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Erratum

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Erratum: Development of a "First-Principles" Water Potential with Flexible Monomers: Dimer Potential Energy Surface, VRT Spectrum, and Second Virial Coefficient

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Supporting Information

This erratum involves the replacement of the Supporting Information for the following paper: "Erratum: Development of a First-Principles Water Potential with Flexible Monomers: Dimer Potential Energy Surface, VRT Spectrum, and Second Virial Coefficient" (*J. Chem. Theory Comput.*, **2014**, *10* (5), 2212−2212, DOI: 10.1021/ct500284p). This is the second erratum regarding the Supporting Information for this paper. The original paper is *J. Chem. Theory Comput.*, **2013**, *9* (12), 5395−5403, DOI: 10.1021/ct400863t.

We have realized that, although all data are correct, some plots of the Supporting Information figures used different scaling factors, which could give a wrong perception of the results. These have been corrected in the Supporting Information accompanying this erratum.

■ ASSOCIATED CONTENT

Supporting Information

Corrected Figures S1−S4, along with the other files included in the original paper's Supporting Information. This material is available free of charge via the Internet at http://pubs.acs.org.

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