

Terephthalic acid–2,2'-dimethyl-1,1'-(butane-1,4-diyl)dibenzimidazole (2/3)

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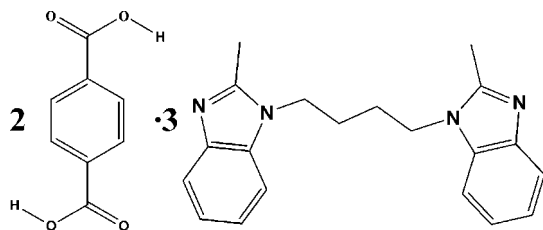
Received 22 November 2007; accepted 8 April 2008

 Key indicators: single-crystal X-ray study; $T = 293$ K; mean $\sigma(\text{C}-\text{C}) = 0.004$ Å; R factor = 0.071; wR factor = 0.164; data-to-parameter ratio = 17.3.

In the crystal structure of the title compound, $2\text{C}_8\text{H}_6\text{O}_4 \cdot 3\text{C}_{20}\text{H}_{22}\text{N}_4$, there are three independent 2,2'-dimethyl-1,1'-(butane-1,4-diyl)dibenzimidazole molecules, each of which lies on an inversion centre. The terephthalic acid molecules are linked to adjacent 2,2'-dimethyl-1,1'-(butane-1,4-diyl)dibenzimidazole molecules *via* $\text{O}-\text{H} \cdots \text{N}$ hydrogen bonds.

Related literature

For related literature, see: Dale & Elsegood (2004); Ma *et al.* (2000).



Experimental

Crystal data

 $2\text{C}_8\text{H}_6\text{O}_4 \cdot 3\text{C}_{20}\text{H}_{22}\text{N}_4$
 $M_r = 1286.88$

 Triclinic, $P\bar{1}$
 $a = 9.225$ (5) Å

 $b = 12.298$ (8) Å

 $c = 14.883$ (7) Å

 $\alpha = 87.17$ (2)°

 $\beta = 87.768$ (16)°

 $\gamma = 89.33$ (2)°

 $V = 1685.0$ (16) Å³
 $Z = 1$

 Mo $K\alpha$ radiation

 $\mu = 0.08$ mm⁻¹
 $T = 293$ (2) K

 $0.49 \times 0.38 \times 0.26$ mm

Data collection

Rigaku R-Axis RAPID diffractometer

 Absorption correction: multi-scan (*ABSCOR*; Higashi, 1995)

 $T_{\min} = 0.963$, $T_{\max} = 0.977$

16449 measured reflections

7573 independent reflections

 3739 reflections with $I > 2\sigma(I)$
 $R_{\text{int}} = 0.055$

Refinement

 $R[F^2 > 2\sigma(F^2)] = 0.071$
 $wR(F^2) = 0.163$
 $S = 1.01$

7573 reflections

437 parameters

H-atom parameters constrained

 $\Delta\rho_{\max} = 0.23$ e Å⁻³
 $\Delta\rho_{\min} = -0.21$ e Å⁻³
Table 1

Hydrogen-bond geometry (Å, °).

$D-H \cdots A$	$D-H$	$H \cdots A$	$D \cdots A$	$D-H \cdots A$
$\text{O1}-\text{H1} \cdots \text{N1}^{\text{i}}$	0.82	1.83	2.609 (3)	157
$\text{O3}-\text{H3A} \cdots \text{N3}^{\text{ii}}$	0.82	1.87	2.668 (3)	165

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

Data collection: *PROCESS-AUTO* (Rigaku, 1998); cell refinement: *PROCESS-AUTO*; data reduction: *PROCESS-AUTO*; program(s) used to solve structure: *SHELXS97* (Sheldrick, 2008); program(s) used to refine structure: *SHELXL97* (Sheldrick, 2008); molecular graphics: *SHELXTL-Plus* (Sheldrick, 2008); software used to prepare material for publication: *SHELXL97*.

The authors thank the Analysis and Testing Foundation of Northeast Normal University for support.

Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: CS2066).

References

- Dale, S. H. & Elsegood, M. R. J. (2004). *Acta Cryst.* **C60**, o444–o448.
 Higashi, T. (1995). *ABSCOR*. Rigaku Corporation, Tokyo, Japan.
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 Rigaku (1998). *PROCESS-AUTO*. Rigaku Corporation, Tokyo, Japan.
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supplementary materials

Acta Cryst. (2008). E64, o838 [doi:10.1107/S1600536808009665]

Terephthalic acid-2,2'-dimethyl-1,1'-(butane-1,4-diyl)dibenzimidazole (2/3)

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Comment

The title compound is composed of terephthalic acid and 2,2'-dimethyl-1,1'-(butane-1,4-diyl)dibenzimidazole (Fig. 1). The bond distances and angles are normal (Dale & Elsegood, 2004; Ma *et al.*, 2000). There are three independent 2,2'-dimethyl-1,1'-(butane-1,4-diyl)dibenzimidazole molecules, which extend in different directions, sitting around independent symmetry centres. The terephthalic acids are hydrogen bonded to the 2,2'-dimethyl-1,1'-(butane-1,4-diyl)dibenzimidazole (Table 1) *via* O—H \cdots N hydrogen bonds into the chains (Fig. 2).

Experimental

A mixture of terephthalic acid (0.033 g, 0.2 mmol), 2,2'-dimethyl-1,1'-(butane-1,4-diyl)dibenzimidazole (0.064 g, 0.2 mmol) in 1:4 (*v/v*) EtOH/H₂O (10 ml) was placed in a Teflon-lined autoclave. The mixture was heated at 430 K for 1 days, and then it was gradually cooled to room temperature at a rate of 10 K.h⁻¹. Colorless single crystals of the title compound were obtained.

Refinement

All H-atoms bound to carbon were refined using a riding model with $d(\text{C—H}) = 0.93 \text{ \AA}$, $U_{\text{iso}} = 1.2U_{\text{eq}}(\text{C})$ for aromatic, 0.97 \AA , $U_{\text{iso}} = 1.5U_{\text{eq}}(\text{C})$ for CH₂ atoms and 0.96 \AA , $U_{\text{iso}} = 1.5U_{\text{eq}}(\text{C})$ for CH₃ atoms. Hydroxyl H atoms were refined using a riding model with $d(\text{O—H}) = 0.82 \text{ \AA}$, $U_{\text{iso}} = 1.5U_{\text{eq}}(\text{O})$.

Figures

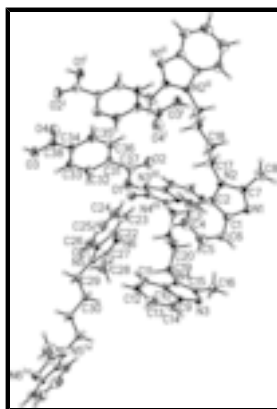


Fig. 1. A view of the title clathrate. The displacement ellipsoids are drawn at the 30% probability level. (i) $-x, -y, -z + 1$. (ii) $-x, -y + 1, -z + 1$. (iii) $-x + 1, -y + 1, -z$. (iv) $-x + 2, -y, -z$.

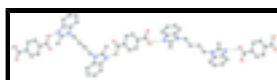


Fig. 2. A one-dimensional chain linked by hydrogen bonds. The H-bonds are shown as dashed lines.

Terephthalic acid–2,2'-dimethyl-1,1'-(butane-1,4-diyl)dibenzimidazole (2/3)

Crystal data

$2\text{C}_8\text{H}_6\text{O}_4 \cdot 3\text{C}_{20}\text{H}_{22}\text{N}_4$	$Z = 1$
$M_r = 1286.88$	$F_{000} = 682$
Triclinic, $P\bar{1}$	$D_x = 1.268 \text{ Mg m}^{-3}$
Hall symbol: -P 1	Mo $K\alpha$ radiation
$a = 9.225 (5) \text{ \AA}$	$\lambda = 0.71069 \text{ \AA}$
$b = 12.298 (8) \text{ \AA}$	Cell parameters from 7573 reflections
$c = 14.883 (7) \text{ \AA}$	$\theta = 3.0\text{--}27.5^\circ$
$\alpha = 87.17 (2)^\circ$	$\mu = 0.08 \text{ mm}^{-1}$
$\beta = 87.768 (16)^\circ$	$T = 293 (2) \text{ K}$
$\gamma = 89.33 (2)^\circ$	Plate, colourless
$V = 1685.0 (16) \text{ \AA}^3$	$0.49 \times 0.38 \times 0.26 \text{ mm}$

Data collection

Rigaku R-Axis RAPID diffractometer	7573 independent reflections
Radiation source: rotation-anode tube	3739 reflections with $I > 2\sigma(I)$
Monochromator: graphite	$R_{\text{int}} = 0.055$
Detector resolution: $10.0 \text{ pixels mm}^{-1}$	$\theta_{\text{max}} = 27.5^\circ$
$T = 293(2) \text{ K}$	$\theta_{\text{min}} = 3.0^\circ$
ω scans	$h = -11 \rightarrow 10$
Absorption correction: multi-scan (ABSCOR; Higashi, 1995)	$k = -15 \rightarrow 15$
$T_{\text{min}} = 0.963$, $T_{\text{max}} = 0.977$	$l = -19 \rightarrow 19$
16449 measured reflections	

Refinement

Refinement on F^2	Hydrogen site location: inferred from neighbouring sites
Least-squares matrix: full	H-atom parameters constrained
$R[F^2 > 2\sigma(F^2)] = 0.071$	$w = 1/[\sigma^2(F_o^2) + (0.0674P)^2 + 0.1144P]$
$wR(F^2) = 0.163$	where $P = (F_o^2 + 2F_c^2)/3$
$S = 1.01$	$(\Delta/\sigma)_{\text{max}} < 0.001$
7573 reflections	$\Delta\rho_{\text{max}} = 0.23 \text{ e \AA}^{-3}$
437 parameters	$\Delta\rho_{\text{min}} = -0.21 \text{ e \AA}^{-3}$
Primary atom site location: structure-invariant direct methods	Extinction correction: SHELXL97 (Sheldrick, 2008), $F_c^* = kF_c[1 + 0.001x F_c^2 \lambda^3 / \sin(2\theta)]^{-1/4}$
Secondary atom site location: difference Fourier map	Extinction coefficient: 0.0101 (15)

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	0.4952 (3)	0.6452 (2)	0.39701 (16)	0.0443 (6)
C2	0.4004 (3)	0.5588 (2)	0.41249 (16)	0.0436 (6)
C3	0.4031 (3)	0.4701 (2)	0.35876 (18)	0.0565 (7)
H3	0.3405	0.4119	0.3701	0.068*
C4	0.5025 (4)	0.4726 (3)	0.2882 (2)	0.0728 (10)
H4	0.5074	0.4144	0.2505	0.087*
C5	0.5956 (4)	0.5585 (3)	0.2710 (2)	0.0713 (9)
H5	0.6609	0.5570	0.2218	0.086*
C6	0.5948 (3)	0.6460 (3)	0.32422 (18)	0.0592 (8)
H6	0.6581	0.7037	0.3123	0.071*
C7	0.3629 (3)	0.6799 (2)	0.51495 (18)	0.0481 (7)
C8	0.2974 (3)	0.7342 (3)	0.5937 (2)	0.0756 (10)
H8A	0.3136	0.6901	0.6475	0.113*
H8B	0.1949	0.7435	0.5864	0.113*
H8C	0.3410	0.8041	0.5982	0.113*
C9	0.9763 (2)	0.4712 (2)	0.16737 (16)	0.0383 (6)
C10	0.8777 (2)	0.4183 (2)	0.11664 (15)	0.0368 (6)
C11	0.8522 (3)	0.3085 (2)	0.12836 (18)	0.0521 (7)
H11	0.7845	0.2741	0.0951	0.063*
C12	0.9318 (3)	0.2525 (2)	0.1917 (2)	0.0611 (8)
H12	0.9170	0.1782	0.2021	0.073*
C13	1.0343 (3)	0.3039 (3)	0.24080 (18)	0.0621 (8)
H13	1.0883	0.2627	0.2819	0.074*
C14	1.0571 (3)	0.4128 (2)	0.22996 (17)	0.0529 (7)
H14	1.1247	0.4469	0.2635	0.063*
C15	0.8801 (2)	0.5945 (2)	0.08040 (16)	0.0416 (6)
C16	0.8408 (3)	0.7004 (2)	0.0362 (2)	0.0604 (8)
H16A	0.8748	0.7024	-0.0256	0.091*
H16B	0.8848	0.7580	0.0664	0.091*
H16C	0.7373	0.7096	0.0392	0.091*
C17	0.1997 (2)	0.5149 (2)	0.52713 (18)	0.0519 (7)
H17A	0.1880	0.5270	0.5909	0.062*
H17B	0.2251	0.4388	0.5207	0.062*
C18	0.0580 (2)	0.5394 (2)	0.48259 (17)	0.0511 (7)
H18A	0.0717	0.5344	0.4180	0.061*
H18B	0.0269	0.6132	0.4944	0.061*
C19	0.7073 (2)	0.4862 (2)	-0.00500 (16)	0.0469 (7)
H19A	0.7158	0.4139	-0.0281	0.056*
H19B	0.7226	0.5386	-0.0553	0.056*
C20	0.5564 (2)	0.5015 (2)	0.03514 (15)	0.0410 (6)
H20A	0.5369	0.4444	0.0812	0.049*
H20B	0.5502	0.5708	0.0635	0.049*
C21	0.6099 (3)	0.0780 (2)	0.09565 (19)	0.0499 (7)
C22	0.5086 (3)	0.1624 (2)	0.08891 (18)	0.0463 (6)
C23	0.4243 (3)	0.1903 (3)	0.1628 (2)	0.0667 (8)

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H23	0.3561	0.2464	0.1585	0.080*
C24	0.4434 (4)	0.1336 (4)	0.2417 (3)	0.0917 (12)
H24	0.3872	0.1516	0.2922	0.110*
C25	0.5439 (5)	0.0499 (4)	0.2497 (3)	0.0946 (13)
H25	0.5544	0.0133	0.3053	0.114*
C26	0.6290 (4)	0.0197 (3)	0.1766 (3)	0.0764 (10)
H26	0.6960	-0.0372	0.1814	0.092*
C27	0.6126 (3)	0.1528 (2)	-0.0416 (2)	0.0569 (7)
C28	0.6558 (4)	0.1745 (4)	-0.1375 (2)	0.0991 (13)
H28A	0.7523	0.2031	-0.1420	0.149*
H28B	0.5899	0.2266	-0.1641	0.149*
H28C	0.6531	0.1081	-0.1685	0.149*
C29	0.7922 (3)	-0.0014 (3)	-0.0138 (3)	0.0739 (10)
H29A	0.7951	-0.0081	-0.0785	0.089*
H29B	0.7731	-0.0729	0.0143	0.089*
C30	0.9373 (3)	0.0365 (2)	0.0140 (3)	0.0741 (10)
H30A	0.9350	0.0407	0.0790	0.089*
H30B	0.9541	0.1093	-0.0121	0.089*
C31	0.1777 (3)	0.0473 (2)	0.42478 (16)	0.0420 (6)
C32	0.2506 (3)	-0.0499 (2)	0.41235 (17)	0.0493 (7)
H32	0.3417	-0.0616	0.4358	0.059*
C33	0.1875 (3)	-0.1296 (2)	0.36497 (17)	0.0490 (7)
H33	0.2373	-0.1947	0.3563	0.059*
C34	0.0517 (3)	-0.1139 (2)	0.33031 (16)	0.0449 (6)
C35	-0.0224 (3)	-0.0178 (2)	0.34588 (18)	0.0504 (7)
H35	-0.1151	-0.0069	0.3245	0.060*
C36	0.0403 (3)	0.0617 (2)	0.39292 (17)	0.0507 (7)
H36	-0.0107	0.1258	0.4033	0.061*
C37	0.2471 (3)	0.1394 (2)	0.46745 (18)	0.0497 (7)
C38	-0.0130 (3)	-0.1938 (2)	0.27258 (19)	0.0518 (7)
N1	0.4693 (2)	0.72048 (18)	0.46235 (15)	0.0498 (6)
N2	0.3174 (2)	0.58270 (18)	0.48811 (13)	0.0442 (5)
N3	0.9751 (2)	0.58143 (17)	0.14404 (14)	0.0445 (5)
N4	0.81999 (19)	0.49945 (16)	0.05991 (13)	0.0393 (5)
N5	0.6745 (2)	0.07253 (18)	0.01097 (17)	0.0530 (6)
N6	0.5129 (2)	0.20832 (19)	0.00184 (15)	0.0567 (6)
O1	0.3721 (2)	0.11390 (16)	0.50186 (14)	0.0674 (6)
H1	0.4049	0.1675	0.5244	0.081*
O2	0.1950 (2)	0.22979 (17)	0.46676 (16)	0.0793 (7)
O3	0.0626 (2)	-0.28314 (16)	0.26517 (13)	0.0667 (6)
H3A	0.0214	-0.3237	0.2325	0.080*
O4	-0.1239 (2)	-0.17682 (18)	0.23391 (17)	0.0881 (8)

Atomic displacement parameters (\AA^2)

	U^{11}	U^{22}	U^{33}	U^{12}	U^{13}	U^{23}
C1	0.0354 (13)	0.0509 (17)	0.0473 (14)	-0.0004 (12)	-0.0090 (12)	-0.0023 (12)
C2	0.0352 (13)	0.0505 (17)	0.0462 (14)	-0.0008 (12)	-0.0111 (12)	-0.0054 (12)

C3	0.0593 (17)	0.0555 (19)	0.0567 (17)	-0.0048 (14)	-0.0188 (15)	-0.0092 (14)
C4	0.086 (2)	0.082 (3)	0.0530 (18)	0.008 (2)	-0.0140 (18)	-0.0234 (17)
C5	0.067 (2)	0.101 (3)	0.0458 (16)	0.007 (2)	0.0036 (15)	-0.0110 (18)
C6	0.0447 (15)	0.077 (2)	0.0550 (17)	-0.0084 (14)	-0.0038 (14)	0.0042 (16)
C7	0.0386 (14)	0.0508 (18)	0.0563 (16)	0.0061 (13)	-0.0086 (13)	-0.0128 (13)
C8	0.073 (2)	0.083 (3)	0.073 (2)	0.0173 (18)	-0.0005 (18)	-0.0272 (18)
C9	0.0306 (12)	0.0401 (15)	0.0448 (13)	0.0012 (10)	-0.0050 (11)	-0.0063 (11)
C10	0.0303 (11)	0.0393 (15)	0.0413 (13)	0.0020 (10)	-0.0022 (11)	-0.0058 (11)
C11	0.0502 (15)	0.0467 (18)	0.0603 (16)	-0.0047 (13)	-0.0004 (14)	-0.0116 (13)
C12	0.074 (2)	0.0376 (17)	0.0702 (19)	0.0023 (15)	0.0017 (17)	0.0047 (14)
C13	0.0659 (19)	0.064 (2)	0.0545 (17)	0.0139 (16)	-0.0068 (15)	0.0121 (15)
C14	0.0465 (15)	0.062 (2)	0.0512 (15)	0.0058 (14)	-0.0139 (13)	-0.0031 (14)
C15	0.0321 (12)	0.0407 (15)	0.0520 (14)	-0.0006 (11)	-0.0035 (12)	-0.0019 (12)
C16	0.0524 (16)	0.0495 (18)	0.079 (2)	0.0006 (13)	-0.0095 (16)	0.0099 (15)
C17	0.0347 (13)	0.067 (2)	0.0529 (15)	-0.0090 (13)	-0.0040 (12)	0.0059 (13)
C18	0.0348 (13)	0.0659 (19)	0.0522 (15)	-0.0069 (12)	-0.0052 (12)	0.0051 (13)
C19	0.0336 (13)	0.0634 (19)	0.0452 (14)	-0.0010 (12)	-0.0112 (12)	-0.0094 (12)
C20	0.0348 (12)	0.0491 (16)	0.0399 (13)	-0.0020 (11)	-0.0075 (11)	-0.0051 (11)
C21	0.0346 (13)	0.0450 (17)	0.0704 (19)	-0.0064 (11)	-0.0049 (14)	-0.0025 (14)
C22	0.0335 (13)	0.0458 (16)	0.0601 (17)	-0.0004 (11)	-0.0031 (13)	-0.0071 (13)
C23	0.0556 (18)	0.074 (2)	0.071 (2)	0.0005 (16)	0.0074 (17)	-0.0108 (17)
C24	0.081 (3)	0.118 (4)	0.075 (3)	-0.013 (2)	0.014 (2)	-0.008 (2)
C25	0.092 (3)	0.117 (4)	0.073 (2)	-0.026 (3)	-0.008 (2)	0.029 (2)
C26	0.064 (2)	0.060 (2)	0.104 (3)	-0.0069 (16)	-0.018 (2)	0.021 (2)
C27	0.0495 (16)	0.0558 (19)	0.0654 (18)	-0.0039 (14)	0.0031 (15)	-0.0061 (15)
C28	0.099 (3)	0.130 (4)	0.068 (2)	-0.011 (3)	0.016 (2)	-0.008 (2)
C29	0.0360 (14)	0.053 (2)	0.135 (3)	0.0018 (13)	0.0092 (17)	-0.0380 (19)
C30	0.0364 (14)	0.053 (2)	0.135 (3)	-0.0005 (13)	0.0082 (17)	-0.0372 (19)
C31	0.0445 (14)	0.0370 (15)	0.0443 (14)	-0.0032 (11)	-0.0018 (12)	0.0001 (11)
C32	0.0450 (14)	0.0439 (17)	0.0599 (16)	0.0015 (12)	-0.0139 (13)	-0.0032 (13)
C33	0.0487 (15)	0.0383 (16)	0.0607 (16)	0.0024 (12)	-0.0113 (14)	-0.0041 (12)
C34	0.0452 (14)	0.0405 (15)	0.0488 (14)	-0.0043 (12)	-0.0051 (12)	0.0039 (12)
C35	0.0419 (14)	0.0489 (18)	0.0606 (16)	-0.0020 (12)	-0.0092 (13)	0.0005 (13)
C36	0.0465 (15)	0.0448 (17)	0.0608 (16)	0.0066 (12)	-0.0019 (14)	-0.0033 (13)
C37	0.0494 (15)	0.0449 (18)	0.0549 (16)	-0.0024 (13)	-0.0004 (14)	-0.0028 (13)
C38	0.0482 (15)	0.0445 (17)	0.0636 (17)	-0.0044 (13)	-0.0095 (14)	-0.0044 (13)
N1	0.0382 (12)	0.0499 (14)	0.0623 (14)	-0.0034 (10)	-0.0078 (11)	-0.0085 (11)
N2	0.0304 (10)	0.0540 (14)	0.0484 (12)	-0.0042 (10)	-0.0031 (10)	-0.0044 (10)
N3	0.0361 (11)	0.0408 (13)	0.0579 (13)	-0.0021 (9)	-0.0122 (10)	-0.0067 (10)
N4	0.0298 (10)	0.0412 (13)	0.0478 (11)	-0.0028 (9)	-0.0090 (9)	-0.0061 (9)
N5	0.0352 (11)	0.0444 (14)	0.0801 (16)	-0.0001 (10)	0.0049 (12)	-0.0153 (12)
N6	0.0486 (13)	0.0566 (16)	0.0644 (15)	0.0090 (11)	-0.0013 (12)	-0.0023 (12)
O1	0.0630 (12)	0.0492 (13)	0.0933 (15)	-0.0026 (10)	-0.0258 (12)	-0.0187 (10)
O2	0.0744 (14)	0.0431 (13)	0.1239 (19)	0.0057 (11)	-0.0263 (13)	-0.0218 (12)
O3	0.0662 (12)	0.0494 (13)	0.0883 (14)	0.0003 (10)	-0.0303 (11)	-0.0199 (11)
O4	0.0714 (14)	0.0772 (17)	0.1216 (19)	0.0137 (12)	-0.0481 (15)	-0.0336 (14)

supplementary materials

Geometric parameters (Å, °)

C1—N1	1.388 (3)	C20—C20 ⁱⁱ	1.506 (4)
C1—C2	1.388 (4)	C20—H20A	0.9700
C1—C6	1.393 (4)	C20—H20B	0.9700
C2—N2	1.378 (3)	C21—N5	1.377 (3)
C2—C3	1.384 (4)	C21—C26	1.387 (4)
C3—C4	1.367 (4)	C21—C22	1.392 (3)
C3—H3	0.9300	C22—C23	1.377 (4)
C4—C5	1.377 (5)	C22—N6	1.387 (3)
C4—H4	0.9300	C23—C24	1.353 (5)
C5—C6	1.367 (4)	C23—H23	0.9300
C5—H5	0.9300	C24—C25	1.381 (5)
C6—H6	0.9300	C24—H24	0.9300
C7—N1	1.319 (3)	C25—C26	1.380 (5)
C7—N2	1.355 (3)	C25—H25	0.9300
C7—C8	1.482 (4)	C26—H26	0.9300
C8—H8A	0.9600	C27—N6	1.308 (3)
C8—H8B	0.9600	C27—N5	1.363 (3)
C8—H8C	0.9600	C27—C28	1.479 (4)
C9—N3	1.382 (3)	C28—H28A	0.9600
C9—C14	1.384 (3)	C28—H28B	0.9600
C9—C10	1.391 (3)	C28—H28C	0.9600
C10—C11	1.375 (4)	C29—N5	1.456 (3)
C10—N4	1.390 (3)	C29—C30	1.502 (4)
C11—C12	1.371 (4)	C29—H29A	0.9700
C11—H11	0.9300	C29—H29B	0.9700
C12—C13	1.393 (4)	C30—C30 ⁱⁱⁱ	1.515 (5)
C12—H12	0.9300	C30—H30A	0.9700
C13—C14	1.359 (4)	C30—H30B	0.9700
C13—H13	0.9300	C31—C36	1.377 (3)
C14—H14	0.9300	C31—C32	1.381 (3)
C15—N3	1.318 (3)	C31—C37	1.488 (4)
C15—N4	1.353 (3)	C32—C33	1.382 (4)
C15—C16	1.477 (4)	C32—H32	0.9300
C16—H16A	0.9600	C33—C34	1.382 (3)
C16—H16B	0.9600	C33—H33	0.9300
C16—H16C	0.9600	C34—C35	1.383 (4)
C17—N2	1.462 (3)	C34—C38	1.482 (4)
C17—C18	1.510 (3)	C35—C36	1.376 (4)
C17—H17A	0.9700	C35—H35	0.9300
C17—H17B	0.9700	C36—H36	0.9300
C18—C18 ⁱ	1.511 (5)	C37—O2	1.206 (3)
C18—H18A	0.9700	C37—O1	1.308 (3)
C18—H18B	0.9700	C38—O4	1.205 (3)
C19—N4	1.463 (3)	C38—O3	1.301 (3)
C19—C20	1.507 (3)	O1—H1	0.8199

C19—H19A	0.9700	O3—H3A	0.8197
C19—H19B	0.9700		
N1—C1—C2	109.2 (2)	N5—C21—C26	133.3 (3)
N1—C1—C6	130.6 (3)	N5—C21—C22	105.5 (2)
C2—C1—C6	120.2 (3)	C26—C21—C22	121.2 (3)
N2—C2—C3	132.0 (3)	C23—C22—N6	129.6 (2)
N2—C2—C1	105.9 (2)	C23—C22—C21	120.7 (3)
C3—C2—C1	122.1 (3)	N6—C22—C21	109.7 (2)
C4—C3—C2	116.5 (3)	C24—C23—C22	118.1 (3)
C4—C3—H3	121.8	C24—C23—H23	121.0
C2—C3—H3	121.8	C22—C23—H23	121.0
C3—C4—C5	122.1 (3)	C23—C24—C25	122.0 (4)
C3—C4—H4	118.9	C23—C24—H24	119.0
C5—C4—H4	118.9	C25—C24—H24	119.0
C6—C5—C4	121.8 (3)	C26—C25—C24	121.1 (3)
C6—C5—H5	119.1	C26—C25—H25	119.4
C4—C5—H5	119.1	C24—C25—H25	119.4
C5—C6—C1	117.2 (3)	C25—C26—C21	117.0 (3)
C5—C6—H6	121.4	C25—C26—H26	121.5
C1—C6—H6	121.4	C21—C26—H26	121.5
N1—C7—N2	112.3 (2)	N6—C27—N5	113.0 (3)
N1—C7—C8	124.5 (3)	N6—C27—C28	124.1 (3)
N2—C7—C8	123.2 (3)	N5—C27—C28	122.9 (3)
C7—C8—H8A	109.5	C27—C28—H28A	109.5
C7—C8—H8B	109.5	C27—C28—H28B	109.5
H8A—C8—H8B	109.5	H28A—C28—H28B	109.5
C7—C8—H8C	109.5	C27—C28—H28C	109.5
H8A—C8—H8C	109.5	H28A—C28—H28C	109.5
H8B—C8—H8C	109.5	H28B—C28—H28C	109.5
N3—C9—C14	130.0 (2)	N5—C29—C30	112.6 (2)
N3—C9—C10	109.79 (19)	N5—C29—H29A	109.1
C14—C9—C10	120.2 (2)	C30—C29—H29A	109.1
C11—C10—N4	132.6 (2)	N5—C29—H29B	109.1
C11—C10—C9	122.3 (2)	C30—C29—H29B	109.1
N4—C10—C9	105.0 (2)	H29A—C29—H29B	107.8
C12—C11—C10	116.4 (2)	C29—C30—C30 ⁱⁱⁱ	113.8 (3)
C12—C11—H11	121.8	C29—C30—H30A	108.8
C10—C11—H11	121.8	C30 ⁱⁱⁱ —C30—H30A	108.8
C11—C12—C13	121.8 (3)	C29—C30—H30B	108.8
C11—C12—H12	119.1	C30 ⁱⁱⁱ —C30—H30B	108.8
C13—C12—H12	119.1	H30A—C30—H30B	107.7
C14—C13—C12	121.4 (2)	C36—C31—C32	119.4 (2)
C14—C13—H13	119.3	C36—C31—C37	119.0 (2)
C12—C13—H13	119.3	C32—C31—C37	121.6 (2)
C13—C14—C9	117.8 (2)	C31—C32—C33	119.8 (2)
C13—C14—H14	121.1	C31—C32—H32	120.1
C9—C14—H14	121.1	C33—C32—H32	120.1
N3—C15—N4	112.5 (2)	C34—C33—C32	120.9 (2)

supplementary materials

N3—C15—C16	124.6 (2)	C34—C33—H33	119.6
N4—C15—C16	122.9 (2)	C32—C33—H33	119.6
C15—C16—H16A	109.5	C33—C34—C35	118.8 (2)
C15—C16—H16B	109.5	C33—C34—C38	122.0 (2)
H16A—C16—H16B	109.5	C35—C34—C38	119.1 (2)
C15—C16—H16C	109.5	C36—C35—C34	120.3 (2)
H16A—C16—H16C	109.5	C36—C35—H35	119.8
H16B—C16—H16C	109.5	C34—C35—H35	119.8
N2—C17—C18	112.0 (2)	C35—C36—C31	120.7 (2)
N2—C17—H17A	109.2	C35—C36—H36	119.6
C18—C17—H17A	109.2	C31—C36—H36	119.6
N2—C17—H17B	109.2	O2—C37—O1	123.6 (3)
C18—C17—H17B	109.2	O2—C37—C31	122.7 (2)
H17A—C17—H17B	107.9	O1—C37—C31	113.7 (2)
C17—C18—C18 ⁱ	110.9 (3)	O4—C38—O3	122.7 (3)
C17—C18—H18A	109.5	O4—C38—C34	123.2 (3)
C18 ⁱ —C18—H18A	109.5	O3—C38—C34	114.1 (2)
C17—C18—H18B	109.5	C7—N1—C1	105.6 (2)
C18 ⁱ —C18—H18B	109.5	C7—N2—C2	107.1 (2)
H18A—C18—H18B	108.0	C7—N2—C17	128.4 (2)
N4—C19—C20	112.81 (19)	C2—N2—C17	124.5 (2)
N4—C19—H19A	109.0	C15—N3—C9	105.54 (19)
C20—C19—H19A	109.0	C15—N4—C10	107.14 (18)
N4—C19—H19B	109.0	C15—N4—C19	126.2 (2)
C20—C19—H19B	109.0	C10—N4—C19	126.5 (2)
H19A—C19—H19B	107.8	C27—N5—C21	106.6 (2)
C20 ⁱⁱ —C20—C19	111.6 (2)	C27—N5—C29	127.9 (3)
C20 ⁱⁱ —C20—H20A	109.3	C21—N5—C29	125.4 (3)
C19—C20—H20A	109.3	C27—N6—C22	105.1 (2)
C20 ⁱⁱ —C20—H20B	109.3	C37—O1—H1	109.3
C19—C20—H20B	109.3	C38—O3—H3A	109.4
H20A—C20—H20B	108.0		

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

Hydrogen-bond geometry ($\text{\AA}, ^\circ$)

$D-H\cdots A$	$D-H$	$H\cdots A$	$D\cdots A$	$D-H\cdots A$
O1—H1 \cdots N1 ^{iv}	0.82	1.83	2.609 (3)	157
O3—H3A \cdots N3 ^v	0.82	1.87	2.668 (3)	165

Symmetry codes: (iv) $-x+1, -y+1, -z+1$; (v) $x-1, y-1, z$.

Fig. 1

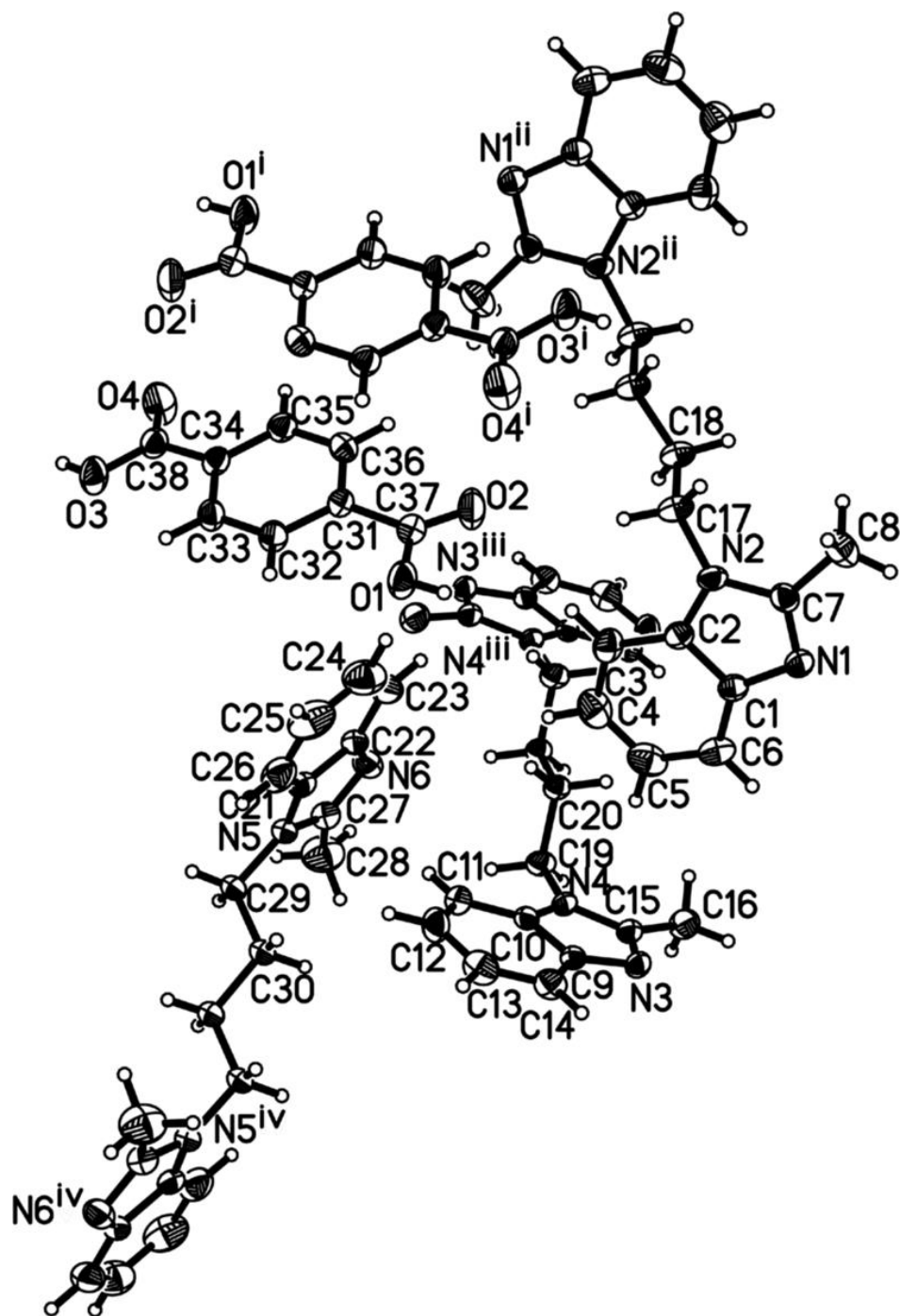


Fig. 2

