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Addition/Correction

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Correction to "Molecular Dynamics Simulation for the Dynamics and Kinetics of Folding Peptides in the Gas Phase"

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In Table 1, page 12940, the value of σ for $V^{(2)}_{G}$ is 50 instead of 500, and the depth of the ion−He interaction potential, ε = 3.94, is in units of meV. The value in kcal mol^{-1} is 0.0909.

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