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Tomasz Włodarski

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- born November 6th 1984
- citizen of the Republic of Poland

Education

- 2008 - 2013 **University of Warsaw**, Warsaw, Poland
Inter-Faculty Interdisciplinary Doctoral Studies in Natural Sciences and Mathematics
PhD studies in Bioinformatics and Biophysics
Thesis: "Application of computational biophysics and bioinformatics to multiscale biological problems" (advisors: Prof. Paweł Golik and Prof. Marek Niezgódka)
- 2003 - 2008 **Jagiellonian University**, Krakow, Poland
M. Sc. in Biophysics, 2008
Thesis: "Study of interactions of fluorescent dyes (Col-F and Sulphorodamine B) with collagens"
(advisors: Prof. Jerzy Dobrucki and Prof. Marta Pasenkiewicz-Gierula)

Work Experience

- 09/2023 - ... **Institute of Biochemistry and Biophysics** Polish Academy of Sciences, Warsaw, Poland
- 06/2013 - 08/2023 **University College London**, Institute of Structural and Molecular Biology, London, UK
- 10/2016 - 04/2019 **University of Cambridge**, Chemistry Department, Cambridge, UK

Joint EMBO Postdoctoral Fellow in Prof. John Christodoulou (UCL) and Prof. Michele Vendruscolo (Univ. of Cambridge) groups.

Scientific Experience

- 11/2011 – 02/2012 **University of Cambridge, Chemistry Department**, Cambridge, UK
visiting student in Prof. Michele Vendruscolo group
(*EMBO Short Term Fellowship*)
- 07/2011 **Max F. Perutz Laboratories**, Vienna, Austria
visiting student in Computational Biophysics of Macromolecules Group
(*advisor: Dr. Bojan Zagrovic*)
- 02/2010 **University of Texas Medical Branch at Galveston**, Galveston, TX, USA
visiting student in Bioinformatics and System Biology Group
Project: “Analysis of sequencing data of human fragile sites”
(*advisor: Prof. Maga Rowicka*)
- 08/2009 **Mediterranean Institute for Life Sciences**, Split, Croatia
Summer studentship in Computational Biophysics of Macromolecules Group
Project: “Conformational selection vs induced fit in protein-protein binding”
(*advisor: Dr. Bojan Zagrovic*)
- 07/2008 – 08/2008 **Mediterranean Institute for Life Sciences**, Split, Croatia
FEBS Summer Scholarship in Computational Biophysics of Macromolecules Group
Project: “Conformational selection vs induced fit in protein-protein binding”
(*advisor: Dr. Bojan Zagrovic*)
- 07/2007 – 08/2007 **University of Illinois in Urbana-Champaign**, Urbana-Champaign, IL, USA
Summer studentship in Theoretical and Computational Biophysics Group
Project: “Computational study of solvation and hydrophobic effect around a simple molecular compound”
(*advisor: Prof. Klaus Schulten*)

Research Interest:

- protein folding and missfolding
- computational protein design
- molecular dynamics simulations
- integrative structural biology
- protein structure and dynamics
- structural bioinformatics
- machine learning

Honors and Fellowships:

- EMBO Long Term Fellowship (2013)
- EMBO Short Term Fellowship (2011)
- Foundation for Polish Science Scholarship - START 2011
- FEBS Summer Scholarship (2008)
- FEBS prize for the best Summer Scholarship report in 2008

Research Grants:

- **POLONEZ BIS** grant from the National Science Centre for the project: “Co-translational protein folding in the light of ribosome evolution”. (2023-2025)

Computational Grants:

- **LUMI EUROHPC-JU**, Regular Access 2022, EU: Computational microscopy: in-cell co-translational folding and misfolding in health and disease (**co-Principal Investigator**)
- **ARCHER2** HecBiosim Grant 2022, UK: “Structures of oncogenic Ras protein folding intermediates on the ribosome as novel drug targets” (**co-Principal Investigator**)
- **EPSRC**, Access to High Performance Computing Call, 2022, UK: “Atomistic insights into co-translational protein folding on the ribosome using enhanced sampling molecular dynamics simulations” (**co-Principal Investigator**)
- **ARCHER2** HecBiosim Grant 2022, UK: “Determining the structural basis of co-translational protein folding by integrating state-of-the-art MD simulations with experimental data.” (**co-Principal Investigator**)
- **EPSRC**, Access to High Performance Computing Call, 2021, UK: “Computational studies of co-translational protein folding on the ribosome” (**co-Principal Investigator**)
- **ARCHER2** HecBiosim Grant 2021, UK: “Applying coarse-grained molecular dynamics simulations to study ribosome interactions: from nascent polypeptides to small molecule drugs” (**co-Principal Investigator**)
- **ICM Okeanos** Computational grant, PL (2019 - 2021): “Studying in atomic details interactions between protein nascent chain and ribosome during protein biosynthesis” (**Principal Investigator**)
- **ICM Okeanos** - Grand Challenges, PL (2017): “Integrative structural biology approach to study protein folding and misfolding on the ribosome” (**Principal Investigator**)
- **ICM Okeanos** - Grand Challenges, PL (2016): “Studying protein folding and misfolding during biosynthesis on the ribosome using a computational microscope.” (**Principal Investigator**)
- **ARCHER** Leadership Grant, UK (2015-2017): “Computational studies of co-translational protein folding and mis-folding on the ribosome” (**co-Principal Investigator**)

Editorial Board:

- Review Editor in *Frontiers in Molecular Biosciences* (Biological Modeling and Simulation)

Other Experience:

- EMBO Laboratory Leadership Course for Postdocs (2018)
- Polonium Foundation Scientific Content Manager (2017 - 2021)

Publications:

1. Chan, S.H.S., **Włodarski, T.**, Streit, J., Cassaignau, A.M.E., Woodburn, L., Ahn, M., Waudby, C.A., Budisa, N., Cabrita, L.D., Christodoulou, J. 2022 “*The ribosome stabilises partially folded intermediates of a nascent multi-domain protein*” **Nat. Chem.** *14*, 1165–1173. <https://doi.org/10.1038/s41557-022-01004-0>
2. **Włodarski, T.**, Ahn, M., Mitropoulou A., Chan, S.H.S., Sidhu, H., Plessa, E., Becker, T.A., Waudby, C.A., Beckmann, R., Cassaignau, A.M.E., Cabrita, L.D., Christodoulou, J. 2022 “*Modulating co-translational protein folding by rational design and ribosome engineering*” **Nat. Commun.** *13*, 4243. <https://doi.org/10.1038/s41467-022-31906-z>
3. **Włodarski, T.**, Deckert, A., Cassaignau, A.M.E., Wang, X., Chan, S.H.S., Waudby, C.A., Kirkpatrick, J.P., Vendruscolo, M., Cabrita, L.D., Christodoulou, J., 2021. *Common sequence motifs of nascent chains engage the ribosome surface and trigger factor.* **Proc Natl Acad Sci USA** *118*. doi:10.1073/pnas.2103015118
4. Cassaignau, A.M.E., **Włodarski, T.**, Chan, S.H.S., Woodburn, L.F., Bukvin, I.V., Streit, J.O., Cabrita, L.D., Waudby, C.A., Christodoulou, J., 2021. *Interactions between nascent proteins and the ribosome surface inhibit co-translational folding.* **Nat. Chem.** *13*, 1214–1220. <https://doi.org/10.1038/s41557-021-00796-x>
5. Burrige, C., Waudby, C.A., **Włodarski, T.**, Cassaignau, A.M.E., Cabrita, L.D., Christodoulou, J., 2021. *Nascent chain dynamics and ribosome interactions within folded ribosome-nascent chain complexes observed by NMR spectroscopy.* **Chem. Sci.** *12*, 13120–13126. doi:10.1039/d1sc04313g
6. Waudby, C.A., **Włodarski, T.**, Karyadi, M.-E., Cassaignau, A.M.E., Chan, S.H.S., Wentink, A.S., Schmidt-Engler, J.M., Camilloni, C., Vendruscolo, M., Cabrita, L.D., Christodoulou, J., 2018. *Systematic mapping of free energy landscapes of a growing filamin domain during biosynthesis.* **Proc Natl Acad Sci USA** *115*, 9744–9749. doi:10.1073/pnas.1716252115
7. Redondo, R.A.F., de Vlarar, H.P., **Włodarski, T.**, Bollback, J.P., 2017. *Evolutionary interplay between structure, energy and epistasis in the coat protein of the ϕ X174 phage family.* **J. R. Soc. Interface** *14*. doi:10.1098/rsif.2016.0139
8. Deckert, A., Waudby, C.A., **Włodarski, T.**, Wentink, A.S., Wang, X., Kirkpatrick, J.P., Paton, J.F.S., Camilloni, C., Kukic, P., Dobson, C.M., Vendruscolo, M., Cabrita, L.D., Christodoulou, J., 2016. *Structural characterization of the interaction of α -synuclein nascent chains with the ribosomal surface and trigger factor.* **Proc Natl Acad Sci USA** *113*, 5012–5017. doi:10.1073/pnas.1519124113
9. Cabrita, L.D., Cassaignau, A.M.E., Launay, H.M.M., Waudby, C.A., **Włodarski, T.**, Camilloni, C., Karyadi, M.-E., Robertson, A.L., Wang, X., Wentink, A.S., Goodsell, L., Woolhead, C.A., Vendruscolo, M., Dobson, C.M., Christodoulou, J., 2016. *A structural ensemble of a ribosome-nascent chain complex during cotranslational protein folding.* **Nat. Struct. Mol. Biol.** *23*, 278–285. doi:10.1038/nsmb.3182
10. **Włodarski, T.**, Kutner, J., Towpik, J., Knizewski, L., Rychlewski, L., Kudlicki, A., Rowicka, M., Dziembowski, A., Ginalski, K., 2011. *Comprehensive structural and substrate specificity classification of the *Saccharomyces cerevisiae* methyltransferase.* **PLoS ONE** *6*, e23168. doi:10.1371/journal.pone.0023168

11. **Włodarski, T.**, Zagrovic, B., 2009. *Conformational selection and induced fit mechanism underlie specificity in noncovalent interactions with ubiquitin*. **Proc Natl Acad Sci USA** 106, 19346–19351. doi:10.1073/pnas.0906966106

Preprints:

1. Javed, A., **Włodarski, T.**, Cassaignau, AME., Cabrita, LD., Christodoulou, J., Orlova, EV. „*Visualising nascent chain dynamics at the ribosome exit tunnel by cryo-electron microscopy*” doi: <https://doi.org/10.1101/722611>

Languages

- **Polish** - native
- **English** - fluent
- **Russian** – basic
- **German** – basic