

# Robert Vãjcha

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/1249814/publications.pdf>

Version: 2024-02-01

68  
papers

4,756  
citations

101543

36  
h-index

95266

68  
g-index

74  
all docs

74  
docs citations

74  
times ranked

5733  
citing authors

| #  | ARTICLE  | IF   | CITATIONS |
|----|--|------|-----------|
| 1  | Magainin 2 and PGLa in Bacterial Membrane Mimics III: Membrane Fusion and Disruption. <i>Biophysical Journal</i> , 2022, , .   | 0.5  | 4         |
| 2  | Synthesis and Profiling of Highly Selective Inhibitors of Methyltransferase DOT1L Based on Carbocyclic C-Nucleosides. <i>Journal of Medicinal Chemistry</i> , 2022, 65, 5701-5723. | 6.4  | 5         |
| 3  | Capsid Structure of <i>Leishmania</i> RNA Virus 1. <i>Journal of Virology</i> , 2021, 95, .  | 3.4  | 10        |
| 4  | Phosphorylation-induced changes in the PDZ domain of Dishevelled 3. <i>Scientific Reports</i> , 2021, 11, 1484.  | 3.3  | 2         |
| 5  | Capsid opening enables genome release of iflaviruses. <i>Science Advances</i> , 2021, 7, .   | 10.3 | 13        |
| 6  | Selecting Collective Variables and Free-Energy Methods for Peptide Translocation across Membranes. <i>Journal of Chemical Information and Modeling</i> , 2021, 61, 819-830.        | 5.4  | 22        |
| 7  | Advances in Molecular Understanding of $\alpha$ -Helical Membrane-Active Peptides. <i>Accounts of Chemical Research</i> , 2021, 54, 2196-2204.                                     | 15.6 | 47        |
| 8  | Enhanced translocation of amphiphilic peptides across membranes by transmembrane proteins. <i>Biophysical Journal</i> , 2021, 120, 2296-2305.                                      | 0.5  | 7         |
| 9  | The impact of the glycan headgroup on the nanoscopic segregation of gangliosides. <i>Biophysical Journal</i> , 2021, 120, 5530-5543.   | 0.5  | 8         |
| 10 | Cargo Release from Nonenveloped Viruses and Virus-like Nanoparticles: Capsid Rupture or Pore Formation. <i>ACS Nano</i> , 2021, 15, 19233-19243.                                   | 14.6 | 7         |
| 11 | Magainin 2 and PGLa in Bacterial Membrane Mimics II: Membrane Fusion and Sponge Phase Formation. <i>Biophysical Journal</i> , 2020, 118, 612-623.                                  | 0.5  | 25        |
| 12 | Effect of membrane composition on DivIVA-membrane interaction. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2020, 1862, 183144.   | 2.6  | 3         |
| 13 | Yeast Spt6 Reads Multiple Phosphorylation Patterns of RNA Polymerase II C-Terminal Domain In Vitro. <i>Journal of Molecular Biology</i> , 2020, 432, 4092-4107.                    | 4.2  | 6         |
| 14 | Effect of Helical Kink on Peptide Translocation across Phospholipid Membranes. <i>Journal of Physical Chemistry B</i> , 2020, 124, 5940-5947.                                      | 2.6  | 13        |
| 15 | Effect of helical kink in antimicrobial peptides on membrane pore formation. <i>ELife</i> , 2020, 9, .   | 6.0  | 39        |
| 16 | Magainin 2 and PGLa in Bacterial Membrane Mimics I: Peptide-Peptide and Lipid-Peptide Interactions. <i>Biophysical Journal</i> , 2019, 117, 1858-1869.                             | 0.5  | 30        |
| 17 | Dishevelled-3 conformation dynamics analyzed by FRET-based biosensors reveals a key role of casein kinase 1. <i>Nature Communications</i> , 2019, 10, 1804.                        | 12.8 | 20        |
| 18 | Enterovirus particles expel capsid pentamers to enable genome release. <i>Nature Communications</i> , 2019, 10, 1138.  | 12.8 | 33        |

| #  | ARTICLE   | IF   | CITATIONS |
|----|---|------|-----------|
| 19 | Aggregate Size Dependence of Amyloid Adsorption onto Charged Interfaces. <i>Langmuir</i> , 2018, 34, 1266-1273.   | 3.5  | 5         |
| 20 | Synergism of Antimicrobial Frog Peptides Couples to Membrane Intrinsic Curvature Strain. <i>Biophysical Journal</i> , 2018, 114, 1945-1954.   | 0.5  | 55        |
| 21 | Optimal Hydrophobicity and Reorientation of Amphiphilic Peptides Translocating through Membrane. <i>Biophysical Journal</i> , 2018, 115, 1045-1054.                                       | 0.5  | 29        |
| 22 | Design of Multivalent Inhibitors for Preventing Cellular Uptake. <i>Scientific Reports</i> , 2017, 7, 11689.  | 3.3  | 9         |
| 23 | Self-assembled clusters of patchy rod-like molecules. <i>Soft Matter</i> , 2017, 13, 7492-7497.   | 2.7  | 8         |
| 24 | GM 1 $\beta$ Gangliosid hemmt die $\beta$ -Amyloid-Oligomerisation, wahrend Sphingomyelin diese initiiert. <i>Angewandte Chemie</i> , 2016, 128, 9557-9562.                              | 2.0  | 1         |
| 25 | GM <sub>1</sub> Ganglioside Inhibits $\beta$ -Amyloid Oligomerization Induced by Sphingomyelin. <i>Angewandte Chemie - International Edition</i> , 2016, 55, 9411-9415.                   | 13.8 | 86        |
| 26 | Optimal conditions for opening of membrane pore by amphiphilic peptides. <i>Journal of Chemical Physics</i> , 2015, 143, 243115.  | 3.0  | 21        |
| 27 | Influence of ligand distribution on uptake efficiency. <i>Soft Matter</i> , 2015, 11, 2726-2730.  | 2.7  | 49        |
| 28 | Simulations Suggest Possible Novel Membrane Pore Structure. <i>Langmuir</i> , 2014, 30, 1304-1310.  | 3.5  | 19        |
| 29 | Surface Effects on Aggregation Kinetics of Amyloidogenic Peptides. <i>Journal of the American Chemical Society</i> , 2014, 136, 11776-11782.  | 13.7 | 158       |
| 30 | Stability of Bicelles: A Simulation Study. <i>Langmuir</i> , 2014, 30, 4229-4235.   | 3.5  | 14        |
| 31 | Composition- and Size-Controlled Cyclic Self-Assembly by Solvent- and C <sub>60</sub> -Responsive Self-Sorting. <i>Journal of the American Chemical Society</i> , 2013, 135, 15263-15268. | 13.7 | 30        |
| 32 | <i>Faunus</i> – a flexible framework for Monte Carlo simulation. <i>Molecular Simulation</i> , 2013, 39, 1233-1239.   | 2.0  | 44        |
| 33 | Connecting Macroscopic Observables and Microscopic Assembly Events in Amyloid Formation Using Coarse Grained Simulations. <i>PLoS Computational Biology</i> , 2012, 8, e1002692.          | 3.2  | 63        |
| 34 | Aqueous Solutions at the Interface with Phospholipid Bilayers. <i>Accounts of Chemical Research</i> , 2012, 45, 74-82.  | 15.6 | 100       |
| 35 | Intracellular Release of Endocytosed Nanoparticles Upon a Change of Ligand-Receptor Interaction. <i>ACS Nano</i> , 2012, 6, 10598-10605.  | 14.6 | 55        |
| 36 | Sodium Dodecyl Sulfate at Water-Hydrophobic Interfaces: A Simulation Study. <i>Journal of Physical Chemistry B</i> , 2012, 116, 11936-11942.  | 2.6  | 31        |

| #  | ARTICLE   | IF   | CITATIONS |
|----|---|------|-----------|
| 37 | Charge Transfer between Water Molecules As the Possible Origin of the Observed Charging at the Surface of Pure Water. <i>Journal of Physical Chemistry Letters</i> , 2012, 3, 107-111.                            | 4.6  | 101       |
| 38 | Ions at Hydrophobic Aqueous Interfaces: Molecular Dynamics with Effective Polarization. <i>Journal of Physical Chemistry Letters</i> , 2012, 3, 2087-2091.  | 4.6  | 89        |
| 39 | Study of the betulin molecule in a water environment; ab initio and molecular simulation calculations. <i>Journal of Molecular Modeling</i> , 2012, 18, 367-376.  | 1.8  | 5         |
| 40 | Relation between Molecular Shape and the Morphology of Self-Assembling Aggregates: A Simulation Study. <i>Biophysical Journal</i> , 2011, 101, 1432-1439.   | 0.5  | 62        |
| 41 | The Orientation and Charge of Water at the Hydrophobic Oil Droplet-Water Interface. <i>Journal of the American Chemical Society</i> , 2011, 133, 10204-10210.   | 13.7 | 213       |
| 42 | Receptor-Mediated Endocytosis of Nanoparticles of Various Shapes. <i>Nano Letters</i> , 2011, 11, 5391-5395.  | 9.1  | 441       |
| 43 | Three-dimensional energy profile measurement of a molecular ion beam by coincidence momentum imaging compared to a retarding field analyzer. <i>Journal of Instrumentation</i> , 2010, 5, P10006-P10006.          | 1.2  | 0         |
| 44 | Dielectric Interpretation of Specificity of Ion Pairing in Water. <i>Journal of Physical Chemistry Letters</i> , 2010, 1, 300-303.  | 4.6  | 41        |
| 45 | Comment on "An explanation for the charge on water's surface" by A. Gray-Weale and J. K. Beattie, <i>Phys. Chem. Chem. Phys.</i> , 2009, 11, 10994. <i>Physical Chemistry Chemical Physics</i> , 2010, 12, 14362. | 2.8  | 12        |
| 46 | Mechanism of Interaction of Monovalent Ions with Phosphatidylcholine Lipid Membranes. <i>Journal of Physical Chemistry B</i> , 2010, 114, 9504-9509.  | 2.6  | 89        |
| 47 | Large variations in the propensity of aqueous oxychlorine anions for the solution/vapor interface. <i>Journal of Chemical Physics</i> , 2009, 131, 124706.  | 3.0  | 24        |
| 48 | Hofmeister series and specific interactions of charged headgroups with aqueous ions. <i>Advances in Colloid and Interface Science</i> , 2009, 146, 42-47.   | 14.7 | 378       |
| 49 | Reply to comments on <i>Frontiers Article "Behavior of hydroxide at the water/vapor interface"</i> . <i>Chemical Physics Letters</i> , 2009, 481, 19-21.  | 2.6  | 22        |
| 50 | Behavior of hydroxide at the water/vapor interface. <i>Chemical Physics Letters</i> , 2009, 474, 241-247.   | 2.6  | 110       |
| 51 | Effects of Alkali Cations and Halide Anions on the DOPC Lipid Membrane. <i>Journal of Physical Chemistry A</i> , 2009, 113, 7235-7243.  | 2.5  | 144       |
| 52 | Ion specific effects of sodium and potassium on the catalytic activity of HIV-1 protease. <i>Physical Chemistry Chemical Physics</i> , 2009, 11, 7599.  | 2.8  | 36        |
| 53 | Molecular Model of a Cell Plasma Membrane With an Asymmetric Multicomponent Composition: Water Permeation and Ion Effects. <i>Biophysical Journal</i> , 2009, 96, 4493-4501.                                      | 0.5  | 75        |
| 54 | Benchmarking Polarizable Molecular Dynamics Simulations of Aqueous Sodium Hydroxide by Diffraction Measurements. <i>Journal of Physical Chemistry A</i> , 2009, 113, 4022-4027.                                   | 2.5  | 22        |

| #  | ARTICLE   | IF  | CITATIONS |
|----|---|-----|-----------|
| 55 | Specific Ion Binding to Macromolecules: Effects of Hydrophobicity and Ion Pairing. <i>Langmuir</i> , 2008, 24, 3387-3391.   | 3.5 | 106       |
| 56 | Response to Comment on Autoionization at the surface of neat water: is the top layer pH neutral, basic, or acidic? by J. K. Beattie, <i>Phys. Chem. Chem. Phys.</i> , 2007, 9, DOI: 10.1039/b713702h. <i>Physical Chemistry Chemical Physics</i> , 2008, 10, 332-333.   | 2.8 | 37        |
| 57 | Cation-Specific Interactions with Carboxylate in Amino Acid and Acetate Aqueous Solutions: X-ray Absorption and <i>ab initio</i> Calculations. <i>Journal of Physical Chemistry B</i> , 2008, 112, 12567-12570.   | 2.6 | 149       |
| 58 | Biomolecular simulations of membranes: Physical properties from different force fields. <i>Journal of Chemical Physics</i> , 2008, 128, 125103.   | 3.0 | 242       |
| 59 | Hydronium and hydroxide at the interface between water and hydrophobic media. <i>Physical Chemistry Chemical Physics</i> , 2008, 10, 4975.  | 2.8 | 68        |
| 60 | Adsorption of Aromatic Hydrocarbons and Ozone at Environmental Aqueous Surfaces. <i>Journal of Physical Chemistry A</i> , 2008, 112, 4942-4950.   | 2.5 | 38        |
| 61 | Water Structuring and Hydroxide Ion Binding at the Interface between Water and Hydrophobic Walls of Varying Rigidity and van der Waals Interactions. <i>Journal of Physical Chemistry C</i> , 2008, 112, 7689-7692.   | 3.1 | 53        |
| 62 | Water surface is acidic. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2007, 104, 7342-7347.   | 7.1 | 332       |
| 63 | Ion Pairing as a Possible Clue for Discriminating between Sodium and Potassium in Biological and Other Complex Environments. <i>Journal of Physical Chemistry B</i> , 2007, 111, 14077-14079.   | 2.6 | 80        |
| 64 | Autoionization at the surface of neat water: is the top layer pH neutral, basic, or acidic?. <i>Physical Chemistry Chemical Physics</i> , 2007, 9, 4736.  | 2.8 | 151       |
| 65 | Adsorption of polycyclic aromatic hydrocarbons at the air-water interface: Molecular dynamics simulations and experimental atmospheric observations. <i>Physical Chemistry Chemical Physics</i> , 2006, 8, 4461-4467.   | 2.8 | 70        |
| 66 | Propensity for the Air/Water Interface and Ion Pairing in Magnesium Acetate vs Magnesium Nitrate Solutions: A Molecular Dynamics Simulations and Surface Tension Measurements. <i>Journal of Physical Chemistry B</i> , 2006, 110, 15939-15944.   | 2.6 | 86        |
| 67 | Quantification and rationalization of the higher affinity of sodium over potassium to protein surfaces. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2006, 103, 15440-15444.  | 7.1 | 212       |
| 68 | Adsorption of Atmospherically Relevant Gases at the Air/Water Interface: Free Energy Profiles of Aqueous Solvation of N <sub>2</sub> , O <sub>2</sub> , O <sub>3</sub> , OH, H <sub>2</sub> O, HO <sub>2</sub> , and H <sub>2</sub> O <sub>2</sub> . <i>Journal of Physical Chemistry A</i> , 2004, 108, 11573-11579. | 2.5 | 195       |