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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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E. van Genderen,[a,b] M. T. B. Clabbers,[a,b] P. P. Das,[c] A. Stewart,[d] I. Nederlof,[a,e] K. C. Barentsen,[a] Q. Portillo,[c,f] N. S. Pannu,[a] S. Nicolopoulos,[c] T. Gruene[g]* and J. P. Abrahams[a,b,g]*

[a]Biophysical Structural Chemistry, Leiden University, Einsteinweg 55, 2333 CC Leiden, The Netherlands, [b]Center for Cellular Imaging and NanoAnalytics (C-CINA), Biozentrum, University of Basel, CH-4058 Basel, Switzerland, [c]Nanomegas SPRL, Boulevard Edmond Machtens 79, B 1080, Brussels, Belgium, [d]Department of Physics and Energy, Materials and Surface Science Institute (MSSI), University of Limerick, Limerick, Ireland, [e]Amsterdam Scientific Instruments, Postbus 41882, 1009 DB Amsterdam, The Netherlands, [f]Centres Científics i Tecnològics de la Universitat de Barcelona, University of Barcelona, Carrer de Lluís Solé i Sabaris, 1-3, Barcelona, Spain, and [g]Biology and Chemistry, Laboratory of Biomolecular Research, Paul Scherrer Institute (PSI), 5232 Villigen, Switzerland. *Correspondence e-mail: tim.gruene@psi.ch, jp.abrahams@unibas.ch

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 $R_{\rm complete}$, $R1$ 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 statistics | | |
| $R_{\rm complete}$† | 31.8 | 37.7 |
| $R1$ (%) | 27.9 | 34.1 |
| $wR2$ (%) | 55.2 | 60.1 |

† Luebben & Gruene (2015).

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## References

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