

# Christoph Held

## List of Publications by Year in descending order

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113  
papers

4,070  
citations

101543

36  
h-index

133252

59  
g-index

114  
all docs

114  
docs citations

114  
times ranked

2937  
citing authors

#	ARTICLE	IF	CITATIONS
1	Solvent Selection for the Extraction of 2-Phenylethanol from Aqueous Phases: Density and Viscosity Studies. <i>Journal of Chemical &amp; Engineering Data</i> , 2022, 67, 1893-1904.	1.9	3
2	Interfacial Properties of Deep Eutectic Solvents by Density Gradient Theory. <i>Industrial &amp; Engineering Chemistry Research</i> , 2022, 61, 2580-2591.	3.7	6
3	The melting properties of D- $\alpha$ -glucose, D- $\beta$ -fructose, D-sucrose, D- $\alpha$ -galactose, and D- $\alpha$ -xylose and their solubility in water: A revision. <i>Food Biophysics</i> , 2022, 17, 181-197.	3.0	3
4	Calculation of Multiphase Equilibria Containing Mixed Solvents and Mixed Electrolytes: General Formulation and Case Studies. <i>Journal of Chemical &amp; Engineering Data</i> , 2022, 67, 1972-1984.	1.9	10
5	Solubility of Electrolytes in Organic Solvents: Solvent-Specific Effects and Ion-Specific Effects. <i>Journal of Chemical &amp; Engineering Data</i> , 2022, 67, 2706-2718.	1.9	19
6	Prediction of pH in multiphase multicomponent systems with ePC-SAFT advanced. <i>Chemical Communications</i> , 2022, 58, 8436-8439.	4.1	9
7	Perspective: Increasing blue carbon around Antarctica is an ecosystem service of considerable societal and economic value worth protecting. <i>Global Change Biology</i> , 2021, 27, 5-12.	9.5	35
8	Measurement and PC-SAFT Modeling of the Solubility of Gallic Acid in Aqueous Mixtures of Deep Eutectic Solvents. <i>Journal of Chemical &amp; Engineering Data</i> , 2021, 66, 958-967.	1.9	17
9	Unravelling the nature of citric acid: arginine:water mixtures: the bifunctional role of water. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 1706-1717.	2.8	20
10	Melting Properties of Peptides and Their Solubility in Water. Part 2: Di- and Tripeptides Based on Glycine, Alanine, Leucine, Proline, and Serine. <i>Industrial &amp; Engineering Chemistry Research</i> , 2021, 60, 4693-4704.	3.7	13
11	New thermodynamic activity-based approach allows predicting the feasibility of glycolysis. <i>Scientific Reports</i> , 2021, 11, 6125.	3.3	4
12	Predicting Vapor-Liquid Equilibria for Sour-Gas Absorption in Aqueous Mixtures of Chemical and Physical Solvents or Ionic Liquids with ePC-SAFT. <i>Industrial &amp; Engineering Chemistry Research</i> , 2021, 60, 6327-6336.	3.7	14
13	Prediction of salting-out in liquid-liquid two-phase systems with ePC-SAFT: Effect of the Born term and of a concentration-dependent dielectric constant. <i>Zeitschrift Fur Anorganische Und Allgemeine Chemie</i> , 2021, 647, 1305-1314.	1.2	19
14	Boosting the kinetic efficiency of formate dehydrogenase by combining the effects of temperature, high pressure and co-solvent mixtures. <i>Colloids and Surfaces B: Biointerfaces</i> , 2021, 208, 112127.	5.0	17
15	Extremely Low Vapor Pressure Data as Access to PC-SAFT Parameter Estimation for Ionic Liquids and Modeling of Precursor Solubility in Ionic Liquids. <i>ChemistryOpen</i> , 2021, 10, 216-226.	1.9	7
16	Tetrahydrothiophene-Based Ionic Liquids: Synthesis and Thermodynamic Characterizations. <i>ChemistryOpen</i> , 2021, 10, 153-163.	1.9	8
17	Protecting Antarctic blue carbon: as marine ice retreats can the law fill the gap?. <i>Climate Policy</i> , 2020, 20, 149-162.	5.1	14
18	Modeling the CO <sub>2</sub> Solubility in Aqueous Electrolyte Solutions Using ePC-SAFT. <i>Journal of Chemical &amp; Engineering Data</i> , 2020, 65, 5768-5777.	1.9	33

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19	Thermodynamics and Kinetics of Glycolytic Reactions. Part I: Kinetic Modeling Based on Irreversible Thermodynamics and Validation by Calorimetry. <i>International Journal of Molecular Sciences</i> , 2020, 21, 8341.	4.1	4
20	Vapor Pressure Assessment of Sulfolane-Based Eutectic Solvents: Experimental, PC-SAFT, and Molecular Dynamics. <i>Journal of Physical Chemistry B</i> , 2020, 124, 10386-10397.	2.6	12
21	Promising Thiolanium Ionic Liquid for Extraction of Aromatics from Aliphatics: Experiments and Modeling. <i>Industrial &amp; Engineering Chemistry Research</i> , 2020, 59, 15707-15717.	3.7	16
22	Partitioning of water-soluble vitamins in biodegradable aqueous two-phase systems: Electrolyte perturbed chain statistical associating fluid theory predictions and experimental validation. <i>AIChE Journal</i> , 2020, 66, e16984.	3.6	9
23	Thermodynamic $E^{\text{g}}$ Models and Equations of State for Electrolytes in a Water-Poor Medium: A Review. <i>Journal of Chemical &amp; Engineering Data</i> , 2020, 65, 5073-5082.	1.9	21
24	Thermodynamics and Kinetics of Glycolytic Reactions. Part II: Influence of Cytosolic Conditions on Thermodynamic State Variables and Kinetic Parameters. <i>International Journal of Molecular Sciences</i> , 2020, 21, 7921.	4.1	4
25	5-Hydroxymethylfurfural Synthesis in Nonaqueous Two-Phase Systems (NTPS) – PC-SAFT Predictions and Validation. <i>Organic Process Research and Development</i> , 2020, 24, 1052-1062.	2.7	8
26	High-Pressure-Mediated Thiourea-Organocatalyzed Asymmetric Michael Addition to (Hetero)aromatic Nitroolefins: Prediction of Reaction Parameters by PC-SAFT Modelling. <i>ChemPlusChem</i> , 2020, 85, 1292-1296.	2.8	4
27	Influence of cytosolic conditions on the reaction equilibrium and the reaction enthalpy of the enolase reaction accessed by calorimetry and van 't Hoff. <i>Biochimica Et Biophysica Acta - General Subjects</i> , 2020, 1864, 129675.	2.4	5
28	Molecular aspects of lipid metabolism in the midgut gland of the brown shrimp <i>Crangon crangon</i> . <i>Comparative Biochemistry and Physiology - B Biochemistry and Molecular Biology</i> , 2020, 248-249, 110465.	1.6	1
29	Standard Gibbs energy of metabolic reactions: IV. Triosephosphate isomerase reaction. <i>Biophysical Chemistry</i> , 2020, 258, 106330.	2.8	10
30	Standard Gibbs energy of metabolic reactions: V. Enolase reaction. <i>Biochimica Et Biophysica Acta - Proteins and Proteomics</i> , 2020, 1868, 140365.	2.3	10
31	Phylogenetic relationship within Cumacea (Crustacea: Peracarida) and genetic variability of two Antarctic species of the family Leuconidae. <i>Scientia Marina</i> , 2020, 84, 385-392.	0.6	3
32	Ancient globetrotters' connectivity and putative native ranges of two cosmopolitan biofouling amphipods. <i>PeerJ</i> , 2020, 8, e9613.	2.0	14
33	Density variations of TMAO solutions in the kilobar range: Experiments, PC-SAFT predictions, and molecular dynamics simulations. <i>Biophysical Chemistry</i> , 2019, 253, 106222.	2.8	8
34	Thermodynamic properties of aqueous osmolyte solutions at high-pressure conditions. <i>Biophysical Chemistry</i> , 2019, 253, 106211.	2.8	7
35	Catalytic Low-Temperature Dehydration of Fructose to 5-Hydroxymethylfurfural Using Acidic Deep Eutectic Solvents and Polyoxometalate Catalysts. <i>Frontiers in Chemistry</i> , 2019, 7, 661.	3.6	44
36	Solubility Enhancement of Vitamins in Water in the Presence of Covitamins: Measurements and ePC-SAFT Predictions. <i>Industrial &amp; Engineering Chemistry Research</i> , 2019, 58, 21761-21771.	3.7	12

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37	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. <i>Biophysical Chemistry</i> , 2019, 254, 106260.	2.8	10
38	Determination of the Total Vapor Pressure of Hydrophobic Deep Eutectic Solvents: Experiments and Perturbed-Chain Statistical Associating Fluid Theory Modeling. <i>ACS Sustainable Chemistry and Engineering</i> , 2019, 7, 4047-4057.	6.7	69
39	Cosolvent and pressure effects on enzyme-catalysed hydrolysis reactions. <i>Biophysical Chemistry</i> , 2019, 252, 106209.	2.8	10
40	Toward Thermodynamic Predictions of Aqueous Vitamin Solubility: An Activity Coefficient-Based Approach. <i>Industrial &amp; Engineering Chemistry Research</i> , 2019, 58, 7362-7369.	3.7	39
41	Incorporating a concentration-dependent dielectric constant into ePC-SAFT. An application to binary mixtures containing ionic liquids. <i>Fluid Phase Equilibria</i> , 2019, 492, 26-33.	2.5	48
42	Thermophysical properties of glyceline-water mixtures investigated by molecular modelling. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 6467-6476.	2.8	38
43	Thermodynamic Properties of Systems Comprising Esters: Experimental Data and Modeling with PC-SAFT and SAFT- $\Gamma^3$ Mie. <i>Industrial &amp; Engineering Chemistry Research</i> , 2019, 58, 6841-6849.	3.7	9
44	Guanidinium Cation Effect on the Water Activity of Ternary (S)Aminopentanedioic Acid Sodium Salt Solutions at 298.15 and 310.15 K. <i>Journal of Chemical &amp; Engineering Data</i> , 2019, 64, 1256-1264.	1.9	2
45	Modeling Thermodynamic Derivative Properties and Gas Solubility of Ionic Liquids with ePC-SAFT. <i>Industrial &amp; Engineering Chemistry Research</i> , 2019, 58, 8401-8417.	3.7	33
46	Simultaneous Prediction of Cosolvent Influence on Reaction Equilibrium and Michaelis Constants of Enzyme-Catalyzed Ketone Reductions. <i>ACS Omega</i> , 2019, 4, 6264-6272.	3.5	7
47	<i>110th Anniversary:</i> Distribution Coefficients of Furfural and 5-Hydroxymethylfurfural in Hydrophobic Deep Eutectic Solvent + Water Systems: Experiments and Perturbed-Chain Statistical Associating Fluid Theory Predictions. <i>Industrial &amp; Engineering Chemistry Research</i> , 2019, 58, 4240-4247.	3.7	42
48	Aggregation control of Ru and Ir nanoparticles by tunable aryl alkyl imidazolium ionic liquids. <i>Nanoscale</i> , 2019, 11, 4073-4082.	5.6	26
49	Combined co-solvent and pressure effect on kinetics of a peptide hydrolysis: an activity-based approach. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 22224-22229.	2.8	22
50	Liquid-Liquid Equilibria for Separation of Alcohols from Esters Using Deep Eutectic Solvents Based on Choline Chloride: Experimental Study and Thermodynamic Modeling. <i>Journal of Chemical &amp; Engineering Data</i> , 2019, 64, 6049-6059.	1.9	20
51	Predicting the high concentration co-solvent influence on the reaction equilibria of the ADH-catalyzed reduction of acetophenone. <i>Journal of Chemical Thermodynamics</i> , 2019, 128, 275-282.	2.0	20
52	Mutual Influence of Furfural and Furancarboxylic Acids on Their Solubility in Aqueous Solutions: Experiments and Perturbed-Chain Statistical Associating Fluid Theory (PC-SAFT) Predictions. <i>Journal of Chemical &amp; Engineering Data</i> , 2018, 63, 1460-1470.	1.9	12
53	Standard Gibbs Energy of Metabolic Reactions: III The 3-Phosphoglycerate Kinase Reaction. <i>ACS Omega</i> , 2018, 3, 1783-1790.	3.5	16
54	Effect of different organic salts on amino acids partition behaviour in PEG-salt ATPS. <i>Fluid Phase Equilibria</i> , 2018, 456, 84-91.	2.5	20

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55	Tunable Hydrophobic Eutectic Solvents Based on Terpenes and Monocarboxylic Acids. ACS Sustainable Chemistry and Engineering, 2018, 6, 8836-8846.	6.7	207
56	Prediction and Experimental Validation of Co-solvent Influence on Michaelis Constants: A Thermodynamic Activity-Based Approach. Chemistry - A European Journal, 2018, 24, 16418-16425.	3.3	19
57	Inducing the Alternative Oxidase Forms Part of the Molecular Strategy of Anoxic Survival in Freshwater Bivalves. Frontiers in Physiology, 2018, 9, 100.	2.8	25
58	The Role of Polyfunctionality in the Formation of [Ch]Cl-Carboxylic Acid-Based Deep Eutectic Solvents. Industrial & Engineering Chemistry Research, 2018, 57, 11195-11209.	3.7	46
59	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	88
60	Thermodynamics of enzyme-catalyzed esterifications: I. Succinic acid esterification with ethanol. Applied Microbiology and Biotechnology, 2017, 101, 5973-5984.	3.6	16
61	Standard Gibbs energy of metabolic reactions: II. Glucose-6-phosphatase reaction and ATP hydrolysis. Biophysical Chemistry, 2017, 223, 30-38.	2.8	32
62	Thermodynamics of enzyme-catalyzed esterifications: II. Levulinic acid esterification with short-chain alcohols. Applied Microbiology and Biotechnology, 2017, 101, 7509-7521.	3.6	27
63	Crowders and Cosolvents – Major Contributors to the Cellular Milieu and Efficient Means to Counteract Environmental Stresses. ChemPhysChem, 2017, 18, 2951-2972.	2.1	82
64	Phase Behavior of Binary Mixtures Containing Succinic Acid or Its Esters. Journal of Chemical & Engineering Data, 2017, 62, 1983-1993.	1.9	10
65	Measuring and Predicting the Extraction Behavior of Biogenic Formic Acid in Biphasic Aqueous/Organic Reaction Mixtures. ACS Omega, 2017, 2, 8982-8989.	3.5	12
66	Influence of Salts on the Partitioning of 5-Hydroxymethylfurfural in Water/MIBK. Journal of Physical Chemistry B, 2016, 120, 3797-3808.	2.6	57
67	Thermodynamics of the alanine aminotransferase reaction. Fluid Phase Equilibria, 2016, 422, 87-98.	2.5	32
68	Benzoic Acid and Chlorobenzoic Acids: Thermodynamic Study of the Pure Compounds and Binary Mixtures With Water. Journal of Pharmaceutical Sciences, 2016, 105, 1050-1058.	3.3	35
69	Modelling interfacial properties of ionic liquids with ePC-SAFT combined with density gradient theory. Molecular Physics, 2016, 114, 2492-2499.	1.7	14
70	Standard Gibbs Energy of Metabolic Reactions: I. Hexokinase Reaction. Biochemistry, 2016, 55, 5665-5674.	2.5	23
71	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.	1.9	10
72	Thermodynamics of Bioreactions. Annual Review of Chemical and Biomolecular Engineering, 2016, 7, 395-414.	6.8	42

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73	Modeling the density of ionic liquids with ePC-SAFT. <i>Fluid Phase Equilibria</i> , 2016, 410, 9-22.	2.5	39
74	A Distinct Mitochondrial Genome with DUI-Like Inheritance in the Ocean Quahog <i>Arctica islandica</i> . <i>Molecular Biology and Evolution</i> , 2016, 33, 375-383.	8.9	20
75	PC-SAFT Modeling of CO <sub>2</sub> Solubilities in Deep Eutectic Solvents. <i>Journal of Physical Chemistry B</i> , 2016, 120, 2300-2310.	2.6	110
76	Thermodynamics of a model biological reaction: A comprehensive combined experimental and theoretical study. <i>Fluid Phase Equilibria</i> , 2016, 422, 99-110.	2.5	21
77	Thermochemistry of halogenobenzoic acids as an access to PC-SAFT solubility modeling. <i>Fluid Phase Equilibria</i> , 2016, 409, 399-407.	2.5	20
78	Activity coefficients at infinite dilution for different alcohols and ketones in [EMpy][ESO4]: Experimental data and modeling with PC-SAFT. <i>Fluid Phase Equilibria</i> , 2016, 424, 32-40.	2.5	12
79	Compatible solutes: Thermodynamic properties relevant for effective protection against osmotic stress. <i>Fluid Phase Equilibria</i> , 2016, 407, 224-235.	2.5	42
80	Measurement and PC-SAFT modelling of three-phase behaviour. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 1800-1810.	2.8	25
81	Predicting CO <sub>2</sub> solubility in aqueous N-methyldiethanolamine solutions with ePC-SAFT. <i>Fluid Phase Equilibria</i> , 2015, 393, 91-100.	2.5	42
82	Modeling thermodynamic derivative properties of ionic liquids with ePC-SAFT. <i>Fluid Phase Equilibria</i> , 2015, 405, 73-82.	2.5	43
83	Separation Performance of BioRenewable Deep Eutectic Solvents. <i>Industrial &amp; Engineering Chemistry Research</i> , 2015, 54, 3498-3504.	3.7	97
84	Benchmark Thermochemistry for Biologically Relevant Adenine and Cytosine. A Combined Experimental and Theoretical Study. <i>Journal of Physical Chemistry A</i> , 2015, 119, 9680-9691.	2.5	56
85	Effect of salts on the solubility of ionic liquids in water: experimental and electrolyte Perturbed-Chain Statistical Associating Fluid Theory. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 32044-32052.	2.8	22
86	Isolation and characterization of 11 microsatellite markers from the squat lobster <i>Munida gregaria</i> (Decapoda: Galatheidae) around the Falkland Islands/Islas Malvinas. <i>Conservation Genetics Resources</i> , 2015, 7, 147-149.	0.8	1
87	Telomere-independent ageing in the longest-lived non-colonial animal, <i>Arctica islandica</i> . <i>Experimental Gerontology</i> , 2014, 51, 38-45.	2.8	29
88	Measuring and modeling aqueous electrolyte/amino-acid solutions with ePC-SAFT. <i>Journal of Chemical Thermodynamics</i> , 2014, 68, 1-12.	2.0	97
89	Modeling imidazolium-based ionic liquids with ePC-SAFT. Part II. Application to H <sub>2</sub> S and synthesis-gas components. <i>Fluid Phase Equilibria</i> , 2014, 363, 59-65.	2.5	65
90	Vapor-Liquid Equilibria of Water + Alkylimidazolium-Based Ionic Liquids: Measurements and Perturbed-Chain Statistical Associating Fluid Theory Modeling. <i>Industrial &amp; Engineering Chemistry Research</i> , 2014, 53, 3737-3748.	3.7	82

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91	A thermodynamic investigation of the glucose-6-phosphate isomerization. <i>Biophysical Chemistry</i> , 2014, 195, 22-31.	2.8	29
92	Density of Mixtures Containing Sugars and Ionic Liquids: Experimental Data and PC-SAFT Modeling. <i>Journal of Chemical &amp; Engineering Data</i> , 2014, 59, 2942-2954.	1.9	36
93	Thermodynamic Properties of Aqueous Glucose-Urea-Salt Systems. <i>Journal of Solution Chemistry</i> , 2014, 43, 1110-1131.	1.2	12
94	ePC-SAFT revised. <i>Chemical Engineering Research and Design</i> , 2014, 92, 2884-2897.	5.6	158
95	Solubility of Sugars and Sugar Alcohols in Ionic Liquids: Measurement and PC-SAFT Modeling. <i>Journal of Physical Chemistry B</i> , 2013, 117, 9980-9995.	2.6	67
96	Isolation and characterization of 10 polymorphic loci for the giant Antarctic isopod, <i>Glyptonotus antarcticus</i> . <i>Conservation Genetics Resources</i> , 2013, 5, 963-965.	0.8	0
97	Liquid-Liquid Equilibria of 1-Butanol/Water/IL Systems. <i>Industrial &amp; Engineering Chemistry Research</i> , 2013, 52, 18472-18481.	3.7	90
98	The role of activity coefficients in bioreaction equilibria: Thermodynamics of methyl ferulate hydrolysis. <i>Biophysical Chemistry</i> , 2013, 173-174, 21-30.	2.8	29
99	Modeling thermodynamic properties of aqueous single-solute and multi-solute sugar solutions with PC-SAFT. <i>AIChE Journal</i> , 2013, 59, 4794-4805.	3.6	57
100	Molecular Interactions in 1-Butanol + IL Solutions by Measuring and Modeling Activity Coefficients. <i>Journal of Physical Chemistry B</i> , 2013, 117, 3173-3185.	2.6	41
101	Modeling imidazolium-based ionic liquids with ePC-SAFT. <i>Fluid Phase Equilibria</i> , 2012, 335, 64-73.	2.5	128
102	Measuring and modeling alcohol/salt systems. <i>Chemical Engineering Science</i> , 2012, 68, 328-339.	3.8	64
103	Thermodynamic properties of aqueous salt containing urea solutions. <i>Fluid Phase Equilibria</i> , 2012, 325, 71-79.	2.5	34
104	Measuring and Modeling Activity Coefficients in Aqueous Amino-Acid Solutions. <i>Industrial &amp; Engineering Chemistry Research</i> , 2011, 50, 131-141.	3.7	129
105	Modeling pH and Solubilities in Aqueous Multisolute Amino Acid Solutions. <i>Industrial &amp; Engineering Chemistry Research</i> , 2011, 50, 3503-3509.	3.7	45
106	Long-distance island hopping without dispersal stages: transportation across major zoogeographic barriers in a Southern Ocean isopod. <i>Die Naturwissenschaften</i> , 2010, 97, 583-594.	1.6	80
107	Compatible solutes: Thermodynamic properties and biological impact of ectoines and prolines. <i>Biophysical Chemistry</i> , 2010, 152, 28-39.	2.8	90
108	Measurement and Modeling Solubility of Aqueous Multisolute Amino-Acid Solutions. <i>Industrial &amp; Engineering Chemistry Research</i> , 2010, 49, 1395-1401.	3.7	69

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109	Modeling aqueous electrolyte solutions. Part 2. Weak electrolytes. Fluid Phase Equilibria, 2009, 279, 141-148.	2.5	90
110	Modelling Specific Ion Effects in Engineering Science. , 2009, , 85-115.		0
111	Encounter of lithodid crab <i>Paralomis birsteini</i> on the continental slope off Antarctica, sampled by ROV. Polar Biology, 2008, 31, 1143-1148.	1.2	41
112	Modeling aqueous electrolyte solutions. Fluid Phase Equilibria, 2008, 270, 87-96.	2.5	173
113	Cryptic speciation in the giant Antarctic isopod <i>Glyptonotus antarcticus</i> (Isopoda,) Tj ETQq1 1 0.784314 rgBT/Over 0.6 121	0.6	121