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Erratum: Fast Calculations of Electrostatic Solvation Free Energy from Reconstructed Solvent Density using proximal Radial Distribution Functions

Bin Lin¹, Ka-Yiu Wong², Char Hu¹, Hironori Kokubo², and B. Montgomery Pettitt^{1,2}

¹Program in Structural and Computational Biology and Molecular Biophysics, Baylor College of Medicine, Houston, TX 77030 USA

²Department of Chemistry, University of Houston, Houston, TX 77204-5003 USA

In *J. Phys. Chem. Lett.* 2011, **2**, 1626–1632, two numbers in table 1 for the electrostatic solvation free energies (ESFE) of deca-alanine using the pRDF3 reconstruction method were incorrect and should be changed. The ESFEs for pRDF3 for I3 and H were –79.2 and –55.1 and should be –83.5 and –58.3 respectively. This changes none of the conclusions and improves the numerical consistency within the hierarchy of approximations.

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