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Retraction of articles by T. Liu *et al.*

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A series of 29 papers by Liu *et al.* are retracted.

As a result of problems with the data sets and incorrect atom assignments, 29 papers by Liu *et al.* are retracted. Full details of all the articles are given in Table 1.

Table 1
Details of articles to be retracted, in order of publication.

Title	Reference	DOI	Refcode
Tetrakis(pyrazine-κN)bis(thiocyanato-κN)manganese(II)	Liu & Xie (2007*a*)	10.1107/S1600536807026852	EDUMAS
(Dihydroxyglyoxime-κ^2N,N')bis(1,10-phenanthroline-κ^2N,N')copper(II) dinitrate dihydrate	Liu, Wang, Wang & Xie (2007*b*)	10.1107/S1600536807028255	EDUVAB
Tetrakis(pyrazine-κN)bis(thiocyanato-κN)zinc(II)	Liu & Xie (2007*b*)	10.1107/S1600536807028735	RIGQAA
Tetrakis(μ-2-pyridyloxyacetato)bis[(1,10-phenanthroline)(2-pyridyloxyacetato)-lanthanum(III)]	Liu, Wang, Wang & Xie (2007*c*)	10.1107/S1600536807030917	UDUMIQ
Polymeric $KNOF_2$	Liu Wang, Wang & Xie (2007*a*)	10.1107/S1600536807027195	ICSD 240891
(Dihydroxyglyoxime-κ^2N,N')bis(1,10-phenanthroline-κ^2N,N')cobalt(II) dinitrate dihydrate	Liu, Wang, Wang & Xie (2007*d*)	10.1107/S1600536807031224	WIHJED
Tetrakis(μ-2-pyridyloxyacetato)bis[(1,10-phenanthroline)(2-pyridyloxyacetato)-praseodymium(III)]	Liu, Wang, Wang & Xie (2007*e*)	10.1107/S1600536807032679	WIHQEK
Tetrakis[μ-(2-pyridyloxy)acetato-$\kappa^2O:O'$]bis{(1,10-phenanthroline-κ^2N,N')-[(2-pyridyloxy)acetato-κO]neodymium(III)}	Liu, Wang, Wang & Xie (2007*f*)	10.1107/S1600536807035349	TIGDAP
(Dihydroxyglyoxime-κ^2N,N')bis(1,10-phenanthroline-κ^2N,N')manganese(II) dinitrate dihydrate	Liu, Wang, Wang & Xie (2007*g*)	10.1107/S1600536807035076	TIGDET
2-Amino-3,5-dinitrobenzoic acid–ammonia (1/1)	Liu & Zhu (2007*j*)	10.1107/S1600536807040068	KIKQAX
2-Hydroxy-3,5-dinitrobenzamide monohydrate	Liu & Zhu (2007*k*)	10.1107/S1600536807039712	KIKQEB
2-(1-Hydroxy-2-pyridyl)acetamide monohydrate	Liu & Zhu (2007*l*)	10.1107/S1600536807040652	CIKQOD
Bis(2,2′-bipyridine-κN,N′)bis(thiocyanato-κN)iron(II)	Liu & Zhu (2007*a*)	10.1107/S1600536807043486	XIFXOA
catena-Poly[hexakis(μ_2-anilinoacetamide)bis(1,10-phenanthroline)disamarium(III]	Liu & Zhu (2007*b*)	10.1107/S1600536807045485	XILNAI
3-Hydroxy-2,4,6-trinitropyridine monohydrate	Liu & Zhu (2007*m*)	10.1107/S1600536807045230	PILNOO
catena-Poly[hexakis(μ_2-anilinoacetamide)bis(1,10-phenanthroline)-dipraseodymium(III)]	Liu & Zhu (2007*c*)	10.1107/S1600536807047733	SILZET
catena-Poly[[tetra-μ-anilinoacetamidato-bis(1,10-phenanthroline)dicerium(III)]-di-μ-anilinoacetamidato]	Liu & Zhu (2007*d*)	10.1107/S1600536807050969	GIMZOS
Tetrakis(pyridine-κN)bis(thiocyanato-κN)chromium(II)	Liu & Zhu (2007*e*)	10.1107/S1600536807051756	WINFAB
2-Ammonio-3-carboxy-5-nitrobenzoate monohydrate	Liu & Zhu (2007*n*)	10.1107/S1600536807048477	GINFEP
2-(Benzoylhydrazinocarbonyl)benzoic acid	Liu & Zhu (2007*o*)	10.1107/S160053680705204X	TINZIA
Tetrakis(pyridine-κN)bis(thiocyanato-κN)vanadium(II)	Liu & Zhu (2007*f*)	10.1107/S1600536807054529	HIPZIQ
catena-Poly[[(nitrato-κO)(1,10-phenanthroline-κ^2N,N')nickel(II)]-μ-acetamido-$\kappa^2O:N$]	Liu & Zhu (2007*g*)	10.1107/S1600536807056504	XIRGIP
catena-Poly[[(nitrato-κO)(1,10-phenanthroline-κ^2N,N')copper(II)]-μ-acetamido-$\kappa^2O:N$]	Liu & Zhu (2007*h*)	10.1107/S1600536807059077	HIQROP
catena-Poly[[(nitrato-κO)(1,10-phenanthroline-κ^2N,N')cobalt(II)]-μ-acetamidato-$\kappa^2O:N$]	Liu & Zhu (2007*i*)	10.1107/S1600536807060631	YIQMER
N′-Benzoyl-4-nitronicotinohydrazide	Liu & Zhu (2007*p*)	10.1107/S1600536807053068	CIPVON
N′-(3-Nitro-4-pyridylcarbonyl)pyridine-4-carbohydrazide	Liu & Zhu (2007*q*)	10.1107/S1600536807054876	RIRWEV

 doi:10.1107/S1600536809049976

Table 1 (continued)

Title	Reference	DOI	Refcode
Ethylenediammonium sulfate	Liu & Zhu (2007*r*)	10.1107/S1600536807056280	ETDAMS03
Ethylenediammonium perchlorate	Liu & Zhu (2007*s*)	10.1107/S1600536807059909	HIRYEN
catena-Poly[[(nitrato-κO)(1,10-phenanthroline-κ^2N,N')manganese(II)]-μ-nitrato-$\kappa^2O:O'$]	Liu & Zhu (2008)	10.1107/S160053680706254X	MIRROV

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N′-Benzoyl-4-nitronicotinohydrazide

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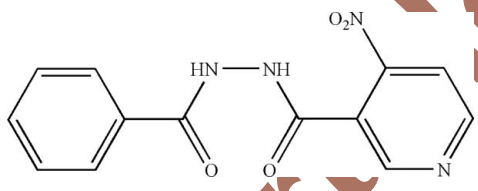
Key indicators: single-crystal X-ray study; T = 273 K; mean σ(C–C) = 0.004 Å; R factor = 0.049; wR factor = 0.139; data-to-parameter ratio = 13.5.

In the crystal structure of the title compound, $C_{13}H_{10}N_4O_4$, all bond lengths and angles are within normal ranges and the benzene and pyridine rings are oriented at a dihedral angle of 7.988 (4)°. N—H···O hydrogen bonds result in the formation of a supramolecular network.

Related literature

For general background, see: Desiraju (1995, 1997); Braga *et al.* (1998); Li *et al.* (2005); Liu *et al.* (2004); Pan & Xu (2004); Wu *et al.* (2003). For bond-length data, see: Allen *et al.* (1987).



Experimental

Crystal data

$C_{13}H_{10}N_4O_4$
M_r = 286.25
Monoclinic, $P2_1/n$
a = 8.290 (2) Å
b = 12.7981 (13) Å
c = 13.0073 (11) Å
β = 92.804 (3)°
V = 1378.4 (4) Å^3
Z = 4
Mo $K\alpha$ radiation
μ = 0.11 mm^{-1}
T = 273 (2) K
0.28 × 0.17 × 0.15 mm

Data collection

Bruker APEXII area-detector diffractometer
Absorption correction: multi-scan (*SADABS*; Sheldrick, 1996)
T_{min} = 0.971, T_{max} = 0.984
8484 measured reflections
2586 independent reflections
1276 reflections with $I > 2\sigma(I)$
R_{int} = 0.043

Refinement

$R[F^2 > 2\sigma(F^2)]$ = 0.049
$wR(F^2)$ = 0.139
S = 1.01
2586 reflections
191 parameters
H-atom parameters constrained
$\Delta\rho_{max}$ = 0.38 e Å^{-3}
$\Delta\rho_{min}$ = −0.26 e Å^{-3}

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

D—H···A	D—H	H···A	D···A	D—H···A
N4—H4A···O1^{i}	0.86	2.11	2.903 (3)	153
N3—H3A···O2^{ii}	0.86	1.98	2.831 (3)	170

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

Data collection: *APEX2* (Bruker, 2005); cell refinement: *SAINT* (Siemens, 1996); data reduction: *SAINT*; program(s) used to solve structure: *SHELXS97* (Sheldrick, 1997); program(s) used to refine structure: *SHELXL97* (Sheldrick, 1997); molecular graphics: *SHELXTL* (Siemens, 1996); software used to prepare material for publication: *SHELXTL*.

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Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: AT2444).

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supplementary materials

Acta Cryst. (2007). E**63**, o4527 [**doi:10.1107/S1600536807053068**]

N'-Benzoyl-4-nitronicotinohydrazide

T. Liu and J. Y. Zhu

Comment

In the synthesis of crystal structures by design, the assembly of molecular units in predefined arrangements is a key goal (Desiraju, 1995, 1997; Braga *et al.*, 1998). Due to pyridine groups are one of the most important classes of biological ligands, the coordination of metal-pyridine groups complexes are of critical importance in biological systems, organic materials and coordination chemistry. Recently, pyridine groups with variable coordination modes have been used to construct metal-organic supramolecular structures (Wu *et al.*, 2003; Pan & Xu, 2004; Liu *et al.*, 2004; Li *et al.*, 2005). We originally attempted to synthesize complexes featuring Cu metal chains by reaction of the copper(II) ion with *N*-(4-nitropyridine-3-carboxamido)-benzamide ligand. Unfortunately, we obtained only the title compound, (I), and we report herein its crystal structure.

In the molecule of (I) (Fig. 1), the ligand bond lengths and angles are within normal ranges (Allen *et al.*, 1987). In the crystal structure, the N—H···O hydrogen bonds (Table 1, Fig. 2) result in the formation of a supramolecular network structure.

Experimental

Crystals of the title compound were synthesized using hydrothermal method in a 23 ml Teflon-lined Parr bomb, which was then sealed. Copper dinitrate hexahydrate (147.8 mg, 0.5 mmol), *N*-(4-nitropyridine-3-carboxamido)-benzamide (286.3 mg, 1 mmol) and distilled water (5 g) were placed into the bomb and sealed. The bomb was then heated under autogenous pressure up to 453 K over the course of 7 d and allowed to cool at room temperature for 24 h. Upon opening the bomb, a clear colourless solution was decanted from small colourless crystals. These crystals were washed with distilled water followed by ethanol, and allowed to air-dry at room temperature.

Refinement

The H atoms were positioned geometrically, with N—H = 0.86 Å (for NH) and C—H = 0.93 Å for aromatic H, and constrained to ride on their parent atoms, with U_{iso}(H) = 1.2U_{eq}(C,*N*).

Figures

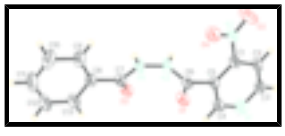

Fig. 1. The molecular structure of the title molecule, with the atom-numbering scheme. Displacement ellipsoids are drawn at the 30% probability level.

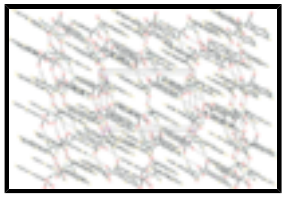

Fig. 2. A packing diagram of (I). Hydrogen bonds are shown as dashed lines.

N'-Benzoyl-4-nitronicotinohydrazide

Crystal data

$C_{13}H_{10}N_4O_4$

$M_r = 286.25$

Monoclinic, $P2_1/n$

Hall symbol: -P 2yn

$a = 8.290\ (2)$ Å

$b = 12.7981\ (13)$ Å

$c = 13.0073\ (11)$ Å

$\beta = 92.804\ (3)^\circ$

$V = 1378.4\ (4)$ Å^3

$Z = 4$

$F_{000} = 592$

$D_x = 1.379$ Mg m^{-3}

Mo $K\alpha$ radiation

$\lambda = 0.71073$ Å

Cell parameters from 1769 reflections

$\theta = 2.3–21.2^\circ$

$\mu = 0.11$ mm^{-1}

$T = 273\ (2)$ K

Prism, colourless

0.28 × 0.17 × 0.15 mm

Data collection

Bruker APEXII area-detector diffractometer

Radiation source: fine-focus sealed tube

Monochromator: graphite

$T = 273(2)$ K

φ and ω scans

Absorption correction: multi-scan (SADABS; Sheldrick, 1996)

$T_{min} = 0.971$, $T_{max} = 0.984$

8484 measured reflections

2586 independent reflections

1276 reflections with $I > 2\sigma(I)$

$R_{int} = 0.043$

$\theta_{max} = 26.0^\circ$

$\theta_{min} = 2.2^\circ$

$h = -10\rightarrow10$

$k = -15\rightarrow15$

$l = -16\rightarrow15$

Refinement

Refinement on F^2

Least-squares matrix: full

$R[F^2 > 2\sigma(F^2)] = 0.049$

$wR(F^2) = 0.139$

$S = 1.01$

2586 reflections

191 parameters

Hydrogen site location: inferred from neighbouring sites

H-atom parameters constrained

$w = 1/[\sigma^2(F_o^2) + (0.0593P)^2 + 0.15P]$ where $P = (F_o^2 + 2F_c^2)/3$

$(\Delta/\sigma)_{max} < 0.001$

$\Delta\rho_{max} = 0.38$ e Å^{-3}

$\Delta\rho_{min} = -0.26$ e Å^{-3}

Extinction correction: SHELXL97 (Sheldrick, 1997), $Fc^* = kFc[1+0.001xFc^2\lambda^3/\sin(2\theta)]^{-1/4}$

Primary atom site location: structure-invariant direct methods
Secondary atom site location: difference Fourier map
Extinction coefficient: 0.0085 (16)

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 $F^2 > 2\mathrm{sigma}(F^2)$ 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 (Å^2)

	x	*y*	*z*	U_{iso}*/U_{eq}
O3	0.7539 (4)	−0.37730 (18)	0.5475 (2)	0.1317 (10)
O4	0.8686 (3)	−0.23377 (16)	0.51348 (15)	0.0856 (7)
O1	0.9288 (2)	−0.01114 (13)	0.62696 (13)	0.0636 (5)
O2	0.5682 (2)	0.15823 (13)	0.51217 (15)	0.0682 (6)
N3	0.7074 (3)	−0.02935 (15)	0.51904 (16)	0.0626 (6)
H3A	0.6180	−0.0619	0.5061	0.075*
N4	0.7559 (3)	0.04934 (15)	0.45320 (16)	0.0632 (6)
H4A	0.8303	0.0384	0.4107	0.076*
N2	0.7830 (3)	−0.28636 (18)	0.56606 (18)	0.0714 (7)
C1	0.7145 (3)	−0.24006 (17)	0.65641 (18)	0.0529 (6)
C2	0.6407 (3)	−0.30418 (18)	0.7262 (2)	0.0609 (7)
H2	0.6335	−0.3756	0.7138	0.073*
C3	0.5793 (3)	−0.2641 (2)	0.8121 (2)	0.0631 (7)
H3	0.5313	−0.3082	0.8585	0.076*
N1	0.5872 (3)	−0.1581 (2)	0.8315 (2)	0.0929 (8)
C4	0.6614 (3)	−0.0939 (2)	0.76155 (19)	0.0611 (7)
H4	0.6682	−0.0226	0.7749	0.073*
C5	0.7258 (3)	−0.13250 (17)	0.67237 (18)	0.0508 (6)
C6	0.8002 (3)	−0.05385 (17)	0.6022 (2)	0.0512 (6)
C7	0.6825 (3)	0.14358 (19)	0.45821 (19)	0.0537 (6)
C8	0.7451 (3)	0.22485 (19)	0.38880 (19)	0.0563 (7)
C9	0.8229 (3)	0.2020 (2)	0.3001 (2)	0.0704 (8)
H9	0.8404	0.1328	0.2822	0.085*
C10	0.8756 (4)	0.2827 (3)	0.2370 (3)	0.0913 (10)
H10	0.9276	0.2677	0.1770	0.110*
C11	0.8490 (5)	0.3840 (3)	0.2656 (4)	0.1109 (14)
H11	0.8825	0.4382	0.2240	0.133*
C12	0.7744 (5)	0.4065 (2)	0.3537 (4)	0.1046 (13)
H12	0.7616	0.4758	0.3730	0.126*
C13	0.7176 (4)	0.3280 (2)	0.4150 (3)	0.0803 (9)

H13	0.6615	0.3442	0.4731	0.096*

Atomic displacement parameters (Å^2)

	U^{11}	U^{22}	U^{33}	U^{12}	U^{13}	U^{23}
O3	0.216 (3)	0.0550 (14)	0.130 (2)	−0.0236 (15)	0.0768 (18)	−0.0343 (13)
O4	0.1122 (17)	0.0719 (13)	0.0763 (14)	0.0048 (12)	0.0411 (12)	−0.0029 (11)
O1	0.0679 (12)	0.0460 (10)	0.0778 (12)	−0.0083 (9)	0.0122 (9)	−0.0001 (8)
O2	0.0728 (13)	0.0579 (12)	0.0753 (13)	−0.0045 (9)	0.0167 (10)	0.0018 (9)
N3	0.0641 (14)	0.0513 (12)	0.0728 (15)	−0.0121 (11)	0.0072 (12)	0.0217 (11)
N4	0.0730 (15)	0.0479 (13)	0.0704 (14)	−0.0018 (11)	0.0212 (11)	0.0171 (10)
N2	0.0987 (19)	0.0471 (14)	0.0704 (16)	0.0043 (13)	0.0260 (14)	−0.0045 (12)
C1	0.0622 (16)	0.0395 (13)	0.0576 (15)	0.0054 (12)	0.0094 (12)	0.0006 (11)
C2	0.0760 (19)	0.0363 (13)	0.0711 (18)	0.0008 (12)	0.0104 (14)	0.0096 (12)
C3	0.082 (2)	0.0484 (16)	0.0599 (17)	0.0002 (14)	0.0158 (14)	0.0135 (13)
N1	0.110 (2)	0.084 (2)	0.0866 (18)	0.0032 (16)	0.0166 (15)	0.0102 (14)
C4	0.0780 (19)	0.0450 (14)	0.0621 (17)	−0.0015 (13)	0.0217 (14)	0.0005 (12)
C5	0.0562 (16)	0.0389 (13)	0.0578 (16)	0.0009 (11)	0.0089 (12)	0.0045 (11)
C6	0.0586 (16)	0.0337 (13)	0.0627 (17)	0.0002 (13)	0.0171 (13)	−0.0005 (11)
C7	0.0573 (17)	0.0486 (16)	0.0551 (16)	−0.0056 (13)	0.0019 (13)	0.0028 (12)
C8	0.0595 (17)	0.0479 (15)	0.0610 (17)	−0.0084 (12)	−0.0009 (13)	0.0109 (12)
C9	0.073 (2)	0.0670 (18)	0.071 (2)	−0.0065 (15)	0.0033 (15)	0.0219 (15)
C10	0.084 (2)	0.110 (3)	0.080 (2)	−0.015 (2)	0.0005 (17)	0.041 (2)
C11	0.123 (3)	0.086 (3)	0.119 (3)	−0.041 (2)	−0.030 (3)	0.057 (3)
C12	0.149 (4)	0.0486 (19)	0.113 (3)	−0.026 (2)	−0.025 (3)	0.023 (2)
C13	0.098 (2)	0.0501 (18)	0.091 (2)	−0.0077 (15)	−0.0154 (17)	0.0086 (15)

Geometric parameters (Å, °)

O3—N2	1.211 (3)	N1—C4	1.392 (3)
O4—N2	1.213 (3)	C4—C5	1.391 (3)
O1—C6	1.227 (3)	C4—H4	0.9300
O2—C7	1.220 (3)	C5—C6	1.511 (3)
N3—C6	1.333 (3)	C7—C8	1.487 (3)
N3—N4	1.394 (3)	C8—C9	1.380 (4)
N3—H3A	0.8600	C8—C13	1.385 (4)
N4—C7	1.354 (3)	C9—C10	1.402 (4)
N4—H4A	0.8600	C9—H9	0.9300
N2—C1	1.456 (3)	C10—C11	1.370 (6)
C1—C2	1.388 (3)	C10—H10	0.9300
C1—C5	1.395 (3)	C11—C12	1.360 (5)
C2—C3	1.351 (4)	C11—H11	0.9300
C2—H2	0.9300	C12—C13	1.379 (5)
C3—N1	1.381 (3)	C12—H12	0.9300
C3—H3	0.9300	C13—H13	0.9300
C6—N3—N4	119.7 (2)	O1—C6—N3	124.6 (2)
C6—N3—H3A	120.1	O1—C6—C5	120.9 (2)
N4—N3—H3A	120.1	N3—C6—C5	114.2 (2)

C7—N4—N3	118.0 (2)	O2—C7—N4	121.9 (2)
C7—N4—H4A	121.0	O2—C7—C8	123.1 (2)
N3—N4—H4A	121.0	N4—C7—C8	115.0 (2)
O4—N2—O3	122.5 (2)	C9—C8—C13	119.8 (2)
O4—N2—C1	119.4 (2)	C9—C8—C7	123.4 (2)
O3—N2—C1	118.1 (2)	C13—C8—C7	116.8 (3)
C2—C1—C5	121.0 (2)	C8—C9—C10	120.4 (3)
C2—C1—N2	119.3 (2)	C8—C9—H9	119.8
C5—C1—N2	119.7 (2)	C10—C9—H9	119.8
C3—C2—C1	120.8 (2)	C11—C10—C9	118.6 (3)
C3—C2—H2	119.6	C11—C10—H10	120.7
C1—C2—H2	119.6	C9—C10—H10	120.7
C2—C3—N1	120.5 (2)	C10—C11—C12	121.0 (3)
C2—C3—H3	119.7	C10—C11—H11	119.5
N1—C3—H3	119.7	C12—C11—H11	119.5
C3—N1—C4	118.6 (2)	C11—C12—C13	121.0 (3)
C5—C4—N1	122.4 (2)	C11—C12—H12	119.5
C5—C4—H4	118.8	C13—C12—H12	119.5
N1—C4—H4	118.8	C12—C13—C8	119.1 (4)
C4—C5—C1	116.7 (2)	C12—C13—H13	120.4
C4—C5—C6	116.9 (2)	C8—C13—H13	120.4
C1—C5—C6	126.5 (2)		

Hydrogen-bond geometry (Å, °)

D—H···*A*	*D*—H	H···*A*	*D*···*A*	*D*—H···*A*
N4—H4A···O1^{i}	0.86	2.11	2.903 (3)	153
N3—H3A···O2^{ii}	0.86	1.98	2.831 (3)	170

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

Fig. 1

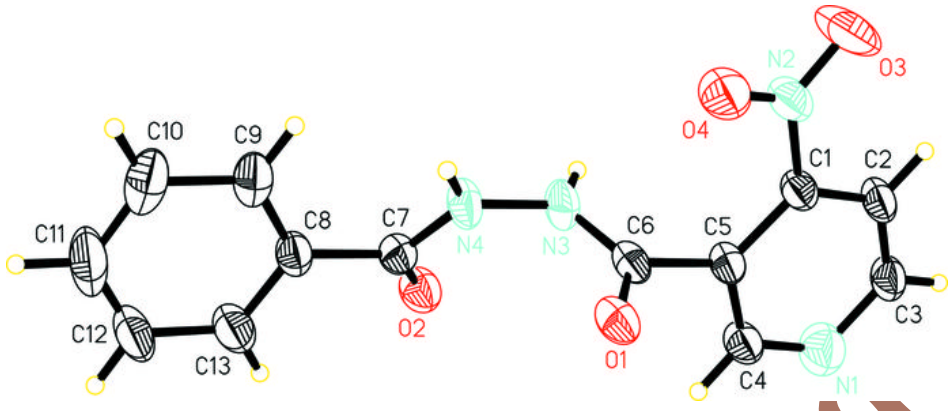

Fig. 2