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Gabapentinium picrate

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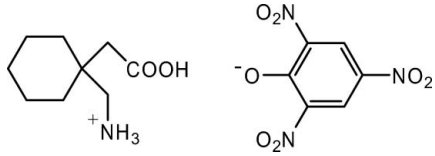
Key indicators: single-crystal X-ray study; T = 296 K; mean σ(C–C) = 0.004 Å; R factor = 0.062; wR factor = 0.182; data-to-parameter ratio = 12.3.

The title compound {systematic name: [1-(carboxymethyl)-cyclohexyl]methanaminium 2,4,6-trinitrophenolate}, $C_9H_{18}NO_2^+\cdot C_6H_2N_3O_7^-$, was synthesized from picric acid and gabapentin. The crystal packing is stabilized by intramolecular N—H···O═N and N—H···O—Ph hydrogen bonds. An O—H···O interaction is also present.

Related literature

For background, see: Bryans & Wustrow (1999). For related structures, see: Ibers (2001); Swamy *et al.* (2007) and references cited therein.



Experimental

Crystal data

$C_9H_{18}NO_2^+\cdot C_6H_2N_3O_7^-$
M_r = 400.35
Monoclinic, $P2_1/n$
a = 11.576 (2) Å
b = 7.7312 (16) Å
c = 19.973 (4) Å
β = 91.425 (2)°
V = 1787.0 (6) Å^3
Z = 4
Mo $K\alpha$ radiation
μ = 0.12 mm^{-1}
T = 296 K
0.25 × 0.25 × 0.20 mm

Data collection

Bruker SMART CCD area-detector diffractometer
Absorption correction: multi-scan (*SADABS*; Sheldrick, 2004)
T_{min} = 0.970, T_{max} = 0.976
8899 measured reflections
3150 independent reflections
2408 reflections with $I > 2\sigma(I)$
R_{int} = 0.034

Refinement

$R[F^2 > 2\sigma(F^2)]$ = 0.062
$wR(F^2)$ = 0.182
S = 1.07
3150 reflections
256 parameters
H-atom parameters constrained
$\Delta\rho_{max}$ = 1.06 e Å^{-3}
$\Delta\rho_{min}$ = −0.55 e Å^{-3}

Table 1
Hydrogen-bond geometry (Å, °).

D—H···A	D—H	H···A	D···A	D—H···A
O9—H9···O8^{i}	0.82	1.86	2.672 (3)	174
N4—H4C···O4^{ii}	0.89	2.32	2.957 (4)	128
N4—H4C···O3^{ii}	0.89	2.06	2.862 (3)	149
N4—H4B···O2^{iii}	0.89	2.41	2.894 (3)	114
N4—H4B···O4^{i}	0.89	2.23	3.063 (3)	155
N4—H4A···O3^{iii}	0.89	2.21	3.081 (3)	166

Symmetry codes: (i) $-x+1, -y+2, -z+1$; (ii) $x+\frac{1}{2}, -y+\frac{3}{2}, z+\frac{1}{2}$; (iii) $-x+1, -y+1, -z+1$.

Data collection: *SMART* (Bruker, 2001); cell refinement: *SAINT* (Bruker, 2001); data reduction: *SAINT*; program(s) used to solve structure: *SHELXS97* (Sheldrick, 2008); program(s) used to refine structure: *SHELXL97* (Sheldrick, 2008); molecular graphics: *SHELXTL* (Sheldrick, 2008); software used to prepare material for publication: *SHELXTL*.

LM thanks University of Mysore for research facilities.

Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: CS2105).

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Swamy, M. T., Ashok, M. A., Yathirajan, H. S., Narayana, B. & Bolte, M. (2007). *Acta Cryst.* E**63**, o4919.

supporting information

Acta Cryst. (2009). E**65**, o783 [doi:10.1107/S1600536809008952]

Gabapentinium picrate

Hongqi Li, H. S. Yathirajan, L. Mallesha, K. N. Mohana and B. Narayana

S1. Comment

Gabapentin, (1-(aminomethyl) cyclohexane acetic acid; Neurontin), is a novel anticonvulsant agent and has therapeutically beneficial effects against chronic pain states and anxiety (Bryans & Wustrow, 1999). Gabapentin is a zwitterion in the solid state (Ibers, 2001). On the other hand, picric acid forms salts or charge-transfer complexes with many organic compounds and we have reported crystal structures of a number of picrate complexes with organic compounds of pharmaceutical importance *viz*. desipraminium picrate (Swamy *et al.*, 2007). The present paper reports the crystal structure of the title compound, (1-(carboxymethyl)cyclohexyl)methanaminium 2,4,6-trinitrophenolate.

S2. Experimental

The title compound was synthesized by mixing solutions of picric acid (4.59 g, 0.01 mol) in 20 ml of distilled water and gabapentin (1.72 g, 0.01 mol) in 20 ml of distilled water and the resulting solution was stirred well for 10 min. A yellow precipitate of gabapentinium picrate was formed almost instantaneously after stirring. The so formed yellow complex was filtered off, washed with distilled water and dried *in vacuo* over anhydrous calcium chloride. The compound was purified by successive recrystallization from methanol (yield 92%). Single crystals for X-ray studies were grown by slow evaporation of a methanol solution. Analysis (%) found (calculated) for $C_{15}H_{20}N_4O_9$: C: 44.68 (45.00); H: 5.11 (5.04); N: 13.73 (13.99)%.

S3. Refinement

All H atoms were placed at calculated positions and refined using a riding model approximation, with C—H = 0.93–0.97 Å and with U_{iso}(H) = 1.2U_{eq}(C).

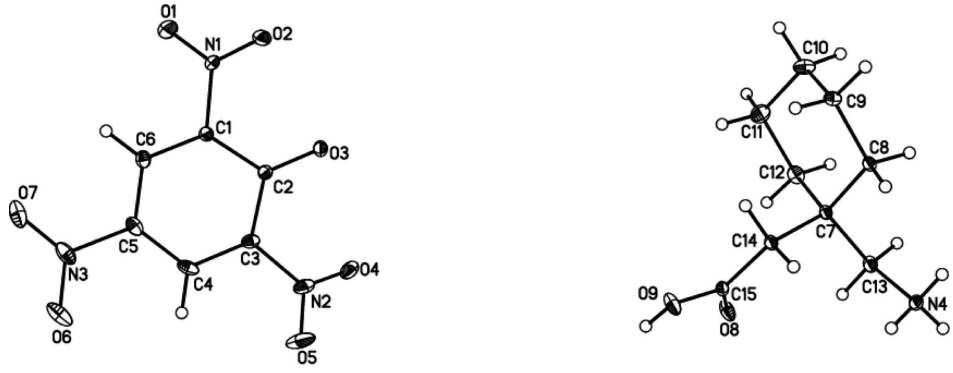

Figure 1

A view of the asymmetric unit of the title compound. Displacement ellipsoids are drawn at the 30% probability level and H atoms are shown as small spheres of arbitrary radii.

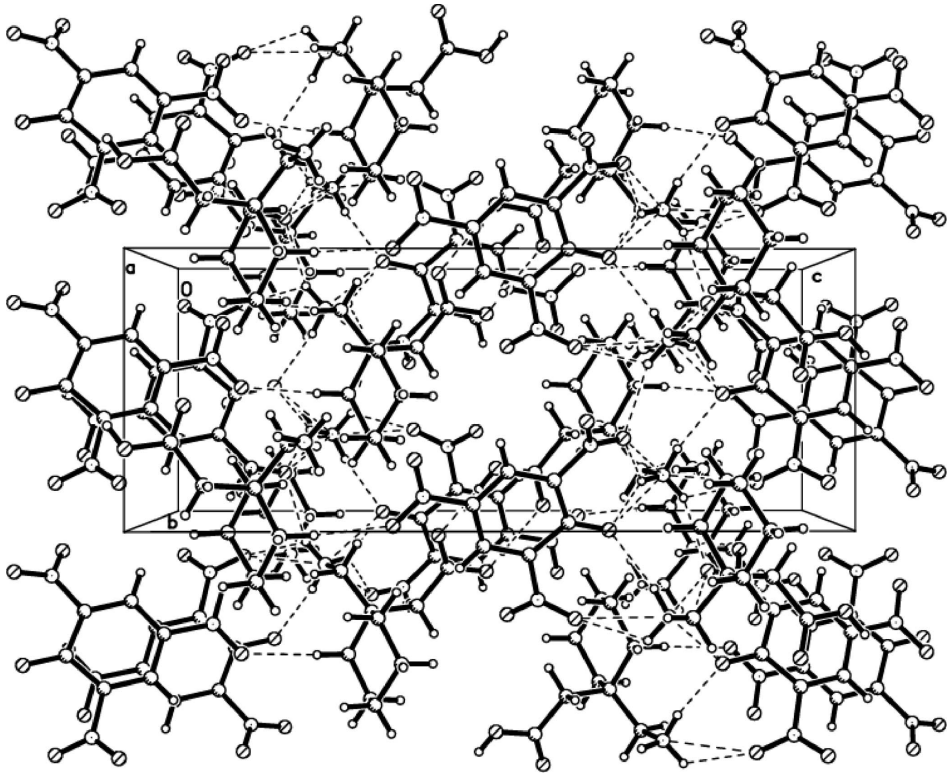

Figure 2

Packing in the crystal structure of the title compound, viewed along the *a* axis

[1-(Carboxymethyl)cyclohexyl]methanaminium 2,4,6-trinitrophenolate

Crystal data

$C_9H_{18}NO_2^+·C_6H_2N_3O_7^-$
M_r = 400.35
Monoclinic, $P2_1/n$
Hall symbol: -P 2yn
a = 11.576 (2) Å
b = 7.7312 (16) Å
c = 19.973 (4) Å
β = 91.425 (2)°
V = 1787.0 (6) Å^3
Z = 4
$F(000)$ = 840
D_x = 1.488 Mg m^{-3}
Melting point = 431–434 K
Mo $K\alpha$ radiation, λ = 0.71073 Å
Cell parameters from 2666 reflections
θ = 2.8–25.1°
μ = 0.12 mm^{-1}
T = 296 K
Block, yellow
0.25 × 0.25 × 0.20 mm

Data collection

Bruker SMART CCD area-detector diffractometer
Radiation source: fine-focus sealed tube
Graphite monochromator
φ and ω scans
Absorption correction: multi-scan (*SADABS*; Sheldrick, 2004)
T_{min} = 0.970, T_{max} = 0.976
8899 measured reflections
3150 independent reflections
2408 reflections with $I > 2\sigma(I)$
R_{int} = 0.034
θ_{max} = 25.1°, θ_{min} = 2.8°
h = −12→13
k = −7→9
l = −23→23

Refinement

Refinement on F^2
Least-squares matrix: full
$R[F^2 > 2\sigma(F^2)] = 0.062$
$wR(F^2) = 0.182$
$S = 1.07$
3150 reflections
256 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
$w = 1/[\sigma^2(F_o^2) + (0.083P)^2 + 1.5949P]$ where $P = (F_o^2 + 2F_c^2)/3$
$(\Delta/\sigma)_{max} < 0.001$
$\Delta\rho_{max} = 1.06$ e Å^{-3}
$\Delta\rho_{min} = -0.55$ e Å^{-3}
Extinction correction: *SHELXL97* (Sheldrick, 2008), $Fc^*=kFc[1+0.001xFc^2\lambda^3/\sin(2\theta)]^{-1/4}$
Extinction coefficient: 0.075 (5)

Fractional atomic coordinates and isotropic or equivalent isotropic displacement parameters (Å²)

	x	*y*	*z*	U_{iso}*/U_{eq}
C1	0.2919 (2)	0.3970 (4)	0.05392 (13)	0.0375 (6)
C2	0.3067 (2)	0.5162 (3)	0.10930 (13)	0.0362 (6)
C3	0.3639 (2)	0.6728 (3)	0.08832 (16)	0.0432 (7)
C4	0.3949 (2)	0.7066 (4)	0.02348 (17)	0.0499 (8)
H4	0.4323	0.8093	0.0131	0.060*
C5	0.3701 (2)	0.5869 (4)	−0.02563 (15)	0.0466 (8)
C6	0.3199 (2)	0.4315 (4)	−0.01060 (14)	0.0433 (7)
H6	0.3052	0.3503	−0.0441	0.052*
C7	0.4810 (2)	0.6530 (3)	0.67642 (12)	0.0328 (6)
C8	0.5192 (2)	0.4847 (3)	0.71209 (13)	0.0372 (6)
H8A	0.6030	0.4811	0.7144	0.045*
H8B	0.4920	0.4871	0.7576	0.045*
C9	0.4752 (3)	0.3205 (4)	0.67810 (17)	0.0521 (8)
H9A	0.5112	0.3084	0.6350	0.063*
H9B	0.4972	0.2213	0.7053	0.063*
C10	0.3458 (3)	0.3223 (5)	0.6679 (2)	0.0727 (11)
H10A	0.3093	0.3214	0.7111	0.087*
H10B	0.3216	0.2192	0.6437	0.087*
C11	0.3074 (3)	0.4818 (5)	0.6289 (2)	0.0664 (10)
H11A	0.2237	0.4833	0.6247	0.080*
H11B	0.3381	0.4771	0.5842	0.080*
C12	0.3490 (3)	0.6467 (4)	0.66370 (16)	0.0502 (8)
H12A	0.3109	0.6569	0.7062	0.060*
H12B	0.3261	0.7453	0.6364	0.060*
C13	0.5043 (3)	0.8069 (4)	0.72276 (14)	0.0429 (7)
H13A	0.4866	0.9122	0.6982	0.051*
H13B	0.4517	0.8003	0.7597	0.051*
C14	0.5472 (2)	0.6665 (3)	0.61064 (12)	0.0354 (6)
H14A	0.6290	0.6669	0.6222	0.042*
H14B	0.5319	0.5615	0.5853	0.042*
C15	0.5244 (2)	0.8160 (3)	0.56474 (13)	0.0365 (6)
N1	0.2450 (3)	0.2259 (3)	0.06681 (13)	0.0530 (7)

N2	0.3944 (3)	0.8022 (3)	0.13862 (17)	0.0582 (8)
N3	0.4040 (2)	0.6182 (5)	−0.09448 (17)	0.0633 (9)
N4	0.6251 (2)	0.8199 (3)	0.75081 (12)	0.0490 (7)
H4A	0.6443	0.7210	0.7710	0.074*
H4B	0.6292	0.9060	0.7803	0.074*
H4C	0.6734	0.8404	0.7178	0.074*
O1	0.1815 (4)	0.1631 (5)	0.02309 (16)	0.1380 (18)
O2	0.2579 (2)	0.1585 (3)	0.11974 (12)	0.0681 (8)
O3	0.2754 (2)	0.4861 (3)	0.16754 (10)	0.0492 (6)
O4	0.3311 (2)	0.8242 (3)	0.18624 (13)	0.0638 (7)
O5	0.4804 (3)	0.8876 (4)	0.1315 (2)	0.1066 (12)
O6	0.4584 (2)	0.7510 (5)	−0.10551 (16)	0.0911 (10)
O7	0.3773 (3)	0.5105 (5)	−0.13704 (14)	0.0893 (10)
O8	0.4688 (2)	0.9447 (3)	0.57786 (10)	0.0609 (7)
O9	0.5759 (2)	0.7959 (3)	0.50733 (11)	0.0616 (7)
H9	0.5576	0.8754	0.4820	0.092*

Atomic displacement parameters (Å²)

	U^{11}	U^{22}	U^{33}	U^{12}	U^{13}	U^{23}
C1	0.0436 (15)	0.0344 (14)	0.0344 (14)	−0.0022 (12)	−0.0046 (11)	0.0041 (11)
C2	0.0400 (14)	0.0318 (14)	0.0364 (15)	0.0020 (11)	−0.0054 (11)	0.0019 (11)
C3	0.0420 (15)	0.0297 (14)	0.0575 (18)	−0.0008 (12)	−0.0075 (13)	0.0030 (13)
C4	0.0384 (16)	0.0382 (16)	0.073 (2)	0.0036 (13)	0.0079 (14)	0.0239 (16)
C5	0.0394 (15)	0.0549 (19)	0.0457 (16)	0.0112 (14)	0.0059 (12)	0.0156 (15)
C6	0.0428 (16)	0.0514 (18)	0.0356 (15)	0.0031 (13)	−0.0026 (12)	0.0021 (13)
C7	0.0415 (14)	0.0297 (13)	0.0271 (13)	0.0035 (11)	0.0006 (10)	0.0015 (10)
C8	0.0490 (16)	0.0323 (14)	0.0301 (14)	−0.0001 (12)	−0.0014 (11)	0.0044 (11)
C9	0.073 (2)	0.0313 (16)	0.0520 (18)	−0.0097 (15)	−0.0040 (15)	0.0041 (13)
C10	0.076 (3)	0.058 (2)	0.084 (3)	−0.031 (2)	0.002 (2)	0.004 (2)
C11	0.0460 (19)	0.078 (3)	0.074 (2)	−0.0127 (18)	−0.0084 (16)	−0.001 (2)
C12	0.0416 (16)	0.059 (2)	0.0505 (18)	0.0065 (14)	0.0054 (13)	0.0072 (15)
C13	0.0624 (19)	0.0337 (15)	0.0325 (14)	0.0073 (13)	0.0012 (13)	−0.0009 (11)
C14	0.0466 (15)	0.0311 (14)	0.0285 (13)	0.0016 (11)	0.0016 (11)	0.0015 (10)
C15	0.0480 (16)	0.0346 (15)	0.0268 (13)	−0.0010 (12)	−0.0009 (11)	0.0005 (11)
N1	0.0781 (19)	0.0435 (15)	0.0369 (14)	−0.0202 (13)	−0.0054 (12)	−0.0054 (12)
N2	0.0603 (17)	0.0320 (14)	0.082 (2)	−0.0054 (13)	−0.0139 (16)	−0.0025 (14)
N3	0.0510 (16)	0.078 (2)	0.0612 (19)	0.0195 (16)	0.0141 (14)	0.0320 (18)
N4	0.0808 (19)	0.0321 (13)	0.0337 (13)	−0.0104 (12)	−0.0085 (12)	−0.0038 (10)
O1	0.240 (5)	0.111 (3)	0.0613 (19)	−0.105 (3)	−0.031 (2)	0.0042 (18)
O2	0.0992 (19)	0.0492 (14)	0.0547 (15)	−0.0240 (13)	−0.0203 (13)	0.0170 (11)
O3	0.0769 (15)	0.0367 (11)	0.0340 (11)	−0.0009 (10)	0.0009 (10)	0.0000 (8)
O4	0.0849 (18)	0.0401 (13)	0.0655 (16)	0.0009 (12)	−0.0174 (14)	−0.0124 (11)
O5	0.089 (2)	0.072 (2)	0.159 (3)	−0.0438 (18)	0.005 (2)	−0.029 (2)
O6	0.0774 (19)	0.101 (2)	0.096 (2)	0.0071 (17)	0.0278 (16)	0.0562 (19)
O7	0.106 (2)	0.116 (3)	0.0471 (15)	0.016 (2)	0.0227 (15)	0.0115 (17)
O8	0.1037 (19)	0.0422 (13)	0.0375 (12)	0.0242 (13)	0.0137 (11)	0.0090 (10)
O9	0.0925 (18)	0.0544 (14)	0.0389 (12)	0.0224 (13)	0.0198 (11)	0.0148 (10)

Geometric parameters (Å, º)

C1—C6	1.363 (4)	C10—H10B	0.9700
C1—C2	1.447 (4)	C11—C12	1.524 (5)
C1—N1	1.455 (4)	C11—H11A	0.9700
C2—O3	1.249 (3)	C11—H11B	0.9700
C2—C3	1.447 (4)	C12—H12A	0.9700
C3—C4	1.377 (4)	C12—H12B	0.9700
C3—N2	1.455 (4)	C13—N4	1.497 (4)
C4—C5	1.374 (5)	C13—H13A	0.9700
C4—H4	0.9300	C13—H13B	0.9700
C5—C6	1.371 (4)	C14—C15	1.495 (4)
C5—N3	1.460 (4)	C14—H14A	0.9700
C6—H6	0.9300	C14—H14B	0.9700
C7—C13	1.527 (4)	C15—O8	1.217 (3)
C7—C14	1.541 (3)	C15—O9	1.314 (3)
C7—C8	1.543 (3)	N1—O2	1.185 (3)
C7—C12	1.544 (4)	N1—O1	1.228 (4)
C8—C9	1.522 (4)	N2—O5	1.205 (4)
C8—H8A	0.9700	N2—O4	1.227 (4)
C8—H8B	0.9700	N3—O7	1.225 (4)
C9—C10	1.506 (5)	N3—O6	1.227 (4)
C9—H9A	0.9700	N4—H4A	0.8900
C9—H9B	0.9700	N4—H4B	0.8900
C10—C11	1.520 (5)	N4—H4C	0.8900
C10—H10A	0.9700	O9—H9	0.8200
C6—C1—C2	124.9 (3)	C10—C11—H11A	109.4
C6—C1—N1	116.3 (3)	C12—C11—H11A	109.4
C2—C1—N1	118.8 (2)	C10—C11—H11B	109.4
O3—C2—C1	124.2 (2)	C12—C11—H11B	109.4
O3—C2—C3	124.8 (3)	H11A—C11—H11B	108.0
C1—C2—C3	111.0 (2)	C11—C12—C7	113.7 (3)
C4—C3—C2	124.2 (3)	C11—C12—H12A	108.8
C4—C3—N2	117.0 (3)	C7—C12—H12A	108.8
C2—C3—N2	118.7 (3)	C11—C12—H12B	108.8
C5—C4—C3	119.3 (3)	C7—C12—H12B	108.8
C5—C4—H4	120.4	H12A—C12—H12B	107.7
C3—C4—H4	120.4	N4—C13—C7	115.4 (2)
C6—C5—C4	121.1 (3)	N4—C13—H13A	108.4
C6—C5—N3	118.5 (3)	C7—C13—H13A	108.4
C4—C5—N3	120.3 (3)	N4—C13—H13B	108.4
C1—C6—C5	119.4 (3)	C7—C13—H13B	108.4
C1—C6—H6	120.3	H13A—C13—H13B	107.5
C5—C6—H6	120.3	C15—C14—C7	119.4 (2)
C13—C7—C14	112.4 (2)	C15—C14—H14A	107.5
C13—C7—C8	109.4 (2)	C7—C14—H14A	107.5
C14—C7—C8	107.9 (2)	C15—C14—H14B	107.5

C13—C7—C12	106.5 (2)	C7—C14—H14B	107.5
C14—C7—C12	111.9 (2)	H14A—C14—H14B	107.0
C8—C7—C12	108.7 (2)	O8—C15—O9	122.6 (2)
C9—C8—C7	114.1 (2)	O8—C15—C14	125.9 (2)
C9—C8—H8A	108.7	O9—C15—C14	111.5 (2)
C7—C8—H8A	108.7	O2—N1—O1	121.3 (3)
C9—C8—H8B	108.7	O2—N1—C1	121.2 (2)
C7—C8—H8B	108.7	O1—N1—C1	117.0 (3)
H8A—C8—H8B	107.6	O5—N2—O4	121.8 (3)
C10—C9—C8	111.9 (3)	O5—N2—C3	118.9 (3)
C10—C9—H9A	109.2	O4—N2—C3	119.4 (3)
C8—C9—H9A	109.2	O7—N3—O6	124.5 (3)
C10—C9—H9B	109.2	O7—N3—C5	118.1 (3)
C8—C9—H9B	109.2	O6—N3—C5	117.3 (4)
H9A—C9—H9B	107.9	C13—N4—H4A	109.5
C9—C10—C11	110.7 (3)	C13—N4—H4B	109.5
C9—C10—H10A	109.5	H4A—N4—H4B	109.5
C11—C10—H10A	109.5	C13—N4—H4C	109.5
C9—C10—H10B	109.5	H4A—N4—H4C	109.5
C11—C10—H10B	109.5	H4B—N4—H4C	109.5
H10A—C10—H10B	108.1	C15—O9—H9	109.5
C10—C11—C12	111.0 (3)		

Hydrogen-bond geometry (Å, º)

D—H···*A*	*D*—H	H···*A*	*D*···*A*	*D*—H···*A*
O9—H9···O8^{i}	0.82	1.86	2.672 (3)	174
N4—H4*C*···O4^{ii}	0.89	2.32	2.957 (4)	128
N4—H4*C*···O3^{ii}	0.89	2.06	2.862 (3)	149
N4—H4*B*···O2^{iii}	0.89	2.41	2.894 (3)	114
N4—H4*B*···O4^{i}	0.89	2.23	3.063 (3)	155
N4—H4*A*···O3^{iii}	0.89	2.21	3.081 (3)	166

Symmetry codes: (i) −*x*+1, −*y*+2, −*z*+1; (ii) *x*+1/2, −*y*+3/2, *z*+1/2; (iii) −*x*+1, −*y*+1, −*z*+1.