

Xiaodong Pang

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/471389/publications.pdf>

Version: 2024-02-01

34
papers

1,487
citations

567281

15
h-index

552781

26
g-index

34
all docs

34
docs citations

34
times ranked

2598
citing authors

#	ARTICLE	IF	CITATIONS
1	Genetic factors associated with cancer racial disparity – an integrative study across twenty-one cancer types. <i>Molecular Oncology</i> , 2020, 14, 2775-2786.	4.6	8
2	Maximizing the reusability of gene expression data by predicting missing metadata. <i>PLoS Computational Biology</i> , 2020, 16, e1007450.	3.2	4
3	Maximizing the reusability of gene expression data by predicting missing metadata. , 2020, 16, e1007450.		0
4	Maximizing the reusability of gene expression data by predicting missing metadata. , 2020, 16, e1007450.		0
5	Maximizing the reusability of gene expression data by predicting missing metadata. , 2020, 16, e1007450.		0
6	Maximizing the reusability of gene expression data by predicting missing metadata. , 2020, 16, e1007450.		0
7	Maximizing the reusability of gene expression data by predicting missing metadata. , 2020, 16, e1007450.		0
8	Maximizing the reusability of gene expression data by predicting missing metadata. , 2020, 16, e1007450.		0
9	The novel BET–CBP/p300 dual inhibitor NEO2734 is active in SPOP mutant and wild-type prostate cancer. <i>EMBO Molecular Medicine</i> , 2019, 11, e10659.	6.9	56
10	Influenza A M2 Channel Clustering at High Protein/Lipid Ratios: Viral Budding Implications. <i>Biophysical Journal</i> , 2019, 116, 1075-1084.	0.5	33
11	Electrostatic Interactions in Protein Structure, Folding, Binding, and Condensation. <i>Chemical Reviews</i> , 2018, 118, 1691-1741.	47.7	577
12	Rate Constants and Mechanisms of Protein–Ligand Binding. <i>Annual Review of Biophysics</i> , 2017, 46, 105-130.	10.0	49
13	The dock-and-coalesce mechanism for the association of a <sc>WASP</sc> disordered region with the Cdc42 <sc>GTPase</sc>. <i>FEBS Journal</i> , 2017, 284, 3381-3391.	4.7	23
14	Structural modeling for the open state of an NMDA receptor. <i>Journal of Structural Biology</i> , 2017, 200, 369-375.	2.8	13
15	Mechanism and rate constants of the Cdc42 GTPase binding with intrinsically disordered effectors. <i>Proteins: Structure, Function and Bioinformatics</i> , 2016, 84, 674-685.	2.6	17
16	Computational full electron structure study of biological activity in Cyclophilin A. <i>Journal of Biomolecular Structure and Dynamics</i> , 2016, 34, 870-876.	3.5	0
17	Two Pathways Mediate Interdomain Allosteric Regulation in Pin1. <i>Structure</i> , 2015, 23, 237-247.	3.3	70
18	The full electron structure of the FKBP12/FK506 complex. <i>Journal of Biomolecular Structure and Dynamics</i> , 2015, 33, 388-394.	3.5	3

#	ARTICLE	IF	CITATIONS
19	Disorder-to-Order Transition of an Active-Site Loop Mediates the Allosteric Activation of Sortase A. <i>Biophysical Journal</i> , 2015, 109, 1706-1715.	0.5	26
20	Truncated ERG Oncoproteins from TMPRSS2-ERG Fusions Are Resistant to SPOP-Mediated Proteasome Degradation. <i>Molecular Cell</i> , 2015, 59, 904-916.	9.7	129
21	Distinct mechanisms of a phosphotyrosyl peptide binding to two SH2 domains. <i>Journal of Theoretical and Computational Chemistry</i> , 2014, 13, 1440003.	1.8	4
22	Design Rules for Selective Binding of Nuclear Localization Signals to Minor Site of Importin β . <i>PLoS ONE</i> , 2014, 9, e91025.	2.5	13
23	Molecular Structure of RADA16-I Designer Self-Assembling Peptide Nanofibers. <i>ACS Nano</i> , 2013, 7, 7562-7572.	14.6	124
24	Activation of signaling receptors: do ligands bind to receptor monomer, dimer, or both?. <i>BMC Biophysics</i> , 2013, 6, 7.	4.4	7
25	Solid-State NMR Evidence for β^2 -Hairpin Structure within MAX8 Designer Peptide Nanofibers. <i>Biophysical Journal</i> , 2013, 105, 222-230.	0.5	24
26	Poisson-Boltzmann Calculations: van der Waals or Molecular Surface?. <i>Communications in Computational Physics</i> , 2013, 13, 1-12.	1.7	41
27	A Common Model for Cytokine Receptor Activation: Combined Scissor-Like Rotation and Self-Rotation of Receptor Dimer Induced by Class I Cytokine. <i>PLoS Computational Biology</i> , 2012, 8, e1002427.	3.2	24
28	Prediction and Dissection of Widely-Varying Association Rate Constants of Actin-Binding Proteins. <i>PLoS Computational Biology</i> , 2012, 8, e1002696.	3.2	13
29	Rate constants and mechanisms of intrinsically disordered proteins binding to structured targets. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 10466.	2.8	75
30	Rationalizing 5000-Fold Differences in Receptor-Binding Rate Constants of Four Cytokines. <i>Biophysical Journal</i> , 2011, 101, 1175-1183.	0.5	20
31	Automated Prediction of Protein Association Rate Constants. <i>Structure</i> , 2011, 19, 1744-1751.	3.3	111
32	Discovery of a potent peptidic cyclophilin A inhibitor Trp-Gly-Pro. <i>European Journal of Medicinal Chemistry</i> , 2011, 46, 1701-1705.	5.5	10
33	A Mathematical Model for Peptide Inhibitor Design. <i>Journal of Computational Biology</i> , 2010, 17, 1081-1093.	1.6	4
34	Two Rules on the Protein-Ligand Interaction. <i>Nature Precedings</i> , 0, , .	0.1	9