

Correction to “Accurate and Numerically Efficient r^2 SCAN Meta-Generalized Gradient Approximation”

James W. Furness,* Aaron D. Kaplan, Jinliang Ning, John P. Perdew, and Jianwei Sun*

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The value of d_{p2} was incorrectly given in the main text. The correct value, for both the r^2 SCAN exchange and correlation energy functionals, is $d_{p2} = 0.361$ as given in the Supporting Information. We regret the error and thank F. Tran for bringing this to our attention.

The authors were made aware at the time of publication of calculations using the TASK exchange functional that demonstrate increased numeric efficiency [Hofmann, F. ; Kümmel, S. *J. Chem. Phys.* **2020**, *153*, 114106; DOI: [10.1063/5.0023657](https://doi.org/10.1063/5.0023657)]. We therefore regret our assertion that the modifications TASK proposes are not expected to increase numerical efficiency and instead refer interested readers to their work. We thank S. Kümmel for bringing this to our attention.

Neither error affects the ensuing material in the original publication nor its conclusions.

AUTHOR INFORMATION

Corresponding Authors

James W. Furness; orcid.org/0000-0003-3146-0977;

Email: jfurness@tulane.edu

Jianwei Sun; orcid.org/0000-0002-2361-6823;

Email: jsun@tulane.edu

Authors

Aaron D. Kaplan; orcid.org/0000-0003-3439-4856

Jinliang Ning; orcid.org/0000-0002-3691-5291

John P. Perdew

Complete contact information is available at:

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