Mohammed Gollapalli

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	Synthesis of benzothiazole derivatives as a potent α-glucosidase inhibitor. Bioorganic Chemistry, 2019, 85, 33-48.	4.1	54
2	Synthesis of Bis-indolylmethane sulfonohydrazides derivatives as potent α-Glucosidase inhibitors. Bioorganic Chemistry, 2018, 80, 112-120.	4.1	49
3	Rainfall Prediction System Using Machine Learning Fusion for Smart Cities. Sensors, 2022, 22, 3504.	3.8	47
4	Supervised Machine Learning Empowered Multifactorial Genetic Inheritance Disorder Prediction. Computational Intelligence and Neuroscience, 2022, 2022, 1-10.	1.7	47
5	Synthesis, in vitro urease inhibitory potential and molecular docking study of Benzimidazole analogues. Bioorganic Chemistry, 2019, 89, 103024.	4.1	45
6	A novel stacking ensemble for detecting three types of diabetes mellitus using a Saudi Arabian dataset: Pre-diabetes, T1DM, and T2DM. Computers in Biology and Medicine, 2022, 147, 105757.	7.0	35
7	Synthesis of new arylhydrazide bearing Schiff bases/thiazolidinone: α-Amylase, urease activities and their molecular docking studies. Bioorganic Chemistry, 2019, 91, 103112.	4.1	33
8	Synthesis of novel quinoline-based thiadiazole, evaluation of their antileishmanial potential and molecular docking studies. Bioorganic Chemistry, 2019, 85, 109-116.	4.1	25
9	Exploring indole-based-thiadiazole derivatives as potent acetylcholinesterase and butyrylcholinesterase enzyme inhibitors. International Journal of Biological Macromolecules, 2021, 188, 1025-1036.	7.5	20
10	Synthesis, anti-leishmanial and molecular docking study of bis-indole derivatives. BMC Chemistry, 2019, 13, 102.	3.8	18
11	Advance Genome Disorder Prediction Model Empowered With Deep Learning. IEEE Access, 2022, 10, 70317-70328.	4.2	15
12	Synthesis of oxadiazole-coupled-thiadiazole derivatives as a potent Î ² -glucuronidase inhibitors and their molecular docking study. Bioorganic and Medicinal Chemistry, 2019, 27, 3145-3155.	3.0	13
13	2,5-Disubstituted thiadiazoles as potent β-glucuronidase inhibitors; Synthesis, in vitro and in silico studies. Bioorganic Chemistry, 2019, 91, 103126.	4.1	12
14	Automated discovery of multi-faceted ontologies for accurate query answering and future semantic reasoning. Data and Knowledge Engineering, 2013, 87, 405-424.	3.4	11
15	Literature Review of Attribute Level and Structure Level Data Linkage Techniques. International Journal of Data Mining & Knowledge Management Process, 2015, 5, 01-20.	0.1	10
16	Indole bearing thiadiazole analogs: synthesis, β-glucuronidase inhibition and molecular docking study. BMC Chemistry, 2019, 13, 14.	3.8	10
17	Ensemble Machine Learning Model to Predict the Waterborne Syndrome. Algorithms, 2022, 15, 93.	2.1	10
18	Synthesis of Thymidine Phosphorylase Inhibitor Based on Quinoxaline Derivatives and Their Molecular Docking Study. Molecules, 2019, 24, 1002.	3.8	9

#	Article	IF	CITATIONS
19	Inhibition potential of phenyl linked benzimidazole-triazolothiadiazole modular hybrids against β-glucuronidase and their interactions thereof. International Journal of Biological Macromolecules, 2020, 161, 355-363.	7.5	9
20	Synthesis, thymidine phosphorylase, angiogenic inhibition and molecular docking study of isoquinoline derivatives. Bioorganic Chemistry, 2019, 89, 102999.	4.1	8
21	Ontology Guided Data Linkage Framework for Discovering Meaningful Data Facts. Lecture Notes in Computer Science, 2011, , 252-265.	1.3	7
22	Saudi Rural Breast Cancer Prevention Framework. , 2015, , .		6
23	Approximate Record Matching Using Hash Grams. , 2011, , .		5
24	Synthesis of Chromen-4-One-Oxadiazole Substituted Analogs as Potent Î ² -Glucuronidase Inhibitors. Molecules, 2019, 24, 1528.	3.8	5
25	The Saudi Recycling eCommerce Framework. , 2015, , .		1