# Oleg Ursu

List of Publications by Year
in descending order

Source: https://exaly.com/author-pdf/7466574/publications.pdf

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|---|---|---|---|
| 17 papers | 2,784 citations | 687363 13 h-index | 794594 19 g-index |
| 20 all docs | 20 docs citations | 20 times ranked | 5092 citing authors |

| # | Article | IF | Citations |
|---|---|---|---|
| 1 | Machine learning prediction and tau-based screening identifies potential Alzheimerâ€™s disease genes relevant to immunity. Communications Biology, 2022, 5, 125. | 4.4 | 18 |
| 2 | PepSeA: Peptide Sequence Alignment and Visualization Tools to Enable Lead Optimization. Journal of Chemical Information and Modeling, 2022, 62, 1259-1267. | 5.4 | 3 |
| 3 | DrugCentral 2021 supports drug discovery and repositioning. Nucleic Acids Research, 2021, 49, D1160-D1169. | 14.5 | 129 |
| 4 | Unexplored therapeutic opportunities in the human genome. Nature Reviews Drug Discovery, 2018, 17, 317-332. | 46.4 | 263 |
| 5 | Chronic obstructive pulmonary disease phenotypes using cluster analysis of electronic medical records. Health Informatics Journal, 2018, 24, 394-409. | 2.1 | 22 |
| 6 | High-Throughput Flow Cytometry Screening of Multidrug Efflux Systems. Methods in Molecular Biology, 2018, 1700, 293-318. | 0.9 | 12 |
| 7 | Pharos: Collating protein information to shed light on the druggable genome. Nucleic Acids Research, 2017, 45, D995-D1002. | 14.5 | 271 |
| 8 | TIN-X: target importance and novelty explorer. Bioinformatics, 2017, 33, 2601-2603. | 4.1 | 27 |
| 9 | A comprehensive map of molecular drug targets. Nature Reviews Drug Discovery, 2017, 16, 19-34. | 46.4 | 1,608 |
| 10 | Formalizing drug indications on the road to therapeutic intent. Journal of the American Medical Informatics Association: JAMIA, 2017, 24, 1169-1172. | 4.4 | 8 |
| 11 | Drug target ontology to classify and integrate drug discovery data. Journal of Biomedical Semantics, 2017, 8, 50. | 1.6 | 63 |
| 12 | Badapple: promiscuity patterns from noisy evidence. Journal of Cheminformatics, 2016, 8, 29. | 6.1 | 85 |
| 13 | Discovery of a specific inhibitor of human GLUT5 by virtual screening and in vitro transport evaluation. Scientific Reports, 2016, 6, 24240. | 3.3 | 45 |
| 14 | Development and validation of a clinical prediction rule for candidemia in hospitalized patients with severe sepsis and septic shock. Journal of Critical Care, 2015, 30, 715-720. | 2.2 | 26 |
| 15 | FRET detection of lymphocyte functionâ€“associated antigen-1 conformational extension. Molecular Biology of the Cell, 2015, 26, 43-54. | 2.1 | 17 |
| 16 | CHEMICAL AND BIOLOGICAL DESCRIPTOR INTEGRATION IMPROVES COMPUTATIONAL MODELING OF RAT TOXICITY. Revue Roumaine De Chimie, 2015, 60, 219-226. | 0.2 | 4 |
| 17 | Model-Free Drug-Likeness from Fragments. Journal of Chemical Information and Modeling, 2010, 50, 1387-1394. | 5.4 | 28 |