

# An Extension and Further Validation of an All-Atomistic Membranes

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Citation Report

#	ARTICLE	IF	CITATIONS
11	Molecular Dynamics Simulations of a Characteristic DPC Micelle in Water. <i>Journal of Chemical Theory and Computation</i> , 2012, 8, 4610-4623.	2.3	62
12	Computer simulation of lipid membranes: Methodology and achievements. <i>Polymer Science - Series C</i> , 2013, 55, 162-180.	0.8	26
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14	Biomolecular Simulations with the Transferable Potentials for Phase Equilibria: Extension to Phospholipids. <i>Journal of Physical Chemistry B</i> , 2013, 117, 9910-9921.	1.2	14
15	Implicit inclusion of atomic polarization in modeling of partitioning between water and lipid bilayers. <i>Physical Chemistry Chemical Physics</i> , 2013, 15, 4677.	1.3	43
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