6583 independent reflections
4644 reflections with $I > 2\sigma(I)$
 $R_{\text{int}} = 0.028$

Refinement

$R[F^2 > 2\sigma(F^2)] = 0.036$
 $wR(F^2) = 0.087$
 $S = 0.98$
6583 reflections
329 parameters
3 restraints
H atoms treated by a mixture of independent and constrained refinement
 $\Delta\rho_{\text{max}} = 0.72$ e Å⁻³
 $\Delta\rho_{\text{min}} = -0.64$ e Å⁻³

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

$D-\text{H}\cdots A$	$D-\text{H}$	$\text{H}\cdots A$	$D\cdots A$	$D-\text{H}\cdots A$
$\text{N1}-\text{H1}\cdots\text{O2}$	0.89 (1)	2.07 (2)	2.680 (2)	125 (2)
$\text{N1}-\text{H1}\cdots\text{Cl2}^i$	0.89 (1)	2.64 (2)	3.3221 (18)	135 (2)
$\text{N2}-\text{H2A}\cdots\text{O4}$	0.89	2.30	3.102 (3)	151
$\text{N2}-\text{H2A}\cdots\text{O3}$	0.89	2.37	2.877 (2)	116
$\text{N2}-\text{H2B}\cdots\text{O4}^{ii}$	0.89	2.04	2.881 (3)	158
$\text{N2}-\text{H2B}\cdots\text{Cl3}^{iii}$	0.89	2.98	3.502 (2)	120
$\text{N2}-\text{H2C}\cdots\text{Cl1}^{iv}$	0.89	2.45	3.306 (2)	162
$\text{O4}-\text{H4B}\cdots\text{Cl4}^{iii}$	0.80 (2)	2.44 (2)	3.2323 (19)	168 (3)
$\text{O4}-\text{H4A}\cdots\text{Cl1}^v$	0.80 (2)	2.72 (2)	3.4165 (19)	147 (3)
$\text{C8}-\text{H8A}\cdots\text{Cl3}$	0.97	2.67	3.474 (2)	140
$\text{C11}-\text{H11}\cdots\text{Cl2}$	0.93	2.82	3.687 (2)	155
$\text{C24}-\text{H24A}\cdots\text{Cl4}^{vi}$	0.97	2.76	3.707 (3)	166

Symmetry codes: (i) $x+1, y, z$; (ii) $-x+1, -y+1, -z$; (iii) $x, y, z-1$; (iv) $-x+1, -y+1, -z+1$; (v) $-x, -y+1, -z+1$; (vi) $x+1, y, z-1$.

Data collection: *CrysAlis PRO* (Oxford Diffraction, 2009); cell refinement: *CrysAlis PRO*; data reduction: *CrysAlis PRO*; program(s) used to solve structure: *SIR97* (Altomare *et al.*, 1999); program(s) used to refine structure: *SHELXTL* (Sheldrick, 2008); molecular graphics: *SHELXTL*; software used to prepare material for publication: *SHELXTL*.

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Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: HB5489).

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supporting information

Acta Cryst. (2010). E66, m791 [doi:10.1107/S1600536810021793]

(2-Methoxybenzyl)(2-methoxybenzylidene)azanium (2-methoxyphenyl)-methanaminium tetrachloridozincate(II) monohydrate

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S1. Comment

As a part of our ongoing investigations in molecular salts containing *meta*-chlorido complexes (Ben Gharbia *et al.*, 2005; Ben Gharbia *et al.*, 2008), we report here the crystal structure of one such compound, $(C_{16}H_{18}NO_2)(C_8H_{12}NO)[ZnCl_4] \cdot H_2O$. The title compound was obtained as a byproduct of the Zn^{2+} and HCl catalyzed condensation of (2-methoxyphenyl) methanamine in water. Subsequent crystallization from the reaction mixture yielded among the main reaction products crystals of the title compound that consist of one *N*-(2-methoxybenzylidene)-1-(2-methoxyphenyl)-methanaminium cation, one (2-methoxyphenyl) methanaminium cation, one $ZnCl_4^{2-}$ anion and one interstitial water molecule (Fig. 1).

The distance N1—C9 [1.273 (2) Å] is substantially shorter than the one of N1—C8 [1.477 (2) Å], indicating the presence of a double bond in the condensation product, thus indicating the nature of the organic molecules in the crystal as indicated in Scheme 1. Preliminary NMR data on the material indicate that the bulk of the reaction product is not identical with the title compound. Further investigations into the nature of the bulk material are under way.

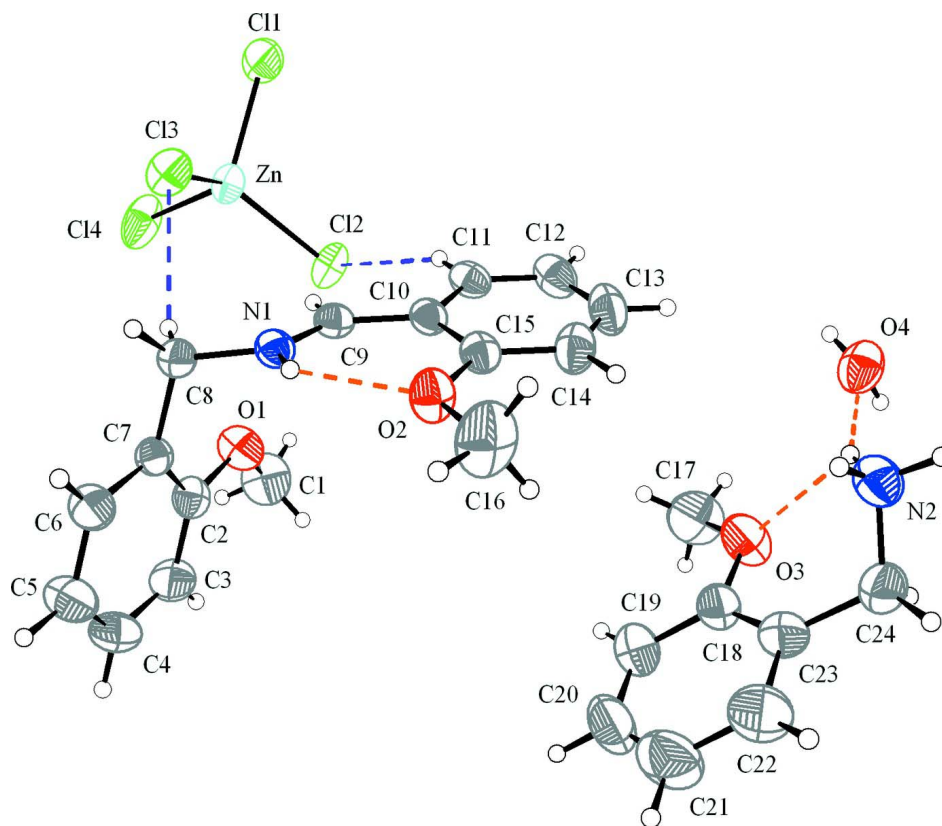
The Cl—Zn—Cl bond angles in the title compound show relatively little distortion from a regular tetrahedron [spread values 104.45 (3)–111.78 (2)] (Gayathri *et al.*, 2008, Hosseinian *et al.*, 2009). Classic N—H \cdots O, O—H \cdots Cl and N—H \cdots Cl hydrogen bonds are observed, which link the two types of organic ammonium cations, the anionic complexes $[ZnCl_4]^{2-}$ and the uncoordinated water molecules into a 3-D hydrogen bonded network, as shown in Fig. 2. Three of the hydrogen bonds are bifurcated: N1—H1 \cdots (Cl2,O2), N2—H2A \cdots (O3,O4) and N2—H2B \cdots (O4,Cl3).

S2. Experimental

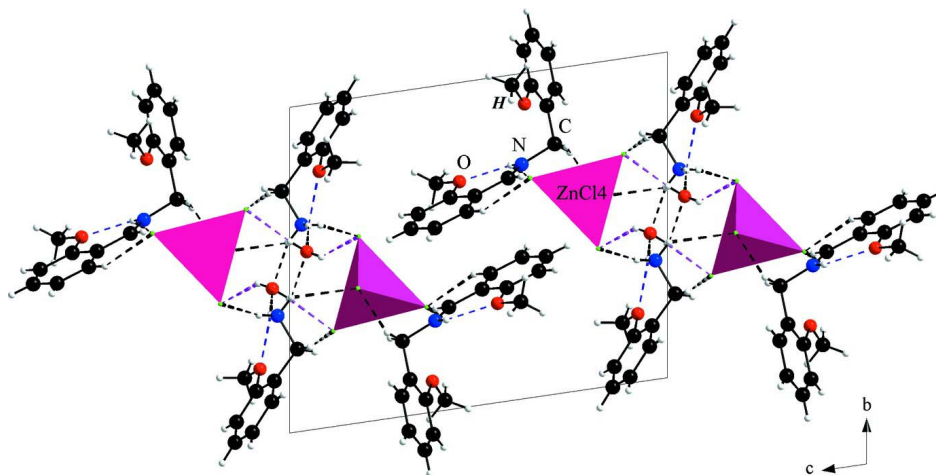
An aqueous solution of (2-methoxyphenyl) methanamine (2-methoxybenzylamine), zinc chloride and 1 M hydrochloric acid in a Petri disk was slowly evaporated at room temperature. A colourless block of (I), which remained stable under normal conditions of temperature and humidity, was isolated as a byproduct of the reaction.

S3. Refinement

C—H and ammonium H atoms were placed in calculated positions with C—H in the range 0.93–0.97 and N—H equal to 0.89 Å. The imminium and the water hydrogen atom positions were refined with N—H and O—H distance restraints of 0.89 (2) and 0.82 (2) Å. The $U_{iso}(H)$ values of all H atoms were constrained to 1.2 or 1.5 times U_{eq} of the respective parent atom.

**Figure 1**

A view of the title compound, showing 50% probability displacement ellipsoids (arbitrary spheres for the H atoms).

**Figure 2**

Projection of the structure along the *a* axis. Hydrogen bonds are denoted by dotted lines.

(2-Methoxybenzyl)(2-methoxybenzylidene)azanium (2-methoxyphenyl)methanaminium tetrachloridozincate(II) monohydrate*Crystal data* $(C_8H_{12}NO)(C_{16}H_{18}NO_2)[ZnCl_4] \cdot H_2O$ $M_r = 619.69$ Triclinic, $P\bar{1}$

Hall symbol: -P 1

 $a = 8.0884$ (9) Å $b = 12.424$ (2) Å $c = 14.678$ (2) Å $\alpha = 98.23$ (1)° $\beta = 97.43$ (1)° $\gamma = 90.29$ (1)° $V = 1447.1$ (3) Å³ $Z = 2$ $F(000) = 640$ $D_x = 1.422$ Mg m⁻³Mo $K\alpha$ radiation, $\lambda = 0.71069$ Å

Cell parameters from 5313 reflections

 $\theta = 3.0$ – 29.2 ° $\mu = 1.25$ mm⁻¹ $T = 293$ K

Block, colourless

 $0.54 \times 0.47 \times 0.25$ mm*Data collection*Oxford Diffraction Xcalibur
diffractometer

Radiation source: fine-focus sealed tube

Graphite monochromator

 ω scans

Absorption correction: analytical

(CrysAlis PRO; Oxford Diffraction, 2009; Clark
& Reid, 1995) $T_{\min} = 0.576$, $T_{\max} = 0.749$

12584 measured reflections

6583 independent reflections

4644 reflections with $I > 2\sigma(I)$ $R_{\text{int}} = 0.028$ $\theta_{\max} = 29.5$ °, $\theta_{\min} = 3.0$ ° $h = -10 \rightarrow 11$ $k = -15 \rightarrow 16$ $l = -19 \rightarrow 19$ *Refinement*Refinement on F^2

Least-squares matrix: full

 $R[F^2 > 2\sigma(F^2)] = 0.036$ $wR(F^2) = 0.087$ $S = 0.98$

6583 reflections

329 parameters

3 restraints

Primary atom site location: structure-invariant
direct methodsSecondary atom site location: difference Fourier
mapHydrogen site location: inferred from
neighbouring sitesH atoms treated by a mixture of independent
and constrained refinement $w = 1/[\sigma^2(F_o^2) + (0.0435P)^2]$ where $P = (F_o^2 + 2F_c^2)/3$ $(\Delta/\sigma)_{\max} = 0.001$ $\Delta\rho_{\max} = 0.72$ e Å⁻³ $\Delta\rho_{\min} = -0.64$ e Å⁻³*Special details*

Geometry. All e.s.d.'s (except the e.s.d. in the dihedral angle between two l.s. planes) are estimated using the full covariance matrix. The cell e.s.d.'s are taken into account individually in the estimation of e.s.d.'s in distances, angles and torsion angles; correlations between e.s.d.'s in cell parameters are only used when they are defined by crystal symmetry. An approximate (isotropic) treatment of cell e.s.d.'s is used for estimating e.s.d.'s involving l.s. planes.

Refinement. Refinement of F^2 against ALL reflections. The weighted R -factor wR and goodness of fit S are based on F^2 , conventional R -factors R are based on F , with F set to zero for negative F^2 . The threshold expression of $F^2 > \sigma(F^2)$ is used only for calculating R -factors(gt) etc. and is not relevant to the choice of reflections for refinement. R -factors based on F^2 are statistically about twice as large as those based on F , and R -factors based on ALL data will be even larger.

Fractional atomic coordinates and isotropic or equivalent isotropic displacement parameters (Å²)

	x	y	z	$U_{\text{iso}}^*/U_{\text{eq}}$
C1	0.2523 (3)	0.0313 (2)	0.58855 (19)	0.0669 (8)

H1A	0.2600	-0.0055	0.5272	0.100*
H1B	0.1579	0.0778	0.5869	0.100*
H1C	0.2392	-0.0214	0.6292	0.100*
C2	0.5461 (3)	0.04191 (18)	0.63917 (14)	0.0377 (5)
C3	0.5657 (3)	-0.06829 (19)	0.61983 (16)	0.0486 (6)
H3	0.4754	-0.1130	0.5916	0.058*
C4	0.7185 (3)	-0.1128 (2)	0.64206 (17)	0.0559 (6)
H4	0.7313	-0.1875	0.6285	0.067*
C5	0.8521 (3)	-0.0475 (2)	0.68406 (17)	0.0535 (6)
H5	0.9549	-0.0779	0.6999	0.064*
C6	0.8333 (3)	0.06297 (19)	0.70261 (14)	0.0421 (5)
H6	0.9247	0.1070	0.7303	0.051*
C7	0.6816 (2)	0.11022 (17)	0.68108 (13)	0.0331 (4)
C8	0.6598 (3)	0.23148 (16)	0.70157 (13)	0.0367 (5)
H8A	0.5629	0.2461	0.7336	0.044*
H8B	0.7568	0.2646	0.7418	0.044*
C9	0.5018 (2)	0.31536 (16)	0.57852 (13)	0.0331 (4)
H9	0.4129	0.3162	0.6129	0.040*
C10	0.4693 (2)	0.35513 (16)	0.49025 (13)	0.0326 (4)
C11	0.3132 (2)	0.39947 (17)	0.46852 (15)	0.0390 (5)
H11	0.2380	0.4048	0.5117	0.047*
C12	0.2683 (3)	0.43548 (18)	0.38452 (16)	0.0485 (6)
H12	0.1641	0.4649	0.3709	0.058*
C13	0.3800 (3)	0.42721 (19)	0.32143 (16)	0.0536 (6)
H13	0.3498	0.4506	0.2643	0.064*
C14	0.5355 (3)	0.3854 (2)	0.34018 (15)	0.0469 (6)
H14	0.6098	0.3814	0.2965	0.056*
C15	0.5808 (2)	0.34898 (17)	0.42502 (14)	0.0360 (5)
C16	0.8537 (3)	0.2994 (3)	0.38782 (19)	0.0764 (9)
H16A	0.8119	0.2554	0.3301	0.115*
H16B	0.9527	0.2678	0.4153	0.115*
H16C	0.8797	0.3714	0.3766	0.115*
C17	0.2513 (3)	0.1890 (3)	0.1188 (2)	0.0783 (9)
H17A	0.2140	0.1149	0.0986	0.117*
H17B	0.1647	0.2370	0.1003	0.117*
H17C	0.2776	0.2010	0.1852	0.117*
C18	0.5273 (3)	0.14345 (19)	0.08828 (15)	0.0467 (6)
C19	0.5371 (4)	0.0578 (2)	0.14046 (17)	0.0616 (7)
H19	0.4496	0.0425	0.1722	0.074*
C20	0.6792 (4)	-0.0042 (2)	0.14433 (19)	0.0741 (8)
H20	0.6864	-0.0621	0.1783	0.089*
C21	0.8083 (4)	0.0192 (3)	0.0987 (2)	0.0772 (9)
H21	0.9038	-0.0221	0.1023	0.093*
C22	0.7973 (3)	0.1043 (2)	0.04706 (19)	0.0641 (7)
H22	0.8862	0.1200	0.0164	0.077*
C23	0.6558 (3)	0.16659 (19)	0.04026 (15)	0.0435 (5)
C24	0.6378 (3)	0.2543 (2)	-0.02067 (16)	0.0512 (6)
H24A	0.7340	0.2540	-0.0539	0.061*

H24B	0.5397	0.2379	-0.0663	0.061*
Cl3	0.40847 (6)	0.41374 (5)	0.81903 (4)	0.04914 (15)
Cl2	0.04753 (7)	0.32552 (5)	0.63718 (3)	0.05224 (16)
Cl1	0.03902 (8)	0.57253 (6)	0.81699 (4)	0.06066 (18)
Cl4	0.03288 (9)	0.29378 (6)	0.88415 (4)	0.0746 (2)
N1	0.6383 (2)	0.27933 (13)	0.61445 (11)	0.0334 (4)
H1	0.729 (2)	0.2772 (17)	0.5852 (13)	0.040*
N2	0.6225 (2)	0.36549 (15)	0.03186 (13)	0.0488 (5)
H2A	0.5218	0.3717	0.0501	0.073*
H2B	0.6365	0.4152	-0.0047	0.073*
H2C	0.7000	0.3756	0.0814	0.073*
O1	0.40047 (18)	0.09490 (13)	0.62188 (11)	0.0508 (4)
O2	0.72997 (17)	0.30455 (13)	0.44960 (10)	0.0458 (4)
O3	0.3943 (2)	0.20972 (14)	0.07819 (12)	0.0616 (5)
O4	0.2655 (2)	0.45092 (16)	0.04551 (11)	0.0573 (5)
H4B	0.220 (3)	0.412 (2)	0.0009 (15)	0.086*
H4A	0.208 (3)	0.473 (3)	0.0836 (17)	0.086*
Zn1	0.12911 (3)	0.39816 (2)	0.786811 (15)	0.03902 (9)

Atomic displacement parameters (Å²)

	U^{11}	U^{22}	U^{33}	U^{12}	U^{13}	U^{23}
C1	0.0402 (13)	0.079 (2)	0.0807 (19)	-0.0143 (13)	0.0011 (13)	0.0147 (16)
C2	0.0393 (12)	0.0406 (14)	0.0358 (11)	-0.0040 (10)	0.0081 (9)	0.0113 (10)
C3	0.0548 (14)	0.0362 (14)	0.0524 (14)	-0.0100 (11)	0.0027 (11)	0.0031 (12)
C4	0.0713 (18)	0.0304 (14)	0.0637 (16)	0.0043 (12)	0.0050 (13)	0.0018 (12)
C5	0.0504 (14)	0.0428 (16)	0.0665 (16)	0.0090 (11)	0.0044 (12)	0.0085 (13)
C6	0.0401 (12)	0.0393 (14)	0.0461 (13)	-0.0011 (10)	0.0011 (10)	0.0075 (11)
C7	0.0400 (11)	0.0314 (12)	0.0299 (10)	-0.0007 (9)	0.0076 (8)	0.0080 (9)
C8	0.0462 (12)	0.0327 (12)	0.0311 (10)	0.0012 (9)	0.0046 (9)	0.0050 (9)
C9	0.0349 (11)	0.0276 (11)	0.0367 (11)	-0.0013 (8)	0.0073 (9)	0.0024 (9)
C10	0.0343 (10)	0.0275 (11)	0.0357 (11)	-0.0030 (8)	0.0015 (8)	0.0061 (9)
C11	0.0349 (11)	0.0307 (12)	0.0504 (13)	-0.0035 (9)	0.0017 (10)	0.0061 (10)
C12	0.0416 (12)	0.0413 (14)	0.0599 (15)	-0.0027 (10)	-0.0104 (11)	0.0142 (12)
C13	0.0635 (16)	0.0510 (16)	0.0454 (13)	-0.0112 (12)	-0.0118 (12)	0.0208 (12)
C14	0.0510 (14)	0.0531 (15)	0.0377 (12)	-0.0075 (11)	0.0052 (10)	0.0118 (11)
C15	0.0355 (11)	0.0327 (12)	0.0393 (11)	-0.0066 (9)	0.0008 (9)	0.0073 (10)
C16	0.0522 (16)	0.116 (3)	0.0707 (18)	0.0130 (16)	0.0307 (14)	0.0248 (18)
C17	0.0612 (17)	0.080 (2)	0.104 (2)	0.0048 (15)	0.0403 (17)	0.0217 (19)
C18	0.0518 (14)	0.0454 (15)	0.0433 (13)	0.0018 (11)	0.0046 (11)	0.0090 (12)
C19	0.0748 (18)	0.0559 (18)	0.0588 (16)	0.0021 (14)	0.0137 (14)	0.0202 (14)
C20	0.101 (2)	0.058 (2)	0.0637 (18)	0.0116 (17)	-0.0055 (17)	0.0245 (16)
C21	0.069 (2)	0.075 (2)	0.084 (2)	0.0259 (16)	-0.0060 (16)	0.0126 (18)
C22	0.0508 (15)	0.066 (2)	0.0751 (18)	0.0077 (13)	0.0082 (13)	0.0080 (16)
C23	0.0402 (12)	0.0427 (14)	0.0457 (13)	0.0002 (10)	0.0025 (10)	0.0019 (11)
C24	0.0579 (15)	0.0516 (16)	0.0472 (13)	-0.0012 (12)	0.0177 (11)	0.0080 (12)
Cl3	0.0335 (3)	0.0638 (4)	0.0473 (3)	0.0050 (3)	0.0046 (2)	-0.0010 (3)
Cl2	0.0449 (3)	0.0768 (5)	0.0322 (3)	-0.0055 (3)	0.0033 (2)	0.0003 (3)

C11	0.0586 (4)	0.0643 (5)	0.0540 (4)	0.0230 (3)	-0.0034 (3)	0.0003 (3)
C14	0.0860 (5)	0.0963 (6)	0.0413 (3)	-0.0483 (4)	0.0023 (3)	0.0165 (4)
N1	0.0369 (9)	0.0298 (10)	0.0347 (9)	-0.0008 (8)	0.0067 (7)	0.0067 (8)
N2	0.0512 (11)	0.0436 (12)	0.0522 (11)	-0.0044 (9)	0.0039 (9)	0.0118 (10)
O1	0.0377 (8)	0.0492 (10)	0.0652 (10)	-0.0031 (7)	0.0023 (7)	0.0120 (8)
O2	0.0374 (8)	0.0595 (11)	0.0444 (8)	0.0057 (7)	0.0109 (7)	0.0159 (8)
O3	0.0493 (10)	0.0612 (12)	0.0850 (13)	0.0119 (8)	0.0259 (9)	0.0309 (10)
O4	0.0568 (11)	0.0688 (13)	0.0442 (10)	-0.0155 (9)	0.0040 (8)	0.0048 (10)
Zn1	0.03518 (14)	0.05046 (18)	0.03133 (14)	-0.00161 (11)	0.00427 (10)	0.00579 (12)

Geometric parameters (Å, °)

C1—O1	1.424 (3)	C16—O2	1.431 (2)
C1—H1A	0.9600	C16—H16A	0.9600
C1—H1B	0.9600	C16—H16B	0.9600
C1—H1C	0.9600	C16—H16C	0.9600
C2—O1	1.365 (2)	C17—O3	1.407 (3)
C2—C3	1.372 (3)	C17—H17A	0.9600
C2—C7	1.402 (3)	C17—H17B	0.9600
C3—C4	1.377 (3)	C17—H17C	0.9600
C3—H3	0.9300	C18—O3	1.364 (3)
C4—C5	1.372 (3)	C18—C23	1.380 (3)
C4—H4	0.9300	C18—C19	1.394 (3)
C5—C6	1.374 (3)	C19—C20	1.386 (4)
C5—H5	0.9300	C19—H19	0.9300
C6—C7	1.382 (3)	C20—C21	1.362 (4)
C6—H6	0.9300	C20—H20	0.9300
C7—C8	1.509 (3)	C21—C22	1.384 (4)
C8—N1	1.477 (2)	C21—H21	0.9300
C8—H8A	0.9700	C22—C23	1.386 (3)
C8—H8B	0.9700	C22—H22	0.9300
C9—N1	1.273 (2)	C23—C24	1.501 (3)
C9—C10	1.445 (3)	C24—N2	1.497 (3)
C9—H9	0.9300	C24—H24A	0.9700
C10—C15	1.392 (3)	C24—H24B	0.9700
C10—C11	1.397 (3)	Cl3—Zn1	2.2495 (6)
C11—C12	1.377 (3)	Cl2—Zn1	2.2639 (6)
C11—H11	0.9300	Cl1—Zn1	2.2903 (8)
C12—C13	1.369 (3)	Cl4—Zn1	2.2664 (7)
C12—H12	0.9300	N1—H1	0.893 (14)
C13—C14	1.373 (3)	N2—H2A	0.8900
C13—H13	0.9300	N2—H2B	0.8900
C14—C15	1.391 (3)	N2—H2C	0.8900
C14—H14	0.9300	O4—H4B	0.801 (17)
C15—O2	1.357 (2)	O4—H4A	0.795 (17)
O1—C1—H1A	109.5	H16A—C16—H16B	109.5
O1—C1—H1B	109.5	O2—C16—H16C	109.5

H1A—C1—H1B	109.5	H16A—C16—H16C	109.5
O1—C1—H1C	109.5	H16B—C16—H16C	109.5
H1A—C1—H1C	109.5	O3—C17—H17A	109.5
H1B—C1—H1C	109.5	O3—C17—H17B	109.5
O1—C2—C3	125.4 (2)	H17A—C17—H17B	109.5
O1—C2—C7	114.2 (2)	O3—C17—H17C	109.5
C3—C2—C7	120.4 (2)	H17A—C17—H17C	109.5
C2—C3—C4	120.2 (2)	H17B—C17—H17C	109.5
C2—C3—H3	119.9	O3—C18—C23	114.11 (19)
C4—C3—H3	119.9	O3—C18—C19	124.9 (2)
C5—C4—C3	120.3 (2)	C23—C18—C19	121.0 (2)
C5—C4—H4	119.9	C20—C19—C18	119.1 (2)
C3—C4—H4	119.9	C20—C19—H19	120.4
C4—C5—C6	119.6 (2)	C18—C19—H19	120.4
C4—C5—H5	120.2	C21—C20—C19	120.4 (3)
C6—C5—H5	120.2	C21—C20—H20	119.8
C5—C6—C7	121.5 (2)	C19—C20—H20	119.8
C5—C6—H6	119.3	C20—C21—C22	120.1 (3)
C7—C6—H6	119.3	C20—C21—H21	120.0
C6—C7—C2	118.0 (2)	C22—C21—H21	120.0
C6—C7—C8	121.79 (19)	C21—C22—C23	121.0 (3)
C2—C7—C8	120.23 (18)	C21—C22—H22	119.5
N1—C8—C7	110.25 (16)	C23—C22—H22	119.5
N1—C8—H8A	109.6	C18—C23—C22	118.4 (2)
C7—C8—H8A	109.6	C18—C23—C24	120.1 (2)
N1—C8—H8B	109.6	C22—C23—C24	121.5 (2)
C7—C8—H8B	109.6	N2—C24—C23	113.39 (18)
H8A—C8—H8B	108.1	N2—C24—H24A	108.9
N1—C9—C10	127.33 (18)	C23—C24—H24A	108.9
N1—C9—H9	116.3	N2—C24—H24B	108.9
C10—C9—H9	116.3	C23—C24—H24B	108.9
C15—C10—C11	118.40 (18)	H24A—C24—H24B	107.7
C15—C10—C9	124.36 (18)	C9—N1—C8	124.65 (17)
C11—C10—C9	117.20 (18)	C9—N1—H1	120.6 (13)
C12—C11—C10	121.4 (2)	C8—N1—H1	114.6 (13)
C12—C11—H11	119.3	C24—N2—H2A	109.5
C10—C11—H11	119.3	C24—N2—H2B	109.5
C13—C12—C11	118.8 (2)	H2A—N2—H2B	109.5
C13—C12—H12	120.6	C24—N2—H2C	109.5
C11—C12—H12	120.6	H2A—N2—H2C	109.5
C12—C13—C14	121.8 (2)	H2B—N2—H2C	109.5
C12—C13—H13	119.1	C2—O1—C1	118.18 (19)
C14—C13—H13	119.1	C15—O2—C16	119.14 (17)
C13—C14—C15	119.4 (2)	C18—O3—C17	119.08 (19)
C13—C14—H14	120.3	H4B—O4—H4A	115 (3)
C15—C14—H14	120.3	Cl3—Zn1—Cl2	111.78 (2)
O2—C15—C14	123.99 (19)	Cl3—Zn1—Cl4	108.77 (3)
O2—C15—C10	115.83 (17)	Cl2—Zn1—Cl4	110.35 (3)

C14—C15—C10	120.17 (19)	C13—Zn1—C11	104.45 (3)
O2—C16—H16A	109.5	C12—Zn1—C11	111.19 (3)
O2—C16—H16B	109.5	C14—Zn1—C11	110.14 (3)
O1—C2—C3—C4	-178.8 (2)	C11—C10—C15—C14	-0.7 (3)
C7—C2—C3—C4	0.5 (3)	C9—C10—C15—C14	177.2 (2)
C2—C3—C4—C5	0.3 (4)	O3—C18—C19—C20	179.4 (2)
C3—C4—C5—C6	-1.0 (4)	C23—C18—C19—C20	0.5 (4)
C4—C5—C6—C7	0.9 (3)	C18—C19—C20—C21	0.8 (4)
C5—C6—C7—C2	-0.2 (3)	C19—C20—C21—C22	-0.9 (5)
C5—C6—C7—C8	179.77 (18)	C20—C21—C22—C23	-0.4 (4)
O1—C2—C7—C6	178.82 (17)	O3—C18—C23—C22	179.3 (2)
C3—C2—C7—C6	-0.5 (3)	C19—C18—C23—C22	-1.7 (4)
O1—C2—C7—C8	-1.1 (3)	O3—C18—C23—C24	-3.2 (3)
C3—C2—C7—C8	179.54 (18)	C19—C18—C23—C24	175.8 (2)
C6—C7—C8—N1	108.7 (2)	C21—C22—C23—C18	1.6 (4)
C2—C7—C8—N1	-71.3 (2)	C21—C22—C23—C24	-175.8 (2)
N1—C9—C10—C15	7.4 (3)	C18—C23—C24—N2	64.7 (3)
N1—C9—C10—C11	-174.7 (2)	C22—C23—C24—N2	-117.9 (2)
C15—C10—C11—C12	0.7 (3)	C10—C9—N1—C8	-174.76 (19)
C9—C10—C11—C12	-177.4 (2)	C7—C8—N1—C9	108.3 (2)
C10—C11—C12—C13	0.0 (3)	C3—C2—O1—C1	5.8 (3)
C11—C12—C13—C14	-0.8 (4)	C7—C2—O1—C1	-173.46 (18)
C12—C13—C14—C15	0.8 (4)	C14—C15—O2—C16	3.6 (3)
C13—C14—C15—O2	178.8 (2)	C10—C15—O2—C16	-177.5 (2)
C13—C14—C15—C10	-0.1 (3)	C23—C18—O3—C17	175.7 (2)
C11—C10—C15—O2	-179.57 (19)	C19—C18—O3—C17	-3.3 (4)
C9—C10—C15—O2	-1.7 (3)		

Hydrogen-bond geometry (\AA , $^\circ$)

<i>D</i> —H \cdots <i>A</i>	<i>D</i> —H	H \cdots <i>A</i>	<i>D</i> \cdots <i>A</i>	<i>D</i> —H \cdots <i>A</i>
N1—H1 \cdots O2	0.89 (1)	2.07 (2)	2.680 (2)	125 (2)
N1—H1 \cdots Cl2 ⁱ	0.89 (1)	2.64 (2)	3.3221 (18)	135 (2)
N2—H2A \cdots O4	0.89	2.30	3.102 (3)	151
N2—H2A \cdots O3	0.89	2.37	2.877 (2)	116
N2—H2B \cdots O4 ⁱⁱ	0.89	2.04	2.881 (3)	158
N2—H2B \cdots Cl3 ⁱⁱⁱ	0.89	2.98	3.502 (2)	120
N2—H2C \cdots Cl1 ^{iv}	0.89	2.45	3.306 (2)	162
O4—H4B \cdots Cl4 ⁱⁱⁱ	0.80 (2)	2.44 (2)	3.2323 (19)	168 (3)
O4—H4A \cdots Cl1 ^v	0.80 (2)	2.72 (2)	3.4165 (19)	147 (3)
C8—H8A \cdots Cl3	0.97	2.67	3.474 (2)	140
C11—H11 \cdots Cl2	0.93	2.82	3.687 (2)	155
C24—H24A \cdots Cl4 ^{vi}	0.97	2.76	3.707 (3)	166

Symmetry codes: (i) $x+1, y, z$; (ii) $-x+1, -y+1, -z$; (iii) $x, y, z-1$; (iv) $-x+1, -y+1, -z+1$; (v) $-x, -y+1, -z+1$; (vi) $x+1, y, z-1$.