## Patrick F J Fuchs

List of Publications by Year in descending order

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Version: 2024-02-01

448610 425179 1,668 35 19 34 citations g-index h-index papers 36 36 36 2596 docs citations times ranked citing authors all docs

#	Article	IF	CITATIONS
1	How to best estimate the viscosity of lipid bilayers. Biophysical Chemistry, 2022, 281, 106732.	1.5	8
2	Structural Bases for the Involvement of Phosphatidylinositol-4,5-bisphosphate in the Internalization of the Cell-Penetrating Peptide Penetratin. ACS Chemical Biology, 2022, 17, 1427-1439.	1.6	6
3	Molecular Study of Ultrasound-Triggered Release of Fluorescein from Liposomes. Langmuir, 2021, 37, 3868-3881.	1.6	7
4	Inverse Conformational Selection in Lipid–Protein Binding. Journal of the American Chemical Society, 2021, 143, 13701-13709.	6.6	16
5	buildH: Build hydrogen atoms from united-atom molecular dynamics of lipids and calculate the order parameters. Journal of Open Source Software, 2021, 6, 3521.	2.0	O
6	On the Effect of the Various Assumptions and Approximations used in Molecular Simulations on the Properties of Bioâ€Molecular Systems: Overview and Perspective on Issues. ChemPhysChem, 2021, 22, 264-282.	1.0	12
7	Cooperation of Conical and Polyunsaturated Lipids to Regulate Initiation and Processing of Membrane Fusion. Frontiers in Molecular Biosciences, 2021, 8, 763115.	1.6	11
8	CX3CL1 homo-oligomerization drives cell-to-cell adherence. Scientific Reports, 2020, 10, 9069.	1.6	13
9	Biophysical Insight on the Membrane Insertion of an Arginine-Rich Cell-Penetrating Peptide. International Journal of Molecular Sciences, 2019, 20, 4441.	1.8	14
10	Influence of the Treatment of Nonbonded Interactions on the Thermodynamic and Transport Properties of Pure Liquids Calculated Using the 2016H66 Force Field. Journal of Chemical Theory and Computation, 2019, 15, 1806-1826.	2.3	28
11	The heptad repeat domain $1$ of Mitofusin has membrane destabilization function in mitochondrial fusion. EMBO Reports, 2018, $19$ , .	2.0	39
12	Effect of Dimethyl Sulfoxide on the Binding of 1-Adamantane Carboxylic Acid to $\hat{l}^2$ - and $\hat{l}^3$ -Cyclodextrins. ACS Omega, 2018, 3, 1014-1021.	1.6	12
13	Meet-U: Educating through research immersion. PLoS Computational Biology, 2018, 14, e1005992.	1.5	4
14	Interdigitation between Triglycerides and Lipids Modulates Surface Properties of Lipid Droplets. Biophysical Journal, 2017, 112, 1417-1430.	0.2	102
15	Simulating Bilayers of Nonionic Surfactants with the GROMOS-Compatible 2016H66 Force Field. Langmuir, 2017, 33, 10225-10238.	1.6	12
16	Targeting surface voids to counter membrane disorders in lipointoxication-related diseases. Journal of Cell Science, 2016, 129, 2368-81.	1.2	5
17	A GROMOS-Compatible Force Field for Small Organic Molecules in the Condensed Phase: The 2016H66 Parameter Set. Journal of Chemical Theory and Computation, 2016, 12, 3825-3850.	2.3	118
18	Nuclear pore targeting of the yeast Pom33 nucleoporin depends on karyopherin- and lipid-binding. Journal of Cell Science, 2015, 128, 305-16.	1.2	29

#	Article	IF	CITATIONS
19	Toward Atomistic Resolution Structure of Phosphatidylcholine Headgroup and Glycerol Backbone at Different Ambient Conditions. Journal of Physical Chemistry B, 2015, 119, 15075-15088.	1.2	109
20	Insights on pH-dependent conformational changes of mosquito odorant binding proteins by molecular dynamics simulations. Journal of Biomolecular Structure and Dynamics, 2014, 32, 1742-1751.	2.0	10
21	Coarse-Grained Simulations of the HIV-1 Matrix Protein Anchoring: Revisiting Its Assembly on Membrane Domains. Biophysical Journal, 2014, 106, 577-585.	0.2	71
22	Modeling and molecular dynamics simulations of the V33 variant of the integrin subunit $\hat{l}^2$ 3: Structural comparison with the L33 (HPA-1a) and P33 (HPA-1b) variants. Biochimie, 2014, 105, 84-90.	1.3	20
23	Conical Lipids in Flat Bilayers Induce Packing Defects Similar to that Induced by Positive Curvature. Biophysical Journal, 2013, 104, 585-593.	0.2	149
24	Amphipathic Lipid Packing Sensor Motifs: Probing Bilayer Defects with Hydrophobic Residues. Biophysical Journal, 2013, 104, 575-584.	0.2	171
25	A GROMOS Parameter Set for Vicinal Diether Functions: Properties of Polyethyleneoxide and Polyethyleneglycol. Journal of Chemical Theory and Computation, 2012, 8, 3943-3963.	2.3	61
26	Modeling and Molecular Dynamics of HPA-1a and -1b Polymorphisms: Effects on the Structure of the $\hat{l}^2$ 3 Subunit of the $\hat{l}^2$ 3 Integrin. PLoS ONE, 2012, 7, e47304.	1.1	33
27	Amphipathic-Lipid-Packing-Sensor interactions with lipids assessed by atomistic molecular dynamics. Biochimica Et Biophysica Acta - Biomembranes, 2011, 1808, 2119-2127.	1.4	35
28	New Interaction Parameters for Oxygen Compounds in the GROMOS Force Field: Improved Pure-Liquid and Solvation Properties for Alcohols, Ethers, Aldehydes, Ketones, Carboxylic Acids, and Esters. Journal of Chemical Theory and Computation, 2011, 7, 1016-1031.	2.3	112
29	Molecular Dynamics of Membrane Peptides and Proteins: Principles and Comparison to Experimental Data. Methods in Molecular Biology, 2010, 654, 403-421.	0.4	4
30	Interpretation of 2H-NMR Experiments on the Orientation of the Transmembrane Helix WALP23 by Computer Simulations. Biophysical Journal, 2010, 99, 1455-1464.	0.2	43
31	On the Orientation of a Designed Transmembrane Peptide:  Toward the Right Tilt Angle?. Journal of the American Chemical Society, 2007, 129, 15174-15181.	6.6	96
32	Kinetics and Thermodynamics of Type VIII $\hat{I}^2$ -Turn Formation: A CD, NMR, and Microsecond Explicit Molecular Dynamics Study of the GDNP Tetrapeptide. Biophysical Journal, 2006, 90, 2745-2759.	0.2	44
33	High accuracy prediction of $\hat{l}^2$ -turns and their types using propensities and multiple alignments. Proteins: Structure, Function and Bioinformatics, 2005, 59, 828-839.	1.5	101
34	NMR and molecular dynamics studies of an autoimmune myelin basic protein peptide and its antagonist. FEBS Journal, 2004, 271, 3399-3413.	0.2	20
35	Conformational Dependence of Collagenase (Matrix Metalloproteinase-1) Up-regulation by Elastin Peptides in Cultured Fibroblasts. Journal of Biological Chemistry, 2001, 276, 5222-5227.	1.6	153