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catena-Poly[zinc(II)-bis[μ-2-(2,4-dichlorophenoxy)acetato]]

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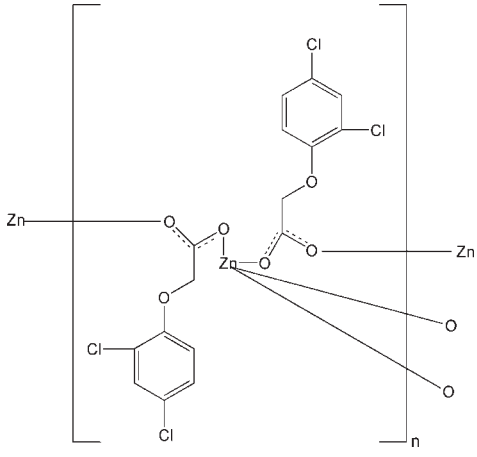
Key indicators: single-crystal X-ray study; T = 296 K; mean σ(C–C) = 0.004 Å; R factor = 0.030; wR factor = 0.091; data-to-parameter ratio = 13.3.

The title polymeric compound, $[Zn(C_8H_5Cl_2O_3)_2]_n$, was prepared by reaction of zinc(II) chloride with 2,4-dichlorophenoxyacetic acid and sodium hydroxide under hydrothermal conditions. The Zn^{II} atom is coordinated in a distorted tetrahedral environment by four O atoms from four 2,4-dichlorophenoxyacetate ligands. Each ligand bridges two Zn^{II} atoms, forming a polymeric chain along the *a* axis. Adjacent chains are connected *via* C—H···Cl hydrogen bonds.

Related literature

For metal-organic coordination polymers, see: Qin *et al.* (2009); Huang *et al.* (2008); Reineke *et al.* (1999); Xiong *et al.* (2002).



Experimental

Crystal data

$[Zn(C_8H_5Cl_2O_3)_2]$
M_r = 505.43
Triclinic, $P\overline{1}$
a = 4.7322 (10) Å
b = 10.459 (2) Å
c = 18.979 (4) Å
α = 79.340 (2)°
β = 89.838 (2)°
γ = 82.847 (3)°
V = 915.8 (3) Å^3
Z = 2
Mo $K\alpha$ radiation
μ = 1.96 mm^{-1}
T = 296 K
0.56 × 0.21 × 0.16 mm

Data collection

Bruker SMART APEXII CCD diffractometer
Absorption correction: multi-scan (*SADABS*; Sheldrick, 1996)
T_{min} = 0.617, T_{max} = 0.731
4737 measured reflections
3255 independent reflections
2849 reflections with $I > 2\sigma(I)$
R_{int} = 0.021

Refinement

$R[F^2 > 2\sigma(F^2)]$ = 0.030
$wR(F^2)$ = 0.091
S = 0.94
3255 reflections
244 parameters
H-atom parameters constrained
$\Delta\rho_{max}$ = 0.34 e Å^{-3}
$\Delta\rho_{min}$ = −0.39 e Å^{-3}

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

D—H···A	D—H	H···A	D···A	D—H···A
C2—H2B···Cl2^{i}	0.97	2.73	3.610 (3)	151

Symmetry code: (i) $-x+1, -y+2, -z+1$.

Data collection: *APEX2* (Bruker, 2004); cell refinement: *SAINT* (Bruker, 2004); data reduction: *SAINT*; program(s) used to solve structure: *SHELXS97* (Sheldrick, 2008); program(s) used to refine structure: *SHELXL97* (Sheldrick, 2008); molecular graphics: *SHELXTL* (Sheldrick, 2008); software used to prepare material for publication: *SHELXTL*.

The author acknowledges South China Normal University for supporting this work.

Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: CI5084).

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supporting information

Acta Cryst. (2010). E**66**, m621 [https://doi.org/10.1107/S1600536810015977]

catena-Poly[zinc(II)-bis[μ-2-(2,4-dichlorophenoxy)acetato]]

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S1. Experimental

A mixture of $ZnCl_2$ (0.068 g, 0.5 mmol), H_2O (6 ml), 2,4-dichlorophenoxyacetic acid (0.22 g, 1 mmol) was stirred vigorously for 10 min and then sodium hydroxide solution (1.0 mol/L) was added to adjust the pH value to 7.0. The mixture was then sealed in a 25 ml Teflon-lined stainless-steel autoclave. The autoclave was heated at 413 K for 3 d and then slowly cooled to room temperature at 6 K/h. The product was collected by filtration, washed with water and air-dried. Colourless needle-shaped crystals were obtained in ca. 42.3% yield based on Zn.

S2. Refinement

H atoms were positioned geometrically [C–H = 0.93–0.97 Å] and refined using a riding model, with U_{iso}(H) = 1.2U_{eq}(C).

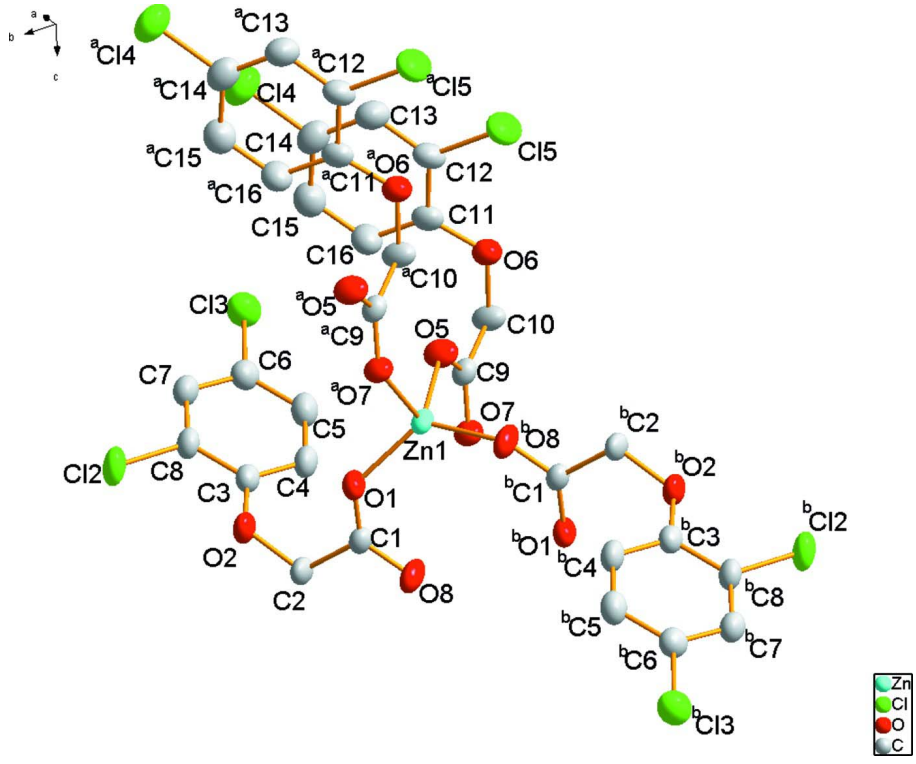

Figure 1

Part of a polymeric chain in the title compound. Displacement ellipsoids are drawn at the 50% probability level. Symmetry codes: (a) 1+x, y, z; (b) 1-x, 1-y, 1-z.

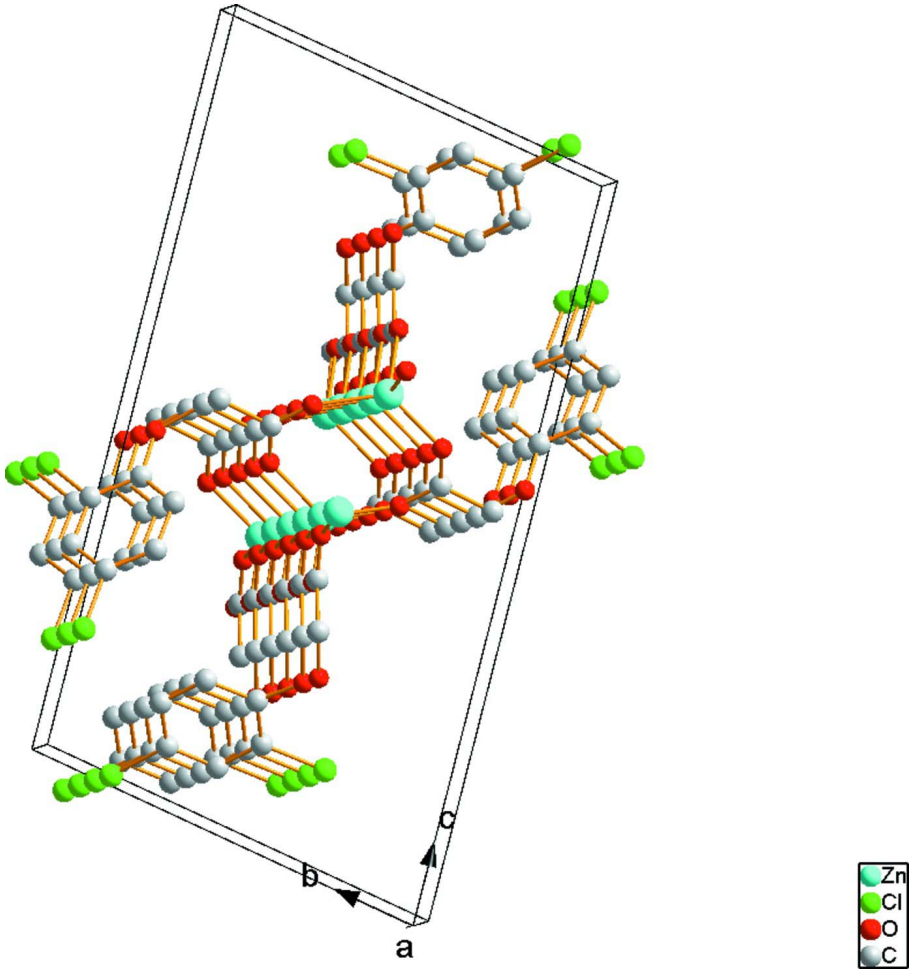

Figure 2

Part of the crystal packing, showing a polymeric chain viewed along the *a* axis. H atoms have been omitted for clarity.

***catena*-Poly[zinc(II)-bis[*μ*-2-(2,4-dichlorophenoxy)acetato]]**

Crystal data

$[Zn(C_8H_5Cl_2O_3)_2]$
$M_r = 505.43$
Triclinic, $P\overline{1}$
Hall symbol: -P 1
$a = 4.7322\ (10)$ Å
$b = 10.459\ (2)$ Å
$c = 18.979\ (4)$ Å
$\alpha = 79.340\ (2)°$
$\beta = 89.838\ (2)°$
$\gamma = 82.847\ (3)°$
$V = 915.8\ (3)$ Å^3
$Z = 2$
$F(000) = 504$
$D_x = 1.833$ Mg m^{-3}
Mo $K\alpha$ radiation, $\lambda = 0.71073$ Å
Cell parameters from 2642 reflections
$\theta = 2.5–27.7°$
$\mu = 1.96$ mm^{-1}
$T = 296$ K
Needle, colourless
0.56 × 0.21 × 0.16 mm

Data collection

Bruker SMART APEXII CCD diffractometer
Radiation source: fine-focus sealed tube
Graphite monochromator
ω scans
Absorption correction: multi-scan (*SADABS*; Sheldrick, 1996)
$T_{min} = 0.617$, $T_{max} = 0.731$

4737 measured reflections
3255 independent reflections
2849 reflections with $I > 2\sigma(I)$
$R_{int} = 0.021$
$\theta_{max} = 25.2°$, $\theta_{min} = 2.0°$
$h = -5 \rightarrow 5$
$k = -12 \rightarrow 8$
$l = -22 \rightarrow 21$

Refinement

Refinement on F^2
Least-squares matrix: full
$R[F^2 > 2\sigma(F^2)] = 0.030$
$wR(F^2) = 0.091$
$S = 0.94$
3255 reflections
244 parameters
0 restraints
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
$w = 1/[\sigma^2(F_o^2) + (0.0663P)^2]$ where $P = (F_o^2 + 2F_c^2)/3$
$(\Delta/\sigma)_{max} = 0.024$
$\Delta\rho_{max} = 0.34$ e Å^{-3}
$\Delta\rho_{min} = -0.39$ e Å^{-3}

Special details

Geometry. All esds (except the esd in the dihedral angle between two l.s. planes) are estimated using the full covariance matrix. The cell esds are taken into account individually in the estimation of esds in distances, angles and torsion angles; correlations between esds in cell parameters are only used when they are defined by crystal symmetry. An approximate (isotropic) treatment of cell esds is used for estimating esds 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 (Å^2)

	x	*y*	*z*	U_{iso}*/U_{eq}
Zn1	0.66929 (6)	0.53465 (3)	0.409892 (14)	0.02622 (12)
Cl2	0.46808 (18)	1.18539 (7)	0.33205 (4)	0.0500 (2)
Cl3	−0.18417 (19)	1.02366 (8)	0.14274 (4)	0.0538 (2)
O1	0.5602 (4)	0.69621 (17)	0.44411 (10)	0.0354 (4)
C2	0.2487 (7)	0.8626 (2)	0.48725 (14)	0.0368 (6)
H2A	0.0425	0.8686	0.4858	0.044*
H2B	0.3034	0.8854	0.5321	0.044*
C1	0.3707 (5)	0.7213 (2)	0.48733 (13)	0.0283 (5)
O2	0.3338 (4)	0.95654 (17)	0.43021 (10)	0.0381 (4)
C8	0.2572 (6)	1.0724 (2)	0.31164 (15)	0.0344 (6)
C3	0.2031 (6)	0.9673 (2)	0.36461 (14)	0.0336 (6)
C7	0.1426 (6)	1.0904 (3)	0.24305 (15)	0.0374 (6)
H7	0.1846	1.1593	0.2077	0.045*
C4	0.0218 (6)	0.8825 (3)	0.34807 (15)	0.0370 (6)
H4	−0.0198	0.8129	0.3830	0.044*
C6	−0.0369 (6)	1.0027 (3)	0.22834 (15)	0.0381 (6)
C5	−0.0982 (6)	0.9003 (3)	0.27995 (16)	0.0409 (7)
H5	−0.2199	0.8429	0.2693	0.049*
O5	0.4663 (4)	0.5250 (2)	0.32238 (10)	0.0379 (4)
C9	0.2002 (5)	0.5268 (2)	0.32394 (13)	0.0285 (5)
C10	0.0513 (6)	0.4985 (3)	0.25952 (14)	0.0365 (6)

H10A	−0.0637	0.4283	0.2752	0.044*
H10B	−0.0762	0.5759	0.2383	0.044*
O6	0.2391 (4)	0.46218 (19)	0.20660 (10)	0.0394 (5)
C11	0.3544 (6)	0.5571 (3)	0.16035 (14)	0.0335 (6)
C12	0.5536 (6)	0.5123 (3)	0.11352 (14)	0.0371 (6)
C14	0.6058 (7)	0.7315 (3)	0.05874 (16)	0.0440 (7)
C15	0.4104 (7)	0.7786 (3)	0.10480 (18)	0.0481 (7)
H15	0.3623	0.8684	0.1018	0.058*
C13	0.6794 (7)	0.5995 (3)	0.06313 (15)	0.0457 (7)
H13	0.8135	0.5685	0.0323	0.055*
C16	0.2865 (7)	0.6908 (3)	0.15540 (17)	0.0452 (7)
H16	0.1552	0.7224	0.1866	0.054*
Cl5	0.6472 (2)	0.34585 (8)	0.11910 (5)	0.0616 (3)
Cl4	0.7620 (2)	0.84005 (10)	−0.00545 (5)	0.0660 (3)
O7	0.0604 (4)	0.54895 (18)	0.37772 (9)	0.0331 (4)
O8	0.2660 (4)	0.64018 (18)	0.53311 (11)	0.0449 (5)

Atomic displacement parameters (Å²)

	U^{11}	U^{22}	U^{33}	U^{12}	U^{13}	U^{23}
Zn1	0.02447 (18)	0.02421 (18)	0.02907 (18)	−0.00534 (12)	0.00625 (12)	−0.00117 (12)
Cl2	0.0636 (5)	0.0330 (4)	0.0527 (5)	−0.0225 (4)	−0.0055 (4)	0.0049 (3)
Cl3	0.0667 (5)	0.0505 (5)	0.0438 (4)	−0.0079 (4)	−0.0080 (4)	−0.0074 (3)
O1	0.0368 (10)	0.0253 (9)	0.0450 (11)	−0.0078 (8)	0.0112 (9)	−0.0070 (8)
C2	0.0508 (17)	0.0239 (13)	0.0339 (14)	−0.0042 (12)	0.0067 (12)	−0.0011 (11)
C1	0.0325 (14)	0.0239 (12)	0.0274 (13)	−0.0053 (10)	−0.0012 (10)	−0.0002 (10)
O2	0.0522 (12)	0.0239 (9)	0.0361 (10)	−0.0091 (8)	0.0002 (9)	0.0025 (8)
C8	0.0348 (15)	0.0244 (13)	0.0427 (15)	−0.0077 (11)	0.0036 (12)	−0.0007 (11)
C3	0.0385 (15)	0.0225 (13)	0.0383 (14)	−0.0022 (11)	0.0054 (12)	−0.0030 (11)
C7	0.0443 (17)	0.0260 (13)	0.0390 (15)	−0.0049 (12)	0.0062 (12)	0.0018 (11)
C4	0.0425 (16)	0.0223 (13)	0.0433 (15)	−0.0063 (11)	0.0061 (12)	0.0026 (11)
C6	0.0412 (16)	0.0328 (14)	0.0388 (15)	0.0008 (12)	0.0007 (12)	−0.0063 (12)
C5	0.0425 (16)	0.0291 (14)	0.0518 (17)	−0.0075 (12)	0.0023 (13)	−0.0070 (12)
O5	0.0200 (9)	0.0564 (12)	0.0392 (10)	−0.0081 (8)	0.0039 (8)	−0.0116 (9)
C9	0.0257 (13)	0.0277 (13)	0.0306 (13)	−0.0048 (10)	0.0023 (10)	−0.0011 (10)
C10	0.0277 (14)	0.0499 (17)	0.0334 (14)	−0.0084 (12)	0.0027 (11)	−0.0091 (12)
O6	0.0429 (11)	0.0428 (11)	0.0352 (10)	−0.0103 (9)	0.0099 (8)	−0.0117 (9)
C11	0.0323 (14)	0.0419 (15)	0.0277 (13)	−0.0034 (12)	−0.0028 (11)	−0.0107 (11)
C12	0.0402 (16)	0.0382 (15)	0.0320 (14)	0.0038 (12)	0.0012 (11)	−0.0105 (12)
C14	0.0419 (17)	0.0481 (18)	0.0391 (16)	−0.0075 (14)	0.0001 (13)	0.0007 (13)
C15	0.0489 (18)	0.0373 (16)	0.0570 (19)	−0.0018 (14)	−0.0011 (15)	−0.0080 (14)
C13	0.0485 (18)	0.0522 (19)	0.0352 (15)	0.0003 (15)	0.0085 (13)	−0.0096 (14)
C16	0.0443 (17)	0.0421 (17)	0.0506 (17)	−0.0006 (14)	0.0137 (14)	−0.0152 (14)
Cl5	0.0829 (7)	0.0394 (4)	0.0605 (5)	0.0066 (4)	0.0191 (5)	−0.0135 (4)
Cl4	0.0775 (6)	0.0650 (6)	0.0514 (5)	−0.0220 (5)	0.0086 (4)	0.0081 (4)
O7	0.0256 (9)	0.0427 (11)	0.0335 (10)	−0.0108 (8)	0.0070 (8)	−0.0095 (8)
O8	0.0527 (13)	0.0266 (10)	0.0495 (12)	−0.0031 (9)	0.0156 (10)	0.0071 (9)

Geometric parameters (Å, º)

Zn1—O1	1.9315 (18)	C5—H5	0.93
Zn1—O8^{i}	1.9335 (18)	O5—C9	1.258 (3)
Zn1—O5	1.9458 (18)	C9—O7	1.256 (3)
Zn1—O7^{ii}	1.9622 (18)	C9—C10	1.507 (3)
Cl2—C8	1.735 (3)	C10—O6	1.412 (3)
Cl3—C6	1.735 (3)	C10—H10A	0.97
O1—C1	1.250 (3)	C10—H10B	0.97
C2—O2	1.415 (3)	O6—C11	1.365 (3)
C2—C1	1.518 (3)	C11—C16	1.382 (4)
C2—H2A	0.97	C11—C12	1.392 (4)
C2—H2B	0.97	C12—C13	1.382 (4)
C1—O8	1.245 (3)	C12—Cl5	1.726 (3)
O2—C3	1.372 (3)	C14—C13	1.369 (4)
C8—C3	1.395 (4)	C14—C15	1.380 (5)
C8—C7	1.383 (4)	C14—Cl4	1.735 (3)
C3—C4	1.383 (4)	C15—C16	1.383 (4)
C7—C6	1.389 (4)	C15—H15	0.93
C7—H7	0.93	C13—H13	0.93
C4—C5	1.385 (4)	C16—H16	0.93
C4—H4	0.93	O7—Zn1^{iii}	1.9622 (17)
C6—C5	1.371 (4)	O8—Zn1^{i}	1.9335 (18)
O1—Zn1—O8^{i}	126.97 (8)	C6—C5—H5	120.1
O1—Zn1—O5	113.18 (8)	C4—C5—H5	120.1
O8^{i}—Zn1—O5	107.88 (9)	C9—O5—Zn1	118.61 (16)
O1—Zn1—O7^{ii}	103.55 (8)	O7—C9—O5	121.7 (2)
O8^{i}—Zn1—O7^{ii}	98.21 (8)	O7—C9—C10	120.3 (2)
O5—Zn1—O7^{ii}	103.04 (7)	O5—C9—C10	118.0 (2)
C1—O1—Zn1	128.96 (17)	O6—C10—C9	113.7 (2)
O2—C2—C1	115.7 (2)	O6—C10—H10A	108.8
O2—C2—H2A	108.4	C9—C10—H10A	108.8
C1—C2—H2A	108.4	O6—C10—H10B	108.8
O2—C2—H2B	108.4	C9—C10—H10B	108.8
C1—C2—H2B	108.4	H10A—C10—H10B	107.7
H2A—C2—H2B	107.4	C11—O6—C10	119.4 (2)
O8—C1—O1	126.4 (2)	O6—C11—C16	125.9 (2)
O8—C1—C2	113.8 (2)	O6—C11—C12	115.7 (2)
O1—C1—C2	119.8 (2)	C16—C11—C12	118.4 (3)
C3—O2—C2	117.3 (2)	C13—C12—C11	120.8 (3)
C3—C8—C7	121.6 (2)	C13—C12—Cl5	119.6 (2)
C3—C8—Cl2	119.6 (2)	C11—C12—Cl5	119.6 (2)
C7—C8—Cl2	118.8 (2)	C13—C14—C15	120.7 (3)
O2—C3—C8	116.8 (2)	C13—C14—Cl4	119.3 (2)
O2—C3—C4	124.7 (2)	C15—C14—Cl4	120.0 (3)
C8—C3—C4	118.5 (3)	C16—C15—C14	119.4 (3)
C8—C7—C6	118.2 (3)	C16—C15—H15	120.3

C8—C7—H7	120.9	C14—C15—H15	120.3
C6—C7—H7	120.9	C14—C13—C12	119.7 (3)
C3—C4—C5	120.6 (3)	C14—C13—H13	120.2
C3—C4—H4	119.7	C12—C13—H13	120.2
C5—C4—H4	119.7	C15—C16—C11	121.1 (3)
C5—C6—C7	121.3 (3)	C15—C16—H16	119.5
C5—C6—Cl3	119.7 (2)	C11—C16—H16	119.5
C7—C6—Cl3	119.0 (2)	C9—O7—Zn1^{iii}	135.66 (16)
C6—C5—C4	119.8 (3)	C1—O8—Zn1^{i}	144.34 (19)

Symmetry codes: (i) −*x*+1, −*y*+1, −*z*+1; (ii) *x*+1, *y*, *z*; (iii) *x*−1, *y*, *z*.

Hydrogen-bond geometry (Å, º)

D—H···*A*	*D*—H	H···*A*	*D*···*A*	*D*—H···*A*
C2—H2*B*···Cl2^{iv}	0.97	2.73	3.610 (3)	151

Symmetry code: (iv) −*x*+1, −*y*+2, −*z*+1.