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A four-armed Schiff base: 6,6',6'',6'''-tetramethoxy-2,2',2'',2'''-[methane-tetrayltetrakis(methylenenitrilomethylidene)]tetraphenol

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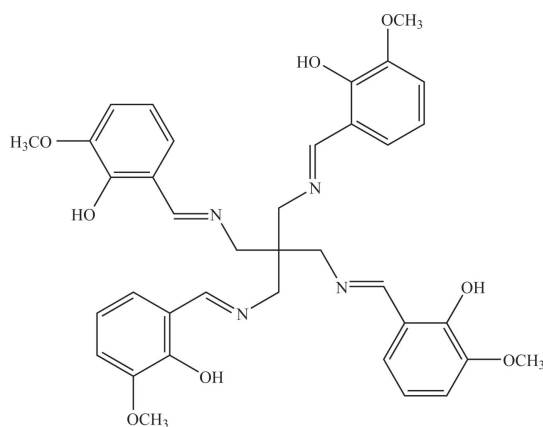
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 Key indicators: single-crystal X-ray study; $T = 273$ K; mean $\sigma(\text{C}-\text{C}) = 0.004$ Å; R factor = 0.054; wR factor = 0.146; data-to-parameter ratio = 14.0.

In the structure of the title compound, $\text{C}_{37}\text{H}_{40}\text{N}_4\text{O}_8$, pentaerythrityltetramine is bonded to four *o*-vanillin molecules, forming a four-armed Schiff base molecule. These molecules are connected by intermolecular $\text{C}-\text{H}\cdots\text{O}$ hydrogen bonds. Intramolecular $\text{C}-\text{H}\cdots\text{N}$ and $\text{O}-\text{H}\cdots\text{N}$ hydrogen bonds are also present.

Related literature

 For related literature, see: Tanaka *et al.* (2006); Bernardo *et al.* (1996); Laliberte *et al.* (2004).


Experimental

Crystal data

 $\text{C}_{37}\text{H}_{40}\text{N}_4\text{O}_8$
 $M_r = 668.73$

 Triclinic, $P\bar{1}$
 $a = 11.3464$ (11) Å
 $b = 12.4437$ (12) Å
 $c = 13.0523$ (14) Å
 $\alpha = 75.861$ (7)°
 $\beta = 88.893$ (7)°
 $\gamma = 78.385$ (7)°

 $V = 1749.6$ (3) Å³
 $Z = 2$
 Mo $K\alpha$ radiation
 $\mu = 0.09$ mm⁻¹
 $T = 273$ (2) K
 $0.18 \times 0.16 \times 0.15$ mm

Data collection

 Bruker APEXII CCD area-detector diffractometer
 Absorption correction: multi-scan (SADABS; Bruker, 2007)
 $T_{\min} = 0.984$, $T_{\max} = 0.987$

 16510 measured reflections
 6221 independent reflections
 3077 reflections with $I > 2\sigma(I)$
 $R_{\text{int}} = 0.058$

Refinement

 $R[F^2 > 2\sigma(F^2)] = 0.054$
 $wR(F^2) = 0.145$
 $S = 0.94$
 6194 reflections

 443 parameters
 H-atom parameters constrained
 $\Delta\rho_{\text{max}} = 0.18$ e Å⁻³
 $\Delta\rho_{\text{min}} = -0.19$ e Å⁻³

Table 1

Hydrogen-bond geometry (Å, °).

$D-\text{H}\cdots A$	$D-\text{H}$	$\text{H}\cdots A$	$D\cdots A$	$D-\text{H}\cdots A$
$\text{O}2-\text{H}2\cdots\text{N}1$	0.82	1.87	2.592 (3)	147
$\text{O}4-\text{H}4\cdots\text{N}4$	0.82	1.84	2.567 (2)	147
$\text{O}6-\text{H}6\cdots\text{N}3$	0.82	1.89	2.623 (3)	148
$\text{O}7-\text{H}7\cdots\text{N}2$	0.82	1.88	2.612 (3)	148
$\text{C}27-\text{H}27\text{B}\cdots\text{N}1$	0.97	2.56	2.936 (3)	103
$\text{C}26-\text{H}26\cdots\text{O}5^i$	0.93	2.58	3.309 (3)	136
$\text{C}28-\text{H}28\text{C}\cdots\text{O}2^i$	0.96	2.57	3.387 (3)	143
$\text{C}12-\text{H}12\cdots\text{O}6^{\text{ii}}$	0.93	2.59	3.400 (3)	146
$\text{C}32-\text{H}32\cdots\text{O}3^{\text{iii}}$	0.93	2.42	3.289 (3)	155
$\text{C}36-\text{H}36\text{B}\cdots\text{O}8^{\text{iv}}$	0.97	2.55	3.399 (3)	146

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

Data collection: APEX2 (Bruker, 2007); cell refinement: SAINT (Bruker, 2007); data reduction: SAINT (Bruker, 2007); 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).

Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: OM2247).

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supporting information

Acta Cryst. (2008). E64, o1455 [doi:10.1107/S1600536808020618]

A four-armed Schiff base: 6,6',6'',6'''-tetramethoxy-2,2',2'',2'''-[methane-tetrayltetrakis(methylenenitrilomethylidyne)]tetraphenol

Guang-Qi Jiang, Jie Cai, Yun-Qian Zhang and Qian-Jun Zhang

S1. Comment

Schiff bases have been intensively investigated during the past decades, mainly because of their strong coordination capability and diverse biological activities, such as antibacterial, antitumor activities, *etc.* (Tanaka *et al.*, 2006; Bernardo *et al.*, 1996). Pentaerythrityl tetramine and its congeners act as versatile building blocks in supramolecular chemistry and have been widely employed for the synthesis of various spiro crown ethers, spiro-macrocyclic ligands *etc.* (Laliberte *et al.*, 2004). However, to the best of our knowledge, few Schiff bases and their complexes derived from pentaerythrityl tetramine have been reported. Herein, we present the synthesis and crystal structure of the multidentate Schiff base ligand derived from *o*-vanillin and pentaerythrityl tetramine.

The molecular structure is shown in Fig. 1. The atoms C9, C18, C27, and C36 form a distorted tetrahedral geometry about the central atom C37. The structure reveals four intramolecular O—H \cdots N hydrogen bonds and one C—H \cdots N hydrogen bond as well as five intermolecular C—H \cdots O hydrogen bonds (Table 1). A view of the packing is given in Fig. 2. The molecules are also connected by C—H \cdots π interactions, with C9—H9B \cdots Cg(1) angle of 138°, C9 \cdots Cg(1) distance of 3.595 (3) Å and C14—H14 \cdots Cg(4) angle of 157°, C14 \cdots Cg(4) distance of 3.614 (3) Å, where Cg(1) and Cg(4) are the centroids of the phenyl ring formed by atoms C2, C3, C4, C5, C6, C7 and C29, C30, C31, C32, C33, C34 at (1 - x, -y, 1 - z) and (1 + x, y, z), respectively.

S2. Experimental

A solution of pentaerythrityl tetramine (0.13 g, 1 mmol) in ethanol (20 ml) was added dropwise while stirring to a solution of *o*-vanillin (0.68 g, 4 mmol) in ethanol (20 ml), the mixture was stirred for about 2 h at room temperature. The resulting solution was filtered, and after slow evaporation of the solvent for several days, crystals suitable for X-ray diffraction were obtained in a yield of 46.4%.

S3. Refinement

All H atoms were placed in calculated positions and refined as riding, with C—H = 0.93–0.97 Å, and with $U_{\text{iso}}(\text{H}) = 1.2\text{--}1.5 U_{\text{eq}}(\text{C}, \text{O})$.

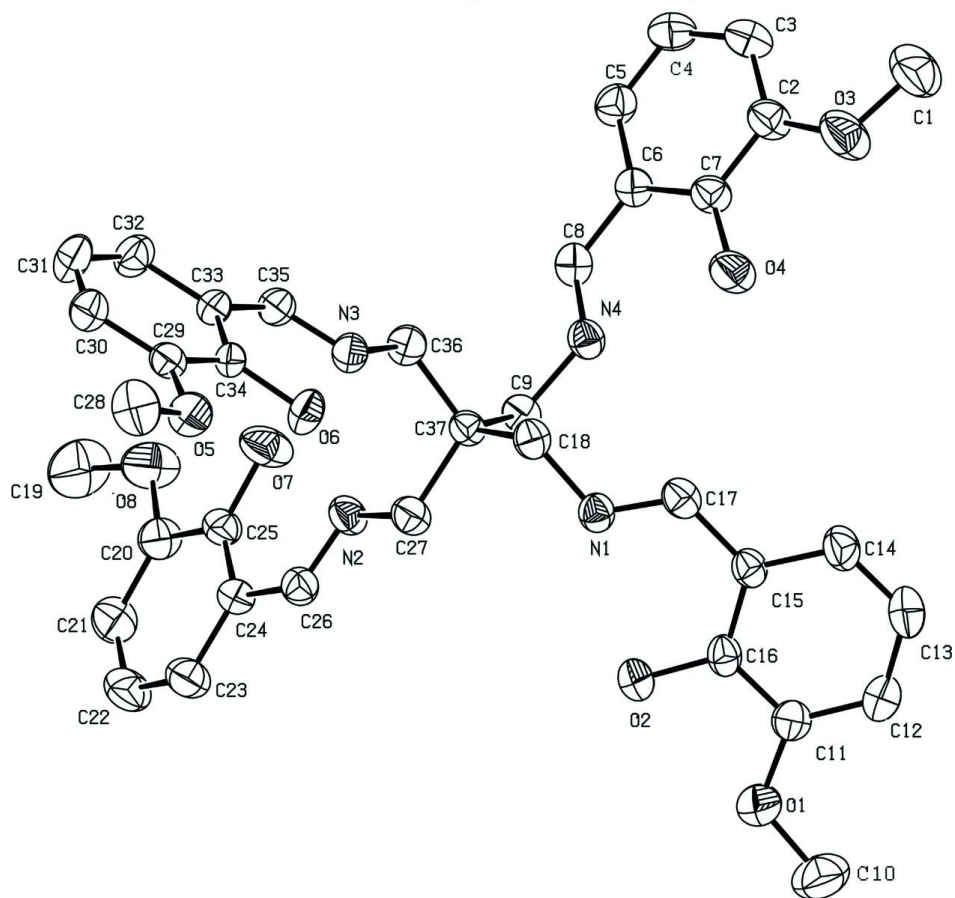


Figure 1

The molecular structure shown with 50% probability displacement ellipsoids.

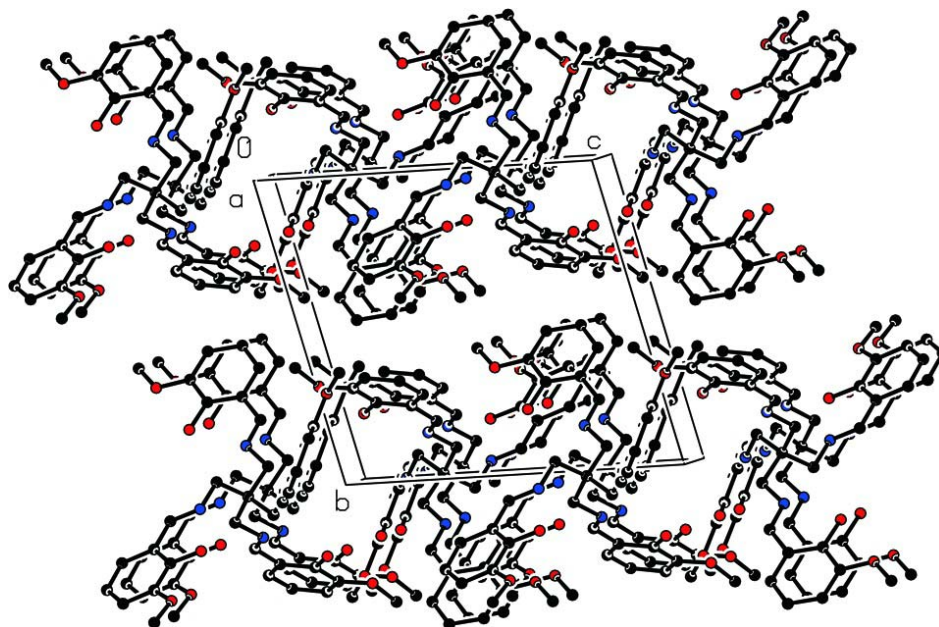


Figure 2

One-dimensional chains of the Schiff base perpendicular to the *bc* plane, showing the packing of the compound. Hydrogen bonds and C—H... π interactions are shown as dashed lines. H atoms have been omitted for clarity.

6,6',6'',6'''-tetramethoxy-2,2',2'',2'''-[methanetetrayl]tetrakis(methylenenitrilomethylidyne)]tetraphenol

Crystal data

$C_{37}H_{40}N_4O_8$

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Triclinic, $P\bar{1}$

Hall symbol: $-P\ 1$

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$c = 13.0523$ (14) Å

$\alpha = 75.861$ (7)°

$\beta = 88.893$ (7)°

$\gamma = 78.385$ (7)°

$V = 1749.6$ (3) Å³

$Z = 2$

$F(000) = 708$

$D_x = 1.269$ Mg m⁻³

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

Cell parameters from 16510 reflections

$\theta = 1.0$ – 25.1 °

$\mu = 0.09$ mm⁻¹

$T = 273$ K

Prism, yellow

$0.18 \times 0.16 \times 0.15$ mm

Data collection

Bruker APEXII CCD area-detector
diffractometer

Radiation source: fine-focus sealed tube

Graphite monochromator

φ and ω scans

Absorption correction: multi-scan
(*SADABS*; Bruker, 2007)

$T_{\min} = 0.984$, $T_{\max} = 0.987$

16510 measured reflections

6221 independent reflections

3077 reflections with $I > 2\sigma(I)$

$R_{\text{int}} = 0.058$

$\theta_{\max} = 25.1$ °, $\theta_{\min} = 1.7$ °

$h = -13 \rightarrow 13$

$k = -14 \rightarrow 14$

$l = -13 \rightarrow 15$

*Refinement*Refinement on F^2

Least-squares matrix: full

 $R[F^2 > 2\sigma(F^2)] = 0.054$ $wR(F^2) = 0.145$ $S = 0.94$

6194 reflections

443 parameters

0 restraints

Primary atom site location: structure-invariant
direct methodsSecondary atom site location: difference Fourier
mapHydrogen site location: inferred from
neighbouring sites

H-atom parameters constrained

 $w = 1/[\sigma^2(F_o^2) + (0.0583P)^2]$ where $P = (F_o^2 + 2F_c^2)/3$ $(\Delta/\sigma)_{\max} < 0.001$ $\Delta\rho_{\max} = 0.18 \text{ e } \text{\AA}^{-3}$ $\Delta\rho_{\min} = -0.19 \text{ e } \text{\AA}^{-3}$ Extinction correction: *SHELXL97* (Sheldrick,
2008), $F_c^* = kF_c[1 + 0.001x F_c^2 \lambda^3 / \sin(2\theta)]^{-1/4}$

Extinction coefficient: 0.0093 (14)

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 (\AA^2)

	<i>x</i>	<i>y</i>	<i>z</i>	$U_{\text{iso}}^*/U_{\text{eq}}$
C1	-0.6783 (3)	0.6040 (3)	0.6346 (3)	0.1205 (14)
H1A	-0.7350	0.5856	0.5914	0.181*
H1B	-0.7208	0.6457	0.6816	0.181*
H1C	-0.6285	0.5355	0.6752	0.181*
C2	-0.5190 (3)	0.7057 (2)	0.6164 (2)	0.0661 (8)
C3	-0.4819 (3)	0.6681 (2)	0.7207 (2)	0.0689 (8)
H3	-0.5142	0.6112	0.7650	0.083*
C4	-0.3970 (3)	0.7138 (2)	0.7606 (2)	0.0705 (8)
H4A	-0.3729	0.6879	0.8315	0.085*
C5	-0.3483 (2)	0.7971 (2)	0.6962 (2)	0.0637 (7)
H5	-0.2922	0.8283	0.7239	0.076*
C6	-0.3823 (2)	0.8355 (2)	0.5891 (2)	0.0504 (6)
C7	-0.4681 (2)	0.7891 (2)	0.5489 (2)	0.0561 (7)
C8	-0.3299 (2)	0.9225 (2)	0.5203 (2)	0.0556 (7)
H8	-0.2759	0.9550	0.5492	0.067*
C9	-0.2957 (2)	1.0390 (2)	0.35387 (19)	0.0541 (7)
H9A	-0.2517	1.0717	0.3972	0.065*
H9B	-0.3559	1.0994	0.3116	0.065*
C10	-0.7806 (3)	1.3762 (3)	-0.1285 (3)	0.1135 (13)
H10A	-0.7593	1.4442	-0.1701	0.170*
H10B	-0.8450	1.3953	-0.0832	0.170*
H10C	-0.8063	1.3352	-0.1746	0.170*
C11	-0.6932 (2)	1.2070 (2)	-0.0018 (2)	0.0585 (7)

C12	-0.7994 (2)	1.1700 (3)	0.0167 (2)	0.0672 (8)
H12	-0.8709	1.2161	-0.0145	0.081*
C13	-0.8005 (3)	1.0645 (3)	0.0816 (3)	0.0773 (9)
H13	-0.8725	1.0393	0.0929	0.093*
C14	-0.6962 (2)	0.9971 (2)	0.1292 (2)	0.0680 (8)
H14	-0.6979	0.9259	0.1725	0.082*
C15	-0.5875 (2)	1.0334 (2)	0.11382 (19)	0.0511 (7)
C16	-0.5865 (2)	1.1396 (2)	0.0472 (2)	0.0527 (7)
C17	-0.4787 (2)	0.9634 (2)	0.16885 (19)	0.0567 (7)
H17	-0.4822	0.8927	0.2121	0.068*
C18	-0.2728 (2)	0.9217 (2)	0.21916 (19)	0.0540 (7)
H18A	-0.2174	0.8917	0.1707	0.065*
H18B	-0.2971	0.8584	0.2683	0.065*
C19	0.1639 (4)	1.4133 (3)	0.5082 (3)	0.1264 (14)
H19A	0.1696	1.3906	0.5840	0.190*
H19B	0.1277	1.4923	0.4858	0.190*
H19C	0.2429	1.4004	0.4803	0.190*
C20	0.0741 (2)	1.3692 (2)	0.3635 (2)	0.0667 (8)
C21	0.1075 (3)	1.4562 (2)	0.2895 (3)	0.0770 (9)
H21	0.1441	1.5075	0.3116	0.092*
C22	0.0876 (3)	1.4686 (2)	0.1828 (3)	0.0822 (9)
H22	0.1117	1.5272	0.1334	0.099*
C23	0.0320 (2)	1.3938 (2)	0.1502 (2)	0.0727 (8)
H23	0.0183	1.4023	0.0784	0.087*
C24	-0.0044 (2)	1.3052 (2)	0.2232 (2)	0.0509 (7)
C25	0.0163 (2)	1.2936 (2)	0.3301 (2)	0.0581 (7)
C26	-0.0634 (2)	1.2277 (2)	0.1863 (2)	0.0547 (7)
H26	-0.0732	1.2373	0.1138	0.066*
C27	-0.1624 (2)	1.0776 (2)	0.19919 (19)	0.0591 (7)
H27A	-0.1064	1.0425	0.1541	0.071*
H27B	-0.2297	1.1260	0.1550	0.071*
C28	0.3376 (3)	0.6069 (3)	-0.0230 (2)	0.0883 (10)
H28A	0.3091	0.6053	-0.0912	0.132*
H28B	0.3572	0.5312	0.0209	0.132*
H28C	0.4082	0.6394	-0.0308	0.132*
C29	0.2713 (2)	0.6854 (2)	0.1230 (2)	0.0519 (7)
C30	0.3835 (2)	0.6521 (2)	0.1727 (2)	0.0666 (8)
H30	0.4489	0.6186	0.1394	0.080*
C31	0.3984 (3)	0.6687 (3)	0.2720 (3)	0.0782 (9)
H31	0.4743	0.6462	0.3052	0.094*
C32	0.3036 (2)	0.7176 (2)	0.3220 (2)	0.0675 (8)
H32	0.3149	0.7272	0.3894	0.081*
C33	0.1896 (2)	0.7534 (2)	0.27244 (19)	0.0493 (6)
C34	0.1742 (2)	0.7371 (2)	0.1721 (2)	0.0472 (6)
C35	0.0911 (2)	0.8099 (2)	0.3253 (2)	0.0543 (7)
H35	0.1062	0.8195	0.3919	0.065*
C36	-0.1053 (2)	0.9013 (2)	0.3478 (2)	0.0583 (7)
H36A	-0.0667	0.9410	0.3885	0.070*

H36B	-0.1378	0.8435	0.3973	0.070*
C37	-0.2085 (2)	0.9849 (2)	0.28059 (18)	0.0475 (6)
N1	-0.37817 (18)	0.99528 (17)	0.16006 (15)	0.0527 (5)
N2	-0.10224 (18)	1.14746 (18)	0.24823 (16)	0.0586 (6)
N3	-0.01463 (19)	0.84665 (17)	0.28493 (15)	0.0544 (6)
N4	-0.35536 (17)	0.95591 (17)	0.42188 (17)	0.0542 (6)
O1	-0.68132 (17)	1.30936 (18)	-0.06743 (17)	0.0877 (6)
O2	-0.48359 (14)	1.17979 (14)	0.02791 (14)	0.0690 (5)
H2	-0.4272	1.1331	0.0611	0.104*
O3	-0.6060 (2)	0.6698 (2)	0.56997 (16)	0.1045 (8)
O4	-0.50509 (17)	0.82276 (16)	0.44657 (14)	0.0791 (6)
H4	-0.4688	0.8710	0.4150	0.119*
O5	0.24695 (16)	0.67307 (15)	0.02443 (14)	0.0692 (5)
O6	0.06588 (14)	0.77182 (15)	0.12026 (13)	0.0641 (5)
H6	0.0164	0.8014	0.1570	0.096*
O7	-0.01615 (18)	1.20882 (16)	0.40428 (14)	0.0870 (7)
H7	-0.0487	1.1704	0.3754	0.131*
O8	0.0914 (2)	1.34865 (17)	0.47009 (17)	0.1004 (7)

Atomic displacement parameters (Å²)

	U^{11}	U^{22}	U^{33}	U^{12}	U^{13}	U^{23}
C1	0.130 (3)	0.122 (3)	0.118 (3)	-0.077 (3)	0.002 (3)	-0.005 (3)
C2	0.080 (2)	0.0671 (19)	0.0570 (19)	-0.0297 (17)	0.0043 (16)	-0.0143 (16)
C3	0.091 (2)	0.0606 (19)	0.0545 (19)	-0.0213 (17)	0.0114 (17)	-0.0093 (16)
C4	0.087 (2)	0.071 (2)	0.0492 (17)	-0.0113 (17)	0.0006 (16)	-0.0090 (16)
C5	0.0650 (18)	0.076 (2)	0.0539 (18)	-0.0153 (16)	-0.0037 (14)	-0.0208 (16)
C6	0.0527 (15)	0.0560 (16)	0.0465 (16)	-0.0115 (13)	0.0051 (13)	-0.0200 (14)
C7	0.0656 (17)	0.0617 (18)	0.0441 (16)	-0.0180 (15)	0.0041 (14)	-0.0149 (14)
C8	0.0527 (16)	0.0623 (18)	0.0592 (18)	-0.0135 (14)	0.0050 (14)	-0.0273 (15)
C9	0.0607 (16)	0.0543 (16)	0.0512 (16)	-0.0170 (14)	0.0047 (14)	-0.0160 (14)
C10	0.117 (3)	0.082 (2)	0.121 (3)	-0.005 (2)	-0.033 (3)	0.005 (2)
C11	0.0613 (18)	0.0593 (19)	0.0583 (18)	-0.0161 (15)	-0.0014 (15)	-0.0173 (15)
C12	0.0532 (17)	0.076 (2)	0.075 (2)	-0.0103 (15)	-0.0054 (15)	-0.0256 (18)
C13	0.0562 (19)	0.088 (2)	0.097 (2)	-0.0291 (18)	0.0082 (18)	-0.030 (2)
C14	0.0591 (18)	0.0647 (19)	0.086 (2)	-0.0271 (16)	0.0036 (16)	-0.0172 (17)
C15	0.0527 (16)	0.0530 (16)	0.0543 (16)	-0.0208 (13)	0.0028 (13)	-0.0176 (14)
C16	0.0481 (16)	0.0660 (18)	0.0525 (16)	-0.0204 (14)	0.0039 (13)	-0.0236 (14)
C17	0.0643 (18)	0.0588 (17)	0.0540 (17)	-0.0225 (15)	0.0071 (14)	-0.0192 (14)
C18	0.0551 (16)	0.0551 (16)	0.0571 (16)	-0.0163 (14)	0.0036 (13)	-0.0197 (14)
C19	0.164 (4)	0.111 (3)	0.124 (3)	-0.049 (3)	-0.046 (3)	-0.046 (3)
C20	0.076 (2)	0.0602 (19)	0.067 (2)	-0.0202 (16)	-0.0088 (16)	-0.0167 (17)
C21	0.084 (2)	0.0555 (19)	0.097 (3)	-0.0239 (17)	-0.0023 (19)	-0.0208 (19)
C22	0.096 (2)	0.064 (2)	0.088 (3)	-0.0369 (18)	0.005 (2)	-0.0059 (19)
C23	0.084 (2)	0.070 (2)	0.0649 (19)	-0.0292 (17)	0.0038 (16)	-0.0058 (17)
C24	0.0545 (15)	0.0473 (16)	0.0500 (17)	-0.0126 (13)	0.0033 (13)	-0.0084 (13)
C25	0.0637 (17)	0.0510 (17)	0.0593 (18)	-0.0201 (14)	-0.0005 (14)	-0.0060 (15)
C26	0.0529 (16)	0.0597 (17)	0.0521 (16)	-0.0138 (14)	0.0036 (13)	-0.0132 (14)

C27	0.0687 (17)	0.0664 (17)	0.0486 (16)	-0.0295 (15)	-0.0004 (13)	-0.0130 (14)
C28	0.087 (2)	0.094 (2)	0.097 (2)	-0.0184 (19)	0.023 (2)	-0.050 (2)
C29	0.0546 (17)	0.0526 (16)	0.0486 (16)	-0.0155 (13)	0.0055 (14)	-0.0091 (13)
C30	0.0521 (18)	0.0663 (19)	0.076 (2)	-0.0059 (14)	0.0028 (16)	-0.0117 (17)
C31	0.0482 (18)	0.095 (2)	0.087 (2)	-0.0025 (16)	-0.0133 (17)	-0.021 (2)
C32	0.0580 (18)	0.084 (2)	0.0589 (18)	-0.0105 (16)	-0.0107 (15)	-0.0164 (16)
C33	0.0467 (15)	0.0541 (16)	0.0467 (15)	-0.0128 (13)	-0.0049 (13)	-0.0088 (13)
C34	0.0379 (14)	0.0488 (15)	0.0529 (16)	-0.0134 (12)	-0.0039 (13)	-0.0048 (13)
C35	0.0574 (17)	0.0608 (17)	0.0476 (16)	-0.0183 (14)	-0.0043 (14)	-0.0130 (14)
C36	0.0566 (16)	0.0681 (18)	0.0516 (16)	-0.0115 (14)	0.0032 (14)	-0.0183 (14)
C37	0.0524 (15)	0.0493 (15)	0.0433 (15)	-0.0156 (13)	0.0027 (13)	-0.0120 (13)
N1	0.0511 (13)	0.0596 (14)	0.0513 (13)	-0.0157 (11)	0.0025 (11)	-0.0175 (11)
N2	0.0651 (14)	0.0654 (15)	0.0520 (13)	-0.0311 (12)	0.0027 (11)	-0.0123 (12)
N3	0.0477 (13)	0.0636 (14)	0.0534 (13)	-0.0122 (11)	0.0016 (11)	-0.0162 (11)
N4	0.0533 (13)	0.0612 (14)	0.0519 (14)	-0.0156 (11)	0.0086 (11)	-0.0182 (12)
O1	0.0758 (14)	0.0767 (14)	0.0974 (16)	-0.0167 (12)	-0.0118 (12)	0.0055 (13)
O2	0.0564 (11)	0.0705 (12)	0.0794 (13)	-0.0261 (10)	0.0020 (10)	-0.0061 (10)
O3	0.1308 (19)	0.1312 (19)	0.0715 (14)	-0.0868 (17)	0.0009 (13)	-0.0134 (14)
O4	0.0976 (14)	0.0991 (15)	0.0501 (12)	-0.0504 (12)	-0.0052 (10)	-0.0104 (11)
O5	0.0691 (12)	0.0804 (14)	0.0596 (12)	-0.0096 (11)	0.0072 (10)	-0.0249 (11)
O6	0.0455 (10)	0.0858 (13)	0.0635 (11)	-0.0079 (9)	-0.0075 (9)	-0.0268 (10)
O7	0.1317 (18)	0.0847 (14)	0.0563 (12)	-0.0599 (14)	-0.0076 (12)	-0.0072 (11)
O8	0.149 (2)	0.0852 (15)	0.0793 (16)	-0.0509 (15)	-0.0278 (14)	-0.0184 (13)

Geometric parameters (Å, °)

C1—O3	1.401 (3)	C19—H19B	0.9600
C1—H1A	0.9600	C19—H19C	0.9600
C1—H1B	0.9600	C20—O8	1.362 (3)
C1—H1C	0.9600	C20—C21	1.375 (3)
C2—O3	1.369 (3)	C20—C25	1.402 (4)
C2—C3	1.371 (3)	C21—C22	1.381 (4)
C2—C7	1.400 (3)	C21—H21	0.9300
C3—C4	1.381 (4)	C22—C23	1.373 (4)
C3—H3	0.9300	C22—H22	0.9300
C4—C5	1.369 (3)	C23—C24	1.396 (3)
C4—H4A	0.9300	C23—H23	0.9300
C5—C6	1.397 (3)	C24—C25	1.387 (3)
C5—H5	0.9300	C24—C26	1.453 (3)
C6—C7	1.397 (3)	C25—O7	1.353 (3)
C6—C8	1.447 (3)	C26—N2	1.270 (3)
C7—O4	1.346 (3)	C26—H26	0.9300
C8—N4	1.270 (3)	C27—N2	1.469 (3)
C8—H8	0.9300	C27—C37	1.535 (3)
C9—N4	1.457 (3)	C27—H27A	0.9700
C9—C37	1.535 (3)	C27—H27B	0.9700
C9—H9A	0.9700	C28—O5	1.415 (3)
C9—H9B	0.9700	C28—H28A	0.9600

C10—O1	1.390 (3)	C28—H28B	0.9600
C10—H10A	0.9600	C28—H28C	0.9600
C10—H10B	0.9600	C29—O5	1.372 (3)
C10—H10C	0.9600	C29—C30	1.378 (3)
C11—C12	1.371 (3)	C29—C34	1.389 (3)
C11—O1	1.380 (3)	C30—C31	1.381 (4)
C11—C16	1.392 (3)	C30—H30	0.9300
C12—C13	1.379 (4)	C31—C32	1.363 (4)
C12—H12	0.9300	C31—H31	0.9300
C13—C14	1.365 (4)	C32—C33	1.397 (3)
C13—H13	0.9300	C32—H32	0.9300
C14—C15	1.391 (3)	C33—C34	1.393 (3)
C14—H14	0.9300	C33—C35	1.452 (3)
C15—C16	1.396 (3)	C34—O6	1.353 (3)
C15—C17	1.447 (3)	C35—N3	1.270 (3)
C16—O2	1.356 (3)	C35—H35	0.9300
C17—N1	1.275 (3)	C36—N3	1.468 (3)
C17—H17	0.9300	C36—C37	1.525 (3)
C18—N1	1.454 (3)	C36—H36A	0.9700
C18—C37	1.534 (3)	C36—H36B	0.9700
C18—H18A	0.9700	O2—H2	0.8200
C18—H18B	0.9700	O4—H4	0.8200
C19—O8	1.432 (3)	O6—H6	0.8200
C19—H19A	0.9600	O7—H7	0.8200
O3—C1—H1A	109.5	C20—C21—C22	121.1 (3)
O3—C1—H1B	109.5	C20—C21—H21	119.5
H1A—C1—H1B	109.5	C22—C21—H21	119.5
O3—C1—H1C	109.5	C23—C22—C21	119.4 (3)
H1A—C1—H1C	109.5	C23—C22—H22	120.3
H1B—C1—H1C	109.5	C21—C22—H22	120.3
O3—C2—C3	125.3 (3)	C22—C23—C24	121.1 (3)
O3—C2—C7	114.8 (2)	C22—C23—H23	119.5
C3—C2—C7	119.8 (3)	C24—C23—H23	119.4
C2—C3—C4	120.6 (3)	C25—C24—C23	118.9 (2)
C2—C3—H3	119.7	C25—C24—C26	121.4 (2)
C4—C3—H3	119.7	C23—C24—C26	119.8 (2)
C5—C4—C3	120.3 (3)	O7—C25—C24	121.4 (2)
C5—C4—H4A	119.8	O7—C25—C20	118.3 (3)
C3—C4—H4A	119.8	C24—C25—C20	120.2 (3)
C4—C5—C6	120.4 (3)	N2—C26—C24	123.1 (2)
C4—C5—H5	119.8	N2—C26—H26	118.4
C6—C5—H5	119.8	C24—C26—H26	118.4
C7—C6—C5	119.1 (2)	N2—C27—C37	112.89 (19)
C7—C6—C8	120.2 (2)	N2—C27—H27A	109.0
C5—C6—C8	120.7 (2)	C37—C27—H27A	109.0
O4—C7—C6	122.0 (2)	N2—C27—H27B	109.0
O4—C7—C2	118.3 (2)	C37—C27—H27B	109.0

C6—C7—C2	119.6 (2)	H27A—C27—H27B	107.8
N4—C8—C6	122.1 (2)	O5—C28—H28A	109.5
N4—C8—H8	118.9	O5—C28—H28B	109.5
C6—C8—H8	118.9	H28A—C28—H28B	109.5
N4—C9—C37	111.28 (19)	O5—C28—H28C	109.5
N4—C9—H9A	109.4	H28A—C28—H28C	109.5
C37—C9—H9A	109.4	H28B—C28—H28C	109.5
N4—C9—H9B	109.4	O5—C29—C30	124.4 (3)
C37—C9—H9B	109.4	O5—C29—C34	115.6 (2)
H9A—C9—H9B	108.0	C30—C29—C34	120.0 (3)
O1—C10—H10A	109.5	C29—C30—C31	119.8 (3)
O1—C10—H10B	109.5	C29—C30—H30	120.1
H10A—C10—H10B	109.5	C31—C30—H30	120.1
O1—C10—H10C	109.5	C32—C31—C30	120.9 (3)
H10A—C10—H10C	109.5	C32—C31—H31	119.5
H10B—C10—H10C	109.5	C30—C31—H31	119.5
C12—C11—O1	125.0 (3)	C31—C32—C33	120.2 (3)
C12—C11—C16	120.1 (3)	C31—C32—H32	119.9
O1—C11—C16	114.9 (2)	C33—C32—H32	119.9
C11—C12—C13	120.2 (3)	C34—C33—C32	119.1 (3)
C11—C12—H12	119.9	C34—C33—C35	121.7 (2)
C13—C12—H12	119.9	C32—C33—C35	119.2 (2)
C14—C13—C12	120.3 (3)	O6—C34—C29	118.8 (2)
C14—C13—H13	119.9	O6—C34—C33	121.2 (2)
C12—C13—H13	119.9	C29—C34—C33	120.0 (2)
C13—C14—C15	120.8 (3)	N3—C35—C33	123.2 (2)
C13—C14—H14	119.6	N3—C35—H35	118.4
C15—C14—H14	119.6	C33—C35—H35	118.4
C14—C15—C16	118.7 (2)	N3—C36—C37	113.1 (2)
C14—C15—C17	120.3 (2)	N3—C36—H36A	109.0
C16—C15—C17	120.9 (2)	C37—C36—H36A	109.0
O2—C16—C11	118.4 (2)	N3—C36—H36B	109.0
O2—C16—C15	121.8 (2)	C37—C36—H36B	109.0
C11—C16—C15	119.8 (2)	H36A—C36—H36B	107.8
N1—C17—C15	122.3 (2)	C36—C37—C18	109.1 (2)
N1—C17—H17	118.8	C36—C37—C27	111.42 (19)
C15—C17—H17	118.8	C18—C37—C27	107.39 (18)
N1—C18—C37	111.7 (2)	C36—C37—C9	108.6 (2)
N1—C18—H18A	109.3	C18—C37—C9	110.63 (19)
C37—C18—H18A	109.3	C27—C37—C9	109.64 (19)
N1—C18—H18B	109.3	C17—N1—C18	119.5 (2)
C37—C18—H18B	109.3	C26—N2—C27	116.9 (2)
H18A—C18—H18B	107.9	C35—N3—C36	117.2 (2)
O8—C19—H19A	109.5	C8—N4—C9	120.7 (2)
O8—C19—H19B	109.5	C11—O1—C10	118.6 (2)
H19A—C19—H19B	109.5	C16—O2—H2	109.5
O8—C19—H19C	109.5	C2—O3—C1	118.8 (2)
H19A—C19—H19C	109.5	C7—O4—H4	109.5

H19B—C19—H19C	109.5	C29—O5—C28	118.0 (2)
O8—C20—C21	125.5 (3)	C34—O6—H6	109.5
O8—C20—C25	115.1 (3)	C25—O7—H7	109.5
C21—C20—C25	119.3 (3)	C20—O8—C19	117.6 (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>
O2—H2...N1	0.82	1.87	2.592 (3)	147
O4—H4...N4	0.82	1.84	2.567 (2)	147
O6—H6...N3	0.82	1.89	2.623 (3)	148
O7—H7...N2	0.82	1.88	2.612 (3)	148
C27—H27 <i>B</i> ...N1	0.97	2.56	2.936 (3)	103
C26—H26...O5 ⁱ	0.93	2.58	3.309 (3)	136
C28—H28 <i>C</i> ...O2 ⁱ	0.96	2.57	3.387 (3)	143
C12—H12...O6 ⁱⁱ	0.93	2.59	3.400 (3)	146
C32—H32...O3 ⁱⁱⁱ	0.93	2.42	3.289 (3)	155
C36—H36 <i>B</i> ...O8 ^{iv}	0.97	2.55	3.399 (3)	146

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