Linh Tran

List of Publications by Year in descending order

Source: https://exaly.com/author-pdf/1757139/publications.pdf

Version: 2024-02-01

25	388	9	19
papers	citations	h-index	g-index
25	25	25	645
all docs	docs citations	times ranked	citing authors

#	Article	IF	CITATIONS
1	Efficacy of Kinesio Taping Compared to Other Treatment Modalities in Musculoskeletal Disorders: A Systematic Review and Meta-Analysis. Research in Sports Medicine, 2023, 31, 416-439.	1.3	8
2	A systematic review of antimalarial activities of Morinda species. South African Journal of Botany, 2022, 148, 396-406.	2.5	1
3	The F19W mutation reduces the binding affinity of the transmembrane Aβ11–40 trimer to the membrane bilayer. RSC Advances, 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. Acta Tropica, 2021, 216, 105823.	2.0	5
5	Quality assessment tools used in systematic reviews of in vitro studies: A systematic review. BMC Medical Research Methodology, 2021, 21, 101.	3.1	39
6	Efficacy of chalcone and xanthine derivatives on lipase inhibition: A systematic review. Chemical Biology and Drug Design, 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. Asian Journal of Psychiatry, 2020, 48, 101916.	2.0	28
8	Frequent inappropriate use of unweighted summary statistics in systematic reviews of pathogen genotypes or genogroups. Journal of Clinical Epidemiology, 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. Physiotherapy, 2020, 107, 176-188.	0.4	28
10	Morinda morindoides: A systematic review of its therapeutic activities. South African Journal of Botany, 2020, 131, 93-103.	2.5	4
11	Effects of Mulberry on The Central Nervous System: A Literature Review. Current Neuropharmacology, 2020, 19, 193-219.	2.9	9
12	Propafenone effects on the stable structures of A \hat{I}^2 16-22 system. Chemical Physics Letters, 2018, 696, 55-60.	2.6	6
13	Ginsenoside Rh1: A Systematic Review of Its Pharmacological Properties. Planta Medica, 2018, 84, 139-152.	1.3	66
14	Binding Modes of a Glycopeptidomimetic Molecule on $\hat{Al^2}$ Protofibrils: Implication for Its Inhibition Mechanism. ACS Chemical Neuroscience, 2018, 9, 2859-2869.	3.5	10
15	Understanding the Binding Mechanism of Amyloid- $\hat{l^2}$ Inhibitors from Molecular Simulations. Current Pharmaceutical Design, 2018, 24, 3341-3346.	1.9	2
16	Ginsenoside Rk1 bioactivity: a systematic review. PeerJ, 2017, 5, e3993.	2.0	20
17	Effect of Post-Translational Amidation on Islet Amyloid Polypeptide Conformational Ensemble: Implications for Its Aggregation Early Steps. International Journal of Molecular Sciences, 2016, 17, 1896.	4.1	7
18	Structure of ring-shaped $\hat{Al^2}42$ oligomers determined by conformational selection. Scientific Reports, 2016, 6, 21429.	3.3	32

#	Article	IF	CITATION
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- \hat{l}^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 ToFâ€SIMS 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