

List of Publications by Year in descending order

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216
papers

15,930
citations

10389

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113
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228
all docs

228
docs citations

228
times ranked

15077
citing authors

#	ARTICLE	IF	CITATIONS
1	Machine learning in the analysis of biomolecular simulations. <i>Advances in Physics: X</i> , 2022, 7, .	4.1	7
2	A cholesterol analog stabilizes the human β_2 -adrenergic receptor nonlinearly with temperature. <i>Science Signaling</i> , 2022, 15, .	3.6	8
3	N-Glycosylation can selectively block or foster different receptor–ligand binding modes. <i>Scientific Reports</i> , 2021, 11, 5239.	3.3	18
4	Martini 3: a general purpose force field for coarse-grained molecular dynamics. <i>Nature Methods</i> , 2021, 18, 382-388.	19.0	557
5	Seipin traps triacylglycerols to facilitate their nanoscale clustering in the endoplasmic reticulum membrane. <i>PLoS Biology</i> , 2021, 19, e3000998.	5.6	54
6	Complexity of seemingly simple lipid nanodiscs. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2020, 1862, 183420.	2.6	22
7	Cryo-EM structure of the complete and ligand-saturated insulin receptor ectodomain. <i>Journal of Cell Biology</i> , 2020, 219, .	5.2	84
8	Tail-Oxidized Cholesterol Enhances Membrane Permeability for Small Solutes. <i>Langmuir</i> , 2020, 36, 10438-10447.	3.5	24
9	Understanding the Functional Properties of Lipid Heterogeneity in Pulmonary Surfactant Monolayers at the Atomistic Level. <i>Frontiers in Cell and Developmental Biology</i> , 2020, 8, 581016.	3.7	18
10	Membrane-Dependent Binding and Entry Mechanism of Dopamine into Its Receptor. <i>ACS Chemical Neuroscience</i> , 2020, 11, 1914-1924.	3.5	21
11	High-resolution imaging and structure-based predictions reveal functional differences between Niemann-Pick C1 variants. <i>Traffic</i> , 2020, 21, 386-397.	2.7	14
12	Pulmonary Surfactant Lipid Reorganization Induced by the Adsorption of the Oligomeric Surfactant Protein B Complex. <i>Journal of Molecular Biology</i> , 2020, 432, 3251-3268.	4.2	29
13	Crystalline Wax Esters Regulate the Evaporation Resistance of Tear Film Lipid Layers Associated with Dry Eye Syndrome. <i>Journal of Physical Chemistry Letters</i> , 2019, 10, 3893-3898.	4.6	17
14	An efficient auxin-inducible degron system with low basal degradation in human cells. <i>Nature Methods</i> , 2019, 16, 866-869.	19.0	117
15	The Role of Temperature and Lipid Charge on Intake/Uptake of Cationic Gold Nanoparticles into Lipid Bilayers. <i>Small</i> , 2019, 15, e1805046.	10.0	35
16	Multiscale Simulations of Biological Membranes: The Challenge To Understand Biological Phenomena in a Living Substance. <i>Chemical Reviews</i> , 2019, 119, 5607-5774.	47.7	209
17	Understanding the Role of Lipids in Signaling Through Atomistic and Multiscale Simulations of Cell Membranes. <i>Annual Review of Biophysics</i> , 2019, 48, 421-439.	10.0	31
18	The Devil Is in the Details: What Do We Really Track in Single-Particle Tracking Experiments of Diffusion in Biological Membranes?. <i>Journal of Physical Chemistry Letters</i> , 2019, 10, 1005-1011.	4.6	13

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19	Bobbing of Oxysterols: Molecular Mechanism for Translocation of Tail-Oxidized Sterols through Biological Membranes. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 1118-1123.	4.6	24
20	Molecular mechanism for inhibition of twinfilin by phosphoinositides. <i>Journal of Biological Chemistry</i> , 2018, 293, 4818-4829.	3.4	15
21	How cardiolipin peroxidation alters the properties of the inner mitochondrial membrane?. <i>Chemistry and Physics of Lipids</i> , 2018, 214, 15-23.	3.2	35
22	Cholesterol Protects the Oxidized Lipid Bilayer from Water Injury: An All-Atom Molecular Dynamics Study. <i>Journal of Membrane Biology</i> , 2018, 251, 521-534.	2.1	12
23	Atomistic Model for Nearly Quantitative Simulations of Langmuir Monolayers. <i>Langmuir</i> , 2018, 34, 2565-2572.	3.5	53
24	Glucosylceramide modifies the LPS-induced inflammatory response in macrophages and the orientation of the LPS/TLR4 complex in silico. <i>Scientific Reports</i> , 2018, 8, 13600.	3.3	33
25	How to minimize dye-induced perturbations while studying biomembrane structure and dynamics: PEG linkers as a rational alternative. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2018, 1860, 2436-2445.	2.6	31
26	Redox-coupled quinone dynamics in the respiratory complex I. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2018, 115, E8413-E8420.	7.1	84
27	Negatively Charged Gangliosides Promote Membrane Association of Amphipathic Neurotransmitters. <i>Neuroscience</i> , 2018, 384, 214-223.	2.3	17
28	Quantitative Assessment of Methods Used To Obtain Rate Constants from Molecular Dynamics Simulations—Translocation of Cholesterol across Lipid Bilayers. <i>Journal of Chemical Theory and Computation</i> , 2018, 14, 3840-3848.	5.3	18
29	Long-chain GM1 gangliosides alter transmembrane domain registration through interdigitation. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2017, 1859, 870-878.	2.6	20
30	Glycosylation and Lipids Working in Concert Direct CD2 Ectodomain Orientation and Presentation. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 1060-1066.	4.6	22
31	Effects of Membrane PEGylation on Entry and Location of Antifungal Drug Itraconazole and Their Pharmacological Implications. <i>Molecular Pharmaceutics</i> , 2017, 14, 1057-1070.	4.6	19
32	Calcium Assists Dopamine Release by Preventing Aggregation on the Inner Leaflet of Presynaptic Vesicles. <i>ACS Chemical Neuroscience</i> , 2017, 8, 1242-1250.	3.5	21
33	Calcium Directly Regulates Phosphatidylinositol 4,5-Bisphosphate Headgroup Conformation and Recognition. <i>Journal of the American Chemical Society</i> , 2017, 139, 4019-4024.	13.7	87
34	On Atomistic Models for Molecular Oxygen. <i>Journal of Physical Chemistry B</i> , 2017, 121, 518-528.	2.6	19
35	Mechanistic principles underlying regulation of the actin cytoskeleton by phosphoinositides. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2017, 114, E8977-E8986.	7.1	106
36	Diffusion of Integral Membrane Proteins in Protein-Rich Membranes. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 4308-4313.	4.6	65

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37	<i>Glycans</i> Tools for Preparing Carbohydrate Structures for Atomistic Simulations of Glycoproteins, Glycolipids, and Carbohydrate Polymers for GROMACS. <i>Journal of Chemical Information and Modeling</i> , 2017, 57, 2401-2406.	5.4	71
38	Dynamics and energetics of the mammalian phosphatidylinositol transfer protein phospholipid exchange cycle. <i>Journal of Biological Chemistry</i> , 2017, 292, 14438-14455.	3.4	25
39	Nanoscale Membrane Domain Formation Driven by Cholesterol. <i>Scientific Reports</i> , 2017, 7, 1143.	3.3	83
40	The Integrin Receptor in Biologically Relevant Bilayers: Insights from Molecular Dynamics Simulations. <i>Journal of Membrane Biology</i> , 2017, 250, 337-351.	2.1	29
41	Phase Partitioning of GM1 and Its Bodipy-Labeled Analog Determine Their Different Binding to Cholera Toxin. <i>Frontiers in Physiology</i> , 2017, 8, 252.	2.8	34
42	Concerted regulation of npc2 binding to endosomal/lysosomal membranes by bis(monoacylglycero)phosphate and sphingomyelin. <i>PLoS Computational Biology</i> , 2017, 13, e1005831.	3.2	27
43	Excessive aggregation of membrane proteins in the Martini model. <i>PLoS ONE</i> , 2017, 12, e0187936.	2.5	147
44	The role of hydrophobic matching on transmembrane helix packing in cells. <i>Cell Stress</i> , 2017, 1, 90-106.	3.2	37
45	Selective effect of cell membrane on synaptic neurotransmission. <i>Scientific Reports</i> , 2016, 6, 19345.	3.3	48
46	Data including GROMACS input files for atomistic molecular dynamics simulations of mixed, asymmetric bilayers including molecular topologies, equilibrated structures, and force field for lipids compatible with OPLS-AA parameters. <i>Data in Brief</i> , 2016, 7, 1171-1174.	1.0	15
47	Atomistic determinants of co-enzyme Q reduction at the Qi-site of the cytochrome bc1 complex. <i>Scientific Reports</i> , 2016, 6, 33607.	3.3	23
48	Protein Crowding in Lipid Bilayers Gives Rise to Non-Gaussian Anomalous Lateral Diffusion of Phospholipids and Proteins. <i>Physical Review X</i> , 2016, 6, .	8.9	152
49	What Can We Learn about Cholesterol's Transmembrane Distribution Based on Cholesterol-Induced Changes in Membrane Dipole Potential?. <i>Journal of Physical Chemistry Letters</i> , 2016, 7, 4585-4590.	4.6	19
50	Lipid membranes: Theory and simulations bridged to experiments. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2016, 1858, 2251-2253.	2.6	12
51	Distribution and dynamics of quinones in the lipid bilayer mimicking the inner membrane of mitochondria. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2016, 1858, 2116-2122.	2.6	47
52	Interdigitation of long-chain sphingomyelin induces coupling of membrane leaflets in a cholesterol dependent manner. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2016, 1858, 281-288.	2.6	76
53	Role of charged lipids in membrane structures Insight given by simulations. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2016, 1858, 2322-2333.	2.6	63
54	Cholesterol oxidation products and their biological importance. <i>Chemistry and Physics of Lipids</i> , 2016, 199, 144-160.	3.2	130

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55	Mechanism of allosteric regulation of β^2 -adrenergic receptor by cholesterol. <i>ELife</i> , 2016, 5, .	6.0	115
56	Building Synthetic Sterols Computationally â€“ Unlocking the Secrets of Evolution?. <i>Frontiers in Bioengineering and Biotechnology</i> , 2015, 3, 121.	4.1	5
57	PEGylated Liposomes as Carriers of Hydrophobic Porphyrins. <i>Journal of Physical Chemistry B</i> , 2015, 119, 6646-6657.	2.6	47
58	Cholesterol under oxidative stressâ€”How lipid membranes sense oxidation as cholesterol is being replaced by oxysterols. <i>Free Radical Biology and Medicine</i> , 2015, 84, 30-41.	2.9	57
59	How Well Does BODIPY-Cholesteryl Ester Mimic Unlabeled Cholesteryl Esters in High Density Lipoprotein Particles?. <i>Journal of Physical Chemistry B</i> , 2015, 119, 15848-15856.	2.6	4
60	Proton-coupled electron transfer and the role of water molecules in proton pumping by cytochrome <i>c</i> oxidase. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2015, 112, 2040-2045.	7.1	59
61	Biogenesis of Nascent High Density Lipoprotein Particles. <i>Structure</i> , 2015, 23, 1153-1154.	3.3	3
62	How To Minimize Artifacts in Atomistic Simulations of Membrane Proteins, Whose Crystal Structure Is Heavily Engineered: β^2 -Adrenergic Receptor in the Spotlight. <i>Journal of Chemical Theory and Computation</i> , 2015, 11, 3432-3445.	5.3	16
63	Oxidation of Cholesterol Does Not Alter Significantly Its Uptake into High-Density Lipoprotein Particles. <i>Journal of Physical Chemistry B</i> , 2015, 119, 4594-4600.	2.6	6
64	Role of subunit III and its lipids in the molecular mechanism of cytochrome c oxidase. <i>Biochimica Et Biophysica Acta - Bioenergetics</i> , 2015, 1847, 690-697.	1.0	24
65	Sec14-nodulin proteins and the patterning of phosphoinositide landmarks for developmental control of membrane morphogenesis. <i>Molecular Biology of the Cell</i> , 2015, 26, 1764-1781.	2.1	44
66	<i>N</i> -Glycosylation as determinant of epidermal growth factor receptor conformation in membranes. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2015, 112, 4334-4339.	7.1	135
67	PIP2 and Talin Join Forces to Activate Integrin. <i>Journal of Physical Chemistry B</i> , 2015, 119, 12381-12389.	2.6	27
68	Redox-induced activation of the proton pump in the respiratory complex I. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2015, 112, 11571-11576.	7.1	122
69	How endoglucanase enzymes act on cellulose nanofibrils: role of amorphous regions revealed by atomistic simulations. <i>Cellulose</i> , 2015, 22, 2911-2925.	4.9	20
70	Experimental determination and computational interpretation of biophysical properties of lipid bilayers enriched by cholesteryl hemisuccinate. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2015, 1848, 422-432.	2.6	45
71	Enzymatic Oxidation of Cholesterol: Properties and Functional Effects of Cholestenone in Cell Membranes. <i>PLoS ONE</i> , 2014, 9, e103743.	2.5	50
72	Co-Exposure with Fullerene May Strengthen Health Effects of Organic Industrial Chemicals. <i>PLoS ONE</i> , 2014, 9, e114490.	2.5	9

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73	How Anacetrapib Inhibits the Activity of the Cholesteryl Ester Transfer Protein? Perspective through Atomistic Simulations. <i>PLoS Computational Biology</i> , 2014, 10, e1003987.	3.2	17
74	The challenges of understanding glycolipid functions: An open outlook based on molecular simulations. <i>Biochimica Et Biophysica Acta - Molecular and Cell Biology of Lipids</i> , 2014, 1841, 1130-1145.	2.4	35
75	Effect of PEGylation on Drug Entry into Lipid Bilayer. <i>Journal of Physical Chemistry B</i> , 2014, 118, 144-151.	2.6	26
76	How well does cholesteryl hemisuccinate mimic cholesterol in saturated phospholipid bilayers?. <i>Journal of Molecular Modeling</i> , 2014, 20, 2121.	1.8	44
77	How to link pyrene to its host lipid to minimize the extent of membrane perturbations and to optimize pyrene dimer formation. <i>Chemistry and Physics of Lipids</i> , 2014, 177, 19-25.	3.2	7
78	Cholesterol, sphingolipids, and glycolipids: What do we know about their role in raft-like membranes?. <i>Chemistry and Physics of Lipids</i> , 2014, 184, 82-104.	3.2	159
79	Cationic Au Nanoparticle Binding with Plasma Membrane-like Lipid Bilayers: Potential Mechanism for Spontaneous Permeation to Cells Revealed by Atomistic Simulations. <i>Journal of Physical Chemistry C</i> , 2014, 118, 11131-11141.	3.1	69
80	How To Tackle the Issues in Free Energy Simulations of Long Amphiphiles Interacting with Lipid Membranes: Convergence and Local Membrane Deformations. <i>Journal of Physical Chemistry B</i> , 2014, 118, 3572-3581.	2.6	61
81	Atomistic simulations of anionic Au ₁₄₄ (SR) ₆₀ nanoparticles interacting with asymmetric model lipid membranes. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2014, 1838, 2852-2860.	2.6	46
82	Refined OPLS All-Atom Force Field for Saturated Phosphatidylcholine Bilayers at Full Hydration. <i>Journal of Physical Chemistry B</i> , 2014, 118, 4571-4581.	2.6	139
83	Can pyrene probes be used to measure lateral pressure profiles of lipid membranes? Perspective through atomistic simulations. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2014, 1838, 1406-1411.	2.6	16
84	Modeling of Lipid Membranes and Lipoproteins. , 2014, , 299-318.		2
85	Parameterization of the prosthetic redox centers of the bacterial cytochrome bc ₁ complex for atomistic molecular dynamics simulations. <i>Theoretical Chemistry Accounts</i> , 2013, 132, 1.	1.4	16
86	Key role of water in proton transfer at the Qo-site of the cytochrome bc ₁ complex predicted by atomistic molecular dynamics simulations. <i>Biochimica Et Biophysica Acta - Bioenergetics</i> , 2013, 1827, 761-768.	1.0	39
87	Anomalous and normal diffusion of proteins and lipids in crowded lipid membranes. <i>Faraday Discussions</i> , 2013, 161, 397-417.	3.2	170
88	Atomistic simulations indicate cardiolipin to have an integral role in the structure of the cytochrome bc ₁ complex. <i>Biochimica Et Biophysica Acta - Bioenergetics</i> , 2013, 1827, 769-778.	1.0	54
89	Molecular lipidomics of exosomes released by PC-3 prostate cancer cells. <i>Biochimica Et Biophysica Acta - Molecular and Cell Biology of Lipids</i> , 2013, 1831, 1302-1309.	2.4	546
90	Do Lipids Retard the Evaporation of the Tear Fluid?. , 2012, 53, 6442.		49

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91	Interfacial Tension and Surface Pressure of High Density Lipoprotein, Low-Density Lipoprotein, and Related Lipid Droplets. <i>Biophysical Journal</i> , 2012, 103, 1236-1244.	0.5	42
92	Mechanism for translocation of fluoroquinolones across lipid membranes. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2012, 1818, 2563-2571.	2.6	76
93	Role of Neutral Lipids in Tear Fluid Lipid Layer: Coarse-Grained Simulation Study. <i>Langmuir</i> , 2012, 28, 17092-17100.	3.5	27
94	The impact of lipid composition on the stability of the tear fluid lipid layer. <i>Soft Matter</i> , 2012, 8, 5826.	2.7	40
95	Interaction of Hematoporphyrin with Lipid Membranes. <i>Journal of Physical Chemistry B</i> , 2012, 116, 4889-4897.	2.6	36
96	Interaction of C70 fullerene with the Kv1.2 potassium channel. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 12526.	2.8	17
97	Strong preferences of dopamine and L-dopa towards lipid head group: importance of lipid composition and implication for neurotransmitter metabolism. <i>Journal of Neurochemistry</i> , 2012, 122, 681-690.	3.9	51
98	Cationic Dimyristoylphosphatidylcholine and Dioleoyloxytrimethylammonium Propane Lipid Bilayers: Atomistic Insight for Structure and Dynamics. <i>Journal of Physical Chemistry B</i> , 2012, 116, 269-276.	2.6	25
99	Atomistic Simulations of Functional Au ₁₄₄ (SR) ₆₀ Gold Nanoparticles in Aqueous Environment. <i>Journal of Physical Chemistry C</i> , 2012, 116, 9805-9815.	3.1	94
100	Revealing structural and dynamical properties of high density lipoproteins through molecular simulations. <i>Soft Matter</i> , 2012, 8, 1262-1267.	2.7	11
101	Lipid Exchange Mechanism of the Cholesteryl Ester Transfer Protein Clarified by Atomistic and Coarse-grained Simulations. <i>PLoS Computational Biology</i> , 2012, 8, e1002299.	3.2	49
102	Drug-Lipid Membrane Interaction Mechanisms Revealed Through Molecular Simulations. <i>Current Physical Chemistry</i> , 2012, 2, 379-400.	0.2	13
103	Effect of Galactosylceramide on the Dynamics of Cholesterol-Rich Lipid Membranes. <i>Journal of Physical Chemistry B</i> , 2011, 115, 14424-14434.	2.6	17
104	Properties of the Membrane Binding Component of Catechol-O-methyltransferase Revealed by Atomistic Molecular Dynamics Simulations. <i>Journal of Physical Chemistry B</i> , 2011, 115, 13541-13550.	2.6	15
105	Protein Shape Change Has a Major Effect on the Gating Energy of a Mechanosensitive Channel. <i>Biophysical Journal</i> , 2011, 100, 1651-1659.	0.5	48
106	Analysis of Twisting of Cellulose Nanofibrils in Atomistic Molecular Dynamics Simulations. <i>Journal of Physical Chemistry B</i> , 2011, 115, 3747-3755.	2.6	129
107	Low density lipoprotein: structure, dynamics, and interactions of apoB-100 with lipids. <i>Soft Matter</i> , 2011, 7, 8135.	2.7	47
108	A MARTINI Coarse-Grained Model of a Thermoset Polyester Coating. <i>Macromolecules</i> , 2011, 44, 6198-6208.	4.8	66

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109	N- and O-methylation of sphingomyelin markedly affects its membrane properties and interactions with cholesterol. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2011, 1808, 1179-1186.	2.6	35
110	Lipid Simulations: A Perspective on Lipids in Action. <i>Cold Spring Harbor Perspectives in Biology</i> , 2011, 3, a004655-a004655.	5.5	35
111	Coarse-graining polymers with the MARTINI force-field: polystyrene as a benchmark case. <i>Soft Matter</i> , 2011, 7, 698-708.	2.7	216
112	Cholesterol modulates glycolipid conformation and receptor activity. <i>Nature Chemical Biology</i> , 2011, 7, 260-262.	8.0	194
113	Lessons from the biophysics of interfaces: Lung surfactant and tear fluid. <i>Progress in Retinal and Eye Research</i> , 2011, 30, 204-215.	15.5	46
114	Lateral sorting in model membranes by cholesterol-mediated hydrophobic matching. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2011, 108, 16628-16633.	7.1	131
115	Association of Lipidome Remodeling in the Adipocyte Membrane with Acquired Obesity in Humans. <i>PLoS Biology</i> , 2011, 9, e1000623.	5.6	213
116	High Density Lipoprotein Structural Changes and Drug Response in Lipidomic Profiles following the Long-Term Fenofibrate Therapy in the FIELD Substudy. <i>PLoS ONE</i> , 2011, 6, e23589.	2.5	33
117	Defect-Mediated Trafficking across Cell Membranes: Insights from <i>in Silico</i> Modeling. <i>Chemical Reviews</i> , 2010, 110, 6077-6103.	47.7	171
118	Molecular Dynamics Simulations Reveal Fundamental Role of Water As Factor Determining Affinity of Binding of β -Blocker Nebivolol to β -Adrenergic Receptor. <i>Journal of Physical Chemistry B</i> , 2010, 114, 8374-8386.	2.6	35
119	Role of Lipids in Spheroidal High Density Lipoproteins. <i>PLoS Computational Biology</i> , 2010, 6, e1000964.	3.2	81
120	Composition and lipid spatial distribution of HDL particles in subjects with low and high HDL-cholesterol. <i>Journal of Lipid Research</i> , 2010, 51, 2341-2351.	4.2	111
121	Free Volume Theory Applied to Lateral Diffusion in Langmuir Monolayers: Atomistic Simulations for a Protein-Free Model of Lung Surfactant. <i>Langmuir</i> , 2010, 26, 15436-15444.	3.5	42
122	Molecular Organization of the Tear Fluid Lipid Layer. <i>Biophysical Journal</i> , 2010, 99, 2559-2567.	0.5	67
123	Role of Glycolipids in Lipid Rafts: A View through Atomistic Molecular Dynamics Simulations with Galactosylceramide. <i>Journal of Physical Chemistry B</i> , 2010, 114, 7797-7807.	2.6	60
124	Effects of DPH on DPPC~Cholesterol Membranes with Varying Concentrations of Cholesterol: From Local Perturbations to Limitations in Fluorescence Anisotropy Experiments. <i>Journal of Physical Chemistry B</i> , 2010, 114, 2704-2711.	2.6	39
125	The hydrophobic effect and its role in cold denaturation. <i>Cryobiology</i> , 2010, 60, 91-99.	0.7	164
126	Reply to the comment by Graziano on "The hydrophobic effect and its role in cold denaturation". <i>Cryobiology</i> , 2010, 60, 356-357.	0.7	1

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127	Temperature-induced structural transition in-situ in porcine lens " Changes observed in void size distribution. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2010, 1798, 958-965.	2.6	12
128	Membrane Proteins Diffuse as Dynamic Complexes with Lipids. <i>Journal of the American Chemical Society</i> , 2010, 132, 7574-7575.	13.7	157
129	Effect of Sphingomyelin Headgroup Size on Molecular Properties and Interactions with Cholesterol. <i>Biophysical Journal</i> , 2010, 99, 3300-3308.	0.5	75
130	Concerted diffusion of lipids in raft-like membranes. <i>Faraday Discussions</i> , 2010, 144, 411-430.	3.2	92
131	Cholesterol Induces Specific Spatial and Orientational Order in Cholesterol/Phospholipid Membranes. <i>PLoS ONE</i> , 2010, 5, e11162.	2.5	101
132	Calculation of the electrostatic potential of lipid bilayers from molecular dynamics simulations: Methodological issues. <i>Journal of Chemical Physics</i> , 2009, 130, 215107.	3.0	63
133	Desipramine induces disorder in cholesterol-rich membranes: implications for viral trafficking. <i>Physical Biology</i> , 2009, 6, 046004.	1.8	10
134	Conformational Changes and Slow Dynamics through Microsecond Polarized Atomistic Molecular Simulation of an Integral Kv1.2 Ion Channel. <i>PLoS Computational Biology</i> , 2009, 5, e1000289.	3.2	108
135	Systematic coarse graining from structure using internal states: Application to phospholipid/cholesterol bilayer. <i>Journal of Chemical Physics</i> , 2009, 131, 055101.	3.0	51
136	Effects of carbon nanoparticles on lipid membranes: a molecular simulation perspective. <i>Soft Matter</i> , 2009, 5, 4433.	2.7	116
137	Mitochondrial Membranes with Mono- and Divalent Salt: Changes Induced by Salt Ions on Structure and Dynamics. <i>Journal of Physical Chemistry B</i> , 2009, 113, 15513-15521.	2.6	41
138	Why is the $n=2$ Chain of Monounsaturated Glycerophospholipids Usually Unsaturated whereas the $n=1$ Chain Is Saturated? Studies of 1-Stearoyl-2-oleoyl-glycero-3-phosphatidylcholine (SOPC) and 1-Oleoyl-2-stearoyl-glycero-3-phosphatidylcholine (OSPC) Membranes with and without Cholesterol. <i>Journal of Physical Chemistry B</i> , 2009, 113, 8347-8356.	2.6	22
139	Probing Biomembranes with Positrons. <i>Journal of Physical Chemistry B</i> , 2009, 113, 1810-1812.	2.6	31
140	Ion Dynamics in Cationic Lipid Bilayer Systems in Saline Solutions. <i>Journal of Physical Chemistry B</i> , 2009, 113, 9226-9234.	2.6	38
141	Atom-scale molecular interactions in lipid raft mixtures. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2009, 1788, 122-135.	2.6	74
142	Ordering effects of cholesterol and its analogues. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2009, 1788, 97-121.	2.6	506
143	Ceramide-1-Phosphate, in Contrast to Ceramide, Is Not Segregated into Lateral Lipid Domains in Phosphatidylcholine Bilayers. <i>Biophysical Journal</i> , 2009, 96, 2216-2226.	0.5	21
144	Atomistic Simulations of Phosphatidylcholines and Cholesteryl Esters in High-Density Lipoprotein-Sized Lipid Droplet and Trilayer: Clues to Cholesteryl Ester Transport and Storage. <i>Biophysical Journal</i> , 2009, 96, 4099-4108.	0.5	33

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145	Intrinsic Potential of Cell Membranes: Opposite Effects of Lipid Transmembrane Asymmetry and Asymmetric Salt Ion Distribution. <i>Journal of Physical Chemistry B</i> , 2009, 113, 7194-7198.	2.6	23
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147	Asymmetric nature of lateral pressure profiles in supported lipid membranes and its implications for membrane protein functions. <i>Soft Matter</i> , 2009, 5, 3258.	2.7	32
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