Martino Bertoni

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.

20 8,361 8 23 g-index

23 11,621 9.3 5.98 ext. papers ext. citations avg, IF L-index

#	Paper	IF	Citations
20	A community challenge for a pancancer drug mechanism of action inference from perturbational profile data <i>Cell Reports Medicine</i> , 2022 , 3, 100492	18	5
19	Bioactivity descriptors for uncharacterized chemical compounds. <i>Nature Communications</i> , 2021 , 12, 393	3 2 17.4	8
18	Extending the small-molecule similarity principle to all levels of biology with the Chemical Checker. <i>Nature Biotechnology</i> , 2020 , 38, 1087-1096	44.5	29
17	Bioactivity Profile Similarities to Expand the Repertoire of COVID-19 Drugs. <i>Journal of Chemical Information and Modeling</i> , 2020 , 60, 5730-5734	6.1	1
16	Introducing "best single template" models as reference baseline for the Continuous Automated Model Evaluation (CAMEO). <i>Proteins: Structure, Function and Bioinformatics</i> , 2019 , 87, 1378-1387	4.2	17
15	Cover Image, Volume 9, Issue 6. <i>Wiley Interdisciplinary Reviews: Computational Molecular Science</i> , 2019 , 9, e1451	7.9	
14	Formatting biological big data for modern machine learning in drug discovery. <i>Wiley Interdisciplinary Reviews: Computational Molecular Science</i> , 2019 , 9, e1408	7.9	6
13	Modeling of Protein Tertiary and Quaternary Structures Based on Evolutionary Information. <i>Methods in Molecular Biology</i> , 2019 , 1851, 301-316	1.4	6
12	Assessment of protein assembly prediction in CASP12. <i>Proteins: Structure, Function and Bioinformatics</i> , 2018 , 86 Suppl 1, 247-256	4.2	37
11	Continuous Automated Model EvaluatiOn (CAMEO) complementing the critical assessment of structure prediction in CASP12. <i>Proteins: Structure, Function and Bioinformatics</i> , 2018 , 86 Suppl 1, 387-3	19 <mark>8</mark> .2	75
10	DynBench3D, a Web-Resource to Dynamically Generate Benchmark Sets of Large Heteromeric Protein Complexes. <i>Journal of Molecular Biology</i> , 2018 , 430, 4431-4438	6.5	3
9	SWISS-MODEL: homology modelling of protein structures and complexes. <i>Nucleic Acids Research</i> , 2018 , 46, W296-W303	20.1	4367
8	Modeling protein quaternary structure of homo- and hetero-oligomers beyond binary interactions by homology. <i>Scientific Reports</i> , 2017 , 7, 10480	4.9	357
7	A non-deterministic approach to forecasting the trophic evolution of lakes. <i>Journal of Limnology</i> , 2016 , 75,	1.5	2
6	SWISS-MODEL: modelling protein tertiary and quaternary structure using evolutionary information. <i>Nucleic Acids Research</i> , 2014 , 42, W252-8	20.1	3437
5	Gene regulatory networks reconstruction from time series datasets using genetic programming: a comparison between tree-based and graph-based approaches. <i>Genetic Programming and Evolvable Machines</i> , 2013 , 14, 431-455	2	
4	GeNet: A Graph-Based Genetic Programming Framework for the Reverse Engineering of Gene Regulatory Networks. <i>Lecture Notes in Computer Science</i> , 2012 , 97-109	0.9	2

LIST OF PUBLICATIONS

3	Extending the small molecule similarity principle to all levels of biology	1
2	Bioactivity descriptors for uncharacterized compounds	3
1	A Community Challenge for Pancancer Drug Mechanism of Action Inference from Perturbational Profile Data	3