

Rovshan G Sadygov

List of Publications by Year in descending order

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60
papers

5,681
citations

201674

27
h-index

123424

61
g-index

61
all docs

61
docs citations

61
times ranked

6949
citing authors

#	ARTICLE	IF	CITATIONS
1	OUP accepted manuscript. Briefings in Bioinformatics, 2022, , .	6.5	6
2	Using Heavy Mass Isotopomers for Protein Turnover in Heavy Water Metabolic Labeling. Journal of Proteome Research, 2021, 20, 2035-2041.	3.7	7
3	A novel estimator of the interaction matrix in Graphical Gaussian Model of omics data using the entropy of non-equilibrium systems. Bioinformatics, 2021, 37, 837-844.	4.1	2
4	Partial Isotope Profiles Are Sufficient for Protein Turnover Analysis Using Closed-Form Equations of Mass Isotopomer Dynamics. Analytical Chemistry, 2020, 92, 14747-14753.	6.5	12
5	High-Resolution Mass Spectrometry for In Vivo Proteome Dynamics using Heavy Water Metabolic Labeling. International Journal of Molecular Sciences, 2020, 21, 7821.	4.1	5
6	Timepoint Selection Strategy for In Vivo Proteome Dynamics from Heavy Water Metabolic Labeling and LC-MS. Journal of Proteome Research, 2020, 19, 2105-2112.	3.7	8
7	Proteome dynamics from heavy water metabolic labeling and peptide tandem mass spectrometry. International Journal of Mass Spectrometry, 2019, 445, 116194.	1.5	8
8	Calculation of the Protein Turnover Rate Using the Number of Incorporated 2H Atoms and Proteomics Analysis of a Single Labeled Sample. Analytical Chemistry, 2019, 91, 14340-14351.	6.5	9
9	Another look at matrix correlations. Bioinformatics, 2019, 35, 4748-4753.	4.1	2
10	Poisson Model To Generate Isotope Distribution for Biomolecules. Journal of Proteome Research, 2018, 17, 751-758.	3.7	15
11	d2ome, Software for in Vivo Protein Turnover Analysis Using Heavy Water Labeling and LC-MS, Reveals Alterations of Hepatic Proteome Dynamics in a Mouse Model of NAFLD. Journal of Proteome Research, 2018, 17, 3740-3748.	3.7	42
12	Hepatic Mitochondrial Defects in a Nonalcoholic Fatty Liver Disease Mouse Model Are Associated with Increased Degradation of Oxidative Phosphorylation Subunits. Molecular and Cellular Proteomics, 2018, 17, 2371-2386.	3.8	59
13	PPARgamma agonists rescue increased phosphorylation of FGF14 at S226 in the Tg2576 mouse model of Alzheimer's disease. Experimental Neurology, 2017, 295, 1-17.	4.1	35
14	Increased serotransferrin and ceruloplasmin turnover in diet-controlled patients with type 2 diabetes. Free Radical Biology and Medicine, 2017, 113, 461-469.	2.9	41
15	Predicting the protein half-life in tissue from its cellular properties. PLoS ONE, 2017, 12, e0180428.	2.5	11
16	Gaussian Process Modeling of Protein Turnover. Journal of Proteome Research, 2016, 15, 2115-2122.	3.7	25
17	Proteome Dynamics Reveals Pro-Inflammatory Remodeling of Plasma Proteome in a Mouse Model of NAFLD. Journal of Proteome Research, 2016, 15, 3388-3404.	3.7	15
18	Integrative proteomic analysis reveals reprogramming tumor necrosis factor signaling in epithelial mesenchymal transition. Journal of Proteomics, 2016, 148, 126-138.	2.4	29

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19	Tracer-based estimates of protein flux in cases of incomplete product renewal: evidence and implications of heterogeneity in collagen turnover. <i>American Journal of Physiology - Endocrinology and Metabolism</i> , 2015, 309, E115-E121.	3.5	16
20	Using SEQUEST with Theoretically Complete Sequence Databases. <i>Journal of the American Society for Mass Spectrometry</i> , 2015, 26, 1858-1864.	2.8	10
21	Mixed-effects model of epithelialâ€mesenchymal transition reveals rewiring of signaling networks. <i>Cellular Signalling</i> , 2015, 27, 1413-1425.	3.6	25
22	Current Bioinformatics Challenges in Proteome Dynamics using Heavy Water-based Metabolic Labeling. <i>Journal of Data Mining in Genomics & Proteomics</i> , 2014, 05, e112.	0.5	2
23	The Cancer Drug Tamoxifen: A Potential Therapeutic Treatment for Spinal Cord Injury. <i>Journal of Neurotrauma</i> , 2014, 31, 268-283.	3.4	37
24	Cognitive Enhancing Treatment with a PPAR β Agonist Normalizes Dentate Granule Cell Presynaptic Function in Tg2576 APP Mice. <i>Journal of Neuroscience</i> , 2014, 34, 1028-1036.	3.6	48
25	Cardiac mitochondrial proteome dynamics with heavy water reveals stable rate of mitochondrial protein synthesis in heart failure despite decline in mitochondrial oxidative capacity. <i>Journal of Molecular and Cellular Cardiology</i> , 2014, 75, 88-97.	1.9	34
26	Use of singular value decomposition analysis to differentiate phosphorylated precursors in strong cation exchange fractions. <i>Electrophoresis</i> , 2014, 35, 3498-3503.	2.4	5
27	Use of Theoretical Peptide Distributions in Phosphoproteome Analysis. <i>Journal of Proteome Research</i> , 2013, 12, 3207-3214.	3.7	1
28	Assessment of cardiac proteome dynamics with heavy water: slower protein synthesis rates in interfibrillar than subsarcolemmal mitochondria. <i>American Journal of Physiology - Heart and Circulatory Physiology</i> , 2013, 304, H1201-H1214.	3.2	66
29	Detection of Structural and Metabolic Changes in Traumatically Injured Hippocampus by Quantitative Differential Proteomics. <i>Journal of Neurotrauma</i> , 2013, 30, 775-788.	3.4	32
30	High Mass Accuracy Phosphopeptide Identification Using Tandem Mass Spectra. <i>International Journal of Proteomics</i> , 2012, 2012, 1-5.	2.0	4
31	Cognitive Enhancement with Rosiglitazone Links the Hippocampal PPAR β and ERK MAPK Signaling Pathways. <i>Journal of Neuroscience</i> , 2012, 32, 16725-16735.	3.6	99
32	Improved Mass Defect Model for Theoretical Tryptic Peptides. <i>Analytical Chemistry</i> , 2012, 84, 3026-3032.	6.5	11
33	Generalized linear and mixed models for label-free shotgun proteomics. <i>Statistics and Its Interface</i> , 2012, 5, 89-98.	0.3	6
34	Examining Troughs in the Mass Distribution of All Theoretically Possible Tryptic Peptides. <i>Journal of Proteome Research</i> , 2011, 10, 4150-4157.	3.7	25
35	SVM Model for Quality Assessment of Medium Resolution Mass Spectra from ¹⁸ O-Water Labeling Experiments. <i>Journal of Proteome Research</i> , 2011, 10, 2095-2103.	3.7	7
36	A parallel method for enumerating amino acid compositions and masses of all theoretical peptides. <i>BMC Bioinformatics</i> , 2011, 12, 432.	2.6	11

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37	Measuring protein synthesis using metabolic 2H labeling, high-resolution mass spectrometry, and an algorithm. <i>Analytical Biochemistry</i> , 2011, 412, 47-55.	2.4	64
38	Bioinformatics Tools for Mass Spectrometry-Based High-Throughput Quantitative Proteomics Platforms. <i>Current Proteomics</i> , 2011, 8, 125-137.	0.3	6
39	Comparison of Programmatic Approaches for Efficient Accessing to mzML Files. <i>Journal of Data Mining in Genomics & Proteomics</i> , 2011, 02, .	0.5	4
40	Altered Retinoic Acid Metabolism in Diabetic Mouse Kidney Identified by 18O Isotopic Labeling and 2D Mass Spectrometry. <i>PLoS ONE</i> , 2010, 5, e11095.	2.5	45
41	Using Power Spectrum Analysis to Evaluate 18O-Water Labeling Data Acquired from Low Resolution Mass Spectrometers. <i>Journal of Proteome Research</i> , 2010, 9, 4306-4312.	3.7	23
42	A New Probabilistic Database Search Algorithm for ETD Spectra. <i>Journal of Proteome Research</i> , 2009, 8, 3198-3205.	3.7	33
43	Charger: A Combination of Signal Processing and Statistical Learning Algorithms for Precursor Charge-State Determination from Electron-Transfer Dissociation Spectra. <i>Analytical Chemistry</i> , 2008, 80, 376-386.	6.5	18
44	Assigning in vivo carbamylation and acetylation in human lens proteins using tandem mass spectrometry and database searching. <i>International Journal of Mass Spectrometry</i> , 2007, 259, 161-173.	1.5	11
45	ChromAlign: A Two-Step Algorithmic Procedure for Time Alignment of Three-Dimensional LC-MS Chromatographic Surfaces. <i>Analytical Chemistry</i> , 2006, 78, 8207-8217.	6.5	101
46	Central Limit Theorem as an Approximation for Intensity-Based Scoring Function. <i>Analytical Chemistry</i> , 2006, 78, 89-95.	6.5	31
47	Large-scale database searching using tandem mass spectra: Looking up the answer in the back of the book. <i>Nature Methods</i> , 2004, 1, 195-202.	19.0	375
48	MS1, MS2, and SQT—three unified, compact, and easily parsed file formats for the storage of shotgun proteomic spectra and identifications. <i>Rapid Communications in Mass Spectrometry</i> , 2004, 18, 2162-2168.	1.5	350
49	Statistical Models for Protein Validation Using Tandem Mass Spectral Data and Protein Amino Acid Sequence Databases. <i>Analytical Chemistry</i> , 2004, 76, 1664-1671.	6.5	127
50	A Model for Random Sampling and Estimation of Relative Protein Abundance in Shotgun Proteomics. <i>Analytical Chemistry</i> , 2004, 76, 4193-4201.	6.5	2,280
51	A Hypergeometric Probability Model for Protein Identification and Validation Using Tandem Mass Spectral Data and Protein Sequence Databases. <i>Analytical Chemistry</i> , 2003, 75, 3792-3798.	6.5	198
52	A Correlation Algorithm for the Automated Quantitative Analysis of Shotgun Proteomics Data. <i>Analytical Chemistry</i> , 2003, 75, 6912-6921.	6.5	276
53	Shotgun identification of protein modifications from protein complexes and lens tissue. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2002, 99, 7900-7905.	7.1	571
54	Code Developments to Improve the Efficiency of Automated MS/MS Spectra Interpretation. <i>Journal of Proteome Research</i> , 2002, 1, 211-215.	3.7	198

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55	Dynamics of primary charge separation in bacterial photosynthesis using the multilevel Redfield-Davies secular approach. International Journal of Quantum Chemistry, 2002, 87, 254-263.	2.0	3
56	Unusual conical intersections in the Jahn-Teller effect: The electronically excited states of Li ₃ . Journal of Chemical Physics, 1999, 110, 3639-3642.	3.0	38
57	On the adiabatic to diabatic states transformation in the presence of a conical intersection: A most diabatic basis from the solution to a Poisson's equation. I. Journal of Chemical Physics, 1998, 109, 20-25.	3.0	86
58	Electronic structure aspects of the spin-forbidden reaction CH ₃ (X ² A ₂) + N(4S) → HCN(X ¹ Σ ⁺) + H ₂ (X ¹ g _g). Journal of Chemical Physics, 1997, 107, 4994-4999.	3.0	28
59	Resonances in the predissociation of the A ² Π _g state of MgBr. Journal of Chemical Physics, 1997, 106, 4091-4101.	3.0	13
60	A theoretical study of the structure and energetics of stacked dimers of polycyclic aromatic hydrocarbons. Application of INDO 1/S method to singlet excimers of naphthalene and phenanthrene. Chemical Physics Letters, 1994, 225, 441-447.	2.6	30