

Ian Walsh

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/4939230/publications.pdf>

Version: 2024-02-01

38
papers

2,613
citations

236612

25
h-index

301761

39
g-index

43
all docs

43
docs citations

43
times ranked

3498
citing authors

#	ARTICLE	IF	CITATIONS
1	Glycoinformatics Tools for Comprehensive Characterization of Glycans Enzymatically Released from Proteins. <i>Methods in Molecular Biology</i> , 2022, 2370, 3-23.	0.4	0
2	Semi-Automated Glycoproteomic Data Analysis of LC-MS Data Using GlycopeptideGraphMS in Process Development of Monoclonal Antibody Biologics. <i>Frontiers in Chemistry</i> , 2021, 9, 661406.	1.8	3
3	DOME: recommendations for supervised machine learning validation in biology. <i>Nature Methods</i> , 2021, 18, 1122-1127.	9.0	105
4	Clustering and curation of electropherograms: an efficient method for analyzing large cohorts of capillary electrophoresis glycomic profiles for bioprocessing operations. <i>Beilstein Journal of Organic Chemistry</i> , 2020, 16, 2087-2099.	1.3	5
5	Utility of Ion-Mobility Spectrometry for Deducing Branching of Multiply Charged Glycans and Glycopeptides in a High-Throughput Positive ion LC-FLR-IMS-MS Workflow. <i>Analytical Chemistry</i> , 2020, 92, 15323-15335.	3.2	30
6	Experimentally Determined Long Intrinsically Disordered Protein Regions Are Now Abundant in the Protein Data Bank. <i>International Journal of Molecular Sciences</i> , 2020, 21, 4496.	1.8	25
7	GlycanAnalyzer: software for automated interpretation of <i>N</i> -glycan profiles after exoglycosidase digestions. <i>Bioinformatics</i> , 2019, 35, 688-690.	1.8	17
8	Combining Glucose Units, <i>m/z</i> , and Collision Cross Section Values: Multiattribute Data for Increased Accuracy in Automated Glycosphingolipid Glycan Identifications and Its Application in Triple Negative Breast Cancer. <i>Analytical Chemistry</i> , 2019, 91, 9078-9085.	3.2	14
9	A Robust and Versatile Automated Glycoanalytical Technology for Serum Antibodies and Acute Phase Proteins: Ovarian Cancer Case Study. <i>Molecular and Cellular Proteomics</i> , 2019, 18, 2191-2206.	2.5	18
10	GlycoStore: a database of retention properties for glycan analysis. <i>Bioinformatics</i> , 2018, 34, 3231-3232.	1.8	77
11	Plasma N-glycans in colorectal cancer risk. <i>Scientific Reports</i> , 2018, 8, 8655.	1.6	57
12	FELLS: fast estimator of latent local structure. <i>Bioinformatics</i> , 2017, 33, 1889-1891.	1.8	72
13	Serum <i>N</i> -glycome alterations in breast cancer during multimodal treatment and follow-up. <i>Molecular Oncology</i> , 2017, 11, 1361-1379.	2.1	32
14	Quantitative profiling of glycans and glycopeptides: an informatics™ perspective. <i>Current Opinion in Structural Biology</i> , 2016, 40, 70-80.	2.6	29
15	Correct machine learning on protein sequences: a peer-reviewing perspective. <i>Briefings in Bioinformatics</i> , 2016, 17, 831-840.	3.2	53
16	MobiDB 2.0: an improved database of intrinsically disordered and mobile proteins. <i>Nucleic Acids Research</i> , 2015, 43, D315-D320.	6.5	177
17	Comprehensive large-scale assessment of intrinsic protein disorder. <i>Bioinformatics</i> , 2015, 31, 201-208.	1.8	154
18	Toward an accurate prediction of inter-residue distances in proteins using 2D recursive neural networks. <i>BMC Bioinformatics</i> , 2014, 15, 6.	1.2	51

#	ARTICLE	IF	CITATIONS
19	RUBI: rapid proteomic-scale prediction of lysine ubiquitination and factors influencing predictor performance. <i>Amino Acids</i> , 2014, 46, 853-862.	1.2	24
20	RepeatsDB: a database of tandem repeat protein structures. <i>Nucleic Acids Research</i> , 2014, 42, D352-D357.	6.5	53
21	PASTA 2.0: an improved server for protein aggregation prediction. <i>Nucleic Acids Research</i> , 2014, 42, W301-W307.	6.5	349
22	NeEMO: a method using residue interaction networks to improve prediction of protein stability upon mutation. <i>BMC Genomics</i> , 2014, 15, S7.	1.2	83
23	Analysis and consensus of currently available intrinsic protein disorder annotation sources in the MobiDB database. <i>BMC Bioinformatics</i> , 2013, 14, S3.	1.2	30
24	PANADA: Protein Association Network Annotation, Determination and Analysis. <i>PLoS ONE</i> , 2013, 8, e78383.	1.1	7
25	ESpritz: accurate and fast prediction of protein disorder. <i>Bioinformatics</i> , 2012, 28, 503-509.	1.8	445
26	RAPHAEL: recognition, periodicity and insertion assignment of solenoid protein structures. <i>Bioinformatics</i> , 2012, 28, 3257-3264.	1.8	27
27	Blues server: electrostatic properties of wild-type and mutated protein structures. <i>Bioinformatics</i> , 2012, 28, 2189-2190.	1.8	72
28	MobiDB: a comprehensive database of intrinsic protein disorder annotations. <i>Bioinformatics</i> , 2012, 28, 2080-2081.	1.8	142
29	RING: networking interacting residues, evolutionary information and energetics in protein structures. <i>Bioinformatics</i> , 2011, 27, 2003-2005.	1.8	116
30	CSpritz: accurate prediction of protein disorder segments with annotation for homology, secondary structure and linear motifs. <i>Nucleic Acids Research</i> , 2011, 39, W190-W196.	6.5	77
31	In Silico Protein Motif Discovery and Structural Analysis. <i>Methods in Molecular Biology</i> , 2011, 760, 341-353.	0.4	2
32	MOBI: a web server to define and visualize structural mobility in NMR protein ensembles. <i>Bioinformatics</i> , 2010, 26, 2916-2917.	1.8	35
33	Ab initio and homology based prediction of protein domains by recursive neural networks. <i>BMC Bioinformatics</i> , 2009, 10, 195.	1.2	13
34	Ab initio and template-based prediction of multi-class distance maps by two-dimensional recursive neural networks. <i>BMC Structural Biology</i> , 2009, 9, 5.	2.3	44
35	Recursive Neural Networks for Undirected Graphs for Learning Molecular Endpoints. <i>Lecture Notes in Computer Science</i> , 2009, , 391-403.	1.0	6
36	LONG-RANGE INFORMATION AND PHYSICALITY CONSTRAINTS IMPROVE PREDICTED PROTEIN CONTACT MAPS. <i>Journal of Bioinformatics and Computational Biology</i> , 2008, 06, 1001-1020.	0.3	3

#	ARTICLE	IF	CITATIONS
37	A two-stage approach for improved prediction of residue contact maps. BMC Bioinformatics, 2006, 7, 180.	1.2	74
38	Distill: a suite of web servers for the prediction of one-, two- and three-dimensional structural features of proteins. BMC Bioinformatics, 2006, 7, 402.	1.2	85