Supporting Information

A multi-scale molecular dynamics simulation of PMAL assisted siRNA delivery

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Figure S1. Mapping from all-atom model to coarse-grained model

(a) bond-tb-bb

(b) bond-bb-ac11

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Figure S2. Comparison of important parameters obtained from all-atom model and coarse-grained model of PMAL

Figure S3. Distance between COMs of PMAL-SiRNA complex and lipid bilayer with applied force time