

Laurent Hoffer

List of Publications by Year in descending order

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Version: 2024-02-01

16
papers

363
citations

840119

11
h-index

940134

16
g-index

19
all docs

19
docs citations

19
times ranked

510
citing authors

#	ARTICLE	IF	CITATIONS
1	Integrated Strategy for Lead Optimization Based on Fragment Growing: The Diversity-Oriented-Target-Focused-Synthesis Approach. <i>Journal of Medicinal Chemistry</i> , 2018, 61, 5719-5732.	2.9	51
2	Aurone derivatives as promising antibacterial agents against resistant Gram-positive pathogens. <i>European Journal of Medicinal Chemistry</i> , 2019, 165, 133-141.	2.6	46
3	Fragment-Based Drug Design: Computational and Experimental State of the Art. <i>Combinatorial Chemistry and High Throughput Screening</i> , 2011, 14, 500-520.	0.6	37
4	In Silico Fragment-Based Drug Discovery: Setup and Validation of a Fragment-to-Lead Computational Protocol Using S4MPLE. <i>Journal of Chemical Information and Modeling</i> , 2013, 53, 836-851.	2.5	33
5	Chemistry-Driven Hit-to-Lead Optimization Guided by Structure-Based Approaches. <i>Molecular Informatics</i> , 2018, 37, e1800059.	1.4	33
6	Fr-PPICChem: An Academic Compound Library Dedicated to Protein-Protein Interactions. <i>ACS Chemical Biology</i> , 2020, 15, 1566-1574.	1.6	29
7	S4MPLE – Sampler For Multiple Protein-Ligand Entities: Simultaneous Docking of Several Entities. <i>Journal of Chemical Information and Modeling</i> , 2013, 53, 88-102.	2.5	26
8	S4MPLE – Sampler for Multiple Protein-Ligand Entities: Methodology and Rigid-Site Docking Benchmarking. <i>Molecules</i> , 2015, 20, 8997-9028.	1.7	25
9	Optimization of a fragment linking hit toward Dengue and Zika virus NS5 methyltransferases inhibitors. <i>European Journal of Medicinal Chemistry</i> , 2019, 161, 323-333.	2.6	18
10	Synthesis and Evaluation of the Antibacterial Activities of 13-Substituted Berberine Derivatives. <i>Antibiotics</i> , 2020, 9, 381.	1.5	18
11	<i>In silico</i> molecular target prediction unveils mebendazole as a potent MAPK14 inhibitor. <i>Molecular Oncology</i> , 2020, 14, 3083-3099.	2.1	17
12	<i>CovaDOTS: In Silico</i> Chemistry-Driven Tool to Design Covalent Inhibitors Using a Linking Strategy. <i>Journal of Chemical Information and Modeling</i> , 2019, 59, 1472-1485.	2.5	13
13	Temporin-SHa and Its Analogs as Potential Candidates for the Treatment of <i>Helicobacter pylori</i> . <i>Biomolecules</i> , 2019, 9, 598.	1.8	10
14	Fragment-based drug design targeting syntenin PDZ2 domain involved in exosomal release and tumour spread. <i>European Journal of Medicinal Chemistry</i> , 2021, 223, 113601.	2.6	3
15	CRCM5484: A BET-BDII Selective Compound with Differential Anti-leukemic Drug Modulation. <i>Journal of Medicinal Chemistry</i> , 2022, 65, 5660-5674.	2.9	2
16	Discovery of Small-Molecule Inhibitors of the PTK7 ^β -Catenin Interaction Targeting the Wnt Signaling Pathway in Colorectal Cancer. <i>ACS Chemical Biology</i> , 2022, 17, 1061-1072.	1.6	1