# **Special Issue**

# Chemometrics in Analytical Chemistry

## Message from the Guest Editor

Due to the enormous development of computer technology during the last decades, chemometrics has become the leading and preferred methodology for the experimental data analysis, especially in analytical chemistry. A significant interest in chemometric methods is also connected with the availability of opensource software, removing the financial barriers of expensive software packages. Today, chemometric methods are available for every interested researcher equipped with an average computer. Therefore, chemometrics can be present everywhere—from simple experimental designs, through multivariate analysis of collected data, up to huge datasets containing millions of samples or variables. This Special Issue focuses on all aspects of chemometrics in analytical chemistryexperimental design, instrumental data analysis, signal processing, image processing, multivariate data mining, neural networks, genetic algorithms, multi-way methods, and multivariate curve resolution-both in context of new methods and algorithms, as well as novel applications of known approaches. Reviews are also welcome.

## **Guest Editor**

Prof. Dr. Lukasz Komsta

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### Deadline for manuscript submissions

closed (31 December 2022)



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## **About the Journal**

## Message from the Editor-in-Chief

As the premier open access journal dedicated to molecular chemistry, now in its 29th year of publication, the papers published in *Molecules* span from classical synthetic methodology to natural product isolation and characterization, as well as physicochemical studies and the applications of these molecules as pharmaceuticals, catalysts, and novel materials. Pushing the boundaries of the discipline, we invite papers on all major fields of molecular chemistry and multidisciplinary topics bridging chemistry with biology, physics, and materials science, as well as timely reviews and topical issues on cutting-edge fields in all of these areas.

#### Editor-in-Chief

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