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). A72, 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 Rcomplete, 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.

<table>
<thead>
<tr>
<th></th>
<th>Carbamazepine</th>
<th>Nicotinic acid</th>
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<tr>
<td>Rcomplete</td>
<td>31.8</td>
<td>37.7</td>
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<tr>
<td>R1 (%)</td>
<td>27.9</td>
<td>34.1</td>
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<tr>
<td>wR2 (%)</td>
<td>55.2</td>
<td>60.1</td>
</tr>
</tbody>
</table>

† Luebben & Gruene (2015).

Acknowledgements

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References
