Supplementary Information

A Theoretical and Experimental Analysis of the Interfacial Mechanism of Dendrimer-Doxorubicin Complexes Formation

Barbara Jachimska[1], Magdalena Goncerz[1], Paweł Wolski[1], Callum Meldrum[2], Łukasz Lustyk[1], Tomasz Panczyk[1],*

[1] Jerzy Haber Institute of Catalysis and Surface Chemistry Polish Academy of Sciences, Krakow, Poland

[2] Department of Chemical and Process Engineering, University of Strathclyde; 75 Montrose Street, Glasgow G1 1XJ, UK

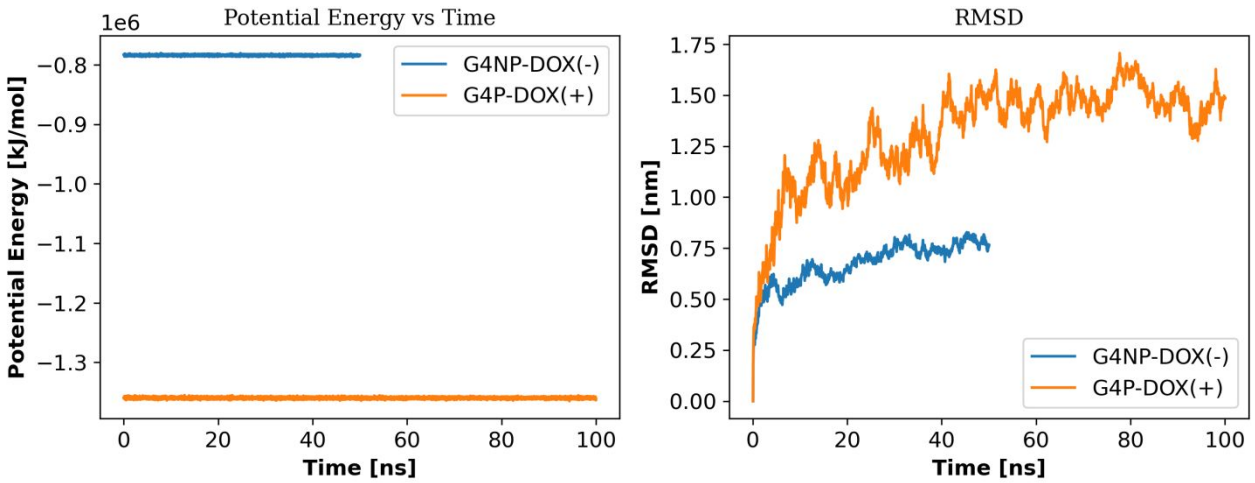

Figure S1. Change in the potential energy (left) and RMSD (right) values as a function of time. The RMSD value was calculated only for the dendrimer using the initial structure as a reference.