

Xiaojiao Yi

List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

8

papers

121

citations

6

h-index

9

g-index

9

ext. papers

173

ext. citations

3.7

avg, IF

2.6

L-index

#	Paper	IF	Citations
8	Global Metabolomic and Lipidomic Analysis Reveal the Synergistic Effect of Bufalin in Combination with Cinobufagin against HepG2 Cells. <i>Journal of Proteome Research</i> , 2020 , 19, 873-883	5.6	5
7	Integration of Transcriptomics and Metabolomics Reveals the Antitumor Mechanism Underlying Shikonin in Colon Cancer. <i>Frontiers in Pharmacology</i> , 2020 , 11, 544647	5.6	6
6	Systems Pharmacology-Based Approach to Comparatively Study the Independent and Synergistic Mechanisms of Danhong Injection and Naoxintong Capsule in Ischemic Stroke Treatment. <i>Evidence-based Complementary and Alternative Medicine</i> , 2019 , 2019, 1056708	2.3	13
5	Investigation of the reverse effect of Danhong injection on doxorubicin-induced cardiotoxicity in H9c2 cells: Insight by LC-MS based non-targeted metabolomic analysis. <i>Journal of Pharmaceutical and Biomedical Analysis</i> , 2018 , 152, 264-270	3.5	17
4	Rapid screening of brain-penetrable antioxidants from natural products by blood-brain barrier specific permeability assay combined with DPPH recognition. <i>Journal of Pharmaceutical and Biomedical Analysis</i> , 2018 , 151, 42-48	3.5	10
3	Chemical profiling and antioxidant evaluation of Yangxinshi Tablet by HPLC-ESI-Q-TOF-MS/MS combined with DPPH assay. <i>Journal of Chromatography B: Analytical Technologies in the Biomedical and Life Sciences</i> , 2017 , 1060, 262-271	3.2	36
2	Immobilized fusion protein affinity chromatography combined with HPLC-ESI-Q-TOF-MS/MS for rapid screening of PPAR ligands from natural products. <i>Talanta</i> , 2017 , 165, 508-515	6.2	14
1	Drug-protein binding of Danhong injection and the potential influence of drug combination with aspirin: Insight by ultrafiltration LC-MS and molecular modeling. <i>Journal of Pharmaceutical and Biomedical Analysis</i> , 2017 , 134, 100-107	3.5	20