

# Larisa Zoranic

## List of Publications by Year in descending order

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Version: 2024-02-01

23  
papers

896  
citations

623734

14  
h-index

642732

23  
g-index

24  
all docs

24  
docs citations

24  
times ranked

1111  
citing authors

| #  | ARTICLE  | IF  | CITATIONS |
|----|--|-----|-----------|
| 1  | Basic Charge Clusters and Predictions of Membrane Protein Topology. <i>Journal of Chemical Information and Computer Sciences</i> , 2002, 42, 620-632.  | 2.8 | 193       |
| 2  | Microstructure of neat alcohols: A molecular dynamics study. <i>Journal of Chemical Physics</i> , 2007, 127, 024502.   | 3.0 | 88        |
| 3  | A comparative Molecular Dynamics study of water-methanol and acetone-methanol mixtures. <i>Journal of Molecular Liquids</i> , 2011, 159, 52-59.  | 4.9 | 79        |
| 4  | Ethanol-water mixtures: ultrasonics, Brillouin scattering and molecular dynamics. <i>Journal of Molecular Liquids</i> , 2011, 164, 66-73.  | 4.9 | 76        |
| 5  | Micro-heterogeneity versus clustering in binary mixtures of ethanol with water or alkanes. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 23971-23979.                                   | 2.8 | 72        |
| 6  | Microstructure of neat alcohols. <i>Physical Review E</i> , 2007, 75, 060502.  | 2.1 | 65        |
| 7  | Antibacterial Activity Affected by the Conformational Flexibility in Glycine-Lysine Based $\alpha$ -Helical Antimicrobial Peptides. <i>Journal of Medicinal Chemistry</i> , 2018, 61, 2924-2936. | 6.4 | 48        |
| 8  | Concentration fluctuations and microheterogeneity in aqueous amide mixtures. <i>Journal of Chemical Physics</i> , 2009, 130, 124315.   | 3.0 | 38        |
| 9  | Structural changes in ethanol-water mixtures: Ultrasonics, Brillouin scattering and molecular dynamics studies. <i>Vibrational Spectroscopy</i> , 2012, 60, 102-106.                             | 2.2 | 34        |
| 10 | A comparison of force fields for ethanol-water mixtures. <i>Molecular Simulation</i> , 2015, 41, 699-712.  | 2.0 | 34        |
| 11 | Simple and complex disorder in binary mixtures with benzene as a common solvent. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 9885-9898.   | 2.8 | 32        |
| 12 | On the Microheterogeneity in Neat and Aqueous Amides: A Molecular Dynamics Study. <i>Journal of Physical Chemistry C</i> , 2007, 111, 15586-15595.   | 3.1 | 24        |
| 13 | PGLa-H tandem-repeat peptides active against multidrug resistant clinical bacterial isolates. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2017, 1859, 228-237.                         | 2.6 | 23        |
| 14 | The microscopic structure of cold aqueous methanol mixtures. <i>Journal of Chemical Physics</i> , 2016, 145, 144502.   | 3.0 | 20        |
| 15 | Anisaxins, helical antimicrobial peptides from marine parasites, kill resistant bacteria by lipid extraction and membrane disruption. <i>Acta Biomaterialia</i> , 2022, 146, 131-144.            | 8.3 | 15        |
| 16 | Molecular dynamics simulations and femtosecond optical Kerr effect spectroscopy of methanol/acetone mixtures. <i>Journal of Molecular Liquids</i> , 2011, 159, 60-69.                            | 4.9 | 12        |
| 17 | A simple two dimensional model of methanol. <i>Journal of Molecular Liquids</i> , 2018, 262, 46-57.  | 4.9 | 8         |
| 18 | Designed peptide with a flexible central motif from ranatuerins adapts its conformation to bacterial membranes. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2018, 1860, 2655-2668.     | 2.6 | 8         |

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|----|---|-----|-----------|
| 19 | Water-like structure with repulsive double-core interactions. <i>Molecular Physics</i> , 2009, 107, 1349-1353.  | 1.7 | 7         |
| 20 | A re-appraisal of the concept of ideal mixtures through a computer simulation study of the methanol-ethanol mixtures. <i>Journal of Chemical Physics</i> , 2016, 145, .           | 3.0 | 7         |
| 21 | Density and energy distribution in water and organic solvents: A molecular dynamics study. <i>Journal of Molecular Liquids</i> , 2007, 136, 199-205.                              | 4.9 | 5         |
| 22 | The structuring in mixtures with acetone as the common solvent. <i>Physics and Chemistry of Liquids</i> , 2020, 58, 184-201.  | 1.2 | 5         |
| 23 | The influence of binary mixtures' structuring on the calculation of Kirkwood-Buff integrals: A molecular dynamics study. <i>Journal of Molecular Liquids</i> , 2021, 324, 114773. | 4.9 | 3         |