Additions & Corrections

Structural Studies of the Polymorphs of Carbamazepine, Its Dihydrate, and Two Solvates

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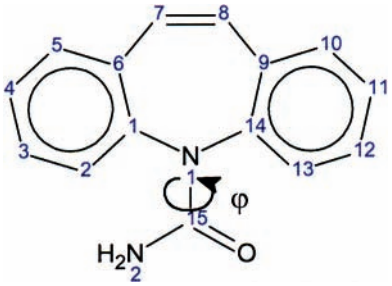
(*Org. Process Res. Dev.* **2005**, *9*, 902–910).

Table 4: There were minor notational errors in the originally published table, but none of the numbers were affected. The corrected version is below.

Table 4. **Some key molecular angles in the forms of carbamazepine[a]**



		intramolecular angles[a]			
structure	database code	φ^b	β^b	γ	δ
triclinic[c]	CBMZPN11 (O1)	−179.6	−17.7	55.0	70.7
	CBMZPN11 (O2)	−176.8	−15.2	54.1	71.2
	CBMZPN11 (O3)	−175.7	−11.8	54.9	69.2
	CBMZPN11 (O4)	−176.7	−15.9	55.1	71.9
P-monoclinic	CBMZPN01	−170.6	+ 2.2	53.4	61.4
	CBMZPN02	−170.4	+ 1.8	53.3	62.9
	CBMZPN10	−170.7	+ 1.6	53.3	61.5
trigonal	CBMZPN03	−176.1	−13.4	55.2	70.3
C-monoclinic	CBMZPN12	−172.4	− 7.5	49.8	68.1
dihydrate	this work	179.6(3)	0.9(5)	55.8(1)	54.5(1)

[a] The angles are defined as follows (see also Figure 6): φ torsion angle C14−N1−C15−N2. β torsion angle C1−N1−C15−N2. γ dihedral angle between benzene rings. δ dihedral angle between N2, O1, C15, N1 and C1, C6, C9, C14 planes. [b] The signs are given in relation to Figure 6 and of the molecule as given in the asymmetric unit in the CIF file. [c] Four independent molecules.

Note Added in Proof: It has come to our notice since submitting this paper that another crystal structure (code FEFNOT01, CCDC-247132) for carbamazepine dihydrate has appeared in the latest version of the CCSD. This perpetuates the disordered orthorhombic case.

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