

Bian Li

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/6357964/publications.pdf>

Version: 2024-02-01

20
papers

527
citations

932766

10
h-index

794141

19
g-index

27
all docs

27
docs citations

27
times ranked

884
citing authors

#	ARTICLE	IF	CITATIONS
1	Micelle-Based Brain-Targeted Drug Delivery Enabled by a Nicotine Acetylcholine Receptor Ligand. <i>Angewandte Chemie - International Edition</i> , 2011, 50, 5482-5485.	7.2	124
2	High-Throughput Reclassification of SCN5A Variants. <i>American Journal of Human Genetics</i> , 2020, 107, 111-123.	2.6	88
3	Predicting changes in protein thermodynamic stability upon point mutation with deep 3D convolutional neural networks. <i>PLoS Computational Biology</i> , 2020, 16, e1008291.	1.5	76
4	Predicting the Functional Impact of KCNQ1 Variants of Unknown Significance. <i>Circulation: Cardiovascular Genetics</i> , 2017, 10, .	5.1	40
5	Finding the needle in the haystack: towards solving the protein-folding problem computationally. <i>Critical Reviews in Biochemistry and Molecular Biology</i> , 2018, 53, 1-28.	2.3	31
6	Contribution of Cotranslational Folding Defects to Membrane Protein Homeostasis. <i>Journal of the American Chemical Society</i> , 2019, 141, 204-215.	6.6	27
7	Distinct Features of Proband With Early Repolarization and Brugada Syndromes Carrying SCN5A Pathogenic Variants. <i>Journal of the American College of Cardiology</i> , 2021, 78, 1603-1617.	1.2	22
8	The 3D mutational constraint on amino acid sites in the human proteome. <i>Nature Communications</i> , 2022, 13, .	5.8	15
9	Molecular Insights into the D1R Agonist and D2R/D3R Antagonist Effects of the Natural Product (α^{S})-Stepholidine: Molecular Modeling and Dynamics Simulations. <i>Journal of Physical Chemistry B</i> , 2012, 116, 8121-8130.	1.2	13
10	CASP11 – An Evaluation of a Modular BCL::Fold-Based Protein Structure Prediction Pipeline. <i>PLoS ONE</i> , 2016, 11, e0152517.	1.1	13
11	Structural determinants of cholesterol recognition in helical integral membrane proteins. <i>Biophysical Journal</i> , 2021, 120, 1592-1604.	0.2	12
12	Improving prediction of helix-helix packing in membrane proteins using predicted contact numbers as restraints. <i>Proteins: Structure, Function and Bioinformatics</i> , 2017, 85, 1212-1221.	1.5	9
13	Dominant negative effects of SCN5A missense variants. <i>Genetics in Medicine</i> , 2022, 24, 1238-1248.	1.1	9
14	Accurate Prediction of Contact Numbers for Multi-Spanning Helical Membrane Proteins. <i>Journal of Chemical Information and Modeling</i> , 2016, 56, 423-434.	2.5	7
15	Interfaces Between Alpha-helical Integral Membrane Proteins: Characterization, Prediction, and Docking. <i>Computational and Structural Biotechnology Journal</i> , 2019, 17, 699-711.	1.9	7
16	Building a Hybrid Physical-Statistical Classifier for Predicting the Effect of Variants Related to Protein-Drug Interactions. <i>Structure</i> , 2019, 27, 1469-1481.e3.	1.6	6
17	A Multitask Deep-Learning Method for Predicting Membrane Associations and Secondary Structures of Proteins. <i>Journal of Proteome Research</i> , 2021, 20, 4089-4100.	1.8	6
18	Discovery of novel small-molecule Src kinase inhibitors via a kinase-focused druglikeness rule and structure-based virtual screening. <i>Molecular Simulation</i> , 2014, 40, 341-348.	0.9	1

#	ARTICLE	IF	CITATIONS
19	Clinical Characteristics and Electrophysiologic Properties of <i>SCN5A</i> Variants in Fever-Induced Brugada Syndrome. SSRN Electronic Journal, 0, , .	0.4	0
20	Integration of Protein Structure and Population-Scale DNA Sequence Data for Disease Gene Discovery and Variant Interpretation. Annual Review of Biomedical Data Science, 2022, 5, .	2.8	0