

# Yiping Li

## List of Publications by Year in descending order

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Version: 2024-02-01

23  
papers

353  
citations

933447

10  
h-index

839539

18  
g-index

23  
all docs

23  
docs citations

23  
times ranked

573  
citing authors

#	ARTICLE	IF	CITATIONS
1	Study on the interaction between nimodipine and five proteinases and the effects of naringin and vitamin C on these interactions by spectroscopic and molecular docking methods. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2022, 272, 120982.	3.9	4
2	Natural Bioactive Molecules as Potential Agents Against SARS-CoV-2. <i>Frontiers in Pharmacology</i> , 2021, 12, 702472.	3.5	8
3	Fluorescence polarization assay for the identification and evaluation of inhibitors at YAPâ€“TEAD proteinâ€“protein interface 3. <i>Analytical Biochemistry</i> , 2019, 586, 113413.	2.4	17
4	Observation of the generation of peroxyxynitrite in mouse liver after acetaminophen overdose with a boronate-based ratiometric fluorescence probe. <i>RSC Advances</i> , 2019, 9, 6510-6514.	3.6	5
5	Polyactin A is a novel and potent immunological adjuvant for peptide-based cancer vaccine. <i>International Immunopharmacology</i> , 2018, 54, 95-102.	3.8	14
6	Synthesis, nitric oxide release, and dipeptidyl peptidase-4 inhibition of sitagliptin derivatives as new multifunctional antidiabetic agents. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2018, 28, 3731-3735.	2.2	7
7	SnCl <sub>4</sub> grafted on silica gel: an efficient catalyst for solvent-free synthesis of coumarins via the Pechmann condensation. <i>Journal of the Iranian Chemical Society</i> , 2017, 14, 737-742.	2.2	8
8	Transbody against virus core protein potently inhibits hepadnavirus replication <i>in vivo</i> : evidence from a duck model of hepatitis B virus. <i>British Journal of Pharmacology</i> , 2017, 174, 2261-2272.	5.4	5
9	A new multifunctional hydroxytyrosol-fenofibrate with antidiabetic, antihyperlipidemic, antioxidant and antiinflammatory action. <i>Biomedicine and Pharmacotherapy</i> , 2017, 95, 1749-1758.	5.6	20
10	Phenylsulfonylfuroxan NO-donor phenols: Synthesis and multifunctional activities evaluation. <i>Bioorganic and Medicinal Chemistry</i> , 2017, 25, 4407-4413.	3.0	7
11	Molecular Dynamics Simulations to Investigate the Binding Mode of the Natural Product Liphagal with Phosphoinositide 3-Kinase I $\alpha$ . <i>Molecules</i> , 2016, 21, 857.	3.8	4
12	Spectroscopy and molecular docking study on the interaction of daidzein and genistein with pepsin. <i>Luminescence</i> , 2016, 31, 1524-1531.	2.9	16
13	Transbody against hepatitis B virus core protein inhibits hepatitis B virus replication <i>in vitro</i> . <i>International Immunopharmacology</i> , 2015, 25, 363-369.	3.8	7
14	Discovery of novel antagonists of glycoprotein IIb/IIIa-mediated platelet aggregation through virtual screening. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2015, 25, 1249-1253.	2.2	10
15	Dissociation Constants and Solubilities of Dalbergin and Nordalbergin in Different Solvents. <i>Journal of Chemical &amp; Engineering Data</i> , 2015, 60, 2585-2593.	1.9	8
16	Oxone-mediated oxidative carbon-heteroatom bond cleavage: synthesis of benzoxazinones from benzoxazoles with I $\alpha$ -oxocarboxylic acids. <i>RSC Advances</i> , 2014, 4, 8720.	3.6	22
17	Definition of the binding mode of phosphoinositide 3-kinase I $\alpha$ -selective inhibitor A-66S through molecular dynamics simulation. <i>Journal of Molecular Modeling</i> , 2014, 20, 2166.	1.8	7
18	Dissociation Constants and Solubilities of Daidzein and Genistein in Different Solvents. <i>Journal of Chemical &amp; Engineering Data</i> , 2014, 59, 1304-1311.	1.9	41

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19	A new series of N2-substituted-5-(p-toluenesulfonylamino)phthalimide analogues as $\alpha$ -glucosidase inhibitors. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2013, 23, 2022-2026.	2.2	17
20	Characterization of molecular recognition of Phosphoinositide-3-kinase $\alpha$ inhibitor through molecular dynamics simulation. <i>Journal of Molecular Modeling</i> , 2012, 18, 1907-1916.	1.8	9
21	Pharmacophore modeling and 3D-QSAR analysis of phosphoinositide 3-kinase p110 $\alpha$ inhibitors. <i>Journal of Molecular Modeling</i> , 2010, 16, 1449-1460.	1.8	32
22	An approach to develop binary chromatographic fingerprints of the total alkaloids from <i>Caulophyllum robustum</i> by high performance liquid chromatography/diode array detector and gas chromatography/mass spectrometry. <i>Journal of Pharmaceutical and Biomedical Analysis</i> , 2007, 43, 1667-1672.	2.8	53
23	Establishment of the model of vascular endothelial cell membrane chromatography and its preliminary application. <i>Science Bulletin</i> , 2007, 52, 922-928.	1.7	32