

Acta Crystallographica Section E

## Structure Reports

Online

ISSN 1600-5368

Ethyl 4-[(3,5-di-*tert*-butyl-2-hydroxybenzyl)amino]benzoate

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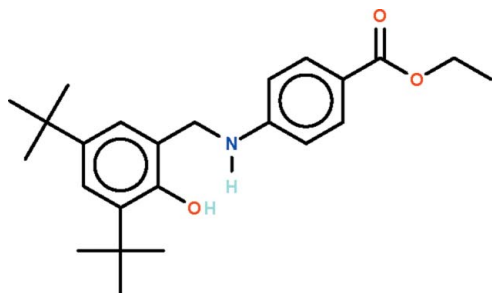
Received 6 October 2010; accepted 11 October 2010

Key indicators: single-crystal X-ray study;  $T = 100$  K; mean  $\sigma(\text{C}-\text{C}) = 0.004$  Å; disorder in main residue;  $R$  factor = 0.058;  $wR$  factor = 0.154; data-to-parameter ratio = 13.9.

The title amine,  $\text{C}_{24}\text{H}_{33}\text{NO}_3$ , has two substituted aromatic rings at either end of the  $-\text{CH}_2\text{NH}-$  linkage [ $\text{C}_{\text{aryl}}-\text{CH}_2-\text{NH}-\text{C}_{\text{aryl}}$  torsion angle =  $77.4(3)^\circ$ ]. The amino and hydroxy groups are hydrogen-bond donors to the carbonyl O atom of an adjacent molecule. These hydrogen bonds give rise to a chain that runs along the  $b$  axis. One of the *tert*-butyl groups is disordered over two positions with a site-occupation factor of 0.834 (6) for the major occupied site.

## Related literature

For the structure of the Schiff-base reactant, see: Shakir *et al.* (2010).



## Experimental

## Crystal data

$\text{C}_{24}\text{H}_{33}\text{NO}_3$   
 $M_r = 383.51$   
 Monoclinic,  $P2_1/c$   
 $a = 17.788(3)$  Å  
 $b = 8.9872(14)$  Å  
 $c = 14.235(2)$  Å  
 $\beta = 101.414(2)^\circ$   
 $V = 2230.7(6)$  Å<sup>3</sup>  
 $Z = 4$   
 Mo  $K\alpha$  radiation  
 $\mu = 0.07$  mm<sup>-1</sup>  
 $T = 100$  K  
 $0.30 \times 0.10 \times 0.03$  mm

## Data collection

Bruker SMART APEX  
 diffractometer  
 17094 measured reflections  
 3928 independent reflections  
 2428 reflections with  $I > 2\sigma(I)$   
 $R_{\text{int}} = 0.092$

## Refinement

$R[F^2 > 2\sigma(F^2)] = 0.058$   
 $wR(F^2) = 0.154$   
 $S = 1.03$   
 3928 reflections  
 283 parameters  
 45 restraints  
 H-atom parameters constrained  
 $\Delta\rho_{\text{max}} = 0.24$  e Å<sup>-3</sup>  
 $\Delta\rho_{\text{min}} = -0.23$  e Å<sup>-3</sup>

Table 1

Hydrogen-bond geometry (Å, °).

$D-H\cdots A$	$D-H$	$H\cdots A$	$D\cdots A$	$D-H\cdots A$
$\text{O1}-\text{H1o}\cdots\text{O2}^i$	0.84	2.13	2.909 (3)	154
$\text{N1}-\text{H1n}\cdots\text{O2}^i$	0.88	2.07	2.827 (3)	143

Symmetry code: (i)  $x, y + 1, z$ .

Data collection: *APEX2* (Bruker, 2009); cell refinement: *SAINTE* (Bruker, 2009); data reduction: *SAINTE*; program(s) used to solve structure: *SHELXS97* (Sheldrick, 2008); program(s) used to refine structure: *SHELXL97* (Sheldrick, 2008); molecular graphics: *X-SEED* (Barbour, 2001); software used to prepare material for publication: *publCIF* (Westrip, 2010).

We thank the University of Malaya for supporting this study.

Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: BT5371).

## References

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**supplementary materials**

*Acta Cryst.* (2010). E66, o2916 [ doi:10.1107/S1600536810040742 ]

## Ethyl 4-[(3,5-di-*tert*-butyl-2-hydroxybenzyl)amino]benzoate

R. M. Shakir, A. Ariffin and S. W. Ng

### Comment

The Schiff base, ethyl 4-(di-3,5-*tert*-butyl-2-hydroxybenzylideneamino)benzoate, is an intensely orange-colored compound whose color can be attributed to the azomethine double-bond; in the crystal structure, the two aromatic rings are aligned at 24.9 (1)° (Shakir *et al.*, 2010). The reduction of this compound affords the title colorless compound (Scheme I, Fig. 1), which is a secondary amine. The amino and hydroxy groups are hydrogen-bond donors to the carbonyl oxygen atom of an adjacent molecule; the hydrogen bonds give rise to a chain that runs along the *b*-axis. One of the *tert*-butyl groups is disordered over two positions with respect to the methyl units in an 83.6 (6):16.4 (6) ratio.

### Experimental

Ethyl 4-aminobenzoate (0.35 g) dissolved in ethanol (5 ml) was added to 3,5-di-*tert*-butyl-2-hydroxybenzaldehyde (0.5 g) dissolved in ethanol (20 ml). Several drops of acetic acid were added. The solution was heated for 3 h. The solvent was evaporated and the product recrystallized from ethanol to yield orange prisms in 80% yield. The compound (0.5 g) was dissolved in methanol-THF (1:1) (20 ml). The solution was cooled to 273 K and three equivalents of sodium tetraborate were added. The mixture was stirred for 4 h until the orange color was discharged. The mixture was filtered and the solvent evaporated. The colorless compound was recrystallized from a hexane-ethanol mixture (yield 90%).

### Refinement

Carbon-bound H-atoms were placed in calculated positions (C—H 0.95–0.99 Å) and were included in the refinement in the riding model approximation, with  $U(\text{H})$  set to 1.2–15 $U(\text{C})$ .

The amino and hydroxy H-atoms were also placed in calculated positions and their displacement parameters were also similarly tied (N—H 0.88 Å and O—H 0.84 Å).

One of the *tert*-butyl groups is disordered over two positions with respect to the methyl units in an 83.4 (6):16.6 (6) ratio; the C—C<sub>methyl</sub> distances were restrained to 1.50±0.01 Å; the anisotropic displacement parameters of the six methyl C-atoms were restrained to be nearly isotropic. By the use of a *SAME* instruction, the bonds involving the C12', C13' and C14' were restrained to the same length as those involving the unprimed ones. In addition, the 1,3-related distances were restrained to be equal in both disordered moieties.

### Figures

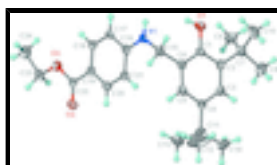


Fig. 1. Anisotropic displacement ellipsoid plot (Barbour, 2001) of  $\text{C}_{24}\text{H}_{33}\text{NO}_3$  at the 70% probability level; hydrogen atoms are drawn as spheres of arbitrary radius. The disorder in one of the *tert*-butyl groups is not shown.

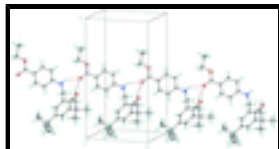


Fig. 2. Hydrogen-bonded chain motif.

## Ethyl 4-[(3,5-di-*tert*-butyl-2-hydroxybenzyl)amino]benzoate

### Crystal data

$C_{24}H_{33}NO_3$	$F(000) = 832$
$M_r = 383.51$	$D_x = 1.142 \text{ Mg m}^{-3}$
Monoclinic, $P2_1/c$	Mo $K\alpha$ radiation, $\lambda = 0.71073 \text{ \AA}$
Hall symbol: -P 2ybc	Cell parameters from 1466 reflections
$a = 17.788 (3) \text{ \AA}$	$\theta = 2.1\text{--}21.3^\circ$
$b = 8.9872 (14) \text{ \AA}$	$\mu = 0.07 \text{ mm}^{-1}$
$c = 14.235 (2) \text{ \AA}$	$T = 100 \text{ K}$
$\beta = 101.414 (2)^\circ$	Prism, colorless
$V = 2230.7 (6) \text{ \AA}^3$	$0.30 \times 0.10 \times 0.03 \text{ mm}$
$Z = 4$	

### Data collection

Bruker SMART APEX diffractometer	2428 reflections with $I > 2\sigma(I)$
Radiation source: fine-focus sealed tube	$R_{\text{int}} = 0.092$
graphite	$\theta_{\text{max}} = 25.0^\circ$ , $\theta_{\text{min}} = 2.3^\circ$
$\omega$ scans	$h = -20 \rightarrow 21$
17094 measured reflections	$k = -10 \rightarrow 10$
3928 independent reflections	$l = -16 \rightarrow 16$

### Refinement

Refinement on $F^2$	Secondary atom site location: difference Fourier map
Least-squares matrix: full	Hydrogen site location: inferred from neighbouring sites
$R[F^2 > 2\sigma(F^2)] = 0.058$	H-atom parameters constrained
$wR(F^2) = 0.154$	$w = 1/[\sigma^2(F_o^2) + (0.0662P)^2 + 0.7154P]$
$S = 1.03$	where $P = (F_o^2 + 2F_c^2)/3$
3928 reflections	$(\Delta/\sigma)_{\text{max}} = 0.001$
283 parameters	$\Delta\rho_{\text{max}} = 0.24 \text{ e \AA}^{-3}$
45 restraints	$\Delta\rho_{\text{min}} = -0.23 \text{ e \AA}^{-3}$
Primary atom site location: structure-invariant direct methods	Extinction correction: <i>SHELXL97</i> (Sheldrick, 2008), $F_c^* = kFc[1 + 0.001xFc^2\lambda^3/\sin(2\theta)]^{-1/4}$
	Extinction coefficient: 0.0079 (12)

Fractional atomic coordinates and isotropic or equivalent isotropic displacement parameters ( $\text{\AA}^2$ )

	x	y	z	$U_{\text{iso}}^*/U_{\text{eq}}$	Occ. (<1)
O1	0.32463 (11)	0.7157 (2)	0.42898 (14)	0.0281 (5)	
H1O	0.3613	0.7145	0.3998	0.042*	
O2	0.46605 (11)	-0.2040 (2)	0.36814 (13)	0.0257 (5)	
O3	0.58429 (11)	-0.1107 (2)	0.41969 (13)	0.0237 (5)	
N1	0.41764 (13)	0.5020 (2)	0.31779 (16)	0.0223 (6)	
H1N	0.4497	0.5690	0.3479	0.027*	
C1	0.26841 (16)	0.6193 (3)	0.38515 (19)	0.0206 (6)	
C2	0.20301 (16)	0.6027 (3)	0.4259 (2)	0.0224 (7)	
C3	0.14725 (16)	0.5024 (3)	0.3809 (2)	0.0246 (7)	
H3	0.1028	0.4890	0.4076	0.029*	
C4	0.15296 (16)	0.4208 (3)	0.2993 (2)	0.0227 (7)	
C5	0.21844 (16)	0.4441 (3)	0.2617 (2)	0.0227 (7)	
H5	0.2236	0.3920	0.2052	0.027*	
C6	0.27635 (15)	0.5401 (3)	0.30322 (19)	0.0197 (6)	
C7	0.19398 (17)	0.6857 (3)	0.5172 (2)	0.0272 (7)	
C8	0.11681 (17)	0.6522 (4)	0.5457 (2)	0.0388 (8)	
H8A	0.1126	0.5451	0.5564	0.058*	
H8B	0.0748	0.6841	0.4942	0.058*	
H8C	0.1138	0.7061	0.6047	0.058*	
C9	0.25655 (17)	0.6366 (4)	0.6014 (2)	0.0332 (8)	
H9A	0.2537	0.5286	0.6097	0.050*	
H9B	0.2490	0.6867	0.6600	0.050*	
H9C	0.3070	0.6631	0.5885	0.050*	
C10	0.1981 (2)	0.8544 (3)	0.5028 (2)	0.0399 (9)	
H10A	0.2468	0.8797	0.4842	0.060*	
H10B	0.1949	0.9054	0.5627	0.060*	
H10C	0.1554	0.8860	0.4522	0.060*	
C11	0.09420 (17)	0.3034 (3)	0.2554 (2)	0.0365 (8)	
C12	0.0664 (3)	0.3407 (7)	0.1489 (3)	0.077 (2)	0.834 (6)
H12A	0.0275	0.2683	0.1199	0.115*	0.834 (6)
H12B	0.1099	0.3368	0.1161	0.115*	0.834 (6)
H12C	0.0442	0.4408	0.1428	0.115*	0.834 (6)
C13	0.1347 (3)	0.1529 (4)	0.2641 (4)	0.0600 (17)	0.834 (6)
H13A	0.0980	0.0750	0.2377	0.090*	0.834 (6)
H13B	0.1557	0.1318	0.3317	0.090*	0.834 (6)
H13C	0.1765	0.1552	0.2284	0.090*	0.834 (6)
C14	0.0263 (2)	0.2976 (5)	0.3027 (4)	0.0452 (13)	0.834 (6)
H14A	-0.0094	0.2209	0.2719	0.068*	0.834 (6)
H14B	0.0005	0.3944	0.2964	0.068*	0.834 (6)
H14C	0.0433	0.2735	0.3707	0.068*	0.834 (6)
C12'	0.1079 (13)	0.210 (2)	0.1736 (14)	0.048 (7)	0.166 (6)
H12D	0.1621	0.2163	0.1694	0.073*	0.166 (6)
H12E	0.0761	0.2472	0.1139	0.073*	0.166 (6)
H12F	0.0944	0.1068	0.1839	0.073*	0.166 (6)
C13'	0.0927 (15)	0.184 (2)	0.3315 (15)	0.074 (9)	0.166 (6)

## supplementary materials

H13D	0.1373	0.1185	0.3355	0.111*	0.166 (6)
H13E	0.0456	0.1248	0.3140	0.111*	0.166 (6)
H13F	0.0941	0.2313	0.3938	0.111*	0.166 (6)
C14'	0.0130 (7)	0.361 (2)	0.2415 (18)	0.048 (7)	0.166 (6)
H14D	-0.0041	0.3589	0.3028	0.072*	0.166 (6)
H14E	-0.0208	0.2991	0.1948	0.072*	0.166 (6)
H14F	0.0114	0.4641	0.2180	0.072*	0.166 (6)
C15	0.34660 (15)	0.5559 (3)	0.25896 (19)	0.0225 (7)	
H15A	0.3374	0.5013	0.1974	0.027*	
H15B	0.3531	0.6623	0.2445	0.027*	
C16	0.43888 (16)	0.3562 (3)	0.33031 (19)	0.0203 (6)	
C17	0.51327 (15)	0.3219 (3)	0.37876 (19)	0.0205 (6)	
H17	0.5482	0.4004	0.4004	0.025*	
C18	0.53670 (16)	0.1770 (3)	0.39565 (19)	0.0204 (6)	
H18	0.5876	0.1565	0.4282	0.024*	
C19	0.48608 (16)	0.0590 (3)	0.36514 (19)	0.0202 (6)	
C20	0.41230 (16)	0.0918 (3)	0.31703 (19)	0.0220 (7)	
H20	0.3774	0.0130	0.2960	0.026*	
C21	0.38877 (16)	0.2380 (3)	0.29915 (18)	0.0208 (6)	
H21	0.3382	0.2581	0.2654	0.025*	
C22	0.50943 (15)	-0.0962 (3)	0.38371 (18)	0.0194 (6)	
C23	0.61300 (17)	-0.2598 (3)	0.4479 (2)	0.0245 (7)	
H23A	0.6072	-0.3259	0.3912	0.029*	
H23B	0.5846	-0.3037	0.4944	0.029*	
C24	0.69625 (17)	-0.2403 (3)	0.4929 (2)	0.0320 (8)	
H24A	0.7188	-0.3374	0.5133	0.048*	
H24B	0.7009	-0.1744	0.5486	0.048*	
H24C	0.7234	-0.1964	0.4460	0.048*	

### Atomic displacement parameters ( $\text{\AA}^2$ )

	$U^{11}$	$U^{22}$	$U^{33}$	$U^{12}$	$U^{13}$	$U^{23}$
O1	0.0285 (12)	0.0281 (11)	0.0296 (12)	-0.0076 (9)	0.0102 (9)	-0.0064 (9)
O2	0.0288 (12)	0.0204 (11)	0.0285 (11)	-0.0036 (9)	0.0072 (9)	-0.0018 (9)
O3	0.0254 (11)	0.0184 (10)	0.0271 (11)	0.0027 (8)	0.0049 (9)	0.0050 (8)
N1	0.0229 (13)	0.0186 (13)	0.0255 (13)	0.0000 (10)	0.0045 (11)	-0.0029 (10)
C1	0.0232 (16)	0.0138 (14)	0.0229 (15)	0.0011 (12)	0.0001 (12)	-0.0007 (12)
C2	0.0201 (16)	0.0216 (15)	0.0254 (15)	0.0033 (12)	0.0044 (12)	0.0012 (12)
C3	0.0204 (16)	0.0246 (16)	0.0295 (17)	0.0006 (12)	0.0068 (13)	0.0009 (13)
C4	0.0191 (16)	0.0217 (15)	0.0252 (16)	0.0030 (12)	-0.0009 (13)	-0.0025 (12)
C5	0.0265 (17)	0.0197 (15)	0.0209 (15)	0.0051 (12)	0.0025 (13)	-0.0006 (12)
C6	0.0225 (16)	0.0176 (14)	0.0189 (14)	0.0036 (12)	0.0035 (12)	0.0029 (12)
C7	0.0275 (17)	0.0290 (17)	0.0277 (16)	0.0003 (13)	0.0114 (14)	-0.0037 (13)
C8	0.0319 (19)	0.051 (2)	0.0372 (19)	-0.0004 (16)	0.0151 (16)	-0.0107 (16)
C9	0.0332 (19)	0.0406 (19)	0.0275 (17)	-0.0082 (15)	0.0101 (14)	-0.0066 (15)
C10	0.047 (2)	0.0295 (18)	0.047 (2)	0.0008 (15)	0.0194 (17)	-0.0112 (16)
C11	0.0273 (18)	0.0364 (19)	0.044 (2)	-0.0023 (15)	0.0038 (16)	-0.0094 (16)
C12	0.063 (3)	0.113 (5)	0.047 (3)	-0.052 (3)	-0.009 (2)	-0.010 (3)

C13	0.045 (3)	0.033 (2)	0.106 (4)	-0.014 (2)	0.025 (3)	-0.032 (3)
C14	0.028 (2)	0.042 (3)	0.068 (3)	-0.0133 (19)	0.016 (2)	-0.016 (2)
C12'	0.043 (10)	0.047 (10)	0.053 (10)	-0.007 (8)	0.004 (8)	-0.026 (8)
C13'	0.078 (12)	0.063 (11)	0.078 (12)	-0.014 (9)	0.007 (9)	-0.005 (9)
C14'	0.046 (10)	0.050 (10)	0.048 (10)	-0.009 (8)	0.007 (8)	-0.012 (8)
C15	0.0259 (17)	0.0199 (15)	0.0218 (15)	0.0045 (12)	0.0047 (13)	0.0009 (12)
C16	0.0259 (16)	0.0195 (15)	0.0174 (14)	0.0004 (12)	0.0090 (12)	-0.0002 (12)
C17	0.0234 (16)	0.0173 (14)	0.0220 (15)	-0.0040 (12)	0.0075 (13)	-0.0046 (12)
C18	0.0201 (15)	0.0225 (15)	0.0191 (14)	0.0006 (12)	0.0054 (12)	-0.0018 (12)
C19	0.0218 (16)	0.0220 (15)	0.0181 (14)	0.0009 (12)	0.0072 (12)	-0.0004 (12)
C20	0.0244 (16)	0.0193 (15)	0.0234 (15)	-0.0033 (12)	0.0075 (13)	-0.0012 (12)
C21	0.0209 (15)	0.0245 (16)	0.0177 (15)	0.0000 (12)	0.0058 (12)	-0.0026 (12)
C22	0.0193 (16)	0.0228 (15)	0.0176 (14)	-0.0010 (13)	0.0070 (12)	-0.0016 (12)
C23	0.0359 (18)	0.0181 (15)	0.0212 (15)	0.0056 (13)	0.0095 (13)	0.0042 (12)
C24	0.0344 (19)	0.0292 (17)	0.0323 (18)	0.0076 (14)	0.0065 (15)	0.0098 (14)

*Geometric parameters (Å, °)*

O1—C1	1.376 (3)	C12—H12A	0.9800
O1—H1O	0.8400	C12—H12B	0.9800
O2—C22	1.231 (3)	C12—H12C	0.9800
O3—C22	1.335 (3)	C13—H13A	0.9800
O3—C23	1.462 (3)	C13—H13B	0.9800
N1—C16	1.366 (3)	C13—H13C	0.9800
N1—C15	1.454 (3)	C14—H14A	0.9800
N1—H1N	0.8800	C14—H14B	0.9800
C1—C6	1.397 (4)	C14—H14C	0.9800
C1—C2	1.407 (4)	C12'—H12D	0.9800
C2—C3	1.398 (4)	C12'—H12E	0.9800
C2—C7	1.534 (4)	C12'—H12F	0.9800
C3—C4	1.395 (4)	C13'—H13D	0.9800
C3—H3	0.9500	C13'—H13E	0.9800
C4—C5	1.391 (4)	C13'—H13F	0.9800
C4—C11	1.529 (4)	C14'—H14D	0.9800
C5—C6	1.384 (4)	C14'—H14E	0.9800
C5—H5	0.9500	C14'—H14F	0.9800
C6—C15	1.513 (4)	C15—H15A	0.9900
C7—C10	1.534 (4)	C15—H15B	0.9900
C7—C9	1.531 (4)	C16—C17	1.400 (4)
C7—C8	1.536 (4)	C16—C21	1.401 (4)
C8—H8A	0.9800	C17—C18	1.374 (4)
C8—H8B	0.9800	C17—H17	0.9500
C8—H8C	0.9800	C18—C19	1.403 (4)
C9—H9A	0.9800	C18—H18	0.9500
C9—H9B	0.9800	C19—C20	1.387 (4)
C9—H9C	0.9800	C19—C22	1.465 (4)
C10—H10A	0.9800	C20—C21	1.387 (4)
C10—H10B	0.9800	C20—H20	0.9500
C10—H10C	0.9800	C21—H21	0.9500

## supplementary materials

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C11—C12'	1.491 (9)	C23—C24	1.503 (4)
C11—C14	1.495 (5)	C23—H23A	0.9900
C11—C14'	1.511 (9)	C23—H23B	0.9900
C11—C13	1.526 (5)	C24—H24A	0.9800
C11—C13'	1.531 (10)	C24—H24B	0.9800
C11—C12	1.535 (5)	C24—H24C	0.9800
C1—O1—H1O	109.5	H13A—C13—H13B	109.5
C22—O3—C23	117.5 (2)	C11—C13—H13C	109.5
C16—N1—C15	125.5 (2)	H13A—C13—H13C	109.5
C16—N1—H1N	117.2	H13B—C13—H13C	109.5
C15—N1—H1N	117.2	C11—C14—H14A	109.5
O1—C1—C6	121.0 (3)	C11—C14—H14B	109.5
O1—C1—C2	117.8 (2)	H14A—C14—H14B	109.5
C6—C1—C2	121.2 (2)	C11—C14—H14C	109.5
C3—C2—C1	116.7 (3)	H14A—C14—H14C	109.5
C3—C2—C7	121.1 (3)	H14B—C14—H14C	109.5
C1—C2—C7	122.1 (2)	C11—C12'—H12D	109.5
C4—C3—C2	123.9 (3)	C11—C12'—H12E	109.5
C4—C3—H3	118.0	H12D—C12'—H12E	109.5
C2—C3—H3	118.0	C11—C12'—H12F	109.5
C5—C4—C3	116.5 (2)	H12D—C12'—H12F	109.5
C5—C4—C11	120.2 (3)	H12E—C12'—H12F	109.5
C3—C4—C11	123.2 (3)	C11—C13'—H13D	109.5
C6—C5—C4	122.6 (3)	C11—C13'—H13E	109.5
C6—C5—H5	118.7	H13D—C13'—H13E	109.5
C4—C5—H5	118.7	C11—C13'—H13F	109.5
C5—C6—C1	119.0 (3)	H13D—C13'—H13F	109.5
C5—C6—C15	119.1 (2)	H13E—C13'—H13F	109.5
C1—C6—C15	121.9 (2)	C11—C14'—H14D	109.5
C10—C7—C9	110.0 (3)	C11—C14'—H14E	109.5
C10—C7—C8	107.4 (3)	H14D—C14'—H14E	109.5
C9—C7—C8	106.6 (2)	C11—C14'—H14F	109.5
C10—C7—C2	110.6 (2)	H14D—C14'—H14F	109.5
C9—C7—C2	110.0 (2)	H14E—C14'—H14F	109.5
C8—C7—C2	112.2 (2)	N1—C15—C6	115.1 (2)
C7—C8—H8A	109.5	N1—C15—H15A	108.5
C7—C8—H8B	109.5	C6—C15—H15A	108.5
H8A—C8—H8B	109.5	N1—C15—H15B	108.5
C7—C8—H8C	109.5	C6—C15—H15B	108.5
H8A—C8—H8C	109.5	H15A—C15—H15B	107.5
H8B—C8—H8C	109.5	N1—C16—C17	119.0 (2)
C7—C9—H9A	109.5	N1—C16—C21	123.0 (2)
C7—C9—H9B	109.5	C17—C16—C21	118.0 (2)
H9A—C9—H9B	109.5	C18—C17—C16	121.3 (2)
C7—C9—H9C	109.5	C18—C17—H17	119.4
H9A—C9—H9C	109.5	C16—C17—H17	119.4
H9B—C9—H9C	109.5	C17—C18—C19	120.6 (3)
C7—C10—H10A	109.5	C17—C18—H18	119.7
C7—C10—H10B	109.5	C19—C18—H18	119.7



H10A—C10—H10B	109.5	C20—C19—C18	118.6 (2)
C7—C10—H10C	109.5	C20—C19—C22	119.9 (2)
H10A—C10—H10C	109.5	C18—C19—C22	121.5 (2)
H10B—C10—H10C	109.5	C19—C20—C21	120.9 (3)
C12'—C11—C14'	113.0 (13)	C19—C20—H20	119.5
C14—C11—C13	110.1 (3)	C21—C20—H20	119.5
C12'—C11—C13'	100.8 (15)	C20—C21—C16	120.7 (3)
C14'—C11—C13'	100.6 (13)	C20—C21—H21	119.7
C12'—C11—C4	120.1 (9)	C16—C21—H21	119.7
C14—C11—C4	112.8 (3)	O2—C22—O3	122.3 (2)
C14'—C11—C4	111.9 (9)	O2—C22—C19	124.9 (2)
C13—C11—C4	107.7 (3)	O3—C22—C19	112.8 (2)
C13'—C11—C4	107.6 (10)	O3—C23—C24	105.7 (2)
C14—C11—C12	108.7 (3)	O3—C23—H23A	110.6
C13—C11—C12	109.1 (4)	C24—C23—H23A	110.6
C4—C11—C12	108.3 (3)	O3—C23—H23B	110.6
C11—C12—H12A	109.5	C24—C23—H23B	110.6
C11—C12—H12B	109.5	H23A—C23—H23B	108.7
H12A—C12—H12B	109.5	C23—C24—H24A	109.5
C11—C12—H12C	109.5	C23—C24—H24B	109.5
H12A—C12—H12C	109.5	H24A—C24—H24B	109.5
H12B—C12—H12C	109.5	C23—C24—H24C	109.5
C11—C13—H13A	109.5	H24A—C24—H24C	109.5
C11—C13—H13B	109.5	H24B—C24—H24C	109.5
O1—C1—C2—C3	-179.0 (2)	C5—C4—C11—C13	60.4 (4)
C6—C1—C2—C3	0.7 (4)	C3—C4—C11—C13	-115.5 (4)
O1—C1—C2—C7	-1.2 (4)	C5—C4—C11—C13'	116.0 (12)
C6—C1—C2—C7	178.5 (2)	C3—C4—C11—C13'	-59.9 (12)
C1—C2—C3—C4	-0.5 (4)	C5—C4—C11—C12	-57.5 (4)
C7—C2—C3—C4	-178.3 (3)	C3—C4—C11—C12	126.6 (4)
C2—C3—C4—C5	-0.6 (4)	C16—N1—C15—C6	77.4 (3)
C2—C3—C4—C11	175.5 (3)	C5—C6—C15—N1	-113.9 (3)
C3—C4—C5—C6	1.5 (4)	C1—C6—C15—N1	65.8 (3)
C11—C4—C5—C6	-174.7 (3)	C15—N1—C16—C17	171.6 (2)
C4—C5—C6—C1	-1.3 (4)	C15—N1—C16—C21	-10.2 (4)
C4—C5—C6—C15	178.4 (2)	N1—C16—C17—C18	178.1 (3)
O1—C1—C6—C5	179.8 (2)	C21—C16—C17—C18	-0.1 (4)
C2—C1—C6—C5	0.2 (4)	C16—C17—C18—C19	-0.5 (4)
O1—C1—C6—C15	0.1 (4)	C17—C18—C19—C20	0.5 (4)
C2—C1—C6—C15	-179.6 (2)	C17—C18—C19—C22	-179.0 (2)
C3—C2—C7—C10	-123.6 (3)	C18—C19—C20—C21	0.1 (4)
C1—C2—C7—C10	58.7 (3)	C22—C19—C20—C21	179.6 (3)
C3—C2—C7—C9	114.7 (3)	C19—C20—C21—C16	-0.7 (4)
C1—C2—C7—C9	-63.1 (3)	N1—C16—C21—C20	-177.4 (3)
C3—C2—C7—C8	-3.8 (4)	C17—C16—C21—C20	0.7 (4)
C1—C2—C7—C8	178.5 (3)	C23—O3—C22—O2	-4.9 (4)
C5—C4—C11—C12'	1.7 (13)	C23—O3—C22—C19	175.4 (2)
C3—C4—C11—C12'	-174.2 (12)	C20—C19—C22—O2	-7.5 (4)
C5—C4—C11—C14	-177.9 (3)	C18—C19—C22—O2	172.0 (3)

## supplementary materials

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C3—C4—C11—C14	6.2 (4)	C20—C19—C22—O3	172.2 (2)
C5—C4—C11—C14'	-134.4 (11)	C18—C19—C22—O3	-8.3 (4)
C3—C4—C11—C14'	49.7 (11)	C22—O3—C23—C24	-175.3 (2)

### *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>
O1—H1 <sub>o</sub> $\cdots$ O2 <sup>i</sup>	0.84	2.13	2.909 (3)	154
N1—H1 <sub>n</sub> $\cdots$ O2 <sup>i</sup>	0.88	2.07	2.827 (3)	143

Symmetry codes: (i) *x*, *y*+1, *z*.

Fig. 1

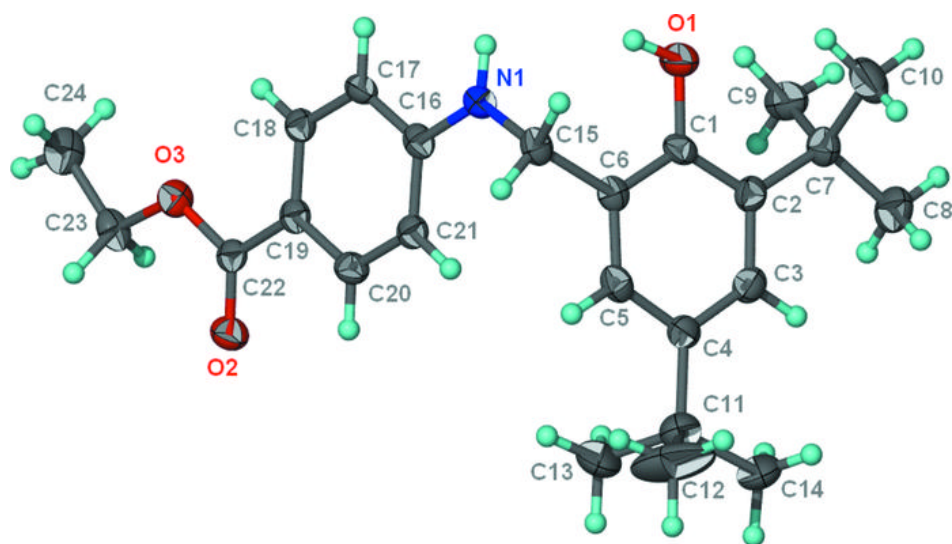


Fig. 2

