# Gabriele Sadowski

## List of Publications by Year in Descending Order



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|---|---|---|---|
| 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 | Predicting Solvent Effects on Homogeneity and Kinetics of the Hydroaminomethylation: A Thermodynamic Approach Using PC-SAFT. *Industrial & Engineering Chemistry Research*, **2022**, 61, 2323-2332 | 3.9 | 0 |
| 306 | Factors Influencing the Crystallization-Onset Time of Metastable ASDs.. *Pharmaceutics*, **2022**, 14, | 6.4 | 2 |
| 305 | Measuring and Modeling Water Sorption in Amorphous Indomethacin and Ritonavir.. *Molecular Pharmaceutics*, **2022**, | 5.6 | 2 |
| 304 | Influence of Temperature and Concentration on the Self-Assembly of Nonionic CE Surfactants: A Light Scattering Study.. *ACS Omega*, **2022**, 7, 7057-7065 | 3.9 | 2 |
| 303 | Influence of Process Temperature and Residence Time on the Manufacturing of Amorphous Solid Dispersions in Hot Melt Extrusion.. *Pharmaceutical Development and Technology*, **2022**, 1-9 | 3.4 | 1 |
| 302 | Water Sorption in Glassy Polyvinylpyrrolidone-Based Polymers.. *Membranes*, **2022**, 12, | 3.8 | 2 |
| 301 | Unraveling the influence of dissolved gases on permeate flux in organic solvent nanofiltration  experimental analysis. *Separation and Purification Technology*, **2022**, 121265 | 8.3 | 0 |
| 300 | Theoretical modeling and prediction of biorelevant solubility of poorly soluble pharmaceuticals. *Chemical Engineering Journal*, **2022**, 444, 136678 | 14.7 | 0 |
| 299 | Predicting the Water Sorption in ASDs. *Pharmaceutics*, **2022**, 14, 1181 | 6.4 | 1 |
| 298 | Generalized Diffusion Relaxation Model for Solvent Sorption in Polymers. *Industrial & Engineering Chemistry Research*, **2021**, 60, 15766-15781 | 3.9 | 8 |
| 297 | Impact of deep eutectic solvents and their constituents on the aqueous solubility of phloroglucinol dihydrate. *Journal of Molecular Liquids*, **2021**, 344, 117932 | 6 | 2 |
| 296 | Solvent Selection in Homogeneous Catalysis Optimization of Kinetics and Reaction Performance. *ACS Catalysis*, **2021**, 11, 590-594 | 13.1 | 6 |
| 295 | Stability of Pharmaceutical Co-Crystals at Humid Conditions Can Be Predicted. *Pharmaceutics*, **2021**, 13, | 6.4 | 5 |
| 294 | Predicting Vapor Liquid Equilibria for Sour-Gas Absorption in Aqueous Mixtures of Chemical and Physical Solvents or Ionic Liquids with ePC-SAFT. *Industrial & Engineering Chemistry Research*, **2021**, 60, 6327-6336 | 3.9 | 6 |
| 293 | Predicting Deliquescence Relative Humidities of Crystals and Crystal Mixtures. *Molecules*, **2021**, 26, | 4.8 | 2 |
| 292 | Swelling and Diffusion in Polymerized Ionic Liquids-Based Hydrogels. *Polymers*, **2021**, 13, | 4.5 | 3 |
| 291 | Thermodynamic Properties of Biogenic Amines and Their Solutions. *Journal of Chemical & Engineering Data*, **2021**, 66, 2822-2831 | 2.8 | 3 |

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| 290 | Co-Crystal Screening by Vapor Sorption of Organic Solvents. *Crystal Growth and Design*, **2021**, 21, 4445-4455 | 4.5 | 1 |
| 289 | Solubility of DNP-amino acids and their partitioning in biodegradable ATPS: Experimental and ePC-SAFT modeling. *Fluid Phase Equilibria*, **2021**, 527, 112830 | 2.5 | 4 |
| 288 | 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 |
| 287 | Combining crystalline and polymeric excipients in API solid dispersions - Opportunity or risk?. *European Journal of Pharmaceutics and Biopharmaceutics*, **2021**, 158, 323-335 | 5.7 | 4 |
| 286 | Measurement and PC-SAFT Modeling of the Solubility of Gallic Acid in Aqueous Mixtures of Deep Eutectic Solvents. *Journal of Chemical & Engineering Data*, **2021**, 66, 958-967 | 2.8 | 8 |
| 285 | PC-SAFT Modeling of Phase Equilibria Relevant for Lipid-Based Drug Delivery Systems. *Journal of Chemical & Engineering Data*, **2021**, 66, 1280-1289 | 2.8 | 2 |
| 284 | Predicting the API partitioning between lipid-based drug delivery systems and water. *International Journal of Pharmaceutics*, **2021**, 595, 120266 | 6.5 | 0 |
| 283 | Insights into influence mechanism of polymeric excipients on dissolution of drug formulations: A molecular interaction-based theoretical model analysis and prediction. *AICHE Journal*, **2021**, 67, e17372 | 3.6 | 0 |
| 282 | Boosting the kinetic efficiency of formate dehydrogenase by combining the effects of temperature, high pressure and co-solvent mixtures. *Colloids and Surfaces B: Biointerfaces*, **2021**, 208, 112127 | 6 | 4 |
| 281 | Solvent mixtures in pharmaceutical development: Maximizing the API solubility and avoiding phase separation. *Fluid Phase Equilibria*, **2021**, 548, 113200 | 2.5 | 0 |
| 280 | Predicting process design spaces for spray drying amorphous solid dispersions. *International Journal of Pharmaceutics: X*, **2021**, 3, 100072 | 3.2 | 2 |
| 279 | Phase behavior of ASDs based on hydroxypropyl cellulose. *International Journal of Pharmaceutics: X* , **2021**, 3, 100070 | 3.2 | 6 |
| 278 | Methodology Based on the Theory of Information to Describe the Representation Ability of the DMC + Alkane Behavior. *Industrial & Engineering Chemistry Research*, **2021**, 60, 1036-1054 | 3.9 | 2 |
| 277 | In-Silico Screening of Lipid-Based Drug Delivery Systems. *Pharmaceutical Research*, **2020**, 37, 249 | 4.5 | 2 |
| 276 | Solubility of Pharmaceutical Ingredients in Natural Edible Oils. *Molecular Pharmaceutics*, **2020**, 17, 2499-2507 | 5.6 | 9 |
| 275 | Thermodynamic Modeling of Solvent-Impact on Phase Separation in Amorphous Solid Dispersions during Drying. *Molecular Pharmaceutics*, **2020**, 17, 2721-2733 | 5.6 | 12 |
| 274 | High-Pressure-Mediated Thiourea-Organocatalyzed Asymmetric Michael Addition to (Hetero)aromatic Nitroolefins: Prediction of Reaction Parameters by PCP-SAFT Modelling. *ChemPlusChem*, **2020**, 85, 1292-1296 | 2.8 | 2 |
| 273 | Hydrate formation in polymer-based pharmaceutical formulations. *Fluid Phase Equilibria*, **2020**, 521, 112677 | 2.5 | 2 |

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| 272 | Influence of cytosolic conditions on the reaction equilibrium and the reaction enthalpy of the enolase reaction accessed by calorimetry and van 't HOFF. *Biochimica Et Biophysica Acta - General Subjects*, **2020**, 1864, 129675 | 4 | 4 |
| 271 | Standard Gibbs energy of metabolic reactions: IV. Triosephosphate isomerase reaction. *Biophysical Chemistry*, **2020**, 258, 106330 | 3.5 | 5 |
| 270 | Standard Gibbs energy of metabolic reactions: V. Enolase reaction. *Biochimica Et Biophysica Acta - Proteins and Proteomics*, **2020**, 1868, 140365 | 4 | 7 |
| 269 | Phase behavior of pharmaceutically relevant polymer/solvent mixtures. *International Journal of Pharmaceutics*, **2020**, 577, 119065 | 6.5 | 14 |
| 268 | Standard Gibbs energy of metabolic reactions: VI. Glyceraldehyde 3-phosphate dehydrogenase reaction. *Fluid Phase Equilibria*, **2020**, 517, 112597 | 2.5 | 3 |
| 267 | The influence of polymeric excipients on desupersaturation profiles of active pharmaceutical ingredients. 1: Polyethylene glycol. *International Journal of Pharmaceutics*, **2020**, 582, 119317 | 6.5 | 4 |
| 266 | The role of molecular interactions on Michaelis constants of α-chymotrypsin catalyzed peptide hydrolyses. *Journal of Chemical Thermodynamics*, **2020**, 148, 106142 | 2.9 | 4 |
| 265 | Correctly Measuring and Predicting Solubilities of Solvates, Hydrates, and Polymorphs. *Crystal Growth and Design*, **2020**, 20, 723-735 | 3.5 | 14 |
| 264 | Modeling the CO2 Solubility in Aqueous Electrolyte Solutions Using ePC-SAFT. *Journal of Chemical & Engineering Data*, **2020**, 65, 5768-5777 | 2.8 | 13 |
| 263 | Application of PC-SAFT and DGT for the Prediction of Self-Assembly. *Journal of Chemical & Engineering Data*, **2020**, 65, 5897-5908 | 2.8 | 2 |
| 262 | Partitioning of water-soluble vitamins in biodegradable aqueous two-phase systems: Electrolyte perturbed-chain statistical associating fluid theory predictions and experimental validation. *AICHE Journal*, **2020**, 66, e16984 | 3.6 | 3 |
| 261 | Viscosity of ASDs at humid conditions. *European Journal of Pharmaceutics and Biopharmaceutics*, **2020**, 154, 387-396 | 5.7 | 2 |
| 260 | The interplay of dissolution, solution crystallization and solid-state transformation of amorphous indomethacin in aqueous solution. *International Journal of Pharmaceutics: X*, **2020**, 2, 100063 | 3.2 | 1 |
| 259 | Modeling of Interfacial Tensions of Long-Chain Molecules and Related Mixtures Using Perturbed Chain-Statistical Associating Fluid Theory and the Density Gradient Theory. *Journal of Chemical & Engineering Data*, **2020**, 65, 1005-1018 | 2.8 | 5 |
| 258 | Protein-protein interactions and water activity coefficients can be used to aid a first excipient choice in protein formulations. *International Journal of Pharmaceutics*, **2019**, 569, 118608 | 6.5 | 4 |
| 257 | Structure and thermodynamics of aqueous urea solutions from ambient to kilobar pressures: From thermodynamic modeling, experiments, and first principles simulations to an accurate force field description. *Biophysical Chemistry*, **2019**, 254, 106260 | 3.5 | 7 |
| 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 | Cosolvent and pressure effects on enzyme-catalysed hydrolysis reactions. *Biophysical Chemistry*, **2019**, 252, 106209 | 3.5 | 7 |

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| 254 | Thermodynamic Activity-Based Solvent Design for Bioreactions. *Trends in Biotechnology*, **2019**, 37, 1038-1041 | 15.1 | 9 |
| 253 | Selecting Excipients Forming Therapeutic Deep Eutectic Systems-A Mechanistic Approach. *Molecular Pharmaceutics*, **2019**, 16, 3091-3099 | 5.6 | 14 |
| 252 | Measurement and Prediction of Excess Properties of Binary Mixtures Methyl Decanoate + an Even-Numbered n-Alkane (C6 C16) at 298.15 K. *Journal of Chemical & Engineering Data*, **2019**, 64, 2816-2825 | 2.8 | 2 |
| 251 | Thermodynamic Approach for Co-crystal Screening. *Crystal Growth and Design*, **2019**, 19, 3253-3264 | 3.5 | 13 |
| 250 | Heterosegmental Modeling of Long-Chain Molecules and Related Mixtures Using PC-SAFT: 2. Associating Compounds. *Industrial & Engineering Chemistry Research*, **2019**, 58, 4625-4643 | 3.9 | 6 |
| 249 | Toward Thermodynamic Predictions of Aqueous Vitamin Solubility: An Activity Coefficient-Based Approach. *Industrial & Engineering Chemistry Research*, **2019**, 58, 7362-7369 | 3.9 | 31 |
| 248 | Second osmotic virial coefficients of therapeutic proteins in the presence of excipient-mixtures can be predicted to aid an efficient formulation design. *Journal of Molecular Liquids*, **2019**, 283, 575-583 | 6 | 5 |
| 247 | Thermodynamic Modeling of Triglycerides using PC-SAFT. *Journal of Chemical & Engineering Data*, **2019**, 64, 1446-1453 | 2.8 | 10 |
| 246 | Reply to Comment on Perturbed-Chain SAFT: An Equation of State Based on a Perturbation Theory for Chain Molecules *Industrial & Engineering Chemistry Research*, **2019**, 58, 5744-5745 | 3.9 | 6 |
| 245 | Thermodynamic Properties of Systems Comprising Esters: Experimental Data and Modeling with PC-SAFT and SAFT- Mie. *Industrial & Engineering Chemistry Research*, **2019**, 58, 6841-6849 | 3.9 | 5 |
| 244 | Simultaneous Prediction of Cosolvent Influence on Reaction Equilibrium and Michaelis Constants of Enzyme-Catalyzed Ketone Reductions. *ACS Omega*, **2019**, 4, 6264-6272 | 3.9 | 5 |
| 243 | Density variations of TMAO solutions in the kilobar range: Experiments, PC-SAFT predictions, and molecular dynamics simulations. *Biophysical Chemistry*, **2019**, 253, 106222 | 3.5 | 7 |
| 242 | Thermodynamic properties of aqueous osmolyte solutions at high-pressure conditions. *Biophysical Chemistry*, **2019**, 253, 106211 | 3.5 | 4 |
| 241 | Highlighting 10 Years of NIST Cooperation and Service to the Thermophysical Properties Data Community. *Journal of Chemical & Engineering Data*, **2019**, 64, 4191-4192 | 2.8 | 2 |
| 240 | Solubility of pharmaceutical ingredients in triglycerides. *European Journal of Pharmaceutics and Biopharmaceutics*, **2019**, 145, 113-120 | 5.7 | 10 |
| 239 | Solubility Enhancement of Vitamins in Water in the Presence of Covitamins: Measurements and ePC-SAFT Predictions. *Industrial & Engineering Chemistry Research*, **2019**, 58, 21761-21771 | 3.9 | 10 |
| 238 | Measurement and Modeling of Lactose Solubility in Aqueous Electrolyte Solutions. *Industrial & Engineering Chemistry Research*, **2019**, 58, 20797-20805 | 3.9 | 7 |
| 237 | Thermodynamic Activity-Based Michaelis Constants **2019**, | | 2 |

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| 236 | 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 |
| 235 | Liquid□liquid Equilibria for Separation of Alcohols from Esters Using Deep Eutectic Solvents Based on Choline Chloride: Experimental Study and Thermodynamic Modeling. *Journal of Chemical & Engineering Data*, **2019**, 64, 6049-6059 | 2.8 | 9 |
| 234 | Heterosegmental Modeling of Long-Chain Molecules and Related Mixtures using PC-SAFT: 1. Polar Compounds. *Industrial & Engineering Chemistry Research*, **2019**, 58, 2551-2574 | 3.9 | 9 |
| 233 | 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 |
| 232 | In-situ determination of crystallization kinetics in ASDs via water sorption experiments. *European Journal of Pharmaceutics and Biopharmaceutics*, **2018**, 127, 183-193 | 5.7 | 10 |
| 231 | 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 |
| 230 | Mutual Influence of Furfural and Furancarboxylic Acids on Their Solubility in Aqueous Solutions: Experiments and Perturbed-Chain Statistical Associating Fluid Theory (PC-SAFT) Predictions. *Journal of Chemical & Engineering Data*, **2018**, 63, 1460-1470 | 2.8 | 10 |
| 229 | Standard Gibbs Energy of Metabolic Reactions: III The 3-Phosphoglycerate Kinase Reaction. *ACS Omega*, **2018**, 3, 1783-1790 | 3.9 | 10 |
| 228 | Physical stability of API/polymer-blend amorphous solid dispersions. *European Journal of Pharmaceutics and Biopharmaceutics*, **2018**, 124, 147-157 | 5.7 | 39 |
| 227 | Influence of Low-Molecular-Weight Excipients on the Phase Behavior of PVPVA64 Amorphous Solid Dispersions. *Pharmaceutical Research*, **2018**, 35, 25 | 4.5 | 20 |
| 226 | Mutual Impact of Phase Separation/Crystallization and Water Sorption in Amorphous Solid Dispersions. *Molecular Pharmaceutics*, **2018**, 15, 669-678 | 5.6 | 17 |
| 225 | 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 |
| 224 | Effect of different organic salts on amino acids partition behaviour in PEG-salt ATPS. *Fluid Phase Equilibria*, **2018**, 456, 84-91 | 2.5 | 10 |
| 223 | 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 |
| 222 | 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 |
| 221 | Thermodynamic prediction of the solvent effect on a transesterification reaction. *Chemical Engineering Science*, **2018**, 176, 264-269 | 4.4 | 8 |
| 220 | Investigating phase separation in amorphous solid dispersions via Raman mapping. *International Journal of Pharmaceutics*, **2018**, 535, 245-252 | 6.5 | 25 |
| 219 | Modeling binary mixtures of n-alkanes and water using PC-SAFT. *Fluid Phase Equilibria*, **2018**, 470, 203-211 | 2.5 | 25 |

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| 218 | Aktivitätskoeffizienten und osmotische Virialkoeffizienten zur Auswahl von Additiven für Antikörper-Formulierungen. *Chemie-Ingenieur-Technik*, **2018**, 90, 1243-1243 | 0.8 | |
| 217 | Choosing Appropriate Solvents for ASD Preparation. *Molecular Pharmaceutics*, **2018**, 15, 5397-5409 | 5.6 | 13 |
| 216 | Phase separation of Pickering-type emulsions – Influence of particles on the interfacial properties of oil/water systems. *Chemie-Ingenieur-Technik*, **2018**, 90, 1321-1322 | 0.8 | |
| 215 | Experimentelle und theoretische Untersuchung der Grenzflächeneigenschaften von wässrigen Systemen. *Chemie-Ingenieur-Technik*, **2018**, 90, 1324-1324 | 0.8 | |
| 214 | 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 |
| 213 | Thermodynamics of enzyme-catalyzed esterifications: I. Succinic acid esterification with ethanol. *Applied Microbiology and Biotechnology*, **2017**, 101, 5973-5984 | 5.7 | 15 |
| 212 | Predicting solvent effects on the 1-dodecene hydroformylation reaction equilibrium. *AICHE Journal* , **2017**, 63, 4576-4585 | 3.6 | 11 |
| 211 | Measuring and Predicting Thermodynamic Limitation of an Alcohol Dehydrogenase Reaction. *Industrial & Engineering Chemistry Research*, **2017**, 56, 5535-5546 | 3.9 | 11 |
| 210 | 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 |
| 209 | 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 |
| 208 | Predicting the Solvent Effect on Esterification Kinetics. *ChemPhysChem*, **2017**, 18, 1977-1980 | 3.2 | 9 |
| 207 | PC-SAFT modeling of CO2 solubilities in hydrophobic deep eutectic solvents. *Fluid Phase Equilibria*, **2017**, 448, 94-98 | 2.5 | 50 |
| 206 | Modeling and analysis of dissolution of paracetamol/Eudragit formulations. *Chemical Engineering Research and Design*, **2017**, 121, 22-31 | 5.5 | 7 |
| 205 | Standard Gibbs energy of metabolic reactions: II. Glucose-6-phosphatase reaction and ATP hydrolysis. *Biophysical Chemistry*, **2017**, 223, 30-38 | 3.5 | 19 |
| 204 | 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 |
| 203 | Thermodynamics of enzyme-catalyzed esterifications: II. Levulinic acid esterification with short-chain alcohols. *Applied Microbiology and Biotechnology*, **2017**, 101, 7509-7521 | 5.7 | 20 |
| 202 | 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 |
| 201 | Moisture-induced phase separation and recrystallization in amorphous solid dispersions. *International Journal of Pharmaceutics*, **2017**, 532, 635-646 | 6.5 | 30 |

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| 200 | Predicting the Solubility of CO2 in Toluene + Ionic Liquid Mixtures with PC-SAFT. *Industrial & Engineering Chemistry Research*, **2017**, 56, 9885-9894 | 3.9 | 12 |
| 199 | Crowders and Cosolvents-Major Contributors to the Cellular Milieu and Efficient Means to Counteract Environmental Stresses. *ChemPhysChem*, **2017**, 18, 2951-2972 | 3.2 | 65 |
| 198 | Phase Behavior of Binary Mixtures Containing Succinic Acid or Its Esters. *Journal of Chemical & Engineering Data*, **2017**, 62, 1983-1993 | 2.8 | 9 |
| 197 | Long-Term Physical Stability of PVP- and PVPVA-Amorphous Solid Dispersions. *Molecular Pharmaceutics*, **2017**, 14, 157-171 | 5.6 | 71 |
| 196 | Solubilization of proteins in aqueous two-phase extraction through combinations of phase-formers and displacement agents. *European Journal of Pharmaceutics and Biopharmaceutics*, **2017**, 112, 38-44 | 5.7 | 8 |
| 195 | Amorphous-Amorphous Phase Separation in API/Polymer Formulations. *Molecules*, **2017**, 22, | 4.8 | 45 |
| 194 | Compatible solutes: Thermodynamic properties relevant for effective protection against osmotic stress. *Fluid Phase Equilibria*, **2016**, 407, 224-235 | 2.5 | 37 |
| 193 | Protein partition coefficients can be estimated efficiently by hybrid shortcut calculations. *Journal of Biotechnology*, **2016**, 233, 151-9 | 3.7 | 5 |
| 192 | Applied catastrophic phase inversion: a continuous non-centrifugal phase separation step in biphasic whole-cell biocatalysis. *Journal of Industrial Microbiology and Biotechnology*, **2016**, 43, 1527-1535 | 4.2 | 9 |
| 191 | Applied Catastrophic Phase Inversion (ACPI) □A Continuous Noncentrifugal Phase Separation in Biphasic Whole-Cell Biocatalysis. *Chemie-Ingenieur-Technik*, **2016**, 88, 1331-1332 | 0.8 | |
| 190 | Thermodynamics of Bioreactions. *Annual Review of Chemical and Biomolecular Engineering*, **2016**, 7, 395-414 | 8.3 | 36 |
| 189 | Recovery of cis,cis-muconic acid from organic phase after reactive extraction. *Separation and Purification Technology*, **2016**, 169, 1-8 | 8.3 | 8 |
| 188 | Polymorphs, Hydrates, Cocrystals, and Cocrystal Hydrates: Thermodynamic Modeling of Theophylline Systems. *Crystal Growth and Design*, **2016**, 16, 4439-4449 | 3.5 | 26 |
| 187 | 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 |
| 186 | VOC Sorption in Stretched Cross-Linked Natural Rubber. *Industrial & Engineering Chemistry Research*, **2016**, 55, 7191-7200 | 3.9 | 4 |
| 185 | Non-monotonic course of protein solubility in aqueous polymer-salt solutions can be modeled using the sol-mxDLVO model. *Biotechnology Journal*, **2016**, 11, 282-9 | 5.6 | |
| 184 | Modeling and prediction of protein solubility using the second osmotic virial coefficient. *Fluid Phase Equilibria*, **2016**, 422, 32-42 | 2.5 | 13 |
| 183 | PC-SAFT Modeling of CO2 Solubilities in Deep Eutectic Solvents. *Journal of Physical Chemistry B*, **2016**, 120, 2300-10 | 3.4 | 78 |

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| 182 | Thermodynamics of a model biological reaction: A comprehensive combined experimental and theoretical study. *Fluid Phase Equilibria*, **2016**, 422, 99-110 | 2.5 | 21 |
| 181 | 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 |
| 180 | Predicting the Solubility of Pharmaceutical Cocrystals in Solvent/Anti-Solvent Mixtures. *Molecules*, **2016**, 21, | 4.8 | 18 |
| 179 | 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 |
| 178 | Drug Release Kinetics and Mechanism from PLGA Formulations. *AICHE Journal*, **2016**, 62, 4055-4065 | 3.6 | 12 |
| 177 | Phase Equilibria for the Hydroesterification of 10-Undecenoic Acid Methyl Ester. *Journal of Chemical & Engineering Data*, **2016**, 61, 3317-3325 | 2.8 | 13 |
| 176 | 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 |
| 175 | Influence of Salts on the Partitioning of 5-Hydroxymethylfurfural in Water/MIBK. *Journal of Physical Chemistry B*, **2016**, 120, 3797-808 | 3.4 | 41 |
| 174 | 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 |
| 173 | Thermodynamics of the alanine aminotransferase reaction. *Fluid Phase Equilibria*, **2016**, 422, 87-98 | 2.5 | 29 |
| 172 | Standard Gibbs Energy of Metabolic Reactions: I. Hexokinase Reaction. *Biochemistry*, **2016**, 55, 5665-5674 | 3.2 | 17 |
| 171 | Cation Effect on the Water Activity of Ternary (S)-Aminobutanedioic Acid Magnesium Salt Solutions at 298.15 and 310.15 K. *Journal of Chemical & Engineering Data*, **2016**, 61, 3190-3199 | 2.8 | 8 |
| 170 | Novel Displacement Agents for Aqueous 2-Phase Extraction Can Be Estimated Based on Hybrid Shortcut Calculations. *Journal of Pharmaceutical Sciences*, **2016**, 105, 3030-3038 | 3.9 | 7 |
| 169 | Inclusion of mPRISM potential for polymer-induced protein interactions enables modeling of second osmotic virial coefficients in aqueous polymer-salt solutions. *Biotechnology Journal*, **2016**, 11, 146-54 | 5.6 | 7 |
| 168 | A Novel Approach for Analyzing the Dissolution Mechanism of Solid Dispersions. *Pharmaceutical Research*, **2015**, 32, 2559-78 | 4.5 | 21 |
| 167 | Thermodynamic Modeling for Efficient Cocrystal Formation. *Crystal Growth and Design*, **2015**, 15, 4406-4416 | 3.5 | 31 |
| 166 | Predicting the Solubility Advantage of Amorphous Pharmaceuticals: A Novel Thermodynamic Approach. *Molecular Pharmaceutics*, **2015**, 12, 2823-33 | 5.6 | 48 |
| 165 | Influence of humidity on the phase behavior of API/polymer formulations. *European Journal of Pharmaceutics and Biopharmaceutics*, **2015**, 94, 352-62 | 5.7 | 50 |

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| 164 | Solubility and Caloric Properties of Cinnarizine. *Journal of Chemical & Engineering Data*, **2015**, 60, 2256-2261 | 2.8 | 17 |
| 163 | Solving Phase Equilibrium Problems by Means of Avoidance-Based Multiobjectivization **2015**, 1159-1171 | | 7 |
| 162 | Phase Equilibria in Systems of Morpholine, Acetonitrile, and n-Alkanes. *Journal of Chemical & Engineering Data*, **2015**, 60, 2098-2103 | 2.8 | 5 |
| 161 | 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 |
| 160 | Partition Coefficients of Pharmaceuticals as Functions of Temperature and pH. *Industrial & Engineering Chemistry Research*, **2015**, 54, 3968-3975 | 3.9 | 16 |
| 159 | Thermodynamic phase behaviour of indomethacin/PLGA formulations. *European Journal of Pharmaceutics and Biopharmaceutics*, **2015**, 93, 88-94 | 5.7 | 34 |
| 158 | Different recycling concepts in the homogeneously catalysed synthesis of terpenyl amines. *Chemical Engineering and Processing: Process Intensification*, **2015**, 98, 22-31 | 3.7 | 9 |
| 157 | Solid Dispersion a Pragmatic Method to Improve the Bioavailability of Poorly Soluble Drugs **2015**, 81-100 | | 0 |
| 156 | Solvent effects on esterification equilibria. *AICHE Journal*, **2015**, 61, 3000-3011 | 3.6 | 21 |
| 155 | Process boundaries of irreversible scCO2 -assisted phase separation in biphasic whole-cell biocatalysis. *Biotechnology and Bioengineering*, **2015**, 112, 2316-23 | 4.9 | 6 |
| 154 | Dissolution of Crystalline Pharmaceuticals: Experimental Investigation and Thermodynamic Modeling. *Industrial & Engineering Chemistry Research*, **2015**, 54, 731-742 | 3.9 | 44 |
| 153 | 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 |
| 152 | Modeling aqueous two-phase systems: III. Polymers and organic salts as ATPS former. *Fluid Phase Equilibria*, **2015**, 387, 178-189 | 2.5 | 24 |
| 151 | Modeling aqueous two-phase systems: I. Polyethylene glycol and inorganic salts as ATPS former. *Fluid Phase Equilibria*, **2014**, 368, 91-103 | 2.5 | 41 |
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