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6-Bromo-1-butyldoline-2,3-dione

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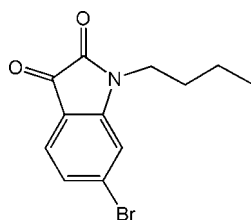
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Key indicators: single-crystal X-ray study; $T = 296$ K; mean $\sigma(\text{C}-\text{C}) = 0.005$ Å; R factor = 0.039; wR factor = 0.070; data-to-parameter ratio = 14.2.

There are two independent molecules in the asymmetric unit of the title compound, $\text{C}_{12}\text{H}_{12}\text{BrNO}_2$. The C—C bond lengths of the two carbonyl C atoms of the five-membered rings are distinctly longer than a normal $\text{Csp}^2-\text{Csp}^2$ single bond. One of the molecules makes parallel self-coupled (inversion) dimers by $\pi-\pi$ interactions with phenyl–phenyl interplanar distances of 3.403 (2) Å. The other molecule also forms self-dimers at longer phenyl–phenyl plane distances [3.649 (2) Å]. In the crystal, a C—H...O interaction is seen.

Related literature

For synthesis and applications, see: Kopka *et al.* (2006); Pirrung *et al.* (2005); Zhou *et al.* (2006). For related crystal structures, see: Goldschmidt & Llewellyn (1950); Palenik *et al.* (1990).



Experimental

Crystal data

$\text{C}_{12}\text{H}_{12}\text{BrNO}_2$
 $M_r = 282.14$
 Monoclinic, $P2_1/c$
 $a = 13.3097$ (2) Å
 $b = 11.8793$ (2) Å
 $c = 16.2238$ (2) Å
 $\beta = 112.340$ (1)°

$V = 2372.62$ (6) Å³
 $Z = 8$
 Mo $K\alpha$ radiation
 $\mu = 3.45$ mm⁻¹
 $T = 296$ (2) K
 $0.37 \times 0.13 \times 0.11$ mm

Data collection

Bruker APEXII CCD diffractometer
 Absorption correction: multi-scan (SADABS; Bruker, 2005)
 $T_{\min} = 0.526$, $T_{\max} = 0.744$
 (expected range = 0.487–0.689)

16713 measured reflections
 5170 independent reflections
 2424 reflections with $I > 2\sigma(I)$
 $R_{\text{int}} = 0.076$

Refinement

$R[F^2 > 2\sigma(F^2)] = 0.039$
 $wR(F^2) = 0.070$
 $S = 0.93$
 5170 reflections
 364 parameters
 1 restraint

H atoms treated by a mixture of independent and constrained refinement
 $\Delta\rho_{\text{max}} = 0.39$ e Å⁻³
 $\Delta\rho_{\text{min}} = -0.49$ e Å⁻³

Table 1

Selected bond lengths (Å).

N1—C1	1.375 (3)	N2—C21	1.373 (3)
N1—C8	1.411 (4)	N2—C28	1.401 (3)
N1—C9	1.425 (4)	N2—C29	1.467 (4)
C1—C2	1.545 (4)	C21—C22	1.558 (4)

Table 2

Hydrogen-bond geometry (Å, °).

$D-H\cdots A$	$D-H$	$H\cdots A$	$D\cdots A$	$D-H\cdots A$
$C7-H7\cdots O3^i$	0.95 (3)	2.44 (3)	3.367 (5)	165 (2)

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

Data collection: APEX2 (Bruker, 2005); cell refinement: SAINT (Bruker, 2005); data reduction: SAINT; program(s) used to solve structure: SHELXL97 (Sheldrick, 2008); program(s) used to refine structure: SHELXL97 (Sheldrick, 2008); molecular graphics: SHELXTL (Sheldrick, 2008); software used to prepare material for publication: WinGX (Farrugia, 1999).

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Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: CS2100).

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supporting information

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6-Bromo-1-butylindoline-2,3-dione

Lei Ji, Qi Fang and Jian-dong Fan

S1. Comment

Isatin, 2,3-indolinedione is traditionally obtained from oxidation of oxindole and indigo blue. Its derivatives have long been used as precursors of medicines and reductive dyes (Zhou *et al.*, 2006; Pirrung *et al.*, 2005; Kopka *et al.*, 2006). The first crystal structure report on isatin was by Goldschmidt and Llewellyn, 1950. Here we report the crystal structure of *N*-butyl-6-bromoisatin. There are two independent molecules in the asymmetric unit with very similar bond parameters. Except for the fused aromatic bond of C3=C8 in one molecule and C23=C28 bond in the other molecule, four other bonds in the five-membered heterocyclic rings can be classified as single C—C bonds and π -conjugated C=N bonds. The C1—C2 (1.545 (4) Å) and C21—C22 (1.558 (4) Å) bond lengths are longer than expected for a C(sp^2)-C(sp^2) single bond. This may be the result of the repulsion of the lone pair electrons of the two oxygen atoms in *cis*-diones (Palenik *et al.*, 1990). Nitrogen atoms and their three bonded carbons are perfectly co-planar, showing that the sp^2 N atoms allocate two p_z electrons for π -bonding. The bond lengths of the π -conjugated N=C bonds (c.f. Table 1) are much shorter than a single C—N bond length. Molecules are packed in dimers (Figure 2). One kind of the independent molecules are linked through a symmetry center to form dimer stack A while two other molecules form another dimer stack B. Mean planes of A and B dimers have a dihedral angle of 85.8 (1)°. The phenyl-phenyl spacing (3.40 (1) Å) in A dimers is considerably shorter than the spacing (3.65 (1) Å) in the B dimers, indicating relatively stronger intermolecular π - π interactions between these A molecules. The intermolecular interactions in dimer A are further strengthened by C—Br short contacts. The C1 \cdots Br1[-x+1, -y+1, -z] and C2 \cdots Br1[-x+1, -y+1, -z] in dimer A are 3.476 (3) Å and 3.538 (3) Å, respectively. By comparison, the corresponding C \cdots Br contacts in dimer B are much longer with the shortest distance being 3.672 (4) Å for C22 \cdots Br2[-x+1, -y+1, -z]. It may be the result of these different C \cdots Br contacts that the C1—C2 bond length in dimer A is marginally shorter than the C21—C22 bond in dimer B. Amongst intermolecular C—H \cdots O hydrogen bonds the strongest is the C7—H7 \cdots O3 one between dimers A and B (see Table 2).

S2. Experimental

6-Bromoisatin (10.5 g) was dissolved in 100 ml DMSO in a three-necked flask. Then KI (3.50 g), cetyltrimethyl ammonium bromide (1.00 g), and KOH (30.0 g) in 20.0 g water were added. Then 40.0 ml of *n*-bromobutane was added drop-wise into the above mixture with stirring. The mixture was stirred at 343 K for 2 days under nitrogen protection. The reaction mixture was washed by water, extracted with CHCl₃, then the chloroform layer was dried by Na₂SO₄. After vaporizing the solvent, the crude product was purified by column chromatography, resulting in 9.5 g (yield 71%) *N*-butyl-6-bromoisatin product. The compound was dissolved in chloroform again. On most of the solvent evaporating at room temperature orange lump like crystals were formed.

S3. Refinement

All H atoms except those on the methyl groups were initially found in difference electron density syntheses and were used in the least-squares refinement. Six H atoms on two terminal methyls could also be located in the difference maps but some H—C bond parameters in the methyl groups became unreasonable. So HFIX 137 instructions were used to restrain the methyl H-positions.

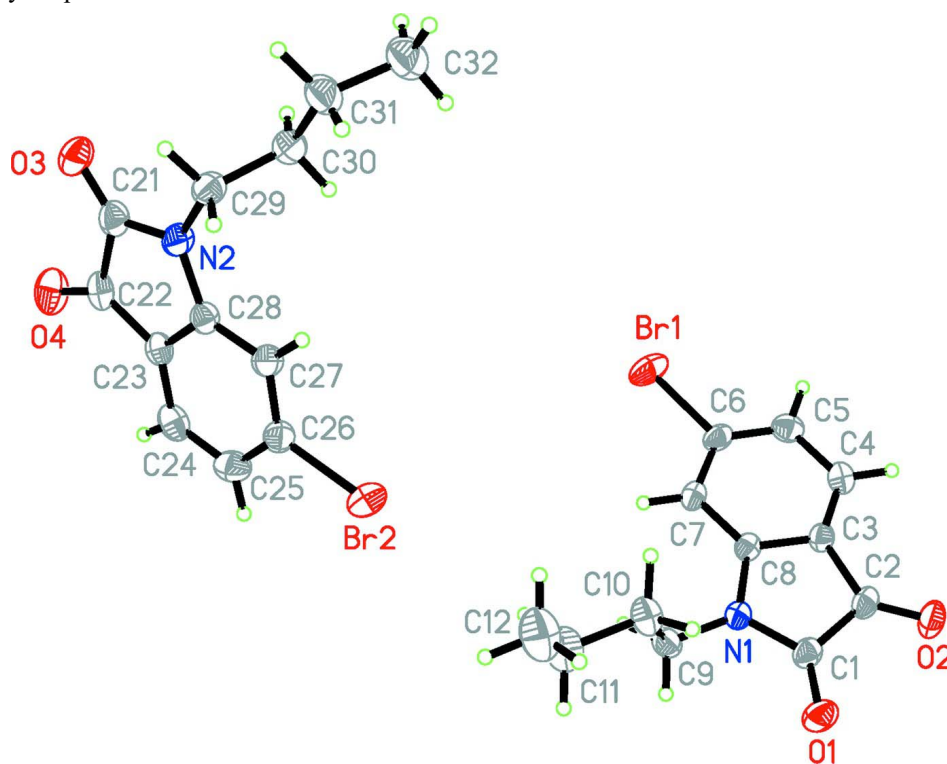


Figure 1

Two *N*-butyl-6-bromoisatin molecules in the asymmetric unit with 30% probability displacement ellipsoids

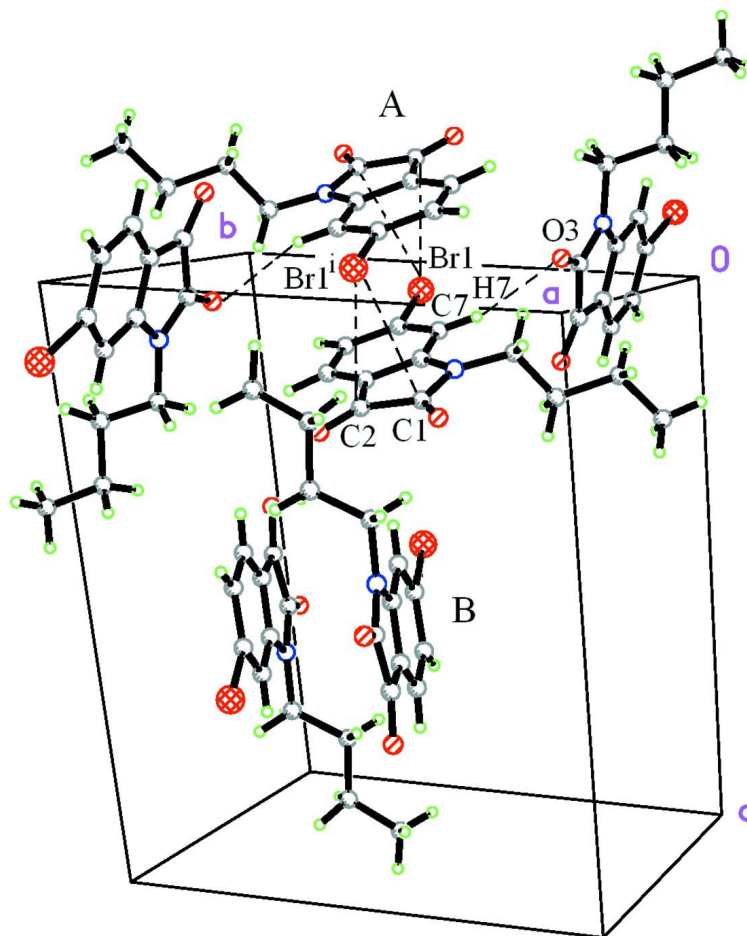


Figure 2

The molecular packing of the crystal showing the dimer structure [symmetry code: (i)-x+1, -y+1, -z]

6-Bromo-1-butylindoline-2,3-dione

Crystal data

$C_{12}H_{12}BrNO_2$

$M_r = 282.14$

Monoclinic, $P2_1/c$

Hall symbol: -P 2ybc

$a = 13.3097$ (2) Å

$b = 11.8793$ (2) Å

$c = 16.2238$ (2) Å

$\beta = 112.340$ (1)°

$V = 2372.62$ (6) Å³

$Z = 8$

$F(000) = 1136$

$D_x = 1.580$ Mg m⁻³

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

Cell parameters from 3677 reflections

$\theta = 2.2$ – 23.8 °

$\mu = 3.45$ mm⁻¹

$T = 296$ K

Plank, orange

$0.37 \times 0.13 \times 0.11$ mm

Data collection

Bruker APEXII CCD
diffractometer

Radiation source: fine-focus sealed tube

Graphite monochromator

Detector resolution: 10.0 pixels mm⁻¹

phi and ω scans

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

$T_{\min} = 0.526$, $T_{\max} = 0.744$

16713 measured reflections

5170 independent reflections

2424 reflections with $I > 2\sigma(I)$

$R_{\text{int}} = 0.076$
 $\theta_{\text{max}} = 27.0^\circ$, $\theta_{\text{min}} = 1.7^\circ$
 $h = -16 \rightarrow 17$

$k = -15 \rightarrow 10$
 $l = -20 \rightarrow 20$

Refinement

Refinement on F^2
 Least-squares matrix: full
 $R[F^2 > 2\sigma(F^2)] = 0.039$
 $wR(F^2) = 0.070$
 $S = 0.93$
 5170 reflections
 364 parameters
 1 restraint
 Primary atom site location: structure-invariant
 direct methods
 Secondary atom site location: difference Fourier
 map

Hydrogen site location: difference Fourier map
 H atoms treated by a mixture of independent
 and constrained refinement
 $w = 1/[\sigma^2(F_o^2) + (0.0158P)^2]$
 where $P = (F_o^2 + 2F_c^2)/3$
 $(\Delta/\sigma)_{\text{max}} = 0.001$
 $\Delta\rho_{\text{max}} = 0.39 \text{ e } \text{Å}^{-3}$
 $\Delta\rho_{\text{min}} = -0.49 \text{ e } \text{Å}^{-3}$
 Extinction correction: *SHELXL97* (Sheldrick,
 2008), $F_c^* = kF_c[1 + 0.001x F_c^2 \lambda^3 / \sin(2\theta)]^{-1/4}$
 Extinction coefficient: 0.00147 (12)

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

	<i>x</i>	<i>y</i>	<i>z</i>	$U_{\text{iso}}^*/U_{\text{eq}}$
Br1	0.25474 (2)	0.52247 (4)	0.03130 (3)	0.07546 (17)
O1	0.82928 (15)	0.3064 (2)	0.18511 (13)	0.0649 (7)
O2	0.82426 (16)	0.5497 (2)	0.22161 (15)	0.0682 (7)
N1	0.64300 (18)	0.3303 (3)	0.12570 (15)	0.0448 (7)
C1	0.7490 (2)	0.3648 (3)	0.1675 (2)	0.0477 (9)
C2	0.7445 (2)	0.4919 (3)	0.18633 (19)	0.0475 (9)
C3	0.6286 (2)	0.5203 (3)	0.15321 (19)	0.0431 (9)
C4	0.5754 (3)	0.6195 (4)	0.1517 (2)	0.0519 (10)
H4	0.616 (2)	0.688 (3)	0.1810 (18)	0.058 (10)*
C5	0.4628 (3)	0.6193 (4)	0.1150 (2)	0.0535 (10)
H5	0.4240 (19)	0.690 (3)	0.1167 (16)	0.047 (9)*
C6	0.4095 (2)	0.5211 (4)	0.0815 (2)	0.0476 (9)
C7	0.4595 (2)	0.4196 (4)	0.0807 (2)	0.0435 (9)
H7	0.4227 (18)	0.351 (2)	0.0576 (16)	0.036 (9)*
C8	0.5720 (2)	0.4213 (3)	0.11826 (18)	0.0382 (8)
C9	0.6102 (3)	0.2182 (3)	0.0965 (3)	0.0504 (10)
H9A	0.538 (2)	0.218 (3)	0.040 (2)	0.073 (10)*
H9B	0.659 (2)	0.185 (3)	0.081 (2)	0.078 (13)*
C10	0.5921 (3)	0.1458 (4)	0.1651 (3)	0.0640 (11)
H10B	0.537 (2)	0.185 (2)	0.1859 (17)	0.052 (9)*
H10A	0.658 (2)	0.147 (3)	0.221 (2)	0.070 (11)*
C11	0.5557 (4)	0.0275 (4)	0.1272 (3)	0.0831 (14)
H11A	0.613 (3)	-0.012 (3)	0.121 (2)	0.100*
H11B	0.477 (3)	0.032 (3)	0.085 (2)	0.100*
C12	0.5496 (3)	-0.0541 (4)	0.1927 (3)	0.1061 (15)
H12A	0.6201	-0.0628	0.2392	0.138 (10)*
H12B	0.5253	-0.1253	0.1641	0.138 (10)*
H12C	0.4993	-0.0276	0.2179	0.138 (10)*
Br2	0.25658 (2)	-0.10901 (4)	0.09621 (3)	0.07726 (17)

O3	-0.32733 (16)	-0.1957 (2)	0.03877 (14)	0.0728 (8)
O4	-0.30063 (18)	-0.2190 (2)	-0.13304 (15)	0.0795 (8)
N2	-0.14314 (18)	-0.1581 (2)	0.08915 (16)	0.0493 (7)
C21	-0.2438 (3)	-0.1837 (3)	0.0266 (2)	0.0550 (10)
C22	-0.2278 (3)	-0.1969 (3)	-0.0632 (2)	0.0538 (9)
C23	-0.1117 (2)	-0.1769 (3)	-0.0404 (2)	0.0446 (8)
C24	-0.0483 (3)	-0.1808 (3)	-0.0900 (3)	0.0552 (10)
H24	-0.084 (2)	-0.193 (3)	-0.1480 (19)	0.060 (11)*
C25	0.0612 (3)	-0.1618 (3)	-0.0495 (3)	0.0599 (10)
H25	0.1067 (18)	-0.164 (2)	-0.0805 (16)	0.035 (8)*
C26	0.1047 (2)	-0.1379 (3)	0.0399 (2)	0.0513 (9)
C27	0.0439 (2)	-0.1332 (3)	0.0925 (2)	0.0501 (10)
H27	0.0727 (17)	-0.120 (2)	0.1503 (16)	0.034 (9)*
C28	-0.0653 (2)	-0.1544 (3)	0.0505 (2)	0.0443 (8)
C29	-0.1191 (3)	-0.1497 (4)	0.1850 (2)	0.0545 (11)
H29A	-0.1731 (19)	-0.188 (2)	0.1991 (16)	0.045 (9)*
H29B	-0.056 (2)	-0.198 (3)	0.2144 (19)	0.068 (11)*
C30	-0.1031 (3)	-0.0315 (4)	0.2208 (3)	0.0592 (11)
H30B	-0.049 (2)	0.010 (3)	0.200 (2)	0.077 (11)*
H30A	-0.162 (2)	0.016 (3)	0.195 (2)	0.067 (12)*
C31	-0.0728 (4)	-0.0318 (4)	0.3219 (3)	0.0771 (13)
H31B	-0.002 (3)	-0.071 (3)	0.355 (2)	0.112 (17)*
H31A	-0.133 (3)	-0.073 (3)	0.333 (2)	0.103 (14)*
C32	-0.0463 (3)	0.0819 (4)	0.3617 (2)	0.0941 (14)
H32A	0.0122	0.1132	0.3483	0.129 (9)*
H32B	-0.0249	0.0766	0.4252	0.129 (9)*
H32C	-0.1090	0.1295	0.3376	0.129 (9)*

Atomic displacement parameters (\AA^2)

	U^{11}	U^{22}	U^{33}	U^{12}	U^{13}	U^{23}
Br1	0.03753 (19)	0.1118 (4)	0.0741 (3)	0.0122 (2)	0.01790 (17)	0.0168 (3)
O1	0.0405 (12)	0.082 (2)	0.0631 (16)	0.0147 (13)	0.0090 (11)	-0.0007 (14)
O2	0.0450 (13)	0.081 (2)	0.0662 (16)	-0.0181 (13)	0.0076 (12)	-0.0127 (15)
N1	0.0369 (14)	0.046 (2)	0.0473 (17)	-0.0027 (15)	0.0110 (12)	-0.0079 (17)
C1	0.0397 (19)	0.060 (3)	0.039 (2)	-0.0006 (19)	0.0107 (16)	0.001 (2)
C2	0.0372 (17)	0.062 (3)	0.038 (2)	-0.0066 (19)	0.0086 (15)	0.003 (2)
C3	0.0376 (17)	0.049 (3)	0.038 (2)	-0.0040 (19)	0.0091 (15)	-0.005 (2)
C4	0.056 (2)	0.045 (3)	0.051 (2)	-0.007 (2)	0.0165 (18)	-0.006 (2)
C5	0.053 (2)	0.051 (3)	0.056 (2)	0.010 (2)	0.0202 (19)	0.002 (2)
C6	0.0317 (16)	0.069 (3)	0.039 (2)	0.006 (2)	0.0095 (15)	0.005 (2)
C7	0.0361 (18)	0.049 (3)	0.042 (2)	-0.008 (2)	0.0114 (16)	-0.003 (2)
C8	0.0375 (17)	0.044 (2)	0.0305 (18)	0.0009 (18)	0.0100 (14)	0.0031 (18)
C9	0.052 (2)	0.043 (3)	0.053 (3)	0.005 (2)	0.016 (2)	-0.007 (2)
C10	0.072 (3)	0.050 (3)	0.064 (3)	-0.009 (2)	0.019 (2)	-0.004 (3)
C11	0.087 (3)	0.075 (4)	0.077 (3)	-0.005 (3)	0.020 (3)	0.003 (3)
C12	0.121 (4)	0.080 (4)	0.107 (4)	-0.018 (3)	0.032 (3)	-0.006 (4)
Br2	0.0453 (2)	0.0909 (4)	0.0969 (3)	0.0008 (2)	0.0285 (2)	-0.0097 (3)

O3	0.0427 (12)	0.090 (2)	0.0790 (17)	-0.0143 (13)	0.0160 (12)	-0.0119 (15)
O4	0.0672 (15)	0.094 (2)	0.0514 (16)	-0.0202 (15)	-0.0066 (13)	-0.0106 (16)
N2	0.0385 (14)	0.070 (2)	0.0342 (16)	-0.0056 (14)	0.0078 (12)	-0.0067 (16)
C21	0.048 (2)	0.059 (3)	0.049 (2)	-0.008 (2)	0.0089 (18)	-0.010 (2)
C22	0.055 (2)	0.042 (3)	0.052 (2)	-0.0093 (19)	0.0065 (18)	-0.004 (2)
C23	0.0489 (19)	0.043 (2)	0.036 (2)	-0.0037 (17)	0.0100 (17)	-0.0023 (19)
C24	0.072 (3)	0.050 (3)	0.036 (2)	-0.003 (2)	0.012 (2)	-0.003 (2)
C25	0.069 (3)	0.063 (3)	0.059 (3)	0.007 (2)	0.036 (2)	0.003 (2)
C26	0.0444 (18)	0.050 (3)	0.057 (2)	0.0006 (17)	0.0175 (18)	-0.001 (2)
C27	0.0423 (19)	0.064 (3)	0.038 (2)	0.0006 (18)	0.0095 (18)	-0.006 (2)
C28	0.0395 (17)	0.050 (2)	0.038 (2)	-0.0023 (16)	0.0089 (16)	-0.0028 (19)
C29	0.048 (2)	0.070 (3)	0.048 (2)	-0.007 (2)	0.0202 (19)	-0.004 (2)
C30	0.057 (2)	0.064 (3)	0.058 (3)	-0.006 (2)	0.023 (2)	-0.008 (3)
C31	0.103 (4)	0.076 (4)	0.065 (3)	-0.021 (3)	0.046 (3)	-0.020 (3)
C32	0.116 (3)	0.101 (4)	0.076 (3)	-0.027 (3)	0.048 (2)	-0.029 (3)

Geometric parameters (Å, °)

Br1—C6	1.906 (3)	Br2—C26	1.907 (3)
O1—C1	1.214 (3)	O3—C21	1.209 (3)
O2—C2	1.210 (3)	O4—C22	1.208 (3)
N1—C1	1.375 (3)	N2—C21	1.373 (3)
N1—C8	1.411 (4)	N2—C28	1.401 (3)
N1—C9	1.425 (4)	N2—C29	1.467 (4)
C1—C2	1.545 (4)	C21—C22	1.558 (4)
C2—C3	1.468 (4)	C22—C23	1.465 (4)
C3—C4	1.370 (4)	C23—C24	1.370 (4)
C3—C8	1.395 (4)	C23—C28	1.391 (4)
C4—C5	1.387 (4)	C24—C25	1.371 (4)
C4—H4	0.99 (3)	C24—H24	0.89 (3)
C5—C6	1.366 (4)	C25—C26	1.372 (4)
C5—H5	1.00 (3)	C25—H25	0.92 (2)
C6—C7	1.379 (4)	C26—C27	1.382 (4)
C7—C8	1.386 (4)	C27—C28	1.374 (4)
C7—H7	0.95 (3)	C27—H27	0.88 (2)
C9—C10	1.496 (5)	C29—C30	1.503 (5)
C9—H9A	1.05 (3)	C29—H29A	0.95 (3)
C9—H9B	0.87 (3)	C29—H29B	0.98 (3)
C10—C11	1.536 (6)	C30—C31	1.532 (5)
C10—H10B	1.03 (3)	C30—H30B	1.04 (3)
C10—H10A	0.99 (3)	C30—H30A	0.93 (3)
C11—C12	1.464 (6)	C31—C32	1.481 (5)
C11—H11A	0.94 (4)	C31—H31B	1.00 (4)
C11—H11B	1.01 (3)	C31—H31A	1.01 (4)
C12—H12A	0.9600	C32—H32A	0.9600
C12—H12B	0.9600	C32—H32B	0.9600
C12—H12C	0.9600	C32—H32C	0.9600

C1—N1—C8	110.0 (3)	C21—N2—C28	110.9 (2)
C1—N1—C9	124.8 (3)	C21—N2—C29	123.9 (3)
C8—N1—C9	125.3 (3)	C28—N2—C29	124.9 (2)
O1—C1—N1	126.3 (3)	O3—C21—N2	127.2 (3)
O1—C1—C2	127.4 (3)	O3—C21—C22	127.0 (3)
N1—C1—C2	106.3 (3)	N2—C21—C22	105.7 (3)
O2—C2—C3	130.8 (4)	O4—C22—C23	131.7 (3)
O2—C2—C1	123.8 (3)	O4—C22—C21	123.5 (3)
C3—C2—C1	105.5 (3)	C23—C22—C21	104.7 (3)
C4—C3—C8	121.5 (3)	C24—C23—C28	120.3 (3)
C4—C3—C2	132.0 (3)	C24—C23—C22	132.2 (3)
C8—C3—C2	106.5 (3)	C28—C23—C22	107.4 (3)
C3—C4—C5	118.2 (4)	C23—C24—C25	119.6 (3)
C3—C4—H4	120.7 (16)	C23—C24—H24	115.6 (18)
C5—C4—H4	120.8 (16)	C25—C24—H24	124.7 (19)
C6—C5—C4	119.1 (4)	C24—C25—C26	119.0 (3)
C6—C5—H5	122.6 (15)	C24—C25—H25	122.3 (15)
C4—C5—H5	118.3 (15)	C26—C25—H25	118.7 (15)
C5—C6—C7	124.8 (3)	C25—C26—C27	123.4 (3)
C5—C6—Br1	118.4 (3)	C25—C26—Br2	119.2 (3)
C7—C6—Br1	116.8 (3)	C27—C26—Br2	117.4 (3)
C6—C7—C8	115.3 (3)	C28—C27—C26	116.3 (3)
C6—C7—H7	125.2 (15)	C28—C27—H27	120.7 (16)
C8—C7—H7	119.5 (15)	C26—C27—H27	123.0 (16)
C7—C8—C3	121.1 (3)	C27—C28—C23	121.4 (3)
C7—C8—N1	127.1 (3)	C27—C28—N2	127.3 (3)
C3—C8—N1	111.8 (3)	C23—C28—N2	111.3 (2)
N1—C9—C10	113.9 (3)	N2—C29—C30	114.5 (3)
N1—C9—H9A	110.8 (17)	N2—C29—H29A	110.0 (15)
C10—C9—H9A	107.0 (17)	C30—C29—H29A	111.5 (17)
N1—C9—H9B	111 (2)	N2—C29—H29B	106.1 (17)
C10—C9—H9B	108 (2)	C30—C29—H29B	112.8 (18)
H9A—C9—H9B	106 (3)	H29A—C29—H29B	101 (2)
C9—C10—C11	110.1 (4)	C29—C30—C31	110.6 (4)
C9—C10—H10B	108.8 (16)	C29—C30—H30B	109.3 (18)
C11—C10—H10B	112.8 (16)	C31—C30—H30B	115.0 (17)
C9—C10—H10A	109.7 (18)	C29—C30—H30A	115 (2)
C11—C10—H10A	114.1 (19)	C31—C30—H30A	109 (2)
H10B—C10—H10A	101 (2)	H30B—C30—H30A	98 (3)
C12—C11—C10	114.3 (4)	C32—C31—C30	112.8 (4)
C12—C11—H11A	91 (3)	C32—C31—H31B	100 (2)
C10—C11—H11A	111 (2)	C30—C31—H31B	113 (2)
C12—C11—H11B	101 (2)	C32—C31—H31A	116 (2)
C10—C11—H11B	108 (2)	C30—C31—H31A	107 (2)
H11A—C11—H11B	129 (3)	H31B—C31—H31A	109 (3)
C11—C12—H12A	109.5	C31—C32—H32A	109.5
C11—C12—H12B	109.5	C31—C32—H32B	109.5
H12A—C12—H12B	109.5	H32A—C32—H32B	109.5

C11—C12—H12C	109.5	C31—C32—H32C	109.5
H12A—C12—H12C	109.5	H32A—C32—H32C	109.5
H12B—C12—H12C	109.5	H32B—C32—H32C	109.5
C8—N1—C1—O1	179.0 (3)	C28—N2—C21—O3	178.2 (4)
C9—N1—C1—O1	0.6 (5)	C29—N2—C21—O3	5.0 (6)
C8—N1—C1—C2	-1.6 (3)	C28—N2—C21—C22	-0.5 (4)
C9—N1—C1—C2	180.0 (3)	C29—N2—C21—C22	-173.6 (3)
O1—C1—C2—O2	0.6 (5)	O3—C21—C22—O4	1.8 (6)
N1—C1—C2—O2	-178.8 (3)	N2—C21—C22—O4	-179.6 (3)
O1—C1—C2—C3	-179.3 (3)	O3—C21—C22—C23	-178.2 (3)
N1—C1—C2—C3	1.3 (3)	N2—C21—C22—C23	0.4 (3)
O2—C2—C3—C4	0.2 (6)	O4—C22—C23—C24	-2.7 (7)
C1—C2—C3—C4	-180.0 (3)	C21—C22—C23—C24	177.3 (4)
O2—C2—C3—C8	179.6 (3)	O4—C22—C23—C28	179.8 (4)
C1—C2—C3—C8	-0.5 (3)	C21—C22—C23—C28	-0.2 (4)
C8—C3—C4—C5	0.7 (5)	C28—C23—C24—C25	-0.3 (5)
C2—C3—C4—C5	180.0 (3)	C22—C23—C24—C25	-177.6 (4)
C3—C4—C5—C6	-0.5 (5)	C23—C24—C25—C26	-0.8 (5)
C4—C5—C6—C7	-0.2 (5)	C24—C25—C26—C27	0.8 (6)
C4—C5—C6—Br1	-179.9 (2)	C24—C25—C26—Br2	-179.0 (3)
C5—C6—C7—C8	0.8 (5)	C25—C26—C27—C28	0.2 (5)
Br1—C6—C7—C8	-179.6 (2)	Br2—C26—C27—C28	-180.0 (2)
C6—C7—C8—C3	-0.6 (4)	C26—C27—C28—C23	-1.3 (5)
C6—C7—C8—N1	-179.6 (3)	C26—C27—C28—N2	177.9 (3)
C4—C3—C8—C7	-0.1 (5)	C24—C23—C28—C27	1.4 (5)
C2—C3—C8—C7	-179.6 (3)	C22—C23—C28—C27	179.3 (3)
C4—C3—C8—N1	179.1 (3)	C24—C23—C28—N2	-177.9 (3)
C2—C3—C8—N1	-0.4 (3)	C22—C23—C28—N2	0.0 (4)
C1—N1—C8—C7	-179.6 (3)	C21—N2—C28—C27	-179.0 (3)
C9—N1—C8—C7	-1.2 (5)	C29—N2—C28—C27	-5.9 (5)
C1—N1—C8—C3	1.3 (3)	C21—N2—C28—C23	0.3 (4)
C9—N1—C8—C3	179.7 (3)	C29—N2—C28—C23	173.4 (3)
C1—N1—C9—C10	89.8 (4)	C21—N2—C29—C30	-106.6 (4)
C8—N1—C9—C10	-88.3 (4)	C28—N2—C29—C30	81.2 (4)
N1—C9—C10—C11	179.3 (3)	N2—C29—C30—C31	-177.1 (3)
C9—C10—C11—C12	172.4 (4)	C29—C30—C31—C32	174.8 (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>
C7—H7 \cdots O3 ⁱ	0.95 (3)	2.44 (3)	3.367 (5)	165 (2)

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