Publisher Correction: Extending machine learning beyond interatomic potentials for predicting molecular properties

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Check for updates

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In the version of this article initially published, there was a typographical error in Fig. 5c, where the top-left formula now reading "$C_{10}H_2$" appeared initially as "$C_{10}H_{12}$". In the bottom section of Fig. 5c, the 4G-HDNNP result now presented was preceded in error by the 3G-HDNNP result from the same source article (Ko, T. W. et al. *Nat. Commun.* https://doi.org/10.1038/s41467-020-20427-2 (2021)). The figure has been updated in the HTML and PDF versions of the article.

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