

Igor F Tsigelny

List of Publications by Year in descending order

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Version: 2024-02-01

53
papers

1,884
citations

304743

22
h-index

276875

41
g-index

58
all docs

58
docs citations

58
times ranked

3062
citing authors

#	ARTICLE	IF	CITATIONS
1	Interactions of large T-Antigen (LT) protein of polyomaviruses with p53 unfold their cancerogenic potential. <i>Journal of Biomolecular Structure and Dynamics</i> , 2022, 40, 5243-5252.	3.5	4
2	Potential SARS-CoV-2 Spike Protein-ACE2 Interface Inhibitors: Repurposing FDA-approved Drugs. <i>Journal of Exploratory Research in Pharmacology</i> , 2022, 7, 17-29.	0.4	5
3	Modeling of mutant superoxide dismutase 1 octamers with cross-linked disulfide bonds. <i>Journal of Molecular Modeling</i> , 2022, 28, 89.	1.8	0
4	<scp>Machineâ€learningâ€based</scp> virtual screening to repurpose drugs for treatment of <i>Candida albicans</i> infection. <i>Mycoses</i> , 2022, 65, 794-805.	4.0	5
5	Finding distinctions between oral cancer and periodontitis using saliva metabolites and machine learning. <i>Oral Diseases</i> , 2021, 27, 484-493.	3.0	32
6	Potential SARS-CoV-2 protease M^{pro} inhibitors: repurposing FDA-approved drugs. <i>Physical Biology</i> , 2021, 18, 025001.	1.8	4
7	Implications of viral infection in cancer development. <i>Biochimica Et Biophysica Acta: Reviews on Cancer</i> , 2021, 1876, 188622.	7.4	5
8	Machine Learning Strategies to Distinguish Oral Cancer from Periodontitis Using Salivary Metabolites. <i>Advances in Intelligent Systems and Computing</i> , 2021, , 511-526.	0.6	3
9	Targeting Epigenetic Regulators Using Machine Learning: Potential Sirtuin 2 Inhibitors. <i>Journal of Computational Biophysics and Chemistry</i> , 2021, 20, 841-851.	1.7	1
10	Deep-learning- and pharmacophore-based prediction of RAGE inhibitors. <i>Physical Biology</i> , 2020, 17, 036003.	1.8	11
11	Bovine leukemia virus relation to human breast cancer: Meta-analysis. <i>Microbial Pathogenesis</i> , 2020, 149, 104417.	2.9	12
12	Prediction of Premature Termination Codon Suppressing Compounds for Treatment of Duchenne Muscular Dystrophy Using Machine Learning. <i>Molecules</i> , 2020, 25, 3886.	3.8	1
13	Human Papillomavirus (HPV) Viral Proteins Substitute for the Impact of Somatic Mutations by Affecting Cancer-Related Genes: Meta-analysis and Perspectives. <i>Journal of Infectiology</i> , 2020, 3, 29-47.	0.8	2
14	Potential COVID-19 papain-like protease PL^{pro} inhibitors: repurposing FDA-approved drugs. <i>PeerJ</i> , 2020, 8, e9965.	2.0	44
15	Recognition of early and late stages of bladder cancer using metabolites and machine learning. <i>Metabolomics</i> , 2019, 15, 94.	3.0	34
16	Polycomb repressive 2 complexâ€™ Molecular mechanisms of function. <i>Protein Science</i> , 2019, 28, 1387-1399.	7.6	57
17	APOBEC-related mutagenesis and neo-peptide hydrophobicity: implications for response to immunotherapy. <i>Oncolmmunology</i> , 2019, 8, 1550341.	4.6	60
18	Artificial intelligence in drug combination therapy. <i>Briefings in Bioinformatics</i> , 2019, 20, 1434-1448.	6.5	60

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19	Next-Generation Sequencing of Circulating Tumor DNA Reveals Frequent Alterations in Advanced Hepatocellular Carcinoma. <i>Oncologist</i> , 2018, 23, 586-593.	3.7	75
20	Cripto stabilizes GRP78 on the cell membrane. <i>Protein Science</i> , 2018, 27, 653-661.	7.6	13
21	High expression of PD-1 ligands is associated with kataegis mutational signature and APOBEC3 alterations. <i>OncoImmunology</i> , 2017, 6, e1284719.	4.6	64
22	Multiple spatially related pharmacophores define small molecule inhibitors of OLIG2 in glioblastoma. <i>Oncotarget</i> , 2017, 8, 22370-22384.	1.8	23
23	A <i>de novo</i> compound targeting α -synuclein improves deficits in models of Parkinson's disease. <i>Brain</i> , 2016, 139, 3217-3236.	7.6	122
24	The copper transporter 1 (CTR1) is required to maintain the stability of copper transporter 2 (CTR2). <i>Metallomics</i> , 2015, 7, 1477-1487.	2.4	18
25	Hierarchical control of coherent gene clusters defines the molecular mechanisms of glioblastoma. <i>Molecular BioSystems</i> , 2015, 11, 1012-1028.	2.9	0
26	Development of a pharmacophore model for the catecholamine release-inhibitory peptide catestatin: Virtual screening and functional testing identify novel small molecule therapeutics of hypertension. <i>Bioorganic and Medicinal Chemistry</i> , 2013, 21, 5855-5869.	3.0	13
27	An All-Atom Model of the Structure of Human Copper Transporter 1. <i>Cell Biochemistry and Biophysics</i> , 2012, 63, 223-234.	1.8	40
28	Role of α -synuclein penetration into the membrane in the mechanisms of oligomer pore formation. <i>FEBS Journal</i> , 2012, 279, 1000-1013.	4.7	146
29	Conformational Changes of the Multispecific Transporter Organic Anion Transporter 1 (OAT1/SLC22A6) Suggests a Molecular Mechanism for Initial Stages of Drug and Metabolite Transport. <i>Cell Biochemistry and Biophysics</i> , 2011, 61, 251-259.	1.8	18
30	Elucidation of common pharmacophores from analysis of targeted metabolites transported by the multispecific drug transporter Organic anion transporter1 (Oat1). <i>Bioorganic and Medicinal Chemistry</i> , 2011, 19, 3320-3340.	3.0	14
31	Interaction of Organic Cations with Organic Anion Transporters. <i>Journal of Biological Chemistry</i> , 2009, 284, 31422-31430.	3.4	58
32	Role of Synucleins in Alzheimer's Disease. <i>Neurotoxicity Research</i> , 2009, 16, 306-317.	2.7	73
33	Mechanism of alpha-synuclein oligomerization and membrane interaction: theoretical approach to unstructured proteins studies. <i>Nanomedicine: Nanotechnology, Biology, and Medicine</i> , 2008, 4, 350-357.	3.3	48
34	MAPAS: a tool for predicting membrane-contacting protein surfaces. <i>Nature Methods</i> , 2008, 5, 119-119.	19.0	19
35	MODELING OF GLYCEROL-3-PHOSPHATE TRANSPORTER SUGGESTS A POTENTIAL 'TILT' MECHANISM INVOLVED IN ITS FUNCTION. <i>Journal of Bioinformatics and Computational Biology</i> , 2008, 06, 885-904.	0.8	12
36	Analysis of Metagene Portraits Reveals Distinct Transitions During Kidney Organogenesis. <i>Science Signaling</i> , 2008, 1, ra16.	3.6	28

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37	Mechanisms of Hybrid Oligomer Formation in the Pathogenesis of Combined Alzheimer's and Parkinson's Diseases. PLoS ONE, 2008, 3, e3135.	2.5	233
38	Dynamics of β -synuclein aggregation and inhibition of pore-like oligomer development by β ² -synuclein. FEBS Journal, 2007, 274, 1862-1877.	4.7	149
39	Structure-activity relationships and determinants of selectivity for congeners of the selective β ⁷ partial agonist 3-(2,4-dimethoxybenzylidene)-anabaseine (DMXBA or GTS21) with the ACh binding proteins (AChBPs). FASEB Journal, 2006, 20, A244.		0
40	Identification of Molecular Determinants That Modulate Trafficking of β ^{F508} CFTR, the Mutant ABC Transporter Associated With Cystic Fibrosis. Cell Biochemistry and Biophysics, 2005, 42, 041-054.	1.8	8
41	Complex Dynamics of Chaperone-Protein Interactions Under Cellular Stress. Cell Biochemistry and Biophysics, 2004, 40, 263-276.	1.8	12
42	Conformational preferences and activities of peptides from the catecholamine release-inhibitory (catestatin) region of chromogranin A. Regulatory Peptides, 2004, 118, 75-87.	1.9	27
43	Finding needles in haystacks: Reranking DOT results by using shape complementarity, cluster analysis, and biological information. Proteins: Structure, Function and Bioinformatics, 2003, 52, 33-40.	2.6	20
44	Hidden Markov Models-based system (HMMSPECTR) for detecting structural homologies on the basis of sequential information. Protein Engineering, Design and Selection, 2002, 15, 347-352.	2.1	6
45	Delineation of Selective Cyclic GMP-Dependent Protein Kinase β ¹ Substrate and Inhibitor Peptides Based on Combinatorial Peptide Libraries on Paper. , 1999, 82, 373-387.		52
46	600 ps Molecular dynamics reveals stable substructures and flexible hinge points in cAMP dependent protein kinase. , 1999, 50, 513-524.		49
47	Kinetic Analyses of Mutations in the Glycine-Rich Loop of cAMP-Dependent Protein Kinase. Biochemistry, 1998, 37, 7708-7715.	2.5	82
48	Solution structure of synthetic peptide inhibitor and substrate of cAMP-dependent protein kinase. A study by 2D ¹ H NMR and molecular dynamics. Chemical Biology and Drug Design, 1997, 49, 210-220.	1.1	7
49	Examination of an active-site electrostatic node in the cAMP-dependent protein kinase catalytic subunit. Protein Science, 1996, 5, 1316-1324.	7.6	36
50	Catalytic subunit of cAMP-dependent protein kinase: Electrostatic features and peptide recognition. Biopolymers, 1996, 39, 353-365.	2.4	15
51	Catalytic subunit of cAMP-dependent protein kinase: Electrostatic features and peptide recognition. Biopolymers, 1996, 39, 353-365.	2.4	9
52	Theoretical analysis of the structure of the peptide fasciculin and its docking to acetylcholinesterase. Protein Science, 1995, 4, 703-715.	7.6	23
53	Adding a positive charge at residue 46 of Drosophila alcohol dehydrogenase increases cofactor specificity for NADP ⁺ . FEBS Letters, 1994, 356, 81-85.	2.8	26