Vladimir R Perović

List of Publications by Year in descending order

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Version: 2024-02-01

47 papers 1,024 citations

16 h-index 30 g-index

54 all docs

54 docs citations 54 times ranked 1343 citing authors

#	Article	IF	CITATIONS
1	Identification of SARSâ€CoVâ€2 Papainâ€like Protease (PLpro) Inhibitors Using Combined Computational Approach**. ChemistryOpen, 2022, 11, e202100248.	1.9	8
2	In Silico Screening of Natural Compounds for Candidates 5HT6 Receptor Antagonists against Alzheimer's Disease. Molecules, 2022, 27, 2626.	3.8	4
3	Alignment-free method for functional annotation of amino acid substitutions: Application on epigenetic factors involved in hematologic malignancies. PLoS ONE, 2021, 16, e0244948.	2.5	O
4	DiNGO: standalone application for Gene Ontology and Human Phenotype Ontology term enrichment analysis. Bioinformatics, 2020, 36, 1981-1982.	4.1	2
5	Biological Rationale for the Repurposing of BCG Vaccine against SARS-CoV-2. Journal of Proteome Research, 2020, 19, 4649-4654.	3.7	11
6	Drug Repurposing for Candidate SARS-CoV-2 Main Protease Inhibitors by a Novel In Silico Method. Molecules, 2020, 25, 3830.	3.8	49
7	Tally-2.0: upgraded validator of tandem repeat detection in protein sequences. Bioinformatics, 2020, 36, 3260-3262.	4.1	2
8	Automated feature engineering improves prediction of protein–protein interactions. Amino Acids, 2019, 51, 1187-1200.	2.7	18
9	Virtual Screen for Repurposing of Drugs for Candidate Influenza a M2 Ion-Channel Inhibitors. Frontiers in Cellular and Infection Microbiology, 2019, 9, 67.	3.9	19
10	The CAFA challenge reports improved protein function prediction and new functional annotations for hundreds of genes through experimental screens. Genome Biology, 2019, 20, 244.	8.8	261
11	Mapping of Protein-Protein Interactions: Web-Based Resources for Revealing Interactomes. Current Medicinal Chemistry, 2019, 26, 3890-3910.	2.4	11
12	Ibuprofen as a template molecule for drug design against Ebola virus. Frontiers in Bioscience - Landmark, 2018, 23, 947-953.	3.0	23
13	IDPpi: Protein-Protein Interaction Analyses of Human Intrinsically Disordered Proteins. Scientific Reports, 2018, 8, 10563.	3.3	18
14	TRI_tool: a web-tool for prediction of protein–protein interactions in human transcriptional regulation. Bioinformatics, 2017, 33, 289-291.	4.1	17
15	Identification of Candidate Allosteric Modulators of the M1 Muscarinic Acetylcholine Receptor Which May Improve Vagus Nerve Stimulation in Chronic Tinnitus. Frontiers in Neuroscience, 2017, 11, 636.	2.8	8
16	A simple method for calculation of basic molecular properties of nutrients and their use as a criterion for a healthy diet. F1000Research, 2017, 6, 13.	1.6	2
17	In silico Therapeutics for Neurogenic Hypertension and Vasovagal Syncope. Frontiers in Neuroscience, 2016, 9, 520.	2.8	2
18	Arginase Flavonoid Anti-Leishmanial in Silico Inhibitors Flagged against Anti-Targets. Molecules, 2016, 21, 589.	3.8	24

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19	A combined ligand- and structure-based approach for the identification of rilmenidine-derived compounds which synergize the antitumor effects of doxorubicin. Bioorganic and Medicinal Chemistry, 2016, 24, 3174-3183.	3.0	15
20	Common molecular mechanism of the hepatic lesion and the cardiac parasympathetic regulation in chronic hepatitis C infection: a critical role for the muscarinic receptor type 3. BMC Bioinformatics, 2016, 17, 139.	2.6	5
21	Rilmenidine suppresses proliferation and promotes apoptosis via the mitochondrial pathway in human leukemic K562 cells. European Journal of Pharmaceutical Sciences, 2016, 81, 172-180.	4.0	11
22	Predicted Enhanced Human Propensity of Current Avian-Like H1N1 Swine Influenza Virus from China. PLoS ONE, 2016, 11, e0165451.	2.5	2
23	Role of genetic markers in sport and recreational physical activity. FiziÄka Kultura, 2016, 70, 5-13.	0.2	0
24	Simple Chemoinformatics Criterion Using Electron Donor-Acceptor Molecular Characteristics for Selection of Antibiotics Against Multi-Drug-Resistant Bacteria. Discoveries, 2016, 4, e64.	2.3	0
25	Evolution of 2014/15 H3N2 Influenza Viruses Circulating in US: Consequences for Vaccine Effectiveness and Possible New Pandemic. Frontiers in Microbiology, 2015, 6, 1456.	3.5	13
26	In silico analysis suggests interaction between Ebola virus and the extracellular matrix. Frontiers in Microbiology, 2015, 6, 135.	3.5	24
27	Improving attrition rates in Ebola virus drug discovery. Expert Opinion on Drug Discovery, 2015, 10, 1025-1032.	5.0	3
28	Virtual screen for repurposing approved and experimental drugs for candidate inhibitors of EBOLA virus infection. F1000Research, 2015, 4, 34.	1.6	41
29	In silico analysis suggests repurposing of ibuprofen for prevention and treatment of EBOLA virus disease. F1000Research, 2015, 4, 104.	1.6	23
30	Natural Products as Promising Therapeutics for Treatment of Influenza Disease. Current Pharmaceutical Design, 2015, 21, 5573-5588.	1.9	27
31	Influenza vaccine as prevention for cardiovascular diseases: Possible molecular mechanism. Vaccine, 2014, 32, 6569-6575.	3.8	51
32	Novel algorithm for phylogenetic analysis of proteins: application to analysis of the evolution of H5N1 influenza viruses. Journal of Mathematical Chemistry, 2013, 51, 2238-2255.	1.5	10
33	New in silico and conventional in vitro approaches to advance HIV drug discovery and design. Expert Opinion on Drug Discovery, 2013, 8, 83-92.	5.0	5
34	Novel Phylogenetic Algorithm to Monitor Human Tropism in Egyptian H5N1-HPAIV Reveals Evolution toward Efficient Human-to-Human Transmission. PLoS ONE, 2013, 8, e61572.	2.5	33
35	Feature-Based Classification of Amino Acid Substitutions outside Conserved Functional Protein Domains. Scientific World Journal, The, 2013, 2013, 1-10.	2.1	5
36	Simple and General Criterion for "in silico―Screening of Candidate HIV Drugs. Current Pharmaceutical Biotechnology, 2013, 14, 561-569.	1.6	13

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37	Assessment of Hepatitis C Virus Protein Sequences with Regard to Interferon/Ribavirin Combination Therapy Response in Patients with HCV Genotype 1b. Protein Journal, 2012, 31, 129-136.	1.6	8
38	The role of long-range intermolecular interactions in discovery of new drugs. Expert Opinion on Drug Discovery, 2011, 6, 1263-1270.	5.0	26
39	Characterization of conserved properties of hemagglutinin of H5N1 and human influenza viruses: possible consequences for therapy and infection control. BMC Structural Biology, 2009, 9, 21.	2.3	67
40	Identification of hemagglutinin structural domain and polymorphisms which may modulate swine H1N1 interactions with human receptor. BMC Structural Biology, 2009, 9, 62.	2.3	62
41	Lipoprotein lipase: A bioinformatics criterion for assessment of mutations as a risk factor for cardiovascular disease. Proteins: Structure, Function and Bioinformatics, 2008, 70, 855-862.	2.6	15
42	Discovery of New Therapeutic Targets by the Informational Spectrum Method. Current Protein and Peptide Science, 2008, 9, 493-506.	1.4	34
43	Novel Virtual Screening Protocol Based on the Combined Use of Molecular Modeling and Electron-lon Interaction Potential Techniques To Design HIV-1 Integrase Inhibitors. Journal of Chemical Information and Modeling, 2007, 47, 1536-1544.	5.4	32
44	In Silico Criterion for Prediction of Effects of p53 Gene Missense Mutations on p53-Mdm2 Feedback Loop. Protein and Peptide Letters, 2006, 13, 807-814.	0.9	2
45	Evolution of SARS-CoV-2 virus and assessment of the effectiveness of COVID-19 vaccine. F1000Research, 0, 10, 28.	1.6	2
46	Prediction of the effectiveness of COVID-19 vaccine candidates. F1000Research, 0, 9, 365.	1.6	5
47	Virtual screen for repurposing approved and experimental drugs for candidate inhibitors of EBOLA virus infection. F1000Research, 0, 4, 34.	1.6	7