

Ulrich Zachariae

List of Publications by Year in descending order

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64
papers

2,962
citations

201674

27
h-index

182427

51
g-index

77
all docs

77
docs citations

77
times ranked

4203
citing authors

#	ARTICLE	IF	CITATIONS
1	Ion permeation in K ⁺ channels occurs by direct Coulomb knock-on. <i>Science</i> , 2014, 346, 352-355.	12.6	271
2	Computational Electrophysiology: The Molecular Dynamics of Ion Channel Permeation and Selectivity in Atomistic Detail. <i>Biophysical Journal</i> , 2011, 101, 809-817.	0.5	214
3	Direct knock-on of desolvated ions governs strict ion selectivity in K ⁺ channels. <i>Nature Chemistry</i> , 2018, 10, 813-820.	13.6	170
4	Structure of the MacAB-TolC ABC-type tripartite multidrug efflux pump. <i>Nature Microbiology</i> , 2017, 2, 17070.	13.3	140
5	Mechanisms of Anion Conduction by Coupled Glutamate Transporters. <i>Cell</i> , 2015, 160, 542-553.	28.9	114
6	Position of Transmembrane Helix 6 Determines Receptor G Protein Coupling Specificity. <i>Journal of the American Chemical Society</i> , 2014, 136, 11244-11247.	13.7	105
7	Î²-Barrel Mobility Underlies Closure of the Voltage-Dependent Anion Channel. <i>Structure</i> , 2012, 20, 1540-1549.	3.3	104
8	Crystal structure and functional mechanism of a human antimicrobial membrane channel. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2013, 110, 4586-4591.	7.1	104
9	Functional dynamics in the voltage-dependent anion channel. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2010, 107, 22546-22551.	7.1	97
10	GPCRmd uncovers the dynamics of the 3D-GPCRome. <i>Nature Methods</i> , 2020, 17, 777-787.	19.0	90
11	Quantitative Structural Analysis of Importin-Î² Flexibility: Paradigm for Solenoid Protein Structures. <i>Structure</i> , 2010, 18, 1171-1183.	3.3	89
12	Molecular Determinants for Nuclear Import of Influenza A PB2 by Importin Î± Isoforms 3 and 7. <i>Structure</i> , 2015, 23, 374-384.	3.3	87
13	Intracellular Transfer of Na ⁺ in an Active-State G-Protein-Coupled Receptor. <i>Structure</i> , 2018, 26, 171-180.e2.	3.3	77
14	Interfacial self-assembly of a bacterial hydrophobin. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2015, 112, 5419-5424.	7.1	68
15	Toward a Consensus Model of the hERG Potassium Channel. <i>ChemMedChem</i> , 2010, 5, 455-467.	3.2	66
16	An Unusual Hydrophobic Core Confers Extreme Flexibility to HEAT Repeat Proteins. <i>Biophysical Journal</i> , 2010, 99, 1596-1603.	0.5	66
17	Structural basis for cooperativity of CRM1 export complex formation. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2013, 110, 960-965.	7.1	64
18	Insights into the function of ion channels by computational electrophysiology simulations. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2016, 1858, 1741-1752.	2.6	60

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19	The Persistent Question of Potassium Channel Permeation Mechanisms. <i>Journal of Molecular Biology</i> , 2021, 433, 167002.	4.2	55
20	The Molecular Mechanism of Toxin-Induced Conformational Changes in a Potassium Channel: Relation to C-Type Inactivation. <i>Structure</i> , 2008, 16, 747-754.	3.3	52
21	Importin- β : Structural and Dynamic Determinants of a Molecular Spring. <i>Structure</i> , 2008, 16, 906-915.	3.3	49
22	Structural Mechanisms of Voltage Sensing in G α Protein-Coupled Receptors. <i>Structure</i> , 2016, 24, 997-1007.	3.3	48
23	Side Chain Flexibilities in the Human Ether-a-go-go Related Gene Potassium Channel (hERG) Together with Matched-Pair Binding Studies Suggest a New Binding Mode for Channel Blockers. <i>Journal of Medicinal Chemistry</i> , 2009, 52, 4266-4276.	6.4	44
24	Vacuolar Protein Sorting: Two Different Functional States of the AAA-ATPase Vps4p. <i>Journal of Molecular Biology</i> , 2008, 377, 352-363.	4.2	41
25	Crystallographic analysis of <i>Neisseria meningitidis</i> PorB extracellular loops potentially implicated in TLR2 recognition. <i>Journal of Structural Biology</i> , 2014, 185, 440-447.	2.8	32
26	Membrane potentials regulating GPCRs: insights from experiments and molecular dynamics simulations. <i>Current Opinion in Pharmacology</i> , 2016, 30, 44-50.	3.5	32
27	Ten Years of High Resolution Structural Research on the Voltage Dependent Anion Channel (VDAC) – Recent Developments and Future Directions. <i>Frontiers in Physiology</i> , 2018, 9, 108.	2.8	32
28	Quantifying Disorder through Conditional Entropy: An Application to Fluid Mixing. <i>PLoS ONE</i> , 2013, 8, e65617.	2.5	32
29	The O-GlcNAc Transferase Intellectual Disability Mutation L254F Distorts the TPR Helix. <i>Cell Chemical Biology</i> , 2018, 25, 513-518.e4.	5.2	30
30	Electrostatic properties of the anion selective porin Omp32 from <i>Delftia acidovorans</i> and of the arginine cluster of bacterial porins. <i>Protein Science</i> , 2002, 11, 1309-1319.	7.6	29
31	High Resolution Crystal Structures and Molecular Dynamics Studies Reveal Substrate Binding in the Porin Omp32. <i>Journal of Biological Chemistry</i> , 2006, 281, 7413-7420.	3.4	29
32	Picosecond Time-Resolved FRET in the Fluorescent Protein from <i>Discosoma Red</i> (wt-DsRed). <i>ChemPhysChem</i> , 2001, 2, 325-328.	2.1	28
33	The Bacterial Hydrophobin BslA is a Switchable Ellipsoidal Janus Nanocolloid. <i>Langmuir</i> , 2015, 31, 11558-11563.	3.5	28
34	The lipid environment determines the activity of the <i>Escherichia coli</i> ammonium transporter AmtB. <i>FASEB Journal</i> , 2019, 33, 1989-1999.	0.5	28
35	A Highly Strained Nuclear Conformation of the Exportin Cse1p Revealed by Molecular Dynamics Simulations. <i>Structure</i> , 2006, 14, 1469-1478.	3.3	27
36	Multistep Mechanism of Chloride Translocation in a Strongly Anion-Selective Porin Channel. <i>Biophysical Journal</i> , 2003, 85, 954-962.	0.5	26

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37	A Designed Inhibitor of a CLC Antiporter Blocks Function through a Unique Binding Mode. <i>Chemistry and Biology</i> , 2012, 19, 1460-1470.	6.0	25
38	Conserved structure and domain organization among bacterial Slc26 transporters. <i>Biochemical Journal</i> , 2014, 463, 297-307.	3.7	25
39	A two-lane mechanism for selective biological ammonium transport. <i>ELife</i> , 2020, 9, .	6.0	23
40	Proton uptake associated with the reduction of the primary quinone QA influences the binding site of the secondary quinone QB in <i>Rhodospseudomonas viridis</i> photosynthetic reaction centers. <i>Biochimica Et Biophysica Acta - Bioenergetics</i> , 2001, 1505, 280-290.	1.0	21
41	A Molecular Switch Driving Inactivation in the Cardiac K ⁺ Channel hERG. <i>PLoS ONE</i> , 2012, 7, e41023.	2.5	19
42	Structural Determinants and Mechanism of Mammalian CRM1 Allostery. <i>Structure</i> , 2013, 21, 1350-1360.	3.3	17
43	Improved 3D continuum calculations of ion flux through membrane channels. <i>European Biophysics Journal</i> , 2003, 32, 689-702.	2.2	16
44	Insights into the ion-coupling mechanism in the MATE transporter NorM-VC. <i>Physical Biology</i> , 2017, 14, 045009.	1.8	16
45	On the ion coupling mechanism of the MATE transporter ClbM. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2020, 1862, 183137.	2.6	16
46	Positively selected modifications in the pore of TbAQP2 allow pentamidine to enter <i>Trypanosoma brucei</i> . <i>ELife</i> , 2020, 9, .	6.0	16
47	Merging In-Solution X-ray and Neutron Scattering Data Allows Fine Structural Analysis of Membrane-Protein Detergent Complexes. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 3910-3914.	4.6	14
48	Hydrophilic Linkers and Polar Contacts Affect Aggregation of FG Repeat Peptides. <i>Biophysical Journal</i> , 2010, 98, 2653-2661.	0.5	13
49	Identification of a cation transport pathway in <i>Neisseria meningitidis</i> PorB. <i>Proteins: Structure, Function and Bioinformatics</i> , 2013, 81, 830-840.	2.6	13
50	Universal Relaxation Governs the Nonequilibrium Elasticity of Biomolecules. <i>Physical Review Letters</i> , 2012, 109, 118304.	7.8	12
51	Effect of Protonation State on the Stability of Amyloid Oligomers Assembled from TTR(105-115). <i>Journal of Physical Chemistry Letters</i> , 2013, 4, 1233-1238.	4.6	12
52	Role of Structural Dynamics at the Receptor G Protein Interface for Signal Transduction. <i>PLoS ONE</i> , 2015, 10, e0143399.	2.5	12
53	High-resolution experimental and computational electrophysiology reveals weak \hat{I}^2 -lactam binding events in the porin PorB. <i>Scientific Reports</i> , 2019, 9, 1264.	3.3	12
54	An antibiotic-resistance conferring mutation in a neisserial porin: Structure, ion flux, and ampicillin binding. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2021, 1863, 183601.	2.6	9

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55	Data-Driven Derivation of Molecular Substructures That Enhance Drug Activity in Gram-Negative Bacteria. <i>Journal of Medicinal Chemistry</i> , 2022, 65, 6088-6099.	6.4	8
56	A role for loop G in the $\hat{2}1$ strand in GABA _A receptor activation. <i>Journal of Physiology</i> , 2016, 594, 5555-5571.	2.9	7
57	Disease related single point mutations alter the global dynamics of a tetratricopeptide (TPR) $\hat{1}\pm$ -solenoid domain. <i>Journal of Structural Biology</i> , 2020, 209, 107405.	2.8	7
58	Modulation of the <i>Neisseria gonorrhoeae</i> drug efflux conduit MtrE. <i>Scientific Reports</i> , 2017, 7, 17091.	3.3	6
59	Coexistence of Ammonium Transporter and Channel Mechanisms in Amt-Mep-Rh Twin-His Variants Impairs the Filamentation Signaling Capacity of Fungal Mep2 Transceptors. <i>MBio</i> , 2022, 13, e0291321.	4.1	6
60	Computational Electrophysiology on Vdac-1 reveals Mechanism of Anion Flux. <i>Biophysical Journal</i> , 2011, 100, 267a.	0.5	0
61	Structural Determinants of Conformational Flexibility and Long-Range Allostery of the CRM1 Export Complex. <i>Biophysical Journal</i> , 2013, 104, 68a.	0.5	0
62	Channel Crystal Structure and Antimicrobial Mechanism of Dermcidin from Human Skin. <i>Biophysical Journal</i> , 2013, 104, 241a.	0.5	0
63	Direct Contacts of K ⁺ Ions in the Selectivity Filter Enable the High Conductance of K ⁺ Channels. <i>Biophysical Journal</i> , 2015, 108, 129a.	0.5	0
64	Characterisation of the Rh50 protein from the ammonia-oxidising bacterium <i>Nitrosomonas europaea</i> . <i>Access Microbiology</i> , 2022, 4, .	0.5	0