

Ahmet Bakan

List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

18
papers

1,972
citations

15
h-index

18
g-index

18
ext. papers

2,329
ext. citations

9.6
avg, IF

4.85
L-index

#	Paper	IF	Citations
18	Allostery as Structure-Encoded Collective Dynamics 2020 , 125-141		1
17	Inhibition of Peroxidase Activity of Cytochrome c: De Novo Compound Discovery and Validation. <i>Molecular Pharmacology</i> , 2015 , 88, 421-7	4.3	19
16	Synthesis and biological evaluation of 3-aminoisoquinolin-1(2H)-one based inhibitors of the dual-specificity phosphatase Cdc25B. <i>Bioorganic and Medicinal Chemistry</i> , 2015 , 23, 2810-8	3.4	28
15	Designing inhibitors of cytochrome c/cardiolipin peroxidase complexes: mitochondria-targeted imidazole-substituted fatty acids. <i>Free Radical Biology and Medicine</i> , 2014 , 71, 221-230	7.8	33
14	Evo and ProDy for bridging protein sequence evolution and structural dynamics. <i>Bioinformatics</i> , 2014 , 30, 2681-3	7.2	130
13	Investigational inhibitors of PTP4A3 phosphatase as antineoplastic agents. <i>Expert Opinion on Investigational Drugs</i> , 2014 , 23, 661-73	5.9	15
12	In vivo structure-activity relationship studies support allosteric targeting of a dual specificity phosphatase. <i>ChemBioChem</i> , 2014 , 15, 1436-45	3.8	34
11	Druggability Assessment of Allosteric Proteins by Dynamics Simulations in the Presence of Probe Molecules. <i>Journal of Chemical Theory and Computation</i> , 2012 , 8, 2435-2447	6.4	113
10	Coupling between catalytic loop motions and enzyme global dynamics. <i>PLoS Computational Biology</i> , 2012 , 8, e1002705	5	35
9	ProDy: protein dynamics inferred from theory and experiments. <i>Bioinformatics</i> , 2011 , 27, 1575-7	7.2	626
8	Pre-existing soft modes of motion uniquely defined by native contact topology facilitate ligand binding to proteins. <i>Protein Science</i> , 2011 , 20, 1645-58	6.3	74
7	Computational generation inhibitor-bound conformers of p38 MAP kinase and comparison with experiments. <i>Pacific Symposium on Biocomputing Pacific Symposium on Biocomputing</i> , 2011 , 181-92	1.3	17
6	Sodium-coupled Secondary Transporters: Insights from Structure-based Computations 2011 , 199-229		2
5	Normal mode analysis of biomolecular structures: functional mechanisms of membrane proteins. <i>Chemical Reviews</i> , 2010 , 110, 1463-97	68.1	383
4	The intrinsic dynamics of enzymes plays a dominant role in determining the structural changes induced upon inhibitor binding. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2009 , 106, 14349-54	11.5	219
3	Zebrafish chemical screening reveals an inhibitor of Dusp6 that expands cardiac cell lineages. <i>Nature Chemical Biology</i> , 2009 , 5, 680-7	11.7	190
2	Toward a molecular understanding of the interaction of dual specificity phosphatases with substrates: insights from structure-based modeling and high throughput screening. <i>Current Medicinal Chemistry</i> , 2008 , 15, 2536-44	4.3	31

- 1 Structurally unique inhibitors of human mitogen-activated protein kinase phosphatase-1 identified in a pyrrole carboxamide library. *Journal of Pharmacology and Experimental Therapeutics*, **2007**, 322, 940-7 22