

Ariane Nunes-Alves

List of Publications by Year in descending order

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Version: 2024-02-01

18
papers

428
citations

1039880

9
h-index

887953

17
g-index

23
all docs

23
docs citations

23
times ranked

415
citing authors

#	ARTICLE	IF	CITATIONS
1	The Non-phosphorylating Glyceraldehyde-3-Phosphate Dehydrogenase GapN Is a Potential New Drug Target in <i>Streptococcus pyogenes</i> . <i>Frontiers in Microbiology</i> , 2022, 13, 802427.	1.5	2
2	Ligand unbinding mechanisms and kinetics for T4 lysozyme mutants from \bar{I} ,RAMD simulations. <i>Current Research in Structural Biology</i> , 2021, 3, 106-111.	1.1	31
3	A Protocol to Use Comparative Binding Energy Analysis to Estimate Drug-Target Residence Time. <i>Methods in Molecular Biology</i> , 2021, 2266, 171-186.	0.4	1
4	Line-FRAP, A Versatile Method to Measure Diffusion Rates In Vitro and In Vivo. <i>Journal of Molecular Biology</i> , 2021, 433, 166898.	2.0	14
5	Structure-kinetic relationship reveals the mechanism of selectivity of FAK inhibitors over PYK2. <i>Cell Chemical Biology</i> , 2021, 28, 686-698.e7.	2.5	36
6	Prediction of the Drug-Target Binding Kinetics for Flexible Proteins by Comparative Binding Energy Analysis. <i>Journal of Chemical Information and Modeling</i> , 2021, 61, 3708-3721.	2.5	12
7	Molecular Dynamics Simulations in Drug Discovery and Pharmaceutical Development. <i>Processes</i> , 2021, 9, 71.	1.3	162
8	From Brazil to Germany: Challenges and Advantages. <i>Journal of Chemical Information and Modeling</i> , 2020, 60, 449-451.	2.5	1
9	Recent progress in molecular simulation methods for drug binding kinetics. <i>Current Opinion in Structural Biology</i> , 2020, 64, 126-133.	2.6	61
10	Structural characterization of an Arf dimer interface: molecular mechanism of Arf-dependent membrane scission. <i>FEBS Letters</i> , 2020, 594, 2240-2253.	1.3	12
11	Myotubularin-related protein 7 activates peroxisome proliferator-activated receptor-gamma. <i>Oncogenesis</i> , 2020, 9, 59.	2.1	6
12	Impact of the <i>Journal of Chemical Information and Modeling</i> Special Issue on Women in Computational Chemistry. <i>Journal of Chemical Information and Modeling</i> , 2020, 60, 3328-3330.	2.5	5
13	What Makes a Paper Be Highly Cited? 60 Years of the <i>Journal of Chemical Information and Modeling</i> . <i>Journal of Chemical Information and Modeling</i> , 2020, 60, 5866-5867.	2.5	1
14	Escape of a Small Molecule from Inside T4 Lysozyme by Multiple Pathways. <i>Biophysical Journal</i> , 2018, 114, 1058-1066.	0.2	44
15	Mechanical Unfolding of Macromolecules Coupled to Bond Dissociation. <i>Journal of Chemical Theory and Computation</i> , 2018, 14, 282-290.	2.3	18
16	Ligand-Receptor Affinities Computed by an Adapted Linear Interaction Model for Continuum Electrostatics and by Protein Conformational Averaging. <i>Journal of Chemical Information and Modeling</i> , 2014, 54, 2309-2319.	2.5	17
17	Tobacco Nitrosamine N-nitrosornicotine as Inhibitor of Neuronal Nicotinic Acetylcholine Receptors. <i>Journal of Molecular Neuroscience</i> , 2013, 49, 52-61.	1.1	5
18	Neurotransmitters as Main Players in the Neural Differentiation and Fate Determination Game. , 2010, , 115-134.		0