

Jure Borisek

List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

24
papers

171
citations

9
h-index

12
g-index

27
ext. papers

323
ext. citations

8.1
avg, IF

3.67
L-index

#	Paper	IF	Citations
24	An oomycete NLP cytolysin forms transient small pores in lipid membranes.. <i>Science Advances</i> , 2022 , 8, eabj9406	14.3	1
23	Molecular Basis of SARS-CoV-2 Nsp1-Induced Immune Translational Shutdown as Revealed by All-Atom Simulations. <i>Journal of Physical Chemistry Letters</i> , 2021 , 12, 11745-11750	6.4	1
22	Senescent cells as promising targets to tackle age-related diseases. <i>Ageing Research Reviews</i> , 2021 , 66, 101251	12	9
21	An Expanded Two-Zn ²⁺ -Ion Motif Orchestrates Pre-mRNA Maturation in the 3'-End Processing Endonuclease Machinery. <i>ACS Catalysis</i> , 2021 , 11, 4319-4326	13.1	3
20	Nep1-like proteins as a target for plant pathogen control. <i>PLoS Pathogens</i> , 2021 , 17, e1009477	7.6	2
19	Computing Metal-Binding Proteins for Therapeutic Benefit. <i>ChemMedChem</i> , 2021 , 16, 2034-2049	3.7	0
18	Allosteric Cross-Talk among Spike ^T Receptor-Binding Domain Mutations of the SARS-CoV-2 South African Variant Triggers an Effective Hijacking of Human Cell Receptor. <i>Journal of Physical Chemistry Letters</i> , 2021 , 12, 5987-5993	6.4	9
17	All-Atom Simulations Reveal a Key Interaction Network in the HLA-E/NKG2A/CD94 Immune Complex Fine-Tuned by the Nonameric Peptide. <i>Journal of Chemical Information and Modeling</i> , 2021 , 61, 3593-3603	6.1	0
16	Atomic-Level Mechanism of Pre-mRNA Splicing in Health and Disease. <i>Accounts of Chemical Research</i> , 2021 , 54, 144-154	24.3	4
15	Domain sliding of two Staphylococcus aureus N-acetylglucosaminidases enables their substrate-binding prior to its catalysis. <i>Communications Biology</i> , 2020 , 3, 178	6.7	4
14	All-Atom Simulations Decrypt the Molecular Terms of RNA Catalysis in the Exon-Ligation Step of the Spliceosome. <i>ACS Catalysis</i> , 2020 , 10, 5328-5334	13.1	12
13	Decrypting the Information Exchange Pathways across the Spliceosome Machinery. <i>Journal of the American Chemical Society</i> , 2020 , 142, 8403-8411	16.4	13
12	Peptide Inhibitors of Bacterial Protein Synthesis with Broad Spectrum and SbmA-Independent Bactericidal Activity against Clinical Pathogens. <i>Journal of Medicinal Chemistry</i> , 2020 , 63, 9590-9602	8.3	11
11	Exploiting Cryo-EM Structural Information and All-Atom Simulations To Decrypt the Molecular Mechanism of Splicing Modulators. <i>Journal of Chemical Information and Modeling</i> , 2020 , 60, 2510-2521	6.1	9
10	Molecular basis for functional diversity among microbial Nep1-like proteins. <i>PLoS Pathogens</i> , 2019 , 15, e1007951	7.6	10
9	Disclosing the Impact of Carcinogenic SF3b Mutations on Pre-mRNA Recognition Via All-Atom Simulations. <i>Biomolecules</i> , 2019 , 9,	5.9	10
8	A Water-Assisted Catalytic Mechanism in Glycoside Hydrolases Demonstrated on the Staphylococcus aureus Autolysin E. <i>ACS Catalysis</i> , 2018 , 8, 4334-4345	13.1	11

7	Discovery of (phenylureido)piperidinyl benzamides as prospective inhibitors of bacterial autolysin E from <i>Staphylococcus aureus</i> . <i>Journal of Enzyme Inhibition and Medicinal Chemistry</i> , 2018 , 33, 1239-1247	5.6	3
6	The CWB2 Cell Wall-Anchoring Module Is Revealed by the Crystal Structures of the <i>Clostridium difficile</i> Cell Wall Proteins Cwp8 and Cwp6. <i>Structure</i> , 2017 , 25, 514-521	5.2	23
5	The Comparison of Docking Search Algorithms and Scoring Functions 2017 , 820-849		
4	Comparison of in silico tools for binding site prediction applied for structure-based design of autolysin inhibitors. <i>SAR and QSAR in Environmental Research</i> , 2016 , 27, 573-87	3.5	4
3	The Comparison of Docking Search Algorithms and Scoring Functions. <i>Advances in Medical Technologies and Clinical Practice Book Series</i> , 2016 , 99-127	0.3	4
2	Development of N-(Functionalized benzoyl)-homocycloleucyl-glycinonitriles as Potent Cathepsin K Inhibitors. <i>Journal of Medicinal Chemistry</i> , 2015 , 58, 6928-37	8.3	19
1	Mechanistic interpretation of artificial neural network-based QSAR model for prediction of cathepsin K inhibition potency. <i>Journal of Chemometrics</i> , 2014 , 28, 272-281	1.6	9