

catena-Poly[[*(4-formylbenzoato-κO¹)-(isonicotinamide-κN¹)zinc(II)-μ-4-formylbenzoato-κ²O¹:O^{1'}*]]

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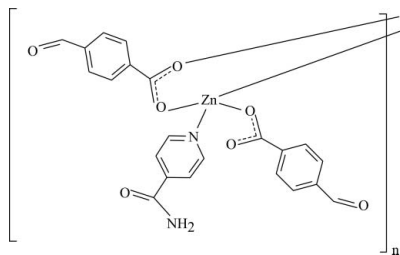
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Key indicators: single-crystal X-ray study; $T = 100$ K; mean $\sigma(\text{C}-\text{C}) = 0.002$ Å; R factor = 0.028; wR factor = 0.080; data-to-parameter ratio = 16.2.

In the title compound, $[\text{Zn}(\text{C}_8\text{H}_5\text{O}_3)_2(\text{C}_6\text{H}_6\text{N}_2\text{O})]_n$, the Zn^{II} ion is tetrahedrally coordinated by two formylbenzoate (FB) and one isonicotinamide (INA) ligands while symmetry-related FB ligands bridge adjacent Zn^{II} ions, forming polymeric chains along the b axis. The carboxylate groups in the two FB ions are twisted away from the attached benzene ring by 9.07 (2) and 26.2 (2)°. The two benzene rings of the FB ions are oriented at a dihedral angle of 81.30 (5)°. In the crystal, adjacent polymeric chains interact *via* $\text{N}-\text{H}\cdots\text{O}$ and $\text{C}-\text{H}\cdots\text{O}$ hydrogen bonds, $\pi-\pi$ contacts between the formylbenzoate rings [centroid-centroid distance = 3.7736 (8) Å] and weak $\text{C}-\text{H}\cdots\pi$ interactions, forming a three-dimensional network.

Related literature

For general background to niacin, see: Krishnamachari (1974). For the crystal structure of *N,N*-diethylnicotinamide, see: Bigoli *et al.* (1972). For related structures, see: Hökelek & Necefoğlu (1996); Hökelek *et al.* (2009).



Experimental

Crystal data

$[\text{Zn}(\text{C}_8\text{H}_5\text{O}_3)_2(\text{C}_6\text{H}_6\text{N}_2\text{O})]$

$M_r = 485.74$

Monoclinic, $P2_1/n$

$a = 13.3143$ (2) Å

$b = 6.7857$ (1) Å

$c = 21.3927$ (3) Å

$\beta = 91.458$ (1)°

$V = 1932.14$ (5) Å³

$Z = 4$

Mo $K\alpha$ radiation

$\mu = 1.32$ mm⁻¹

$T = 100$ K

$0.22 \times 0.12 \times 0.08$ mm

Data collection

Bruker Kappa APEXII CCD area-detector diffractometer

Absorption correction: multi-scan (*SADABS*; Bruker, 2005)

$T_{\text{min}} = 0.829$, $T_{\text{max}} = 0.903$

17841 measured reflections

4812 independent reflections

4086 reflections with $I > 2\sigma(I)$

$R_{\text{int}} = 0.072$

Refinement

$R[F^2 > 2\sigma(F^2)] = 0.028$

$wR(F^2) = 0.080$

$S = 1.10$

4812 reflections

297 parameters

H atoms treated by a mixture of independent and constrained refinement

$\Delta\rho_{\text{max}} = 0.47$ e Å⁻³

$\Delta\rho_{\text{min}} = -0.39$ e Å⁻³

Table 1

Selected bond lengths (Å).

Zn1—O1	1.9153 (11)	Zn1—O4	1.9450 (10)
Zn1—O3	1.9723 (11)	Zn1—N1	2.0270 (12)

Table 2

Hydrogen-bond geometry (Å, °).

$D-\text{H}\cdots A$	$D-\text{H}$	$\text{H}\cdots A$	$D\cdots A$	$D-\text{H}\cdots A$
$\text{N2}-\text{H2A}\cdots\text{O2}^{\text{i}}$	0.86	2.08	2.9242 (17)	165
$\text{N2}-\text{H2B}\cdots\text{O2}^{\text{ii}}$	0.86	2.11	2.9439 (17)	163
$\text{C4}-\text{H4}\cdots\text{O5}^{\text{iii}}$	0.93	2.41	3.298 (2)	160
$\text{C6}-\text{H6}\cdots\text{O7}^{\text{iv}}$	0.93	2.50	3.223 (2)	135
$\text{C15}-\text{H15}\cdots\text{O6}^{\text{iv}}$	0.93	2.32	3.2049 (19)	159
$\text{C16}-\text{H16}\cdots\text{O2}^{\text{ii}}$	0.93	2.44	3.3541 (18)	169
$\text{C3}-\text{H3}\cdots\text{Cg1}$	0.93	2.73	3.6332 (17)	163

Symmetry codes: (i) $-x + 2, -y, -z$; (ii) $x, y - 1, z$; (iii) $x + \frac{1}{2}, -y + \frac{1}{2}, z + \frac{1}{2}$; (iv) $-x + \frac{5}{2}, y - \frac{1}{2}, -z + \frac{1}{2}$. Cg1 is the centroid of the C9–C14 ring.

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: *ORTEP-3 for Windows* (Farrugia, 1997); software used to prepare material for publication: *WinGX* (Farrugia, 1999) and *PLATON* (Spek, 2009).

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Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: CI2939).

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supporting information

Acta Cryst. (2009). E65, m1399–m1400 [https://doi.org/10.1107/S160053680904241X]

catena-Poly[[[(4-formylbenzoato- κ O¹)(isonicotinamide- κ N¹)zinc(II)]- μ -4-formylbenzoato- κ^2 O¹:O¹]

Tuncer Hökelek, Filiz Yılmaz, Barış Tercan, Mustafa Sertçelik and Hacali Necefoğlu

S1. Comment

As a part of our ongoing investigation on transition metal complexes of nicotinamide (NA), one form of niacin (Krishnamachari, 1974), and/or the nicotinic acid derivative *N,N*-diethylnicotinamide (DENA), an important respiratory stimulant (Bigoli *et al.*, 1972), the title compound was synthesized and its crystal structure is reported herein.

In the crystal structure of the title compound, each Zn^{II} ion is coordinated by two formylbenzoate (FB) and one isonicotinamide (INA) ligands (Fig. 1), while symmetry related FB ligands bridge the Zn^{II} ions forming polymeric chains along the *b* axis (Fig.2). The structures of similar complexes of Zn^{II} ion, [Zn₂(C₁₀H₁₄N₂O)₂(C₇H₅O₃)₄].2H₂O (Hökelek & Necefoğlu, 1996) and [Zn(C₉H₁₀NO₂)₂(C₆H₆N₂O)(H₂O)₂] (Hökelek *et al.*, 2009) have also been reported.

The average Zn—O bond length (Table 1) is 1.9442 (11) Å and the Zn1 atom is displaced out of the least-squares planes of the carboxylate groups (O1/C1/O2) and (O3/C8/O4*) by 0.687 (5) Å and 0.703 (2) Å, respectively. The O1/C1/O2 and O3/C8/O4* carboxylate planes form dihedral angles of 9.07 (2)° and 26.2 (2)°, respectively, with benzene rings A(C2-C7) and B(C9-C14), while the angles between rings A, B and C (N1/C15-C19) are A/B = 81.30 (5), A/C = 63.17 (5) and B/C = 46.11 (5)°.

In the crystal structure, N—H \cdots O and C—H \cdots O hydrogen bonds (Table 2) link adjacent chains into a three-dimensional network. In addition, π – π contacts between symmetry related A(C2-C7) formylbenzoate rings at (x, y, z) and (5/2-x, -1/2+y, 1/2-z)/(5/2-x, 1/2+y, 1/2-z) with a centroid-to-centroid distance of 3.7736 (8) Å, and weak C—H \cdots π interaction (Table 2) involving the B(C9-C14) ring stabilize the structure.

S2. Experimental

The title compound was prepared by the reaction of ZnSO₄.H₂O (0.90 g, 5 mmol) in H₂O (25 ml) and INA (1.22 g, 10 mmol) in H₂O (40 ml) with sodium 4-formylbenzoate (1.72 g, 10 mmol) in H₂O (50 ml). The mixture was filtered and set aside to crystallize at ambient temperature for several days, giving colourless single crystals.

S3. Refinement

Atoms H21 and H22 (for methine) were located in a difference Fourier map and refined isotropically. The remaining H atoms were positioned geometrically with N-H = 0.86 Å (for NH₂) and C-H = 0.93 Å for aromatic H atoms and constrained to ride on their parent atoms, with $U_{\text{iso}}(\text{H}) = 1.2U_{\text{eq}}(\text{C,N})$.

catena-Poly[[[4-formylbenzoato- κO^1](isonicotinamide- κN^1)zinc(II)]- μ -4-formylbenzoato- $\kappa^2 O^1:O^1$]

Crystal data

[Zn(C₈H₅O₃)₂(C₆H₆N₂O)] $M_r = 485.74$ Monoclinic, $P2_1/n$

Hall symbol: -P 2yn

 $a = 13.3143$ (2) Å $b = 6.7857$ (1) Å $c = 21.3927$ (3) Å $\beta = 91.458$ (1)° $V = 1932.14$ (5) Å³ $Z = 4$ $F(000) = 992$ $D_x = 1.670$ Mg m⁻³Mo $K\alpha$ radiation, $\lambda = 0.71073$ Å

Cell parameters from 9448 reflections

 $\theta = 3.1$ – 28.4 ° $\mu = 1.32$ mm⁻¹ $T = 100$ K

Needle, colourless

 $0.22 \times 0.12 \times 0.08$ mm

Data collection

Bruker Kappa APEXII CCD area-detector
diffractometer

Radiation source: fine-focus sealed tube

Graphite monochromator

 φ and ω scansAbsorption correction: multi-scan
(SADABS; Bruker, 2005) $T_{\min} = 0.829$, $T_{\max} = 0.903$

17841 measured reflections

4812 independent reflections

4086 reflections with $I > 2\sigma(I)$ $R_{\text{int}} = 0.072$ $\theta_{\max} = 28.4$ °, $\theta_{\min} = 1.8$ ° $h = -17 \rightarrow 14$ $k = -8 \rightarrow 9$ $l = -28 \rightarrow 28$

Refinement

Refinement on F^2

Least-squares matrix: full

 $R[F^2 > 2\sigma(F^2)] = 0.028$ $wR(F^2) = 0.080$ $S = 1.10$

4812 reflections

297 parameters

0 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 $w = 1/[\sigma^2(F_o^2) + (0.0394P)^2 + 0.0092P]$ where $P = (F_o^2 + 2F_c^2)/3$ $(\Delta/\sigma)_{\max} = 0.001$ $\Delta\rho_{\max} = 0.47$ e Å⁻³ $\Delta\rho_{\min} = -0.39$ 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	$U_{\text{iso}}^*/U_{\text{eq}}$
Zn1	0.816160 (13)	0.51300 (2)	0.187880 (8)	0.01058 (7)
O1	0.95052 (8)	0.60767 (16)	0.20123 (5)	0.0149 (2)
O2	1.00643 (8)	0.55706 (16)	0.10502 (5)	0.0149 (2)

O3	0.77863 (8)	0.37932 (15)	0.26575 (5)	0.0134 (2)
O4	0.70741 (8)	0.67550 (14)	0.15404 (5)	0.0137 (2)
O5	0.83582 (11)	-0.07543 (19)	-0.07713 (6)	0.0312 (3)
O6	1.49773 (9)	0.61367 (17)	0.23858 (5)	0.0199 (3)
O7	1.10801 (11)	0.97953 (16)	0.45862 (6)	0.0236 (3)
N1	0.81632 (9)	0.29546 (17)	0.12295 (6)	0.0112 (3)
N2	0.91501 (11)	-0.28715 (18)	-0.01158 (6)	0.0182 (3)
H2A	0.9269	-0.3695	-0.0410	0.022*
H2B	0.9348	-0.3125	0.0262	0.022*
C1	1.02044 (11)	0.5821 (2)	0.16202 (7)	0.0110 (3)
C2	1.12564 (11)	0.5862 (2)	0.18971 (7)	0.0109 (3)
C3	1.14051 (12)	0.5879 (2)	0.25423 (7)	0.0126 (3)
H3	1.0857	0.5867	0.2803	0.015*
C4	1.23802 (12)	0.5913 (2)	0.27978 (7)	0.0131 (3)
H4	1.2487	0.5906	0.3229	0.016*
C5	1.31866 (11)	0.5957 (2)	0.24033 (7)	0.0122 (3)
C6	1.30379 (12)	0.5942 (2)	0.17551 (7)	0.0143 (3)
H6	1.3585	0.5979	0.1494	0.017*
C7	1.20746 (12)	0.5872 (2)	0.15043 (7)	0.0134 (3)
H7	1.1971	0.5832	0.1073	0.016*
C8	0.81930 (11)	0.3293 (2)	0.31722 (7)	0.0115 (3)
C9	0.89783 (12)	0.4527 (2)	0.34954 (7)	0.0114 (3)
C10	0.96360 (12)	0.3649 (2)	0.39317 (7)	0.0149 (3)
H10	0.9603	0.2299	0.4004	0.018*
C11	1.03354 (13)	0.4785 (2)	0.42548 (8)	0.0162 (3)
H11	1.0792	0.4188	0.4531	0.019*
C12	1.03609 (12)	0.6823 (2)	0.41700 (7)	0.0140 (3)
C13	0.97104 (12)	0.7698 (2)	0.37288 (7)	0.0138 (3)
H13	0.9730	0.9053	0.3666	0.017*
C14	0.90409 (12)	0.6556 (2)	0.33872 (7)	0.0122 (3)
H14	0.8627	0.7138	0.3083	0.015*
C15	0.86757 (11)	0.1289 (2)	0.13520 (7)	0.0125 (3)
H15	0.8924	0.1074	0.1757	0.015*
C16	0.88489 (13)	-0.0119 (2)	0.09003 (8)	0.0139 (3)
H16	0.9213	-0.1251	0.1000	0.017*
C17	0.84719 (12)	0.0179 (2)	0.02948 (8)	0.0126 (3)
C18	0.79165 (12)	0.1872 (2)	0.01756 (7)	0.0147 (3)
H18	0.7639	0.2097	-0.0221	0.018*
C19	0.77770 (12)	0.3220 (2)	0.06473 (7)	0.0136 (3)
H19	0.7404	0.4350	0.0561	0.016*
C20	0.86597 (12)	-0.1204 (2)	-0.02441 (7)	0.0157 (3)
C21	1.42128 (13)	0.6036 (2)	0.26818 (8)	0.0161 (3)
H21	1.4266 (13)	0.600 (3)	0.3157 (9)	0.028 (5)*
C22	1.10601 (13)	0.8010 (2)	0.45686 (8)	0.0188 (4)
H22	1.1561 (13)	0.714 (3)	0.4842 (9)	0.020 (5)*

Atomic displacement parameters (\AA^2)

	U^{11}	U^{22}	U^{33}	U^{12}	U^{13}	U^{23}
Zn1	0.01001 (11)	0.01290 (10)	0.00882 (11)	0.00080 (6)	-0.00023 (7)	-0.00111 (6)
O1	0.0092 (6)	0.0225 (6)	0.0129 (5)	-0.0007 (4)	0.0003 (4)	-0.0045 (4)
O2	0.0144 (6)	0.0195 (5)	0.0106 (5)	0.0021 (5)	-0.0019 (5)	-0.0020 (4)
O3	0.0120 (6)	0.0184 (5)	0.0098 (5)	-0.0013 (4)	-0.0021 (4)	0.0016 (4)
O4	0.0134 (6)	0.0141 (5)	0.0135 (5)	0.0030 (4)	-0.0013 (5)	-0.0022 (4)
O5	0.0524 (9)	0.0310 (7)	0.0097 (6)	0.0163 (7)	-0.0092 (6)	-0.0050 (5)
O6	0.0130 (6)	0.0295 (6)	0.0172 (6)	-0.0013 (5)	-0.0005 (5)	0.0005 (5)
O7	0.0279 (8)	0.0248 (6)	0.0181 (7)	-0.0107 (5)	0.0007 (6)	-0.0038 (5)
N1	0.0094 (7)	0.0145 (6)	0.0096 (6)	-0.0011 (5)	-0.0002 (5)	0.0003 (5)
N2	0.0256 (8)	0.0193 (6)	0.0094 (6)	0.0055 (6)	-0.0018 (6)	-0.0037 (5)
C1	0.0111 (8)	0.0103 (6)	0.0115 (7)	0.0008 (6)	-0.0008 (6)	-0.0005 (6)
C2	0.0111 (8)	0.0102 (6)	0.0113 (7)	0.0000 (6)	-0.0016 (6)	0.0001 (6)
C3	0.0128 (8)	0.0145 (7)	0.0106 (7)	0.0009 (6)	0.0017 (6)	0.0001 (6)
C4	0.0164 (8)	0.0143 (7)	0.0085 (7)	0.0000 (6)	-0.0008 (6)	-0.0002 (6)
C5	0.0121 (8)	0.0110 (7)	0.0135 (8)	0.0003 (6)	-0.0016 (6)	0.0002 (6)
C6	0.0113 (8)	0.0190 (7)	0.0127 (7)	-0.0008 (6)	0.0028 (6)	-0.0003 (6)
C7	0.0143 (8)	0.0183 (7)	0.0076 (7)	0.0006 (6)	-0.0008 (6)	0.0006 (6)
C8	0.0096 (8)	0.0138 (7)	0.0113 (7)	0.0029 (6)	0.0015 (6)	-0.0023 (6)
C9	0.0098 (8)	0.0154 (7)	0.0091 (7)	-0.0002 (6)	0.0015 (6)	-0.0020 (6)
C10	0.0173 (9)	0.0137 (7)	0.0137 (8)	0.0002 (6)	-0.0001 (7)	-0.0003 (6)
C11	0.0144 (9)	0.0206 (8)	0.0134 (8)	0.0021 (6)	-0.0043 (7)	-0.0003 (6)
C12	0.0113 (8)	0.0191 (7)	0.0117 (7)	-0.0021 (6)	-0.0004 (6)	-0.0029 (6)
C13	0.0149 (8)	0.0136 (7)	0.0129 (7)	-0.0015 (6)	0.0041 (6)	-0.0024 (6)
C14	0.0111 (8)	0.0164 (7)	0.0090 (7)	0.0020 (6)	0.0012 (6)	0.0004 (6)
C15	0.0116 (8)	0.0171 (7)	0.0087 (7)	0.0004 (6)	-0.0023 (6)	0.0008 (6)
C16	0.0157 (9)	0.0140 (7)	0.0120 (8)	0.0026 (6)	-0.0017 (7)	-0.0003 (5)
C17	0.0147 (8)	0.0143 (7)	0.0088 (8)	-0.0022 (6)	0.0006 (6)	-0.0003 (5)
C18	0.0169 (9)	0.0176 (7)	0.0094 (7)	-0.0006 (6)	-0.0029 (6)	0.0014 (6)
C19	0.0130 (8)	0.0150 (7)	0.0128 (7)	0.0019 (6)	-0.0011 (6)	0.0029 (6)
C20	0.0180 (9)	0.0185 (7)	0.0106 (7)	0.0001 (6)	-0.0013 (7)	-0.0024 (6)
C21	0.0173 (9)	0.0171 (7)	0.0139 (8)	0.0003 (6)	-0.0024 (7)	-0.0003 (6)
C22	0.0174 (9)	0.0260 (8)	0.0131 (8)	-0.0046 (7)	0.0007 (7)	-0.0018 (7)

Geometric parameters (\AA , $^\circ$)

Zn1—O1	1.9153 (11)	C7—C6	1.378 (2)
Zn1—O3	1.9723 (11)	C7—H7	0.93
Zn1—O4	1.9450 (10)	C8—O4 ⁱⁱ	1.2669 (18)
Zn1—N1	2.0270 (12)	C8—C9	1.495 (2)
O1—C1	1.2807 (18)	C9—C10	1.397 (2)
O2—C1	1.2405 (18)	C9—C14	1.3990 (19)
O3—C8	1.2611 (17)	C10—H10	0.93
O4—C8 ⁱ	1.2669 (18)	C11—C10	1.381 (2)
O5—C20	1.2261 (18)	C11—H11	0.93
O6—C21	1.214 (2)	C12—C11	1.395 (2)

O7—C22	1.2124 (19)	C12—C13	1.397 (2)
N1—C15	1.3427 (18)	C12—C22	1.484 (2)
N1—C19	1.3474 (18)	C13—C14	1.377 (2)
N2—C20	1.332 (2)	C13—H13	0.93
N2—H2A	0.86	C14—H14	0.93
N2—H2B	0.86	C15—H15	0.93
C2—C1	1.507 (2)	C16—C15	1.382 (2)
C2—C3	1.390 (2)	C16—C17	1.392 (2)
C2—C7	1.393 (2)	C16—H16	0.93
C3—C4	1.396 (2)	C18—C17	1.386 (2)
C3—H3	0.93	C18—H18	0.93
C4—H4	0.93	C19—C18	1.378 (2)
C5—C4	1.383 (2)	C19—H19	0.93
C5—C6	1.396 (2)	C20—C17	1.512 (2)
C5—C21	1.478 (2)	C21—H21	1.018 (19)
C6—H6	0.93	C22—H22	1.057 (17)
O1—Zn1—O3	106.50 (4)	C14—C9—C8	121.23 (13)
O1—Zn1—O4	123.30 (5)	C9—C10—H10	120.0
O1—Zn1—N1	109.22 (5)	C11—C10—C9	119.95 (14)
O3—Zn1—N1	104.39 (5)	C11—C10—H10	120.0
O4—Zn1—O3	111.89 (5)	C10—C11—C12	120.45 (15)
O4—Zn1—N1	99.88 (5)	C10—C11—H11	119.8
C1—O1—Zn1	123.27 (10)	C12—C11—H11	119.8
C8—O3—Zn1	138.48 (10)	C11—C12—C13	119.50 (14)
C8 ⁱ —O4—Zn1	120.13 (10)	C11—C12—C22	118.70 (15)
C15—N1—Zn1	119.26 (10)	C13—C12—C22	121.77 (14)
C15—N1—C19	118.22 (13)	C12—C13—H13	120.0
C19—N1—Zn1	121.95 (10)	C14—C13—C12	120.09 (14)
C20—N2—H2A	120.0	C14—C13—H13	120.0
C20—N2—H2B	120.0	C9—C14—H14	119.8
H2A—N2—H2B	120.0	C13—C14—C9	120.42 (14)
O1—C1—C2	115.07 (13)	C13—C14—H14	119.8
O2—C1—O1	124.67 (14)	N1—C15—C16	122.57 (14)
O2—C1—C2	120.25 (14)	N1—C15—H15	118.7
C3—C2—C1	119.87 (14)	C16—C15—H15	118.7
C3—C2—C7	120.36 (14)	C15—C16—C17	119.19 (14)
C7—C2—C1	119.76 (13)	C15—C16—H16	120.4
C2—C3—C4	119.77 (14)	C17—C16—H16	120.4
C2—C3—H3	120.1	C16—C17—C20	123.82 (13)
C4—C3—H3	120.1	C18—C17—C16	118.00 (14)
C3—C4—H4	120.3	C18—C17—C20	118.16 (14)
C5—C4—C3	119.36 (14)	C17—C18—H18	120.1
C5—C4—H4	120.3	C19—C18—C17	119.77 (14)
C4—C5—C6	120.90 (14)	C19—C18—H18	120.1
C4—C5—C21	118.62 (14)	N1—C19—C18	122.20 (14)
C6—C5—C21	120.48 (14)	N1—C19—H19	118.9
C5—C6—H6	120.2	C18—C19—H19	118.9

C7—C6—C5	119.60 (14)	O5—C20—N2	123.26 (15)
C7—C6—H6	120.2	O5—C20—C17	119.35 (14)
C2—C7—H7	120.0	N2—C20—C17	117.39 (13)
C6—C7—C2	119.98 (14)	O6—C21—C5	124.78 (15)
C6—C7—H7	120.0	O6—C21—H21	119.0 (11)
O3—C8—O4 ⁱⁱ	121.72 (14)	C5—C21—H21	116.3 (11)
O3—C8—C9	122.19 (13)	O7—C22—C12	124.98 (17)
O4 ⁱⁱ —C8—C9	116.01 (13)	O7—C22—H22	121.9 (9)
C10—C9—C8	119.27 (13)	C12—C22—H22	113.1 (9)
C10—C9—C14	119.45 (14)		
O3—Zn1—O1—C1	132.27 (11)	C21—C5—C4—C3	178.69 (13)
O4—Zn1—O1—C1	-96.44 (12)	C4—C5—C6—C7	-0.3 (2)
N1—Zn1—O1—C1	20.07 (13)	C21—C5—C6—C7	-179.85 (14)
O1—Zn1—O3—C8	-6.56 (15)	C4—C5—C21—O6	-177.81 (15)
O4—Zn1—O3—C8	-143.97 (14)	C6—C5—C21—O6	1.7 (2)
N1—Zn1—O3—C8	108.94 (15)	C2—C7—C6—C5	1.5 (2)
O1—Zn1—O4—C8 ⁱ	-60.90 (13)	O3—C8—C9—C10	-158.23 (15)
O3—Zn1—O4—C8 ⁱ	68.17 (12)	O3—C8—C9—C14	24.3 (2)
N1—Zn1—O4—C8 ⁱ	178.16 (11)	O4 ⁱⁱ —C8—C9—C10	24.8 (2)
O1—Zn1—N1—C15	62.86 (12)	O4 ⁱⁱ —C8—C9—C14	-152.61 (15)
O1—Zn1—N1—C19	-108.32 (12)	C8—C9—C10—C11	-177.06 (15)
O3—Zn1—N1—C15	-50.72 (12)	C14—C9—C10—C11	0.4 (2)
O3—Zn1—N1—C19	138.10 (12)	C8—C9—C14—C13	174.28 (14)
O4—Zn1—N1—C15	-166.53 (11)	C10—C9—C14—C13	-3.2 (2)
O4—Zn1—N1—C19	22.29 (13)	C12—C11—C10—C9	2.9 (3)
Zn1—O1—C1—O2	25.4 (2)	C13—C12—C11—C10	-3.6 (3)
Zn1—O1—C1—C2	-155.21 (9)	C22—C12—C11—C10	174.41 (16)
Zn1—O3—C8—O4 ⁱⁱ	-147.48 (12)	C11—C12—C13—C14	0.9 (2)
Zn1—O3—C8—C9	35.8 (2)	C22—C12—C13—C14	-177.08 (15)
Zn1—N1—C15—C16	-169.24 (12)	C11—C12—C22—O7	-171.97 (17)
C19—N1—C15—C16	2.3 (2)	C13—C12—C22—O7	6.0 (3)
Zn1—N1—C19—C18	169.38 (12)	C12—C13—C14—C9	2.5 (2)
C15—N1—C19—C18	-1.9 (2)	C17—C16—C15—N1	-0.7 (2)
C3—C2—C1—O1	9.2 (2)	C15—C16—C17—C18	-1.4 (2)
C3—C2—C1—O2	-171.42 (14)	C15—C16—C17—C20	176.70 (15)
C7—C2—C1—O1	-171.18 (13)	C19—C18—C17—C16	1.7 (2)
C7—C2—C1—O2	8.2 (2)	C19—C18—C17—C20	-176.45 (14)
C1—C2—C3—C4	179.87 (13)	N1—C19—C18—C17	-0.1 (2)
C7—C2—C3—C4	0.2 (2)	N2—C20—C17—C16	5.3 (2)
C1—C2—C7—C6	178.94 (14)	N2—C20—C17—C18	-176.65 (15)
C3—C2—C7—C6	-1.4 (2)	O5—C20—C17—C16	-174.85 (17)
C2—C3—C4—C5	0.9 (2)	O5—C20—C17—C18	3.2 (2)
C6—C5—C4—C3	-0.8 (2)		

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

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>
N2—H2 <i>A</i> ···O2 ⁱⁱⁱ	0.86	2.08	2.9242 (17)	165
N2—H2 <i>B</i> ···O2 ^{iv}	0.86	2.11	2.9439 (17)	163
C4—H4···O5 ^v	0.93	2.41	3.298 (2)	160
C6—H6···O7 ^{vi}	0.93	2.50	3.223 (2)	135
C15—H15···O6 ^{vi}	0.93	2.32	3.2049 (19)	159
C16—H16···O2 ^{iv}	0.93	2.44	3.3541 (18)	169
C3—H3···C <i>g</i> 1	0.93	2.73	3.6332 (17)	163

Symmetry codes: (iii) $-x+2, -y, -z$; (iv) $x, y-1, z$; (v) $x+1/2, -y+1/2, z+1/2$; (vi) $-x+5/2, y-1/2, -z+1/2$.