## Joaquim Jaumot

## List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

69 3,797 29 61 g-index

70 4,291 5.5 5.86 ext. papers ext. citations avg, IF L-index

| #  | Paper   | IF                 | Citations |
|----|---|--------------------|-----------|
| 69 | Lipidomic analysis of single and combined effects of polyethylene microplastics and polychlorinated biphenyls on human hepatoma cells. <i>Journal of Hazardous Materials</i> , <b>2022</b> , 421, 126777  | 12.8               | 8         |
| 68 | Adverse Effects of Arsenic Uptake in Rice Metabolome and Lipidome Revealed by Untargeted Liquid Chromatography Coupled to Mass Spectrometry (LC-MS) and Regions of Interest Multivariate Curve Resolution. <i>Separations</i> , <b>2022</b> , 9, 79 | 3.1                | О         |
| 67 | Comparison of Multivariate ANOVA-Based Approaches for the Determination of Relevant Variables in Experimentally Designed Metabolomic Studies. <i>Molecules</i> , <b>2022</b> , 27, 3304   | 4.8                | O         |
| 66 | Quantification strategies for two-dimensional liquid chromatography datasets using regions of interest and multivariate curve resolution approaches. <i>Talanta</i> , <b>2022</b> , 247, 123586   | 6.2                |           |
| 65 | Two-Dimensional Liquid Chromatography in Metabolomics and Lipidomics. <i>Neuromethods</i> , <b>2021</b> , 25-47   | 0.4                | 3         |
| 64 | Untangling comprehensive two-dimensional liquid chromatography data sets using regions of interest and multivariate curve resolution approaches. <i>TrAC - Trends in Analytical Chemistry</i> , <b>2021</b> , 137, 116207                           | 14.6               | 6         |
| 63 | MSroi: A pre-processing tool for mass spectrometry-based studies. <i>Chemometrics and Intelligent Laboratory Systems</i> , <b>2021</b> , 215, 104333  | 3.8                | 5         |
| 62 | Mass Spectrometry Imaging: Chemometric Data Analysis <b>2020</b> , 381-394  |                    |           |
| 61 | Interaction of Environmental Pollutants with Microplastics: A Critical Review of Sorption Factors, Bioaccumulation and Ecotoxicological Effects. <i>Toxics</i> , <b>2020</b> , 8,   | 4.7                | 45        |
| 60 | Chemometrics in comprehensive two-dimensional liquid chromatography: A study of the data structure and its multilinear behavior. <i>Chemometrics and Intelligent Laboratory Systems</i> , <b>2020</b> , 201, 1040                                   | ා <sub>ල්ල</sub> 8 | 6         |
| 59 | Partial characterization of the lipidome of the cold-water scallop, Chlamys islandica. <i>Environmental Science and Pollution Research</i> , <b>2020</b> , 27, 1475-1484  | 5.1                | 2         |
| 58 | ROIMCR: a powerful analysis strategy for LC-MS metabolomic datasets. <i>BMC Bioinformatics</i> , <b>2019</b> , 20, 256  | 3.6                | 31        |
| 57 | Untargeted lipidomic evaluation of hydric and heat stresses on rice growth. <i>Journal of Chromatography B: Analytical Technologies in the Biomedical and Life Sciences</i> , <b>2019</b> , 1104, 148-156   | 3.2                | 7         |
| 56 | Assessment of endocrine disruptors effects on zebrafish (Danio rerio) embryos by untargeted LC-HRMS metabolomic analysis. <i>Science of the Total Environment</i> , <b>2018</b> , 635, 156-166  | 10.2               | 55        |
| 55 | Chemometric Strategies for Peak Detection and Profiling from Multidimensional Chromatography. <i>Proteomics</i> , <b>2018</b> , 18, e1700327  | 4.8                | 13        |
| 54 | A chemometric approach for characterization of serum transthyretin in familial amyloidotic polyneuropathy type I (FAP-I) by electrospray ionization-ion mobility mass spectrometry. <i>Talanta</i> , <b>2018</b> , 181, 87-94                       | 6.2                | 3         |
| 53 | An untargeted lipidomic strategy combining comprehensive two-dimensional liquid chromatography and chemometric analysis. <i>Journal of Chromatography A</i> , <b>2018</b> , 1568, 80-90   | 4.5                | 25        |

## (2015-2018)

| 52 | Compression and Resolution Tools for the Analysis of Untargeted Metabolomic Data. <i>Comprehensive Analytical Chemistry</i> , <b>2018</b> , 82, 337-368   | 1.9  | 1   |
|----|---|------|-----|
| 51 | Identification of antihypertensive peptides in nutraceuticals by capillary electrophoresis-mass spectrometry. <i>Journal of Chromatography A</i> , <b>2018</b> , 1579, 129-137  | 4.5  | 10  |
| 50 | Introduction to the Data Analysis Relevance in the Omic Era. <i>Comprehensive Analytical Chemistry</i> , <b>2018</b> , 1-12   | 1.9  | 1   |
| 49 | Chemometric evaluation of hydrophilic interaction liquid chromatography stationary phases: resolving complex mixtures of metabolites. <i>Analytical Methods</i> , <b>2017</b> , 9, 774-785  | 3.2  | 8   |
| 48 | The human mitochondrial transcription factor A is a versatile G-quadruplex binding protein. <i>Scientific Reports</i> , <b>2017</b> , 7, 43992  | 4.9  | 29  |
| 47 | Metabolomic analysis of the effects of cadmium and copper treatment in Oryza sativa L. using untargeted liquid chromatography coupled to high resolution mass spectrometry and all-ion fragmentation. <i>Metallomics</i> , <b>2017</b> , 9, 660-675 | 4.5  | 29  |
| 46 | Knowledge integration strategies for untargeted metabolomics based on MCR-ALS analysis of CE-MS and LC-MS data. <i>Analytica Chimica Acta</i> , <b>2017</b> , 978, 10-23  | 6.6  | 35  |
| 45 | Untargeted Comprehensive Two-Dimensional Liquid Chromatography Coupled with High-Resolution Mass Spectrometry Analysis of Rice Metabolome Using Multivariate Curve Resolution. <i>Analytical Chemistry</i> , <b>2017</b> , 89, 7675-7683            | 7.8  | 62  |
| 44 | Analysis of multiple mass spectrometry images from different Phaseolus vulgaris samples by multivariate curve resolution. <i>Talanta</i> , <b>2017</b> , 175, 557-565   | 6.2  | 14  |
| 43 | Metabolic disruption of zebrafish (Danio rerio) embryos by bisphenol A. An integrated metabolomic and transcriptomic approach. <i>Environmental Pollution</i> , <b>2017</b> , 231, 22-36  | 9.3  | 47  |
| 42 | Modelling of Hydrophilic Interaction Liquid Chromatography Stationary Phases Using Chemometric Approaches. <i>Metabolites</i> , <b>2017</b> , 7,  | 5.6  | 6   |
| 41 | Data analysis strategies for targeted and untargeted LC-MS metabolomic studies: Overview and workflow. <i>TrAC - Trends in Analytical Chemistry</i> , <b>2016</b> , 82, 425-442   | 14.6 | 172 |
| 40 | Metabolic profiling for the identification of Huntington biomarkers by on-line solid-phase extraction capillary electrophoresis mass spectrometry combined with advanced data analysis tools. <i>Electrophoresis</i> , <b>2016</b> , 37, 795-808    | 3.6  | 26  |
| 39 | Chemometric analysis of comprehensive LCIIC-MS data: Resolution of triacylglycerol structural isomers in corn oil. <i>Talanta</i> , <b>2016</b> , 160, 624-635  | 6.2  | 27  |
| 38 | Compression strategies for the chemometric analysis of mass spectrometry imaging data. <i>Journal of Chemometrics</i> , <b>2016</b> , 30, 575-588   | 1.6  | 24  |
| 37 | Lipidomic data analysis: tutorial, practical guidelines and applications. <i>Analytica Chimica Acta</i> , <b>2015</b> , 885, 1-16   | 6.6  | 72  |
| 36 | qRT-PCR evaluation of the transcriptional response of zebra mussel to heavy metals. <i>BMC Genomics</i> , <b>2015</b> , 16, 354   | 4.5  | 6   |
| 35 | Phenotypic malignant changes and untargeted lipidomic analysis of long-term exposed prostate cancer cells to endocrine disruptors. <i>Environmental Research</i> , <b>2015</b> , 140, 18-31   | 7.9  | 29  |

| 34 | Evaluation of changes induced in rice metabolome by Cd and Cu exposure using LC-MS with XCMS and MCR-ALS data analysis strategies. <i>Analytical and Bioanalytical Chemistry</i> , <b>2015</b> , 407, 8835-47   | 4.4  | 32  |
|----|---|------|-----|
| 33 | Epithelial-to-mesenchymal transition involves triacylglycerol accumulation in DU145 prostate cancer cells. <i>Molecular BioSystems</i> , <b>2015</b> , 11, 3397-406   |      | 31  |
| 32 | Potential use of multivariate curve resolution for the analysis of mass spectrometry images. <i>Analyst, The</i> , <b>2015</b> , 140, 837-46  | 5    | 34  |
| 31 | MCR-ALS GUI 2.0: New features and applications. <i>Chemometrics and Intelligent Laboratory Systems</i> , <b>2015</b> , 140, 1-12  | 3.8  | 466 |
| 30 | Combination of CE-MS and advanced chemometric methods for high-throughput metabolic profiling. <i>Electrophoresis</i> , <b>2015</b> , 36, 2324-2335   | 3.6  | 22  |
| 29 | Vibrational spectroscopic image analysis of biological material using multivariate curve resolution-alternating least squares (MCR-ALS). <i>Nature Protocols</i> , <b>2015</b> , 10, 217-40   | 18.8 | 190 |
| 28 | Multivariate Curve Resolution (MCR). Solving the mixture analysis problem. <i>Analytical Methods</i> , <b>2014</b> , 6, 4964-4976   | 3.2  | 340 |
| 27 | Modeling strategies for pharmaceutical blend monitoring and end-point determination by near-infrared spectroscopy. <i>International Journal of Pharmaceutics</i> , <b>2014</b> , 473, 219-31  | 6.5  | 26  |
| 26 | Solution equilibria of cytosine- and guanine-rich sequences near the promoter region of the n-myc gene that contain stable hairpins within lateral loops. <i>Biochimica Et Biophysica Acta - General Subjects</i> , <b>2014</b> , 1840, 41-52         | 4    | 31  |
| 25 | Blending process modeling and control by multivariate curve resolution. <i>Talanta</i> , <b>2013</b> , 117, 492-504   | 6.2  | 27  |
| 24 | Combination of chromatographic and chemometric methods to study the interactions between DNA strands. <i>Analytica Chimica Acta</i> , <b>2012</b> , 722, 34-42  | 6.6  | 9   |
| 23 | Porphyrin binding mechanism is altered by protonation at the loops in G-quadruplex DNA formed near the transcriptional activation site of the human c-kit gene. <i>Biochimica Et Biophysica Acta - General Subjects</i> , <b>2012</b> , 1820, 1987-96 | 4    | 19  |
| 22 | Experimental methods for studying the interactions between G-quadruplex structures and ligands. <i>Current Pharmaceutical Design</i> , <b>2012</b> , 18, 1900-16  | 3.3  | 97  |
| 21 | Influence of pH, temperature and the cationic porphyrin TMPyP4 on the stability of the i-motif formed by the 5V(C3TA2)4-3Vsequence of the human telomere. <i>International Journal of Biological Macromolecules</i> , <b>2011</b> , 49, 729-36        | 7.9  | 42  |
| 20 | Chemical equilibria studies using multivariate analysis methods. <i>Analytical and Bioanalytical Chemistry</i> , <b>2011</b> , 399, 1983-97   | 4.4  | 23  |
| 19 | Using principal component analysis to find correlations between loop-related and thermodynamic variables for G-quadruplex-forming sequences. <i>Biochimie</i> , <b>2010</b> , 92, 1016-23   | 4.6  | 6   |
| 18 | Application of multivariate curve resolution to the analysis of yeast genome-wide screens. <i>Chemometrics and Intelligent Laboratory Systems</i> , <b>2010</b> , 104, 53-64  | 3.8  | 23  |
| 17 | MCR-BANDS: A user friendly MATLAB program for the evaluation of rotation ambiguities in Multivariate Curve Resolution. <i>Chemometrics and Intelligent Laboratory Systems</i> , <b>2010</b> , 103, 96-107   | 3.8  | 157 |

## LIST OF PUBLICATIONS

| 16 | MCRC software: A tool for chemometric analysis of two-way chromatographic data. <i>Chemometrics and Intelligent Laboratory Systems</i> , <b>2010</b> , 104, 155-171   | 3.8  | 52  |
|----|---|------|-----|
| 15 | pH-Modulated Watson-Crick duplex-quadruplex equilibria of guanine-rich and cytosine-rich DNA sequences 140 base pairs upstream of the c-kit transcription initiation site. <i>Chemistry - A European Journal</i> , <b>2009</b> , 15, 12663-71 | 4.8  | 37  |
| 14 | Classification of nucleic acids structures by means of the chemometric analysis of circular dichroism spectra. <i>Analytica Chimica Acta</i> , <b>2009</b> , 642, 117-26  | 6.6  | 29  |
| 13 | Targeting the G-quadruplex-forming region near the P1 promoter in the human BCL-2 gene with the cationic porphyrin TMPyP4 and with the complementary C-rich strand. <i>Biochimie</i> , <b>2009</b> , 91, 894-902                              | 4.6  | 36  |
| 12 | Study of the interaction between the G-quadruplex-forming thrombin-binding aptamer and the porphyrin 5,10,15,20-tetrakis-(N-methyl-4-pyridyl)-21,23H-porphyrin tetratosylate. <i>Analytical Biochemistry</i> , <b>2008</b> , 379, 8-15        | 3.1  | 42  |
| 11 | Quality assessment of the results obtained by multivariate curve resolution analysis of multiple runs of gasoline blending processes. <i>Journal of Chemometrics</i> , <b>2006</b> , 20, 54-67  | 1.6  | 25  |
| 10 | Resolution of a structural competition involving dimeric G-quadruplex and its C-rich complementary strand. <i>Nucleic Acids Research</i> , <b>2006</b> , 34, 206-16   | 20.1 | 34  |
| 9  | Exploratory data analysis of DNA microarrays by multivariate curve resolution. <i>Analytical Biochemistry</i> , <b>2006</b> , 358, 76-89  | 3.1  | 28  |
| 8  | A graphical user-friendly interface for MCR-ALS: a new tool for multivariate curve resolution in MATLAB. <i>Chemometrics and Intelligent Laboratory Systems</i> , <b>2005</b> , 76, 101-110   | 3.8  | 852 |
| 7  | Non-negativity constraints for elimination of multiple solutions in fitting of multivariate kinetic models to spectroscopic data. <i>Journal of Chemometrics</i> , <b>2005</b> , 19, 97-106   | 1.6  | 26  |
| 6  | Application of multivariate resolution methods to the study of biochemical and biophysical processes. <i>Analytical Biochemistry</i> , <b>2004</b> , 327, 1-13  | 3.1  | 46  |
| 5  | Noise propagation and error estimations in multivariate curve resolution alternating least squares using resampling methods. <i>Journal of Chemometrics</i> , <b>2004</b> , 18, 327-340   | 1.6  | 47  |
| 4  | Indications towards a stereoselectivity of the salt-induced peptide formation reaction. <i>Inorganica Chimica Acta</i> , <b>2004</b> , 357, 649-656   | 2.7  | 29  |
| 3  | Multivariate curve resolution applied to the analysis and resolution of two-dimensional [1H,15N] NMR reaction spectra. <i>Analytical Chemistry</i> , <b>2004</b> , 76, 7094-101   | 7.8  | 51  |
| 2  | Multivariate resolution of NMR labile signals by means of hard- and soft-modelling methods. <i>Analytica Chimica Acta</i> , <b>2003</b> , 490, 253-264  | 6.6  | 17  |
| 1  | Multivariate curve resolution: a powerful tool for the analysis of conformational transitions in nucleic acids. <i>Nucleic Acids Research</i> , <b>2002</b> , 30, e92   | 20.1 | 61  |