

# Martin Å Ñ-cho

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/6592751/publications.pdf>

Version: 2024-02-01

10  
papers

397  
citations

1163117  
8  
h-index

1474206  
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g-index

12  
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docs citations

12  
times ranked

508  
citing authors

#	ARTICLE	IF	CITATIONS
1	GLORYx: Prediction of the Metabolites Resulting from Phase 1 and Phase 2 Biotransformations of Xenobiotics. Chemical Research in Toxicology, 2021, 34, 286-299.	3.3	51
2	CYPlebrity: Machine learning models for the prediction of inhibitors of cytochrome P450 enzymes. Bioorganic and Medicinal Chemistry, 2021, 46, 116388.	3.0	16
3	GenUI: interactive and extensible open source software platform for de novo molecular generation and cheminformatics. Journal of Cheminformatics, 2021, 13, 73.	6.1	7
4	NERDD: a web portal providing access to <i>in silico</i> tools for drug discovery. Bioinformatics, 2020, 36, 1291-1292.	4.1	46
5	FAME 3: Predicting the Sites of Metabolism in Synthetic Compounds and Natural Products for Phase 1 and Phase 2 Metabolic Enzymes. Journal of Chemical Information and Modeling, 2019, 59, 3400-3412.	5.4	60
6	GLORY: Generator of the Structures of Likely Cytochrome P450 Metabolites Based on Predicted Sites of Metabolism. Frontiers in Chemistry, 2019, 7, 402.	3.6	57
7	Hit Dexter 2.0: Machine-Learning Models for the Prediction of Frequent Hitters. Journal of Chemical Information and Modeling, 2019, 59, 1030-1043.	5.4	70
8	Hit Dexter: A Machineâ€Learning Model for the Prediction of Frequent Hitters. ChemMedChem, 2018, 13, 564-571.	3.2	34
9	FAME 2: Simple and Effective Machine Learning Model of Cytochrome P450 Regioselectivity. Journal of Chemical Information and Modeling, 2017, 57, 1832-1846.	5.4	56
10	Activity-driven exploration of chemical space with morphing. , 2015, , .		0