Ville Alopaeus

List of Publications by Citations

158 papers	**2,436** citations	**23** h-index	**42** g-index
162 ext. papers	**2,790** ext. citations	**3.8** avg, IF	**5.19** L-index

#	Paper	IF	Citations
158	Modelling aerosol transport and virus exposure with numerical simulations in relation to SARS-CoV-2 transmission by inhalation indoors. *Safety Science*, **2020**, 130, 104866	5.8	193
157	Simulation of the population balances for liquid–liquid systems in a nonideal stirred tank. Part 2—Parameter fitting and the use of the multiblock model for dense dispersions. *Chemical Engineering Science*, **2002**, 57, 1815-1825	4.4	158
156	Simulation of the population balances for liquid–liquid systems in a nonideal stirred tank. Part 1 Description and qualitative validation of the model. *Chemical Engineering Science*, **1999**, 54, 5887-5899	4.4	131
155	Modelling local bubble size distributions in agitated vessels. *Chemical Engineering Science*, **2007**, 62, 721-740	4.4	121
154	Validation of bubble breakage, coalescence and mass transfer models for gas–liquid dispersion in agitated vessel. *Chemical Engineering Science*, **2006**, 61, 218-228	4.4	120
153	CFD modeling of radial spreading of flow in trickle-bed reactors due to mechanical and capillary dispersion. *Chemical Engineering Science*, **2009**, 64, 207-218	4.4	75
152	Solubility of Organosolv Lignin in γ-Valerolactone/Water Binary Mixtures. *ChemSusChem*, **2016**, 9, 2939-2947	9.1	52
151	Modeling method for combining fluid dynamics and algal growth in a bubble column photobioreactor. *Chemical Engineering Journal*, **2013**, 229, 559-568	14.7	51
150	The Henry's law constant of N2O and CO2 in aqueous binary and ternary amine solutions (MEA, DEA, DIPA, MDEA, and AMP). *Fluid Phase Equilibria*, **2011**, 311, 59-66	2.5	51
149	Solution of population balances with breakage and agglomeration by high-order moment-conserving method of classes. *Chemical Engineering Science*, **2006**, 61, 6732-6752	4.4	47
148	Modelling mass transfer in an aerated 0.2m3 vessel agitated by Rushton, Phasejet and Combijet impellers. *Chemical Engineering Journal*, **2008**, 142, 95-108	14.7	44
147	Improved Hydrodynamic Model for Wetting Efficiency, Pressure Drop, and Liquid Holdup in Trickle-Bed Reactors. *Industrial & Engineering Chemistry Research*, **2008**, 47, 8436-8444	3.9	36
146	A Model for Chlorine Dioxide Delignification of Chemical Pulp. *Journal of Wood Chemistry and Technology*, **2010**, 30, 230-268	2	32
145	Novel micro-distillation column for process development. *Chemical Engineering Research and Design*, **2009**, 87, 705-710	5.5	32
144	Vapor–liquid Equilibria, Excess Enthalpy, and Density of Aqueous γ-Valerolactone Solutions.. *Journal of Chemical & Engineering Data*, **2016**, 61, 881-890	2.8	30
143	Dynamic modeling of local reaction conditions in an agitated aerobic fermenter. *AICHE Journal*, **2006**, 52, 1673-1689	3.6	28
142	Hydrogen solubility in heavy oil systems: Experiments and modeling. *Fuel*, **2014**, 137, 393-404	7.1	27

141	Solubility of carbon dioxide in aqueous solutions of diisopropanolamine and methyldiethanolamine. *Fluid Phase Equilibria*, **2010**, 293, 101-109	2.5	27
140	Numerical solution of moment-transformed population balance equation with fixed quadrature points. *Chemical Engineering Science*, **2006**, 61, 4919-4929	4.4	27
139	Measurements and modeling of CO2 solubility in 1,8-diazabicyclo-[5.4.0]-undec-7-ene Glycerol solutions. *Fluid Phase Equilibria*, **2014**, 374, 25-36	2.5	25
138	Compartmental modeling of large stirred tank bioreactors with high gas volume fractions. *Chemical Engineering Journal*, **2018**, 334, 2319-2334	14.7	25
137	Experimental and Theoretical Thermodynamic Study of Distillable Ionic Liquid 1,5-Diazabicyclo[4.3.0]non-5-enium Acetate. *Industrial & Engineering Chemistry Research*, **2016**, 55, 10445-10454	3.9	24
136	Catalyst Screening and Kinetic Modeling for CO Production by High Pressure and Temperature Reverse Water Gas Shift for Fischer Tropsch Applications. *Industrial & Engineering Chemistry Research*, **2017**, 56, 13262-13272	3.9	23
135	An Analytical Model for Capillary Pressure Saturation Relation for Gas liquid System in a Packed-Bed of Spherical Particles. *Transport in Porous Media*, **2009**, 77, 17-40	3.1	23
134	Lignin oxidation mechanisms under oxygen delignification conditions. Part 3. Reaction pathways and modeling. *Holzforschung*, **2011**, 65,	2	22
133	A new simple approach for the scale-up of aerated stirred tanks. *Chemical Engineering Research and Design*, **2015**, 95, 150-161	5.5	21
132	Heavy oil characterization method for PC-SAFT. *Fuel*, **2014**, 133, 216-223	7.1	21
131	Characteristics of liquid and tracer dispersion in trickle-bed reactors: Effect on CFD modeling and experimental analyses. *Computers and Chemical Engineering*, **2011**, 35, 41-49	4	21
130	Solution of population balances with growth and nucleation by high order moment-conserving method of classes. *Chemical Engineering Science*, **2007**, 62, 2277-2289	4.4	21
129	Chemical Recovery of Valerolactone/Water Biorefinery. *Industrial & Engineering Chemistry Research*, **2018**, 57, 15147-15158	3.9	20
128	Hydrogen solubility in furfural and 2-propanol: Experiments and modeling. *Journal of Chemical Thermodynamics*, **2017**, 112, 1-6	2.9	19
127	Development of a unique modular distillation column using 3D printing. *Chemical Engineering and Processing: Process Intensification*, **2016**, 109, 136-148	3.7	19
126	Vapor Pressures, Densities, and PC-SAFT Parameters for 11 Bio-compounds. *International Journal of Thermophysics*, **2019**, 40, 1	2.1	19
125	Analysis of stirred tanks with two-zone models. *AICHE Journal*, **2009**, 55, 2545-2552	3.6	19
124	Feasibility of thermal separation in recycling of the distillable ionic liquid [DBNH][OAc] in cellulose fiber production. *Chemical Engineering Research and Design*, **2016**, 114, 287-298	5.5	19

123	Investigation of drag models in CFD modeling and comparison to experiments of liquid-solid fluidized systems. *International Journal of Mineral Processing*, **2012**, 104-105, 58-70		18
122	Measurements of H2S solubility in aqueous diisopropanolamine solutions and vapour pressure of diisopropanolamine. *Fluid Phase Equilibria*, **2013**, 338, 164-171	2.5	17
121	A dynamic model for plug flow reactor state profiles. *Computers and Chemical Engineering*, **2008**, 32, 1494-1506	4	17
120	Approximate high flux corrections for multicomponent mass transfer models and some explicit methods. *Chemical Engineering Science*, **1999**, 54, 4267-4271	4.4	17
119	Application of the compartmental model to the gas-liquid precipitation of CO2-Ca(OH)2 aqueous system in a stirred tank. *AICHE Journal*, **2017**, 63, 378-386	3.6	16
118	Ternary and Binary LLE Measurements for Solvent (4-Methyl-2-pentanone and 2-Methyl-2-butanol) + Furfural + Water between 298 and 401 K. *Journal of Chemical & Engineering Data*, **2016**, 61, 903-911	2.8	16
117	A comprehensive thermodynamic study of heat stable acetic acid salt of monoethanolamine. *International Journal of Greenhouse Gas Control*, **2014**, 22, 313-324	4.2	16
116	Thermodynamics of aqueous solutions of methyldiethanolamine and diisopropanolamine. *Fluid Phase Equilibria*, **2010**, 299, 51-59	2.5	16
115	Modeling of nickel-based hydrotalcite catalyst coated on heat exchanger reactors for CO2 methanation. *Chemical Engineering Journal*, **2018**, 349, 694-707	14.7	16
114	Ternary and binary LLE measurements for solvent (2-methyltetrahydrofuran and cyclopentyl methyl ether) + furfural + water between 298 and 343 K. *Journal of Chemical Thermodynamics*, **2017**, 110, 127-136	2.9	14
113	Vapor-liquid Equilibrium, Excess Molar Enthalpies, and Excess Molar Volumes of Binary Mixtures Containing Methyl Isobutyl Ketone (MIBK) and 2-Butanol, tert-Pentanol, or 2-Ethyl-1-hexanol. *Journal of Chemical & Engineering Data*, **2012**, 57, 3092-3101	2.8	14
112	A calculation method for multicomponent mass transfer coefficient correlations. *Computers and Chemical Engineering*, **1999**, 23, 1177-1182	4	14
111	Detailed Modeling of Kraft Pulping Chemistry. Delignification. *Industrial & Engineering Chemistry Research*, **2020**, 59, 12977-12985	3.9	13
110	Vapor pressure, vapor-liquid equilibria, liquid-liquid equilibria and excess enthalpy of the system consisting of isophorone, furfural, acetic acid and water. *Chemical Engineering Science*, **2018**, 176, 19-34	4.4	13
109	Modeling of Gas-liquid Packed-Bed Reactors with Momentum Equations and Local Interaction Closures. *Industrial & Engineering Chemistry Research*, **2006**, 45, 8189-8198	3.9	13
108	Mass-transfer calculation methods for transient diffusion within particles. *AICHE Journal*, **2000**, 46, 2369-2372	3.6	13
107	Model for Degradation of Galactoglucomannan in Hot Water Extraction Conditions. *Industrial & Engineering Chemistry Research*, **2012**, 51, 10338-10344	3.9	12
106	The moment method for one-dimensional dynamic reactor models with axial dispersion. *Computers and Chemical Engineering*, **2011**, 35, 423-433	4	12

105	The Kinetics and Stoichiometry of the Reaction between Hypochlorous Acid and Chlorous Acid in Mildly Acidic Solutions. *Industrial & Engineering Chemistry Research*, **2009**, 48, 6280-6286	3.9	12
104	Solubility of hydrogen in bio-oil compounds. *Journal of Chemical Thermodynamics*, **2016**, 102, 406-412	2.9	12
103	Population balance based modeling of changes in cellulose molecular weight distribution during ageing. *Cellulose*, **2015**, 22, 151-163	5.5	10
102	Mathematical model of precipitation of magnesium carbonate with carbon dioxide from the magnesium hydroxide slurry. *Computers and Chemical Engineering*, **2016**, 87, 180-189	4	10
101	Dew points of pure DBN and DBU and vapor⎕liquid equilibria of water+DBN and water+DBU systems for cellulose solvent recycling. *Fluid Phase Equilibria*, **2016**, 408, 79-87	2.5	10
100	Physical Properties of 7-Methyl-1,5,7-triazabicyclo[4.4.0]dec-5-ene (mTBD). *International Journal of Thermophysics*, **2019**, 40, 1	2.1	10
99	Control of reflux and reboil flow rates for milli and micro distillation. *Chemical Engineering Research and Design*, **2013**, 91, 753-760	5.5	10
98	Distillable Protic Ionic Liquid 2-(Hydroxy)ethylammonium Acetate (2-HEAA): Density, Vapor Pressure, Vapor⎕liquid Equilibrium, and Solid⎕liquid Equilibrium. *Industrial & Engineering Chemistry Research*, **2014**, 53, 19322-19330	3.9	10
97	Vapor⎕liquid Equilibrium at 350 K, Excess Molar Enthalpies at 298 K, and Excess Molar Volumes at 298 K of Binary Mixtures Containing Ethyl Acetate, Butyl Acetate, and 2-Butanol. *Journal of Chemical & Engineering Data*, **2013**, 58, 1011-1019	2.8	10
96	The role of Donnan effect in kraft cooking liquor impregnation and hot water extraction of wood. *Holzforschung*, **2013**, 67, 511-521	2	10
95	Vapor⎕liquid Equilibrium for Methoxymethane + Methyl Formate, Methoxymethane + Hexane, and Methyl Formate + Methanol. *Journal of Chemical & Engineering Data*, **2011**, 56, 2634-2640	2.8	10
94	Phase equilibria of binary systems of 3-methylthiophene with four different hydrocarbons. *Fluid Phase Equilibria*, **2010**, 288, 155-160	2.5	10
93	Modeling outdoors algal cultivation with compartmental approach. *Chemical Engineering Journal*, **2015**, 259, 945-960	14.7	9
92	Quaternary and ternary LLE measurements for solvent (2-methyltetrahydrofuran and cyclopentyl methyl ether) + furfural + acetic acid + water between 298 and 343 K. *Journal of Chemical Thermodynamics*, **2018**, 119, 61-75	2.9	9
91	Quaternary, Ternary, and Binary LLE Measurements for 2-Methoxy-2-methylpropane + Furfural + Acetic Acid + Water at Temperatures between 298 and 307 K. *Journal of Chemical & Engineering Data*, **2016**, 61, 2458-2466	2.8	9
90	Isothermal vapor⎕lquid equilibrium and excess molar enthalpies of the binary mixtures furfural+methyl isobutyl ketone, +2-butanol and +2-methyl-2-butanol. *Fluid Phase Equilibria*, **2014**, 372, 85-99	2.5	9
89	Measurements and modeling of LLE and HE for (methanol + 2,4,4-trimethyl-1-pentene), and LLE for (water + methanol + 2,4,4-trimethyl-1-pentene). *Journal of Chemical Thermodynamics*, **2015**, 85, 120-128	2.9	9
88	Vapor⎕liquid Equilibrium for Methoxymethane + Thiophene, + Diethylsulfide, + 2-Methyl-2-propanethiol and 1-Hexene, + 1-Propanethiol. *Journal of Chemical & Engineering Data*, **2013**, 58, 956-963	2.8	9

87	Modelling of crystal growth of KDP in a 100dm3 suspension crystallizer using combination of CFD and multiblock model. *Chemical Engineering Research and Design*, **2010**, 88, 1297-1303	5.5	9
86	Solution of population balances by high order moment-conserving method of classes: reconstruction of a non-negative density distribution. *Chemical Engineering Science*, **2008**, 63, 2741-2751	4.4	9
85	Modelling a molten salt thermal energy system ⬚A validation study. *Applied Energy*, **2019**, 233-234, 126-145	15.7	9
84	Modelling and dynamic simulation of a large MSF plant using local phase equilibrium and simultaneous mass, momentum, and energy solver. *Computers and Chemical Engineering*, **2017**, 97, 242-258	4	8
83	Population balance model and experimental validation for reactive dissolution of particle agglomerates. *Computers and Chemical Engineering*, **2018**, 108, 240-249	4	8
82	A modified continuous flow apparatus for gas solubility measurements at high pressure and temperature with camera system. *Fluid Phase Equilibria*, **2014**, 382, 150-157	2.5	8
81	Analysis of concentration polydispersity in mixed liquid⬚quid systems. *Chemical Engineering Research and Design*, **2014**, 92, 612-618	5.5	8
80	Vapor⬚iquid Equilibria, Excess Enthalpy, and Excess Volume of Binary Mixtures Containing an Alcohol (1-Butanol, 2-Butanol, or 2-Methyl-2-butanol) and 2-Ethoxy-2-methylbutane. *Journal of Chemical & Engineering Data*, **2012**, 57, 3502-3509	2.8	8
79	Appropriate Simplifications in Calculation of Mass Transfer in a Multicomponent Rate-Based Distillation Tray Model. *Industrial & Engineering Chemistry Research*, **2000**, 39, 4336-4345	3.9	8
78	Vapor⬚iquid Equilibrium of Ionic Liquid 7-Methyl-1,5,7-triazabicyclo[4.4.0]dec-5-enium Acetate and Its Mixtures with Water. *Journal of Chemical & Engineering Data*, **2020**, 65, 2405-2421	2.8	7
77	Detailed modeling of the kraft pulping chemistry: carbohydrate reactions. *AICHE Journal*, **2020**, 66, e16252	3.6	7
76	The effect of microwave on the crystallization process of magnesium carbonate from aqueous solutions. *Powder Technology*, **2018**, 328, 358-366	5.2	7
75	Comparative study: Absorption enthalpy of carbon dioxide into aqueous diisopropanolamine and monoethanolamine solutions and densities of the carbonated amine solutions. *Fluid Phase Equilibria*, **2014**, 376, 85-95	2.5	7
74	Hydrogen Solubility of Shale Oil Containing Polar Phenolic Compounds. *Industrial & Engineering Chemistry Research*, **2017**, 56, 8738-8747	3.9	7
73	Comparison and validation of CFD models in liquid⬚olid suspensions. *Canadian Journal of Chemical Engineering*, **2011**, 89, 696-706	2.3	7
72	Isobaric vapor⬚iquid equilibrium for binary systems containing benzothiophene. *Fluid Phase Equilibria*, **2011**, 307, 180-184	2.5	7
71	Modeling and Simulation of an Industrial Trickle-Bed Reactor for Benzene Hydrogenation: Model Validation against Plant Data. *Industrial & Engineering Chemistry Research*, **2009**, 48, 1866-1872	3.9	7
70	Phase equilibria on five binary systems containing 1-butanethiol and 3-methylthiophene in hydrocarbons. *Fluid Phase Equilibria*, **2010**, 293, 157-163	2.5	7

69	Utilization of population balances in simulation of liquid-liquid systems in mixed tanks. *Chemical Engineering Communications*, **2003**, 190, 1468-1484	2.2	7
68	Moment based weighted residual method—New numerical tool for a nonlinear multicomponent chromatographic general rate model. *Computers and Chemical Engineering*, **2013**, 53, 153-163	4	6
67	A novel continuous flow apparatus with a video camera system for high pressure phase equilibrium measurements. *Fluid Phase Equilibria*, **2013**, 356, 291-300	2.5	6
66	Hydrogen solubility measurements of analyzed tall oil fractions and a solubility model. *Journal of Chemical Thermodynamics*, **2017**, 105, 15-20	2.9	6
65	Vapor–liquid Equilibrium for Dimethyl Disulfide + Butane, + trans-But-2-ene, + 2-Methylpropane, + 2-Methylpropene, + Ethanol, and 2-Ethoxy-2-methylpropane. *Journal of Chemical & Engineering Data*, **2011**, 56, 2501-2510	2.8	6
64	Phase equilibria on binary systems containing diethyl sulfide. *Fluid Phase Equilibria*, **2011**, 301, 200-205	2.5	6
63	Vapour–liquid equilibrium for the systems diethyl sulphide+1-butene, +cis-2-butene, +2-methylpropane, +2-methylpropene, +n-butane, +trans-2-butene. *Fluid Phase Equilibria*, **2010**, 291, 180-187	2.5	6
62	Vapor–liquid Equilibrium for the Systems trans-2-Butene + Methanol, + 1-Propanol, + 2-Propanol, + 2-Butanol, and + 2-Methyl-2-propanol at 364.5 K. *Journal of Chemical & Engineering Data*, **2008**, 53, 607-612	2.8	6
61	Vapor–liquid Equilibrium for 1-Butene + Methanol, + 1-Propanol, + 2-Propanol, + 2-Butanol, and 2-Methyl-2-propanol (TBA) at 364.5 K. *Journal of Chemical & Engineering Data*, **2008**, 53, 1829-1835	2.8	6
60	A cellular automata model for liquid distribution in trickle bed reactors. *Chemical Engineering Science*, **2006**, 61, 4930-4943	4.4	6
59	On approximate calculation of multicomponent mass transfer fluxes. *Computers and Chemical Engineering*, **2002**, 26, 461-466	4	6
58	A high-order moment-conserving method of classes (HMMC) based population balance model for mechanical flotation cells. *Minerals Engineering*, **2017**, 108, 36-52	4.9	5
57	The use of microplants in process development—Case study of etherification of 2-ethoxy-2-methylbutane. *Chemical Engineering and Processing: Process Intensification*, **2013**, 74, 75-82	3.7	5
56	Measurement of activity coefficient at infinite dilution for some bio-oil components in water and mass transfer study of bubbles in the dilutor. *Fluid Phase Equilibria*, **2015**, 392, 1-11	2.5	5
55	Vapor–liquid Equilibrium, Excess Molar Enthalpies, and Excess Molar Volumes of Binary Mixtures Containing 2-Ethoxy-2-methylpropane or 2-Ethoxy-2-methylbutane and Acetonitrile or Propanenitrile. *Journal of Chemical & Engineering Data*, **2013**, 58, 943-950	2.8	5
54	Isothermal Binary Vapor–liquid Equilibrium for 2-Methylpropane and n-Butane with 1,2-Ethanedithiol and 2-Methyl-2-propanethiol. *Journal of Chemical & Engineering Data*, **2010**, 55, 291-296	2.8	5
53	Vapor–liquid Equilibrium for Tetrahydrothiophene + n-Butane, + trans-2-Butene, + 2-Methylpropane, and + 2-Methylpropene. *Journal of Chemical & Engineering Data*, **2009**, 54, 1311-1317	2.8	5
52	Phase equilibrium measurements for systems containing propanenitrile with tert-butyl ethyl ether and C4-hydrocarbons. *Fluid Phase Equilibria*, **2010**, 299, 148-160	2.5	5

51	Phase equilibria for systems containing dimethyl disulfide and diethyl disulfide with hydrocarbons at 368.15K. *Fluid Phase Equilibria*, **2010**, 293, 175-181	2.5	5
50	Vapor–liquid Equilibrium for the cis-2-Butene + Methanol, + 2-Propanol, + 2-Butanol, + 2-Methyl-2-propanol Systems at 364.5 K. *Journal of Chemical & Engineering Data*, **2008**, 53, 1539-1544	2.8	5
49	Physicochemical Modeling for Hot Water Extraction of Birch Wood. *Industrial & Engineering Chemistry Research*, **2016**, 55, 11062-11073	3.9	5
48	Design of Equilibrium Cells for Phase Equilibria and PVT Measurements in Large Ranges of Temperatures and Pressures. I. Vapor–liquid–liquid Equilibria. *Journal of Chemical & Engineering Data*, **2016**, 61, 2700-2711	2.8	5
47	Isobaric Vapor–liquid Equilibrium of Furfural + –Valerolactone at 30 kPa and Isothermal Liquid–liquid Equilibrium of Carbon Dioxide + –Valerolactone + Water at 298 K. *Journal of Chemical & Engineering Data*, **2018**,	2.8	5
46	Quaternary, Ternary and Binary LLE Measurements for 2-Methoxy-2-methylbutane + Furfural + Acetic Acid + Water at Temperatures between 298 and 341 K. *Journal of Chemical & Engineering Data*, **2016**, 61, 3730-3739	2.8	4
45	Prototyping a calorimeter mixing cell with direct metal laser sintering. *Chemical Engineering Research and Design*, **2016**, 108, 146-151	5.5	4
44	Modeling of Mass Transfer and Reactions in Anisotropic Biomass Particles with Reduced Computational Load. *Industrial & Engineering Chemistry Research*, **2014**, 53, 4096-4103	3.9	4
43	Solubility of carbon monoxide in bio-oil compounds. *Journal of Chemical Thermodynamics*, **2017**, 105, 296-311	2.9	4
42	Chlorate Formation in Chlorine Dioxide Delignification–An Analysis via Elementary Kinetic Modeling. *Journal of Wood Chemistry and Technology*, **2009**, 29, 191-213	2	4
41	Vapor–liquid equilibrium for the binary systems tetrahydrothiophene + toluene and tetrahydrothiophene + o-xylene at 368.15 K and 383.15 K. *Fluid Phase Equilibria*, **2010**, 296, 4-8	2.5	4
40	Physicochemical Modeling for Pressurized Hot Water Extraction of Spruce Wood. *Industrial & Engineering Chemistry Research*, **2018**, 57, 16664-16674	3.9	4
39	A comprehensive study of CO 2 solubility in aqueous 2-HEAA and MEA + 2-HEAA solutions – Measurements and modeling. *International Journal of Greenhouse Gas Control*, **2015**, 42, 296-306	4.2	3
38	Temperature and Pressure Dependence of Density of a Shale Oil and Derived Thermodynamic Properties. *Industrial & Engineering Chemistry Research*, **2018**, 57, 5128-5135	3.9	3
37	110th Anniversary: Critical Properties and High Temperature Vapor Pressures for Furan, 2-Methylfuran, 2-Methoxy-2-methylpropane, 2-Ethoxy-2-methylbutane, n-Hexane, and Ethanol and Bubble Points of Mixtures with a New Apparatus. *Industrial & Engineering Chemistry Research*, **2019**, 58, 22350-22362	3.9	3
36	A new moment analysis method to estimate the characteristic parameters in chromatographic general rate model. *Computers and Chemical Engineering*, **2013**, 55, 50-60	4	3
35	Vapor–liquid Equilibrium for Thiophene + Butane, + trans-But-2-ene, + 2-Methylpropane, and + 2-Methylpropene. *Journal of Chemical & Engineering Data*, **2011**, 56, 614-621	2.8	3
34	Three-Phase CFD-Model for Trickle Bed Reactors. *International Journal of Nonlinear Sciences and Numerical Simulation*, **2012**, 13,	1.8	3

33	Infinite dilution activity coefficient and vapour liquid equilibrium measurements for dimethylsulphide and tetrahydrothiophene with hydrocarbons. *Fluid Phase Equilibria*, **2010**, 295, 17-25	2.5	3
32	Vapor–liquid Equilibrium for the Systems 2-Methylpropane + Methanol, + 2-Propanol, + 2-Butanol, and + 2-Methyl-2-propanol at 364.5 K. *Journal of Chemical & Engineering Data*, **2008**, 53, 913-918	2.8	3
31	Modeling surfactant and drop size dynamics in polydisperse liquid-liquid systems with population balances. *Chemical Engineering Science*, **2022**, 248, 117269	4.4	3
30	CPMAS 13C NMR Analysis of Fully Bleached Eucalypt Pulp Samples: Links to Handsheet Hygroexpansivity and Strength Properties. *Journal of Applied Sciences*, **2009**, 9, 3991-3998	0.3	3
29	Densities, Viscosities, and Thermal Conductivities of the Ionic Liquid 7-Methyl-1,5,7-triazabicyclo[4.4.0]dec-5-enium Acetate and Its Mixtures with Water. *International Journal of Thermophysics*, **2020**, 41, 1	2.1	3
28	Vapor- liquid equilibrium for the n-dodecane + phenol and n-hexadecane + phenol systems at 523 K and 573 K. *Fluid Phase Equilibria*, **2021**, 537, 112991	2.5	3
27	Toward solvent screening in the extractive desulfurization using ionic liquids: QSPR modeling and experimental validations. *Fuel*, **2021**, 302, 121159	7.1	3
26	Micro-scale piloting of a process for production of 2-methoxy-2,4,4-trimethylpentane. *Chemical Engineering and Processing: Process Intensification*, **2017**, 122, 143-154	3.7	2
25	Isothermal and Isobaric Vapor–liquid Equilibrium and Excess Molar Enthalpy of the Binary Mixtures of 2-Methoxy-2-methylpropane + 2-Methyl-2-butanol or + 2-Butanol. *Journal of Chemical & Engineering Data*, **2015**, 60, 2655-2664	2.8	2
24	Application of GaInSn Liquid Metal Alloy Replacing Mercury in a Phase Equilibrium Cell: Vapor Pressures of Toluene, Hexylbenzene, and 2-Ethylnaphthalene. *Journal of Chemical & Engineering Data*, **2020**, 65, 3270-3276	2.8	2
23	Hydrodeoxygenation Model Compounds Heptalactone and Nonalactone: Density from 293 to 473 K and H2 Solubility from 479 to 582 K. *Journal of Chemical & Engineering Data*, **2020**, 65, 2764-2773	2.8	2
22	Analytical time-stepping solution of the discretized population balance equation. *Computers and Chemical Engineering*, **2020**, 135, 106741	4	2
21	Study of CO2 Absorption into Phase Change Solvents MAPA and DEEA. *Journal of Chemical & Engineering Data*, **2017**, 62, 2261-2271	2.8	2
20	Modeling of alkaline extraction chemistry and kinetics of softwood kraft pulp. *Holzforschung*, **2014**, 68, 733-746	2	2
19	Modeling of Catalyst Activity Profiles in Fixed-Bed Reactors with a Moment Transformation Method. *Industrial & Engineering Chemistry Research*, **2008**, 47, 8192-8196	3.9	2
18	Vapor–liquid Equilibrium for 1-Butanol + 1-Butene at (318.4 and 364.5) K and Vapor–liquid Equilibrium of 1-Butanol + 2-Methylpropane, + n-Butane and 1-Butene + 2-Methylpropane at 318.4 K. *Journal of Chemical & Engineering Data*, **2008**, 53, 2454-2461	2.8	2
17	Application-Related Consideration of the Thermal Stability of [mTBDH][OAc] Compared to Amidine-Based Ionic Liquids in the Presence of Various Amounts of Water. *Industrial & Engineering Chemistry Research*, **2022**, 61, 259-268	3.9	2
16	Volumetric Properties of Aqueous Solutions of Zinc Sulfate at Temperatures from 298.15 to 393.15 K and Pressures up to 10 MPa. *Journal of Chemical & Engineering Data*, **2021**, 66, 45-57	2.8	2

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15	Modelling the kraft pulping process on a fibre scale by considering the intrinsic heterogeneous nature of the lignocellulosic feedstock. *Chemical Engineering Journal*, **2022**, 438, 135548	14.7	2
14	CFD analysis of laboratory scale phase equilibrium cell operation. *Review of Scientific Instruments*, **2017**, 88, 105110	1.7	1
13	Approximating Catalyst Effectiveness Factors with Reaction Rate Profiles. *Catalysts*, **2019**, 9, 255	4	1
12	Measurements and modeling for the density of 2-methoxy-2,4,4-trimethylpentane, HE for (methanol + 2-methoxy-2,4,4-trimethylpentane), LLE for (water + 2-methoxy-2,4,4-trimethylpentane) and LLE for (water + methanol + 2-methoxy-2,4,4-trimethylpentane). *Journal of Chemical Thermodynamics*, **2015**, 91, 313-320	2.9	1
11	Vapor–liquid equilibrium measurements of dimethylsulfide, +ethanol, +dimethylether, +methylacetate with a static total pressure method. *Fluid Phase Equilibria*, **2013**, 355, 34-39	2.5	1
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