

Martino Bertoni

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/7525493/publications.pdf>

Version: 2024-02-01

18
papers

13,776
citations

840776

11
h-index

888059

17
g-index

23
all docs

23
docs citations

23
times ranked

27346
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|---|------|-----------|
| 1 | A community challenge for a pancancer drug mechanism of action inference from perturbational profile data. <i>Cell Reports Medicine</i> , 2022, 3, 100492. | 6.5 | 33 |
| 2 | Bioactivity descriptors for uncharacterized chemical compounds. <i>Nature Communications</i> , 2021, 12, 3932. | 12.8 | 44 |
| 3 | Bioactivity Profile Similarities to Expand the Repertoire of COVID-19 Drugs. <i>Journal of Chemical Information and Modeling</i> , 2020, 60, 5730-5734. | 5.4 | 3 |
| 4 | Extending the small-molecule similarity principle to all levels of biology with the Chemical Checker. <i>Nature Biotechnology</i> , 2020, 38, 1087-1096. | 17.5 | 78 |
| 5 | 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. | 2.6 | 31 |
| 6 | Cover Image, Volume 9, Issue 6. <i>Wiley Interdisciplinary Reviews: Computational Molecular Science</i> , 2019, 9, e1451. | 14.6 | 0 |
| 7 | Formatting biological big data for modern machine learning in drug discovery. <i>Wiley Interdisciplinary Reviews: Computational Molecular Science</i> , 2019, 9, e1408. | 14.6 | 17 |
| 8 | Modeling of Protein Tertiary and Quaternary Structures Based on Evolutionary Information. <i>Methods in Molecular Biology</i> , 2019, 1851, 301-316. | 0.9 | 13 |
| 9 | Assessment of protein assembly prediction in CASP12. <i>Proteins: Structure, Function and Bioinformatics</i> , 2018, 86, 247-256. | 2.6 | 54 |
| 10 | Continuous Automated Model EvaluatiOn (CAMEO) complementing the critical assessment of structure prediction in CASP12. <i>Proteins: Structure, Function and Bioinformatics</i> , 2018, 86, 387-398. | 2.6 | 127 |
| 11 | DynBench3D, a Web-Resource to Dynamically Generate Benchmark Sets of Large Heteromeric Protein Complexes. <i>Journal of Molecular Biology</i> , 2018, 430, 4431-4438. | 4.2 | 3 |
| 12 | SWISS-MODEL: homology modelling of protein structures and complexes. <i>Nucleic Acids Research</i> , 2018, 46, W296-W303. | 14.5 | 8,474 |
| 13 | Modeling protein quaternary structure of homo- and hetero-oligomers beyond binary interactions by homology. <i>Scientific Reports</i> , 2017, 7, 10480. | 3.3 | 608 |
| 14 | A non-deterministic approach to forecasting the trophic evolution of lakes. <i>Journal of Limnology</i> , 2016, 75, . | 1.1 | 8 |
| 15 | SWISS-MODEL: modelling protein tertiary and quaternary structure using evolutionary information. <i>Nucleic Acids Research</i> , 2014, 42, W252-W258. | 14.5 | 4,247 |
| 16 | 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. | 2.2 | 1 |
| 17 | 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. | 1.3 | 2 |
| 18 | Extending the Small Molecule Similarity Principle to All Levels of Biology. <i>SSRN Electronic Journal</i> , 0, , . | 0.4 | 0 |