

George Hedger

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/12160995/publications.pdf>

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12
papers

1,172
citations

759055

12
h-index

1199470

12
g-index

14
all docs

14
docs citations

14
times ranked

1779
citing authors

#	ARTICLE	IF	CITATIONS
1	Structural basis of Smoothed regulation by its extracellular domains. <i>Nature</i> , 2016, 535, 517-522.	13.7	300
2	PtdIns(4,5)P2 stabilizes active states of GPCRs and enhances selectivity of G-protein coupling. <i>Nature</i> , 2018, 559, 423-427.	13.7	236
3	Lipid interaction sites on channels, transporters and receptors: Recent insights from molecular dynamics simulations. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2016, 1858, 2390-2400.	1.4	143
4	Convergence and Sampling in Determining Free Energy Landscapes for Membrane Protein Association. <i>Journal of Physical Chemistry B</i> , 2017, 121, 3364-3375.	1.2	93
5	The juxtamembrane regions of human receptor tyrosine kinases exhibit conserved interaction sites with anionic lipids. <i>Scientific Reports</i> , 2015, 5, 9198.	1.6	89
6	Cholesterol Interaction Sites on the Transmembrane Domain of the Hedgehog Signal Transducer and Class F G Protein-Coupled Receptor Smoothed. <i>Structure</i> , 2019, 27, 549-559.e2.	1.6	77
7	Lipid-Loving ANTs: Molecular Simulations of Cardiolipin Interactions and the Organization of the Adenine Nucleotide Translocase in Model Mitochondrial Membranes. <i>Biochemistry</i> , 2016, 55, 6238-6249.	1.2	63
8	Free Energy Landscape of Lipid Interactions with Regulatory Binding Sites on the Transmembrane Domain of the EGF Receptor. <i>Journal of Physical Chemistry B</i> , 2016, 120, 8154-8163.	1.2	60
9	Lipid Interactions of a Ciliary Membrane TRP Channel: Simulation and Structural Studies of Polycystin-2. <i>Structure</i> , 2020, 28, 169-184.e5.	1.6	37
10	Interactions of the EphA2 Kinase Domain with PIPs in Membranes: Implications for Receptor Function. <i>Structure</i> , 2018, 26, 1025-1034.e2.	1.6	33
11	Molecular simulations of glycolipids: Towards mammalian cell membrane models. <i>Biochimie</i> , 2016, 120, 105-109.	1.3	27
12	GPCRs: What Can We Learn from Molecular Dynamics Simulations?. <i>Methods in Molecular Biology</i> , 2018, 1705, 133-158.	0.4	13