

Christopher W Murray

List of Publications by Year in descending order

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14
papers

1,419
citations

759233

12
h-index

1058476

14
g-index

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all docs

14
docs citations

14
times ranked

1917
citing authors

#	ARTICLE	IF	CITATIONS
1	Discovery of (2,4-Dihydroxy-5-isopropylphenyl)-[5-(4-methylpiperazin-1-ylmethyl)-1,3-dihydroisoindol-2-yl]methanone (AT13387), a Novel Inhibitor of the Molecular Chaperone Hsp90 by Fragment Based Drug Design. <i>Journal of Medicinal Chemistry</i> , 2010, 53, 5956-5969.	6.4	230
2	Experiences in fragment-based drug discovery. <i>Trends in Pharmacological Sciences</i> , 2012, 33, 224-232.	8.7	227
3	Design Principles for Fragment Libraries: Maximizing the Value of Learnings from Pharma Fragment-Based Drug Discovery (FBDD) Programs for Use in Academia. <i>Journal of Medicinal Chemistry</i> , 2016, 59, 8189-8206.	6.4	182
4	Fragment-Based Drug Discovery Applied to Hsp90. Discovery of Two Lead Series with High Ligand Efficiency. <i>Journal of Medicinal Chemistry</i> , 2010, 53, 5942-5955.	6.4	173
5	Opportunity Knocks: Organic Chemistry for Fragment-Based Drug Discovery (FBDD). <i>Angewandte Chemie - International Edition</i> , 2016, 55, 488-492.	13.8	151
6	Efficient exploration of chemical space by fragment-based screening. <i>Progress in Biophysics and Molecular Biology</i> , 2014, 116, 82-91.	2.9	127
7	Discovery of an allosteric mechanism for the regulation of HCV NS3 protein function. <i>Nature Chemical Biology</i> , 2012, 8, 920-925.	8.0	103
8	Fragment-to-Lead Medicinal Chemistry Publications in 2015. <i>Journal of Medicinal Chemistry</i> , 2017, 60, 89-99.	6.4	47
9	A successful collaboration between academia, biotech and pharma led to discovery of erdafitinib, a selective FGFR inhibitor recently approved by the FDA. <i>MedChemComm</i> , 2019, 10, 1509-1511.	3.4	47
10	Fragment-Based Discovery of a Potent, Orally Bioavailable Inhibitor That Modulates the Phosphorylation and Catalytic Activity of ERK1/2. <i>Journal of Medicinal Chemistry</i> , 2018, 61, 4978-4992.	6.4	42
11	Fragment-based drug discovery: opportunities for organic synthesis. <i>RSC Medicinal Chemistry</i> , 2021, 12, 321-329.	3.9	35
12	The Fragment Network: A Chemistry Recommendation Engine Built Using a Graph Database. <i>Journal of Medicinal Chemistry</i> , 2017, 60, 6440-6450.	6.4	33
13	ASTX029, a Novel Dual-mechanism ERK Inhibitor, Modulates Both the Phosphorylation and Catalytic Activity of ERK. <i>Molecular Cancer Therapeutics</i> , 2021, 20, 1757-1768.	4.1	13
14	Discovery of ASTX029, A Clinical Candidate Which Modulates the Phosphorylation and Catalytic Activity of ERK1/2. <i>Journal of Medicinal Chemistry</i> , 2021, 64, 12286-12303.	6.4	9