

Mahmud Tareq Hassan Khan

List of Publications by Citations

Source:
<https://exaly.com/author-pdf/10494805/mahmud-tareq-hassan-khan-publications-by-citations.pdf>
Version: 2024-04-05

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.

32 papers	1,853 citations	22 h-index	32 g-index
32 ext. papers	2,120 ext. citations	3.9 avg, IF	4.86 L-index

#	Paper	IF	Citations
32	A comprehensive review on tyrosinase inhibitors. <i>Journal of Enzyme Inhibition and Medicinal Chemistry</i> , 2019 , 34, 279-309	5.6	331
31	Extracts and molecules from medicinal plants against herpes simplex viruses. <i>Antiviral Research</i> , 2005 , 67, 107-19	10.8	200
30	Cholinesterase inhibitory activities of some flavonoid derivatives and chosen xanthone and their molecular docking studies. <i>Chemico-Biological Interactions</i> , 2009 , 181, 383-9	5	156
29	Structure-activity relationships of tyrosinase inhibitory combinatorial library of 2,5-disubstituted-1,3,4-oxadiazole analogues. <i>Bioorganic and Medicinal Chemistry</i> , 2005 , 13, 3385-95	3.4	144
28	Predictions of the ADMET properties of candidate drug molecules utilizing different QSAR/QSPR modelling approaches. <i>Current Drug Metabolism</i> , 2010 , 11, 285-95	3.5	107
27	Tetraketones: a new class of tyrosinase inhibitors. <i>Bioorganic and Medicinal Chemistry</i> , 2006 , 14, 344-51	3.4	83
26	Oxazolones: new tyrosinase inhibitors; synthesis and their structure-activity relationships. <i>Bioorganic and Medicinal Chemistry</i> , 2006 , 14, 6027-33	3.4	79
25	Recent advancements for the evaluation of anti-viral activities of natural products. <i>New Biotechnology</i> , 2009 , 25, 347-68	6.4	71
24	Molecular design of tyrosinase inhibitors: A critical review of promising novel inhibitors from synthetic origins. <i>Pure and Applied Chemistry</i> , 2007 , 79, 2277-2295	2.1	70
23	TOMOCOMD-CARDD descriptors-based virtual screening of tyrosinase inhibitors: evaluation of different classification model combinations using bond-based linear indices. <i>Bioorganic and Medicinal Chemistry</i> , 2007 , 15, 1483-503	3.4	68
22	Dragon method for finding novel tyrosinase inhibitors: Biosilico identification and experimental in vitro assays. <i>European Journal of Medicinal Chemistry</i> , 2007 , 42, 1370-81	6.8	57
21	Molecular interactions of cholinesterases inhibitors using in silico methods: current status and future prospects. <i>New Biotechnology</i> , 2009 , 25, 331-46	6.4	48
20	New tyrosinase inhibitors selected by atomic linear indices-based classification models. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2006 , 16, 324-30	2.9	48
19	Prediction of tyrosinase inhibition activity using atom-based bilinear indices. <i>ChemMedChem</i> , 2007 , 2, 449-78	3.7	43
18	Tyrosinase inhibitors from <i>Rhododendron collettianum</i> and their structure-activity relationship (SAR) studies. <i>Chemical and Pharmaceutical Bulletin</i> , 2004 , 52, 1458-61	1.9	41
17	Ligand-based computer-aided discovery of tyrosinase inhibitors. Applications of the TOMOCOMD-CARDD method to the elucidation of new compounds. <i>Current Pharmaceutical Design</i> , 2010 , 16, 2601-24	3.3	33
16	Microbial transformation of 17 α -ethynyl- and 17 α -ethylsteroids, and tyrosinase inhibitory activity of transformed products. <i>Steroids</i> , 2005 , 70, 798-802	2.8	30

15	Discovery of potent thermolysin inhibitors using structure based virtual screening and binding assays. <i>Journal of Medicinal Chemistry</i> , 2009 , 52, 48-61	8.3	26
14	Ethnomedicines and ethnomedicinal phytophores against herpesviruses. <i>Biotechnology Annual Review</i> , 2008 , 14, 297-348		26
13	Predictive QSAR modeling for the successful predictions of the ADMET properties of candidate drug molecules. <i>Current Drug Discovery Technologies</i> , 2007 , 4, 141-9	1.5	26
12	Tyrosinase inhibition studies of cycloartane and cucurbitane glycosides and their structure-activity relationships. <i>Bioorganic and Medicinal Chemistry</i> , 2006 , 14, 6085-8	3.4	24
11	Atom- and bond-based 2D TOMOCOMD-CARDD approach and ligand-based virtual screening for the drug discovery of new tyrosinase inhibitors. <i>Journal of Biomolecular Screening</i> , 2008 , 13, 1014-24		23
10	Identification of novel quinazolin-4(3H)-ones as inhibitors of thermolysin, the prototype of the M4 family of proteinases. <i>Bioorganic and Medicinal Chemistry</i> , 2010 , 18, 4317-27	3.4	21
9	Expression of Estrogen Receptor Gene in Breast Cancer Cells Treated With Transcription Factor Decoy Is Modulated by Bangladeshi Natural Plant Extracts. <i>Oncology Research</i> , 2005 , 15, 69-79	4.8	21
8	Potentials of phenolic molecules of natural origin and their derivatives as anti-HIV agents. <i>Biotechnology Annual Review</i> , 2007 , 13, 223-64		18
7	Bond-based 2D TOMOCOMD-CARDD approach for drug discovery: aiding decision-making in 'in silico' selection of new lead tyrosinase inhibitors. <i>Journal of Computer-Aided Molecular Design</i> , 2007 , 21, 167-88	4.2	18
6	Synthesis of methyl ether analogues of sildenafil (Viagra) possessing tyrosinase inhibitory potential. <i>Chemistry and Biodiversity</i> , 2005 , 2, 470-6	2.5	17
5	Heterocyclic Compounds Against the Enzyme Tyrosinase Essential for Melanin Production: Biochemical Features of Inhibition 2007 , 119-138		11
4	Tyrosinase inhibitory effect of benzoic acid derivatives and their structure-activity relationships. <i>Journal of Enzyme Inhibition and Medicinal Chemistry</i> , 2010 , 25, 812-7	5.6	8
3	Dipeptide inhibitors of thermolysin and angiotensin I-converting enzyme. <i>Current Topics in Medicinal Chemistry</i> , 2012 , 12, 1748-62	3	3
2	Determinants for psychrophilic and thermophilic features of metallopeptidases of the M4 family. <i>In Silico Biology</i> , 2009 , 9, 105-24	2	2
1	Retrained Classification of Tyrosinase Inhibitors and In Silico Potency Estimation by Using Atom-Type Linear Indices. <i>International Journal of Chemoinformatics and Chemical Engineering</i> , 2012 , 2, 42-144	0.3	