Exploring ion induced folding of a single-stranded DNA oligomer from molecular simulation studies

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Electronic Supplementary Information

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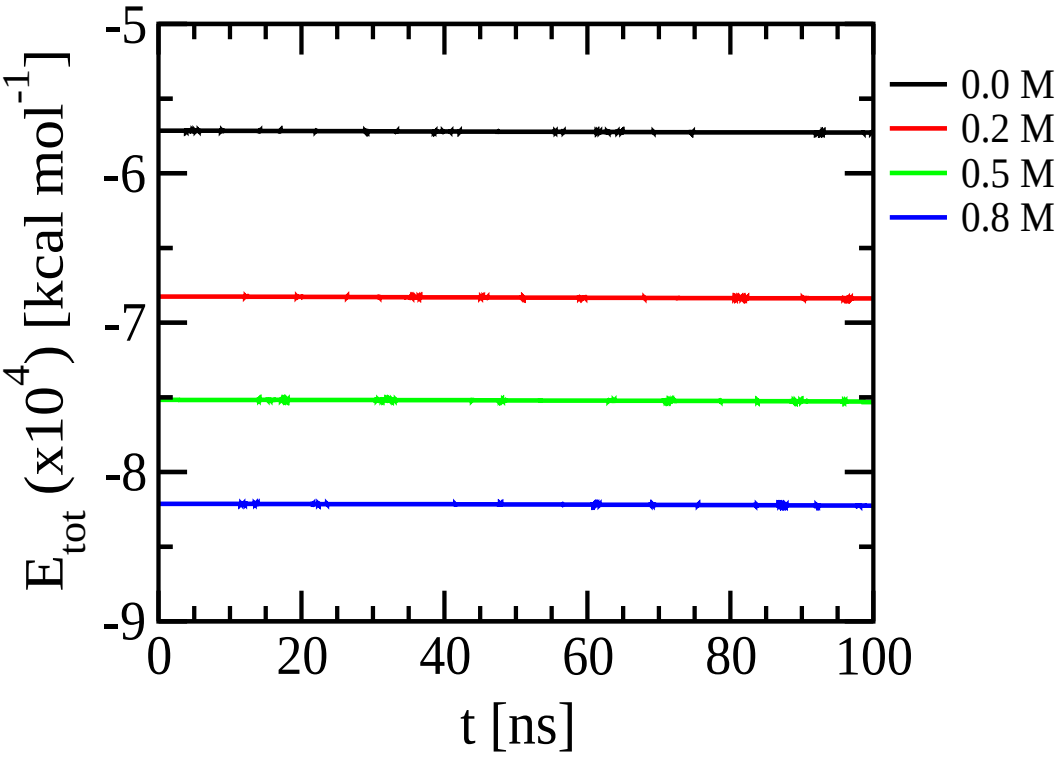

Figure SI-1. Time evolutions of the total energy of different simulation systems as obtained from the last 100 ns of the NVE trajectories.