Deresh Ramjugernath

List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

5,356 312 35 59 h-index g-index citations papers 6,057 3.2 317 5.93 L-index avg, IF ext. citations ext. papers

#	Paper	IF	Citations
312	Purification of Nitrogen Trifluoride by Physical Separation. <i>Fluid Phase Equilibria</i> , 2022 , 113405	2.5	
311	Experimental measurements of CO2 solubility, viscosity, density, sound velocity and evaporation rate for 2-(2-aminoethoxy)ethanol (DGA) + 1-methylpyrrolidin-2-one[(NMP) / water + ionic liquid systems. Fluid Phase Equilibria, 2022, 113475	2.5	
310	The Effect of Cobalt Catalyst Loading at Very High Pressure Plasma-Catalysis in Fischer-Tropsch Synthesis. <i>Catalysts</i> , 2021 , 11, 1324	4	O
309	Investigation of Mixed MEA-Based Solvents Featuring Ionic Liquids and NMP for CO2 Capture: Experimental Measurement of CO2 Solubility and Thermophysical Properties. <i>Journal of Chemical & Engineering Data</i> , 2021 , 66, 899-914	2.8	2
308	Plasma-Catalytic Fischer Tropsch Synthesis at Very High Pressure. Catalysts, 2021, 11, 297	4	2
307	Thermodynamic measurement and modeling of hydrate dissociation for CO2/refrigerant + sucrose/fructose/glucose solutions. <i>AICHE Journal</i> , 2021 , 67, e17379	3.6	2
306	Gas hydrate concentration measurements on sucrose solutions using a new pilot test rig. <i>AICHE Journal</i> , 2020 , 66, e16281	3.6	4
305	Effect of temperature on molecular interactions between tri(butyl)methylphosphonium methylsulfate and furfural. <i>Journal of Chemical Thermodynamics</i> , 2020 , 149, 106150	2.9	3
304	Separation of thiophene from octane/hexadecane with ionic liquids in ternary liquid-liquid phase equilibrium. <i>Fluid Phase Equilibria</i> , 2020 , 509, 112467	2.5	8
303	Estimation of Pure Component Properties, Part 5: Estimation of the Thermal Conductivity of Nonelectrolyte Organic Liquids via Group Contributions. <i>Journal of Chemical & Data</i> , 2020 , 65, 1300-1312	2.8	4
302	Critical analysis of the effect of transport phenomena and operational parameters on the performance of an intermediate-scale surface fluorination reactor. <i>Journal of Fluorine Chemistry</i> , 2020 , 237, 109617	2.1	
301	Measurement and Modeling of the Solubility of Tetrafluoromethane in Either Perfluoroheptane or Perfluorodecalin. <i>Journal of Chemical & Engineering Data</i> , 2020 , 65, 4862-4868	2.8	1
300	Experimental measurements and thermodynamic modelling of hydrate phase equilibrium conditions for CF4+TBAB aqueous solutions. <i>Chemical Engineering Communications</i> , 2020 , 207, 185-193	2.2	1
299	The distribution coefficients of Y3+ and Eu3+ between HNO3 and HDEHP. <i>Minerals Engineering</i> , 2020 , 153, 106285	4.9	
298	Solubility Data for Roflumilast and Maraviroc in Various Solvents between T = (278.2B23.2) K. <i>Journal of Chemical & Data</i> , 2019, 64, 4599-4604	2.8	1
297	Isothermal Vaporliquid Equilibrium Data for the Binary Systems Consisting of 1,1,2,3,3,3-Hexafluoro-1-propene and Either Methylcyclohexane, Cyclohexane, n-Hexane, 2-Methyltetrahydrofuran, or 2,2,3,3,4,4,4-Heptafluoro-1-butanol. <i>Journal of Chemical & Company</i> 2010, 2010	2.8	О
296	Engineering Data, 2019 , 64, 5232-5237 Hydrate Dissociation Data for the Systems (CO2/CH4/Ar) + Water with (TBAF/TBAA/TBPB/TBANO3 and Cyclopentane). <i>Journal of Chemical & Data</i> , 2019, 64, 2542-2549	2.8	1

(2019-2019)

295	Phase Stability Conditions for Clathrate Hydrates Formation in CO2 + (NaCl or CaCl2 or MgCl2) + Cyclopentane + Water Systems: Experimental Measurements and Thermodynamic Modeling. Journal of Chemical & Chemical & Camp; Engineering Data, 2019, 64, 4638-4646	2.8	7	
294	Experimental study of carbon dioxide gas hydrate formation in the presence of zwitterionic compounds. <i>Journal of Chemical Thermodynamics</i> , 2019 , 137, 94-100	2.9	3	
293	Investigation of temperature and composition dependence of molecular interactions between phosphonium-based ionic liquid + N, N-dimethylformamide: A study of thermophysical properties. <i>Journal of Molecular Liquids</i> , 2019 , 291, 110987	6	3	
292	Vapor Liquid Equilibrium Data for 2,3-Pentanedione + (Acetaldehyde or Acetone) at (100, 150, and 200) kPa. <i>Journal of Chemical & Data</i> , 2019, 64, 2388-2394	2.8	2	
291	Phase stability conditions for clathrate hydrate formation in (fluorinated refrigerant + water + single and mixed electrolytes + cyclopentane) systems: Experimental measurements and thermodynamic modelling. <i>Journal of Chemical Thermodynamics</i> , 2019 , 136, 59-76	2.9	10	
290	Isothermal Vaporliquid Equilibrium Measurements for Alcohol + Water/n-Hexane Azeotropic Systems Using Both Dynamic and Automated Static-Synthetic Methods. <i>Journal of Chemical & Engineering Data</i> , 2019 , 64, 2657-2670	2.8	2	
289	Ternary liquid-liquid phase equilibria of {ionic liquid + thiophene + (octane/hexadecane)}. <i>Journal of Chemical Thermodynamics</i> , 2019 , 134, 157-163	2.9	8	
288	Characterization of C7+ fraction properties of crude oils and gas-condensates using data driven models. <i>Petroleum Science and Technology</i> , 2019 , 37, 1516-1522	1.4	3	
287	Evaluation of wax disappearance temperatures in hydrocarbon fluids using soft computing approaches. <i>Petroleum Science and Technology</i> , 2019 , 37, 829-836	1.4	3	
286	Experimental study and modeling of the kinetics of gas hydrate formation for acetylene, ethylene, propane and propylene in the presence and absence of SDS. <i>Petroleum Science and Technology</i> , 2019 , 37, 506-512	1.4	2	
285	Investigation into the use of gas hydrate technology for the treatment of vinasse. <i>Fluid Phase Equilibria</i> , 2019 , 492, 67-77	2.5	5	
284	State of the art and kinetics of refrigerant hydrate formation. <i>International Journal of Refrigeration</i> , 2019 , 98, 410-427	3.8	17	
283	Solid Liquid Equilibrium Measurements for Posaconazole and Voriconazole in Several Solvents between T = 278.2 and 323.2 K Using Differential Thermal Analysis/Thermal Gravimetric Analysis. <i>Journal of Chemical & Data</i> , 2019, 64, 3367-3374	2.8	5	
282	Ligand Free Heterogeneous Sonogashira Cross-Coupling Reaction over an in Situ Organoiodine Capsulized Palladium Anchored to a Perovskite Catalyst. <i>ACS Sustainable Chemistry and Engineering</i> , 2019 , 7, 12697-12706	8.3	7	
281	An insight into thermodynamic consistency of hydrogen sulfide sulfur content data at isobaric conditions. <i>Petroleum Science and Technology</i> , 2019 , 37, 763-769	1.4	1	
2 80	Fabrication, physical and optical properties of functionalized cellulose based polymethylmethacrylate nanocomposites. <i>Microsystem Technologies</i> , 2019 , 1	1.7		
279	VLE measurements and modelling for the binary systems of (CF41+ C6F14) and (CF4+ C8F18). <i>Fluid Phase Equilibria</i> , 2019 , 485, 146-152	2.5	2	
278	Vapor l iquid Equilibrium Measurements of Ether Alcohol Blends for Investigation on Reformulated Gas. <i>Journal of Chemical & Engineering Data</i> , 2019 , 64, 115-123	2.8	2	

277	Effect of temperature on intermolecular interactions between the organic solvents: Insights from density and excess volume. <i>Journal of Chemical Thermodynamics</i> , 2019 , 132, 461-469	2.9	5
276	Isothermal vapour-liquid equilibrium data for the binary systems of CHF 3 with (n-nonane, n-decane, or n-undecane) and C 2 F 6 with (n-nonane or n-decane). <i>Fluid Phase Equilibria</i> , 2018 , 464, 64-	78 ^{.5}	2
275	Valorisation of chicken feather barbs: Utilisation in yarn production and technical textile applications. <i>Sustainable Chemistry and Pharmacy</i> , 2018 , 8, 38-49	3.9	17
274	Valorisation of waste chicken feathers: Optimisation of decontamination and pre-treatment with bleaching agents using response surface methodology. <i>Sustainable Chemistry and Pharmacy</i> , 2018 , 8, 21-37	3.9	18
273	Valorisation of mango seed via extraction of starch: preliminary techno-economic analysis. <i>Clean Technologies and Environmental Policy</i> , 2018 , 20, 81-94	4.3	15
272	Recent Advances in Chitosan-Based Films for Novel Biosensor 2018 , 137-161		3
271	Experimental Phase Equilibrium for the Binary System ofn-Pentane +2-Propanol Using a New Equilibrium Cell and the Static Total Pressure Method. <i>Journal of Chemical & Data</i> , 2018, 63, 732-740	2.8	1
270	Application of Decafluorobiphenyl (DFBP) Moiety as a Linker in Bioconjugation. <i>Bioconjugate Chemistry</i> , 2018 , 29, 225-233	6.3	5
269	Development of a Computational Tool for the Analysis and Synthesis of Crystallization Processes. Organic Process Research and Development, 2018 , 22, 219-227	3.9	1
268	Isothermal vapour-liquid equilibrium data for binary systems of (CHF 3 or C 2 F 6) with (1-hexene or 3-methylpentane). <i>Journal of Chemical Thermodynamics</i> , 2018 , 121, 79-90	2.9	2
267	Data-driven modeling for determination of asphaltene stability condition in oil system. <i>Petroleum Science and Technology</i> , 2018 , 36, 726-731	1.4	O
266	Isothermal Vaporlliquid Equilibrium Data for Binary Mixtures of Hexafluoroethane (R116) + n-Pentane or n-Hexane at Two Temperatures, 288 and 296 K. <i>Journal of Chemical & Data</i> , 2018, 63, 1228-1233	2.8	2
265	Kinetic study of hydrate formation for argon + TBAB + SDS aqueous solution system. <i>Journal of Chemical Thermodynamics</i> , 2018 , 116, 121-129	2.9	13
264	Beneficiation of pulp and paper mill sludge: production and characterisation of functionalised crystalline nanocellulose. <i>Clean Technologies and Environmental Policy</i> , 2018 , 20, 1835-1845	4.3	14
263	Modeling of Trifluoromethane (R-23) or Hexafluoroethane (R-116) and Alkane Binary Mixtures using the Group-Contribution with Association Equation of State. <i>Industrial & Engineering Chemistry Research</i> , 2018 , 57, 10640-10648	3.9	0
262	Isothermal Vaporliquid Equilibrium Data for Binary Systems of CHF3 or C2F6 with Methylcyclohexane or Toluene. <i>Journal of Chemical & Engineering Data</i> , 2018 , 63, 2114-2126	2.8	1
261	Model evaluation for the prediction of solubility of active pharmaceutical ingredients (APIs) to guide solid-liquid separator design. <i>Asian Journal of Pharmaceutical Sciences</i> , 2018 , 13, 265-278	9	6
260	Optimisation of surfactant decontamination and pre-treatment of waste chicken feathers by using response surface methodology. <i>Waste Management</i> , 2018 , 72, 371-388	8.6	9

259	Can 2-methyl-2-butene and isoprene form clathrate hydrates?. <i>Petroleum Science and Technology</i> , 2018 , 36, 1696-1702	1.4		
258	Perfluorophenyl Derivatives as Unsymmetrical Linkers for Solid Phase Conjugation. <i>Frontiers in Chemistry</i> , 2018 , 6, 589	5	2	
257	PII Data and Modeling for Propan-1-ol + n-Octane or n-Nonane or n-Decane from 313.15 K to 363.15 K and 1 MPa to 20 MPa. <i>Journal of Chemical & Engineering Data</i> , 2018, 63, 4136-4156	2.8	6	
256	Isothermal Vaporliquid Equilibrium for the 2-Pentanone (MPK) + 2-Methyl Propan-1-ol Binary Mixture. <i>Journal of Chemical & Engineering Data</i> , 2018 , 63, 4076-4084	2.8	2	
255	Valorisation of avocado seeds: extraction and characterisation of starch for textile applications. <i>Clean Technologies and Environmental Policy</i> , 2018 , 20, 2135-2154	4.3	13	
254	Experimental measurement of carbon dioxide solubility in 1-methylpyrrolidin-2-one (NMP) 1-butyl-3-methyl-1H-imidazol-3-ium tetrafluoroborate ([bmim][BF4]) mixtures using a new static-synthetic cell. <i>Fluid Phase Equilibria</i> , 2018 , 477, 62-77	2.5	3	
253	Valorisation of chicken feathers: Characterisation of thermal, mechanical and electrical properties. <i>Sustainable Chemistry and Pharmacy</i> , 2018 , 9, 27-34	3.9	25	
252	Crystal structure of (S)-tert-butyl-(1-hydroxypropan-2-yl)carbamate, C8H17NO3. <i>Zeitschrift Fur Kristallographie - New Crystal Structures</i> , 2018 , 233, 49-50	0.2		
251	Thermodynamics and activity coefficients at infinite dilution for organic solutes in the ionic liquid 1-butyl-1-methylpyrrolidinium dicyanamide. <i>Fluid Phase Equilibria</i> , 2018 , 473, 175-182	2.5	9	
250	Effect of temperature on density, sound velocity, and their derived properties for the binary systems glycerol with water or alcohols. <i>Journal of Chemical Thermodynamics</i> , 2017 , 109, 124-136	2.9	23	
249	Experimental determination of the critical loci for R-23 + (n-propane or n-hexane) and R-116 + n-propane binary mixtures. <i>Journal of Chemical Thermodynamics</i> , 2017 , 108, 84-96	2.9	7	
248	Microwave synthesis, biological evaluation and docking studies of 2-substituted methyl 1-(4-fluorophenyl)-1H-benzimidazole-5-carboxylates. <i>Medicinal Chemistry Research</i> , 2017 , 26, 484-498	2.2	5	
247	Design of a continuous gas-phase process for the production of hexafluoropropene oxide. <i>Chemical Engineering Research and Design</i> , 2017 , 119, 93-100	5.5		
246	Valorisation of chicken feathers: Characterisation of physical properties and morphological structure. <i>Journal of Cleaner Production</i> , 2017 , 149, 349-365	10.3	44	
245	Influence of gravitational potential on the thermodynamic stability of pure and mixed clathrate hydrates. <i>European Physical Journal B</i> , 2017 , 90, 1	1.2	1	
244	Experimental solubility data for prednisolone and hydrocortisone in various solvents between (293.2 and 328.2) K by employing combined DTA/TGA. <i>Journal of Molecular Liquids</i> , 2017 , 240, 303-312	6	4	
243	Vaporlliquid Equilibrium for Methyl Isobutyl Ketone (MIBK) + (1-Propanol or 2-Propanol) Binary Mixtures. <i>Journal of Chemical & Engineering Data</i> , 2017 , 62, 2014-2020	2.8	4	
242	Synthesis, in vitro antimicrobial, antioxidant, and antidiabetic activities of thiazolidinequinoxaline derivatives with amino acid side chains. <i>Medicinal Chemistry Research</i> , 2017 , 26, 2141-2151	2.2	10	

241	Carbon Dioxide to Energy: Killing Two Birds with One Stone 2017 , 93-103		0
240	Phosphonium Salts in Asymmetric Catalysis: A Journey in a Decade's Extensive Research Work. <i>Advanced Synthesis and Catalysis</i> , 2017 , 359, 3676-3706	5.6	58
239	Experimental Solubility Data for Binary Mixtures of Ethane and 2,2,4-Trimethylpentane at Pressures up to 6 MPa Using a New Variable-Volume Sapphire Cell. <i>Journal of Chemical & Engineering Data</i> , 2017 , 62, 3915-3920	2.8	4
238	Investigation of conventional and non-conventional hydrogen bonds: a comparison of fluorine-substituted and non-fluorine substituted compounds. <i>Monatshefte Fil Chemie</i> , 2017 , 148, 2061-	2068	4
237	Valorisation of chicken feathers: a review on recycling and recovery routedurrent status and future prospects. <i>Clean Technologies and Environmental Policy</i> , 2017 , 19, 2363-2378	4.3	55
236	Quality of Component- and Group-Interaction-Based Regression of Binary Vaporliquid Equilibrium Data. <i>Industrial & Data. Industrial & Data. Indust</i>	3.9	
235	Isothermal Vaporlliquid Equilibrium Data for the Hexafluoroethane (R116) +n-Butane System at Temperatures from 273 to 323 K. <i>Journal of Chemical & Data</i> , 2017, 62, 3483-3487	2.8	3
234	Phase stability conditions of carbon dioxide and methane clathrate hydrates in the presence of KBr, CaBr 2 , MgCl 2 , HCOONa, and HCOOK aqueous solutions: Experimental measurements and thermodynamic modelling. <i>Journal of Chemical Thermodynamics</i> , 2017 , 115, 307-317	2.9	11
233	Phase equilibrium and critical point data for ethylene and chlorodifluoromethane binary mixtures using a new Btatic-analyticDapparatus. <i>Fluid Phase Equilibria</i> , 2017 , 451, 106-113	2.5	6
232	Binary Vaporliquid Equilibrium Data for Perfluorooctane with Light Gases (Oxygen, Nitrogen, and Methane). <i>Journal of Chemical & Data</i> , 2017, 62, 4301-4309	2.8	1
231	Influence of temperature on thermophysical properties of tri(butyl)methylphosphonium methyl sulfate + N-methyl-2-pyrrolidone. <i>Journal of Molecular Liquids</i> , 2017 , 242, 375-381	6	6
230	Valorisation of chicken feathers: Characterisation of chemical properties. <i>Waste Management</i> , 2017 , 68, 626-635	8.6	60
229	Valorisation of chicken feathers: Application in paper production. <i>Journal of Cleaner Production</i> , 2017 , 164, 1324-1331	10.3	21
228	Influence of fluorination on barrier properties of polymers: Insights from Monte Carlo simulations of eicosanes + methane. <i>European Physical Journal E</i> , 2017 , 40, 12	1.5	1
227	The influence of temperature and composition on the density, viscosity and excess properties of aqueous mixtures of carboxylic-based ionic liquids. <i>Journal of Chemical Thermodynamics</i> , 2017 , 109, 71-50.	8 1 .9	20
226	Fluid-phase-equilibrium prediction of fluorocompound-containing binary systems with the predictive E-PPR78 model. <i>International Journal of Refrigeration</i> , 2017 , 73, 65-90	3.8	22
225	Activity coefficients at infinite dilution of hydrocarbons in glycols: Experimental data and thermodynamic modeling with the GCA-EoS. <i>Journal of Chemical Thermodynamics</i> , 2017 , 105, 226-237	2.9	5
224	Experimental solubility of diosgenin and estriol in various solvents between T=(293.2B28.2)K. Journal of Chemical Thermodynamics, 2017, 106, 199-207	2.9	6

223	Isothermal vapour-liquid equilibrium data for the binary systems 2-propanone + (2-butanol or propanoic acid). <i>Fluid Phase Equilibria</i> , 2017 , 433, 119-125	2.5	5	
222	Screening of environmental friendly ionic liquid as a solvent for the different types of separations problem: Insight from activity coefficients at infinite dilution measurement using (gas + liquid) chromatography technique. <i>Journal of Chemical Thermodynamics</i> , 2016 , 92, 35-42	2.9	13	
221	Isothermal (vapour + liquid) equilibrium data for binary systems of (n-hexane + CO2 or CHF3). Journal of Chemical Thermodynamics, 2016 , 94, 31-42	2.9	11	
220	Experimental measurement and thermodynamic modelling of hydrate phase equilibrium conditions for krypton + n -butyl ammonium bromide aqueous solution. <i>Journal of Supercritical Fluids</i> , 2016 , 107, 676-681	4.2	8	
219	A kinetic study of the selective production of difluoromethoxymethane from chlorodifluoromethane. <i>Journal of the Taiwan Institute of Chemical Engineers</i> , 2016 , 66, 70-79	5.3		
218	Phase Equilibria for Perfluoroethane + (n-Perfluorohexane or n-Perfluorooctane) Binary Systems: Measurement and Modeling. <i>Journal of Chemical & Data, 2016</i> , 61, 3363-3370	2.8	5	
217	Binary vapour-liquid equilibrium data for C7 and C9 straight-chain perfluorocarbons with ethylene. <i>Fluid Phase Equilibria</i> , 2016 , 429, 37-44	2.5	3	
216	Synthesis and structure elucidation using 2D NMR and thermal coefficient investigation on amino acid tethered quinoxalines. <i>Magnetic Resonance in Chemistry</i> , 2016 , 54, 921-929	2.1	O	
215	Vapor-liquid equilibria, density and sound velocity measurements of (water or methanol or ethanol + 1,3-propanediol) binary systems at different temperatures. <i>Thermochimica Acta</i> , 2016 , 642, 111-123	2.9	10	
214	Antioxidant properties, computational studies and corrosion inhibition potential of 3-hydroxy-1-(2-hydroxyphenyl)-5-(phenyl)-2,4-pentadien-1-one analogues. <i>Journal of Molecular Liquids</i> , 2016 , 223, 819-827	6	1	
213	Extraction of 2-phenylethanol (PEA) from aqueous phases using tetracyanoborate-based ionic liquids. <i>Journal of Molecular Liquids</i> , 2016 , 224, 1124-1130	6	10	
212	Development of a novel approach for modeling acid gas solubility in alkanolamine aqueous solution. <i>Journal of Natural Gas Science and Engineering</i> , 2016 , 34, 112-123	4.6	4	
211	Experimental Clathrate Hydrate Dissociation Data for Systems Comprising Refrigerant + CaCl2 Aqueous Solutions. <i>Journal of Chemical & Engineering Data</i> , 2016 , 61, 827-836	2.8	13	
210	A corresponding states-based method for the estimation of natural gas compressibility factors. <i>Journal of Molecular Liquids</i> , 2016 , 216, 25-34	6	17	
209	Thermodynamic stability conditions of clathrate hydrates for refrigerant (R134a or R410a or R507) with MgCl 2 aqueous solution. <i>Fluid Phase Equilibria</i> , 2016 , 413, 92-98	2.5	17	
208	On the prediction of critical temperatures of ionic liquids: Model development and evaluation. <i>Fluid Phase Equilibria</i> , 2016 , 411, 24-32	2.5	14	
207	Hydrate phase equilibria for CO2, CH4, or N2lletrabutylphosphonium bromide (TBPB) aqueous solution. <i>Fluid Phase Equilibria</i> , 2016 , 411, 88-92	2.5	15	
206	Experimental solubility for betulin and estrone in various solvents within the temperature range T = (293.2 to 328.2) K. <i>Journal of Chemical Thermodynamics</i> , 2016 , 98, 42-50	2.9	11	

205	Modeling of the vaporization enthalpies of petroleum fractions. Fluid Phase Equilibria, 2016, 412, 228-23	34 .5	5
204	Clathrate hydrate dissociation conditions for refrigerant Bucrose aqueous solution: Experimental measurement and thermodynamic modelling. <i>Fluid Phase Equilibria</i> , 2016 , 413, 99-109	2.5	7
203	Rapid method for the estimation of dew point pressures in gas condensate reservoirs. <i>Journal of the Taiwan Institute of Chemical Engineers</i> , 2016 , 60, 258-266	5.3	22
202	High-pressure phase equilibria data for mixtures involving ethene and perfluoro-n-octane from 293 to 353 K. <i>Fluid Phase Equilibria</i> , 2016 , 408, 33-37	2.5	7
201	Stirred cell ultrafiltration of lignin from black liquor generated from South African kraft mills. <i>South African Journal of Science</i> , 2016 , Volume 112,	1.3	3
200	New Insights into the Kinetics of the Gas-Phase Oxidation of Hexafluoropropene. <i>Progress in Reaction Kinetics and Mechanism</i> , 2016 , 41, 418-427	0.5	1
199	Assessing hydrate formation as a separation process for mixtures of close-boiling point compounds: A modelling study. <i>Journal of Natural Gas Science and Engineering</i> , 2016 , 35, 1405-1415	4.6	5
198	Phase equilibrium data for binary mixtures of carbon dioxide with {1,1,2,3,3,3-hexafluoro-1-propene or 2,2,3-trifluoro-3-(trifluoromethyl)oxirane} at temperatures between (233 and 273) K. <i>Fluid Phase Equilibria</i> , 2016 , 425, 114-119	2.5	2
197	Thermodynamic stability conditions for semi-clathrate hydrates of CO 2 , CH 4 , or N 2 with tetrabutyl ammonium nitrate (TBANO 3) aqueous solution. <i>Journal of Chemical Thermodynamics</i> , 2016 , 96, 52-56	2.9	4
196	Experimental measurements and thermodynamic modeling of the cloud point pressure for solubility of copolymers of vinyl acetate and dibutyl maleate in supercritical CO2. <i>Fluid Phase Equilibria</i> , 2016 , 425, 136-142	2.5	10
195	A group contribution model for prediction of the viscosity with temperature dependency for fluorine-containing ionic liquids. <i>Journal of Fluorine Chemistry</i> , 2016 , 186, 19-27	2.1	10
194	Isothermal vapour-liquid equilibrium data for the binary systems of (CHF3 or C2F6) and n-heptane. <i>Journal of Chemical Thermodynamics</i> , 2016 , 102, 237-247	2.9	5
193	Equilibrium data and GC-PC SAFT predictions for furanic extraction. <i>Fluid Phase Equilibria</i> , 2016 , 430, 57-66	2.5	7
192	GasIlquid mass transfer in a falling film microreactor: Effect of reactor orientation on liquid-side mass transfer coefficient. <i>Chemical Engineering Science</i> , 2016 , 155, 38-44	4.4	18
191	Factors influencing clathrate hydrate stability in equilibrium with liquid water: Insights from information-based statistical analysis. <i>Journal of Molecular Liquids</i> , 2016 , 222, 8-13	6	1
190	Clathrate hydrates modelled with classical density functional theory coupled with a simple lattice gas and van der Waals-Platteeuw theory. <i>Philosophical Magazine</i> , 2016 , 96, 2853-2867	1.6	6
189	Isothermal vaporliquid equilibrium data for the ethene+2,2,3-trifluoro-3-(trifluoromethyl)oxirane binary system between 258 and 308K at pressures up to 4.5MPa. <i>Fluid Phase Equilibria</i> , 2015 , 394, 88-92	2.5	1
188	Prediction of refractive indices of ionic liquids [A quantitative structure-property relationship based model. <i>Journal of the Taiwan Institute of Chemical Engineers</i> , 2015 , 52, 165-180	5.3	19

187	Liquid I liquid Equilibria for Mixtures of Hexadecane and Ethanol with Imidazolium-Based Ionic Liquids. <i>Journal of Solution Chemistry</i> , 2015 , 44, 593-605	1.8	9	
186	Synthesis and in vitro antiplatelet aggregation screening of novel fluorinated diethyl-2-(benzylthio)-2,3-dihydro-1H-imidazole-4,5-dicarboxylate derivatives. <i>Medicinal Chemistry Research</i> , 2015 , 24, 2075-2084	2.2	2	
185	Isothermal phase (vapour+liquid) equilibrium data for binary mixtures of propene (R1270) with either 1,1,2,3,3,3-hexafluoro-1-propene (R1216) or 2,2,3-trifluoro-3-(trifluoromethyl)oxirane in the temperature range of (279 to 318)K. <i>Journal of Chemical Thermodynamics</i> , 2015 , 90, 100-105	2.9	3	
184	Clathrate hydrate dissociation conditions of refrigerants R404A, R406A, R408A and R427A: Experimental measurements and thermodynamic modeling. <i>Journal of Chemical Thermodynamics</i> , 2015 , 90, 193-198	2.9	11	
183	CO2 Solubility in Hybrid Solvents Containing 1-Butyl-3-methylimidazolium Tetrafluoroborate and Mixtures of Alkanolamines. <i>Journal of Chemical & Engineering Data</i> , 2015 , 60, 2380-2391	2.8	13	
182	Isothermal Vapor□iquid Equilibrium Data for the 1,1,2,3,3,3-Hexafluoroprop-1-ene +1,1,2,2,3,3,4,4-Octafluorocyclobutane Binary System: Measurement and Modeling from (292 to 352) K and Pressures up to 2.6 MPa. <i>Journal of Chemical & Data</i> , 2015, 60, 966-969	2.8	1	
181	Stereo-selective synthesis, structural and antibacterial studies of novel glycosylated 2 ,3-amino acid analogues. <i>Medicinal Chemistry Research</i> , 2015 , 24, 3174-3193	2.2	3	
180	A universal segment approach for the prediction of the activity coefficient of complex pharmaceuticals in non-electrolyte solvents. <i>Fluid Phase Equilibria</i> , 2015 , 396, 98-110	2.5		
179	Experimental Measurements and Thermodynamic Modeling of Hydrate Dissociation Conditions for the Xenon + TBAB + Water System. <i>Journal of Chemical & Data</i> , 2015, 60, 1324-1330	2.8	8	
178	Copper-catalysed cross-coupling affected by the Smiles rearrangement: a new chapter on diversifying the synthesis of chiral fluorinated 1,4-benzoxazine derivatives. <i>RSC Advances</i> , 2015 , 5, 835	7 <i>6</i> 2835	80	
177	Isothermal Vaporlliquid Equilibrium (VLE) and Vaporlliquidlliquid Equilibrium (VLLE) Data for Two Binary Systems Containing Perfluorohexane with Carbon Monoxide or Hydrogen Sulfide at (293, 313, and 333) K. <i>Journal of Chemical & Engineering Data</i> , 2015 , 60, 2461-2468	2.8	3	
176	Phase equilibria study of binary systems comprising an (ionic liquid+hydrocarbon). <i>Journal of Chemical Thermodynamics</i> , 2015 , 83, 90-96	2.9	7	
175	Solubilities of carbon dioxide and oxygen in the ionic liquids methyl trioctyl ammonium bis(trifluoromethylsulfonyl)imide, 1-butyl-3-methyl imidazolium bis(trifluoromethylsulfonyl)imide, and 1-butyl-3-methyl imidazolium methyl sulfate. <i>Journal of Physical Chemistry B</i> , 2015 , 119, 1503-14	3.4	42	
174	Excess molar volumes of binary mixtures (an ionic liquid + water): A review. <i>Journal of Chemical Thermodynamics</i> , 2015 , 82, 34-46	2.9	82	
173	Experimental study and modeling of the kinetics of refrigerant hydrate formation. <i>Journal of Chemical Thermodynamics</i> , 2015 , 82, 47-52	2.9	29	
172	(Liquid + liquid) equilibria for mixtures of dodecane and ethanol with alkylsulfate-based ionic liquids. <i>Journal of Chemical Thermodynamics</i> , 2015 , 81, 95-100	2.9	12	
171	Experimental measurements and thermodynamic modeling of refrigerant hydrates dissociation conditions. <i>Journal of Chemical Thermodynamics</i> , 2015 , 80, 30-40	2.9	30	
170	Kinetic and thermodynamic behaviour of CF 4 clathrate hydrates. <i>Journal of Chemical Thermodynamics</i> , 2015 , 81, 52-59	2.9	26	

169	Phase equilibrium data for potentially hazardous binary mixtures involving dichlorosilane, trichlorosilane and silicon-tetrachloride. <i>Journal of Chemical Thermodynamics</i> , 2015 , 91, 420-426	2.9	4
168	Phase equilibrium data for mixtures involving 1,1,2,3,3,3-hexafluoro-1-propene with either propane or n-butane between 312 and 343K. <i>Fluid Phase Equilibria</i> , 2015 , 406, 156-162	2.5	6
167	Gas-phase equilibrium constants for the thermally initiated oxidation of hexafluoropropene with molecular oxygen. <i>Journal of Physical Organic Chemistry</i> , 2015 , 28, 460-471	2.1	4
166	Reliable method for the determination of surfactant retention in porous media during chemical flooding oil recovery. <i>Fuel</i> , 2015 , 158, 122-128	7.1	32
165	Influence of unlike dispersive interactions on methane adsorption in graphite: a grand canonical Monte Carlo simulation and classical density functional theory study. <i>European Physical Journal B</i> , 2015 , 88, 1	1.2	7
164	Application of the bio-inspired Krill Herd optimization technique to phase equilibrium calculations. <i>Computers and Chemical Engineering</i> , 2015 , 74, 75-88	4	18
163	Phase Equilibria of Clathrate Hydrates of Ethyne + Propene. <i>Journal of Chemical & Data</i> , 2015 , 60, 217-221	2.8	10
162	Isothermal Vaporlliquid Equilibrium Data for the Binary System 1,1,2,3,3,3-Hexafluoro-1-propene (R1216) + 2,2,3-Trifluoro-3-(trifluoromethyl)oxirane from (268.13 to 308.19) K. <i>Journal of Chemical & Chemical Pata</i> , 2015, 60, 568-573	2.8	5
161	Dry reforming of methane in a tiplip arc discharge reactor at very high pressure. <i>International Journal of Hydrogen Energy</i> , 2015 , 40, 3388-3401	6.7	11
160	A thermodynamic consistency test for experimental isobaric data of wax solubility in gaseous systems. <i>Fluid Phase Equilibria</i> , 2015 , 388, 182-187	2.5	
159	Effect of the alkyl side chain of the 1-alkylpiperidinium-based ionic liquids on desulfurization of fuels. <i>Journal of Chemical Thermodynamics</i> , 2014 , 72, 31-36	2.9	35
158	A chemical structure based model for the determination of speed of sound in ionic liquids. <i>Journal of Molecular Liquids</i> , 2014 , 196, 7-13	6	8
157	A group contribution method for determination of the standard molar chemical exergy of organic compounds. <i>Energy</i> , 2014 , 70, 288-297	7.9	18
156	Development of a group contribution method for the estimation of heat capacities of ionic liquids. Journal of Thermal Analysis and Calorimetry, 2014 , 115, 1863-1882	4.1	36
155	Toward a group contribution method for determination of speed of sound in saturated liquids. Journal of Molecular Liquids, 2014 , 194, 159-165	6	2
154	Determination of the speed of sound in ionic liquids using a least squares support vector machine group contribution method. <i>Fluid Phase Equilibria</i> , 2014 , 367, 188-193	2.5	12
153	Extrapolation/interpolation of infinite dilution, activity coefficient as well as liquid and solid solubility between solvents: Part 1. Alkane solvents. <i>Fluid Phase Equilibria</i> , 2014 , 361, 69-82	2.5	6
152	Measurement of activity coefficients at infinite dilution of organic solutes in the ionic liquid 1-ethyl-3-methylimidazolium 2-(2-methoxyethoxy) ethylsulfate at T = (308.15, 313.15, 323.15 and 333.15) K using gas + liquid chromatography. <i>Journal of Chemical Thermodynamics</i> , 2014 , 70, 245-252	2.9	31

151	A group contribution method for estimating the refractive indices of ionic liquids. <i>Journal of Molecular Liquids</i> , 2014 , 200, 410-415	6	15
150	Experimental Measurement and Thermodynamic Modeling of Hydrate Dissociation Conditions for the Argon + TBAB + Water System. <i>Journal of Chemical & Data</i> , 2014, 59, 3900-3906	2.8	13
149	A chemical structure based model for the estimation of refractive indices of organic compounds. <i>Fluid Phase Equilibria</i> , 2014 , 384, 1-13	2.5	6
148	Experimental Measurements and Thermodynamic Modeling of Clathrate Hydrate Dissociation Conditions for Refrigerants R116, R23, and Their Mixture R508B. <i>Journal of Chemical & Engineering Data</i> , 2014 , 59, 3907-3911	2.8	15
147	Assessment of Potential of Croton gratissimus Oil for Macroscale Production of Biodiesel Based on Thermophysical Properties. <i>Energy & Energy & 2014</i> , 28, 7576-7581	4.1	5
146	On the application of binary correction factors in lattice distortion calculations for methane clathrate hydrates. <i>Philosophical Magazine</i> , 2014 , 94, 974-990	1.6	6
145	Ternary LiquidIliquid Equilibrium Data for the Water + Acetonitrile + {Butan-1-ol or 2-Methylpropan-1-ol} Systems at (303.2, 323.2, 343.2) K and 1 atm. <i>Journal of Chemical & Engineering Data</i> , 2014 , 59, 3820-3824	2.8	10
144	Vapor l iquid Equilibrium (VLE) Data and Thermodynamic Modeling for Binary Systems Containing Perfluorobutane (R610) with Carbon Monoxide or Nitric Oxide at (293, 313, and 333) K. <i>Journal of Chemical & Data</i> , 2014 , 59, 346-354	2.8	4
143	A group contribution model for the prediction of the freezing point of organic compounds. <i>Fluid Phase Equilibria</i> , 2014 , 382, 21-30	2.5	5
142	Vapourliquid equilibrium of carboxylic acidlicohol binary systems: 2-Propanol+butyric acid, 2-butanol+butyric acid and 2-methyl-1-propanol+butyric acid. <i>Fluid Phase Equilibria</i> , 2014 , 380, 18-27	2.5	4
141	Vaporliquid Equilibrium Data for 1-Methyl-2-Pyrrolidone + (1-Butanol or 1-Hexene or Water) Binary Mixtures. <i>Journal of Chemical & Engineering Data</i> , 2014 , 59, 1643-1650	2.8	6
140	Isothermal Vaporliquid Equilibrium Data and Thermodynamic Modeling for Binary Systems of Perfluorobutane (R610) + (Methane or Hydrogen Sulfide) at (293, 313, and 333) K. <i>Journal of Chemical & Data</i> , 2014 , 59, 2865-2871	2.8	3
139	Isothermal Vaporlliquid Equilibrium Data for the 1,1,2,2-Tetrafluoroethene + 1,1,2,3,3,3-Hexafluoroprop-1-ene Binary System: Measurement and Modeling from (248 to 283) K. <i>Journal of Chemical & Data</i> , 2014, 59, 82-88	2.8	7
138	Effect of temperature on density, sound velocity, refractive index and their derived properties for the binary systems (heptanoic acid + propanoic or butanoic acids). <i>Journal of Chemical Thermodynamics</i> , 2014 , 78, 7-15	2.9	27
137	The 3-(4,5-dimethylthiazol-2-yl)-2,5-diphenyl tetrazolium bromide (MTT) assay is a rapid, cheap, screening test for the in vitro anti-tuberculous activity of chalcones. <i>Journal of Microbiological Methods</i> , 2014 , 104, 72-8	2.8	31
136	GAS-PHASE NON-CATALYTIC EPOXIDATION OF HEXAFLUOROPROPENE IN A TUBULAR REACTOR: OPTIMAL REACTION CONDITIONS. <i>Chemical Engineering Communications</i> , 2014 , 201, 1173-1197	2.2	2
135	Vaporliquid Equilibrium Data for Binary Systems of n-Dodecane + {Propan-1-ol, Butan-1-ol, 2-Methylpropan-1-ol} at 40 kPa. <i>Journal of Chemical & Data</i> , 2014, 59, 1710-1713	2.8	
134	Group Contribution Model for the Prediction of Refractive Indices of Organic Compounds. <i>Journal of Chemical & Engineering Data</i> , 2014 , 59, 1930-1943	2.8	12

133	Apparent Molar Volumes of 1-Alkyl (n = 2, 4, 6)-3-methylimidazolium Bromides in a Mixed-Solvent Medium of Acetonitrile + Water at Temperatures of (293.15, 303.15, and 313.15) K. <i>Journal of Chemical & Data</i> , 2014 , 59, 1086-1093	2.8	3
132	Phase Equilibria of Clathrate Hydrates of Ethyne + Propane. <i>Journal of Chemical & Data</i> , 2014 , 59, 2914-2919	2.8	19
131	Experimental Measurements and Thermodynamic Modeling of the Dissociation Conditions of Clathrate Hydrates for (Refrigerant + NaCl + Water) Systems. <i>Journal of Chemical & Data</i> , 2014, 59, 466-475	2.8	54
130	Influence of unlike dispersion interactions in modeling methane clathrate hydrates. <i>Fluid Phase Equilibria</i> , 2014 , 381, 108-115	2.5	15
129	Experimental study on the formation of higher fluorocarbons from CF4 by a tiplip electrical arc discharge at very high pressure. <i>Journal of Fluorine Chemistry</i> , 2014 , 166, 96-103	2.1	2
128	A group contribution method for determination of thermal conductivity of liquid chemicals at atmospheric pressure. <i>Journal of Molecular Liquids</i> , 2014 , 190, 223-230	6	6
127	Vapourllquid equilibrium of propionic acid + caproic acid, isobutyric acid + caproic acid, valeric acid + caproic acid and caproic acid + enanthoic acid binary mixtures. <i>Fluid Phase Equilibria</i> , 2014 , 375, 201-2	0 8 ·5	2
126	Solubility data and modeling for sugar alcohols in ionic liquids. <i>Journal of Chemical Thermodynamics</i> , 2014 , 77, 23-30	2.9	4
125	Solubility of ionic liquids in 2-phenylethanol (PEA) and water. Fluid Phase Equilibria, 2014, 376, 55-63	2.5	14
124	Review of carbon dioxide capture and storage with relevance to the South African power sector. <i>South African Journal of Science</i> , 2014 , 110, 1-12	1.3	7
123	Structure elucidation of a series of fluoro-2-styrylchromones and methoxy-2-styrylchromones using 1D and 2D NMR spectroscopy. <i>Magnetic Resonance in Chemistry</i> , 2014 , 52, 521-9	2.1	4
122	A New Method for the Analysis of Soluble and Insoluble Oxalate in Pulp and Paper Matrices. <i>Journal of Wood Chemistry and Technology</i> , 2014 , 34, 55-66	2	
121	Development of a LSSVM-GC model for estimating the electrical conductivity of ionic liquids. <i>Chemical Engineering Research and Design</i> , 2014 , 92, 66-79	5.5	24
120	Hydrocarbons synthesis from syngas by very high pressure plasma. <i>Chemical Engineering Journal</i> , 2014 , 241, 1-8	14.7	15
119	Phase equilibria of methane clathrate hydrates from Grand Canonical Monte Carlo simulations. <i>Fluid Phase Equilibria</i> , 2014 , 369, 47-54	2.5	35
118	Assessing the ability of force-fields to predict liquid I quid equilibria of ternary systems of light alcohols+water+dodecane by Monte Carlo simulation. Fluid Phase Equilibria, 2014, 368, 65-71	2.5	7
117	Effects of temperature and concentration on interactions in methanol + ethyl acetate and ethanol + methyl acetate or ethyl acetate systems: Insights from apparent molar volume and apparent molar isentropic compressibility study. <i>Thermochimica Acta</i> , 2014 , 577, 87-94	2.9	17
116	Experimental Measurements and Thermodynamic Modeling of the Dissociation Conditions of Clathrate Hydrates for (Refrigerant + NaCl + Water) Systems. <i>Journal of Chemical & Company Engineering</i>	2.8	4

115	A group contribution model for determining the sublimation enthalpy of organic compounds at the standard reference temperature of 298K. <i>Fluid Phase Equilibria</i> , 2013 , 354, 265-285	2.5	17
114	Determination of the normal boiling point of chemical compounds using a quantitative structureproperty relationship strategy: Application to a very large dataset. <i>Fluid Phase Equilibria</i> , 2013 , 354, 250-258	2.5	15
113	A Bon-linear Quantitative structure property relationship for the prediction of electrical conductivity of ionic liquids. <i>Chemical Engineering Science</i> , 2013 , 101, 478-485	4.4	17
112	Estimation of the Heat Capacity of Ionic Liquids: A Quantitative Structure P roperty Relationship Approach. <i>Industrial & Discrete Approach. Industrial & Disc</i>	3.9	19
111	Experimental vapour I quid equilibrium data and modeling for binary mixtures of 1-butene with 1,1,2,3,3,3-hexafluoro-1-propene, 2,2,3-trifluoro-3-(trifluoromethyl)oxirane, or difluoromethane. <i>Journal of Chemical Thermodynamics</i> , 2013 , 61, 18-26	2.9	9
110	Isothermal vaporliquid equilibrium data for the ethylene + 1,1,2,3,3,3-hexafluoro-1-propene binary system between 258 and 308 K at pressures up to 4.56 MPa. <i>Fluid Phase Equilibria</i> , 2013 , 353, 7-14	2.5	12
109	Gas hydrate phase equilibrium in porous media: An assessment test for experimental data. <i>Fluid Phase Equilibria</i> , 2013 , 360, 161-168	2.5	7
108	Vaporliquid Equilibrium Data for the Morpholine-4-carbaldehyde + n-Hexane or n-Heptane Binary Systems Using a Static-Synthetic Apparatus. <i>Journal of Chemical & Engineering Data</i> , 2013 , 58, 2552	2- <mark>2</mark> .8 2- 2 566	3
107	Isothermal vaporlīquid equilibrium of R170 + n-perfluorooctane at 308B38 K: Measurement, equation of state modelling, and molecular simulation. <i>Fluid Phase Equilibria</i> , 2013 , 344, 84-91	2.5	7
106	Activity coefficients at infinite dilution of organic solutes in diethylene glycol and triethylene glycol from gas l lquid chromatography. <i>Journal of Chemical Thermodynamics</i> , 2013 , 65, 120-130	2.9	7
105	Activity coefficients at infinite dilution of organic solutes in N-formylmorpholine and N-methylpyrrolidone from gas[lquid chromatography. <i>Journal of Chemical Thermodynamics</i> , 2013 , 61, 154-160	2.9	8
104	Development of a group contribution method for estimating the thermal decomposition temperature of ionic liquids. <i>Fluid Phase Equilibria</i> , 2013 , 355, 81-86	2.5	19
103	A novel dynamic recirculating apparatus for vapourliquid equilibrium measurements at moderate pressures and temperatures. <i>Fluid Phase Equilibria</i> , 2013 , 358, 121-130	2.5	11
102	Dissociation Data and Thermodynamic Modeling of Clathrate Hydrates of Ethene, Ethyne, and Propene. <i>Journal of Chemical & Engineering Data</i> , 2013 , 58, 3259-3264	2.8	28
101	A group contribution model for determining the vaporization enthalpy of organic compounds at the standard reference temperature of 298K. <i>Fluid Phase Equilibria</i> , 2013 , 360, 279-292	2.5	12
100	An assessment test for phase equilibrium data of water soluble and insoluble clathrate hydrate formers. <i>Fluid Phase Equilibria</i> , 2013 , 360, 68-76	2.5	17
99	Vapourllquid equilibrium (VLE) for the systems furan+n-hexane and furan+toluene. Measurements, data treatment and modeling using molecular models. <i>Fluid Phase Equilibria</i> , 2013 , 337, 234-245	2.5	20
98	A novel static analytical apparatus for phase equilibrium measurements. <i>Fluid Phase Equilibria</i> , 2013 , 338, 188-196	2.5	27

97	Isothermal (vapour+liquid) equilibria for binary mixtures of diisopropyl ether with (methanol, or ethanol, or 1-butanol): Experimental data, correlations, and predictions. <i>Journal of Chemical Thermodynamics</i> , 2013 , 58, 330-339	2.9	7
96	Activity coefficients at infinite dilution of organic solutes in the ionic liquid PEG-5 cocomonium methylsulfate at T = (313.15, 323.15, 333.15, and 343.15) K: Experimental results and COSMO-RS predictions. <i>Journal of Chemical Thermodynamics</i> , 2013 , 58, 322-329	2.9	17
95	Activity coefficients at infinite dilution of organic solutes in the ionic liquid trihexyltetradecylphosphonium bis (trifluoromethylsulfonyl) imide using gaslīquid chromatography at T=(313.15, 333.15, 353.15 and 373.15)K. <i>Journal of Chemical Thermodynamics</i> ,	2.9	14
94	2013, 65, 159-167 Experimental (vapour + liquid) equilibrium data and modelling for binary mixtures of decafluorobutane with propane and 1-butene. <i>Journal of Chemical Thermodynamics</i> , 2013, 67, 134-142	2.9	10
93	Development of a general model for determination of thermal conductivity of liquid chemical compounds at atmospheric pressure. <i>AICHE Journal</i> , 2013 , 59, 1702-1708	3.6	21
92	Monte Carlo simulations of water solubility and structures in poly(difluoromethylene). <i>Molecular Simulation</i> , 2013 , 39, 367-384	2	1
91	Assessment of Carbon Dioxide Dissociation as a New Route for Syngas Production: A Comparative Review and Potential of Plasma-Based Technologies. <i>Energy & Energy & E</i>	4.1	88
90	Isothermal Vaporliquid Equilibrium Data for the Butan-2-one + Methanol or Ethanol Systems Using a Static-Analytic Microcell. <i>Journal of Chemical & Engineering Data</i> , 2013 , 58, 1280-1287	2.8	9
89	Prediction of the solubility of selected pharmaceuticals in water and alcohols with a group contribution method. <i>Journal of Chemical Thermodynamics</i> , 2013 , 62, 118-129	2.9	6
88	Development of a quantitative structurellquid thermal conductivity relationship for pure chemical compounds. <i>Fluid Phase Equilibria</i> , 2013 , 355, 52-80	2.5	15
87	Density, speed of sound, and refractive index measurements for the binary systems (butanoic acid+propanoic acid, or 2-methyl-propanoic acid) at T=(293.15 to 313.15)K. <i>Journal of Chemical Thermodynamics</i> , 2013 , 57, 203-211	2.9	50
86	SolidIlquid equilibria measurements for binary systems comprising (butyric acid+propionic or pentanoic acid) and (heptanoic acid+propionic or butyric or pentanoic or hexanoic acid). <i>Journal of Chemical Thermodynamics</i> , 2013 , 57, 485-492	2.9	14
85	Isothermal Vaporlliquid Equilibrium Data and Modeling for the Ethane (R170) + Perfluoropropane (R218) System at Temperatures from (264 to 308) K. <i>Journal of Chemical & Chemica</i>	2.8	7
84	Phase Equilibria of Clathrate Hydrates of Ethane + Ethene. <i>Journal of Chemical & Data</i> , 2013 , 58, 896-901	2.8	25
83	Synthesis and Evaluation of Novel Fluorinated 2-Styrylchromones as Antibacterial Agents. <i>Journal of Chemistry</i> , 2013 , 2013, 1-13	2.3	3
82	3-Bromo-chroman-4-one. Acta Crystallographica Section E: Structure Reports Online, 2013, 69, o473		
81	3-(3-Nitro-benz-yl)-4H-chromen-4-one. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2013 , 69, o364		1
80	3-(4-Nitro-benz-yl)-4H-chromen-4-one. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2013 , 69, o513		

79	Phase equilibrium measurements for semi-clathrate hydrates of the (CO2+N2+tetra-n-butylammonium bromide) aqueous solution system. <i>Journal of Chemical Thermodynamics</i> , 2012 , 46, 57-61	2.9	91
78	Application of gas hydrate formation in separation processes: A review of experimental studies. <i>Journal of Chemical Thermodynamics</i> , 2012 , 46, 62-71	2.9	403
77	Activity coefficients at infinite dilution of organic solutes in the ionic liquid trihexyltetradecylphosphonium hexafluorophosphate using gas[Iquid chromatography at T=(313.15, 333.15, 353.15, and 363.15)K. <i>Journal of Chemical Thermodynamics</i> , 2012 , 49, 46-53	2.9	15
76	Speeds of Sound in Liquid and Supercritical Hexafluoropropylene (HFP) and Hexafluoropropylene Oxide (HFPO) at Pressures up to 400 MPa. <i>Journal of Chemical & Engineering Data</i> , 2012 , 57, 2568-	-2 37 5	1
75	Quantitative structureproperty relationship for thermal decomposition temperature of ionic liquids. <i>Chemical Engineering Science</i> , 2012 , 84, 557-563	4.4	31
74	Vaporliquid equilibria of ethylene (C2H4) + decafluorobutane (C4F10) at 268\(\bar{\textit{19}}\)98 K from experiment, molecular simulation and the Peng\(\bar{\textbf{R}}\)obinson equation of state. Fluid Phase Equilibria, 2012, 336, 104-112	2.5	10
73	Absorption Data and Modeling of Carbon Dioxide in Aqueous Blends of Bis(2-hydroxyethyl)methylamine (MDEA) and 2,2-Iminodiethanol (DEA): 25 % MDEA + 25 % DEA and 30 % MDEA + 20 % DEA. <i>Journal of Chemical & Description Data</i> , 2012 , 57, 1607-1620	2.8	10
72	Isothermal VaporIliquid Equilibrium Data for the 1,1,2,2-Tetrafluoroethene + 1,1,2,2,3,3,4,4-Octafluorocyclobutane Binary System: Measurement and Modeling from (248 to 283) K. <i>Journal of Chemical & Data</i> , Engineering Data, 2012 , 57, 1978-1983	2.8	7
71	Vaporlliquid Equilibrium Data for Binary Systems of 1H-Pyrrole with Butan-1-ol, Propan-1-ol, or Pentan-1-ol. <i>Journal of Chemical & Engineering Data</i> , 2012 , 57, 2520-2527	2.8	8
70	Vapor l liquid Equilibrium Measurements and Modeling for the Ethane (R-170) + 1,1,2,3,3,3-Hexafluoro-1-propene (R-1216) Binary System. <i>Journal of Chemical & Data</i> , 2012, 57, 2947-2955	2.8	15
69	Kinetics of the Gas-Phase Noncatalytic Oxidation of Hexafluoropropene. <i>Industrial & amp; Engineering Chemistry Research</i> , 2012 , 51, 13961-13972	3.9	2
68	Development of a group contribution method for determination of viscosity of ionic liquids at atmospheric pressure. <i>Chemical Engineering Science</i> , 2012 , 80, 326-333	4.4	53
67	Synthesis and Characterization of Imidazolium Salts Bearing Fluorinated Anions. <i>Zeitschrift Fur Anorganische Und Allgemeine Chemie</i> , 2012 , 638, n/a-n/a	1.3	2
66	Vaporlliquid Equilibrium Data for Binary Systems of 1-Methyl-4-(1-methylethenyl)-cyclohexene + {Ethanol, Propan-1-ol, Propan-2-ol, Butan-1-ol, Pentan-1-ol, or Hexan-1-ol} at 40 kPa. <i>Journal of Chemical & Engineering Data</i> , 2012 , 57, 2053-2058	2.8	10
65	2,4-Dibromo-1,3-dimeth-oxy-5-methyl-benzene. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2012 , 68, o2062		
64	Phase-dependent energy cross-parameters in a monatomic binary fluid system. <i>Molecular Simulation</i> , 2012 , 38, 838-849	2	6
63	2-Acetyl-phenyl (2E)-3-(4-fluoro-phen-yl)acrylate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2012 , 68, o3049		1
62	3-(3-Meth-oxy-benzyl-idene)chroman-4-one. <i>Acta Crystallographica Section E: Structure Reports</i> Online, 2012 , 68, o1006		6

61	(E)-3-(4-Cyclo-hexyl-3-fluoro-benzyl-idene)chroman-4-one. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2012 , 68, o1972		5
60	3-(3,4-Dichloro-benzyl-idene)chroman-4-one. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2012 , 68, o3062		3
59	Phase equilibria study of {N-hexylisoquinolinium bis{(trifluoromethyl)sulfonyl}imide + aromatic hydrocarbons or an alcohol} binary systems. <i>Journal of Physical Chemistry B</i> , 2011 , 115, 4003-10	3.4	21
58	Phase Equilibria of Methane and Carbon Dioxide Clathrate Hydrates in the Presence of Aqueous Solutions of Tributylmethylphosphonium Methylsulfate Ionic Liquid. <i>Journal of Chemical & Engineering Data</i> , 2011 , 56, 3620-3629	2.8	115
57	Liquid Diquid Equilibria of Methanol, Ethanol, and Propan-2-ol with Water and Dodecane. <i>Journal of Chemical & Chemical &</i>	2.8	18
56	Experimental Study of Hydrocarbons Synthesis from Syngas by a Tipllip Electrical Discharge at Very High Pressure. <i>Plasma Chemistry and Plasma Processing</i> , 2011 , 31, 663-679	3.6	13
55	Measurements of activity coefficients at infinite dilution of organic compounds and water in isoquinolinium-based ionic liquid [C8iQuin][NTf2] using GLC. <i>Journal of Chemical Thermodynamics</i> , 2011 , 43, 499-504	2.9	69
54	Activity coefficients at infinite dilution of organic solutes in the ionic liquid trihexyl(tetradecyl)phosphonium tetrafluoroborate using gasIlquid chromatography at T = (313.15, 333.15, 353.15, and 373.15) K. <i>Journal of Chemical Thermodynamics</i> , 2011 , 43, 670-676	2.9	16
53	Activity coefficients at infinite dilution of organic solutes in the ionic liquid 1-butyl-3-methylimidazolium hexafluoroantimonate using gas[]quid chromatography at T= (313.15, 323.15, and 333.15) K. <i>Journal of Chemical Thermodynamics</i> , 2011 , 43, 829-833	2.9	17
52	Determination of activity coefficients at infinite dilution of water and organic solutes (polar and non-polar) in the Ammoeng 100 ionic liquid at T=(308.15, 313.5, 323.15, and 333.15)K. <i>Journal of Chemical Thermodynamics</i> , 2011 , 43, 1178-1184	2.9	12
51	Isothermal Vaporliquid Equilibrium Data for the Perfluorobutane (R610) + Ethane System at Temperatures from (263 to 353) K. <i>Journal of Chemical & Engineering Data</i> , 2011 , 56, 1918-1924	2.8	29
50	Experimental Measurement of Vapor Pressures and Densities at Saturation of Pure Hexafluoropropylene Oxide: Modeling Using a Crossover Equation of State. <i>Industrial & Engineering Chemistry Research</i> , 2011 , 50, 4761-4768	3.9	21
49	Isothermal vaporllquid equilibrium data for the carbon dioxide (R744)+decafluorobutane (R610) system at temperatures from 263 to 353K. <i>Fluid Phase Equilibria</i> , 2011 , 304, 44-51	2.5	25
48	Liquid II quid phase equilibrium of (piperidinium-based ionic liquid+an alcohol) binary systems and modelling with NRHB and PCP-SAFT. <i>Fluid Phase Equilibria</i> , 2011 , 305, 43-52	2.5	64
47	Activity coefficients at infinite dilution of organic solutes in the ionic liquid, methyl(trioctyl)ammonium thiosalicylate, [N1888][TS] by gasIlquid chromatography at T=(303.15, 313.15, and 323.15)K. <i>Journal of Chemical Thermodynamics</i> , 2011 , 43, 754-758	2.9	14
46	6-Hy-droxy-2H-1,3-benzodioxole-5-carbaldehyde. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011 , 67, o2681		
45	[1,4-Phenyl-enebis(methyl-ene)]bis-(tri-phenyl-phospho-nium) bis-(tetra-fluoro-borate). <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011 , 67, o3391		1
44	Gibbs ensemble Monte Carlo simulations of binary vapourliquidliquid equilibrium: application to n-hexaneWater and ethaneBthanol systems. <i>Molecular Simulation</i> , 2010 , 36, 758-762	2	5

43	Phase equilibria and modeling of pyridinium-based ionic liquid solutions. <i>Journal of Physical Chemistry B</i> , 2010 , 114, 15011-7	3.4	26
42	Pure Component and Binary Vaporliquid Equilibrium + Modeling for Hexafluoropropylene and Hexafluoropropylene Oxide with Toluene and Hexafluoroethane. <i>Journal of Chemical & Engineering Data</i> , 2010 , 55, 411-418	2.8	15
41	High-Pressure Vaporliquid Equilibrium Data for (Carbon Dioxide + Cyclopentanol) and (Propane + Cyclopentanol). <i>Journal of Chemical & Data</i> , 2010, 55, 196-200	2.8	2
40	Experimental Measurement of Vapor Pressures and Densities of Pure Hexafluoropropylene. Journal of Chemical & Densities of Pure Hexafluoropropylene.	2.8	41
39	Isothermal Vaporliquid Equilibrium Data for the Hexafluoropropylene (R1216) + Propylene System at Temperatures from (263.17 to 353.14) K. <i>Journal of Chemical & Engineering Data</i> , 2010 , 55, 1636-1639	2.8	21
38	Separation of an Alcohol and a Tetrahydrofuran, Methyl tert-Butyl Ether, or Ethyl tert-Butyl Ether by Solvent Extraction with a Hyperbranched Polymer at T = 298.15 K. <i>Journal of Chemical &</i> Engineering Data, 2010 , 55, 2879-2885	2.8	17
37	Activity coefficients at infinite dilution of organic solutes in the ionic liquid 1-ethyl-3-methylimidazolium trifluoromethanesulfonate using gasIlquid chromatography at T=(313.15, 323.15, and 333.15)K. <i>Journal of Chemical Thermodynamics</i> , 2010 , 42, 78-83	2.9	48
36	Activity coefficients at infinite dilution for solutes in the trioctylmethylammonium bis(trifluoromethylsulfonyl)imide ionic liquid using gas Iquid chromatography. <i>Journal of Chemical Thermodynamics</i> , 2010 , 42, 256-261	2.9	36
35	Activity coefficients at infinite dilution of organic solutes in the ionic liquid 1-octyl-3-methylimidazolium hexafluorophosphate using gasIlquid chromatography at T= (313.15, 323.15, and 333.15) K. <i>Journal of Chemical Thermodynamics</i> , 2010 , 42, 646-650	2.9	21
34	Phase equilibria study of {N-butylquinolinium bis{(trifluoromethyl)sulfonyl}imide+aromatic hydrocarbons, or an alcohol} binary systems. <i>Journal of Chemical Thermodynamics</i> , 2010 , 42, 1180-1186	2.9	21
33	(Solid+liquid) and (liquid+liquid) phase equilibria study and correlation of the binary systems {N-butyl-3-methylpyridinium tosylate+water, or+an alcohol, or+a hydrocarbon}. <i>Fluid Phase Equilibria</i> , 2010 , 294, 89-97	2.5	20
32	Monte Carlo simulations of vaporliquidliquid equilibrium of some ternary petrochemical mixtures. <i>Fluid Phase Equilibria</i> , 2010 , 299, 24-31	2.5	4
31	Determination of critical properties of pure and multi-component mixtures using a BynamicByntheticDapparatus. <i>Journal of Supercritical Fluids</i> , 2010 , 55, 545-553	4.2	41
30	Estimation of pure component properties. Part 4: Estimation of the saturated liquid viscosity of non-electrolyte organic compounds via group contributions and group interactions. <i>Fluid Phase Equilibria</i> , 2009 , 281, 97-119	2.5	27
29	Vaporliquid equilibrium in the n-butane+methanol system, measurement and modeling from 323.2 to 443.2K. <i>Fluid Phase Equilibria</i> , 2009 , 277, 152-161	2.5	24
28	Vapor I lquid equilibrium measurements and modeling of the n-butane + ethanol system from 323 to 423 K. <i>Fluid Phase Equilibria</i> , 2009 , 286, 79-87	2.5	15
27	Activity coefficients at infinite dilution measurements for organic solutes in the ionic liquid N-butyl-4-methylpyridinium tosylate using GLC at T = (328.15, 333.15, 338.15, and 343.15) K. <i>Fluid Phase Equilibria</i> , 2009 , 276, 31-36	2.5	39
26	Isothermal Vaporliquid Equilibrium Data for the Hexafluoroethane (R116) + Propane System at Temperatures from (263 to 323) K. <i>Journal of Chemical & Data,</i> 2009, 54, 1292-1296	2.8	28

25	Vaporliquid Equilibrium for Binary Systems of 2,3-Pentanedione with Diacetyl and Acetone. <i>Journal of Chemical & Diacetyl and Acetone. Journal of Chemical & Diacetyl and Acetone.</i>	2.8	15
24	Determination of Activity Coefficients at Infinite Dilution of Solutes in the Ionic Liquid, Trihexyltetradecylphosphonium Bis(trifluoromethylsulfonyl) Imide, Using Gas[liquid Chromatography at T = (303.15, 308.15, 313.15, and 318.15) K. <i>Journal of Chemical & Description</i>	2.8	39
23	Estimation of pure component properties: Part 3. Estimation of the vapor pressure of non-electrolyte organic compounds via group contributions and group interactions. <i>Fluid Phase Equilibria</i> , 2008 , 269, 117-133	2.5	200
22	A new high-pressure vapour[Iquid equilibrium apparatus. Fluid Phase Equilibria, 2008, 269, 104-112	2.5	29
21	Activity coefficients at infinite dilution measurements for organic solutes in the ionic liquid trihexyltetradecylphosphonium-bis-(2,4,4-trimethylpentyl)-phosphinate using g.l.c. at T=(303.15, 308.15, 313.15, and 318.15)K. <i>Journal of Chemical Thermodynamics</i> , 2008 , 40, 1243-1247	2.9	40
20	Estimation of the vapour pressure of non-electrolyte organic compounds via group contributions and group interactions. <i>Journal of Molecular Liquids</i> , 2008 , 143, 52-63	6	85
19	High pressure vapor []quid equilibrium measurements of carbon dioxide with naphthalene and benzoic acid. <i>Fluid Phase Equilibria</i> , 2007 , 260, 60-64	2.5	11
18	Estimation of pure component properties. Fluid Phase Equilibria, 2007, 252, 1-27	2.5	111
17	Liquid Diquid Equilibria for Ternary Mixtures (an Ionic Liquid + Benzene + Heptane or Hexadecane) at T = 298.2 K and Atmospheric Pressure. <i>Journal of Chemical & Digital Benzening Data</i> , 2006 , 51, 988-99	1 ^{2.8}	100
16	Vaporliquid Equilibrium for Binary Systems of Diacetyl with Methanol and Acetone. <i>Journal of Chemical & Data</i> , 2006 , 51, 2083-2087	2.8	9
15	Monte carlo simulation of carboxylic acid phase equilibria. <i>Journal of Physical Chemistry B</i> , 2006 , 110, 21938-43	3.4	21
14	Membrane distillation of concentrated brines R ole of water activities in the evaluation of driving force. <i>Journal of Membrane Science</i> , 2006 , 280, 937-947	9.6	68
13	New developments in differential ebulliometry: Experimental and theoretical. <i>Journal of Molecular Liquids</i> , 2006 , 125, 45-57	6	7
12	Vaporliquid Equilibrium Measurements of MTBE and TAME with Toluene. <i>Journal of Chemical & Engineering Data</i> , 2005 , 50, 56-59	2.8	6
11	Simulation of 1-alkene and n-alkane binary vapour Ilquid equilibrium using different united-atom transferable force fields. <i>Fluid Phase Equilibria</i> , 2005 , 232, 136-148	2.5	4
10	Vapourllquid equilibrium of carboxylic acid systems: Propionic acid+valeric acid and isobutyric acid+valeric acid. <i>Fluid Phase Equilibria</i> , 2005 , 237, 89-99	2.5	24
9	Estimation of pure component properties: Part 1. Estimation of the normal boiling point of non-electrolyte organic compounds via group contributions and group interactions. <i>Fluid Phase Equilibria</i> , 2004 , 226, 45-63	2.5	192
8	Subatmospheric Vapor Pressure Curves for Propionic Acid, Butyric Acid, Isobutyric Acid, Valeric Acid, Isovaleric Acid, Hexanoic Acid, and Heptanoic Acid. <i>Journal of Chemical & Data</i> 2004 49, 1189-1192	2.8	25

LIST OF PUBLICATIONS

7	Activity coefficients of hydrocarbon solutes at infinite dilution in the ionic liquid, 1-methyl-3-octyl-imidazolium chloride from gasllquid chromatography. <i>Journal of Chemical Thermodynamics</i> , 2003 , 35, 1335-1341	2.9	94
6	Activity Coefficients at Infinite Dilution of Organic Solutes in 1-Hexyl-3-methylimidazolium Hexafluorophosphate from Gasliquid Chromatography. <i>Journal of Chemical & Data</i> , 2003, 48, 708-711	2.8	123
5	Ternary Liquidliquid Equilibria for Mixtures of 1-Methyl-3-octylimidazolium Chloride + an Alkanol + an Alkane at 298.2 K and 1 bar. <i>Journal of Chemical & Engineering Data</i> , 2003 , 48, 904-907	2.8	89
4	Isobaric Vapor liquid Equilibria for the Systems Propionic Acid + Butyric Acid, Isobutyric Acid + Butyric Acid, Butyric Acid + Isovaleric Acid, and Butyric Acid + Hexanoic Acid at 14 kPa. <i>Journal of Chemical & Data</i> , 2002, 47, 603-607	2.8	36
3	Monoethanolamine as an Extractive Solvent for the n-Hexane + Benzene, Cyclohexane + Ethanol, and Acetone + Methanol Binary Systems. <i>Journal of Chemical & Engineering Data</i> , 2002 , 47, 781-787	, 2.8	6
2	Rigorous characterization of static and dynamic apparatus for measuring limiting activity coefficients. <i>Fluid Phase Equilibria</i> , 2001 , 187-188, 473-487	2.5	8
1	Ternary Liquid Liquid Equilibra for Pseudoternary Mixtures Containing an n-Alkane + an Aromatic Hydrocarbon + {N-Methyl-2-pyrrolidinone + a Solvent} at 298.2 K and 1 atm. <i>Journal of Chemical & Engineering Data</i> , 2001 , 46, 1375-1380	2.8	11