

Linh Tran

List of Publications by Year in descending order

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Version: 2024-02-01

25
papers

388
citations

1040018

9
h-index

794568

19
g-index

25
all docs

25
docs citations

25
times ranked

645
citing authors

#	ARTICLE	IF	CITATIONS
1	Efficacy of Kinesio Taping Compared to Other Treatment Modalities in Musculoskeletal Disorders: A Systematic Review and Meta-Analysis. <i>Research in Sports Medicine</i> , 2023, 31, 416-439.	1.3	8
2	A systematic review of antimalarial activities of <i>Morinda</i> species. <i>South African Journal of Botany</i> , 2022, 148, 396-406.	2.5	1
3	The F19W mutation reduces the binding affinity of the transmembrane A β 40 trimer to the membrane bilayer. <i>RSC Advances</i> , 2021, 11, 2664-2676.	3.6	2
4	Role of cytokines produced by T helper immune-modulators in dengue pathogenesis: A systematic review and meta-analysis. <i>Acta Tropica</i> , 2021, 216, 105823.	2.0	5
5	Quality assessment tools used in systematic reviews of in vitro studies: A systematic review. <i>BMC Medical Research Methodology</i> , 2021, 21, 101.	3.1	39
6	Efficacy of chalcone and xanthine derivatives on lipase inhibition: A systematic review. <i>Chemical Biology and Drug Design</i> , 2020, 95, 205-214.	3.2	3
7	Association of breastfeeding status with risk of autism spectrum disorder: A systematic review, dose-response analysis and meta-analysis. <i>Asian Journal of Psychiatry</i> , 2020, 48, 101916.	2.0	28
8	Frequent inappropriate use of unweighted summary statistics in systematic reviews of pathogen genotypes or genogroups. <i>Journal of Clinical Epidemiology</i> , 2020, 119, 26-35.	5.0	0
9	Efficacy of kinesio taping in treatment of shoulder pain and disability: a systematic review and meta-analysis of randomised controlled trials. <i>Physiotherapy</i> , 2020, 107, 176-188.	0.4	28
10	<i>Morinda morindoides</i> : A systematic review of its therapeutic activities. <i>South African Journal of Botany</i> , 2020, 131, 93-103.	2.5	4
11	Effects of Mulberry on The Central Nervous System: A Literature Review. <i>Current Neuropharmacology</i> , 2020, 19, 193-219.	2.9	9
12	Propafenone effects on the stable structures of A β 16-22 system. <i>Chemical Physics Letters</i> , 2018, 696, 55-60.	2.6	6
13	Ginsenoside Rh1: A Systematic Review of Its Pharmacological Properties. <i>Planta Medica</i> , 2018, 84, 139-152.	1.3	66
14	Binding Modes of a Glycopeptidomimetic Molecule on A β Protofibrils: Implication for Its Inhibition Mechanism. <i>ACS Chemical Neuroscience</i> , 2018, 9, 2859-2869.	3.5	10
15	Understanding the Binding Mechanism of Amyloid- β Inhibitors from Molecular Simulations. <i>Current Pharmaceutical Design</i> , 2018, 24, 3341-3346.	1.9	2
16	Ginsenoside Rk1 bioactivity: a systematic review. <i>PeerJ</i> , 2017, 5, e3993.	2.0	20
17	Effect of Post-Translational Amidation on Islet Amyloid Polypeptide Conformational Ensemble: Implications for Its Aggregation Early Steps. <i>International Journal of Molecular Sciences</i> , 2016, 17, 1896.	4.1	7
18	Structure of ring-shaped A β 42 oligomers determined by conformational selection. <i>Scientific Reports</i> , 2016, 6, 21429.	3.3	32

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19	Comparative study of structural models of Leishmania donovani and human GDP-mannose pyrophosphorylases. European Journal of Medicinal Chemistry, 2016, 107, 109-118.	5.5	12
20	Insights into the Conformational Ensemble of Human Islet Amyloid Polypeptide from Molecular Simulations. Current Pharmaceutical Design, 2016, 22, 3601-3607.	1.9	5
21	Exploring the Alzheimer amyloid- β^2 peptide conformational ensemble: A review of molecular dynamics approaches. Peptides, 2015, 69, 86-91.	2.4	69
22	Recent progress and challenges in the computer-aided design of inhibitors for influenza A M2 channel proteins. Medicinal Chemistry Research, 2014, 23, 3607-3616.	2.4	4
23	Using imaging ToF-MS data to determine the cell wall thickness of fibers in wood. Surface and Interface Analysis, 2014, 46, 225-228.	1.8	4
24	Strategy in structure-based drug design for influenza A virus targeting M2 channel proteins. Medicinal Chemistry Research, 2013, 22, 6078-6088.	2.4	4
25	Discovery of Potential M2 Channel Inhibitors Based on the Amantadine Scaffold via Virtual Screening and Pharmacophore Modeling. Molecules, 2011, 16, 10227-10255.	3.8	20