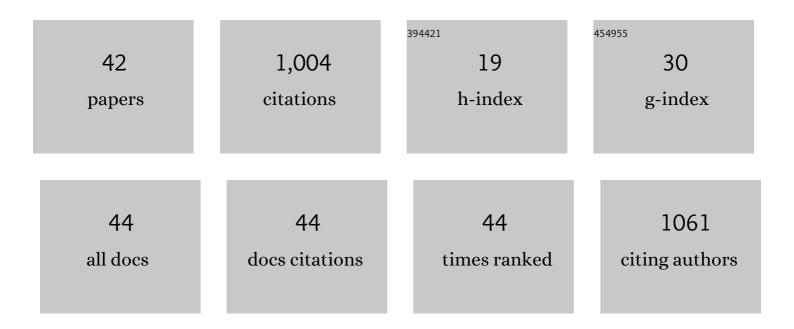
Vladislav B IvaniÅ;tÅ;ev

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	Electrical double layer in ionic liquids: Structural transitions from multilayer to monolayer structure at the interface. Electrochimica Acta, 2013, 110, 762-771.	5.2	110
2	Poly(a)morphic portrait of the electrical double layer in ionic liquids. Electrochemistry Communications, 2014, 48, 61-64.	4.7	64
3	Screening of Ion–Craphene Electrode Interactions by Ionic Liquids: The Effects of Liquid Structure. Journal of Physical Chemistry C, 2014, 118, 5841-5847.	3.1	54
4	Interfaces between Charged Surfaces and Ionic Liquids: Insights from Molecular Simulations. Electrochemical Society Interface, 2014, 23, 65-69.	0.4	45
5	Molecular Response of 1-Butyl-3-Methylimidazolium Dicyanamide Ionic Liquid at the Graphene Electrode Interface Investigated by Sum Frequency Generation Spectroscopy and Molecular Dynamics Simulations. Journal of Physical Chemistry C, 2015, 119, 26009-26019.	3.1	44
6	Molecular dynamics simulation of the behaviour of water in nano-confined ionic liquid–water mixtures. Journal of Physics Condensed Matter, 2016, 28, 464001.	1.8	44
7	Influence of cation chemical composition and structure on the double layer capacitance for Bi(111) room temperature ionic liquid interface. Journal of Electroanalytical Chemistry, 2012, 668, 30-36.	3.8	43
8	On the thickness of the double layer in ionic liquids. Physical Chemistry Chemical Physics, 2018, 20, 10275-10285.	2.8	40
9	Molecular origin of high free energy barriers for alkali metal ion transfer through ionic liquid–graphene electrode interfaces. Physical Chemistry Chemical Physics, 2016, 18, 1302-1310.	2.8	39
10	Restructuring of the electrical double layer in ionic liquids upon charging. Journal of Physics Condensed Matter, 2015, 27, 102101.	1.8	37
11	Predictions of Physicochemical Properties of Ionic Liquids with DFT. Computation, 2016, 4, 25.	2.0	35
12	Molecular dynamics simulation of the structure and interfacial free energy barriers of mixtures of ionic liquids and divalent salts near a graphene wall. Physical Chemistry Chemical Physics, 2017, 19, 846-853.	2.8	33
13	Ionic liquid–metal interface: The origins of capacitance peaks. Electrochimica Acta, 2021, 379, 138148.	5.2	28
14	The nanostructure of a lithium glyme solvate ionic liquid at electrified interfaces. Physical Chemistry Chemical Physics, 2017, 19, 11004-11010.	2.8	27
15	Electrochemical Investigation of 1-Ethyl-3-methylimidazolium Bromide and Tetrafluoroborate Mixture at Bi(111) Electrode Interface. Journal of the Electrochemical Society, 2016, 163, H723-H730.	2.9	26
16	Density Functional Theory Study of Ionic Liquid Adsorption on Circumcoronene Shaped Graphene. Journal of Physical Chemistry C, 2018, 122, 2624-2631.	3.1	26
17	Graphene–Ionic Liquid Interfacial Potential Drop from Density Functional Theory-Based Molecular Dynamics Simulations. Journal of Physical Chemistry C, 2020, 124, 19548-19555.	3.1	24
18	Influence of the electrode potential and in situ STM scanning conditions on the phase boundary structure of the single crystal Bi(1 1 1) 1-butyl-4-methylpyridinium tetrafluoroborate interface. Journal of Electroanalytical Chemistry, 2013, 709, 46-56.	3.8	22

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19	Catalytic CO ₂ /CO Reduction: Gas, Aqueous, and Aprotic Phases. ACS Catalysis, 2022, 12, 2561-2568.	11.2	22
20	Self-interaction error in DFT-based modelling of ionic liquids. Physical Chemistry Chemical Physics, 2016, 18, 2175-2182.	2.8	20
21	Comparative Impedance Study of Cd(0001) Electrode in EMImBF ₄ and KI Aqueous Solution at Different Temperatures. Journal of the Electrochemical Society, 2013, 160, H368-H375.	2.9	18
22	Density functional theory study of the water adsorption at Bi(111) electrode surface. Surface Science, 2010, 604, 1919-1927.	1.9	16
23	Simulation of a Solvate Ionic Liquid at a Polarizable Electrode with a Constant Potential. Journal of Physical Chemistry C, 2019, 123, 3935-3943.	3.1	16
24	Molecular dynamics simulations of novel electrolytes based on mixtures of protic and aprotic ionic liquids at the electrochemical interface: Structure and capacitance of the electric double layer. Electrochimica Acta, 2019, 305, 223-231.	5.2	16
25	Balance of the interfacial interactions of 4,4′-bipyridine at Bi(111) surface. Electrochimica Acta, 2014, 120, 86-95.	5.2	15
26	On the role of the surface charge plane position at Au(hkl)–BMImPF6 interfaces. Electrochimica Acta, 2019, 318, 76-82.	5.2	15
27	A comparative DFT study of the adsorption of H2O molecules at Bi, Hg, and Ga surfaces. Surface Science, 2013, 609, 91-99.	1.9	14
28	Hysteresis in the MD Simulations of Differential Capacitance at the Ionic Liquid–Au Interface. Journal of Physical Chemistry Letters, 2020, 11, 10408-10413.	4.6	14
29	Multifunctional Electrocatalysis on Single-Site Metal Catalysts: A Computational Perspective. Catalysts, 2021, 11, 1165.	3.5	11
30	Adsorption of 4,4′â~'bipyridine on the Cd(0001) single crystal electrode surface. Electrochimica Acta, 2015, 180, 965-976.	5.2	10
31	Performance of SCAN density functional for a set of ionic liquid ion pairs. International Journal of Quantum Chemistry, 2018, 118, e25582.	2.0	10
32	Understanding the Behavior of Fully Non-Toxic Polypyrrole-Gelatin and Polypyrrole-PVdF Soft Actuators with Choline Ionic Liquids. Actuators, 2020, 9, 40.	2.3	10
33	Double layer in ionic liquids: Temperature effect and bilayer model. Journal of Molecular Liquids, 2022, 363, 119747.	4.9	10
34	Interplay between the hydrophilicity of metal electrodes and their interfacial capacitance. Electrochimica Acta, 2016, 210, 615-621.	5.2	8
35	Structure and dynamics of ionic liquids: general discussion. Faraday Discussions, 2018, 206, 291-337.	3.2	8
36	NaRIBaS—A Scripting Framework for Computational Modeling of Nanomaterials and Room Temperature Ionic Liquids in Bulk and Slab. Computation, 2018, 6, 57.	2.0	7

#	Article	IF	CITATIONS
37	Predicting Melting Points of Biofriendly Choline-Based Ionic Liquids with Molecular Dynamics. Applied Sciences (Switzerland), 2019, 9, 5367.	2.5	7
38	Characteristics of Capacitors Based on Ionic Liquids: From Dielectric Polymers to Redox-Active Adsorbed Species. ECS Transactions, 2016, 75, 161-170.	0.5	6
39	Impedance study of adsorption of iodide ions at Cd(0001) and Bi(111) electrode from various solutions with constant ionic strength. Journal of Solid State Electrochemistry, 2010, 14, 555-563.	2.5	4
40	Direct migration of scientific computing experiments to the cloud. , 2013, , .		3
41	Calculation of coreâ€level electron spectra of ionic liquids. International Journal of Quantum Chemistry, 2020, 120, e26247.	2.0	3
42	Ionic liquids at interfaces: general discussion. Faraday Discussions, 2018, 206, 549-586.	3.2	0