

Methyl 1-(7-acetamido-5,8-dimethoxyquinolin-2-yl)-4-methyl- β -carboline-3-carboxylate

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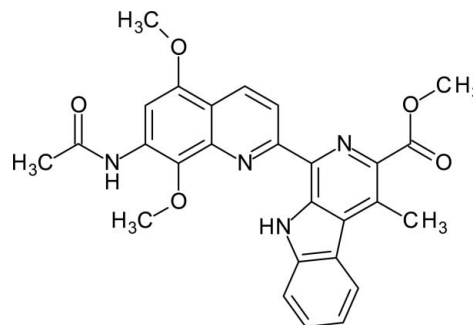
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 Key indicators: single-crystal X-ray study; $T = 193$ K; mean $\sigma(\text{C}-\text{C}) = 0.008$ Å; R factor = 0.083; wR factor = 0.248; data-to-parameter ratio = 13.4.

The title compound, $\text{C}_{27}\text{H}_{24}\text{N}_4\text{O}_5$, is an intermediate in the synthesis of lavendamycin *via* a ruthenium-catalysed [2 + 2 + 2] cycloaddition. An intramolecular hydrogen-bond bridge from the carboline to the quinoline stabilizes a highly planar geometry [maximum deviation = 0.065 (6) Å] for the two rigid units. This hydrogen-bond-stabilized coplanarity has a very close analogy in the structure of the antitumor antibiotic streptonigrin in the solid state and in solution. Intermolecular hydrogen-bond bridges of amides groups along the a axis and π - π stacking interactions [centroid-centroid distance = 3.665 (9) Å] connect molecules arranged in a parallel manner.

Related literature

For metal-catalysed transformations of tethered alkynyl-ynamides to carbolines and other heteroannulated indoles, see: Nissen *et al.* (2011); Dassonneville *et al.* (2010, 2011). For the synthesis of the natural product lavendamycin (systematic name 1-(7-amino-5,8-dioxoquinolin-2-yl)-4-methyl-9H-pyrido[3,4-*b*]indole-3-carboxylic acid) *via* [2 + 2 + 2] cycloaddition, see: Nissen & Detert (2011). For the isolation of lavendamycin from *streptomyces lavendulae*, see: Doyle *et al.* (1981). For the anti-tumor activity of lavendamycin, see: Fang *et al.* (2003). For the preparation of lavendamycin, see: Behforouz *et al.* (1996) Godard *et al.* (1993). For related structures, see: Chiu & Lipscomb (1975); Harding *et al.* (1993).



Experimental

Crystal data

$\text{C}_{27}\text{H}_{24}\text{N}_4\text{O}_5$	$\gamma = 96.85$ (11) $^\circ$
$M_r = 484.50$	$V = 1161$ (4) Å ³
Triclinic, $P\bar{1}$	$Z = 2$
$a = 4.646$ (10) Å	Cu $K\alpha$ radiation
$b = 13.81$ (3) Å	$\mu = 0.80$ mm ⁻¹
$c = 18.768$ (19) Å	$T = 193$ K
$\alpha = 102.06$ (9) $^\circ$	$0.58 \times 0.06 \times 0.03$ mm
$\beta = 95.23$ (10) $^\circ$	

Data collection

Enraf-Nonius CAD-4 diffractometer	2069 reflections with $I > 2\sigma(I)$
4993 measured reflections	$R_{\text{int}} = 0.054$
4412 independent reflections	3 standard reflections every 60 min
	intensity decay: 2%

Refinement

$R[F^2 > 2\sigma(F^2)] = 0.083$	330 parameters
$wR(F^2) = 0.248$	H-atom parameters constrained
$S = 0.99$	$\Delta\rho_{\text{max}} = 0.26$ e Å ⁻³
4412 reflections	$\Delta\rho_{\text{min}} = -0.34$ e Å ⁻³

Table 1

 Hydrogen-bond geometry (Å, $^\circ$).

$D-H\cdots A$	$D-H$	$H\cdots A$	$D\cdots A$	$D-H\cdots A$
N2-H2 \cdots N3	0.88	2.17	2.693 (7)	118
N4-H4 \cdots O5 ⁱ	0.88	1.98	2.824 (8)	160

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

Data collection: *CAD-4 Software* (Enraf-Nonius, 1989); cell refinement: *CAD-4 Software*; data reduction: *CORINC* (Dräger & Gattow, 1971); program(s) used to solve structure: *SIR97* (Altomare *et al.*, 1999); program(s) used to refine structure: *SHELXL97* (Sheldrick, 2008); molecular graphics: *PLATON* (Spek, 2009); software used to prepare material for publication: *PLATON*.

Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: BT5551).

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supporting information

Acta Cryst. (2011). E67, o1497–o1498 [doi:10.1107/S1600536811018794]

Methyl 1-(7-acetamido-5,8-dimethoxyquinolin-2-yl)-4-methyl- β -carboline-3-carboxylate

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S1. Comment

The title compound was prepared as part of a larger project focusing on the synthesis of indolo-annulated heterocycles *via* [2 + 2+2] cycloaddition of alkynyl-ynamides and nitriles or heterocumulenes catalyzed by rhodium or ruthenium (Nissen *et al.*, 2011; Dassonneville *et al.*, 2011) and is a synthetic precursor of the natural product lavendamycin (Nissen & Detert, 2011). The molecule is built up by two rigid and planar units, both are coplanar. This flat structure is stabilized by an intramolecular hydrogen bond from the carboline-NH to the quinoline-N N2—H2 \cdots N3 with a distance of 2.17 Å. Whereas the ester group is coplanar with the carboline core [torsion angle O1—C1—C3—C9 of -0.4 (8)°], the acetamido group is twisted out of the plane of the quinoline framework. The torsion angle C20—N4—C19—C22 amounts to -47.9 (7)°. This torsion can result from the sterical hindrance due to the neighbouring methoxy group, but its ability to act as a hydrogen bridge donor and acceptor is important for the formation of crystals of the title compound. Parallel molecules, arranged in the shape of tilted staples are connected to infinite chains *via* H-bridging along the *a*-axis. The tilt angle of the mean plane of the molecules relative to the *ab*-plane is 54.3°. The second interaction is π - π -stacking, resulting in a distance of only 3.665 (9) Å of the centroids of the pyrrole ring (N2—C6) and the benzo-ring (C6—C14) (symmetry code 1+*X*,*Y*,*Z*). Two molecules, connected *via* a center of inversion fill the unit cell.

S2. Experimental

Pd/C (10%, 47 mg, 44 μ mol, 5 mol%) was added to a suspension of Methyl 1-(5,8-dimethoxy-7-nitroquinolin-2-yl)-4-methyl- β -carboline-3-carboxylate (405 mg, 0.86 mmol) (Nissen & Detert, 2011) in THF (170 ml) and the mixture was stirred 13 h under H₂ atmosphere. Ac₂O (5.0 ml) was added and the mixture was heated to 323 K. After 3 h the solvent was removed *in vacuo*, the residue dissolved in CH₂Cl₂ and filtered through celite. The filtrate was washed with aqueous NaHCO₃ (8%, 8 ml), dried (MgSO₄) and concentrated. Crystallization from chloroform/diethyl ether yielded the title compound (399 mg, 0.82 mmol, 96%) as a bright yellow solid. *M.p.* 467–468 K *R_f*: 0.12 (SiO₂, hexane:ethyl acetate:ethanol 6:3:1).

S3. Refinement

Hydrogen atoms attached to carbons were placed at calculated positions with C—H = 0.95 Å (aromatic) or 0.98–0.99 Å (*sp*³ C-atom). All H atoms were refined in the riding-model approximation with isotropic displacement parameters set at 1.2–1.5 times of the *U*_{eq} of the parent atom.

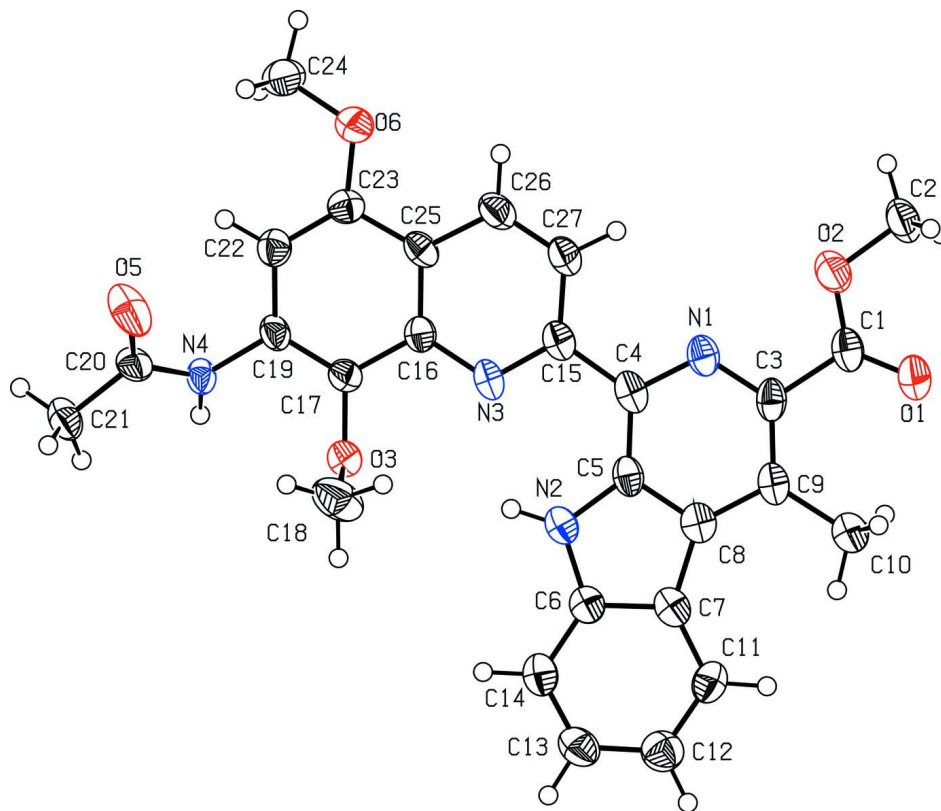


Figure 1

View of compound I. Displacement ellipsoids are drawn at the 50% probability level.

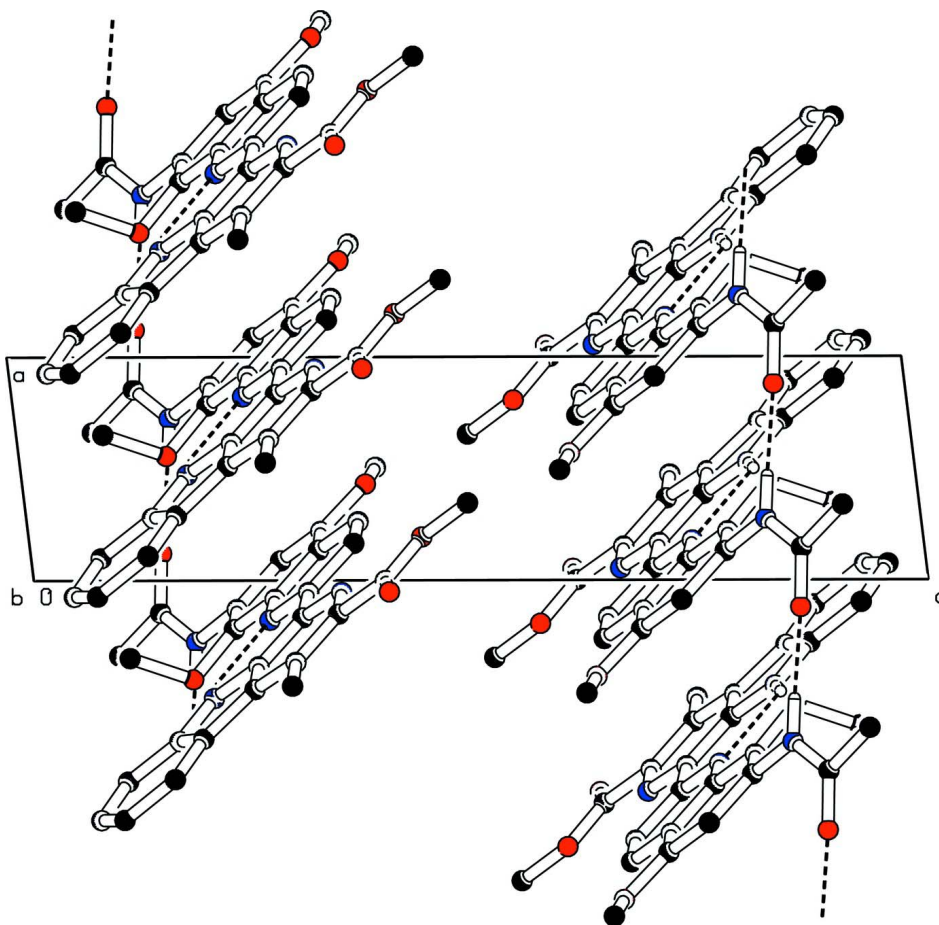


Figure 2

Part of the packing diagram showing the hydrogen bonds. View along *b*-axis.

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Crystal data

$C_{27}H_{24}N_4O_5$

$M_r = 484.50$

Triclinic, $P\bar{1}$

Hall symbol: -P 1

$a = 4.646 (10) \text{ \AA}$

$b = 13.81 (3) \text{ \AA}$

$c = 18.768 (19) \text{ \AA}$

$\alpha = 102.06 (9)^\circ$

$\beta = 95.23 (10)^\circ$

$\gamma = 96.85 (11)^\circ$

$V = 1161 (4) \text{ \AA}^3$

$Z = 2$

$F(000) = 508$

$D_x = 1.386 \text{ Mg m}^{-3}$

Cu $K\alpha$ radiation, $\lambda = 1.54178 \text{ \AA}$

Cell parameters from 25 reflections

$\theta = 10\text{--}25^\circ$

$\mu = 0.80 \text{ mm}^{-1}$

$T = 193 \text{ K}$

10, yellow

$0.58 \times 0.06 \times 0.03 \text{ mm}$

Data collection

Enraf-Nonius CAD-4

diffractometer

Radiation source: rotating anode

Graphite monochromator

$\omega/2\theta$ scans

4993 measured reflections

4412 independent reflections

2069 reflections with $I > 2\sigma(I)$

$R_{\text{int}} = 0.054$

$\theta_{\text{max}} = 70.0^\circ$, $\theta_{\text{min}} = 2.4^\circ$

$h = -5 \rightarrow 0$
 $k = -16 \rightarrow 16$
 $l = -22 \rightarrow 22$

3 standard reflections every 60 min
 intensity decay: 2%

Refinement

Refinement on F^2
 Least-squares matrix: full
 $R[F^2 > 2\sigma(F^2)] = 0.083$
 $wR(F^2) = 0.248$
 $S = 0.99$
 4412 reflections
 330 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
 $w = 1/[\sigma^2(F_o^2) + (0.1159P)^2]$
 where $P = (F_o^2 + 2F_c^2)/3$
 $(\Delta/\sigma)_{\max} < 0.001$
 $\Delta\rho_{\max} = 0.26 \text{ e } \text{\AA}^{-3}$
 $\Delta\rho_{\min} = -0.34 \text{ e } \text{\AA}^{-3}$
 Extinction correction: *SHELXL97* (Sheldrick,
 2008), $F_c^* = kF_c[1 + 0.001x F_c^2 \lambda^3 / \sin(2\theta)]^{-1/4}$
 Extinction coefficient: 0.0010 (5)

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)

	x	y	z	$U_{\text{iso}}^*/U_{\text{eq}}$
O1	0.9484 (9)	0.6415 (3)	0.3958 (2)	0.0631 (12)
O2	1.1937 (9)	0.5349 (3)	0.43953 (19)	0.0530 (11)
O3	0.5569 (7)	-0.0529 (2)	0.16500 (18)	0.0381 (8)
N4	0.7247 (8)	-0.2335 (3)	0.1735 (2)	0.0314 (9)
H4	0.5329	-0.2404	0.1660	0.038*
O5	1.1210 (8)	-0.3053 (3)	0.1464 (2)	0.0691 (14)
O6	1.4309 (8)	-0.0606 (3)	0.38367 (19)	0.0469 (10)
N1	0.9463 (9)	0.3816 (3)	0.3424 (2)	0.0363 (10)
N2	0.4724 (9)	0.2191 (3)	0.1858 (2)	0.0367 (10)
H2	0.4994	0.1561	0.1806	0.044*
N3	0.8225 (9)	0.1206 (3)	0.2585 (2)	0.0349 (10)
C1	0.9992 (12)	0.5579 (4)	0.3929 (3)	0.0392 (12)
C2	1.3460 (13)	0.6178 (4)	0.4953 (3)	0.0506 (15)
H2A	1.4680	0.6626	0.4725	0.076*
H2B	1.4695	0.5926	0.5305	0.076*
H2C	1.2039	0.6545	0.5210	0.076*
C3	0.8488 (11)	0.4675 (3)	0.3360 (3)	0.0362 (12)
C4	0.8328 (11)	0.2958 (3)	0.2955 (3)	0.0361 (12)
C5	0.6143 (11)	0.2973 (3)	0.2397 (3)	0.0341 (11)
C6	0.2800 (11)	0.2556 (4)	0.1412 (3)	0.0352 (11)

C7	0.2906 (11)	0.3594 (4)	0.1682 (3)	0.0355 (11)
C8	0.5118 (11)	0.3860 (3)	0.2323 (3)	0.0344 (11)
C9	0.6334 (11)	0.4754 (3)	0.2817 (3)	0.0368 (12)
C10	0.5268 (13)	0.5721 (4)	0.2747 (3)	0.0508 (15)
H10A	0.4642	0.6039	0.3214	0.076*
H10B	0.3616	0.5585	0.2359	0.076*
H10C	0.6850	0.6169	0.2623	0.076*
C11	0.1095 (11)	0.4101 (4)	0.1309 (3)	0.0416 (13)
H11	0.1105	0.4800	0.1476	0.050*
C12	-0.0723 (12)	0.3573 (4)	0.0691 (3)	0.0490 (14)
H12	-0.1993	0.3914	0.0439	0.059*
C13	-0.0731 (12)	0.2558 (4)	0.0431 (3)	0.0477 (14)
H13	-0.1969	0.2220	-0.0002	0.057*
C14	0.0991 (12)	0.2038 (4)	0.0782 (3)	0.0432 (13)
H14	0.0965	0.1340	0.0605	0.052*
C15	0.9462 (11)	0.2043 (3)	0.3056 (3)	0.0367 (12)
C16	0.9155 (11)	0.0334 (3)	0.2670 (3)	0.0320 (11)
C17	0.7812 (10)	-0.0564 (3)	0.2176 (3)	0.0330 (11)
C18	0.6517 (15)	-0.0421 (5)	0.0971 (3)	0.0645 (19)
H18A	0.7902	0.0195	0.1047	0.097*
H18B	0.4831	-0.0388	0.0627	0.097*
H18C	0.7473	-0.0996	0.0769	0.097*
C19	0.8703 (10)	-0.1450 (3)	0.2232 (3)	0.0323 (11)
C20	0.8550 (10)	-0.3067 (4)	0.1373 (3)	0.0374 (12)
C21	0.6628 (11)	-0.3906 (4)	0.0857 (3)	0.0447 (14)
H21A	0.7272	-0.3984	0.0368	0.067*
H21B	0.4610	-0.3762	0.0833	0.067*
H21C	0.6731	-0.4525	0.1029	0.067*
C22	1.0881 (11)	-0.1513 (4)	0.2785 (3)	0.0373 (12)
H22	1.1437	-0.2143	0.2820	0.045*
C23	1.2197 (10)	-0.0657 (4)	0.3273 (3)	0.0338 (11)
C24	1.5056 (13)	-0.1523 (4)	0.3972 (3)	0.0514 (15)
H24A	1.3311	-0.1930	0.4064	0.077*
H24B	1.6528	-0.1387	0.4403	0.077*
H24C	1.5846	-0.1885	0.3544	0.077*
C25	1.1358 (11)	0.0293 (3)	0.3224 (3)	0.0345 (11)
C26	1.2625 (12)	0.1205 (4)	0.3705 (3)	0.0413 (13)
H26	1.4150	0.1214	0.4080	0.050*
C27	1.1656 (11)	0.2061 (4)	0.3626 (3)	0.0392 (12)
H27	1.2453	0.2676	0.3955	0.047*

Atomic displacement parameters (\AA^2)

	U^{11}	U^{22}	U^{33}	U^{12}	U^{13}	U^{23}
O1	0.068 (3)	0.032 (2)	0.077 (3)	0.006 (2)	-0.016 (2)	-0.004 (2)
O2	0.073 (3)	0.037 (2)	0.042 (2)	-0.003 (2)	-0.006 (2)	0.0036 (17)
O3	0.038 (2)	0.0337 (18)	0.0398 (19)	0.0018 (16)	-0.0003 (17)	0.0052 (15)
N4	0.022 (2)	0.0242 (19)	0.044 (2)	0.0010 (16)	0.0027 (18)	0.0013 (17)

O5	0.0185 (19)	0.067 (3)	0.097 (3)	0.0068 (18)	-0.001 (2)	-0.033 (2)
O6	0.047 (2)	0.039 (2)	0.050 (2)	0.0009 (17)	-0.0097 (18)	0.0100 (17)
N1	0.039 (2)	0.028 (2)	0.040 (2)	-0.0042 (18)	0.009 (2)	0.0057 (18)
N2	0.036 (2)	0.031 (2)	0.038 (2)	-0.0019 (18)	0.002 (2)	0.0024 (18)
N3	0.038 (2)	0.027 (2)	0.034 (2)	-0.0025 (18)	0.0078 (19)	-0.0016 (17)
C1	0.042 (3)	0.029 (3)	0.044 (3)	-0.002 (2)	0.012 (3)	0.003 (2)
C2	0.069 (4)	0.033 (3)	0.041 (3)	-0.006 (3)	0.000 (3)	0.000 (2)
C3	0.047 (3)	0.024 (2)	0.038 (3)	0.001 (2)	0.016 (2)	0.007 (2)
C4	0.040 (3)	0.032 (3)	0.033 (3)	-0.004 (2)	0.013 (2)	0.001 (2)
C5	0.033 (3)	0.027 (2)	0.039 (3)	-0.002 (2)	0.010 (2)	0.003 (2)
C6	0.031 (3)	0.037 (3)	0.037 (3)	0.004 (2)	0.008 (2)	0.005 (2)
C7	0.030 (3)	0.039 (3)	0.035 (3)	0.001 (2)	0.011 (2)	0.002 (2)
C8	0.034 (3)	0.036 (3)	0.034 (3)	0.001 (2)	0.015 (2)	0.008 (2)
C9	0.042 (3)	0.028 (2)	0.038 (3)	0.000 (2)	0.012 (2)	0.002 (2)
C10	0.058 (4)	0.038 (3)	0.050 (3)	0.004 (3)	-0.007 (3)	0.004 (3)
C11	0.041 (3)	0.035 (3)	0.052 (3)	0.010 (2)	0.010 (3)	0.011 (3)
C12	0.041 (3)	0.052 (3)	0.053 (3)	0.004 (3)	0.004 (3)	0.012 (3)
C13	0.039 (3)	0.050 (3)	0.048 (3)	-0.002 (3)	-0.002 (3)	0.004 (3)
C14	0.045 (3)	0.035 (3)	0.047 (3)	0.001 (2)	0.007 (3)	0.005 (2)
C15	0.040 (3)	0.031 (3)	0.036 (3)	-0.002 (2)	0.006 (2)	0.002 (2)
C16	0.033 (3)	0.028 (2)	0.034 (3)	0.001 (2)	0.005 (2)	0.006 (2)
C17	0.029 (3)	0.033 (3)	0.034 (3)	0.002 (2)	0.000 (2)	0.005 (2)
C18	0.075 (5)	0.071 (4)	0.040 (3)	-0.016 (4)	-0.004 (3)	0.016 (3)
C19	0.028 (3)	0.025 (2)	0.039 (3)	-0.003 (2)	0.001 (2)	0.002 (2)
C20	0.020 (2)	0.040 (3)	0.046 (3)	0.004 (2)	-0.003 (2)	-0.001 (2)
C21	0.032 (3)	0.039 (3)	0.054 (3)	0.002 (2)	0.002 (3)	-0.009 (3)
C22	0.038 (3)	0.032 (3)	0.041 (3)	0.003 (2)	0.005 (2)	0.008 (2)
C23	0.026 (2)	0.038 (3)	0.037 (3)	0.001 (2)	0.003 (2)	0.011 (2)
C24	0.047 (3)	0.049 (3)	0.057 (4)	0.003 (3)	-0.008 (3)	0.020 (3)
C25	0.032 (3)	0.034 (3)	0.033 (3)	-0.006 (2)	0.002 (2)	0.003 (2)
C26	0.040 (3)	0.038 (3)	0.038 (3)	-0.007 (2)	-0.004 (2)	0.004 (2)
C27	0.043 (3)	0.031 (3)	0.037 (3)	-0.003 (2)	0.002 (2)	0.001 (2)

Geometric parameters (Å, °)

O1—C1	1.198 (6)	C16—C17	1.420 (7)
O2—C1	1.313 (6)	C17—C19	1.359 (7)
O2—C2	1.442 (6)	C19—C22	1.406 (7)
O3—C17	1.380 (6)	C20—C21	1.485 (7)
O3—C18	1.417 (7)	C22—C23	1.373 (7)
N4—C20	1.336 (6)	C23—C25	1.431 (7)
N4—C19	1.425 (6)	C25—C26	1.414 (7)
O5—C20	1.229 (6)	C26—C27	1.346 (7)
O6—C23	1.360 (6)	N2—H2	0.8800
O6—C24	1.419 (6)	N4—H4	0.8800
N1—C4	1.337 (6)	C2—H2A	0.9800
N1—C3	1.342 (6)	C2—H2B	0.9800
N2—C5	1.370 (6)	C2—H2C	0.9800

N2—C6	1.380 (6)	C10—H10A	0.9800
N3—C15	1.330 (6)	C10—H10B	0.9800
N3—C16	1.362 (6)	C10—H10C	0.9800
C1—C3	1.512 (7)	C11—H11	0.9500
C3—C9	1.392 (7)	C12—H12	0.9500
C4—C5	1.396 (7)	C13—H13	0.9500
C4—C15	1.467 (7)	C14—H14	0.9500
C5—C8	1.395 (7)	C18—H18A	0.9800
C6—C14	1.392 (7)	C18—H18B	0.9800
C6—C7	1.410 (7)	C18—H18C	0.9800
C7—C11	1.388 (7)	C21—H21A	0.9800
C7—C8	1.462 (7)	C21—H21B	0.9800
C8—C9	1.400 (7)	C21—H21C	0.9800
C9—C10	1.505 (7)	C22—H22	0.9500
C11—C12	1.383 (7)	C24—H24A	0.9800
C12—C13	1.383 (8)	C24—H24B	0.9800
C13—C14	1.352 (7)	C24—H24C	0.9800
C15—C27	1.402 (7)	C26—H26	0.9500
C16—C25	1.406 (7)	C27—H27	0.9500
C1—O2—C2	115.5 (4)	C16—C25—C26	117.4 (5)
C17—O3—C18	113.6 (4)	C16—C25—C23	119.0 (4)
C20—N4—C19	125.5 (4)	C26—C25—C23	123.6 (5)
C23—O6—C24	117.3 (4)	C27—C26—C25	119.5 (5)
C4—N1—C3	120.2 (5)	C26—C27—C15	119.7 (5)
C5—N2—C6	108.6 (4)	C6—N2—H2	126.00
C15—N3—C16	117.6 (4)	C5—N2—H2	126.00
O1—C1—O2	123.1 (5)	C19—N4—H4	117.00
O1—C1—C3	124.4 (5)	C20—N4—H4	117.00
O2—C1—C3	112.5 (4)	O2—C2—H2A	110.00
N1—C3—C9	124.4 (4)	O2—C2—H2B	109.00
N1—C3—C1	113.9 (5)	O2—C2—H2C	109.00
C9—C3—C1	121.7 (4)	H2A—C2—H2B	109.00
N1—C4—C5	118.9 (5)	H2A—C2—H2C	109.00
N1—C4—C15	117.8 (5)	H2B—C2—H2C	109.00
C5—C4—C15	123.2 (4)	C9—C10—H10A	109.00
N2—C5—C8	110.2 (4)	C9—C10—H10B	109.00
N2—C5—C4	128.5 (5)	C9—C10—H10C	110.00
C8—C5—C4	121.4 (4)	H10A—C10—H10B	109.00
N2—C6—C14	128.4 (5)	H10A—C10—H10C	109.00
N2—C6—C7	109.4 (4)	H10B—C10—H10C	109.00
C14—C6—C7	122.2 (5)	C7—C11—H11	120.00
C11—C7—C6	118.2 (5)	C12—C11—H11	121.00
C11—C7—C8	136.1 (5)	C11—C12—H12	119.00
C6—C7—C8	105.7 (4)	C13—C12—H12	119.00
C5—C8—C9	119.0 (5)	C12—C13—H13	119.00
C5—C8—C7	106.1 (4)	C14—C13—H13	119.00
C9—C8—C7	134.8 (5)	C6—C14—H14	121.00

C3—C9—C8	116.0 (5)	C13—C14—H14	121.00
C3—C9—C10	124.0 (4)	O3—C18—H18A	110.00
C8—C9—C10	120.0 (5)	O3—C18—H18B	109.00
C12—C11—C7	119.0 (5)	O3—C18—H18C	109.00
C11—C12—C13	121.4 (5)	H18A—C18—H18B	109.00
C14—C13—C12	121.3 (5)	H18A—C18—H18C	109.00
C13—C14—C6	118.0 (5)	H18B—C18—H18C	109.00
N3—C15—C27	123.1 (5)	C20—C21—H21A	109.00
N3—C15—C4	115.5 (5)	C20—C21—H21B	109.00
C27—C15—C4	121.4 (4)	C20—C21—H21C	110.00
N3—C16—C25	122.7 (4)	H21A—C21—H21B	109.00
N3—C16—C17	117.9 (4)	H21A—C21—H21C	109.00
C25—C16—C17	119.4 (4)	H21B—C21—H21C	109.00
C19—C17—O3	120.5 (4)	C19—C22—H22	120.00
C19—C17—C16	120.0 (4)	C23—C22—H22	120.00
O3—C17—C16	119.5 (4)	O6—C24—H24A	109.00
C17—C19—C22	121.7 (4)	O6—C24—H24B	109.00
C17—C19—N4	118.0 (4)	O6—C24—H24C	109.00
C22—C19—N4	120.2 (4)	H24A—C24—H24B	109.00
O5—C20—N4	121.8 (5)	H24A—C24—H24C	109.00
O5—C20—C21	121.4 (5)	H24B—C24—H24C	110.00
N4—C20—C21	116.8 (4)	C25—C26—H26	120.00
C23—C22—C19	119.5 (5)	C27—C26—H26	120.00
O6—C23—C22	125.9 (5)	C15—C27—H27	120.00
O6—C23—C25	113.7 (4)	C26—C27—H27	120.00
C22—C23—C25	120.5 (5)		
C2—O2—C1—O1	-0.2 (8)	N2—C6—C14—C13	179.9 (5)
C2—O2—C1—C3	179.9 (4)	C7—C6—C14—C13	-0.7 (8)
C4—N1—C3—C9	-1.0 (7)	C16—N3—C15—C27	0.1 (7)
C4—N1—C3—C1	-179.9 (4)	C16—N3—C15—C4	-178.3 (4)
O1—C1—C3—N1	178.5 (5)	N1—C4—C15—N3	178.7 (4)
O2—C1—C3—N1	-1.6 (6)	C5—C4—C15—N3	-1.4 (7)
O1—C1—C3—C9	-0.4 (8)	N1—C4—C15—C27	0.3 (7)
O2—C1—C3—C9	179.5 (5)	C5—C4—C15—C27	-179.8 (5)
C3—N1—C4—C5	0.6 (7)	C15—N3—C16—C25	-0.6 (7)
C3—N1—C4—C15	-179.5 (4)	C15—N3—C16—C17	179.1 (5)
C6—N2—C5—C8	1.5 (5)	C18—O3—C17—C19	-88.9 (6)
C6—N2—C5—C4	-178.0 (5)	C18—O3—C17—C16	92.3 (6)
N1—C4—C5—N2	179.1 (5)	N3—C16—C17—C19	179.0 (4)
C15—C4—C5—N2	-0.8 (8)	C25—C16—C17—C19	-1.4 (7)
N1—C4—C5—C8	-0.4 (7)	N3—C16—C17—O3	-2.3 (7)
C15—C4—C5—C8	179.7 (4)	C25—C16—C17—O3	177.4 (4)
C5—N2—C6—C14	177.4 (5)	O3—C17—C19—C22	-176.8 (4)
C5—N2—C6—C7	-2.0 (5)	C16—C17—C19—C22	2.0 (7)
N2—C6—C7—C11	-179.4 (4)	O3—C17—C19—N4	-0.2 (7)
C14—C6—C7—C11	1.1 (7)	C16—C17—C19—N4	178.6 (4)
N2—C6—C7—C8	1.7 (5)	C20—N4—C19—C17	135.4 (5)

C14—C6—C7—C8	-177.8 (5)	C20—N4—C19—C22	-47.9 (7)
N2—C5—C8—C9	-178.9 (4)	C19—N4—C20—O5	2.7 (8)
C4—C5—C8—C9	0.6 (7)	C19—N4—C20—C21	-177.5 (4)
N2—C5—C8—C7	-0.4 (5)	C17—C19—C22—C23	-1.4 (7)
C4—C5—C8—C7	179.1 (4)	N4—C19—C22—C23	-178.0 (5)
C11—C7—C8—C5	-179.3 (6)	C24—O6—C23—C22	-5.6 (7)
C6—C7—C8—C5	-0.8 (5)	C24—O6—C23—C25	173.7 (4)
C11—C7—C8—C9	-1.1 (10)	C19—C22—C23—O6	179.5 (5)
C6—C7—C8—C9	177.4 (5)	C19—C22—C23—C25	0.2 (7)
N1—C3—C9—C8	1.2 (7)	N3—C16—C25—C26	-0.1 (7)
C1—C3—C9—C8	179.9 (5)	C17—C16—C25—C26	-179.8 (5)
N1—C3—C9—C10	-179.7 (5)	N3—C16—C25—C23	179.9 (4)
C1—C3—C9—C10	-0.9 (8)	C17—C16—C25—C23	0.2 (7)
C5—C8—C9—C3	-1.0 (7)	O6—C23—C25—C16	-179.0 (4)
C7—C8—C9—C3	-179.0 (5)	C22—C23—C25—C16	0.3 (7)
C5—C8—C9—C10	179.9 (5)	O6—C23—C25—C26	1.0 (7)
C7—C8—C9—C10	1.9 (8)	C22—C23—C25—C26	-179.7 (5)
C6—C7—C11—C12	-0.1 (7)	C16—C25—C26—C27	1.4 (7)
C8—C7—C11—C12	178.3 (5)	C23—C25—C26—C27	-178.7 (5)
C7—C11—C12—C13	-1.2 (8)	C25—C26—C27—C15	-1.8 (8)
C11—C12—C13—C14	1.6 (9)	N3—C15—C27—C26	1.1 (8)
C12—C13—C14—C6	-0.6 (8)	C4—C15—C27—C26	179.5 (5)

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...N3	0.88	2.17	2.693 (7)	118
N4—H4...O5 ⁱ	0.88	1.98	2.824 (8)	160

Symmetry code: (i) *x*-1, *y*, *z*.