# Template and instructions for writing contribution abstracts to MIB 2023 

First Author ${ }^{1}$, Second Author ${ }^{1}$, Third Author ${ }^{2}$<br>${ }^{1}$ First Afilliation, ${ }^{2}$ Second Affiliation

This document is a template for abstracts submitted to the conference Modeling Interactions in Biomolecules 2023. The LaTeX and word-processor (DOC and RTF) source files can be found on the conferece web site, http://physics.mff.cuni.cz/kchfo/MIB23/contribution.htm.

The entire abstract should fit on a single page, including text, tables, figures, keywords, and references. The page layout and spacing of the template should not be changed, and excessive additional formatting (changing fonts, bold and italic text) is strongly discouraged. The entire text should be written in 12 pt size.

|  | Column 1 | Column 2 | Column 3 |
| :---: | :---: | :---: | :---: |
| Row 1 | $\alpha$ | $\beta$ | $\gamma$ |
| Row 2 | $\alpha$ | $\beta$ | $\gamma$ |

Table 1: Table caption
Any tables and figures should usually span the entire width of the text and be centered on the page. Table and figure captions should also be written in 12 pt size, centered below the table or figure.

If a table or figure needs to be wrapped by text (due to an unconventional shape, or to fit more text on the page), it should take up less than half the width of the page. The abstract, including figures, must be greyscale only. Figures should be large enough so that all details are clearly visible, and include only embedded graphic files (no other inline documents) in one of the following formats: BMP, JPG, GIF, PNG, SVG, EPS, PS, PDF.

Equations should span the entire width of the text and occur


Figure 1: Figure caption where they are referred to. They should be written in the same size as the main text and be numbered in parentheses, for example:

$$
\begin{equation*}
M_{k l m} \approx A e^{i \omega t} \frac{1-x}{\sin ^{2} x} \tag{1}
\end{equation*}
$$

Acknowledgements and a maximum of four references should be placed at the bottom of the page as shown below. References should be listed in the order in which they appear in the text.

Acknowledgements: Support from EU grant No. 12345 is acknowledged. Calculations were performed at the Wroclaw Centre for Networking and Supercomputing.
[1] Jackson, M.; Webster, A.B. Curr. Org. Chem. 2000, 11 (5), 123-456.

