

Bradley C Doak

List of Publications by Year in descending order

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Version: 2024-02-01

25
papers

1,761
citations

623734

14
h-index

610901

24
g-index

30
all docs

30
docs citations

30
times ranked

2195
citing authors

#	ARTICLE	IF	CITATIONS
1	Selective Binding of Small Molecules to <i>Vibrio cholerae</i> DsbA Offers a Starting Point for the Design of Novel Antibacterials. <i>ChemMedChem</i> , 2022, 17, .	3.2	3
2	Methyl probes in proteins for determining ligand binding mode in weak protein-ligand complexes. <i>Scientific Reports</i> , 2022, 12, .	3.3	3
3	Drug Syntheses Beyond the Rule of 5. <i>Chemistry - A European Journal</i> , 2020, 26, 49-88.	3.3	36
4	Frontispiece: Drug Syntheses Beyond the Rule of 5. <i>Chemistry - A European Journal</i> , 2020, 26, .	3.3	1
5	NMR fragment screening reveals a novel small molecule binding site near the catalytic surface of the disulfide dithiol oxidoreductase enzyme DsbA from <i>Burkholderia pseudomallei</i> . <i>Journal of Biomolecular NMR</i> , 2020, 74, 595-611.	2.8	7
6	Rapid Elaboration of Fragments into Leads by X-ray Crystallographic Screening of Parallel Chemical Libraries (REFIL _X). <i>Journal of Medicinal Chemistry</i> , 2020, 63, 6863-6875.	6.4	16
7	A ligand-induced structural change in fatty acid-binding protein 1 is associated with potentiation of peroxisome proliferator-activated receptor β agonists. <i>Journal of Biological Chemistry</i> , 2019, 294, 3720-3734.	3.4	17
8	Impact of Dynamically Exposed Polarity on Permeability and Solubility of Chameleonic Drugs Beyond the Rule of 5. <i>Journal of Medicinal Chemistry</i> , 2018, 61, 4189-4202.	6.4	150
9	Design of a Fragment-Screening Library. <i>Methods in Enzymology</i> , 2018, 610, 97-115.	1.0	7
10	Cyclophilin Succumbs to a Macrocyclic Chameleon. <i>Journal of Medicinal Chemistry</i> , 2018, 61, 9469-9472.	6.4	3
11	Structural and biochemical insights into the disulfide reductase mechanism of DsbD, an essential enzyme for neisserial pathogens. <i>Journal of Biological Chemistry</i> , 2018, 293, 16559-16571.	3.4	10
12	Opportunities and guidelines for discovery of orally absorbed drugs in beyond rule of 5 space. <i>Current Opinion in Chemical Biology</i> , 2018, 44, 23-29.	6.1	64
13	Applications of NMR Spectroscopy in FBDD. , 2018, , 2211-2231.		2
14	Drug discovery beyond the rule of 5 - Opportunities and challenges. <i>Expert Opinion on Drug Discovery</i> , 2017, 12, 115-119.	5.0	75
15	Applications of NMR Spectroscopy in FBDD. , 2017, , 1-22.		0
16	Cell permeability beyond the rule of 5. <i>Advanced Drug Delivery Reviews</i> , 2016, 101, 42-61.	13.7	196
17	The ways and means of fragment-based drug design. , 2016, 167, 28-37.		67
18	Structural and conformational determinants of macrocycle cell permeability. <i>Nature Chemical Biology</i> , 2016, 12, 1065-1074.	8.0	152

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19	Determination of ligand binding modes in weak protein–ligand complexes using sparse NMR data. <i>Journal of Biomolecular NMR</i> , 2016, 66, 195-208.	2.8	19
20	How Beyond Rule of 5 Drugs and Clinical Candidates Bind to Their Targets. <i>Journal of Medicinal Chemistry</i> , 2016, 59, 2312-2327.	6.4	248
21	Application of Fragment-Based Screening to the Design of Inhibitors of <i>Escherichia coli</i> DsbA. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 2179-2184.	13.8	46
22	Oral Druggable Space beyond the Rule of 5: Insights from Drugs and Clinical Candidates. <i>Chemistry and Biology</i> , 2014, 21, 1115-1142.	6.0	523
23	Design and Evaluation of the Performance of an NMR Screening Fragment Library. <i>Australian Journal of Chemistry</i> , 2013, 66, 1465.	0.9	33
24	Synthesis of Unsymmetrical 1,1-Di-Disubstituted Bis(1,2,3-triazole)s Using Monosilylbutadiynes. <i>Organic Letters</i> , 2011, 13, 537-539.	4.6	43
25	Structure and Function of the Oxidoreductase DsbA1 from <i>Neisseria meningitidis</i> . <i>Journal of Molecular Biology</i> , 2009, 394, 931-943.	4.2	36