



Crystal structure of 2-(*p*-tolyl)-6-(tri-fluoromethyl)benzo[*b*]thiophene-3-carbonitrile

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Received 30 April 2015; accepted 5 May 2015

Edited by W. T. A. Harrison, University of Aberdeen, Scotland

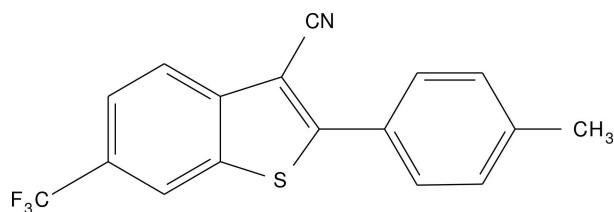
In the title compound, C₁₇H₁₀F₃NS, the dihedral angle between the fused benzothiophene ring system (r.m.s. deviation = 0.042 Å) and the benzene ring is 29.78 (11)°. The crystal structure features C—H···F and very weak C—H···N hydrogen bonds, which generate (001) sheets.

Keywords: crystal structure; benzo[*b*]thiophene; hydrogen bonding.

CCDC reference: 1063141

1. Related literature

For background to benzothiophene derivatives, see: Bettinetti *et al.* (2002); Roberts & Hartley (2004).



2. Experimental

2.1. Crystal data

C ₁₇ H ₁₀ F ₃ NS	$V = 1425.03 (10) \text{ \AA}^3$
$M_r = 317.33$	$Z = 4$
Monoclinic, $P2_1/c$	Cu $K\alpha$ radiation
$a = 13.7576 (5) \text{ \AA}$	$\mu = 2.29 \text{ mm}^{-1}$
$b = 14.5343 (6) \text{ \AA}$	$T = 293 \text{ K}$
$c = 7.1353 (3) \text{ \AA}$	$0.30 \times 0.27 \times 0.25 \text{ mm}$
$\beta = 92.817 (3)^\circ$	

2.2. Data collection

Bruker X8 Proteum diffractometer	7045 measured reflections
Absorption correction: multi-scan (SADABS; Bruker, 2013)	2316 independent reflections
$T_{\min} = 0.546$, $T_{\max} = 0.598$	1860 reflections with $I > 2\sigma(I)$
	$R_{\text{int}} = 0.063$

2.3. Refinement

$R[F^2 > 2\sigma(F^2)] = 0.053$	200 parameters
$wR(F^2) = 0.141$	H-atom parameters constrained
$S = 1.06$	$\Delta\rho_{\max} = 0.43 \text{ e \AA}^{-3}$
2316 reflections	$\Delta\rho_{\min} = -0.40 \text{ e \AA}^{-3}$

Table 1

Hydrogen-bond geometry (Å, °).

$D-H\cdots A$	$D-H$	$H\cdots A$	$D\cdots A$	$D-H\cdots A$
C3—H3···N11 ⁱ	0.93	2.62	3.411 (4)	143
C22—H22C···F15 ⁱⁱ	0.96	2.45	3.375 (4)	162

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

Data collection: *APEX2* (Bruker, 2013); cell refinement: *SAINT* (Bruker, 2013); data reduction: *SAINT*; program(s) used to solve structure: *SHELXS97* (Sheldrick, 2008); program(s) used to refine structure: *SHELXL97* (Sheldrick, 2008); molecular graphics: *PLATON* (Spek, 2009); software used to prepare material for publication: *PLATON*.

Acknowledgements

The authors are thankful to Institution of Excellence, University of Mysore, Mysore, for providing the single-crystal X-ray diffraction facility.

Supporting information for this paper is available from the IUCr electronic archives (Reference: HB7416).

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supporting information

Acta Cryst. (2015). E71, o382 [doi:10.1107/S2056989015008671]

Crystal structure of 2-(*p*-tolyl)-6-(trifluoromethyl)benzo[*b*]thiophene-3-carbonitrile

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S1. Comment

Benzo[*b*]thiophene derivatives are important heterocyclic compounds because of their various applications in medicinal chemistry. They represent an important heterocyclic core and are shown to display a range of promising pharmacological properties such as antipsychotic, antidepressive, antithrombolytic, dopamine receptor antagonist and 5-lipoxygenase inhibitor. Number of 2-arylbenzo[*b*]thiophene derivatives have indeed, these sulfur heterocycles are essential components of clinically important drugs such as Clopidogrel (Bettinetti *et al.*, 2002), Raloxifene (Roberts & Hartley, 2004) and Zileuton.

S2. Experimental

To a solution of 2-(2-chlorophenyl)acetonitrile (1.0 mmol), methyl benzodithioate (1.0 mmol) in DMF (2 ml), K₃PO₄ (2.0 mmol), pivalic acid (1.5 mmol), cuprous iodide (0.2 mmol) were added. The mixture was stirred at 80°C and progress was monitored by TLC. When the dithioesters could no longer be detected, the reaction mixture was extracted with EtOAc (3 × 10 ml). The organic layer was dried over anhydrous Na₂SO₄. The solvent was then removed under reduced pressure and the residue was purified by silica gel chromatography. White solid single crystals were obtained from slow evaporation of its solvent.

S3. Refinement

All H atoms were positioned geometrically and allowed to ride on their parent atom, with C–H = 0.93–0.97 Å, and with $U_{\text{iso}}(\text{H}) = 1.2\text{--}1.5U_{\text{eq}}(\text{C})$.

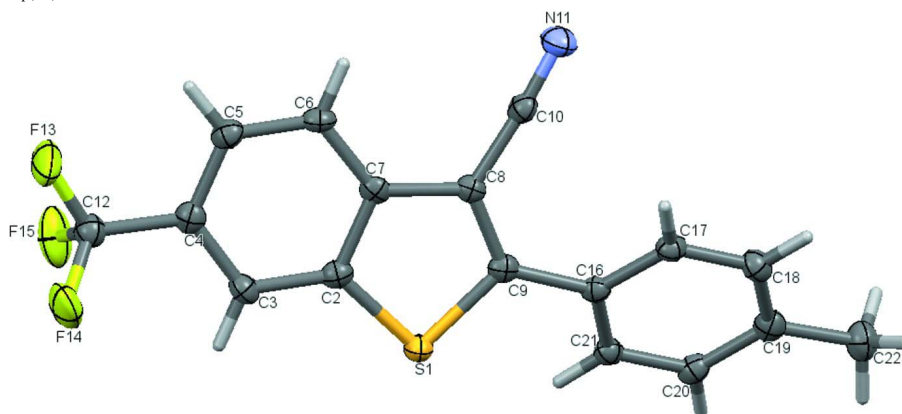
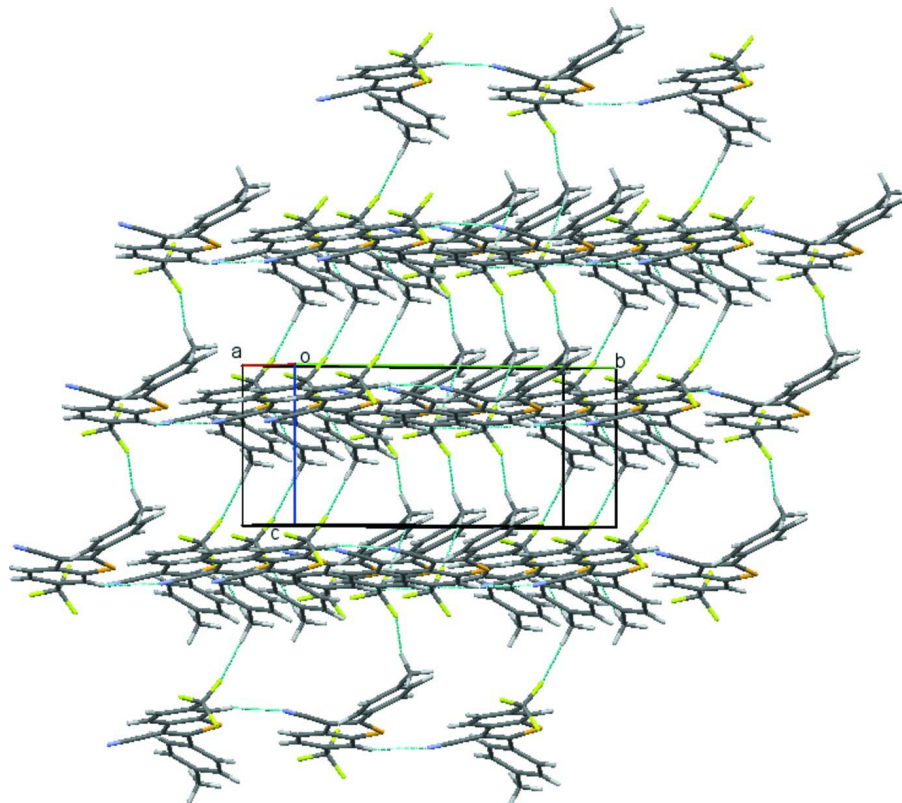


Figure 1

A view of the title compound with displacement ellipsoids drawn at the 50% probability level.

**Figure 2**

A view along the *a* axis of the crystal packing of the title compound. Hydrogen bonds are shown as dashed lines (see Table 2 for details).

2-(*p*-Tolyl)-6-(trifluoromethyl)benzo[*b*]thiophene-3-carbonitrile

Crystal data

$C_{17}H_{10}F_3NS$

$M_r = 317.33$

Monoclinic, $P2_1/c$

Hall symbol: $-P\ 2_1/c$

$a = 13.7576$ (5) Å

$b = 14.5343$ (6) Å

$c = 7.1353$ (3) Å

$\beta = 92.817$ (3)°

$V = 1425.03$ (10) Å³

$Z = 4$

$F(000) = 648$

$D_x = 1.479$ Mg m⁻³

Cu $K\alpha$ radiation, $\lambda = 1.54178$ Å

Cell parameters from 2316 reflections

$\theta = 6.4\text{--}64.7^\circ$

$\mu = 2.29$ mm⁻¹

$T = 293$ K

Block, colourless

$0.30 \times 0.27 \times 0.25$ mm

Data collection

Bruker X8 Proteum
diffractometer

Radiation source: Rotating Anode

Graphite monochromator

Detector resolution: 18.4 pixels mm⁻¹

φ and ω scans

Absorption correction: multi-scan
(*SADABS*; Bruker, 2013)

$T_{\min} = 0.546$, $T_{\max} = 0.598$

7045 measured reflections

2316 independent reflections

1860 reflections with $I > 2\sigma(I)$

$R_{\text{int}} = 0.063$
 $\theta_{\text{max}} = 64.7^\circ$, $\theta_{\text{min}} = 6.4^\circ$
 $h = -15 \rightarrow 15$

$k = -15 \rightarrow 16$
 $l = -8 \rightarrow 6$

Refinement

Refinement on F^2
 Least-squares matrix: full
 $R[F^2 > 2\sigma(F^2)] = 0.053$
 $wR(F^2) = 0.141$
 $S = 1.06$
 2316 reflections
 200 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
 $w = 1/[\sigma^2(F_o^2) + (0.0726P)^2 + 0.2459P]$
 where $P = (F_o^2 + 2F_c^2)/3$
 $(\Delta/\sigma)_{\text{max}} < 0.001$
 $\Delta\rho_{\text{max}} = 0.43 \text{ e } \text{\AA}^{-3}$
 $\Delta\rho_{\text{min}} = -0.40 \text{ e } \text{\AA}^{-3}$

Special details

Geometry. Bond distances, angles *etc.* have been calculated using the rounded fractional coordinates. All su's are estimated from the variances of the (full) variance-covariance matrix. The cell e.s.d.'s are taken into account in the estimation of distances, angles and torsion angles

Refinement. Refinement on F^2 for ALL reflections except those flagged by the user for potential systematic errors. Weighted R -factors wR and all goodnesses 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 observed criterion of $F^2 > \sigma(F^2)$ is used only for calculating $-R$ -factor-obs *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	$U_{\text{iso}}^*/U_{\text{eq}}$
S1	0.51842 (5)	0.17833 (5)	0.23250 (10)	0.0214 (2)
F13	0.10569 (13)	-0.01265 (15)	0.0632 (4)	0.0518 (8)
F14	0.14077 (13)	0.11887 (16)	-0.0469 (3)	0.0450 (7)
F15	0.11752 (13)	0.10360 (17)	0.2439 (3)	0.0484 (8)
N11	0.64047 (19)	-0.14957 (19)	0.3600 (4)	0.0297 (9)
C2	0.4251 (2)	0.0996 (2)	0.1944 (4)	0.0202 (9)
C3	0.3272 (2)	0.1178 (2)	0.1485 (4)	0.0212 (9)
C4	0.2634 (2)	0.0444 (2)	0.1373 (4)	0.0210 (9)
C5	0.2957 (2)	-0.0459 (2)	0.1679 (4)	0.0217 (9)
C6	0.3920 (2)	-0.0643 (2)	0.2103 (4)	0.0203 (8)
C7	0.4583 (2)	0.0087 (2)	0.2249 (4)	0.0178 (8)
C8	0.5601 (2)	0.0069 (2)	0.2816 (4)	0.0174 (8)
C9	0.6018 (2)	0.0928 (2)	0.2949 (4)	0.0183 (8)
C10	0.6079 (2)	-0.0785 (2)	0.3267 (4)	0.0204 (9)
C12	0.1574 (2)	0.0636 (2)	0.0993 (4)	0.0253 (9)
C16	0.7014 (2)	0.1180 (2)	0.3613 (4)	0.0182 (8)
C17	0.7798 (2)	0.0588 (2)	0.3430 (4)	0.0198 (9)
C18	0.8717 (2)	0.0826 (2)	0.4144 (4)	0.0236 (9)
C19	0.8890 (2)	0.1655 (2)	0.5074 (4)	0.0235 (9)
C20	0.8107 (2)	0.2255 (2)	0.5220 (4)	0.0217 (9)
C21	0.7190 (2)	0.2025 (2)	0.4506 (4)	0.0194 (8)
C22	0.9878 (2)	0.1887 (2)	0.5939 (5)	0.0306 (10)

H3	0.30560	0.17760	0.12600	0.0250*
H5	0.25120	-0.09410	0.15940	0.0260*
H6	0.41310	-0.12450	0.22910	0.0240*
H17	0.77040	0.00260	0.28230	0.0240*
H18	0.92320	0.04210	0.39990	0.0280*
H20	0.82050	0.28200	0.58110	0.0260*
H21	0.66800	0.24390	0.46190	0.0230*
H22A	0.99640	0.15870	0.71340	0.0460*
H22B	0.99300	0.25410	0.61070	0.0460*
H22C	1.03700	0.16810	0.51280	0.0460*

Atomic displacement parameters (Å²)

	U^{11}	U^{22}	U^{33}	U^{12}	U^{13}	U^{23}
S1	0.0194 (4)	0.0164 (4)	0.0281 (4)	-0.0001 (3)	-0.0009 (3)	0.0010 (3)
F13	0.0215 (9)	0.0385 (13)	0.0941 (17)	-0.0028 (9)	-0.0107 (10)	-0.0047 (12)
F14	0.0293 (10)	0.0607 (15)	0.0442 (12)	0.0123 (10)	-0.0052 (9)	0.0178 (11)
F15	0.0285 (10)	0.0782 (17)	0.0390 (11)	0.0161 (10)	0.0059 (9)	-0.0160 (11)
N11	0.0321 (15)	0.0196 (16)	0.0369 (16)	0.0022 (12)	-0.0027 (12)	0.0019 (13)
C2	0.0237 (15)	0.0198 (16)	0.0171 (14)	-0.0014 (12)	0.0024 (11)	-0.0003 (13)
C3	0.0214 (14)	0.0211 (17)	0.0210 (14)	0.0051 (12)	0.0011 (12)	-0.0005 (13)
C4	0.0223 (15)	0.0258 (17)	0.0149 (13)	-0.0007 (13)	0.0005 (11)	-0.0005 (13)
C5	0.0244 (15)	0.0211 (17)	0.0196 (14)	-0.0038 (12)	-0.0001 (12)	-0.0010 (13)
C6	0.0268 (15)	0.0145 (15)	0.0198 (14)	-0.0004 (12)	0.0026 (12)	0.0003 (12)
C7	0.0216 (14)	0.0207 (16)	0.0111 (13)	0.0011 (12)	0.0023 (11)	0.0016 (12)
C8	0.0203 (14)	0.0170 (15)	0.0149 (13)	0.0012 (12)	0.0017 (11)	-0.0007 (12)
C9	0.0220 (14)	0.0191 (16)	0.0142 (13)	0.0006 (12)	0.0041 (11)	-0.0006 (12)
C10	0.0199 (14)	0.0220 (18)	0.0191 (15)	-0.0039 (13)	-0.0011 (12)	-0.0006 (13)
C12	0.0240 (15)	0.0254 (18)	0.0265 (16)	0.0002 (13)	0.0002 (13)	-0.0017 (14)
C16	0.0206 (14)	0.0174 (16)	0.0166 (13)	-0.0004 (12)	0.0005 (11)	0.0015 (12)
C17	0.0219 (15)	0.0169 (16)	0.0209 (14)	-0.0001 (12)	0.0034 (12)	-0.0013 (13)
C18	0.0229 (15)	0.0240 (17)	0.0241 (15)	0.0042 (13)	0.0026 (12)	-0.0007 (13)
C19	0.0237 (15)	0.0293 (18)	0.0175 (14)	-0.0019 (13)	0.0003 (12)	0.0033 (13)
C20	0.0240 (15)	0.0223 (17)	0.0187 (14)	-0.0044 (13)	0.0011 (12)	-0.0009 (13)
C21	0.0204 (14)	0.0186 (16)	0.0195 (14)	0.0003 (12)	0.0032 (11)	0.0022 (13)
C22	0.0251 (16)	0.037 (2)	0.0295 (16)	0.0014 (14)	-0.0015 (13)	-0.0015 (16)

Geometric parameters (Å, °)

S1—C2	1.731 (3)	C16—C17	1.391 (4)
S1—C9	1.735 (3)	C16—C21	1.399 (4)
F13—C12	1.335 (4)	C17—C18	1.384 (4)
F14—C12	1.328 (4)	C18—C19	1.390 (4)
F15—C12	1.326 (4)	C19—C20	1.394 (4)
N11—C10	1.146 (4)	C19—C22	1.503 (4)
C2—C3	1.396 (4)	C20—C21	1.378 (4)
C2—C7	1.411 (4)	C3—H3	0.9300
C3—C4	1.381 (4)	C5—H5	0.9300

C4—C5	1.399 (4)	C6—H6	0.9300
C4—C12	1.496 (4)	C17—H17	0.9300
C5—C6	1.371 (4)	C18—H18	0.9300
C6—C7	1.400 (4)	C20—H20	0.9300
C7—C8	1.439 (4)	C21—H21	0.9300
C8—C9	1.375 (4)	C22—H22A	0.9600
C8—C10	1.434 (4)	C22—H22B	0.9600
C9—C16	1.474 (4)	C22—H22C	0.9600
C2—S1—C9	92.44 (14)	C17—C16—C21	117.9 (3)
S1—C2—C3	127.7 (2)	C16—C17—C18	120.6 (3)
S1—C2—C7	111.3 (2)	C17—C18—C19	121.6 (3)
C3—C2—C7	121.0 (3)	C18—C19—C20	117.6 (3)
C2—C3—C4	118.1 (3)	C18—C19—C22	121.4 (3)
C3—C4—C5	121.3 (3)	C20—C19—C22	121.0 (3)
C3—C4—C12	118.5 (3)	C19—C20—C21	121.2 (3)
C5—C4—C12	120.2 (3)	C16—C21—C20	121.1 (3)
C4—C5—C6	120.9 (3)	C2—C3—H3	121.00
C5—C6—C7	119.2 (3)	C4—C3—H3	121.00
C2—C7—C6	119.6 (3)	C4—C5—H5	120.00
C2—C7—C8	111.3 (2)	C6—C5—H5	120.00
C6—C7—C8	129.0 (3)	C5—C6—H6	120.00
C7—C8—C9	113.6 (3)	C7—C6—H6	120.00
C7—C8—C10	120.6 (3)	C16—C17—H17	120.00
C9—C8—C10	125.8 (3)	C18—C17—H17	120.00
S1—C9—C8	111.4 (2)	C17—C18—H18	119.00
S1—C9—C16	119.8 (2)	C19—C18—H18	119.00
C8—C9—C16	128.7 (3)	C19—C20—H20	119.00
N11—C10—C8	175.6 (3)	C21—C20—H20	119.00
F13—C12—F14	106.3 (2)	C16—C21—H21	120.00
F13—C12—F15	106.1 (2)	C20—C21—H21	119.00
F13—C12—C4	112.7 (2)	C19—C22—H22A	109.00
F14—C12—F15	106.5 (2)	C19—C22—H22B	109.00
F14—C12—C4	112.6 (2)	C19—C22—H22C	109.00
F15—C12—C4	112.2 (2)	H22A—C22—H22B	109.00
C9—C16—C17	121.9 (3)	H22A—C22—H22C	110.00
C9—C16—C21	120.1 (3)	H22B—C22—H22C	109.00
C9—S1—C2—C3	175.6 (3)	C2—C7—C8—C9	0.5 (4)
C9—S1—C2—C7	-1.4 (2)	C2—C7—C8—C10	177.7 (3)
C2—S1—C9—C8	1.7 (2)	C6—C7—C8—C9	-175.0 (3)
C2—S1—C9—C16	-175.3 (2)	C6—C7—C8—C10	2.2 (5)
S1—C2—C3—C4	-175.5 (2)	C7—C8—C9—S1	-1.6 (3)
C7—C2—C3—C4	1.2 (4)	C7—C8—C9—C16	175.2 (3)
S1—C2—C7—C6	176.7 (2)	C10—C8—C9—S1	-178.6 (2)
S1—C2—C7—C8	0.8 (3)	C10—C8—C9—C16	-1.8 (5)
C3—C2—C7—C6	-0.5 (4)	S1—C9—C16—C17	-154.1 (2)
C3—C2—C7—C8	-176.4 (3)	S1—C9—C16—C21	27.8 (4)

C2—C3—C4—C5	-1.0 (4)	C8—C9—C16—C17	29.4 (5)
C2—C3—C4—C12	176.5 (3)	C8—C9—C16—C21	-148.7 (3)
C3—C4—C5—C6	0.1 (4)	C9—C16—C17—C18	-177.0 (3)
C12—C4—C5—C6	-177.4 (3)	C21—C16—C17—C18	1.2 (4)
C3—C4—C12—F13	170.3 (3)	C9—C16—C21—C20	176.8 (3)
C3—C4—C12—F14	50.1 (4)	C17—C16—C21—C20	-1.5 (4)
C3—C4—C12—F15	-70.0 (3)	C16—C17—C18—C19	0.4 (4)
C5—C4—C12—F13	-12.2 (4)	C17—C18—C19—C20	-1.8 (4)
C5—C4—C12—F14	-132.4 (3)	C17—C18—C19—C22	176.6 (3)
C5—C4—C12—F15	107.5 (3)	C18—C19—C20—C21	1.5 (4)
C4—C5—C6—C7	0.7 (4)	C22—C19—C20—C21	-176.9 (3)
C5—C6—C7—C2	-0.5 (4)	C19—C20—C21—C16	0.1 (4)
C5—C6—C7—C8	174.6 (3)		

Hydrogen-bond geometry (Å, °)

<i>D—H...A</i>	<i>D—H</i>	<i>H...A</i>	<i>D...A</i>	<i>D—H...A</i>
C3—H3...N11 ⁱ	0.93	2.62	3.411 (4)	143
C22—H22C...F15 ⁱⁱ	0.96	2.45	3.375 (4)	162

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