

# Josefa Fernandez

## 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.

229  
papers

6,384  
citations

41  
h-index

65  
g-index

243  
ext. papers

6,812  
ext. citations

3.2  
avg. IF

5.84  
L-index

#	Paper	IF	Citations
229	Improvement of the lubrication performance of an ester base oil with coated ferrite nanoadditives for different material pairs. <i>Journal of Molecular Liquids</i> , <b>2022</b> , 350, 118550	6	1
228	ZnO nanoparticles coated with oleic acid as additives for a polyalphaolefin lubricant. <i>Journal of Molecular Liquids</i> , <b>2022</b> , 348, 118401	6	2
227	Comparison between thermophysical and tribological properties of two engine lubricant additives: electrochemically exfoliated graphene and molybdenum disulfide nanoplatelets. <i>Nanotechnology</i> , <b>2021</b> , 32, 025701	3.4	9
226	Tribological properties of hexagonal boron nitride nanoparticles or graphene nanoplatelets blended with an ionic liquid as additives of an ester base oil. <i>Lubrication Science</i> , <b>2021</b> , 33, 269-278	1.3	0
225	Tribological synergies among chemical-modified graphene oxide nanomaterials and a phosphonium ionic liquid as additives of a biolubricant. <i>Journal of Molecular Liquids</i> , <b>2021</b> , 116885	6	3
224	Effect of hydrophobic phenomena over the volumetric behavior of aqueous ionic liquid solutions. <i>Journal of Molecular Liquids</i> , <b>2021</b> , 333, 115962	6	1
223	Tribological properties of graphene nanoplatelets or boron nitride nanoparticles as additives of a polyalphaolefin base oil. <i>Journal of Molecular Liquids</i> , <b>2021</b> , 333, 115911	6	7
222	Hybrid combinations of graphene nanoplatelets and phosphonium ionic liquids as lubricant additives for a polyalphaolefin. <i>Journal of Molecular Liquids</i> , <b>2021</b> , 336, 116266	6	7
221	Lubricant properties of trimethylolpropane trioleate biodegradable oil: High pressure density and viscosity, film thickness, Stribeck curves and influence of nanoadditives. <i>Journal of Molecular Liquids</i> , <b>2021</b> , 335, 116410	6	2
220	Heat capacity, density, surface tension, and contact angle for polyalphaolefins and ester lubricants. <i>Thermochimica Acta</i> , <b>2021</b> , 703, 178994	2.9	5
219	Double hybrid lubricant additives consisting of a phosphonium ionic liquid and graphene nanoplatelets/hexagonal boron nitride nanoparticles. <i>Tribology International</i> , <b>2021</b> , 163, 107189	4.9	5
218	Modeling of the Production of Lipid Microparticles Using PGSS Technique. <i>Molecules</i> , <b>2020</b> , 25,	4.8	4
217	Thermal conductivity of ionic liquids under pressure. <i>Fluid Phase Equilibria</i> , <b>2020</b> , 515, 112573	2.5	9
216	Synergistic effects of hexagonal boron nitride nanoparticles and phosphonium ionic liquids as hybrid lubricant additives. <i>Journal of Molecular Liquids</i> , <b>2020</b> , 311, 113343	6	15
215	Synergy between boron nitride or graphene nanoplatelets and tri(butyl)ethylphosphonium diethylphosphate ionic liquid as lubricant additives of triisotridecyltrimellitate oil. <i>Journal of Molecular Liquids</i> , <b>2020</b> , 301, 112442	6	14
214	Tribological performance of silicon nitride and carbon black Ionanofluids based on 1-ethyl-3-methylimidazolium methanesulfonate. <i>Journal of Molecular Liquids</i> , <b>2020</b> , 319, 114335	6	8
213	Tribological Behavior of Nanolubricants Based on Coated Magnetic Nanoparticles and Trimethylolpropane Trioleate Base Oil. <i>Nanomaterials</i> , <b>2020</b> , 10,	5.4	16

212	Tribological and Thermophysical Properties of Environmentally-Friendly Lubricants Based on Trimethylolpropane Trioleate with Hexagonal Boron Nitride Nanoparticles as an Additive. <i>Coatings</i> , <b>2019</b> , 9, 509	2.9	11
211	High pressure viscosity behaviour of tris(2-ethylhexyl) trimellitate up to 150 MPa. <i>Journal of Chemical Thermodynamics</i> , <b>2019</b> , 138, 159-166	2.9	7
210	Experimental Convection Heat Transfer Analysis of a Nano-Enhanced Industrial Coolant. <i>Nanomaterials</i> , <b>2019</b> , 9,	5.4	13
209	Functionalized graphene nanoplatelet nanofluids based on a commercial industrial antifreeze for the thermal performance enhancement of wind turbines. <i>Applied Thermal Engineering</i> , <b>2019</b> , 152, 113-125	5.8	25
208	Thermophysical properties of polyalphaolefin oil modified with nanoadditives. <i>Journal of Chemical Thermodynamics</i> , <b>2019</b> , 131, 192-205	2.9	16
207	International Standard for viscosity at temperatures up to 473 K and pressures below 200 MPa (IUPAC Technical Report). <i>Pure and Applied Chemistry</i> , <b>2019</b> , 91, 161-172	2.1	5
206	High Pressure Characterization of the Viscous and Volumetric Behavior of Three Transmission Oils. <i>Industrial &amp; Engineering Chemistry Research</i> , <b>2019</b> , 58, 1732-1742	3.9	2
205	Tribological properties of dispersions based on reduced graphene oxide sheets and trimethylolpropane trioleate or PAO 40 oils. <i>Journal of Molecular Liquids</i> , <b>2019</b> , 274, 568-576	6	25
204	Excess thermodynamic functions in aqueous systems containing soluble fullerene derivatives. <i>Journal of Molecular Liquids</i> , <b>2018</b> , 256, 305-311	6	14
203	Isobaric heat capacity of nanostructured liquids with potential use as lubricants. <i>Journal of Chemical Thermodynamics</i> , <b>2018</b> , 123, 107-116	2.9	9
202	Effect of ZrO <sub>2</sub> nanoparticles on thermophysical and rheological properties of three synthetic oils. <i>Journal of Molecular Liquids</i> , <b>2018</b> , 262, 126-138	6	21
201	Viscosity-pressure dependence for nanostructured ionic liquids. Experimental values for butyltrimethylammonium and 1-butyl-3-methylpyridinium bis(trifluoromethylsulfonyl)imide. <i>Journal of Chemical Thermodynamics</i> , <b>2018</b> , 121, 27-38	2.9	9
200	Determination of derived volumetric properties and heat capacities at high pressures using two density scaling based equations of state. Application to dipentaerythritol hexa(3,5,5-trimethylhexanoate). <i>Physical Chemistry Chemical Physics</i> , <b>2018</b> , 20, 3531-3542	3.6	5
199	Speed of sound in ionic liquids with a common ion as a function of pressure and temperature. <i>Journal of Chemical Thermodynamics</i> , <b>2018</b> , 116, 235-240	2.9	18
198	Thermal stability of aprotic ionic liquids as potential lubricants. Comparison with synthetic oil bases. <i>Journal of Chemical Thermodynamics</i> , <b>2018</b> , 116, 185-196	2.9	29
197	Pressure and temperature dependence of light fullerenes solubility in n-heptane. <i>Journal of Molecular Liquids</i> , <b>2018</b> , 268, 569-577	6	1
196	Thermophysical and tribological properties of dispersions based on graphene and a trimethylolpropane trioleate oil. <i>Journal of Molecular Liquids</i> , <b>2018</b> , 268, 854-866	6	22
195	High pressure densities of two nanostructured liquids based on the bis(trifluoromethylsulfonyl)imide anion from (278 to 398) K and up to 120 MPa. <i>Journal of Chemical Thermodynamics</i> , <b>2018</b> , 118, 67-76	2.9	6

194	Thermal stability of some imidazolium [NTF 2 ] ionic liquids: Isothermal and dynamic kinetic study through thermogravimetric procedures. <i>Journal of Chemical Thermodynamics</i> , <b>2017</b> , 112, 105-113	2.9	27
193	In Pursuit of a High-Temperature, High-Pressure, High-Viscosity Standard: The Case of Tris(2-ethylhexyl) Trimellitate. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2017</b> , 62, 2884-2895	2.8	18
192	Pressure dependence of the solubility of light fullerenes in n -nonane. <i>Journal of Chemical Thermodynamics</i> , <b>2017</b> , 112, 259-266	2.9	3
191	High Pressure Rheological Behavior of 1-Ethyl-3-methylimidazolium n-Hexylsulfate and Trihexyl(tetradecyl)phosphonium Tris(pentafluoroethyl)trifluorophosphate. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2017</b> , 62, 2927-2936	2.8	4
190	PEG 400-Based Phase Change Materials Nano-Enhanced with Functionalized Graphene Nanoplatelets. <i>Nanomaterials</i> , <b>2017</b> , 8,	5.4	36
189	Volumetric Behavior of Some Motor and Gear-Boxes Oils at High Pressure: Compressibility Estimation at EHL Conditions. <i>Industrial &amp; Engineering Chemistry Research</i> , <b>2017</b> , 56, 10877-10885	3.9	3
188	Tribological Behavior of Two 1-Ethyl-3-methylimidazolium Alkyl Sulfates as Neat Lubricants for a Steel/Steel Contact. <i>Tribology Transactions</i> , <b>2017</b> , 60, 729-738	1.8	3
187	Volumetric behaviour of six ionic liquids from T = (278 to 398) K and up to 120 MPa. <i>Journal of Chemical Thermodynamics</i> , <b>2016</b> , 93, 24-33	2.9	21
186	Volumetric behavior of the ternary system (methyl tert-butyl ether + methylbenzene + butan-1-ol) and its binary sub-system (methyl tert-butyl ether + butan-1-ol) within the temperature range (298.15 to 328.15) K. <i>Journal of Chemical Thermodynamics</i> , <b>2015</b> , 90, 59-70	2.9	6
185	On the viscosity of two 1-butyl-1-methylpyrrolidinium ionic liquids: Effect of the temperature and pressure. <i>Journal of Chemical Thermodynamics</i> , <b>2015</b> , 87, 43-51	2.9	18
184	Density and isothermal compressibility for two trialkylimidazolium-based ionic liquids at temperatures from (278 to 398) K and up to 120 MPa. <i>Journal of Chemical Thermodynamics</i> , <b>2015</b> , 81, 124-130	2.9	18
183	Thermodynamic scaling of the shear viscosity of Mie n-6 fluids and their binary mixtures. <i>Journal of Chemical Physics</i> , <b>2015</b> , 142, 174501	3.9	15
182	Pressure dependence of the solubility of light fullerenes in 1-hexanol from 298.15K to 363.15K. <i>Journal of Molecular Liquids</i> , <b>2015</b> , 209, 71-76	6	2
181	Krytox GPL102 Oil as Reference Fluid for High Viscosities: High Pressure Volumetric Properties, Heat Capacities, and Thermal Conductivities. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2015</b> , 60, 3660-3669	2.8	8
180	Thermodynamic Properties of Dichloromethane, Bromochloromethane, and Dibromomethane under Elevated Pressure: Experimental Results and SAFT-VR Mie Predictions. <i>Industrial &amp; Engineering Chemistry Research</i> , <b>2015</b> , 54, 720-730	3.9	11
179	Interactions and structure of ionic liquids on graphene and carbon nanotubes surfaces. <i>RSC Advances</i> , <b>2014</b> , 4, 18017-18024	3.7	61
178	High pressure density and solubility for the CO <sub>2</sub> +1-ethyl-3-methylimidazolium ethylsulfate system. <i>Journal of Supercritical Fluids</i> , <b>2014</b> , 88, 46-55	4.2	20
177	Viscosity measurements for squalane at high pressures to 350MPa from T=(293.15 to 363.15)K. <i>Journal of Chemical Thermodynamics</i> , <b>2014</b> , 69, 201-208	2.9	44

176	Reference Correlations for the Density and Viscosity of Squalane from 273 to 473 K at Pressures to 200 MPa. <i>Journal of Physical and Chemical Reference Data</i> , <b>2014</b> , 43, 013104	4.3	32
175	Density and viscosity of three (2,2,2-trifluoroethanol+1-butyl-3-methylimidazolium) ionic liquid binary systems. <i>Journal of Chemical Thermodynamics</i> , <b>2014</b> , 70, 101-110	2.9	80
174	Pressure-viscosity behaviour and film thickness in elastohydrodynamic regime of lubrication of ionic liquids and other base oils. <i>Lubrication Science</i> , <b>2014</b> , 26, 449-462	1.3	18
173	Bulk and liquid-vapor interface of pyrrolidinium-based ionic liquids: a molecular simulation study. <i>Journal of Physical Chemistry B</i> , <b>2014</b> , 118, 731-42	3.4	47
172	Compressibilities and Viscosities of Reference, Vegetable, and Synthetic Gear Lubricants. <i>Industrial &amp; Engineering Chemistry Research</i> , <b>2014</b> , 53, 4499-4510	3.9	6
171	Ionic liquids based on phosphonium cations as neat lubricants or lubricant additives for a steel/steel contact. <i>ACS Applied Materials &amp; Interfaces</i> , <b>2014</b> , 6, 13115-28	9.5	102
170	Correlation and Prediction of Dense Fluid Transport Coefficients. IX. Ionic Liquids. <i>International Journal of Thermophysics</i> , <b>2014</b> , 35, 812-829	2.1	19
169	Tribo-chemical reactions of anion in pyrrolidinium salts for steel/steel contact. <i>Tribology International</i> , <b>2014</b> , 77, 160-170	4.9	18
168	Long-term thermal stability of some 1-butyl-1-methylpyrrolidinium ionic liquids. <i>Journal of Chemical Thermodynamics</i> , <b>2014</b> , 74, 51-57	2.9	40
167	Tribological Properties of Two Bis(trifluoromethylsulfonyl)imide-Based Ionic Liquids on Steel/Steel Contact. <i>Tribology Transactions</i> , <b>2014</b> , 57, 637-646	1.8	11
166	Ionic liquids as hydraulic fluids: comparison of several properties with those of conventional oils. <i>Lubrication Science</i> , <b>2014</b> , 26, 488-499	1.3	20
165	Phase equilibrium of two CO <sub>2</sub> + biodegradable oil systems up to 72MPa. <i>Journal of Supercritical Fluids</i> , <b>2014</b> , 91, 90-97	4.2	4
164	Friction and anti-wear properties of two tris(pentafluoroethyl)trifluorophosphate ionic liquids as neat lubricants. <i>Tribology International</i> , <b>2014</b> , 70, 104-111	4.9	57
163	High pressure viscosity characterization of four vegetable and mineral hydraulic oils. <i>Industrial Crops and Products</i> , <b>2014</b> , 54, 281-290	5.9	16
162	Excess volumes and excess heat capacities for alkanediol+water systems in the temperature interval (283.15-313.15)K. <i>Fluid Phase Equilibria</i> , <b>2013</b> , 356, 1-10	2.5	31
161	Reference Correlation of the Viscosity of Squalane from 273 to 373 K at 0.1 MPa. <i>Journal of Physical and Chemical Reference Data</i> , <b>2013</b> , 42, 033101	4.3	34
160	Long-term thermal stability of five imidazolium ionic liquids. <i>Journal of Chemical Thermodynamics</i> , <b>2013</b> , 65, 184-190	2.9	66
159	Influence of the pressure, temperature, cation and anion on the volumetric properties of ionic liquids: New experimental values for two salts. <i>Journal of Chemical Thermodynamics</i> , <b>2013</b> , 58, 440-448	2.9	33

158	Experimental measurements and modeling of CO <sub>2</sub> solubility in sunflower, castor and rapeseed oils. <i>Journal of Supercritical Fluids</i> , <b>2013</b> , 82, 191-199	4.2	4
157	Pressure dependence on the viscosities of 1-butyl-2,3-dimethylimidazolium bis(trifluoromethylsulfonyl)imide and two tris(pentafluoroethyl)trifluorophosphate based ionic liquids: New measurements and modelling. <i>Journal of Chemical Thermodynamics</i> , <b>2013</b> , 62, 162-169	2.9	32
156	Carbon dioxide solubility in reference and vegetable lubricants developed for two stroke engines. <i>Journal of Supercritical Fluids</i> , <b>2012</b> , 68, 123-130	4.2	9
155	Experimental density and viscosity measurements of di(2ethylhexyl)sebacate at high pressure. <i>Journal of Chemical Thermodynamics</i> , <b>2012</b> , 44, 38-43	2.9	36
154	High pressure volumetric properties of 1-ethyl-3-methylimidazolium ethylsulfate and 1-(2-methoxyethyl)-1-methyl-pyrrolidinium bis(trifluoromethylsulfonyl)imide. <i>Journal of Chemical Thermodynamics</i> , <b>2012</b> , 48, 213-220	2.9	44
153	An experimental setup for isobaric heat capacities for viscous fluids at high pressure: Squalane, bis(2-ethylhexyl) sebacate and bis(2-ethylhexyl) phthalate. <i>Journal of Chemical Thermodynamics</i> , <b>2012</b> , 49, 75-80	2.9	8
152	Pressure Viscosity Coefficients for Polyalkylene Glycol Oils and Other Ester or Ionic Lubricants. <i>Tribology Letters</i> , <b>2012</b> , 45, 89-100	2.8	31
151	On the density scaling of pVT data and transport properties for molecular and ionic liquids. <i>Journal of Chemical Physics</i> , <b>2012</b> , 136, 214502	3.9	15
150	Using molecular simulation to understand the structure of [C <sub>2</sub> C <sub>1</sub> im] <sup>+</sup> -alkylsulfate ionic liquids: bulk and liquid-vapor interfaces. <i>Journal of Physical Chemistry B</i> , <b>2012</b> , 116, 14159-70	3.4	27
149	Effect of the pressure on the viscosities of ionic liquids: Experimental values for 1-ethyl-3-methylimidazolium ethylsulfate and two bis(trifluoromethyl-sulfonyl)imide salts. <i>Journal of Chemical Thermodynamics</i> , <b>2012</b> , 54, 302-309	2.9	35
148	Density scaling of the transport properties of molecular and ionic liquids. <i>Journal of Chemical Physics</i> , <b>2011</b> , 134, 144507	3.9	85
147	Influence of Molecular Structure on Densities and Viscosities of Several Ionic Liquids. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2011</b> , 56, 4984-4999	2.8	137
146	Compressibilities and viscosities of reference and vegetable oils for their use as hydraulic fluids and lubricants. <i>Green Chemistry</i> , <b>2011</b> , 13, 1293	10	45
145	Vapor-Pressure Measurements and Modeling of Dipentaerythritol Ester Lubricants. <i>Industrial &amp; Engineering Chemistry Research</i> , <b>2011</b> , 50, 4231-4237	3.9	6
144	On the isobaric thermal expansivity of liquids. <i>Journal of Chemical Physics</i> , <b>2011</b> , 134, 094502	3.9	31
143	High pressure densities of carbon dioxide+dipentaerythritol hexaheptanoate: New experimental setup and volumetric behavior. <i>Journal of Supercritical Fluids</i> , <b>2011</b> , 58, 189-197	4.2	10
142	Scaling of the viscosity of the Lennard-Jones chain fluid model, argon, and some normal alkanes. <i>Journal of Chemical Physics</i> , <b>2011</b> , 134, 064505	3.9	76
141	Effect of Water on the Viscosities and Densities of 1-Butyl-3-methylimidazolium Dicyanamide and 1-Butyl-3-methylimidazolium Tricyanomethane at Atmospheric Pressure. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2010</b> , 55, 645-652	2.8	200

140	How Pressure Affects the Dynamic Viscosities of Two Poly(propylene glycol) Dimethyl Ether Lubricants. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2010</b> , 55, 4088-4094	2.8	20
139	Isobaric Thermal Expansivity of Highly Polar Nitrogen Compounds at Temperatures from (278.15 to 348.15) K and at Pressures from (5 to 55) MPa. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2010</b> , 55, 1537-1541	2.8	5
138	Influence of the Molecular Structure on the Volumetric Properties and Viscosities of Dialkyl Adipates (Dimethyl, Diethyl, and Diisobutyl Adipates). <i>Journal of Chemical &amp; Engineering Data</i> , <b>2010</b> , 55, 3697-3703	2.8	23
137	Solubilities of Carbon Dioxide in a Dipentaerythritol Ester and in a Polyether. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2010</b> , 55, 5483-5488	2.8	8
136	Experimental Dynamic Viscosities of Dipentaerythritol Ester Lubricants at High Pressure. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2010</b> , 55, 3216-3223	2.8	18
135	Pressure and Temperature Dependence of Isobaric Heat Capacity for [Emim][BF <sub>4</sub> ], [Bmim][BF <sub>4</sub> ], [Hmim][BF <sub>4</sub> ], and [Omim][BF <sub>4</sub> ]. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2010</b> , 55, 600-604	2.8	53
134	Isobaric Thermal Expansivity for Nonpolar Compounds. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2010</b> , 55, 2173-2179	2.8	16
133	Unusual Behavior of the Thermodynamic Response Functions of Ionic Liquids. <i>Journal of Physical Chemistry Letters</i> , <b>2010</b> , 1, 211-214	6.4	41
132	Excess molar enthalpies of dichloropropane + n-alkane mixtures. <i>Journal of Thermal Analysis and Calorimetry</i> , <b>2010</b> , 101, 1121-1125	4.1	2
131	Temperature and pressure dependences of volumetric properties of two poly(propylene glycol) dimethyl ether lubricants. <i>Journal of Chemical Thermodynamics</i> , <b>2010</b> , 42, 84-89	2.9	27
130	Solubility of carbon dioxide in pentaerythritol ester oils. New data and modeling using the PC-SAFT model. <i>Journal of Supercritical Fluids</i> , <b>2010</b> , 55, 62-70	4.2	17
129	Compressed liquid densities of two dipentaerythritol esters. <i>Fluid Phase Equilibria</i> , <b>2010</b> , 296, 30-36	2.5	13
128	Isobaric thermal expansivity of the binary system 1-hexanol+n-hexane as a function of temperature and pressure. <i>Fluid Phase Equilibria</i> , <b>2009</b> , 276, 1-6	2.5	14
127	Excess molar properties for binary systems of alkylimidazolium-based ionic liquids + nitromethane. Experimental results and ERAS-model calculations. <i>Journal of Chemical Thermodynamics</i> , <b>2009</b> , 41, 334-347	2.9	105
126	Excess enthalpy, density, and heat capacity for binary systems of alkylimidazolium-based ionic liquids + water. <i>Journal of Chemical Thermodynamics</i> , <b>2009</b> , 41, 161-166	2.9	162
125	Automated densimetric system: Measurements and uncertainties for compressed fluids. <i>Journal of Chemical Thermodynamics</i> , <b>2009</b> , 41, 632-638	2.9	109
124	Study of the effects of pressure on the viscosity and density of diisodecyl phthalate. <i>Journal of Chemical Thermodynamics</i> , <b>2009</b> , 41, 1007-1015	2.9	53
123	Solubility of HFC134a (1,1,1,2-Tetrafluoroethane) in Two Dialkyl Carbonates. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2009</b> , 54, 2609-2615	2.8	7

122	Thermophysical Characterization of Liquids Using Precise Density and Isobaric Heat Capacity Measurements As a Function of Pressure. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2009</b> , 54, 904-915	2.8	49
121	Solubility of Carbon Dioxide in Two Pentaerythritol Ester Oils between (283 and 333) K. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2008</b> , 53, 1854-1861	2.8	39
120	Relationship between viscosity coefficients and volumetric properties using a scaling concept for molecular and ionic liquids. <i>Journal of Physical Chemistry B</i> , <b>2008</b> , 112, 5563-74	3.4	79
119	Excess properties for binary systems ionic liquid+ethanol: Experimental results and theoretical description using the ERAS model. <i>Fluid Phase Equilibria</i> , <b>2008</b> , 274, 59-67	2.5	145
118	The Pressure-Viscosity Coefficient of Several Ionic Liquids. <i>Tribology Letters</i> , <b>2008</b> , 31, 107-118	2.8	89
117	Viscosities for Ionic Liquid Binary Mixtures with a Common Ion. <i>Journal of Solution Chemistry</i> , <b>2008</b> , 37, 677-688	1.8	98
116	High-pressure viscosity and density of carbon dioxide + pentaerythritol ester mixtures: Measurements and modeling. <i>AIChE Journal</i> , <b>2008</b> , 54, 1625-1636	3.6	16
115	Density and refractive index in mixtures of ionic liquids and organic solvents: Correlations and predictions. <i>Journal of Chemical Thermodynamics</i> , <b>2008</b> , 40, 949-956	2.9	120
114	Viscosity and density measurements for carbon dioxide + pentaerythritol ester lubricant mixtures at low lubricant concentration. <i>Journal of Supercritical Fluids</i> , <b>2008</b> , 44, 172-185	4.2	44
113	Phase and volumetric behavior of binary systems containing carbon dioxide and lubricants for transcritical refrigeration cycles. <i>Journal of Supercritical Fluids</i> , <b>2008</b> , 45, 261-271	4.2	13
112	Experimental and PC-SAFT volumetric and phase behavior of carbon dioxide+PAG or POE lubricant systems. <i>Journal of Supercritical Fluids</i> , <b>2008</b> , 47, 8-16	4.2	16
111	Density and Heat Capacity as a Function of Temperature for Binary Mixtures of 1-Butyl-3-methylpyridinium Tetrafluoroborate + Water, + Ethanol, and + Nitromethane. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2007</b> , 52, 2261-2265	2.8	60
110	Heat capacity of associated systems. Experimental data and application of a two-state model to pure liquids and mixtures. <i>Journal of Physical Chemistry B</i> , <b>2007</b> , 111, 1119-28	3.4	48
109	Dynamic Viscosity under Pressure for Mixtures of Pentaerythritol Ester Lubricants with 32 Viscosity Grade: Measurements and Modeling. <i>Industrial &amp; Engineering Chemistry Research</i> , <b>2007</b> , 46, 1826-1833	3.9	19
108	Density Measurements under Pressure for Mixtures of Pentaerythritol Ester Lubricants. Analysis of a Density-Viscosity Relationship. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2007</b> , 52, 1429-1436	2.8	29
107	Vapor pressure measurements in the range 10 <sup>5</sup> Pa to 1 Pa of four pentaerythritol esters: Density and vapor-liquid equilibria modeling of ester lubricants. <i>Fluid Phase Equilibria</i> , <b>2007</b> , 260, 248-261	2.5	41
106	Volumetric properties of 1-iodoperfluorohexane+n-octane binary system at several temperatures. <i>Journal of Thermal Analysis and Calorimetry</i> , <b>2007</b> , 87, 179-187	4.1	3
105	General friction theory viscosity model for the PC-SAFT equation of state. <i>AIChE Journal</i> , <b>2006</b> , 52, 1600-1610	3.6	63



104	P <sup>∞</sup> Measurements and Equation of State (EoS) Predictions of Ester Lubricants up to 45 MPa. <i>Industrial &amp; Engineering Chemistry Research</i> , <b>2006</b> , 45, 1172-1182	3.9	41
103	Relationship between Viscosity Coefficients and Volumetric Properties: Measurements and Modeling for Pentaerythritol Esters. <i>Industrial &amp; Engineering Chemistry Research</i> , <b>2006</b> , 45, 9171-9183	3.9	28
102	Thermodynamic Properties of Imidazolium-Based Ionic Liquids: Densities, Heat Capacities, and Enthalpies of Fusion of [bmim][PF <sub>6</sub> ] and [bmim][NTf <sub>2</sub> ]. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2006</b> , 51, 1856-1859	2.8	240
101	High-Pressure Characterization of Dynamic Viscosity and Derived Properties for Squalane and Two Pentaerythritol Ester Lubricants: Pentaerythritol Tetra-2-ethylhexanoate and Pentaerythritol Tetranonanoate. <i>Industrial &amp; Engineering Chemistry Research</i> , <b>2006</b> , 45, 2394-2404	3.9	56
100	Temperature and Pressure Dependences of Thermophysical Properties of Some Ethylene Glycol Dimethyl Ethers from Ultrasonic Measurements. <i>International Journal of Thermophysics</i> , <b>2006</b> , 27, 1354-1372	2.7	18
99	Volumetric behaviour of the environmentally compatible lubricants pentaerythritol tetraheptanoate and pentaerythritol tetranonanoate at high pressures. <i>Green Chemistry</i> , <b>2005</b> , 7, 775	1.0	50
98	Experimental Dynamic Viscosities of 2,3-Dimethylpentane up to 60 MPa and from (303.15 to 353.15) K Using a Rolling-Ball Viscometer. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2005</b> , 50, 849-855	2.8	46
97	Compressed Liquid Densities of Squalane and Pentaerythritol Tetra(2-ethylhexanoate). <i>Journal of Chemical &amp; Engineering Data</i> , <b>2005</b> , 50, 939-946	2.8	96
96	Density measurements under pressure for the binary system (ethanol+methylcyclohexane). <i>Journal of Chemical Thermodynamics</i> , <b>2005</b> , 37, 1294-1304	2.9	22
95	Phase and viscosity behaviour of refrigerant-lubricant mixtures. <i>International Journal of Refrigeration</i> , <b>2005</b> , 28, 714-724	3.8	33
94	Influence of the molecular structure on the viscosity of some alkoxyethanols. <i>Fluid Phase Equilibria</i> , <b>2005</b> , 236, 229-236	2.5	19
93	Volumetric properties under pressure for the binary system ethanol+toluene. <i>Fluid Phase Equilibria</i> , <b>2005</b> , 235, 139-151	2.5	56
92	Thermodynamic properties of 1,1,1,2-tetrafluoroethane + polyether mixtures up to 60 MPa. <i>Canadian Journal of Chemistry</i> , <b>2004</b> , 82, 1271-1279	0.9	1
91	Influence of the number of CH <sub>2</sub> ?CH <sub>2</sub> ?O groups on the viscosity of polyethylene glycol dimethyl ethers at high pressure. <i>Fluid Phase Equilibria</i> , <b>2004</b> , 222-223, 331-338	2.5	15
90	Prediction of the pressure dependence on the thermodynamic properties of dialkyl carbonate + alkane mixtures using Nitta-Chao model. <i>Fluid Phase Equilibria</i> , <b>2004</b> , 217, 165-173	2.5	4
89	Experimental and Predicted Solubilities of HFC134a (1,1,1,2-Tetrafluoroethane) in Polyethers. <i>Industrial &amp; Engineering Chemistry Research</i> , <b>2004</b> , 43, 1523-1529	3.9	17
88	Heat Capacities, Densities, and Speeds of Sound for {(1,5-Dichloropentane or 1,6-Dichlorohexane) + Dodecane}. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2004</b> , 49, 333-338	2.8	21
87	Dynamic Viscosity for HFC-134a + Polyether Mixtures up to 373.15 K and 140 MPa at Low Polyether Concentration. Measurements and Modeling. <i>Industrial &amp; Engineering Chemistry Research</i> , <b>2004</b> , 43, 804-814	3.9	29

86	$p$ - $x$ Data for the Dimethyl Carbonate + Decane System. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2004</b> , 49, 923-927	2.8	34
85	Phase Equilibria, PVT Behavior, and Critical Phenomena in Carbon Dioxide + n-Alkane Mixtures Using the Perturbed-Chain Statistical Associating Fluid Theory Approach. <i>Industrial &amp; Engineering Chemistry Research</i> , <b>2004</b> , 43, 8345-8353	3.9	46
84	$p$ - $T$ Measurements and EoS Predictions of Glycol Ethers from (283.15 to 353.15) K at Pressures up to 25 MPa. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2004</b> , 49, 1400-1405	2.8	18
83	High-Pressure Volumetric Properties of Three Monoethylene Glycol Alkyl Ethers. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2004</b> , 49, 1344-1349	2.8	17
82	Liquid Density Measurements of Diethylene Glycol Monoalkyl Ethers as a Function of Temperature and Pressure. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2004</b> , 49, 376-379	2.8	26
81	Volumetric Properties of Binary Tetraethylene Glycol Dimethyl Ether + Heptane Mixtures between (278.15 and 353.15) K and up to 25 MPa. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2003</b> , 48, 1271-1278	2.8	17
80	Modeling of Gas Solubility Data for HFCs-Lubricant Oil Binary Systems by Means of the SRK Equation of State. <i>International Journal of Thermophysics</i> , <b>2003</b> , 24, 1043-1060	2.1	15
79	Viscosity measurements and correlations for 1,1,1,2-tetrafluoroethane (HFC-134a) up to 140 MPa. <i>Fluid Phase Equilibria</i> , <b>2003</b> , 210, 21-32	2.5	14
78	Sakou-Mu-Braunsnitz equation of state for modelling phase equilibria and high-pressure PVT of mixtures containing dialkyl carbonate and alkane. <i>Fluid Phase Equilibria</i> , <b>2003</b> , 210, 77-89	2.5	2
77	Experimental densities and dynamic viscosities of organic carbonate + n-alkane or p-xylene systems at 298.15 K. <i>Fluid Phase Equilibria</i> , <b>2003</b> , 204, 233-243	2.5	29
76	Phase equilibria and pVT predictions for alkyl carbonate + n-alkane systems using equations of state. <i>Fluid Phase Equilibria</i> , <b>2003</b> , 212, 111-128	2.5	14
75	Thermophysical Properties of Two Poly(alkylene glycol) Derivative Lubricants from High Pressure Acoustic Measurements. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2003</b> , 48, 1208-1213	2.8	11
74	Volumetric Properties of Monoethylene Glycol Dimethyl Ether and Diethylene Glycol Dimethyl Ether up to 60 MPa. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2003</b> , 48, 1044-1049	2.8	30
73	Pressure and temperature dependence of the excess thermodynamic properties of binary dimethyl carbonate + n-octane mixtures. <i>Canadian Journal of Chemistry</i> , <b>2003</b> , 81, 840-849	0.9	6
72	UNIFAC calculation of thermodynamic properties of binary 1-chloroalkane + alkane and 1,1-dichloroalkane + alkane mixtures: Comparison with Nitta-Chao and DISQUAC predictions. <i>Canadian Journal of Chemistry</i> , <b>2003</b> , 81, 392-405	0.9	2
71	$p$ - $x$ measurements for HFC-134a + triethylene glycol dimethylether system. <i>Fluid Phase Equilibria</i> , <b>2002</b> , 199, 185-195	2.5	31
70	High pressure viscosity and density modeling of two polyethers and two dialkyl carbonates. <i>Fluid Phase Equilibria</i> , <b>2002</b> , 199, 249-263	2.5	39
69	( $p$ , $V_m$ , $T$ , $x$ ) measurements of dimethyl carbonate + octane binary mixtures: II. Excess molar volumes. <i>Fluid Phase Equilibria</i> , <b>2002</b> , 199, 135-145	2.5	18

68	Modelling of PVT for some poly alkylene glycol lubricants using Sakou-Mu-Braunsnitz EOS. <i>Fluid Phase Equilibria</i> , <b>2002</b> , 199, 23-31	2.5	4
67	Determination of Excess Molar Enthalpies of 1,1-Dichloroalkane + 1-Butanol or 1-Heptanol Mixtures at 298.15 K. Analysis and Comparison with Predicted Values of UNIFAC. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2002</b> , 47, 411-415	2.8	9
66	High-Pressure Volumetric Behavior of 1,1,1,2-Tetrafluoroethane + (1,2,5,8,11,14-Pentaoxapentadecane (TEGDME) Mixtures. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2002</b> , 47, 233-238	2.8	20
65	Excess molar enthalpies of mixtures of methyl derivatives of polyethylene glycol with 1-alkanol at 298.15 K and 101.3 kPa. <i>Canadian Journal of Chemistry</i> , <b>2002</b> , 80, 462-466	0.9	14
64	(p, Vm, T, x) measurements of dimethyl carbonate+octane binary mixtures. <i>Fluid Phase Equilibria</i> , <b>2001</b> , 186, 235-255	2.5	89
63	Reply to the letter to the editor by J. Gmehling and J. Lohmann about the paper Analysis of the molecular interactions of organic anhydride + alkane binary mixtures using the Nitta-Chao model [Fluid Phase Equilib. 170 (2000) 69-85]. <i>Fluid Phase Equilibria</i> , <b>2001</b> , 189, 197-201	2.5	
62	High-Pressure Measurements of the Viscosity and Density of Two Polyethers and Two Dialkyl Carbonates. <i>International Journal of Thermophysics</i> , <b>2001</b> , 22, 749-768	2.1	125
61	Thermodynamics of Rubber Elasticity. <i>Journal of Chemical Education</i> , <b>2001</b> , 78, 263	2.4	13
60	Modelling thermodynamic properties of iodoalkane + alkane systems using group contribution models. <i>Physical Chemistry Chemical Physics</i> , <b>2001</b> , 3, 5006	3.6	2
59	Temperature dependence of the excess molar volume of (dimethyl carbonate, or diethyl carbonate+ toluene) from T= 278.15 K to 323.15 K. <i>Journal of Chemical Thermodynamics</i> , <b>2000</b> , 32, 743-754	2.9	35
58	Analysis of the molecular interactions of organic anhydride+alkane binary mixtures using the Nitta-Chao model. <i>Fluid Phase Equilibria</i> , <b>2000</b> , 170, 69-85	2.5	4
57	pVT Measurements of Polyethylene Glycol Dimethylethers Between 278.15 and 328.15 K at Pressures to 12 MPa. <i>International Journal of Thermophysics</i> , <b>2000</b> , 21, 831-851	2.1	49
56	Thermophysical properties of methanol+some polyethylene glycol dimethyl ether by UNIFAC and DISQUAC group-contribution models for absorption heat pumps. <i>Fluid Phase Equilibria</i> , <b>1999</b> , 155, 327-337	2.5	13
55	Pressure and Temperature Dependence of Excess Enthalpies of Methanol + Tetraethylene Glycol Dimethyl Ether and Methanol + Polyethylene Glycol Dimethyl Ether 250. <i>Journal of Chemical &amp; Engineering Data</i> , <b>1999</b> , 44, 1409-1413	2.8	7
54	Excess Properties of Some Methanol + Amide Systems Proposed as Working Fluids for Absorption Machines. <i>Journal of Chemical &amp; Engineering Data</i> , <b>1999</b> , 44, 309-313	2.8	17
53	Excess molar volumes of liquid 1-bromoalkane + alkane mixtures. Nitta-Chao characterization of the bromine-bromine and bromine-methylene interactions in binary 1-bromoalkane + alkane mixtures. <i>Canadian Journal of Chemistry</i> , <b>1999</b> , 77, 299-307	0.9	4
52	Excess molar enthalpies of (1,1-Dichloroalkane + pentanol, or hexanol) at T=298.15 K. <i>Journal of Chemical Thermodynamics</i> , <b>1998</b> , 30, 1061-1068	2.9	13
51	Characteristic parameters of the Tassios, Larsen and Gmehling versions of the UNIFAC model for enthalpies of mixing in organic anhydrides + N-alkanes mixtures. <i>Thermochimica Acta</i> , <b>1998</b> , 317, 59-64	2.9	5

50	Estimation of parameters of Nitta-Chao model for ester+1-alkanol mixtures. <i>Fluid Phase Equilibria</i> , <b>1998</b> , 148, 49-68	2.5	20
49	Experimental excess volumes of organic carbonate+alkane systems. Estimation of the parameters of the Nitta-Chao model for this kind of binary mixture. <i>Journal of the Chemical Society, Faraday Transactions</i> , <b>1998</b> , 94, 1707-1712		39
48	Estimation of parameters of Nitta-Chao model for linear monoether + 1-alkanol mixtures. <i>Fluid Phase Equilibria</i> , <b>1997</b> , 133, 57-72	2.5	13
47	Experimental and predicted excess enthalpies of the working pairs (methanol or trifluoroethanol + polyglycol ethers) for absorption cycles. <i>Fluid Phase Equilibria</i> , <b>1997</b> , 133, 229-238	2.5	26
46	Orthobaric surface tension of (tetrachloromethane + tetradecafluoromethylcyclohexane) in the critical region at temperatures from T= 303 K to T= 315 K. <i>Journal of Chemical Thermodynamics</i> , <b>1996</b> , 28, 767-778	2.9	6
45	Prediction of enthalpies of mixing and vapor-liquid equilibria for mixtures containing organic carbonates + n-alkanes using several versions of the unifac model. <i>Thermochimica Acta</i> , <b>1996</b> , 286, 321-332	2.9	20
44	Estimation of the Nitta-Chao Parameters for the ether-group. Ether + n-alkane mixtures. <i>Fluid Phase Equilibria</i> , <b>1995</b> , 110, 53-71	2.5	12
43	Experimental and predicted excess enthalpies of the 2,2,2-trifluoroethanol-water-tetraethylene glycol dimethyl ether ternary system using binary mixing data. <i>Journal of the Chemical Society, Faraday Transactions</i> , <b>1995</b> , 91, 2071-2079		20
42	Excess molar volumes of binary mixtures with 2-pentanone and 1-chloroalkanes at 298.15 and 308.15 K. <i>Canadian Journal of Chemistry</i> , <b>1995</b> , 73, 139-145	0.9	12
41	Analysis of Excess Molar Volumes of the Ternary Systems Containing a Propyl Alkanoate and Two Alkanes with Some Predictive Methods. <i>Physics and Chemistry of Liquids</i> , <b>1995</b> , 30, 141-150	1.5	3
40	Prediction of density and excess volume for the ternary mixture: (water + 2,2,2-trifluoroethanol + 2,5,8,11,14-pentaoxapentadecane) from experimental binary values at temperatures from 283.15 K to 333.15 K. <i>Journal of Chemical Thermodynamics</i> , <b>1995</b> , 27, 281-292	2.9	22
39	Analysis of the intramolecular proximity effect on dichloroalkane + alkane mixtures using Nitta-Chao model. <i>Fluid Phase Equilibria</i> , <b>1995</b> , 110, 31-51	2.5	16
38	Excess volumes and viscosities of the ternary system water-trifluoroethanol-tetraethylene glycol dimethyl ether at 303.15 K. <i>Thermochimica Acta</i> , <b>1995</b> , 259, 57-70	2.9	14
37	Densities, viscosities, and excess properties of trifluoroethanol-water, tetraethylene glycol dimethylether-water, and trifluoroethanol-tetraethylene glycol dimethylether at 303.15 K. <i>International Journal of Thermophysics</i> , <b>1994</b> , 15, 661-674	2.1	23
36	Estimation of DISQUAC interaction parameters for low molecular mass analogues of polymers: Chloroalkane + ester mixtures. <i>Journal of Solution Chemistry</i> , <b>1994</b> , 23, 135-152	1.8	26
35	Intramolecular-proximity effect on the excess enthalpies of (a dichloroalkane + an alkan-2-one). <i>Journal of Chemical Thermodynamics</i> , <b>1994</b> , 26, 53-59	2.9	11
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32	Measurements and analysis of the excess enthalpies of some dichloroalkane + 2-ketone systems using UNIFAC group-contribution model. <i>Canadian Journal of Chemistry</i> , <b>1994</b> , 72, 304-307	0.9	2
31	VE of binary mixtures containing 3-pentanone or 3-heptanone with 1-chloroalkanes at 298.15 K. <i>Journal of Chemical &amp; Engineering Data</i> , <b>1994</b> , 39, 172-174	2.8	11
30	Estimation of Nitta-Chao Interaction Parameters for Low Molecular Mass Analogues of Polymers. 1-Chloroalkane + n-alkane mixtures. <i>Zeitschrift Fur Elektrotechnik Und Elektrochemie</i> , <b>1993</b> , 97, 106-113		10
29	Excess molar volumes of (1-chloropentane or 1-chlorohexane + heptane + decane) at the temperature 298.15 K. <i>Journal of Chemical Thermodynamics</i> , <b>1993</b> , 25, 1041-1047	2.9	6
28	Excess molar enthalpies of some examples of (a dichloroalkane+a ket-2-one) at the temperature 298.15 K. <i>Journal of Chemical Thermodynamics</i> , <b>1993</b> , 25, 1127-1132	2.9	11
27	Excess molar volumes at the temperature 298.15 K of $\{x_1C_2H_5CO_2(CH_2)_2CH_3 + x_2Cl(CH_2)_5CH_3 + (1-x_1-x_2)c-C_6H_{12}\}$ , $\{x_1C_2H_5CO_2(CH_2)_2CH_3 + x_2Cl(CH_2)_5CH_3 + (1-x_1-x_2)CH_3(CH_2)_4CH_3\}$ , and $\{x_1C_2H_5CO_2(CH_2)_2CH_3 + x_2c-C_6H_{12} + (1-x_1-x_2)CH_3(CH_2)_4CH_3\}$ . <i>Journal of Chemical Thermodynamics</i> , <b>1993</b> , 25, 1133-1139	2.9	21
26	Excess molar enthalpies at the temperature 298.15 K of (an n-alkyl formate + an n-alkanol) IV. $\{xHCO_2(CH_2)_iCH_3 + (1-x)C_jH_{2j+1}OH\}$ (i = 0 to 3 and j = 1 and 2). <i>Journal of Chemical Thermodynamics</i> , <b>1992</b> , 24, 809-814	2.9	13
25	Excess enthalpies of some 2-alkanone + 1-chloroalkane binary mixtures at 25 and 35°C. <i>Journal of Solution Chemistry</i> , <b>1991</b> , 20, 115-124	1.8	23
24	Measurements and Analysis of Excess Enthalpies of Ester + n-Alkane Using the UNIFAC Model. <i>Zeitschrift Fur Elektrotechnik Und Elektrochemie</i> , <b>1991</b> , 95, 128-135		33
23	Excess Molar Volumes of Binary Mixtures of Propyl Ethanoate with some n-Alkanes at 298.15 K and 308.15 K. <i>Physics and Chemistry of Liquids</i> , <b>1991</b> , 24, 13-20	1.5	8
22	Thermodynamic properties of binary mixtures of 2-hexanone with n-alkanes at 35°C. <i>Journal of Solution Chemistry</i> , <b>1990</b> , 19, 1095-1102	1.8	18
21	Analysis of thermodynamic properties of 1-alkanol + n-alkane mixtures using the nitta-chao group contribution model. <i>Fluid Phase Equilibria</i> , <b>1990</b> , 55, 293-308	2.5	46
20	Analysis of excess enthalpies of ethyl formate + n-alkane or 1-alkanol with two group contribution models. <i>Fluid Phase Equilibria</i> , <b>1990</b> , 56, 219-234	2.5	37
19	Excess molar volumes of (methyl butanoate + n-heptane + n-decane) and of (butyl butanoate + n-heptane + n-decane) at 298.15 K. <i>Journal of Chemical Thermodynamics</i> , <b>1990</b> , 22, 865-871	2.9	18
18	Thermodynamic properties of (a propyl ester + an n-alkane) at 298.15 K II. $\{xC_3H_7CO_2C_3H_7 + (1-x)n-C_mH_{2m+2}\}$ , (m = 6 to 10). <i>Journal of Chemical Thermodynamics</i> , <b>1990</b> , 22, 263-268	2.9	20
17	Darc analysis of binary mixtures. Excess enthalpies of ketone + alkane and ketone + alcohol systems. <i>Thermochimica Acta</i> , <b>1989</b> , 156, 21-26	2.9	
16	Thermodynamic properties of (a propyl ester + an n-alkane) at 298.15 K I. $\{xC_2H_5CO_2C_3H_7 + (1-x)n-C_nH_{2n+2}\}$ , (n = 6 to 10). <i>Journal of Chemical Thermodynamics</i> , <b>1989</b> , 21, 1017-1022	2.9	41
15	Analysis of excess enthalpies of ester+ 1-chloroalkanes with two group contribution models: primary parameters. <i>Fluid Phase Equilibria</i> , <b>1988</b> , 43, 295-316	2.5	39

14	Isobaric expansivities of the binary mixtures $C_3H_7(OH) + C_nH_{2n+2}$ ( $n = 11, 12$ ) between 288.15 and 318.15 K. <i>Thermochimica Acta</i> , <b>1988</b> , 131, 57-64	2.9	24
13	Excess molar enthalpies at 298.15 K of (an n-alkyl formate + an n-alkanol) III. $\{xHCO_2CH_3 + (1-x)C_nH_{2n+1}OH\}$ , ( $n = 3$ to 10). <i>Journal of Chemical Thermodynamics</i> , <b>1988</b> , 20, 1315-1319	2.9	16
12	Excess molar volumes of (ethyl formate or ethyl acetate + 1-chloroalkane) at 298.15 K. <i>Journal of Chemical &amp; Engineering Data</i> , <b>1987</b> , 32, 464-466	2.8	59
11	Excess Gibbs energies and excess volumes of some alcohol-methyl ester binary mixtures. <i>Journal of Chemical &amp; Engineering Data</i> , <b>1987</b> , 32, 17-22	2.8	12
10	Excess molar enthalpies at 298.15 K of (an n-alkyl formate + an n-alkanol) II. , ( $n = 3$ to 10). <i>Journal of Chemical Thermodynamics</i> , <b>1987</b> , 19, 147-150	2.9	8
9	Vapor-liquid equilibrium of the binary mixtures $C_nH_{2n+1}OH$ ( $n = 2,3,4$ ) + Propyl Ethanoate and + Ethyl Propanoate. <i>Canadian Journal of Chemical Engineering</i> , <b>1987</b> , 65, 982-990	2.3	22
8	Excess molar enthalpies of butyl acetate + an n-alkane at 298.15 K. <i>Fluid Phase Equilibria</i> , <b>1986</b> , 28, 183-189	2.9	20
7	Excess molar enthalpies at 298.15 K of (an n-alkyl formate + an n-alkanol) I. $\{xHCO_2(CH_2)_3CH_3 + (1-x)C_nH_{2n+1}OH\}$ , ( $n = 3$ to 10). <i>Journal of Chemical Thermodynamics</i> , <b>1986</b> , 18, 1003-1006	2.9	7
6	Excess volumes of (toluene + each of several isomers of hexanol) at 298.15 K. <i>Journal of Chemical Thermodynamics</i> , <b>1986</b> , 18, 419-422	2.9	5
5	Excess thermodynamics functions of 1-propanol + methyl propanoate and 1-propanol + methyl butanoate systems. <i>Fluid Phase Equilibria</i> , <b>1985</b> , 20, 145-153	2.5	22
4	Excess enthalpies of 1-heptanol + n-alkane and di-n-propylamine + normal alcohol mixtures at 298.15 K. <i>Journal of Chemical &amp; Engineering Data</i> , <b>1985</b> , 30, 321-323	2.8	43
3	Excess enthalpies of some ester + alcohol binary mixtures. <i>Journal of Chemical &amp; Engineering Data</i> , <b>1985</b> , 30, 318-320	2.8	24
2	Excess enthalpies of (methyl propanoate or methyl pentanoate + 1-alkanol) at 298.15 K. <i>Journal of Chemical Thermodynamics</i> , <b>1983</b> , 15, 1145-1149	2.9	8
1	Excess enthalpies of (secondary amine + alcohol) at 298.15 K. <i>Journal of Chemical Thermodynamics</i> , <b>1983</b> , 15, 581-584	2.9	40