

Shaun M Kandathil

List of Publications by Year in descending order

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Version: 2024-02-01

18
papers

1,635
citations

758635

12
h-index

996533

15
g-index

21
all docs

21
docs citations

21
times ranked

1469
citing authors

#	ARTICLE	IF	CITATIONS
1	A guide to machine learning for biologists. <i>Nature Reviews Molecular Cell Biology</i> , 2022, 23, 40-55.	16.1	626
2	A review of the chemistry and pharmacology of the date fruits (<i>Phoenix dactylifera</i> L.). <i>Food Research International</i> , 2011, 44, 1812-1822.	2.9	320
3	High precision in protein contact prediction using fully convolutional neural networks and minimal sequence features. <i>Bioinformatics</i> , 2018, 34, 3308-3315.	1.8	157
4	Deep learning extends de novo protein modelling coverage of genomes using iteratively predicted structural constraints. <i>Nature Communications</i> , 2019, 10, 3977.	5.8	144
5	Prediction of interresidue contacts with DeepMetaPSICOV in CASP13. <i>Proteins: Structure, Function and Bioinformatics</i> , 2019, 87, 1092-1099.	1.5	102
6	Recent developments in deep learning applied to protein structure prediction. <i>Proteins: Structure, Function and Bioinformatics</i> , 2019, 87, 1179-1189.	1.5	56
7	Accuracy and tractability of a kriging model of intramolecular polarizable multipolar electrostatics and its application to histidine. <i>Journal of Computational Chemistry</i> , 2013, 34, 1850-1861.	1.5	47
8	Generating, Maintaining, and Exploiting Diversity in a Memetic Algorithm for Protein Structure Prediction. <i>Evolutionary Computation</i> , 2016, 24, 577-607.	2.3	38
9	Ultrafast end-to-end protein structure prediction enables high-throughput exploration of uncharacterized proteins. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2022, 119, .	3.3	26
10	The prediction of atomic kinetic energies from coordinates of surrounding atoms using kriging machine learning. <i>Theoretical Chemistry Accounts</i> , 2014, 133, 1.	0.5	20
11	Accurate prediction of polarised high order electrostatic interactions for hydrogen bonded complexes using the machine learning method kriging. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2015, 136, 32-41.	2.0	19
12	Toward a detailed understanding of search trajectories in fragment assembly approaches to protein structure prediction. <i>Proteins: Structure, Function and Bioinformatics</i> , 2016, 84, 411-426.	1.5	18
13	Improved fragment-based protein structure prediction by redesign of search heuristics. <i>Scientific Reports</i> , 2018, 8, 13694.	1.6	12
14	Proton tunnelling and promoting vibrations during the oxidation of ascorbate by ferricyanide?. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 2256.	1.3	10
15	Uncommon mutational profiles of metastatic colorectal cancer detected during routine genotyping using next generation sequencing. <i>Scientific Reports</i> , 2019, 9, 7083.	1.6	5
16	Using Machine Learning to Explore the Relevance of Local and Global Features During Conformational Search in Rosetta. , 2015, , .		1
17	On heuristic bias in fragment-assembly methods for protein structure prediction. , 2017, , .		0
18	Reliable Generation of Native-Like Decoys Limits Predictive Ability in Fragment-Based Protein Structure Prediction. <i>Biomolecules</i> , 2019, 9, 612.	1.8	0