Erratum: Nuclear Charge Radii of Silicon Isotopes [Phys. Rev. Lett. 132, 162502 (2024)]

Kristian König, Julian C. Berengut, Anastasia Borschevsky, Alex Brinson, B. Alex Brown, Adam Dockery, Serdar Elhatisari, Ephraim Eliav, Ronald F. Garcia Ruiz, Jason D. Holt, Bai-Shan Hu, Jonas Karthein, Dean Lee, Yuan-Zhuo Ma, Ulf-G. Meißner, Kei Minamisono, Alexander V. Oleynichenko, Skyy V. Pineda, Sergey D. Prosnyak, Marten L. Reitsma, Leonid V. Skripnikov, Adam Vernon, and Andréi Zaitsevskii

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In this Letter, we did not clarify the method used in the density functional theory calculations. The calculated charge radii using two functionals, NL3* [1] and SVmin [2], shown in Figs. 2 and 5, were taken from the database available at [3]. This website was not cited in the published version of this Letter. We would like to thank Witold Nazarewicz for pointing this out and highlighting that the calculations published at [3] do not represent the latest results that can be obtained with density functional theory.

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