

Mikko J Vainio

List of Publications by Year in descending order

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11
papers

852
citations

1040056

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1281871

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11
times ranked

1424
citing authors

#	ARTICLE	IF	CITATIONS
1	Algorithmic Analysis of Cahn-Ingold-Prelog Rules of Stereochemistry: Proposals for Revised Rules and a Guide for Machine Implementation. <i>Journal of Chemical Information and Modeling</i> , 2018, 58, 1755-1765.	5.4	10
2	Scaffold Hopping by Fragment Replacement. <i>Journal of Chemical Information and Modeling</i> , 2013, 53, 1825-1835.	5.4	31
3	Big pharma screening collections: more of the same or unique libraries? The AstraZeneca-Bayer Pharma AG case. <i>Drug Discovery Today</i> , 2013, 18, 1014-1024.	6.4	58
4	Automated Recycling of Chemistry for Virtual Screening and Library Design. <i>Journal of Chemical Information and Modeling</i> , 2012, 52, 1777-1786.	5.4	18
5	Accurate conformation-dependent molecular electrostatic potentials for high-throughput <i>in silico</i> drug discovery. <i>Journal of Computational Chemistry</i> , 2010, 31, 1722-1732.	3.3	80
6	The Binding of Synthetic Retinoids to Lipocalin β_2 -Lactoglobulins. <i>Journal of Medicinal Chemistry</i> , 2010, 53, 514-518.	6.4	9
7	ShaEP: Molecular Overlay Based on Shape and Electrostatic Potential. <i>Journal of Chemical Information and Modeling</i> , 2009, 49, 492-502.	5.4	177
8	Binding of Phenolic Compounds and Their Derivatives to Bovine and Reindeer β_2 -Lactoglobulin. <i>Journal of Agricultural and Food Chemistry</i> , 2008, 56, 7721-7729.	5.2	85
9	Generating Conformer Ensembles Using a Multiobjective Genetic Algorithm. <i>Journal of Chemical Information and Modeling</i> , 2007, 47, 2462-2474.	5.4	326
10	Similarity Based Virtual Screening: A Tool for Targeted Library Design. <i>Journal of Medicinal Chemistry</i> , 2006, 49, 2353-2356.	6.4	35
11	McQSAR: A Multiconformational Quantitative Structure-Activity Relationship Engine Driven by Genetic Algorithms. <i>Journal of Chemical Information and Modeling</i> , 2005, 45, 1953-1961.	5.4	23