

Aqua(4-hydroxybenzoato- κ O)- (4-hydroxybenzoato- κ^2 O,O')(1,10- phenanthroline- κ^2 N,N')zinc(II) monohydrate

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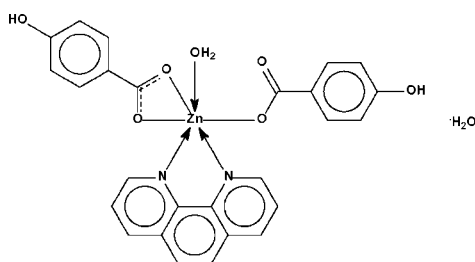
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Key indicators: single-crystal X-ray study; $T = 295$ K; mean $\sigma(\text{C}-\text{C}) = 0.005$ Å; R factor = 0.044; wR factor = 0.131; data-to-parameter ratio = 16.5.

The Zn atom in the title compound, $[\text{Zn}(\text{C}_7\text{H}_5\text{O}_3)_2(\text{C}_{12}\text{H}_8\text{N}_2)(\text{H}_2\text{O})] \cdot \text{H}_2\text{O}$, exists in a distorted *cis*- ZnN_2O_4 octahedral coordination geometry. One of the 4-hydroxybenzoate anions chelates in a bidentate manner whereas the other is monodentate. The complex molecules are linked through the uncoordinated water molecules into a hydrogen-bonded sheet structure.

Related literature

For related zinc bis(4-hydroxybenzoate) structures containing an *N*-heterocycle, see: Hökelek & Necefoglu (1996); Nadzhafov *et al.* (1981); Necefoglu *et al.* (2002); Wang & Okabe (2005); Zheng *et al.* (2006).



Experimental

Crystal data

 $[\text{Zn}(\text{C}_7\text{H}_5\text{O}_3)_2(\text{C}_{12}\text{H}_8\text{N}_2)(\text{H}_2\text{O})] \cdot \text{H}_2\text{O}$
 $M_r = 555.83$

 Monoclinic, $P2_1/c$
 $a = 11.1169$ (5) Å

 $b = 19.738$ (1) Å

 $c = 11.5503$ (6) Å

 $\beta = 106.298$ (1)°

 $V = 2432.5$ (2) Å³
 $Z = 4$

 Mo $K\alpha$ radiation

 $\mu = 1.06$ mm⁻¹
 $T = 295$ (2) K

 $0.30 \times 0.24 \times 0.18$ mm

Data collection

Rigaku R-Axis RAPID diffractometer
Absorption correction: multi-scan (ABSCOR; Higashi, 1995)
 $T_{\min} = 0.674$, $T_{\max} = 0.832$

23098 measured reflections
5541 independent reflections
3924 reflections with $I > 2\sigma(I)$
 $R_{\text{int}} = 0.036$

Refinement

$R[F^2 > 2\sigma(F^2)] = 0.044$
 $wR(F^2) = 0.131$
 $S = 1.08$
5541 reflections

336 parameters
H-atom parameters constrained
 $\Delta\rho_{\text{max}} = 0.67$ e Å⁻³
 $\Delta\rho_{\text{min}} = -0.88$ e Å⁻³

Table 1

Selected bond lengths (Å).

Zn1—O1	2.025 (2)	Zn1—O1w	2.143 (2)
Zn1—O2	2.470 (2)	Zn1—N1	2.075 (2)
Zn1—O4	2.069 (2)	Zn1—N2	2.155 (2)

Table 2

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>
O3—H3o...O2 ⁱ	0.85	1.81	2.619 (3)	157
O6—H6o...O4 ⁱⁱ	0.85	1.90	2.742 (3)	171
O1w—H1w1...O2w ⁱⁱⁱ	0.85	2.01	2.819 (4)	158
O1w—H1w2...O5	0.85	1.84	2.599 (3)	148
O2w—H2w1...O5	0.85	2.04	2.801 (4)	148

Symmetry codes: (i) $x, -y + \frac{1}{2}, z - \frac{1}{2}$; (ii) $x, -y + \frac{3}{2}, z - \frac{1}{2}$; (iii) $-x + 1, -y + 1, -z + 1$.

Data collection: *RAPID-AUTO* (Rigaku, 1998); cell refinement: *RAPID-AUTO*; data reduction: *CrystalStructure* (Rigaku/MS, 2002); 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).

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Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: HB2688).

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Acta Cryst. (2008). E64, m422 [doi:10.1107/S1600536808002249]

Aqua(4-hydroxybenzoato- κO)(4-hydroxybenzoato- $\kappa^2 O, O'$)(1,10-phenanthroline- $\kappa^2 N, N'$)zinc(II) monohydrate

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Comment

The title compound, (I), extends the range of adducts of zinc bis(4-hydroxybenzoate) with *N*-heterocycles: For the *N,N*-diethylnicotinamide adduct, see: Hökelek & Necefouglu (1996). For the pyridine adduct, see; Nadzhafov *et al.* (1981). For the nicotinamide adduct, see: Necefoğlu *et al.* (2002). For the bis(2-pyridyl)amine adduct, see: Wang & Okabe (2005). For the benzimidazole adduct, see: Zheng *et al.* (2006).

The Zn atom in (I) adopts a distorted *cis*-ZnN₂O₄ coordination geometry. One of the 4-hydroxybenzoate anions chelates in an anisobidentate manner whereas the other is unidentate (Table 1, Fig. 1).

In the crystal, the complex molecules are linked through the uncoordinated water molecules into a hydrogen-bonded sheet structure (Table 2).

Experimental

Zinc diacetate dihydrate (1 mmol), 1,10-phenanthroline (2 mmol) and 4-hydroxybenzoic acid (2 mmol) were dissolved in aqueous ethanol. Colourless blocks of (I) were isolated after several days.

Refinement

The carbon-bound and hydroxyl H atoms were placed in calculated positions (C–H = 0.93, O–H = 0.85 Å) and refined as riding with $U_{\text{iso}}(\text{H})$ 1.2–1.5 $U_{\text{eq}}(\text{C}, \text{O})$. The water H atoms were placed in chemically reasonable positions with O–H = 0.85 Å on the basis of likely hydrogen bonding interactions and refined as riding with $U_{\text{iso}}(\text{H}) = 1.2U_{\text{eq}}(\text{O})$.

Figures

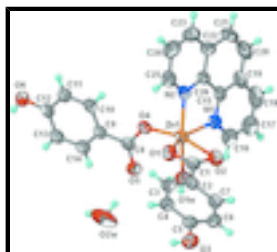


Fig. 1. View of the molecular structure of (I) with displacement ellipsoids drawn at the 50% probability level (H atoms drawn as spheres of arbitrary radius).

supplementary materials

Aquabis(4-hydroxybenzoato- κ O)(4-hydroxybenzoato- κ^2 O, O')(1,10-phenanthroline- κ^2 N, N')zinc(II) monohydrate

Crystal data

$[\text{Zn}(\text{C}_7\text{H}_5\text{O}_3)_2(\text{C}_{12}\text{H}_8\text{N}_2)(\text{H}_2\text{O})]\cdot\text{H}_2\text{O}$	$F(000) = 1144$
$M_r = 555.83$	$D_x = 1.518 \text{ Mg m}^{-3}$
Monoclinic, $P2_1/c$	Mo $K\alpha$ radiation, $\lambda = 0.71073 \text{ \AA}$
Hall symbol: -P 2ybc	Cell parameters from 10359 reflections
$a = 11.1169 (5) \text{ \AA}$	$\theta = 3.1\text{--}27.5^\circ$
$b = 19.738 (1) \text{ \AA}$	$\mu = 1.06 \text{ mm}^{-1}$
$c = 11.5503 (6) \text{ \AA}$	$T = 295 \text{ K}$
$\beta = 106.298 (1)^\circ$	Block, colorless
$V = 2432.5 (2) \text{ \AA}^3$	$0.30 \times 0.24 \times 0.18 \text{ mm}$
$Z = 4$	

Data collection

Rigaku R-AXIS RAPID diffractometer	5541 independent reflections
Radiation source: fine-focus sealed tube graphite	3924 reflections with $I > 2\sigma(I)$
Detector resolution: $10.000 \text{ pixels mm}^{-1}$	$R_{\text{int}} = 0.036$
ω scans	$\theta_{\text{max}} = 27.5^\circ$, $\theta_{\text{min}} = 3.1^\circ$
Absorption correction: multi-scan (ABSCOR; Higashi, 1995)	$h = -14 \rightarrow 14$
$T_{\text{min}} = 0.674$, $T_{\text{max}} = 0.832$	$k = -25 \rightarrow 25$
23098 measured reflections	$l = -14 \rightarrow 14$

Refinement

Refinement on F^2	Primary atom site location: structure-invariant direct methods
Least-squares matrix: full	Secondary atom site location: difference Fourier map
$R[F^2 > 2\sigma(F^2)] = 0.044$	Hydrogen site location: inferred from neighbouring sites
$wR(F^2) = 0.131$	H-atom parameters constrained
$S = 1.08$	$w = 1/[\sigma^2(F_o^2) + (0.0632P)^2 + 0.8555P]$
5541 reflections	where $P = (F_o^2 + 2F_c^2)/3$
336 parameters	$(\Delta/\sigma)_{\text{max}} = 0.001$
0 restraints	$\Delta\rho_{\text{max}} = 0.67 \text{ e \AA}^{-3}$
	$\Delta\rho_{\text{min}} = -0.88 \text{ e \AA}^{-3}$

Fractional atomic coordinates and isotropic or equivalent isotropic displacement parameters (\AA^2)

x	y	z	$U_{\text{iso}}^*/U_{\text{eq}}$
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Zn1	0.23086 (3)	0.517846 (16)	0.72570 (3)	0.05090 (13)
O1	0.2083 (2)	0.43474 (11)	0.6190 (2)	0.0681 (6)
O2	0.2863 (2)	0.40272 (11)	0.8070 (2)	0.0663 (6)
O3	0.3803 (2)	0.13818 (12)	0.5298 (2)	0.0729 (6)
H3o	0.3548	0.1357	0.4533	0.109*
O4	0.19613 (18)	0.60020 (9)	0.61024 (17)	0.0507 (4)
O5	0.3820 (2)	0.61864 (13)	0.5776 (2)	0.0725 (6)
O6	0.03862 (19)	0.81782 (11)	0.19250 (19)	0.0612 (5)
H6o	0.0924	0.8392	0.1669	0.092*
O1w	0.4267 (2)	0.52299 (12)	0.7387 (2)	0.0749 (7)
H1w1	0.4483	0.4874	0.7080	0.112*
H1w2	0.4409	0.5576	0.7007	0.112*
O2w	0.5117 (3)	0.5725 (2)	0.4171 (4)	0.1517 (16)
H2w1	0.4623	0.5704	0.4614	0.182*
H2w2	0.4914	0.6047	0.3664	0.182*
N1	0.2575 (2)	0.55676 (12)	0.8980 (2)	0.0524 (6)
N2	0.0408 (2)	0.52319 (11)	0.7369 (2)	0.0478 (5)
C1	0.2581 (3)	0.38959 (15)	0.6961 (3)	0.0555 (7)
C2	0.2870 (3)	0.32281 (14)	0.6520 (3)	0.0498 (6)
C3	0.2367 (3)	0.30452 (15)	0.5320 (3)	0.0559 (7)
H3	0.1830	0.3342	0.4795	0.067*
C4	0.2653 (3)	0.24292 (16)	0.4895 (3)	0.0589 (7)
H4	0.2291	0.2308	0.4095	0.071*
C5	0.3479 (3)	0.19928 (15)	0.5664 (3)	0.0549 (7)
C6	0.4005 (3)	0.21700 (15)	0.6863 (3)	0.0574 (7)
H6	0.4563	0.1879	0.7381	0.069*
C7	0.3693 (3)	0.27787 (15)	0.7279 (3)	0.0555 (7)
H7	0.4039	0.2893	0.8085	0.067*
C8	0.2667 (3)	0.62967 (14)	0.5549 (3)	0.0495 (6)
C9	0.2074 (3)	0.67968 (14)	0.4592 (2)	0.0480 (6)
C10	0.0776 (3)	0.68671 (15)	0.4180 (3)	0.0539 (7)
H10	0.0269	0.6598	0.4508	0.065*
C11	0.0234 (3)	0.73289 (16)	0.3295 (3)	0.0562 (7)
H11	-0.0634	0.7370	0.3031	0.067*
C12	0.0982 (3)	0.77342 (14)	0.2796 (2)	0.0509 (6)
C13	0.2280 (3)	0.76739 (15)	0.3202 (3)	0.0562 (7)
H13	0.2786	0.7945	0.2874	0.067*
C14	0.2815 (3)	0.72096 (14)	0.4094 (3)	0.0543 (7)
H14	0.3683	0.7172	0.4366	0.065*
C15	0.1503 (3)	0.56991 (13)	0.9281 (2)	0.0501 (6)
C16	0.3666 (3)	0.57100 (18)	0.9768 (3)	0.0701 (9)
H16	0.4399	0.5617	0.9560	0.084*
C17	0.3757 (4)	0.5985 (2)	1.0862 (4)	0.0874 (12)
H17	0.4542	0.6070	1.1392	0.105*
C18	0.2708 (5)	0.6137 (2)	1.1192 (3)	0.0868 (12)
H18	0.2774	0.6333	1.1940	0.104*
C19	0.1501 (4)	0.59957 (16)	1.0383 (3)	0.0669 (9)
C20	0.0339 (5)	0.61090 (19)	1.0621 (4)	0.0853 (12)
H20	0.0331	0.6306	1.1350	0.102*

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C21	-0.0760 (4)	0.5941 (2)	0.9824 (4)	0.0843 (12)
H21	-0.1505	0.6015	1.0022	0.101*
C22	-0.0799 (3)	0.56479 (16)	0.8669 (3)	0.0640 (8)
C23	-0.1891 (4)	0.5470 (2)	0.7791 (4)	0.0847 (12)
H23	-0.2668	0.5547	0.7923	0.102*
C24	-0.1823 (4)	0.5185 (2)	0.6742 (4)	0.0856 (12)
H24	-0.2551	0.5069	0.6150	0.103*
C25	-0.0643 (3)	0.50667 (18)	0.6557 (3)	0.0654 (8)
H25	-0.0604	0.4866	0.5840	0.078*
C26	0.0332 (3)	0.55205 (13)	0.8411 (2)	0.0472 (6)

Atomic displacement parameters (\AA^2)

	U^{11}	U^{22}	U^{33}	U^{12}	U^{13}	U^{23}
Zn1	0.0592 (2)	0.0422 (2)	0.0576 (2)	0.00058 (14)	0.02675 (17)	0.00114 (14)
O1	0.0915 (16)	0.0451 (11)	0.0708 (14)	0.0159 (11)	0.0279 (12)	0.0032 (10)
O2	0.0923 (16)	0.0487 (12)	0.0629 (13)	0.0041 (11)	0.0302 (12)	-0.0055 (10)
O3	0.0831 (16)	0.0597 (13)	0.0764 (14)	0.0212 (12)	0.0235 (13)	-0.0112 (12)
O4	0.0592 (11)	0.0410 (10)	0.0603 (11)	0.0033 (8)	0.0306 (9)	0.0069 (8)
O5	0.0613 (13)	0.0814 (16)	0.0829 (15)	0.0100 (12)	0.0336 (12)	0.0261 (13)
O6	0.0682 (13)	0.0559 (12)	0.0626 (12)	-0.0031 (10)	0.0232 (10)	0.0126 (10)
O1w	0.0624 (14)	0.0749 (16)	0.0959 (18)	0.0095 (11)	0.0359 (13)	0.0233 (13)
O2w	0.140 (3)	0.187 (4)	0.162 (3)	-0.004 (3)	0.099 (3)	-0.052 (3)
N1	0.0549 (14)	0.0441 (13)	0.0575 (13)	-0.0065 (10)	0.0145 (12)	-0.0008 (11)
N2	0.0518 (13)	0.0453 (12)	0.0484 (12)	-0.0067 (10)	0.0174 (11)	0.0025 (10)
C1	0.0639 (17)	0.0443 (15)	0.0639 (18)	-0.0001 (13)	0.0271 (15)	0.0002 (14)
C2	0.0550 (15)	0.0416 (14)	0.0586 (16)	0.0015 (12)	0.0257 (13)	0.0001 (12)
C3	0.0608 (17)	0.0522 (17)	0.0568 (16)	0.0093 (14)	0.0200 (14)	0.0030 (13)
C4	0.0648 (18)	0.0591 (18)	0.0544 (16)	0.0071 (15)	0.0193 (15)	-0.0058 (14)
C5	0.0582 (16)	0.0458 (15)	0.0656 (18)	0.0065 (13)	0.0254 (14)	-0.0047 (14)
C6	0.0605 (17)	0.0492 (16)	0.0620 (17)	0.0121 (13)	0.0162 (14)	0.0017 (14)
C7	0.0614 (17)	0.0499 (16)	0.0568 (16)	0.0013 (13)	0.0193 (14)	-0.0018 (13)
C8	0.0573 (16)	0.0418 (14)	0.0552 (15)	-0.0020 (12)	0.0254 (14)	-0.0023 (12)
C9	0.0580 (16)	0.0409 (14)	0.0505 (14)	-0.0050 (12)	0.0244 (13)	-0.0025 (12)
C10	0.0591 (17)	0.0531 (16)	0.0557 (16)	-0.0110 (13)	0.0263 (14)	0.0028 (13)
C11	0.0554 (16)	0.0565 (17)	0.0588 (17)	-0.0053 (13)	0.0197 (14)	0.0047 (14)
C12	0.0654 (17)	0.0408 (14)	0.0511 (15)	-0.0042 (12)	0.0241 (14)	-0.0017 (12)
C13	0.0648 (18)	0.0468 (15)	0.0663 (18)	-0.0070 (13)	0.0336 (15)	0.0051 (14)
C14	0.0562 (16)	0.0483 (16)	0.0630 (17)	-0.0055 (13)	0.0242 (14)	0.0016 (13)
C15	0.0711 (18)	0.0350 (13)	0.0475 (14)	-0.0020 (12)	0.0220 (14)	0.0025 (11)
C16	0.068 (2)	0.065 (2)	0.071 (2)	-0.0107 (17)	0.0100 (17)	-0.0038 (17)
C17	0.089 (3)	0.088 (3)	0.077 (2)	-0.018 (2)	0.009 (2)	-0.005 (2)
C18	0.136 (4)	0.065 (2)	0.0498 (18)	-0.021 (2)	0.011 (2)	-0.0128 (16)
C19	0.105 (3)	0.0459 (17)	0.0577 (18)	0.0020 (17)	0.0366 (19)	-0.0015 (14)
C20	0.138 (4)	0.063 (2)	0.071 (2)	0.011 (2)	0.055 (3)	-0.0033 (18)
C21	0.105 (3)	0.068 (2)	0.108 (3)	0.026 (2)	0.076 (3)	0.017 (2)
C22	0.0641 (19)	0.0555 (18)	0.082 (2)	0.0096 (15)	0.0358 (18)	0.0179 (16)
C23	0.062 (2)	0.085 (3)	0.115 (3)	0.0119 (19)	0.039 (2)	0.032 (3)

C24	0.056 (2)	0.092 (3)	0.099 (3)	-0.0126 (19)	0.005 (2)	0.020 (2)
C25	0.069 (2)	0.066 (2)	0.0607 (18)	-0.0145 (16)	0.0163 (16)	0.0010 (15)
C26	0.0586 (16)	0.0359 (13)	0.0534 (15)	0.0016 (11)	0.0259 (13)	0.0065 (11)

Geometric parameters (Å, °)

Zn1—O1	2.025 (2)	C7—H7	0.9300
Zn1—O2	2.470 (2)	C8—C9	1.490 (4)
Zn1—O4	2.069 (2)	C9—C14	1.392 (4)
Zn1—O1w	2.143 (2)	C9—C10	1.394 (4)
Zn1—N1	2.075 (2)	C10—C11	1.376 (4)
Zn1—N2	2.155 (2)	C10—H10	0.9300
O1—C1	1.272 (4)	C11—C12	1.391 (4)
O2—C1	1.257 (4)	C11—H11	0.9300
O3—C5	1.359 (3)	C12—C13	1.392 (4)
O3—H3o	0.8501	C13—C14	1.383 (4)
O4—C8	1.282 (3)	C13—H13	0.9300
O5—C8	1.253 (3)	C14—H14	0.9300
O6—C12	1.358 (3)	C15—C19	1.402 (4)
O6—H6o	0.8501	C15—C26	1.447 (4)
O1w—H1w1	0.8500	C16—C17	1.353 (6)
O1w—H1w2	0.8501	C16—H16	0.9300
O2w—H2w1	0.8499	C17—C18	1.358 (6)
O2w—H2w2	0.8500	C17—H17	0.9300
N1—C16	1.326 (4)	C18—C19	1.431 (5)
N1—C15	1.357 (4)	C18—H18	0.9300
N2—C25	1.318 (4)	C19—C20	1.412 (6)
N2—C26	1.356 (3)	C20—C21	1.349 (6)
C1—C2	1.481 (4)	C20—H20	0.9300
C2—C3	1.388 (4)	C21—C22	1.443 (5)
C2—C7	1.394 (4)	C21—H21	0.9300
C3—C4	1.381 (4)	C22—C23	1.391 (5)
C3—H3	0.9300	C22—C26	1.394 (4)
C4—C5	1.384 (4)	C23—C24	1.357 (6)
C4—H4	0.9300	C23—H23	0.9300
C5—C6	1.389 (4)	C24—C25	1.406 (6)
C6—C7	1.374 (4)	C24—H24	0.9300
C6—H6	0.9300	C25—H25	0.9300
O1—Zn1—O4	105.98 (8)	O4—C8—C9	117.7 (2)
O1—Zn1—N1	147.52 (9)	C14—C9—C10	118.4 (3)
O4—Zn1—N1	106.14 (8)	C14—C9—C8	120.3 (3)
O1—Zn1—O1w	91.75 (10)	C10—C9—C8	121.3 (2)
O4—Zn1—O1w	90.57 (8)	C11—C10—C9	121.1 (3)
N1—Zn1—O1w	92.66 (10)	C11—C10—H10	119.5
O1—Zn1—N2	96.91 (9)	C9—C10—H10	119.5
O4—Zn1—N2	89.51 (8)	C10—C11—C12	120.0 (3)
N1—Zn1—N2	78.65 (9)	C10—C11—H11	120.0
O1w—Zn1—N2	170.97 (10)	C12—C11—H11	120.0
O1—Zn1—O2	57.39 (8)	O6—C12—C11	117.0 (3)

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O4—Zn1—O2	161.68 (7)	O6—C12—C13	123.3 (3)
N1—Zn1—O2	91.30 (8)	C11—C12—C13	119.7 (3)
O1w—Zn1—O2	82.91 (8)	C14—C13—C12	119.7 (3)
N2—Zn1—O2	99.59 (8)	C14—C13—H13	120.1
C1—O1—Zn1	100.74 (19)	C12—C13—H13	120.1
C1—O2—Zn1	80.68 (17)	C13—C14—C9	121.1 (3)
C5—O3—H3o	109.5	C13—C14—H14	119.5
C8—O4—Zn1	130.40 (18)	C9—C14—H14	119.5
C12—O6—H6o	109.5	N1—C15—C19	122.5 (3)
Zn1—O1w—H1w1	109.4	N1—C15—C26	117.3 (2)
Zn1—O1w—H1w2	109.5	C19—C15—C26	120.1 (3)
H1w1—O1w—H1w2	109.5	N1—C16—C17	122.7 (4)
H2w1—O2w—H2w2	111.0	N1—C16—H16	118.7
C16—N1—C15	118.9 (3)	C17—C16—H16	118.7
C16—N1—Zn1	126.4 (2)	C16—C17—C18	120.4 (4)
C15—N1—Zn1	114.65 (18)	C16—C17—H17	119.8
C25—N2—C26	118.2 (3)	C18—C17—H17	119.8
C25—N2—Zn1	129.3 (2)	C17—C18—C19	119.6 (3)
C26—N2—Zn1	112.33 (18)	C17—C18—H18	120.2
O2—C1—O1	120.3 (3)	C19—C18—H18	120.2
O2—C1—C2	121.2 (3)	C15—C19—C20	118.5 (3)
O1—C1—C2	118.5 (3)	C15—C19—C18	115.8 (3)
O2—C1—Zn1	70.63 (16)	C20—C19—C18	125.7 (3)
O1—C1—Zn1	50.35 (14)	C21—C20—C19	122.1 (3)
C2—C1—Zn1	164.2 (2)	C21—C20—H20	118.9
C3—C2—C7	118.2 (3)	C19—C20—H20	118.9
C3—C2—C1	120.6 (3)	C20—C21—C22	121.0 (3)
C7—C2—C1	121.2 (3)	C20—C21—H21	119.5
C4—C3—C2	121.0 (3)	C22—C21—H21	119.5
C4—C3—H3	119.5	C23—C22—C26	117.0 (3)
C2—C3—H3	119.5	C23—C22—C21	124.7 (3)
C3—C4—C5	119.8 (3)	C26—C22—C21	118.3 (3)
C3—C4—H4	120.1	C24—C23—C22	120.0 (3)
C5—C4—H4	120.1	C24—C23—H23	120.0
O3—C5—C4	122.6 (3)	C22—C23—H23	120.0
O3—C5—C6	117.3 (3)	C23—C24—C25	119.5 (4)
C4—C5—C6	120.0 (3)	C23—C24—H24	120.2
C7—C6—C5	119.5 (3)	C25—C24—H24	120.2
C7—C6—H6	120.3	N2—C25—C24	122.0 (4)
C5—C6—H6	120.3	N2—C25—H25	119.0
C6—C7—C2	121.5 (3)	C24—C25—H25	119.0
C6—C7—H7	119.3	N2—C26—C22	123.4 (3)
C2—C7—H7	119.3	N2—C26—C15	116.8 (2)
O5—C8—O4	123.4 (3)	C22—C26—C15	119.8 (3)
O5—C8—C9	118.9 (2)		
O4—Zn1—O1—C1	166.06 (18)	C1—C2—C7—C6	-177.4 (3)
N1—Zn1—O1—C1	-22.8 (3)	Zn1—O4—C8—O5	-11.7 (4)
O1w—Zn1—O1—C1	75.0 (2)	Zn1—O4—C8—C9	168.48 (17)
N2—Zn1—O1—C1	-102.5 (2)	O5—C8—C9—C14	-9.9 (4)

O2—Zn1—O1—C1	-5.47 (18)	O4—C8—C9—C14	169.9 (3)
O1—Zn1—O2—C1	5.51 (18)	O5—C8—C9—C10	170.5 (3)
O4—Zn1—O2—C1	-21.3 (3)	O4—C8—C9—C10	-9.7 (4)
N1—Zn1—O2—C1	176.32 (19)	C14—C9—C10—C11	0.5 (4)
O1w—Zn1—O2—C1	-91.16 (19)	C8—C9—C10—C11	-179.9 (3)
N2—Zn1—O2—C1	97.61 (19)	C9—C10—C11—C12	0.1 (5)
O1—Zn1—O4—C8	-80.1 (2)	C10—C11—C12—O6	179.9 (3)
N1—Zn1—O4—C8	104.8 (2)	C10—C11—C12—C13	-0.5 (4)
O1w—Zn1—O4—C8	11.9 (2)	O6—C12—C13—C14	179.9 (3)
N2—Zn1—O4—C8	-177.1 (2)	C11—C12—C13—C14	0.3 (4)
O2—Zn1—O4—C8	-56.9 (4)	C12—C13—C14—C9	0.3 (5)
C1—Zn1—O4—C8	-70.8 (3)	C10—C9—C14—C13	-0.7 (4)
O1—Zn1—N1—C16	92.9 (3)	C8—C9—C14—C13	179.7 (3)
O4—Zn1—N1—C16	-95.9 (3)	C16—N1—C15—C19	1.6 (4)
O1w—Zn1—N1—C16	-4.6 (3)	Zn1—N1—C15—C19	-176.6 (2)
N2—Zn1—N1—C16	177.9 (3)	C16—N1—C15—C26	-178.4 (3)
O2—Zn1—N1—C16	78.4 (3)	Zn1—N1—C15—C26	3.4 (3)
C1—Zn1—N1—C16	80.4 (3)	C15—N1—C16—C17	-0.3 (5)
O1—Zn1—N1—C15	-89.0 (3)	Zn1—N1—C16—C17	177.7 (3)
O4—Zn1—N1—C15	82.14 (19)	N1—C16—C17—C18	-1.0 (6)
O1w—Zn1—N1—C15	173.51 (19)	C16—C17—C18—C19	1.0 (6)
N2—Zn1—N1—C15	-4.01 (18)	N1—C15—C19—C20	-179.6 (3)
O2—Zn1—N1—C15	-103.54 (19)	C26—C15—C19—C20	0.5 (4)
C1—Zn1—N1—C15	-101.5 (2)	N1—C15—C19—C18	-1.5 (4)
O1—Zn1—N2—C25	-33.1 (3)	C26—C15—C19—C18	178.5 (3)
O4—Zn1—N2—C25	73.0 (3)	C17—C18—C19—C15	0.2 (5)
N1—Zn1—N2—C25	179.5 (3)	C17—C18—C19—C20	178.1 (4)
O2—Zn1—N2—C25	-91.1 (3)	C15—C19—C20—C21	0.1 (5)
C1—Zn1—N2—C25	-62.0 (3)	C18—C19—C20—C21	-177.8 (4)
O1—Zn1—N2—C26	151.55 (17)	C19—C20—C21—C22	-1.4 (6)
O4—Zn1—N2—C26	-102.41 (18)	C20—C21—C22—C23	-178.9 (4)
N1—Zn1—N2—C26	4.16 (17)	C20—C21—C22—C26	2.2 (5)
O2—Zn1—N2—C26	93.56 (17)	C26—C22—C23—C24	0.2 (5)
C1—Zn1—N2—C26	122.68 (18)	C21—C22—C23—C24	-178.7 (4)
Zn1—O2—C1—O1	-8.6 (3)	C22—C23—C24—C25	0.5 (6)
Zn1—O2—C1—C2	168.4 (3)	C26—N2—C25—C24	0.4 (5)
Zn1—O1—C1—O2	10.5 (3)	Zn1—N2—C25—C24	-174.8 (2)
Zn1—O1—C1—C2	-166.6 (2)	C23—C24—C25—N2	-0.8 (6)
O2—C1—C2—C3	168.8 (3)	C25—N2—C26—C22	0.3 (4)
O1—C1—C2—C3	-14.2 (4)	Zn1—N2—C26—C22	176.3 (2)
O2—C1—C2—C7	-14.0 (4)	C25—N2—C26—C15	-179.7 (3)
O1—C1—C2—C7	163.1 (3)	Zn1—N2—C26—C15	-3.7 (3)
C7—C2—C3—C4	1.3 (4)	C23—C22—C26—N2	-0.6 (4)
C1—C2—C3—C4	178.7 (3)	C21—C22—C26—N2	178.3 (3)
C2—C3—C4—C5	-1.8 (5)	C23—C22—C26—C15	179.4 (3)
C3—C4—C5—O3	-179.6 (3)	C21—C22—C26—C15	-1.7 (4)
C3—C4—C5—C6	1.0 (5)	N1—C15—C26—N2	0.4 (4)
O3—C5—C6—C7	-179.1 (3)	C19—C15—C26—N2	-179.6 (2)
C4—C5—C6—C7	0.3 (5)	N1—C15—C26—C22	-179.6 (2)

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C5—C6—C7—C2	-0.8 (5)	C19—C15—C26—C22	0.4 (4)
C3—C2—C7—C6	0.0 (4)		

Hydrogen-bond geometry (Å, °)

<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>
O3—H3o \cdots O2 ⁱ	0.85	1.81	2.619 (3)	157
O6—H6o \cdots O4 ⁱⁱ	0.85	1.90	2.742 (3)	171
O1w—H1w1 \cdots O2w ⁱⁱⁱ	0.85	2.01	2.819 (4)	158
O1w—H1w2 \cdots O5	0.85	1.84	2.599 (3)	148
O2w—H2w1 \cdots O5	0.85	2.04	2.801 (4)	148

Symmetry codes: (i) $x, -y+1/2, z-1/2$; (ii) $x, -y+3/2, z-1/2$; (iii) $-x+1, -y+1, -z+1$.

Fig. 1

