

Mushtaque S Shaikh

List of Publications by Citations

Source: <https://exaly.com/author-pdf/8784062/mushtaque-s-shaikh-publications-by-citations.pdf>

Version: 2024-04-27

This document has been generated based on the publications and citations recorded by exaly.com. For the latest version of this publication list, visit the link given above.

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

367
citations

12
h-index

18
g-index

18
ext. papers

388
ext. citations

3.1
avg, IF

3.05
L-index

#	Paper	IF	Citations
18	3,5-Diformylboron dipyrromethenes as fluorescent pH sensors. <i>Inorganic Chemistry</i> , 2011 , 50, 4392-400	5.1	84
17	Synthesis and studies of covalently linked meso-furyl boron-dipyrromethene-ferrocene conjugates. <i>Journal of Organometallic Chemistry</i> , 2012 , 697, 65-73	2.3	32
16	3D-QSAR studies of Dipeptidyl peptidase IV inhibitors using a docking based alignment. <i>Journal of Molecular Modeling</i> , 2007 , 13, 1047-71	2	30
15	Screening for in vitro antimycobacterial activity and three-dimensional quantitative structure-activity relationship (3D-QSAR) study of 4-(arylamino)coumarin derivatives. <i>Chemical Biology and Drug Design</i> , 2010 , 76, 412-24	2.9	28
14	Synthesis and electronic properties of meso-furyl boron dipyrromethenes. <i>Inorganica Chimica Acta</i> , 2012 , 383, 257-266	2.7	27
13	Synthesis, anti-tuberculosis activity and 3D-QSAR study of amino acid conjugates of 4-(adamantan-1-yl) group containing quinolines. <i>European Journal of Medicinal Chemistry</i> , 2009 , 44, 2017-29	6.8	22
12	Synthesis and photophysical properties of covalently linked boron dipyrromethene dyads. <i>Dyes and Pigments</i> , 2012 , 94, 66-73	4.6	19
11	Molecular docking and 3D-QSAR studies of HIV-1 protease inhibitors. <i>Journal of Molecular Modeling</i> , 2010 , 16, 1251-68	2	17
10	Design, synthesis, and evaluation of 4-(substituted)phenyl-2-thioxo-3,4-dihydro-1H-chromino[4,3-d]pyrimidin-5-one and 4-(substituted)phenyl-3,4-dihydro-1H-chromino[4,3-d]pyrimidine-2,5-dione analogs as anti-tubercular agents. <i>Medicinal Chemistry Research</i> , 2014 , 22, 2511-2525	2.2	15
9	Drug-DNA Interaction Studies of Acridone-Based Derivatives. <i>Nucleosides, Nucleotides and Nucleic Acids</i> , 2015 , 34, 309-31	1.4	14
8	Molecular modeling studies, synthesis, and biological evaluation of Plasmodium falciparum enoyl-acyl carrier protein reductase (PFENR) inhibitors. <i>Molecular Diversity</i> , 2009 , 13, 501-17	3.1	14
7	Synthesis and properties of meso-unsubstituted 3-pyrrolyl boron dipyrromethene. <i>Journal of Fluorescence</i> , 2013 , 23, 519-25	2.4	12
6	Molecular modeling studies, synthesis and biological evaluation of derivatives of N-phenylbenzamide as Plasmodium falciparum dihydroorotate dehydrogenase (PFDHODH) inhibitors. <i>Medicinal Chemistry Research</i> , 2011 , 20, 321-332	2.2	12
5	Synthesis of 1-(2,6-dichlorophenyl)-3-methylene-1,3-dihydro-indol-2-one derivatives and in vitro anticancer evaluation against SW620 colon cancer cell line. <i>European Journal of Medicinal Chemistry</i> , 2009 , 44, 1355-62	6.8	11
4	Quantifying ligand-receptor interactions for gorge-spanning acetylcholinesterase inhibitors for the treatment of Alzheimer's disease. <i>Journal of Biomolecular Structure and Dynamics</i> , 2015 , 33, 1107-25	3.6	10
3	Molecular Mechanics Force Fields and their Applications in Drug Design. <i>Anti-Infective Agents in Medicinal Chemistry</i> , 2009 , 8, 128-150		10
2	Self nanoemulsifying granules (SNEGs) of meloxicam: preparation, characterization, molecular modeling and evaluation of in vivo anti-inflammatory activity. <i>Drug Development and Industrial Pharmacy</i> , 2017 , 43, 600-610	3.6	7

- 1 De novo design of 7-aminocoumarin derivatives as novel falcipain-3 inhibitors. *Journal of Molecular Modeling*, **2012**, 18, 1481-93 2 3