## Bernd Rupp

List of Publications by Year in descending order

Source: https://exaly.com/author-pdf/2130363/publications.pdf

Version: 2024-02-01

11	268	1040056	1372567
papers	citations	h-index	g-index
11	11	11	651
all docs	docs citations	times ranked	citing authors

#	Article	IF	CITATIONS
1	Design of chemical libraries with potentially bioactive molecules applying a maximum common substructure concept. Molecular Diversity, 2010, 14, 401-408.	3.9	69
2	Molecular design of two sterol $14\hat{l}_{\pm}$ -demethylase homology models and their interactions with the azole antifungals ketoconazole and bifonazole. Journal of Computer-Aided Molecular Design, 2005, 19, 149-163.	2.9	38
3	Anchor Side Chains of Short Peptide Fragments Trigger Ligand-Exchange of Class II MHC Molecules. PLoS ONE, 2008, 3, e1814.	2.5	34
4	Characterization of Structural Features Controlling the Receptiveness of Empty Class II MHC Molecules. PLoS ONE, 2011, 6, e18662.	2.5	31
5	Design of a Generalâ€Purpose European Compound Screening Library for EUâ€OPENSCREEN. ChemMedChem, 2014, 9, 2309-2326.	3.2	29
6	Chemoenzymatic Synthesis of a Glycolipid Library and Elucidation of the Antigenic Epitope for Construction of a Vaccine Against Lyme Disease. Chemistry - A European Journal, 2010, 16, 3536-3544.	3.3	20
7	Chronic oral LOAEL prediction by using a commercially available computational QSAR tool. Archives of Toxicology, 2010, 84, 681-688.	4.2	17
8	Substrate Transport Activation Is Mediated through Second Periplasmic Loop of Transmembrane Protein MalF in Maltose Transport Complex of Escherichia coli. Journal of Biological Chemistry, 2012, 287, 17040-17049.	3.4	16
9	Structural Characterization and Ligand/Inhibitor Identification Provide Functional Insights into the Mycobacterium tuberculosis Cytochrome P450 CYP126A1. Journal of Biological Chemistry, 2017, 292, 1310-1329.	3.4	13
10	11th German Conference on Chemoinformatics (GCC 2015). Journal of Cheminformatics, 2016, 8, 18.	6.1	1
11	DACS: from compound collections to rationally designed HTS library. Journal of Cheminformatics, 2014, 6, .	6.1	0