

Andreas Lange

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/9034439/publications.pdf>

Version: 2024-02-01

16
papers

1,561
citations

623188

14
h-index

940134

16
g-index

26
all docs

26
docs citations

26
times ranked

2587
citing authors

#	ARTICLE	IF	CITATIONS
1	Heterologous expression of naturally evolved putative <i>de novo</i> proteins with chaperones. <i>Protein Science</i> , 2022, 31, .	3.1	8
2	Structural and functional characterization of a putative de novo gene in <i>Drosophila</i> . <i>Nature Communications</i> , 2021, 12, 1667.	5.8	40
3	Structure and function of naturally evolved de novo proteins. <i>Current Opinion in Structural Biology</i> , 2021, 68, 175-183.	2.6	49
4	Scaffold Effects on Halogen Bonding Strength. <i>Journal of Chemical Information and Modeling</i> , 2019, 59, 885-894.	2.5	24
5	<p>Small-Molecule Intervention At The Dimerization Interface Of Survivin By Novel Rigidized Scaffolds</p>. <i>Drug Design, Development and Therapy</i> , 2019, Volume 13, 4247-4263.	2.0	9
6	Tri- and Tetrasubstituted Pyridinylimidazoles as Covalent Inhibitors of c-Jun N-Terminal Kinase 3. <i>Journal of Medicinal Chemistry</i> , 2017, 60, 594-607.	2.9	46
7	Using Surface Scans for the Evaluation of Halogen Bonds toward the Side Chains of Aspartate, Asparagine, Glutamate, and Glutamine. <i>Journal of Chemical Information and Modeling</i> , 2016, 56, 1373-1383.	2.5	20
8	Fluorescence polarization-based assays for detecting compounds binding to inactive c-Jun N-terminal kinase 3 and p38 β mitogen-activated protein kinase. <i>Analytical Biochemistry</i> , 2016, 503, 28-40.	1.1	22
9	Evaluating the Potential of Halogen Bonding in Molecular Design: Automated Scaffold Decoration Using the New Scoring Function XBScore. <i>Journal of Chemical Information and Modeling</i> , 2015, 55, 687-699.	2.5	33
10	Targeting the Gatekeeper MET146 of C-Jun N-Terminal Kinase 3 Induces a Bivalent Halogen/Chalcogen Bond. <i>Journal of the American Chemical Society</i> , 2015, 137, 14640-14652.	6.6	73
11	Machine Learning Estimates of Natural Product Conformational Energies. <i>PLoS Computational Biology</i> , 2014, 10, e1003400.	1.5	30
12	Halogen-enriched fragment libraries as chemical probes for harnessing halogen bonding in fragment-based lead discovery. <i>Future Medicinal Chemistry</i> , 2014, 6, 617-639.	1.1	38
13	Targeting Histidine Side Chains in Molecular Design through Nitrogen- π -Halogen Bonds. <i>Journal of Chemical Information and Modeling</i> , 2013, 53, 3178-3189.	2.5	23
14	Principles and Applications of Halogen Bonding in Medicinal Chemistry and Chemical Biology. <i>Journal of Medicinal Chemistry</i> , 2013, 56, 1363-1388.	2.9	1,002
15	Using halogen bonds to address the protein backbone: a systematic evaluation. <i>Journal of Computer-Aided Molecular Design</i> , 2012, 26, 935-945.	1.3	86
16	Addressing Methionine in Molecular Design through Directed Sulfur- π -Halogen Bonds. <i>Journal of Chemical Theory and Computation</i> , 2011, 7, 2307-2315.	2.3	49