Joseph L Baker

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	Electroluminescent Zinc(II) Bis(8-hydroxyquinoline):Â Structural Effects on Electronic States and Device Performance. Journal of the American Chemical Society, 2002, 124, 6119-6125.	13.7	260
2	Quantitative PCR assays for mouse enteric flora reveal strain-dependent differences in composition that are influenced by the microenvironment. Mammalian Genome, 2006, 17, 1093-1104.	2.2	124
3	Quantitative determination of mechanical stability in the novel coronavirus spike protein. Nanoscale, 2020, 12, 16409-16413.	5.6	49
4	Characterization of Structural and Energetic Differences between Conformations of the SARS-CoV-2 Spike Protein. Materials, 2020, 13, 5362.	2.9	46
5	Twelve Transmembrane Helices Form the Functional Core of Mammalian MATE1 (Multidrug and Toxin) Tj ETQq1 3	l 0.78431 3.4	4 rgBT /Ove
6	Cations Stiffen Actin Filaments by Adhering a Key Structural Element to Adjacent Subunits. Journal of Physical Chemistry B, 2016, 120, 4558-4567.	2.6	39
7	Electrostatic Interactions between the Bni1p Formin FH2 Domain and Actin Influence Actin Filament Nucleation. Structure, 2015, 23, 68-79.	3.3	24
8	Steered Molecular Dynamics Simulations of a Type IV Pilus Probe Initial Stages of a Force-Induced Conformational Transition. PLoS Computational Biology, 2013, 9, e1003032.	3.2	22
9	Simulations of substrate transport in the multidrug transporter EmrD. Proteins: Structure, Function and Bioinformatics, 2012, 80, 1620-1632.	2.6	20
10	Scope and efficacy of the broad-spectrum topical antiseptic choline geranate. PLoS ONE, 2019, 14, e0222211.	2.5	16
11	Probing late neutrino mass properties with supernova neutrinos. Physical Review D, 2007, 76, .	4.7	15
12	Effects of ATP and Actin-Filament Binding on the Dynamics of the Myosin II S1 Domain. Biophysical Journal, 2013, 105, 1624-1634.	0.5	13
13	Nearest-Neighbor-Atom Core-Hole Transfer in Isolated Molecules. Physical Review Letters, 2004, 92, 223002.	7.8	12
14	Network visualization of conformational sampling during molecular dynamics simulation. Journal of Molecular Graphics and Modelling, 2013, 46, 140-149.	2.4	11
15	Influence of the ionic liquid [C 4 mpy][Tf 2 N] on the structure of the miniprotein Trp-cage. Journal of Molecular Graphics and Modelling, 2015, 62, 202-212.	2.4	10
16	Theory of Change to Practice: How Experimentalist Teaching Enabled Faculty to Navigate the COVID-19 Disruption. Journal of Chemical Education, 2020, 97, 2788-2792.	2.3	8
17	Influence of an ionic liquid on the conformational sampling of Xaa-Pro dipeptides. Journal of Molecular Liquids, 2017, 227, 66-75.	4.9	7
18	The ionic liquid [C ₄ mpy][Tf ₂ N] induces bound-like structure in the intrinsically disordered protein FlgM. Physical Chemistry Chemical Physics, 2019, 21, 17950-17958.	2.8	7

JOSEPH L BAKER

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19	Impact of an alpha helix and a cysteine–cysteine disulfide bond on the resistance of bacterial adhesion pili to stress. Proceedings of the National Academy of Sciences of the United States of America, 2021, 118, .	7.1	6
20	Molecular simulations of lactose-bound and unbound forms of the FaeG adhesin reveal critical amino acids involved in sugar binding. Journal of Molecular Graphics and Modelling, 2016, 70, 100-108.	2.4	3
21	Density, Enthalpy of Vaporization and Local Structure of Neat N-Alkane Liquids. Liquids, 2021, 1, 47-59.	2.5	3
22	Long-ranged heterogeneous structure in aqueous solutions of the deep eutectic solvent choline and geranate at the liquid–vapor interface. Physical Chemistry Chemical Physics, 2022, 24, 13720-13729.	2.8	3
23	Unveiling molecular interactions that stabilize bacterial adhesion pili. Biophysical Journal, 2022, 121, 2096-2106.	0.5	2
24	Assessing the Stability of Biological Fibrils by Molecular-Scale Simulations. Methods in Molecular Biology, 2022, 2340, 357-378.	0.9	1