PaweÅ, SledÅº

List of Publications by Year
in descending order

14	1,205	933447 10	1058476 14
papers	citations	h-index	g-index

17	17	17	1917
all docs	docs citations	times ranked	citing authors

#	Article	IF	Citations
1	Protein structure-based drug design: from docking to molecular dynamics. Current Opinion in Structural Biology, 2018, 48, 93-102.	5.7	405
2	Structural insights into the molecular mechanism of the m6A writer complex. ELife, 2016, 5, .	6.0	386
3	Smallâ€Molecule Inhibitors of METTL3, the Major Human Epitranscriptomic Writer. ChemMedChem, 2020, 15, 744-748.	3.2	106
4	Using Ligandâ€Mapping Simulations to Design a Ligand Selectively Targeting a Cryptic Surface Pocket of Poloâ€Like Kinase 1. Angewandte Chemie - International Edition, 2012, 51, 10078-10081.	13.8	71
5	From Crystal Packing to Molecular Recognition: Prediction and Discovery of a Binding Site on the Surface of Poloâ€Like Kinase 1. Angewandte Chemie - International Edition, 2011, 50, 4003-4006.	13.8	57
6	Molecular basis for cytoplasmic <scp>RNA</scp> surveillance by uridylationâ€triggered decay in Drosophila. EMBO Journal, 2016, 35, 2417-2434.	7.8	50
7	A Reader-Based Assay for m⁶A Writers and Erasers. Analytical Chemistry, 2019, 91, 3078-3084.	6.5	36
8	Highâ€Throughput Interrogation of Ligand Binding Mode Using a Fluorescenceâ€Based Assay. Angewandte Chemie - International Edition, 2012, 51, 7680-7683.	13.8	30
9	Flexible Binding of m⁶A Reader Protein YTHDC1 to Its Preferred RNA Motif. Journal of Chemical Theory and Computation, 2019, 15, 7004-7014.	5.3	18
10	Selectively Disrupting m⁶A-Dependent Proteinâ€“RNA Interactions with Fragments. ACS Chemical Biology, 2020, 15, 618-625.	3.4	17
11	Binding Motifs in the CBP Bromodomain: An Analysis of 20 Crystal Structures of Complexes with Small Molecules. ACS Medicinal Chemistry Letters, 2018, 9, 929-934.	2.8	8
12	Hitting a Moving Target: Simulation and Crystallography Study of ATAD2 Bromodomain Blockers. ACS Medicinal Chemistry Letters, 2020, 11, 1573-1580.	2.8	8
13	Understanding the mechanism of action of pyrrolo[3,2-b]quinoxaline-derivatives as kinase inhibitors. RSC Medicinal Chemistry, 2020, 11, 665-675.	3.9	4
14	Ligand retargeting by binding site analogy. European Journal of Medicinal Chemistry, 2019, 175, 107-113.	5.5	2