

# Ahmet Bakan

## List of Publications by Citations

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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	ProDy: protein dynamics inferred from theory and experiments. <i>Bioinformatics</i> , <b>2011</b> , 27, 1575-7	7.2	626
17	Normal mode analysis of biomolecular structures: functional mechanisms of membrane proteins. <i>Chemical Reviews</i> , <b>2010</b> , 110, 1463-97	68.1	383
16	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> , <b>2009</b> , 106, 14349-54	11.5	219
15	Zebrafish chemical screening reveals an inhibitor of Dusp6 that expands cardiac cell lineages. <i>Nature Chemical Biology</i> , <b>2009</b> , 5, 680-7	11.7	190
14	Evol and ProDy for bridging protein sequence evolution and structural dynamics. <i>Bioinformatics</i> , <b>2014</b> , 30, 2681-3	7.2	130
13	Druggability Assessment of Allosteric Proteins by Dynamics Simulations in the Presence of Probe Molecules. <i>Journal of Chemical Theory and Computation</i> , <b>2012</b> , 8, 2435-2447	6.4	113
12	Pre-existing soft modes of motion uniquely defined by native contact topology facilitate ligand binding to proteins. <i>Protein Science</i> , <b>2011</b> , 20, 1645-58	6.3	74
11	Coupling between catalytic loop motions and enzyme global dynamics. <i>PLoS Computational Biology</i> , <b>2012</b> , 8, e1002705	5	35
10	In vivo structure-activity relationship studies support allosteric targeting of a dual specificity phosphatase. <i>ChemBioChem</i> , <b>2014</b> , 15, 1436-45	3.8	34
9	Designing inhibitors of cytochrome c/cardiolipin peroxidase complexes: mitochondria-targeted imidazole-substituted fatty acids. <i>Free Radical Biology and Medicine</i> , <b>2014</b> , 71, 221-230	7.8	33
8	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> , <b>2008</b> , 15, 2536-44	4.3	31
7	Synthesis and biological evaluation of 3-aminoisoquinolin-1(2H)-one based inhibitors of the dual-specificity phosphatase Cdc25B. <i>Bioorganic and Medicinal Chemistry</i> , <b>2015</b> , 23, 2810-8	3.4	28
6	Structurally unique inhibitors of human mitogen-activated protein kinase phosphatase-1 identified in a pyrrole carboxamide library. <i>Journal of Pharmacology and Experimental Therapeutics</i> , <b>2007</b> , 322, 940-7	4.7	22
5	Inhibition of Peroxidase Activity of Cytochrome c: De Novo Compound Discovery and Validation. <i>Molecular Pharmacology</i> , <b>2015</b> , 88, 421-7	4.3	19
4	Computational generation inhibitor-bound conformers of p38 MAP kinase and comparison with experiments. <i>Pacific Symposium on Biocomputing Pacific Symposium on Biocomputing</i> , <b>2011</b> , 181-92	1.3	17
3	Investigational inhibitors of PTP4A3 phosphatase as antineoplastic agents. <i>Expert Opinion on Investigational Drugs</i> , <b>2014</b> , 23, 661-73	5.9	15
2	Sodium-coupled Secondary Transporters: Insights from Structure-based Computations <b>2011</b> , 199-229		2

1 Allosteric as Structure-Encoded Collective Dynamics **2020**, 125-141

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