

Zdeněk Wagner

List of Publications by Year in descending order

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papers

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687363

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citing authors

#	ARTICLE	IF	CITATIONS
1	Carbon Nanotube-Based Ionanofluids for Efficient Energy Storage: Thermophysical Propertiesâ€™ Determination and Advanced Data Analysis. <i>Industrial & Engineering Chemistry Research</i> , 2021, 60, 7714-7728.	3.7	2
2	Density and sound velocity measurement by an Anton Paar DSA 5000 density meter: Precision and long-time stability. <i>Journal of Molecular Liquids</i> , 2021, 329, 115547.	4.9	16
3	Thermal properties of novel oligoether-substituted ionic liquids and the influence of alkyl-substituent isomery. <i>Fluid Phase Equilibria</i> , 2020, 514, 112561.	2.5	5
4	Comparison of prediction methods of heat capacity of ionic liquids with selected experimental data by means of advanced data analysis. <i>Thermochimica Acta</i> , 2020, 690, 178602.	2.7	1
5	Thermochemical Properties of Menthol and Terpineol. <i>Journal of Solution Chemistry</i> , 2020, 49, 1267-1278.	1.2	6
6	Thermochemical Properties of Selected Terpenes. <i>Journal of Solution Chemistry</i> , 2020, 49, 1137-1153.	1.2	7
7	Heat capacity of 1-hexadecyl-3-methylimidazolium based ionic liquids in solid and liquid phase. <i>Journal of Molecular Liquids</i> , 2020, 305, 112847.	4.9	10
8	Phase transitions in higher-melting imidazolium-based ionic liquids: Experiments and advanced data analysis. <i>Journal of Molecular Liquids</i> , 2019, 292, 111222.	4.9	9
9	Densities, Vapor Pressures, and Surface Tensions of Selected Terpenes. <i>Journal of Solution Chemistry</i> , 2019, 48, 1147-1166.	1.2	11
10	Ionic Liquids as Thermal Energy Storage Materials: On the Importance of Reliable Data Analysis in Assessing Thermodynamic Data. <i>Journal of Solution Chemistry</i> , 2019, 48, 949-961.	1.2	12
11	Group Contribution Method for Evaluation of Volumetric Properties of Ionic Liquids Using Experimental Data Recommended by Mathematical Gnostics. <i>Industrial & Engineering Chemistry Research</i> , 2017, 56, 6827-6840.	3.7	10
12	Thermal properties of 1-alkyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide ionic liquids with linear, branched and cyclic alkyl substituents. <i>Fluid Phase Equilibria</i> , 2017, 443, 32-43.	2.5	9
13	Influence of the alkyl side chain length on the thermophysical properties of chiral ionic liquids with a (1R,2S,5R)-(-)-menthol substituent and data analysis by means of mathematical gnostics. <i>Journal of Molecular Liquids</i> , 2017, 242, 336-348.	4.9	15
14	Liquid Phase Behavior in Systems of 1-Butyl-3-alkylimidazolium bis{(trifluoromethyl)sulfonyl}imide Ionic Liquids with Water: Influence of the Structure of the C5 Alkyl Substituent. <i>Journal of Solution Chemistry</i> , 2017, 46, 1456-1474.	1.2	11
15	Advanced Analysis of Isobaric Heat Capacities by Mathematical Gnostics. <i>Journal of Solution Chemistry</i> , 2017, 46, 1836-1853.	1.2	13
16	Number Concentrations and Modal Structure of Indoor/Outdoor Fine Particles in Four European Cities. <i>Aerosol and Air Quality Research</i> , 2017, 17, 131-146.	2.1	11
17	Volumetric, acoustic and optical properties for binary mixtures of nitroethane with chloroalkane at temperatures between 298.15 K and 318.15 K. Comparison with theories. <i>Journal of Molecular Liquids</i> , 2016, 223, 790-804.	4.9	4
18	New Method Based on the UNIFACâ€™VISCO Model for the Estimation of Ionic Liquids Viscosity Using the Experimental Data Recommended by Mathematical Gnostics. <i>Journal of Chemical & Engineering Data</i> , 2016, 61, 3908-3921.	1.9	13

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19	Evaluation of a Conceptual Model for Gas-Particle Partitioning of Polycyclic Aromatic Hydrocarbons Using Polyparameter Linear Free Energy Relationships. <i>Environmental Science & Technology</i> , 2016, 50, 12312-12319.	10.0	46
20	New arrangement of dynamic permeation method for determination of gas separation ability of ionic liquids. <i>Separation and Purification Technology</i> , 2015, 147, 1-8.	7.9	9
21	Thermal Properties of Alkyl-triethylammonium bis(trifluoromethyl)sulfonyl imide Ionic Liquids. <i>Journal of Solution Chemistry</i> , 2015, 44, 790-810.	1.2	27
22	Solid-liquid equilibria in systems [C _x mim][Tf ₂ N] with diethylamine. <i>Pure and Applied Chemistry</i> , 2015, 87, 453-460.	1.9	5
23	Speeds of sound, isentropic compressibilities and refractive indices for some binary mixtures of nitromethane with chloroalkane at temperatures from 298.15 to 318.15K. Comparison with theories. <i>Fluid Phase Equilibria</i> , 2015, 385, 105-119.	2.5	15
24	Branched and cyclic alkyl groups in imidazolium-based ionic liquids: Molecular organization and physico-chemical properties. <i>Fluid Phase Equilibria</i> , 2014, 371, 41-49.	2.5	34
25	Using Partial Least-Squares Regression in Multivariate UV Spectroscopic Analysis of Mixtures of Imidazolium-Based Ionic Liquids and 1-Methylimidazole for Measurements of Liquid-Liquid Equilibria. <i>Journal of Solution Chemistry</i> , 2012, 41, 2164-2172.	1.2	2
26	Mutual Solubilities of Ammonium-Based Ionic Liquids with Water and with Water/Methanol Mixture. <i>Procedia Engineering</i> , 2012, 42, 1229-1241.	1.2	15
27	Phase Behaviour, Interactions, and Structural Studies of (Amines+Ionic Liquids) Binary Mixtures. <i>ChemPhysChem</i> , 2012, 13, 1825-1835.	2.1	24
28	Semi-empirical model of toluene transport in polyethylene membranes based on the data using a new type of apparatus for determining gas permeability, diffusivity and solubility. <i>Chemical Engineering Science</i> , 2011, 66, 5566-5574.	3.8	11
29	Dynamics of Atmospheric Aerosol Number Size Distributions in the Eastern Mediterranean During the AERO-Project. <i>Water, Air, and Soil Pollution</i> , 2011, 214, 133-146.	2.4	7
30	Thermodynamic description of liquid-liquid equilibria in systems 1-ethyl-3-methylimidazolium ethylsulfate+C7-hydrocarbons by polymer-solution models. <i>Fluid Phase Equilibria</i> , 2009, 284, 80-85.	2.5	16
31	Comparison of Two Approaches to Modeling Atmospheric Aerosol Particle Size Distributions. <i>Aerosol and Air Quality Research</i> , 2008, 8, 392-410.	2.1	2
32	Liquid-Liquid Equilibrium in Binary System [bmim][PF ₆] + 1-Butanol. <i>Journal of Chemical & Engineering Data</i> , 2006, 51, 2126-2131.	1.9	28
33	Vapour-liquid equilibrium at high pressure in the system containing carbon dioxide and propyl acetate. <i>Fluid Phase Equilibria</i> , 1995, 110, 175-182.	2.5	19
34	Possibility of pore size determination in separation layer of ceramic membrane using permeation method. <i>Journal of Membrane Science</i> , 1995, 103, 151-157.	8.2	9
35	Vapour-liquid equilibrium in the carbon dioxide-ethyl propanoate system at pressures from 2 to 9 MPa and temperatures from 303 to 323 K. <i>Fluid Phase Equilibria</i> , 1995, 112, 125-129.	2.5	11
36	Vapour-liquid equilibrium in the carbon dioxide-ethyl acetate system at high pressure. <i>Fluid Phase Equilibria</i> , 1994, 97, 119-126.	2.5	33

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37	Vapour-liquid equilibrium in the carbon dioxide \leftrightarrow p-cymene system at high pressure. Fluid Phase Equilibria, 1993, 90, 135-141.	2.5	8
38	Vapour-liquid equilibrium in the sulphur hexafluoride \leftrightarrow n-pentane system at high pressure. Fluid Phase Equilibria, 1990, 54, 35-45.	2.5	8
39	Nonclassical behaviour of binary mixtures in gas-liquid critical region and its quantitative description. Collection of Czechoslovak Chemical Communications, 1989, 54, 2863-2867.	1.0	0
40	High-pressure vapour-liquid equilibrium in systems containing carbon dioxide, 1-hexene, and n-hexane. Fluid Phase Equilibria, 1987, 33, 109-123.	2.5	89
41	Relation of the temperature derivative of heat of vaporization to the difference of heat capacities along the saturated vapour pressure curve. Collection of Czechoslovak Chemical Communications, 1981, 46, 2446-2454.	1.0	1