



<https://doi.org/10.1038/s41467-020-16706-7>

OPEN

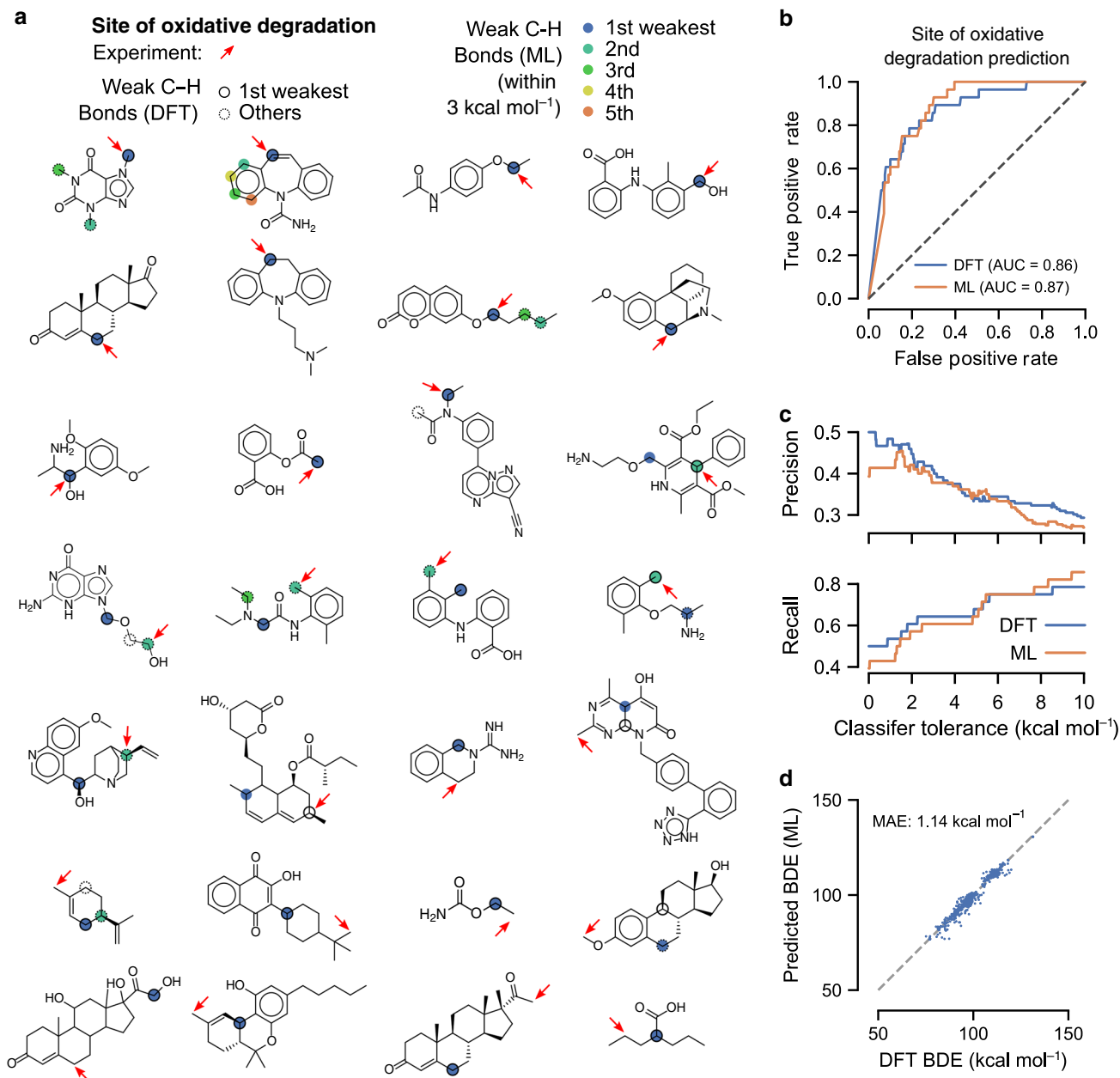
# Publisher Correction: Prediction of organic homolytic bond dissociation enthalpies at near chemical accuracy with sub-second computational cost

Peter C. St. John , Yanfei Guan , Yeonjoon Kim , Seonah Kim  & Robert S. Paton 

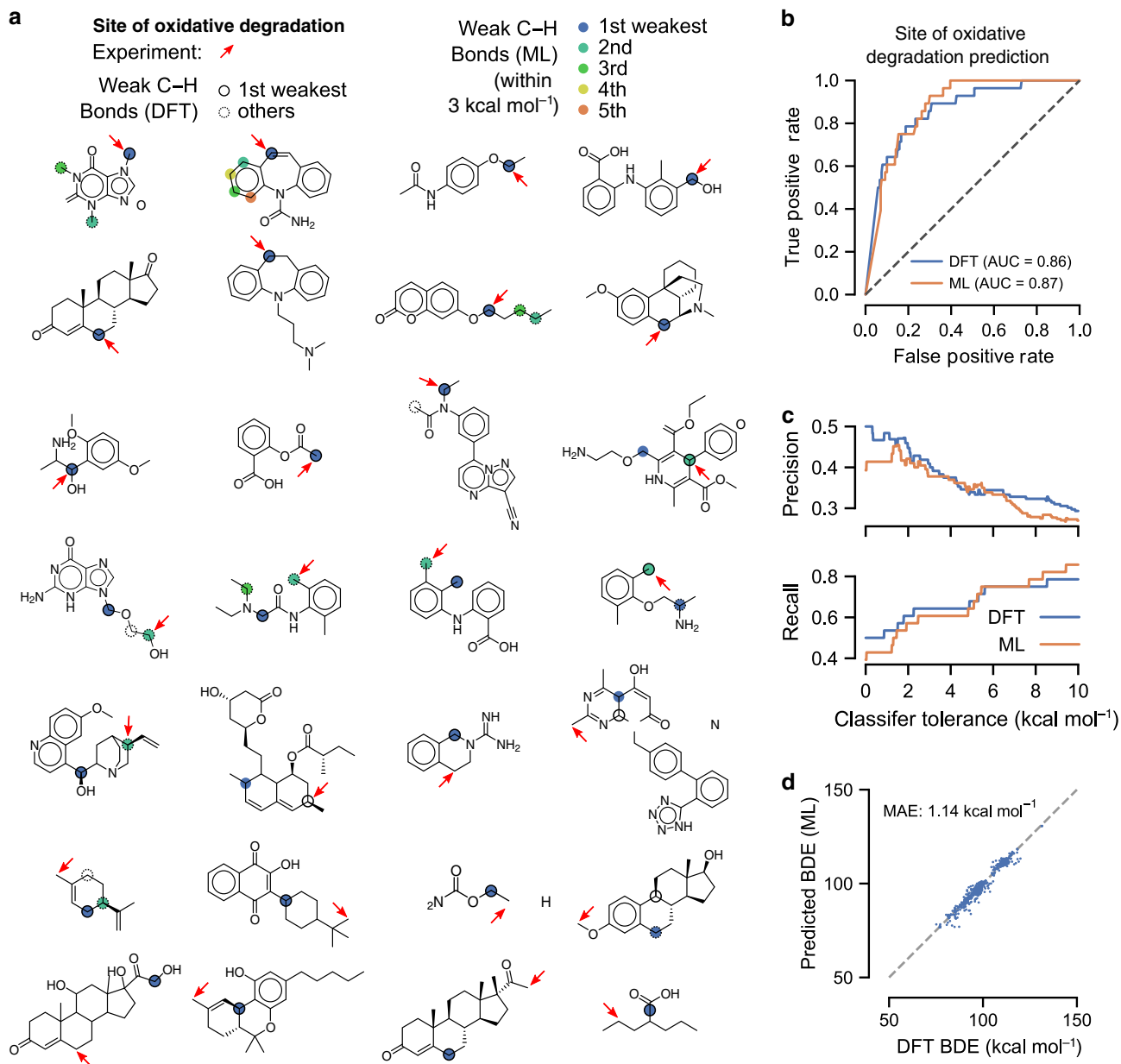
Correction to: *Nature Communications* <https://doi.org/10.1038/s41467-020-16201-z>, published online 11 May 2020.

The original version of this Article contained some errors in Fig. 7, in which there were four structures where one of the atoms was inadvertently displaced to the right of the structures.

The correct version of Fig. 7 is



which replaces the previous incorrect version from the original pdf of the paper.



This has been corrected in the PDF and HTML versions of the Article.

Published online: 11 June 2020

**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/>.

© The Author(s) 2020