

# Vladimir R PeroviÄ

## List of Publications by Year in descending order

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Version: 2024-02-01

47  
papers

1,024  
citations

516710

16  
h-index

454955

30  
g-index

54  
all docs

54  
docs citations

54  
times ranked

1343  
citing authors

#	ARTICLE	IF	CITATIONS
1	The CAFA challenge reports improved protein function prediction and new functional annotations for hundreds of genes through experimental screens. <i>Genome Biology</i> , 2019, 20, 244.	8.8	261
2	Characterization of conserved properties of hemagglutinin of H5N1 and human influenza viruses: possible consequences for therapy and infection control. <i>BMC Structural Biology</i> , 2009, 9, 21.	2.3	67
3	Identification of hemagglutinin structural domain and polymorphisms which may modulate swine H1N1 interactions with human receptor. <i>BMC Structural Biology</i> , 2009, 9, 62.	2.3	62
4	Influenza vaccine as prevention for cardiovascular diseases: Possible molecular mechanism. <i>Vaccine</i> , 2014, 32, 6569-6575.	3.8	51
5	Drug Repurposing for Candidate SARS-CoV-2 Main Protease Inhibitors by a Novel In Silico Method. <i>Molecules</i> , 2020, 25, 3830.	3.8	49
6	Virtual screen for repurposing approved and experimental drugs for candidate inhibitors of EBOLA virus infection. <i>F1000Research</i> , 2015, 4, 34.	1.6	41
7	Discovery of New Therapeutic Targets by the Informational Spectrum Method. <i>Current Protein and Peptide Science</i> , 2008, 9, 493-506.	1.4	34
8	Novel Phylogenetic Algorithm to Monitor Human Tropism in Egyptian H5N1-HPAIV Reveals Evolution toward Efficient Human-to-Human Transmission. <i>PLoS ONE</i> , 2013, 8, e61572.	2.5	33
9	Novel Virtual Screening Protocol Based on the Combined Use of Molecular Modeling and Electron-Ion Interaction Potential Techniques To Design HIV-1 Integrase Inhibitors. <i>Journal of Chemical Information and Modeling</i> , 2007, 47, 1536-1544.	5.4	32
10	Natural Products as Promising Therapeutics for Treatment of Influenza Disease. <i>Current Pharmaceutical Design</i> , 2015, 21, 5573-5588.	1.9	27
11	The role of long-range intermolecular interactions in discovery of new drugs. <i>Expert Opinion on Drug Discovery</i> , 2011, 6, 1263-1270.	5.0	26
12	In silico analysis suggests interaction between Ebola virus and the extracellular matrix. <i>Frontiers in Microbiology</i> , 2015, 6, 135.	3.5	24
13	Arginase Flavonoid Anti-Leishmanial in Silico Inhibitors Flagged against Anti-Targets. <i>Molecules</i> , 2016, 21, 589.	3.8	24
14	Ibuprofen as a template molecule for drug design against Ebola virus. <i>Frontiers in Bioscience - Landmark</i> , 2018, 23, 947-953.	3.0	23
15	In silico analysis suggests repurposing of ibuprofen for prevention and treatment of EBOLA virus disease. <i>F1000Research</i> , 2015, 4, 104.	1.6	23
16	Virtual Screen for Repurposing of Drugs for Candidate Influenza a M2 Ion-Channel Inhibitors. <i>Frontiers in Cellular and Infection Microbiology</i> , 2019, 9, 67.	3.9	19
17	IDPpi: Protein-Protein Interaction Analyses of Human Intrinsically Disordered Proteins. <i>Scientific Reports</i> , 2018, 8, 10563.	3.3	18
18	Automated feature engineering improves prediction of proteinâ€“protein interactions. <i>Amino Acids</i> , 2019, 51, 1187-1200.	2.7	18

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19	TRI_tool: a web-tool for prediction of proteinâ€“protein interactions in human transcriptional regulation. <i>Bioinformatics</i> , 2017, 33, 289-291.	4.1	17
20	Lipoprotein lipase: A bioinformatics criterion for assessment of mutations as a risk factor for cardiovascular disease. <i>Proteins: Structure, Function and Bioinformatics</i> , 2008, 70, 855-862.	2.6	15
21	A combined ligand- and structure-based approach for the identification of rilmenidine-derived compounds which synergize the antitumor effects of doxorubicin. <i>Bioorganic and Medicinal Chemistry</i> , 2016, 24, 3174-3183.	3.0	15
22	Evolution of 2014/15 H3N2 Influenza Viruses Circulating in US: Consequences for Vaccine Effectiveness and Possible New Pandemic. <i>Frontiers in Microbiology</i> , 2015, 6, 1456.	3.5	13
23	Simple and General Criterion for â€œin silicoâ€“Screening of Candidate HIV Drugs. <i>Current Pharmaceutical Biotechnology</i> , 2013, 14, 561-569.	1.6	13
24	Rilmenidine suppresses proliferation and promotes apoptosis via the mitochondrial pathway in human leukemic K562 cells. <i>European Journal of Pharmaceutical Sciences</i> , 2016, 81, 172-180.	4.0	11
25	Biological Rationale for the Repurposing of BCG Vaccine against SARS-CoV-2. <i>Journal of Proteome Research</i> , 2020, 19, 4649-4654.	3.7	11
26	Mapping of Protein-Protein Interactions: Web-Based Resources for Revealing Interactomes. <i>Current Medicinal Chemistry</i> , 2019, 26, 3890-3910.	2.4	11
27	Novel algorithm for phylogenetic analysis of proteins: application to analysis of the evolution of H5N1 influenza viruses. <i>Journal of Mathematical Chemistry</i> , 2013, 51, 2238-2255.	1.5	10
28	Assessment of Hepatitis C Virus Protein Sequences with Regard to Interferon/Ribavirin Combination Therapy Response in Patients with HCV Genotype 1b. <i>Protein Journal</i> , 2012, 31, 129-136.	1.6	8
29	Identification of Candidate Allosteric Modulators of the M1 Muscarinic Acetylcholine Receptor Which May Improve Vagus Nerve Stimulation in Chronic Tinnitus. <i>Frontiers in Neuroscience</i> , 2017, 11, 636.	2.8	8
30	Identification of SARSâ€“CoVâ€“2 Papainâ€“like Protease (PLpro) Inhibitors Using Combined Computational Approach**. <i>ChemistryOpen</i> , 2022, 11, e202100248.	1.9	8
31	Virtual screen for repurposing approved and experimental drugs for candidate inhibitors of EBOLA virus infection. <i>F1000Research</i> , 0, 4, 34.	1.6	7
32	New in silico and conventional in vitro approaches to advance HIV drug discovery and design. <i>Expert Opinion on Drug Discovery</i> , 2013, 8, 83-92.	5.0	5
33	Feature-Based Classification of Amino Acid Substitutions outside Conserved Functional Protein Domains. <i>Scientific World Journal, The</i> , 2013, 2013, 1-10.	2.1	5
34	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. <i>BMC Bioinformatics</i> , 2016, 17, 139.	2.6	5
35	Prediction of the effectiveness of COVID-19 vaccine candidates. <i>F1000Research</i> , 0, 9, 365.	1.6	5
36	In Silico Screening of Natural Compounds for Candidates 5HT6 Receptor Antagonists against Alzheimerâ€“s Disease. <i>Molecules</i> , 2022, 27, 2626.	3.8	4

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37	Improving attrition rates in Ebola virus drug discovery. <i>Expert Opinion on Drug Discovery</i> , 2015, 10, 1025-1032.	5.0	3
38	In Silico Criterion for Prediction of Effects of p53 Gene Missense Mutations on p53-Mdm2 Feedback Loop. <i>Protein and Peptide Letters</i> , 2006, 13, 807-814.	0.9	2
39	In silico Therapeutics for Neurogenic Hypertension and Vasovagal Syncope. <i>Frontiers in Neuroscience</i> , 2016, 9, 520.	2.8	2
40	DiNGO: standalone application for Gene Ontology and Human Phenotype Ontology term enrichment analysis. <i>Bioinformatics</i> , 2020, 36, 1981-1982.	4.1	2
41	Tally-2.0: upgraded validator of tandem repeat detection in protein sequences. <i>Bioinformatics</i> , 2020, 36, 3260-3262.	4.1	2
42	Evolution of SARS-CoV-2 virus and assessment of the effectiveness of COVID-19 vaccine. <i>F1000Research</i> , 0, 10, 28.	1.6	2
43	A simple method for calculation of basic molecular properties of nutrients and their use as a criterion for a healthy diet. <i>F1000Research</i> , 2017, 6, 13.	1.6	2
44	Predicted Enhanced Human Propensity of Current Avian-Like H1N1 Swine Influenza Virus from China. <i>PLoS ONE</i> , 2016, 11, e0165451.	2.5	2
45	Alignment-free method for functional annotation of amino acid substitutions: Application on epigenetic factors involved in hematologic malignancies. <i>PLoS ONE</i> , 2021, 16, e0244948.	2.5	0
46	Role of genetic markers in sport and recreational physical activity. <i>FiziÄka Kultura</i> , 2016, 70, 5-13.	0.2	0
47	Simple Chemoinformatics Criterion Using Electron Donor-Acceptor Molecular Characteristics for Selection of Antibiotics Against Multi-Drug-Resistant Bacteria. <i>Discoveries</i> , 2016, 4, e64.	2.3	0