

Erratum: “Perspective: Machine learning potentials for atomistic simulations” [J. Chem. Phys. 145, 170901 (2016)]

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Citation: *The Journal of Chemical Physics* **145**, 219901 (2016); doi: 10.1063/1.4971792

View online: <http://dx.doi.org/10.1063/1.4971792>

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(Received 17 November 2016; accepted 23 November 2016; published online 2 December 2016)

[<http://dx.doi.org/10.1063/1.4971792>]

There is a misprint in the references of the paper.¹ Reference 72 should be as follows: J. S. Elias, N. Artrith, M. Bugnet, L. Giordano, G. A. Botton, A. M. Kolpak, and Y. Shao-Horn, *ACS Catal.* **6**, 1675 (2016).²

¹J. Behler, *J. Chem. Phys.* **145**, 170901 (2016).

²J. S. Elias, N. Artrith, M. Bugnet, L. Giordano, G. A. Botton, A. M. Kolpak, and Y. Shao-Horn, *ACS Catal.* **6**, 1675 (2016).