Supporting information for Exploring the non-monotonic DNA capture behavior in a charged graphene nanopore

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Fig. S1. The RMSD (a, b) and the number of hydrogen bonds (c, d) between two stands of DNA as functions of time in the simulations.



Fig. S2. (a) The number of K^+ ions binding to each base pair. (b) Typical snapshots illustrating the binding of K^+ ions to O1P and O2P atoms, where the red vdw sphere is the oxygen atom in the phosphate group and the pink one is K^+ ion.