

M R Gunner

List of Publications by Year in descending order

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Version: 2024-02-01

52
papers

3,194
citations

236925

25
h-index

189892

50
g-index

53
all docs

53
docs citations

53
times ranked

2690
citing authors

#	ARTICLE	IF	CITATIONS
1	High-resolution cryo-electron microscopy structure of photosystem II from the mesophilic cyanobacterium, <i>Synechocystis</i> sp. PCC 6803. Proceedings of the National Academy of Sciences of the United States of America, 2022, 119, .	7.1	58
2	Comparison of proton transfer paths to the QA and QB sites of the Rb. sphaeroides photosynthetic reaction centers. Photosynthesis Research, 2022, 152, 153-165.	2.9	10
3	Characterizing Protein Protonation Microstates Using Monte Carlo Sampling. Journal of Physical Chemistry B, 2022, 126, 2476-2485.	2.6	9
4	Overview of the SAMPL6 pKa challenge: evaluating small molecule microscopic and macroscopic pKa predictions. Journal of Computer-Aided Molecular Design, 2021, 35, 131-166.	2.9	23
5	Poor Person's pH Simulation of Membrane Proteins. Methods in Molecular Biology, 2021, 2315, 197-217.	0.9	1
6	Evaluation of log $\hat{A}P$, pKa, and log $\hat{A}D$ predictions from the SAMPL7 blind challenge. Journal of Computer-Aided Molecular Design, 2021, 35, 771-802.	2.9	42
7	Protein Motifs for Proton Transfers That Build the Transmembrane Proton Gradient. Frontiers in Chemistry, 2021, 9, 660954.	3.6	15
8	Proton exit pathways surrounding the oxygen evolving complex of photosystem II. Biochimica Et Biophysica Acta - Bioenergetics, 2021, 1862, 148446.	1.0	30
9	Characterizing the Water Wire in the Gramicidin Channel Found by Monte Carlo Sampling Using Continuum Electrostatics and in Molecular Dynamics Trajectories with Conventional or Polarizable Force Fields. Journal of Computational Biophysics and Chemistry, 2021, 20, 111-130.	1.7	11
10	Mesoscopic to Macroscopic Electron Transfer by Hopping in a Crystal Network of Cytochromes. Journal of the American Chemical Society, 2020, 142, 10459-10467.	13.7	13
11	Identifying the proton loading site cluster in the ba cytochrome c oxidase that loads and traps protons. Biochimica Et Biophysica Acta - Bioenergetics, 2020, 1861, 148239.	1.0	13
12	Hydrogen bond network analysis reveals the pathway for the proton transfer in the E-channel of T. thermophilus Complex I. Biochimica Et Biophysica Acta - Bioenergetics, 2020, 1861, 148240.	1.0	20
13	Identification of a Na ⁺ -Binding Site near the Oxygen-Evolving Complex of Spinach Photosystem II. Biochemistry, 2020, 59, 2823-2831.	2.5	5
14	Standard state free energies, not pKas, are ideal for describing small molecule protonation and tautomeric states. Journal of Computer-Aided Molecular Design, 2020, 34, 561-573.	2.9	20
15	Charge Transfer and Chemo-Mechanical Coupling in Respiratory Complex I. Journal of the American Chemical Society, 2020, 142, 9220-9230.	13.7	22
16	Thermodynamics of the S ₂ -to-S ₃ state transition of the oxygen-evolving complex of photosystem II. Physical Chemistry Chemical Physics, 2019, 21, 20840-20848.	2.8	21
17	Relative stability of the S2 isomers of the oxygen evolving complex of photosystem II. Photosynthesis Research, 2019, 141, 331-341.	2.9	18
18	Photosystem II oxygen-evolving complex photoassembly displays an inverse H/D solvent isotope effect under chloride-limiting conditions. Proceedings of the National Academy of Sciences of the United States of America, 2019, 116, 18917-18922.	7.1	41

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19	Network analysis of a proposed exit pathway for protons to the P-side of cytochrome c oxidase. <i>Biochimica Et Biophysica Acta - Bioenergetics</i> , 2018, 1859, 997-1005.	1.0	35
20	L-2-Hydroxyglutarate production arises from noncanonical enzyme function at acidic pH. <i>Nature Chemical Biology</i> , 2017, 13, 494-500.	8.0	190
21	X-ray Free Electron Laser Radiation Damage through the S-State Cycle of the Oxygen-Evolving Complex of Photosystem II. <i>Journal of Physical Chemistry B</i> , 2017, 121, 9382-9388.	2.6	14
22	Two Cl Ions and a Glu Compete for a Helix Cage in the CLC Proton/Cl ⁻ Antiporter. <i>Biophysical Journal</i> , 2017, 113, 1025-1036.	0.5	16
23	The design features cells use to build their transmembrane proton gradient. <i>Physical Biology</i> , 2017, 14, 013001.	1.8	4
24	Unraveling the mechanism of proton translocation in the extracellular half-channel of bacteriorhodopsin. <i>Proteins: Structure, Function and Bioinformatics</i> , 2016, 84, 639-654.	2.6	17
25	Continuum Electrostatics Approaches to Calculating pKas and Ems in Proteins. <i>Methods in Enzymology</i> , 2016, 578, 1-20.	1.0	27
26	Proton-Coupled Electron Transfer During the S-State Transitions of the Oxygen-Evolving Complex of Photosystem II. <i>Journal of Physical Chemistry B</i> , 2015, 119, 7366-7377.	2.6	49
27	Affinity and activity of non-native quinones at the QB site of bacterial photosynthetic reaction centers. <i>Photosynthesis Research</i> , 2014, 120, 181-196.	2.9	6
28	Halorhodopsin pumps Cl ⁻ and bacteriorhodopsin pumps protons by a common mechanism that uses conserved electrostatic interactions. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2014, 111, 16377-16382.	7.1	24
29	Characterizing the proton loading site in cytochrome c oxidase. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2014, 111, 12414-12419.	7.1	54
30	Molecular mechanisms for generating transmembrane proton gradients. <i>Biochimica Et Biophysica Acta - Bioenergetics</i> , 2013, 1827, 892-913.	1.0	37
31	MCCE analysis of the pK _a s of introduced buried acids and bases in staphylococcal nuclease. <i>Proteins: Structure, Function and Bioinformatics</i> , 2011, 79, 3306-3319.	2.6	56
32	The pK _a Cooperative: A collaborative effort to advance structure-based calculations of pK _a values and electrostatic effects in proteins. <i>Proteins: Structure, Function and Bioinformatics</i> , 2011, 79, 3249-3259.	2.6	105
33	The measured and calculated affinity of methyl- and methoxy-substituted benzoquinones for the Q _A site of bacterial reaction centers. <i>Proteins: Structure, Function and Bioinformatics</i> , 2010, 78, 2638-2654.	2.6	11
34	MCCE2: Improving protein pK _a calculations with extensive side chain rotamer sampling. <i>Journal of Computational Chemistry</i> , 2009, 30, 2231-2247.	3.3	192
35	Analysis of the electrochemistry of hemes with E _m s spanning 800 mV. <i>Proteins: Structure, Function and Bioinformatics</i> , 2009, 75, 719-734.	2.6	81
36	Using Multiconformation Continuum Electrostatics to Compare Chloride Binding Motifs in α -Amylase, Human Serum Albumin, and Omp32. <i>Journal of Molecular Biology</i> , 2009, 387, 840-856.	4.2	41

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37	Computational analysis of photosynthetic systems. <i>Photosynthesis Research</i> , 2008, 97, 1-3.	2.9	3
38	Modification of quinone electrochemistry by the proteins in the biological electron transfer chains: examples from photosynthetic reaction centers. <i>Journal of Bioenergetics and Biomembranes</i> , 2008, 40, 509-19.	2.3	57
39	Ligand preference and orientation in α - and β -type heme-binding proteins. <i>Proteins: Structure, Function and Bioinformatics</i> , 2008, 73, 690-704.	2.6	46
40	Factors influencing the energetics of electron and proton transfers in proteins. What can be learned from calculations. <i>Biochimica Et Biophysica Acta - Bioenergetics</i> , 2006, 1757, 942-968.	1.0	89
41	Are Acidic and Basic Groups in Buried Proteins Predicted to be Ionized?. <i>Journal of Molecular Biology</i> , 2005, 348, 1283-1298.	4.2	106
42	Modeling the First Electron Transfer from QA to QB in Reaction Center Proteins from <i>Rb. sphaeroid.</i> <i>ACS Symposium Series</i> , 2004, , 93-106.	0.5	0
43	Characterization of a symmetrized mutant RC with 42 residues from the QA site replacing residues in the Q(B) site. <i>Photosynthesis Research</i> , 2000, 64, 41-52.	2.9	5
44	Backbone Dipoles Generate Positive Potentials in all Proteins: Origins and Implications of the Effect. <i>Biophysical Journal</i> , 2000, 78, 1126-1144.	0.5	82
45	Modeling the Effects of Mutations on the Free Energy of the First Electron Transfer from QA ⁻ to QB in Photosynthetic Reaction Centers. <i>Biochemistry</i> , 2000, 39, 5940-5952.	2.5	56
46	Temperature Dependence of the Free Energy, Enthalpy, and Entropy of P+QA-Charge Recombination in <i>Rhodobacter sphaeroides</i> R-26 Reaction Centers. <i>Journal of Physical Chemistry B</i> , 2000, 104, 8035-8043.	2.6	38
47	Calculated Protein and Proton Motions Coupled to Electron Transfer: A Electron Transfer from QA-to QB in Bacterial Photosynthetic Reaction Centers. <i>Biochemistry</i> , 1999, 38, 8253-8270.	2.5	243
48	Incorporating protein conformational flexibility into the calculation of pH-dependent protein properties. <i>Biophysical Journal</i> , 1997, 72, 2075-2093.	0.5	343
49	The importance of the protein in controlling the electrochemistry of heme metalloproteins: methods of calculation and analysis. <i>Journal of Biological Inorganic Chemistry</i> , 1997, 2, 126-134.	2.6	84
50	Calculated coupling of electron and proton transfer in the photosynthetic reaction center of <i>Rhodospseudomonas viridis</i> . <i>Biophysical Journal</i> , 1996, 70, 2469-2492.	0.5	194
51	On the calculation of pKas in proteins. <i>Proteins: Structure, Function and Bioinformatics</i> , 1993, 15, 252-265.	2.6	514
52	Characterizing the water wire in the Gramicidin channel found by Monte Carlo sampling using continuum electrostatics and in molecular dynamics trajectories with conventional or polarizable force fields. <i>Journal of Theoretical and Computational Chemistry</i> , 0, , 2042001.	1.8	2