

Andris Jankevics

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/8184290/publications.pdf>

Version: 2024-02-01

28
papers

2,508
citations

430754

18
h-index

526166

27
g-index

35
all docs

35
docs citations

35
times ranked

4167
citing authors

#	ARTICLE	IF	CITATIONS
1	Characterization of Monophasic Solvent-Based Tissue Extractions for the Detection of Polar Metabolites and Lipids Applying Ultrahigh-Performance Liquid Chromatography–Mass Spectrometry Clinical Metabolic Phenotyping Assays. <i>Journal of Proteome Research</i> , 2021, 20, 831-840.	1.8	20
2	mzMLb: A Future-Proof Raw Mass Spectrometry Data Format Based on Standards-Compliant mzML and Optimized for Speed and Storage Requirements. <i>Journal of Proteome Research</i> , 2021, 20, 172-183.	1.8	12
3	An improved strategy for analysis of lipid molecules utilising a reversed phase C30 UHPLC column and scheduled MS/MS acquisition. <i>Talanta</i> , 2021, 229, 122262.	2.9	8
4	Assessment of human plasma and urine sample preparation for reproducible and high-throughput UHPLC-MS clinical metabolic phenotyping. <i>Analyst</i> , The, 2020, 145, 6511-6523.	1.7	28
5	Multi-Omics Analysis of Diabetic Heart Disease in the db/db Model Reveals Potential Targets for Treatment by a Longevity-Associated Gene. <i>Cells</i> , 2020, 9, 1283.	1.8	11
6	Metabolic engineering against the arginine microenvironment enhances CAR-T cell proliferation and therapeutic activity. <i>Blood</i> , 2020, 136, 1155-1160.	0.6	84
7	Metabolic characterisation of disturbances in the APOC3/triglyceride-rich lipoprotein pathway through sample-based recall by genotype. <i>Metabolomics</i> , 2020, 16, 69.	1.4	3
8	RankProd 2.0: a refactored bioconductor package for detecting differentially expressed features in molecular profiling datasets. <i>Bioinformatics</i> , 2017, 33, 2774-2775.	1.8	113
9	Chapter 7. Algorithms for MS1-Based Quantitation. <i>New Developments in Mass Spectrometry</i> , 2016, , 133-154.	0.2	0
10	MetAssign: probabilistic annotation of metabolites from LC–MS data using a Bayesian clustering approach. <i>Bioinformatics</i> , 2014, 30, 2764-2771.	1.8	63
11	Metabolic adaptations of <i>Cryptosporidium parvum</i> in relation to differentiation, drug resistance, and drug pressure. <i>Molecular Microbiology</i> , 2013, 90, 428-442.	1.2	48
12	LC-MS METABOLOMICS FROM STUDY DESIGN TO DATA-ANALYSIS – USING A VERSATILE PATHOGEN AS A TEST CASE. <i>Computational and Structural Biotechnology Journal</i> , 2013, 4, e201301002.	1.9	39
13	Bioanalysis Young Investigator Award 2013. <i>Bioanalysis</i> , 2013, 5, 1479-1484.	0.6	0
14	mzMatch–ISO: an R tool for the annotation and relative quantification of isotope-labelled mass spectrometry data. <i>Bioinformatics</i> , 2013, 29, 281-283.	1.8	91
15	Bioanalysis Young Investigator: Announcing our finalists!. <i>Bioanalysis</i> , 2013, 5, 1963-1964.	0.6	1
16	Metabolomics for Secondary Metabolite Research. <i>Metabolites</i> , 2013, 3, 1076-1083.	1.3	59
17	IDEOM: an Excel interface for analysis of LC–MS-based metabolomics data. <i>Bioinformatics</i> , 2012, 28, 1048-1049.	1.8	307
18	Serine is a natural ligand and allosteric activator of pyruvate kinase M2. <i>Nature</i> , 2012, 491, 458-462.	13.7	519

#	ARTICLE	IF	CITATIONS
19	Stable Isotope-Assisted Metabolomics for Network-Wide Metabolic Pathway Elucidation. <i>Analytical Chemistry</i> , 2012, 84, 8442-8447.	3.2	132
20	Separating the wheat from the chaff: a prioritisation pipeline for the analysis of metabolomics datasets. <i>Metabolomics</i> , 2012, 8, 29-36.	1.4	50
21	Metabolomics methods for the synthetic biology of secondary metabolism. <i>FEBS Letters</i> , 2012, 586, 2177-2183.	1.3	63
22	PeakML/mzMatch: A File Format, Java Library, R Library, and Tool-Chain for Mass Spectrometry Data Analysis. <i>Analytical Chemistry</i> , 2011, 83, 2786-2793.	3.2	305
23	Exploring the metabolic state of microorganisms using metabolomics. <i>Bioanalysis</i> , 2011, 3, 2443-2458.	0.6	19
24	Toward Global Metabolomics Analysis with Hydrophilic Interaction Liquid Chromatography–Mass Spectrometry: Improved Metabolite Identification by Retention Time Prediction. <i>Analytical Chemistry</i> , 2011, 83, 8703-8710.	3.2	326
25	Metabolomic analysis of a synthetic metabolic switch in <i>Streptomyces coelicolor</i> . <i>Proteomics</i> , 2011, 11, 4622-4631.	1.3	20
26	Towards an unbiased metabolic profiling of protozoan parasites: optimisation of a <i>Leishmania</i> sampling protocol for HILIC-orbitrap analysis. <i>Analytical and Bioanalytical Chemistry</i> , 2010, 398, 2059-2069.	1.9	48
27	Metabolomics to Unveil and Understand Phenotypic Diversity between Pathogen Populations. <i>PLoS Neglected Tropical Diseases</i> , 2010, 4, e904.	1.3	91
28	Metabolomic studies of experimental diabetic urine samples by ¹ H NMR spectroscopy and LC/MS method. <i>Chemometrics and Intelligent Laboratory Systems</i> , 2009, 97, 11-17.	1.8	19