

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

S ome of the CC-pol dimer energies shown in Figures S1–S4 of the Supporting Information of ref 1 were calculated incorrectly. The correct Figures (S1–S4) are reported in the PDF file in the Supporting Information (zip) of this erratum, along with the correct root-mean square (RMS) errors obtained with the CCpol-8sf, HBB2, HBB2, and MB-pol potentials (Table S1).

ASSOCIATED CONTENT

Supporting Information

Corrected Figures S1—S4 and Table S1, along with all 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.

REFERENCES

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