

Mushtaque S Shaikh

List of Publications by Year in Descending Order

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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	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
17	Drug-DNA Interaction Studies of Acridone-Based Derivatives. <i>Nucleosides, Nucleotides and Nucleic Acids</i> , 2015 , 34, 309-31	1.4	14
16	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
15	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 antitubercular agents. <i>Medicinal Chemistry Research</i> , 2014 , 23, 2564-2575	2.2	15
14	Synthesis and properties of meso-unsubstituted 3-pyrrolyl boron dipyrromethene. <i>Journal of Fluorescence</i> , 2013 , 23, 519-25	2.4	12
13	Synthesis and photophysical properties of covalently linked boron dipyrromethene dyads. <i>Dyes and Pigments</i> , 2012 , 94, 66-73	4.6	19
12	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
11	De novo design of 7-aminocoumarin derivatives as novel falcipain-3 inhibitors. <i>Journal of Molecular Modeling</i> , 2012 , 18, 1481-93	2	3
10	Synthesis and electronic properties of meso-furyl boron dipyrromethenes. <i>Inorganica Chimica Acta</i> , 2012 , 383, 257-266	2.7	27
9	3,5-Diformylboron dipyrromethenes as fluorescent pH sensors. <i>Inorganic Chemistry</i> , 2011 , 50, 4392-400	5.1	84
8	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
7	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
6	Molecular docking and 3D-QSAR studies of HIV-1 protease inhibitors. <i>Journal of Molecular Modeling</i> , 2010 , 16, 1251-68	2	17
5	Molecular Mechanics Force Fields and their Applications in Drug Design. <i>Anti-Infective Agents in Medicinal Chemistry</i> , 2009 , 8, 128-150		10
4	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
3	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
2	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

1 3D-QSAR studies of Dipeptidyl peptidase IV inhibitors using a docking based alignment. *Journal of Molecular Modeling*, **2007**, 13, 1047-71 2 30