## Antoine Daina

## List of Publications by Citations

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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.

10<br/>papers5,043<br/>citations8<br/>h-index11<br/>g-index11<br/>ext. papers7,681<br/>ext. citations9.8<br/>avg, IF6.65<br/>L-index

#	Paper	IF	Citations
10	SwissADME: a free web tool to evaluate pharmacokinetics, drug-likeness and medicinal chemistry friendliness of small molecules. <i>Scientific Reports</i> , <b>2017</b> , 7, 42717	4.9	3362
9	SwissTargetPrediction: updated data and new features for efficient prediction of protein targets of small molecules. <i>Nucleic Acids Research</i> , <b>2019</b> , 47, W357-W364	20.1	631
8	SwissTargetPrediction: a web server for target prediction of bioactive small molecules. <i>Nucleic Acids Research</i> , <b>2014</b> , 42, W32-8	20.1	557
7	iLOGP: a simple, robust, and efficient description of n-octanol/water partition coefficient for drug design using the GB/SA approach. <i>Journal of Chemical Information and Modeling</i> , <b>2014</b> , 54, 3284-301	6.1	294
6	SwissSimilarity: A Web Tool for Low to Ultra High Throughput Ligand-Based Virtual Screening. Journal of Chemical Information and Modeling, <b>2016</b> , 56, 1399-404	6.1	129
5	Attracting cavities for docking. Replacing the rough energy landscape of the protein by a smooth attracting landscape. <i>Journal of Computational Chemistry</i> , <b>2016</b> , 37, 437-47	3.5	27
4	Application of the SwissDrugDesign Online Resources in Virtual Screening. <i>International Journal of Molecular Sciences</i> , <b>2019</b> , 20,	6.3	25
3	Rational Design, Synthesis, and Pharmacological Characterization of Novel Ghrelin Receptor Inverse Agonists as Potential Treatment against Obesity-Related Metabolic Diseases. <i>Journal of Medicinal Chemistry</i> , <b>2018</b> , 61, 11039-11060	8.3	12
2	SwissBioisostere 2021: updated structural, bioactivity and physicochemical data delivered by a reshaped web interface. <i>Nucleic Acids Research</i> , <b>2021</b> ,	20.1	2
1	Computer-Aided Drug Design for Cancer Therapy <b>2021</b> , 386-401		2