Gabriele Sadowski

List of Publications by Citations

307 papers	10,867 citations	49 h-index	95 g-index
327 ext. papers	12,227 ext. citations	3.8 avg, IF	6.81 L-index

#	Paper	IF	Citations
307	Perturbed-Chain SAFT: An Equation of State Based on a Perturbation Theory for Chain Molecules. *Industrial & Engineering Chemistry Research*, **2001**, 40, 1244-1260	3.9	2283
306	Application of the Perturbed-Chain SAFT Equation of State to Associating Systems. *Industrial & Engineering Chemistry Research*, **2002**, 41, 5510-5515	3.9	794
305	Modeling Polymer Systems Using the Perturbed-Chain Statistical Associating Fluid Theory Equation of State. *Industrial & Engineering Chemistry Research*, **2002**, 41, 1084-1093	3.9	309
304	Application of perturbation theory to a hard-chain reference fluid: an equation of state for square-well chains. *Fluid Phase Equilibria*, **2000**, 168, 183-199	2.5	251
303	Modeling of Aqueous Electrolyte Solutions with Perturbed-Chain Statistical Associated Fluid Theory. *Industrial & Engineering Chemistry Research*, **2005**, 44, 3355-3362	3.9	213
302	Modeling Copolymer Systems Using the Perturbed-Chain SAFT Equation of State. *Industrial & Engineering Chemistry Research*, **2003**, 42, 1266-1274	3.9	170
301	Solubility of Amino Acids: Influence of the pH value and the Addition of Alcoholic Cosolvents on Aqueous Solubility. *Industrial & Engineering Chemistry Research*, **2006**, 45, 6578-6584	3.9	160
300	Modeling aqueous electrolyte solutions. *Fluid Phase Equilibria*, **2008**, 270, 87-96	2.5	145
299	Modeling the solubility of pharmaceuticals in pure solvents and solvent mixtures for drug process design. *Journal of Pharmaceutical Sciences*, **2009**, 98, 4205-15	3.9	143
298	Application of the Perturbed-Chain SAFT equation of state to polar systems. *Fluid Phase Equilibria*, **2004**, 217, 233-239	2.5	129
297	Modeling of Polar Systems Using PCP-SAFT: An Approach to Account for Induced-Association Interactions *Journal of Physical Chemistry C*, **2007**, 111, 15544-15553	3.8	121
296	Design of Contact-Active Antimicrobial Acrylate-Based Materials Using Biocidal Macromers. *Advanced Materials*, **2008**, 20, 104-108	24	112
295	ePC-SAFT revised. *Chemical Engineering Research and Design*, **2014**, 92, 2884-2897	5.5	111
294	Measuring and Modeling Activity Coefficients in Aqueous Amino-Acid Solutions. *Industrial & Engineering Chemistry Research*, **2011**, 50, 131-141	3.9	109
293	Modeling imidazolium-based ionic liquids with ePC-SAFT. *Fluid Phase Equilibria*, **2012**, 335, 64-73	2.5	104
292	Thermodynamic phase behavior of API/polymer solid dispersions. *Molecular Pharmaceutics*, **2014**, 11, 2294-304	5.6	102
291	Hydroformylation of 1-Dodecene in the Thermomorphic Solvent System Dimethylformamide/Decane. Phase Behavior Reaction Performance Catalyst Recycling. *Industrial & Engineering Chemistry Research*, **2012**, 51, 10296-10306	3.9	102

290	Thermodynamic modeling of complex systems using PC-SAFT. *Fluid Phase Equilibria*, **2005**, 228-229, 89-98	2.5	99
289	Modeling of polymer phase equilibria using Perturbed-Chain SAFT. *Fluid Phase Equilibria*, **2002**, 194-197, 541-551	2.5	97
288	Modeling of aqueous amino acid and polypeptide solutions with PC-SAFT. *Chemical Engineering and Processing: Process Intensification*, **2008**, 47, 1018-1025	3.7	91
287	Modeling of Polar Systems with the Perturbed-Chain SAFT Equation of State. Investigation of the Performance of Two Polar Terms. *Industrial & Engineering Chemistry Research*, **2005**, 44, 6928-6938	3.9	89
286	Measurement and modelling of high-pressure phase equilibria in the systems polyethyleneglycol (PEG)–propane, PEG–nitrogen and PEG–carbon dioxide. *Journal of Supercritical Fluids*, **2000**, 17, 1-12	4.2	79
285	PC-SAFT Modeling of CO2 Solubilities in Deep Eutectic Solvents. *Journal of Physical Chemistry B*, **2016**, 120, 2300-10	3.4	78
284	Liquid–liquid Equilibria of 1-Butanol/Water/IL Systems. *Industrial & Engineering Chemistry Research*, **2013**, 52, 18472-18481	3.9	77
283	Modeling aqueous electrolyte solutions. Part 2. Weak electrolytes. *Fluid Phase Equilibria*, **2009**, 279, 141-148	2.5	77
282	Solvent-sensitive reversible stress-response of shape memory natural rubber. *ACS Applied Materials & Interfaces*, **2013**, 5, 3504-7	9.5	76
281	Modeling of aqueous poly(oxyethylene) solutions: 1. Atomistic simulations. *Journal of Physical Chemistry B*, **2008**, 112, 2388-98	3.4	75
280	Compatible solutes: Thermodynamic properties and biological impact of ectoines and prolines. *Biophysical Chemistry*, **2010**, 152, 28-39	3.5	73
279	Measuring and modeling aqueous electrolyte/amino-acid solutions with ePC-SAFT. *Journal of Chemical Thermodynamics*, **2014**, 68, 1-12	2.9	72
278	Long-Term Physical Stability of PVP- and PVPVA-Amorphous Solid Dispersions. *Molecular Pharmaceutics*, **2017**, 14, 157-171	5.6	71
277	Vapor–Liquid Equilibria of Water + Alkylimidazolium-Based Ionic Liquids: Measurements and Perturbed-Chain Statistical Associating Fluid Theory Modeling. *Industrial & Engineering Chemistry Research*, **2014**, 53, 3737-3748	3.9	69
276	Separations and Material Processing in Solutions with Dense Gases. *Industrial & Engineering Chemistry Research*, **1998**, 37, 3208-3220	3.9	69
275	Measurement and PC-SAFT modeling of solid-liquid equilibrium of deep eutectic solvents of quaternary ammonium chlorides and carboxylic acids. *Fluid Phase Equilibria*, **2017**, 448, 69-80	2.5	66
274	Crowders and Cosolvents-Major Contributors to the Cellular Milieu and Efficient Means to Counteract Environmental Stresses. *ChemPhysChem*, **2017**, 18, 2951-2972	3.2	65
273	Influence of copolymer composition on the phase behavior of solid dispersions. *Molecular Pharmaceutics*, **2014**, 11, 4189-98	5.6	62

272	Solubility of sugars and sugar alcohols in ionic liquids: measurement and PC-SAFT modeling. *Journal of Physical Chemistry B*, **2013**, 117, 9980-95	3.4	59
271	Measuring and modeling alcohol/salt systems. *Chemical Engineering Science*, **2012**, 68, 328-339	4.4	58
270	Impact of Polymer Type and Relative Humidity on the Long-Term Physical Stability of Amorphous Solid Dispersions. *Molecular Pharmaceutics*, **2017**, 14, 4374-4386	5.6	57
269	Solubility of Pharmaceuticals and Their Salts As a Function of pH. *Industrial & Engineering Chemistry Research*, **2013**, 52, 2721-2731	3.9	57
268	Experimental investigation and prediction of oiling out during crystallization process. *Journal of Crystal Growth*, **2008**, 310, 4163-4168	1.6	57
267	Measurement and Modeling Solubility of Aqueous Multisolute Amino-Acid Solutions. *Industrial & Engineering Chemistry Research*, **2010**, 49, 1395-1401	3.9	56
266	Cloud-point curves of ethylene–(meth)acrylate copolymers in fluid ethene up to high pressures and temperatures—experimental study and PC–SAFT modeling. *Fluid Phase Equilibria*, **2004**, 215, 263-282	2.5	55
265	Modeling imidazolium-based ionic liquids with ePC-SAFT. Part II. Application to H2S and synthesis-gas components. *Fluid Phase Equilibria*, **2014**, 363, 59-65	2.5	54
264	Modeling of aqueous poly(oxyethylene) solutions. 2. Mesoscale simulations. *Journal of Physical Chemistry B*, **2008**, 112, 13561-71	3.4	51
263	Fickian and Non-Fickian Sorption Kinetics of Toluene in Glassy Polystyrene. *Macromolecules*, **2005**, 38, 8408-8417	5.5	51
262	PC-SAFT modeling of CO2 solubilities in hydrophobic deep eutectic solvents. *Fluid Phase Equilibria*, **2017**, 448, 94-98	2.5	50
261	Influence of humidity on the phase behavior of API/polymer formulations. *European Journal of Pharmaceutics and Biopharmaceutics*, **2015**, 94, 352-62	5.7	50
260	Modeling thermodynamic properties of aqueous single-solute and multi-solute sugar solutions with PC-SAFT. *AICHE Journal*, **2013**, 59, 4794-4805	3.6	49
259	Modeling Poly(N-isopropylacrylamide) Hydrogels in Water/Alcohol Mixtures with PC-SAFT. *Macromolecules*, **2012**, 45, 6686-6696	5.5	49
258	Predicting the Solubility Advantage of Amorphous Pharmaceuticals: A Novel Thermodynamic Approach. *Molecular Pharmaceutics*, **2015**, 12, 2823-33	5.6	48
257	Phase equilibria in polydisperse and associating copolymer solutions: Poly(ethene-co-(meth)acrylic acid)–monomer mixtures. *Fluid Phase Equilibria*, **2006**, 241, 113-123	2.5	48
256	Determination of the Total Vapor Pressure of Hydrophobic Deep Eutectic Solvents: Experiments and Perturbed-Chain Statistical Associating Fluid Theory Modeling. *ACS Sustainable Chemistry and Engineering*, **2019**, 7, 4047-4057	8.3	47
255	Finite and infinite dilution activity coefficients in polycarbonate systems. *Fluid Phase Equilibria*, **1997**, 139, 391-403	2.5	46

254	Amorphous-Amorphous Phase Separation in API/Polymer Formulations. *Molecules*, **2017**, 22,	4.8	45
253	Dissolution of Crystalline Pharmaceuticals: Experimental Investigation and Thermodynamic Modeling. *Industrial & Engineering Chemistry Research*, **2015**, 54, 731-742	3.9	44
252	Modeling aqueous two-phase systems: I. Polyethylene glycol and inorganic salts as ATPS former. *Fluid Phase Equilibria*, **2014**, 368, 91-103	2.5	41
251	Modeling pH and Solubilities in Aqueous Multisolute Amino Acid Solutions. *Industrial & Engineering Chemistry Research*, **2011**, 50, 3503-3509	3.9	41
250	Supercritical antisolvent fractionation: measurements in the systems monodisperse and bidisperse polystyrenecyclohexanecarbon dioxide. *Fluid Phase Equilibria*, **1997**, 139, 349-359	2.5	41
249	Influence of Salts on the Partitioning of 5-Hydroxymethylfurfural in Water/MIBK. *Journal of Physical Chemistry B*, **2016**, 120, 3797-808	3.4	41
248	Characterization and Modeling of the Liquid Phase of Deep Eutectic Solvents Based on Fatty Acids/Alcohols and Choline Chloride. *Industrial & Engineering Chemistry Research*, **2017**, 56, 12192-12202	3.9	40
247	Physical stability of API/polymer-blend amorphous solid dispersions. *European Journal of Pharmaceutics and Biopharmaceutics*, **2018**, 124, 147-157	5.7	39
246	Thermophobicity of liquids: heats of transport in mixtures as pure component properties. *Physical Review Letters*, **2012**, 109, 065901	7.4	39
245	Interfacial tension of binary mixtures exhibiting azeotropic behavior: Measurement and modeling with PCP-SAFT combined with Density Gradient Theory. *Fluid Phase Equilibria*, **2014**, 362, 151-162	2.5	38
244	Modeling of solid–liquid equilibria for systems with solid-complex phase formation. *Fluid Phase Equilibria*, **2007**, 260, 98-104	2.5	38
243	Compatible solutes: Thermodynamic properties relevant for effective protection against osmotic stress. *Fluid Phase Equilibria*, **2016**, 407, 224-235	2.5	37
242	Molecular interactions in 1-butanol + IL solutions by measuring and modeling activity coefficients. *Journal of Physical Chemistry B*, **2013**, 117, 3173-85	3.4	37
241	Thermodynamics of Bioreactions. *Annual Review of Chemical and Biomolecular Engineering*, **2016**, 7, 395-414	8.3	36
240	Development of a group contribution method for polymers within the PC-SAFT model. *Fluid Phase Equilibria*, **2012**, 324, 70-79	2.5	35
239	Predicting the Aqueous Solubility of Pharmaceutical Cocrystals As a Function of pH and Temperature. *Crystal Growth and Design*, **2016**, 16, 2726-2740	3.5	35
238	Thermodynamic phase behaviour of indomethacin/PLGA formulations. *European Journal of Pharmaceutics and Biopharmaceutics*, **2015**, 93, 88-94	5.7	34
237	Salt influence on MIBK/water liquid–liquid equilibrium: Measuring and modeling with ePC-SAFT and COSMO-RS. *Fluid Phase Equilibria*, **2016**, 416, 83-93	2.5	34

236	The Role of Polyfunctionality in the Formation of [Ch]Cl-Carboxylic Acid-Based Deep Eutectic Solvents. *Industrial & Engineering Chemistry Research*, **2018**, 57, 11195-11209	3.9	34
235	Environmental memory of polymer networks under stress. *Advanced Materials*, **2014**, 26, 3441-4	24	33
234	Density of Mixtures Containing Sugars and Ionic Liquids: Experimental Data and PC-SAFT Modeling. *Journal of Chemical & Engineering Data*, **2014**, 59, 2942-2954	2.8	33
233	Thermodynamic properties of aqueous salt containing urea solutions. *Fluid Phase Equilibria*, **2012**, 325, 71-79	2.5	32
232	Toward Thermodynamic Predictions of Aqueous Vitamin Solubility: An Activity Coefficient-Based Approach. *Industrial & Engineering Chemistry Research*, **2019**, 58, 7362-7369	3.9	31
231	Thermodynamic Modeling for Efficient Cocrystal Formation. *Crystal Growth and Design*, **2015**, 15, 4406-4416	3.9	31
230	Co-solvent effects on reaction rate and reaction equilibrium of an enzymatic peptide hydrolysis. *Physical Chemistry Chemical Physics*, **2018**, 20, 11317-11326	3.6	31
229	Oil desulfurization using deep eutectic solvents as sustainable and economical extractants via liquid-liquid extraction: Experimental and PC-SAFT predictions. *Fluid Phase Equilibria*, **2018**, 467, 33-44	2.5	31
228	Moisture-induced phase separation and recrystallization in amorphous solid dispersions. *International Journal of Pharmaceutics*, **2017**, 532, 635-646	6.5	30
227	Modelling of high-pressure phase equilibria using the Sako□Wu□Prausnitz equation of state. *Fluid Phase Equilibria*, **1999**, 163, 79-98	2.5	30
226	Phase-equilibrium measurement and modeling of the PMMA/MMA/carbon dioxide ternary system. *Journal of Supercritical Fluids*, **2008**, 46, 218-225	4.2	29
225	Thermodynamics of the alanine aminotransferase reaction. *Fluid Phase Equilibria*, **2016**, 422, 87-98	2.5	29
224	A thermodynamic investigation of the glucose-6-phosphate isomerization. *Biophysical Chemistry*, **2014**, 195, 22-31	3.5	28
223	Influence of electrolytes on liquid-liquid equilibria of water/1-butanol and on the partitioning of 5-hydroxymethylfurfural in water/1-butanol. *Fluid Phase Equilibria*, **2016**, 428, 102-111	2.5	28
222	Modelling of organic-solvent flux through a polyimide membrane. *Journal of Membrane Science*, **2013**, 428, 554-561	9.6	27
221	The role of activity coefficients in bioreaction equilibria: thermodynamics of methyl ferulate hydrolysis. *Biophysical Chemistry*, **2013**, 173-174, 21-30	3.5	27
220	Design of hybrid distillation/melt crystallisation processes for separation of close boiling mixtures. *Chemical Engineering and Processing: Process Intensification*, **2013**, 67, 16-24	3.7	27
219	Modeling the Phase Behavior of PEO□PPO□PEO Surfactants in Carbon Dioxide Using the PC-SAFT Equation of State: Application to Dry Decontamination of Solid Substrates□*Journal of Chemical & Engineering Data*, **2009**, 54, 1551-1559	2.8	27

218	Polymorphs, Hydrates, Cocrystals, and Cocrystal Hydrates: Thermodynamic Modeling of Theophylline Systems. *Crystal Growth and Design*, **2016**, 16, 4439-4449	3.5	26
217	Investigating phase separation in amorphous solid dispersions via Raman mapping. *International Journal of Pharmaceutics*, **2018**, 535, 245-252	6.5	25
216	Modeling binary mixtures of n-alkanes and water using PC-SAFT. *Fluid Phase Equilibria*, **2018**, 470, 203-211	2.5	25
215	High-pressure gas solubility in multicomponent solvent systems for hydroformylation. Part I: Carbon monoxide solubility. *Journal of Supercritical Fluids*, **2013**, 81, 23-32	4.2	24
214	Modeling aqueous two-phase systems: III. Polymers and organic salts as ATPS former. *Fluid Phase Equilibria*, **2015**, 387, 178-189	2.5	24
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212	Solubility calculation of pharmaceutical compounds—A priori parameter estimation using quantum-chemistry. *Fluid Phase Equilibria*, **2010**, 299, 161-170	2.5	23
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209	Solubility, crystallization and oiling-out behavior of PEGDME: 1. Pure-solvent systems. *Fluid Phase Equilibria*, **2010**, 298, 253-261	2.5	22
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207	Modeling of Solid/Fluid Phase Equilibria in Multicomponent Systems at High Pressure. *Chemical Engineering and Technology*, **2001**, 24, 607-612	2	22
206	A Novel Approach for Analyzing the Dissolution Mechanism of Solid Dispersions. *Pharmaceutical Research*, **2015**, 32, 2559-78	4.5	21
205	Thermodynamics of a model biological reaction: A comprehensive combined experimental and theoretical study. *Fluid Phase Equilibria*, **2016**, 422, 99-110	2.5	21
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203	Solvent effects on esterification equilibria. *AICHE Journal*, **2015**, 61, 3000-3011	3.6	21
202	PC-SAFT parameters from ab initio calculations. *Fluid Phase Equilibria*, **2014**, 362, 41-50	2.5	21
201	Calculation of complex phase equilibria of DMF/alkane systems using the PCP-SAFT equation of state. *Chemical Engineering Science*, **2014**, 115, 49-57	4.4	21

200	Efficient phase separation and product recovery in organic-aqueous bioprocessing using supercritical carbon dioxide. *Biotechnology and Bioengineering*, **2010**, 107, 642-51	4.9	21
199	Influence of pH Value and Ionic Liquids on the Solubility of l-Alanine and l-Glutamic Acid in Aqueous Solutions at 30 °C. *Journal of Chemical & Engineering Data*, **2017**, 62, 52-61	2.8	20
198	Thermodynamics of enzyme-catalyzed esterifications: II. Levulinic acid esterification with short-chain alcohols. *Applied Microbiology and Biotechnology*, **2017**, 101, 7509-7521	5.7	20
197	Influence of Low-Molecular-Weight Excipients on the Phase Behavior of PVPVA64 Amorphous Solid Dispersions. *Pharmaceutical Research*, **2018**, 35, 25	4.5	20
196	High-Pressure Phase Behavior of the System PCHC□CHO□CO2 for the Development of a Solvent-Free Alternative toward Polycarbonate Production. *Industrial & Engineering Chemistry Research*, **2005**, 44, 3363-3366	3.9	20
195	Standard Gibbs energy of metabolic reactions: II. Glucose-6-phosphatase reaction and ATP hydrolysis. *Biophysical Chemistry*, **2017**, 223, 30-38	3.5	19
194	Osmotic coefficients of aqueous weak electrolyte solutions: influence of dissociation on data reduction and modeling. *Journal of Physical Chemistry B*, **2012**, 116, 7479-91	3.4	19
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191	High-pressure gas solubility in multicomponent solvent systems for hydroformylation. Part II: Syngas solubility. *Journal of Supercritical Fluids*, **2014**, 88, 74-84	4.2	18
190	Using complex layer melt crystallization models for the optimization of hybrid distillation/melt crystallization processes. *Chemical Engineering and Processing: Process Intensification*, **2014**, 85, 10-23	3.7	18
189	Modeling Growth Rates in Static Layer Melt Crystallization. *Crystal Growth and Design*, **2013**, 13, 5229-5240	3.9	18
188	Liquid□liquid Equilibria of Systems with Linear Aldehydes. Experimental Data and Modeling with PCP-SAFT. *Industrial & Engineering Chemistry Research*, **2012**, 51, 14525-14534	3.9	18
187	Reducing the amount of PCP□SAFT fitting parameters. 2. Associating components. *Fluid Phase Equilibria*, **2012**, 326, 31-44	2.5	18
186	Activity Coefficients of Complex Molecules by Molecular Simulation and Gibbs-Duhem Integration. *Soft Materials*, **2012**, 10, 26-41	1.7	18
185	Predicting the Solubility of Pharmaceutical Cocrystals in Solvent/Anti-Solvent Mixtures. *Molecules*, **2016**, 21,	4.8	18
184	Solubility and Caloric Properties of Cinnarizine. *Journal of Chemical & Engineering Data*, **2015**, 60, 2256-2261	2.8	17
183	Mutual Impact of Phase Separation/Crystallization and Water Sorption in Amorphous Solid Dispersions. *Molecular Pharmaceutics*, **2018**, 15, 669-678	5.6	17

182	Prediction and Experimental Validation of Co-Solvent Influence on Michaelis Constants: A Thermodynamic Activity-Based Approach. *Chemistry - A European Journal*, **2018**, 24, 16418-16425	4.8	17
181	Application of the PC-SAFT equation of state to modeling of solid-liquid equilibria in systems with organic components forming chemical compounds. *Russian Journal of Applied Chemistry*, **2007**, 80, 542-548	0.8	17
180	Standard Gibbs Energy of Metabolic Reactions: I. Hexokinase Reaction. *Biochemistry*, **2016**, 55, 5665-5674	3.2	17
179	Partition Coefficients of Pharmaceuticals as Functions of Temperature and pH. *Industrial & Engineering Chemistry Research*, **2015**, 54, 3968-3975	3.9	16
178	Extension of the PC-SAFT based group contribution method for polymers to aromatic, oxygen- and silicon-based polymers. *Fluid Phase Equilibria*, **2013**, 339, 89-104	2.5	16
177	Predicting the high concentration co-solvent influence on the reaction equilibria of the ADH-catalyzed reduction of acetophenone. *Journal of Chemical Thermodynamics*, **2019**, 128, 275-282	2.9	16
176	Solvent influence on the phase behavior and glass transition of Amorphous Solid Dispersions. *European Journal of Pharmaceutics and Biopharmaceutics*, **2021**, 158, 132-142	5.7	16
175	Thermodynamics of enzyme-catalyzed esterifications: I. Succinic acid esterification with ethanol. *Applied Microbiology and Biotechnology*, **2017**, 101, 5973-5984	5.7	15
174	Reaction Equilibrium of the Transamination of (S)-Phenylethylamine: Experiments and ePC-SAFT Modeling. *Organic Process Research and Development*, **2017**, 21, 976-986	3.9	15
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169	Selecting Excipients Forming Therapeutic Deep Eutectic Systems-A Mechanistic Approach. *Molecular Pharmaceutics*, **2019**, 16, 3091-3099	5.6	14
168	Phase behavior of pharmaceutically relevant polymer/solvent mixtures. *International Journal of Pharmaceutics*, **2020**, 577, 119065	6.5	14
167	Minimal Experimental Data Set Required for Estimating PCP-SAFT Parameters. *Industrial & Engineering Chemistry Research*, **2011**, 50, 11746-11754	3.9	14
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165	Correctly Measuring and Predicting Solubilities of Solvates, Hydrates, and Polymorphs. *Crystal Growth and Design*, **2020**, 20, 723-735	3.5	14

164	Thermodynamic Approach for Co-crystal Screening. *Crystal Growth and Design*, **2019**, 19, 3253-3264	3.5	13
163	The dynamic influence of cells on the formation of stable emulsions in organic-aqueous biotransformations. *Journal of Industrial Microbiology and Biotechnology*, **2015**, 42, 1011-26	4.2	13
162	Modeling and prediction of protein solubility using the second osmotic virial coefficient. *Fluid Phase Equilibria*, **2016**, 422, 32-42	2.5	13
161	Measurement and Modeling of Phase Equilibria in Systems of Acetonitrile,n-Alkanes, and Myrcene. *Industrial & Engineering Chemistry Research*, **2015**, 54, 1153-1160	3.9	13
160	Modelling of high-pressure phase equilibria using the Sako Wu Prausnitz equation of state: I. Pure-components and heavy n-alkane solutions. *Fluid Phase Equilibria*, **1999**, 163, 61-77	2.5	13
159	Modeling the CO2 Solubility in Aqueous Electrolyte Solutions Using ePC-SAFT. *Journal of Chemical & Engineering Data*, **2020**, 65, 5768-5777	2.8	13
158	Phase Equilibria for the Hydroesterification of 10-Undecenoic Acid Methyl Ester. *Journal of Chemical & Engineering Data*, **2016**, 61, 3317-3325	2.8	13
157	Combined co-solvent and pressure effect on kinetics of a peptide hydrolysis: an activity-based approach. *Physical Chemistry Chemical Physics*, **2019**, 21, 22224-22229	3.6	13
156	Choosing Appropriate Solvents for ASD Preparation. *Molecular Pharmaceutics*, **2018**, 15, 5397-5409	5.6	13
155	Thermodynamic Modeling of Solvent-Impact on Phase Separation in Amorphous Solid Dispersions during Drying. *Molecular Pharmaceutics*, **2020**, 17, 2721-2733	5.6	12
154	Thermodynamic Properties of Aqueous Glucose Urea Salt Systems. *Journal of Solution Chemistry*, **2014**, 43, 1110-1131	1.8	12
153	Phase behavior of poly(dimethylsiloxane) poly(ethylene oxide) amphiphilic block and graft copolymers in compressed carbon dioxide. *Journal of Supercritical Fluids*, **2012**, 62, 211-218	4.2	12
152	Predicting the Solubility of CO2 in Toluene + Ionic Liquid Mixtures with PC-SAFT. *Industrial & Engineering Chemistry Research*, **2017**, 56, 9885-9894	3.9	12
151	VOC sorption in glassy polyimides Measurements and modeling. *Journal of Membrane Science*, **2012**, 415-416, 596-607	9.6	12
150	Reducing the amount of PCP-SAFT fitting parameters. 1. Non-polar and dipolar components. *Fluid Phase Equilibria*, **2012**, 326, 21-30	2.5	12
149	pePC-SAFT: Modeling of polyelectrolyte systems. *Fluid Phase Equilibria*, **2010**, 299, 84-93	2.5	12
148	Innovative Verfahren mit komprimierten Gasen. *Chemie-Ingenieur-Technik*, **1997**, 69, 298-311	0.8	12
147	Drug Release Kinetics and Mechanism from PLGA Formulations. *AICHE Journal*, **2016**, 62, 4055-4065	3.6	12

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145	Measuring and Predicting Thermodynamic Limitation of an Alcohol Dehydrogenase Reaction. *Industrial & Engineering Chemistry Research*, **2017**, 56, 5535-5546	3.9	11
144	Predicting the Effect of pH on Stability and Solubility of Polymorphs, Hydrates, and Cocrystals. *Crystal Growth and Design*, **2016**, 16, 4136-4147	3.5	11
143	Toluene Sorption in Poly(styrene) and Poly(vinyl acetate) near the Glass Transition. *Industrial & Engineering Chemistry Research*, **2013**, 52, 8917-8927	3.9	11
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140	Homogeneously catalyzed hydroamination in a Taylor□couette reactor using a thermormorphic multicomponent solvent system. *Chemical Engineering Research and Design*, **2016**, 112, 263-273	5.5	11
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