Sergey M Ivanov

List of Publications by Year in descending order

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SEDCEV M IVANOV

#	Article	IF	CITATIONS
1	Chemo- and bioinformatics resources for inÂsilico drug discovery from medicinal plants beyond their traditional use: a critical review. Natural Product Reports, 2014, 31, 1585-1611.	5.2	104
2	DIGEP-Pred: web service for <i>in silico</i> prediction of drug-induced gene expression profiles based on structural formula. Bioinformatics, 2013, 29, 2062-2063.	1.8	87
3	In silico assessment of adverse drug reactions and associated mechanisms. Drug Discovery Today, 2016, 21, 58-71.	3.2	51
4	ADVERPred–Web Service for Prediction of Adverse Effects of Drugs. Journal of Chemical Information and Modeling, 2018, 58, 8-11.	2.5	50
5	Web Resources for Discovery and Development of New Medicines. Pharmaceutical Chemistry Journal, 2017, 51, 91-99.	0.3	18
6	Comparison of Quantitative and Qualitative (Q)SAR Models Created for the Prediction of Ki and IC50 Values of Antitarget Inhibitors. Frontiers in Pharmacology, 2018, 9, 1136.	1.6	17
7	Combined network pharmacology and virtual reverse pharmacology approaches for identification of potential targets to treat vascular dementia. Scientific Reports, 2020, 10, 257.	1.6	16
8	Identification of Drug-Induced Myocardial Infarction-Related Protein Targets through the Prediction of Drug–Target Interactions and Analysis of Biological Processes. Chemical Research in Toxicology, 2014, 27, 1263-1281.	1.7	14
9	Network-Based Analysis of OMICs Data to Understand the HIV–Host Interaction. Frontiers in Microbiology, 2020, 11, 1314.	1.5	14
10	A Computational Approach for the Prediction of Treatment History and the Effectiveness or Failure of Antiretroviral Therapy. International Journal of Molecular Sciences, 2020, 21, 748.	1.8	13
11	In silicomethod for identification of promising anticancer drug targets. SAR and QSAR in Environmental Research, 2009, 20, 755-766.	1.0	12
12	<i>In Silico</i> Identification of Proteins Associated with Drug-induced Liver Injury Based on the Prediction of Drug-target Interactions. Molecular Informatics, 2017, 36, 1600142.	1.4	12
13	Data and Text Mining Help Identify Key Proteins Involved in the Molecular Mechanisms Shared by SARS-CoV-2 and HIV-1. Molecules, 2020, 25, 2944.	1.7	12
14	Identification of Drug Targets Related to the Induction of Ventricular Tachyarrhythmia Through a Systems Chemical Biology Approach. Toxicological Sciences, 2015, 145, 321-336.	1.4	11
15	Identification of potential drug targets for treatment of refractory epilepsy using network pharmacology. Journal of Bioinformatics and Computational Biology, 2018, 16, 1840002.	0.3	10
16	Attenuation of hyperhomocysteinemia induced vascular dementia by sodium orthovanadate perhaps via PTP1B: Pertinent downstream outcomes. Behavioural Brain Research, 2019, 364, 29-40.	1.2	9
17	Glycogen synthase kinase-3 inhibition as a potential pharmacological target for vascular dementia: In silico and in vivo evidence. Computers in Biology and Medicine, 2019, 108, 305-316.	3.9	8
18	Integral estimation of xenobiotics' toxicity with regard to their metabolism in human organism. Pure and Applied Chemistry, 2017, 89, 1449-1458.	0.9	8

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19	Assessment of the cardiovascular adverse effects of drug-drug interactions through a combined analysis of spontaneous reports and predicted drug-target interactions. PLoS Computational Biology, 2019, 15, e1006851.	1.5	7
20	Computer search for molecular mechanisms of ulcerogenic action of non-steroidal anti-inflammatory drugs. Biochemistry (Moscow) Supplement Series B: Biomedical Chemistry, 2013, 7, 40-45.	0.2	4
21	Automated Extraction of Information From Texts of Scientific Publications: Insights Into HIV Treatment Strategies. Frontiers in Genetics, 2020, 11, 618862.	1.1	4
22	A computational analysis of transcriptional profiles from CD8(+) T lymphocytes reveals potential mechanisms of HIV/AIDS control and progression. Computational and Structural Biotechnology Journal, 2021, 19, 2447-2459.	1.9	4
23	Relationships between the Structure and Severe Drug-Induced Liver Injury for Low, Medium, and High Doses of Drugs. Chemical Research in Toxicology, 2022, 35, 402-411.	1.7	4
24	Computer Prediction of Adverse Drug Effects on the Cardiovascular System. Pharmaceutical Chemistry Journal, 2018, 52, 758-762.	0.3	3
25	Bioinformatics guided rotenone adjuvant kindling in mice as a new animal model of drug-resistant epilepsy. Computers in Biology and Medicine, 2022, 147, 105754.	3.9	2