

## OPEN

**Author Correction: Deaminative chlorination of aminoheterocycles**Clément Ghiazza , Teresa Faber, Alejandro Gómez-Palomino and Josep Cornella Correction to: *Nature Chemistry* <https://doi.org/10.1038/s41557-021-00812-0>, published online 16 December 2021.

In the version of this Article originally published, there was an error in the structure of compound **31** in Table 1 (and pages 25, 158, 159 of the Supplementary Information) and in the structure of compound **58** in Fig. 3 (and pages 41, 252, 253 of the Supplementary Information). In the original version, the nitrogen and sulfur atoms in the benzisothiazole ring present in each structure were the wrong way around. This has now been corrected.

















**Open Access** This article is licensed under a Creative Commons Attribution 4.0 International License, which permits use, sharing, adaptation, distribution and reproduction in any medium or format, as long as you give appropriate credit to the original author(s) and the source, provide a link to the Creative Commons license, and indicate if changes were made. The images or other third party material in this article are included in the article's Creative Commons license, unless indicated otherwise in a credit line to the material. If material is not included in the article's Creative Commons license and your intended use is not permitted by statutory regulation or exceeds the permitted use, you will need to obtain permission directly from the copyright holder. To view a copy of this license, visit <http://creativecommons.org/licenses/by/4.0/>.

Published online: 30 March 2022

<https://doi.org/10.1038/s41557-022-00932-1>

© The Author(s), under exclusive licence to Springer Nature Limited 2022

**Publisher Correction: Porous covalent organic nanotubes and their assembly in loops and toroids**Kalipada Koner , Shayan Karak , Sharath Kandambeth , Suwendu Karak , Neethu Thomas , Luigi Leanza , Claudio Perego , Luca Pesce , Riccardo Capelli , Monika Moun , Monika Bhakar , Thalasseril G. Ajithkumar , Giovanni M. Pavan  and Rahul Banerjee Correction to: *Nature Chemistry* <https://doi.org/10.1038/s41557-022-00908-1>, published online 14 March 2022.

In the version of the article originally published, there was an error in Fig. 1b, where the word “or” was missing between the structures of DMDA and TA. The word has been restored in the HTML and PDF versions of the article.

Published online: 22 April 2022

<https://doi.org/10.1038/s41557-022-00952-x>

© The Author(s), under exclusive licence to Springer Nature Limited 2022