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Ab initio structure determination of nanocrystals of organic pharmaceutical compounds by electron diffraction at room temperature using a Timepix quantum area direct electron detector. Corrigendum

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Corrections are made to Table 1 in the article by van Genderen *et al.* [*Acta Cryst.* (2016), A**72**, 236–242].

In the article by van Genderen *et al.* (2016), the scattering factors used in the refinements and in the CIF file were for X-ray scattering rather than for electron scattering. The correct scattering factors have now been used and the statistics that were affected by this error (model statistics R_{complete} , R1 and wR2) have been recalculated.

This affects six entries in Table 1 of the original publication. The correct values are given here. The corrected CIF and supporting information are also made available.

Table 1 Corrected statistics.

	Carbamazepine	Nicotinic acid
Refinement statisti	cs	
$R_{\text{complete}}^{\dagger}$	31.8	37.7
R1 (%)	27.9	34.1
wR2 (%)	55.2	60.1

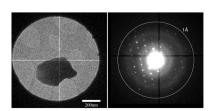
† Luebben & Gruene (2015).

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