

Correction to “Fast Calculations of Electrostatic Solvation Free Energy from Reconstructed Solvent Density Using Proximal Radial Distribution Functions”

Bin Lin, Ka-Yiu Wong, Char Hu, Hironori Kokubo, and B. Montgomery Pettitt*

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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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