

Sebastian Kelm

List of Publications by Year in descending order

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Version: 2024-02-01

29
papers

1,235
citations

471509

17
h-index

552781

26
g-index

34
all docs

34
docs citations

34
times ranked

2044
citing authors

#	ARTICLE	IF	CITATIONS
1	Crystal structure of dopamine D1 receptor in complex with G protein and a non-catechol agonist. Nature Communications, 2021, 12, 3305.	12.8	34
2	The Chemical Synthesis of Knob Domain Antibody Fragments. ACS Chemical Biology, 2021, 16, 1757-1769.	3.4	10
3	Co-evolutionary distance predictions contain flexibility information. Bioinformatics, 2021, , .	4.1	9
4	Computational approaches to therapeutic antibody design: established methods and emerging trends. Briefings in Bioinformatics, 2020, 21, 1549-1567.	6.5	126
5	Structural diversity of B-cell receptor repertoires along the B-cell differentiation axis in humans and mice. PLoS Computational Biology, 2020, 16, e1007636.	3.2	27
6	B-cell epitopes: Discontinuity and conformational analysis. Molecular Immunology, 2019, 114, 643-650.	2.2	28
7	Reconstruction of apo A2A receptor activation pathways reveal ligand-competent intermediates and state-dependent cholesterol hotspots. Scientific Reports, 2019, 9, 14199.	3.3	24
8	The Aminotriazole Antagonist Cmpdâ€1 Stabilises a Distinct Inactive State of the Adenosine 2A Receptor. Angewandte Chemie, 2019, 131, 9499-9503.	2.0	1
9	The Aminotriazole Antagonist Cmpdâ€1 Stabilises a Distinct Inactive State of the Adenosine 2A Receptor. Angewandte Chemie - International Edition, 2019, 58, 9399-9403.	13.8	3
10	Electrostatic interactions modulate the differential aggregation propensities of IgG1 and IgG4P antibodies and inform charged residue substitutions for improved developability. Protein Engineering, Design and Selection, 2019, 32, 277-288.	2.1	19
11	SCALOP: sequence-based antibody canonical loop structure annotation. Bioinformatics, 2019, 35, 1774-1776.	4.1	29
12	Filtering Next-Generation Sequencing of the Ig Gene Repertoire Data Using Antibody Structural Information. Journal of Immunology, 2018, 201, 3694-3704.	0.8	11
13	Observed Antibody Space: A Resource for Data Mining Next-Generation Sequencing of Antibody Repertoires. Journal of Immunology, 2018, 201, 2502-2509.	0.8	165
14	Insight into small molecule binding to the neonatal Fc receptor by X-ray crystallography and 100 kHz magic-angle-spinning NMR. PLoS Biology, 2018, 16, e2006192.	5.6	31
15	Structurally Mapping Antibody Repertoires. Frontiers in Immunology, 2018, 9, 1698.	4.8	36
16	Efficient Sampling for the Prediction of Long and Multidomain Protein Structures. Biophysical Journal, 2018, 114, 574a.	0.5	0
17	A multi-crystal method for extracting obscured crystallographic states from conventionally uninterpretable electron density. Nature Communications, 2017, 8, 15123.	12.8	186
18	Investigating Cotranslational Folding in Membrane Proteins using Fragment-Based Structure Prediction. Biophysical Journal, 2017, 112, 61a.	0.5	1

#	ARTICLE	IF	CITATIONS
19	Sphinx: merging knowledge-based and <i>ab initio</i> approaches to improve protein loop prediction. <i>Bioinformatics</i> , 2017, 33, 1346-1353.	4.1	49
20	Examining the Conservation of Kinks in Alpha Helices. <i>PLoS ONE</i> , 2016, 11, e0157553.	2.5	20
21	SAbPred: a structure-based antibody prediction server. <i>Nucleic Acids Research</i> , 2016, 44, W474-W478.	14.5	155
22	Length-independent structural similarities enrich the antibody CDR canonical class model. <i>MAbs</i> , 2016, 8, 751-760.	5.2	49
23	PANDDAs: multi-dataset methods for finding hits from fragment screening by X-ray crystallography. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2015, 71, s258-s258.	0.1	2
24	Protein Modeling and Structural Prediction. , 2014, , 171-182.		1
25	Fragment-based modeling of membrane protein loops: Successes, failures, and prospects for the future. <i>Proteins: Structure, Function and Bioinformatics</i> , 2014, 82, 175-186.	2.6	15
26	Memoir: template-based structure prediction for membrane proteins. <i>Nucleic Acids Research</i> , 2013, 41, W379-W383.	14.5	38
27	Environment specific substitution tables improve membrane protein alignment. <i>Bioinformatics</i> , 2011, 27, i15-i23.	4.1	20
28	MEDELLER: homology-based coordinate generation for membrane proteins. <i>Bioinformatics</i> , 2010, 26, 2833-2840.	4.1	103
29	iMembrane: homology-based membrane-insertion of proteins. <i>Bioinformatics</i> , 2009, 25, 1086-1088.	4.1	29