SUPPORTING INFORMATION

Diffusion Dynamics of the Li^+ ion on C_{60}: A direct molecular orbital-molecular dynamics Study

Hiroto TACHIKAWA*

Department of Materials Chemistry, Graduate School of Engineering, Hokkaido University, Sapporo 060-8628, Japan

---

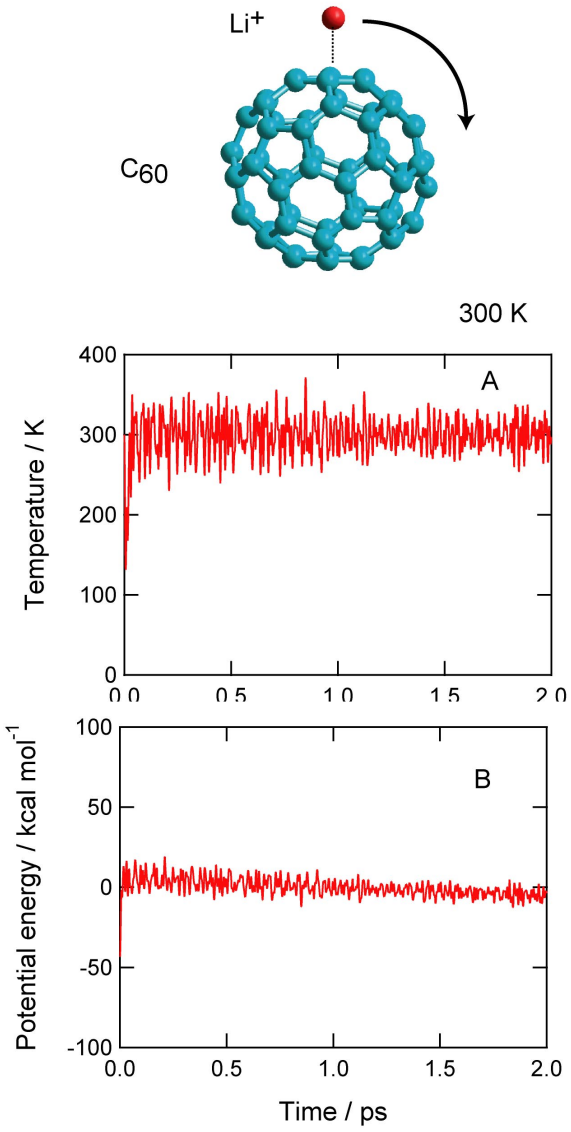

Figure S1. Temperature and potential energy of the Li^+C_{60} system plotted as a function of simulation time. Mean temperature is 300 K.