

Mymoona Akhter

List of Publications by Citations

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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.

71
papers

1,913
citations

19
h-index

42
g-index

82
ext. papers

2,356
ext. citations

3.7
avg, IF

4.84
L-index

#	Paper	IF	Citations
71	Quinoline: A versatile heterocyclic. <i>Saudi Pharmaceutical Journal</i> , 2013 , 21, 1-12	4.4	248
70	The therapeutic voyage of pyrazole and its analogs: A review. <i>European Journal of Medicinal Chemistry</i> , 2016 , 120, 170-201	6.8	235
69	A review exploring biological activities of hydrazones. <i>Journal of Pharmacy and Bioallied Sciences</i> , 2014 , 6, 69-80	1.1	174
68	Therapeutic evolution of benzimidazole derivatives in the last quinquennial period. <i>European Journal of Medicinal Chemistry</i> , 2017 , 126, 705-753	6.8	119
67	Piperazine scaffold: A remarkable tool in generation of diverse pharmacological agents. <i>European Journal of Medicinal Chemistry</i> , 2015 , 102, 487-529	6.8	103
66	Pyrazolines: a biological review. <i>Mini-Reviews in Medicinal Chemistry</i> , 2013 , 13, 921-31	3.2	100
65	Arroylpropionic acid based 2,5-disubstituted-1,3,4-oxadiazoles: synthesis and their anti-inflammatory and analgesic activities. <i>European Journal of Medicinal Chemistry</i> , 2009 , 44, 2372-8	6.8	94
64	Green recipes to quinoline: A review. <i>European Journal of Medicinal Chemistry</i> , 2019 , 164, 121-170	6.8	85
63	Synthesis and evaluation of anticancer activity of some novel 6-aryl-2-(p-sulfamylphenyl)-pyridazin-3(2H)-ones. <i>European Journal of Medicinal Chemistry</i> , 2012 , 49, 304-9 ^{6,8}	6.8	57
62	The therapeutic journey of pyridazinone. <i>European Journal of Medicinal Chemistry</i> , 2016 , 123, 256-281	6.8	45
61	Revealing quinquennial anticancer journey of morpholine: A SAR based review. <i>European Journal of Medicinal Chemistry</i> , 2019 , 167, 324-356	6.8	44
60	Synthesis, pharmacological activity and hydrolytic behavior of glyceride prodrugs of ibuprofen. <i>European Journal of Medicinal Chemistry</i> , 2005 , 40, 371-6	6.8	39
59	Pyrazole-pyrazoline as promising novel antimalarial agents: A mechanistic study. <i>European Journal of Medicinal Chemistry</i> , 2018 , 149, 139-147	6.8	37
58	A Review Exploring Therapeutic Worth of 1,3,4-Oxadiazole Tailored Compounds. <i>Mini-Reviews in Medicinal Chemistry</i> , 2019 , 19, 477-509	3.2	32
57	Synthesis of pyrazole acrylic acid based oxadiazole and amide derivatives as antimalarial and anticancer agents. <i>Bioorganic Chemistry</i> , 2018 , 77, 106-124	5.1	26
56	Synthesis of 6-aminomethyl derivatives of benzopyran-4-one with dual biological properties: anti-inflammatory-analgesic and antimicrobial. <i>European Journal of Medicinal Chemistry</i> , 2009 , 44, 4896-903 ^{6,8}	6.8	22
55	Pharmacophore modeling, 3D-QSAR, docking study and ADME prediction of acyl 1,3,4-thiadiazole amides and sulfonamides as antitubulin agents. <i>Arabian Journal of Chemistry</i> , 2019 , 12, 5000-5018	5.9	22

54	4, 5-Dihydrooxazole-pyrazoline hybrids: Synthesis and their evaluation as potential antimalarial agents. <i>European Journal of Medicinal Chemistry</i> , 2016 , 123, 476-486	6.8	20
53	Unveiling novel diphenyl-1H-pyrazole based acrylates tethered to 1,2,3-triazole as promising apoptosis inducing cytotoxic and anti-inflammatory agents. <i>Bioorganic Chemistry</i> , 2019 , 87, 667-678	5.1	19
52	Hepatoprotective activity of Marrubium vulgare against paracetamol induced toxicity. <i>Journal of Pharmacy Research</i> , 2013 , 7, 565-570		19
51	Recent updates on biological activities of oxadiazoles. <i>Mini-Reviews in Medicinal Chemistry</i> , 2013 , 13, 1027-46	3.2	19
50	Pharmacophore modeling, 3D-QSAR, docking and ADME prediction of quinazoline based EGFR inhibitors. <i>Arabian Journal of Chemistry</i> , 2019 , 12, 4815-4839	5.9	19
49	Synthesis and biological evaluation of 2,5-disubstituted 1,3,4-oxadiazole derivatives with both COX and LOX inhibitory activity. <i>Journal of Enzyme Inhibition and Medicinal Chemistry</i> , 2011 , 26, 767-76	5.6	17
48	Synthesis of some new 3,4-dihydro-2H-1,3-benzoxazines under microwave irradiation in solvent-free conditions and their biological activity. <i>Medicinal Chemistry Research</i> , 2011 , 20, 1147-1153	2.2	17
47	Structure based virtual screening of MDPI database: discovery of structurally diverse and novel DPP-IV inhibitors. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2014 , 24, 3447-51	2.9	15
46	In silico approach for bioremediation of arsenic by structure prediction and docking studies of arsenite oxidase from Pseudomonas stutzeri TS44. <i>International Biodeterioration and Biodegradation</i> , 2017 , 122, 82-91	4.8	14
45	Novel pyrazole-pyrazoline hybrids endowed with thioamide as antimalarial agents: their synthesis and 3D-QSAR studies. <i>Journal of Enzyme Inhibition and Medicinal Chemistry</i> , 2015 , 30, 597-606	5.6	14
44	Synthesis of quinoline attached-furan-2(3H)-ones having anti-inflammatory and antibacterial properties with reduced gastro-intestinal toxicity and lipid peroxidation. <i>Journal of the Serbian Chemical Society</i> , 2011 , 76, 1617-1626	0.9	14
43	Synthesis of novel benzimidazole acrylonitriles for inhibition of Plasmodium falciparum growth by dual target inhibition. <i>Archiv Der Pharmazie</i> , 2018 , 351, 1700251	4.3	14
42	Novel hybrid-pyrrole derivatives: their synthesis, antitubercular evaluation and docking studies. <i>RSC Advances</i> , 2015 , 5, 12807-12820	3.7	13
41	Doxorubicin loaded carboxymethyl Assam bora rice starch coated superparamagnetic iron oxide nanoparticles as potential antitumor cargo. <i>Heliyon</i> , 2019 , 5, e01955	3.6	12
40	Structural comparison of Mtb-DHFR and h-DHFR for design, synthesis and evaluation of selective non-pteridine analogues as antitubercular agents. <i>Bioorganic Chemistry</i> , 2018 , 80, 319-333	5.1	12
39	Targeting malaria and leishmaniasis: Synthesis and pharmacological evaluation of novel pyrazole-1,3,4-oxadiazole hybrids. Part II. <i>Bioorganic Chemistry</i> , 2019 , 89, 102986	5.1	11
38	Synthesis and antimalarial activity of quinoline-substituted furanone derivatives and their identification as selective falcipain-2 inhibitors. <i>Medicinal Chemistry Research</i> , 2015 , 24, 879-890	2.2	11
37	Synthesis, 3D-QSAR and docking studies of pyrimidine nitrile-pyrazoline: a novel class of hybrid antimalarial agents. <i>Medicinal Chemistry Research</i> , 2015 , 24, 1018-1037	2.2	11

36	Synthesis of Hybrids of Dihydropyrimidine and Pyridazinone as potential Anti-Breast Cancer Agents. <i>Mini-Reviews in Medicinal Chemistry</i> , 2018 , 18, 369-379	3.2	11
35	Novel hydrazine derivatives as selective DPP-IV inhibitors: findings from virtual screening and validation through molecular dynamics simulations. <i>Journal of Molecular Modeling</i> , 2014 , 20, 2118	2	10
34	Design, synthesis and pharmacological evaluation of some novel derivatives of 1-[[3-(furan-2-yl)-5-phenyl-4,5-dihydro-1,2-oxazol-4-yl]methyl]-4-methyl piperazine. <i>Arabian Journal of Chemistry</i> , 2017 , 10, 141-149	5.9	10
33	3D-QSAR of amino-substituted pyrido[3,2B]pyrazinones as PDE-5 inhibitors. <i>Medicinal Chemistry Research</i> , 2012 , 21, 202-211	2.2	8
32	Synthesis, Antiinflammatory and Antimicrobial Activity of Some New 1-(3-Phenyl-3,4-Dihydro-2H-1,3-Benzoxazin-6-yl)-Ethanone Derivatives. <i>Indian Journal of Pharmaceutical Sciences</i> , 2011 , 73, 101-4	1.5	8
31	Design and synthesis of pyrazole-pyrazoline hybrids as cancer-associated selective COX-2 inhibitors. <i>Archiv Der Pharmazie</i> , 2021 , 354, e2000116	4.3	8
30	Synthesis, COX-2 inhibition and metabolic stability studies of 6-(4-fluorophenyl)-pyrimidine-5-carbonitrile derivatives as anticancer and anti-inflammatory agents. <i>Journal of Fluorine Chemistry</i> , 2020 , 236, 109579	2.1	7
29	Biochemical characterization of unusual cysteine protease of <i>P. falciparum</i> , metacaspase-2 (MCA-2). <i>Molecular and Biochemical Parasitology</i> , 2018 , 220, 28-41	1.9	7
28	3D-QSAR and molecular docking studies on 3-anilino-4-arylmaleimide derivatives as glycogen synthase kinase-3 inhibitors. <i>Chemical Biology and Drug Design</i> , 2012 , 79, 560-71	2.9	7
27	Molecular interactions of dioxins and DLCs with the xenosensors (PXR and CAR): An in silico risk assessment approach. <i>Journal of Molecular Recognition</i> , 2017 , 30, e2651	2.6	6
26	Pharmacophore based virtual screening, synthesis and SAR of novel inhibitors of Mycobacterium sulfotransferase. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2015 , 25, 701-7	2.9	6
25	Molecular interactions of dioxins and DLCs with the ketosteroid receptors: an in silico risk assessment approach. <i>Toxicology Mechanisms and Methods</i> , 2017 , 27, 151-163	3.6	5
24	Synthesis and biological evaluation of benzimidazole pendant cyanopyrimidine derivatives as anticancer agents. <i>Journal of Heterocyclic Chemistry</i> , 2020 , 57, 3350	1.9	5
23	Benzimidazole based derivatives as anticancer agents: Structure activity relationship analysis for various targets. <i>Journal of Heterocyclic Chemistry</i> ,	1.9	5
22	Dibenzepinones, dibenzoxepines and benzosuberones based p38MAP kinase inhibitors: Their pharmacophore modelling, 3D-QSAR and docking studies. <i>Computers in Biology and Medicine</i> , 2019 , 110, 175-185	7	4
21	Expansion of a novel lead targeting <i>M. tuberculosis</i> DHFR as antitubercular agents. <i>Bioorganic and Medicinal Chemistry</i> , 2019 , 27, 1421-1429	3.4	4
20	Methylene-bearing sulfur-containing cyanopyrimidine derivatives for treatment of cancer: Part-II. <i>Archiv Der Pharmazie</i> , 2020 , 353, e1900333	4.3	4
19	Molecular interactions of bisphenols and analogs with glucocorticoid biosynthetic pathway enzymes: an in silico approach. <i>Toxicology Mechanisms and Methods</i> , 2018 , 28, 45-54	3.6	4

18	Pharmacophore model generation and 3D-QSAR analysis of N-acyl and N-arylpiprazolines for enzymatic and cellular B-Raf kinase inhibition. <i>Medicinal Chemistry Research</i> , 2013 , 22, 2174-2187	2.2	4
17	Malaria: hitches and hopes. <i>Mini-Reviews in Medicinal Chemistry</i> , 2014 , 14, 453-70	3.2	4
16	Identification of novel small molecule non-peptidomimetic inhibitor for prolyl oligopeptidase through and approaches. <i>Journal of Biomolecular Structure and Dynamics</i> , 2020 , 38, 1292-1305	3.6	4
15	1, 2, 4-Oxadiazole Incorporated Ketoprofen Analogues in Search of Safer Non-steroidal Anti-inflammatory Agents: Design, Syntheses, Biological Evaluation and Molecular Docking Studies. <i>Letters in Drug Design and Discovery</i> , 2018 , 15, 590-601	0.8	3
14	Identification of novel selective Mtb-DHFR inhibitors as antitubercular agents through structure-based computational techniques. <i>Archiv Der Pharmazie</i> , 2020 , 353, e1900287	4.3	3
13	Synthesis, ADMET prediction and reverse screening study of 3,4,5-trimethoxy phenyl ring pendant sulfur-containing cyanopyrimidine derivatives as promising apoptosis inducing anticancer agents. <i>Bioorganic Chemistry</i> , 2020 , 104, 104282	5.1	3
12	Identification of novel Mycobacterium tuberculosis dihydrofolate reductase inhibitors through rational drug design. <i>International Journal of Mycobacteriology</i> , 2016 , 5 Suppl 1, S96	0.9	3
11	strategies for probing novel DPP-IV inhibitors as anti-diabetic agents. <i>Journal of Biomolecular Structure and Dynamics</i> , 2021 , 39, 2118-2132	3.6	3
10	NEW AMIDES OF SULPHONAMIDES: SYNTHESIS AND BIOLOGICAL EVALUATION. <i>Journal of the Chilean Chemical Society</i> , 2010 , 55,	2.5	2
9	Mining of potential dipeptidyl peptidase-IV inhibitors as anti-diabetic agents using integrated approaches. <i>Journal of Biomolecular Structure and Dynamics</i> , 2020 , 38, 5349-5361	3.6	2
8	Anti-proliferative and anti-malarial activities of spiroisoxazoline analogues of artemisinin. <i>Archiv Der Pharmazie</i> , 2018 , 352, e1800192	4.3	2
7	Synthesis and anticonvulsant activity of some newer dihydro-pyrimidine-5-carbonitrile derivatives: Part II. <i>Journal of Taibah University Medical Sciences</i> , 2015 , 10, 437-443	1.7	1
6	3D quantitative structure-activity relationship for quinoline, benzimidazole and benzofuran-based analogs as phosphodiesterases IV (PDE-IV) inhibitors. <i>Medicinal Chemistry Research</i> , 2013 , 22, 5153-5166	2.2	1
5	3D QSAR Based Virtual Screening of Pyrido[1,2-a] Benzimidazoles as Potent Antimalarial Agents. <i>Letters in Drug Design and Discovery</i> , 2019 , 16, 301-312	0.8	1
4	Ameliorative effect of rubiadin-loaded nanocarriers in STZ-NA-induced diabetic nephropathy in rats: formulation optimization, molecular docking, and in vivo biological evaluation. <i>Drug Delivery and Translational Research</i> , 2021 , 1	6.2	1
3	A review on synthetic procedures and applications of phosphorus oxychloride (POCl ₃) in the last biennial period (2018-19). <i>Phosphorus, Sulfur and Silicon and the Related Elements</i> , 2021 , 196, 211-274	1	1
2	Synthesis, biological evaluation and docking studies of methylene bearing cyanopyrimidine derivatives possessing a hydrazone moiety as potent Lysine specific demethylase-1 (LSD1) inhibitors: A promising anticancer agents. <i>Bioorganic Chemistry</i> , 2022 , 126, 105885	5.1	1
1	Identification, analysis of deleterious SNPs of the human GSR gene and their effects on the structure and functions of associated proteins and other diseases.. <i>Scientific Reports</i> , 2022 , 12, 5474	4.9	0

