Kai Sundmacher

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.

8,349 46 69 393 h-index g-index citations papers 6.66 9,704 4.7 413 avg, IF L-index ext. citations ext. papers

#	Paper	IF	Citations
393	Sustainability of green solvents Ireview and perspective. <i>Green Chemistry</i> , 2022 , 24, 410-437	10	9
392	Integrated metalBrganic framework and pressure/vacuum swing adsorption process design: Descriptor optimization. <i>AICHE Journal</i> , 2022 , 68, e17524	3.6	2
391	Load-flexible fixed-bed reactors by multi-period design optimization. <i>Chemical Engineering Journal</i> , 2022 , 428, 130771	14.7	1
390	Closed-loop real-time optimization for unsteady operating production systems. <i>Journal of Process Control</i> , 2022 , 113, 80-95	3.9	
389	Increased efficiency of charge-mediated fusion in polymer/lipid hybrid membranes <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2022 , 119, e2122468119	11.5	3
388	Fusion-Induced Growth of Biomimetic Polymersomes: Behavior of Poly(dimethylsiloxane)-Poly(ethylene oxide) Vesicles in Saline Solutions Under High Agitation. <i>Macromolecular Rapid Communications</i> , 2021 , e2100712	4.8	4
387	Integrated computer-aided molecular and process design: Green solvents for the hydroformylation of long-chain olefines. <i>Chemical Engineering Science</i> , 2021 , 249, 117243	4.4	1
386	Computer-aided solvent screening for the fractionation of wet microalgae biomass. <i>Green Chemistry</i> , 2021 , 23, 10014-10029	10	0
385	Comparative Screening of Organic Solvents, Ionic Liquids, and Their Binary Mixtures for Vitamin E Extraction from Deodorizer Distillate. <i>Chemical Engineering and Processing: Process Intensification</i> , 2021 , 108711	3.7	O
384	Light-Powered Reactivation of Flagella and Contraction of Microtubule Networks: Toward Building an Artificial Cell. <i>ACS Synthetic Biology</i> , 2021 , 10, 1490-1504	5.7	2
383	Scale up of Transmembrane NADH Oxidation in Synthetic Giant Vesicles. <i>Bioconjugate Chemistry</i> , 2021 , 32, 897-903	6.3	2
382	Hybrid Data-Driven and Mechanistic Modeling Approaches for Multiscale Material and Process Design. <i>Engineering</i> , 2021 , 7, 1231-1231	9.7	5
381	Bottom-Up Synthesis of Artificial Cells: Recent Highlights and Future Challenges. <i>Annual Review of Chemical and Biomolecular Engineering</i> , 2021 , 12, 287-308	8.9	6
380	Model-based optimal design of phase change ionic liquids for efficient thermal energy storage. <i>Green Energy and Environment</i> , 2021 , 6, 392-404	5.7	8
379	Integrated ionic liquid and rate-based absorption process design for gas separation: Global optimization using hybrid models. <i>AICHE Journal</i> , 2021 , 67, e17340	3.6	8
378	Advances in the HCl gas-phase electrolysis employing an oxygen-depolarized cathode. <i>Electrochimica Acta</i> , 2021 , 365, 137282	6.7	0
377	Evaluation of COSMO-RS for solid[Iquid equilibria prediction of binary eutectic solvent systems. <i>Green Energy and Environment</i> , 2021 , 6, 371-379	5.7	10

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376	Computer-Aided Screening of Deep Eutectic Solvent Systems for the Associative Extraction of Frocopherol from Deodorizer Distillate. <i>Computer Aided Chemical Engineering</i> , 2021 , 50, 341-346	0.6	
375	Neural recommender system for the activity coefficient prediction and UNIFAC model extension of ionic liquid-solute systems. <i>AICHE Journal</i> , 2021 , 67, e17171	3.6	14
374	Rational Design of Ionic Liquid Phase-Change Material for Efficient Thermal Energy Storage. <i>Computer Aided Chemical Engineering</i> , 2021 , 191-196	0.6	0
373	Multi-Period Design Optimization of Flexible Fixed-Bed Reactors by Stoichiometry-Based Model Reduction. <i>Computer Aided Chemical Engineering</i> , 2021 , 947-952	0.6	
372	Power-to-Chemicals: A Superstructure Problem for Sustainable Syngas Production. <i>Mathematics in Industry</i> , 2021 , 145-168	0.2	0
371	Forschungsarbeiten am Institut fil Verfahrenstechnik der Otto-von-Guericke-Universit i l Magdeburg. <i>Chemie-Ingenieur-Technik</i> , 2021 , 93, 345-352	0.8	
370	ECarotene extraction from Dunaliella salina by supercritical CO2. <i>Journal of Applied Phycology</i> , 2021 , 33, 1435-1445	3.2	5
369	Guest Editorial Special Issue on Deep Integration of Artificial Intelligence and Data Science for Process Manufacturing. <i>IEEE Transactions on Neural Networks and Learning Systems</i> , 2021 , 32, 3294-329	5 ^{10.3}	
368	En route to dynamic life processes by SNARE-mediated fusion of polymer and hybrid membranes. <i>Nature Communications</i> , 2021 , 12, 4972	17.4	10
367	Decoupling oxygen and water transport dynamics in polymer electrolyte membrane fuel cells through frequency response methods based on partial pressure perturbations. <i>Electrochimica Acta</i> , 2021 , 390, 138788	6.7	2
366	Non-intrusive Time-POD for Optimal Control of a Fixed-Bed Reactor for CO2 Methanation. <i>IFAC-PapersOnLine</i> , 2021 , 54, 122-127	0.7	
365	Hybrid Semi-parametric Modeling in Separation Processes: A Review. <i>Chemie-Ingenieur-Technik</i> , 2020 , 92, 842-855	0.8	10
364	Model-Based Analysis of the Limiting Mechanisms in the Gas-Phase Oxidation of HCl Employing an Oxygen Depolarized Cathode. <i>Journal of the Electrochemical Society</i> , 2020 , 167, 013537	3.9	2
363	Symmetry Breaking and Emergence of Directional Flows in Minimal Actomyosin Cortices. <i>Cells</i> , 2020 , 9,	7.9	4
362	Light-Driven ATP Regeneration in Diblock/Grafted Hybrid Vesicles. <i>ChemBioChem</i> , 2020 , 21, 2149-2160	3.8	22
361	Selectivity and Sustainability of Electroenzymatic Process for Glucose Conversion to Gluconic Acid. <i>Catalysts</i> , 2020 , 10, 269	4	4
360	Transformation of remnant algal biomass to 5-HMF and levulinic acid: influence of a biphasic solvent system <i>RSC Advances</i> , 2020 , 10, 24753-24763	3.7	17
359	Constructing artificial respiratory chain in polymer compartments: Insights into the interplay between oxidase and the membrane. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2020 , 117, 15006-15017	11.5	22

358	Systematic Green Solvent Selection for the Hydroformylation of Long-Chain Alkenes. <i>ACS Sustainable Chemistry and Engineering</i> , 2020 ,	8.3	2
357	Time-minimal set point transition for nonlinear SISO systems under different constraints. <i>Automatica</i> , 2020 , 114, 108806	5.7	3
356	Spezielle labortechnische Reaktoren: Wendelrohrreaktor. <i>Springer Reference Naturwissenschaften</i> , 2020 , 1289-1320	0.2	
355	Numerical Methods for Coupled Population Balance Systems Applied to the Dynamical Simulation of Crystallization Processes 2020 , 475-518		Ο
354	Systematic Screening of Deep Eutectic Solvents as Sustainable Separation Media Exemplified by the CO2 Capture Process. <i>ACS Sustainable Chemistry and Engineering</i> , 2020 , 8, 8741-8751	8.3	29
353	Screening of functional solvent system for automatic aldehyde and ketone separation in aldol reaction: A combined COSMO-RS and experimental approach. <i>Chemical Engineering Journal</i> , 2020 , 385, 123399	14.7	12
352	Extending the UNIFAC model for ionic liquidBolute systems by combining experimental and computational databases. <i>AICHE Journal</i> , 2020 , 66, e16821	3.6	29
351	Reconstruction and analysis of a carbon-core metabolic network for Dunaliella salina. <i>BMC Bioinformatics</i> , 2020 , 21, 1	3.6	104
350	The FluxMax approach: Simultaneous flux optimization and heat integration by discretization of thermodynamic state space illustrated on methanol synthesis process. <i>Chemical Engineering Science</i> , 2020 , 215, 115382	4.4	6
349	Optimal catalyst particle design for flexible fixed-bed CO2 methanation reactors. <i>Chemical Engineering Journal</i> , 2020 , 387, 123704	14.7	34
349		14.7 5.4	29
	Engineering Journal, 2020, 387, 123704 Computer-aided solvent selection and design for efficient chemical processes. Current Opinion in		
348	Engineering Journal, 2020, 387, 123704 Computer-aided solvent selection and design for efficient chemical processes. Current Opinion in Chemical Engineering, 2020, 27, 35-44 Computer Aided Molecular Design of Green Solvents for the Hydroformylation of Long-Chain	5.4	29
348 347	Computer-aided solvent selection and design for efficient chemical processes. <i>Current Opinion in Chemical Engineering</i> , 2020 , 27, 35-44 Computer Aided Molecular Design of Green Solvents for the Hydroformylation of Long-Chain Olefines. <i>Computer Aided Chemical Engineering</i> , 2020 , 48, 745-750 In silico Screening of Metal-organic Frameworks for Acetylene/ethylene Separation. <i>Computer</i>	5.4	29
348 347 346	Computer-aided solvent selection and design for efficient chemical processes. <i>Current Opinion in Chemical Engineering</i> , 2020 , 27, 35-44 Computer Aided Molecular Design of Green Solvents for the Hydroformylation of Long-Chain Olefines. <i>Computer Aided Chemical Engineering</i> , 2020 , 48, 745-750 In silico Screening of Metal-organic Frameworks for Acetylene/ethylene Separation. <i>Computer Aided Chemical Engineering</i> , 2020 , 895-900 Power-to-Syngas Processes by Reactor-Separator Superstructure Optimization. <i>Computer Aided</i>	5·4 o.6 o.6	29
348 347 346 345	Computer-aided solvent selection and design for efficient chemical processes. Current Opinion in Chemical Engineering, 2020, 27, 35-44 Computer Aided Molecular Design of Green Solvents for the Hydroformylation of Long-Chain Olefines. Computer Aided Chemical Engineering, 2020, 48, 745-750 In silico Screening of Metal-organic Frameworks for Acetylene/ethylene Separation. Computer Aided Chemical Engineering, 2020, 895-900 Power-to-Syngas Processes by Reactor-Separator Superstructure Optimization. Computer Aided Chemical Engineering, 2020, 48, 1387-1392 CO2 methanation process synthesis by superstructure optimization. Journal of CO2 Utilization,	5.4 o.6 o.6	29
348347346345344	Computer-aided solvent selection and design for efficient chemical processes. Current Opinion in Chemical Engineering, 2020, 27, 35-44 Computer Aided Molecular Design of Green Solvents for the Hydroformylation of Long-Chain Olefines. Computer Aided Chemical Engineering, 2020, 48, 745-750 In silico Screening of Metal-organic Frameworks for Acetylene/ethylene Separation. Computer Aided Chemical Engineering, 2020, 895-900 Power-to-Syngas Processes by Reactor-Separator Superstructure Optimization. Computer Aided Chemical Engineering, 2020, 48, 1387-1392 CO2 methanation process synthesis by superstructure optimization. Journal of CO2 Utilization, 2020, 40, 101228 Polymer Electrolyte Fuel Cell Degradation Mechanisms and Their Diagnosis by Frequency Response	5.4 o.6 o.6 o.6	29 1 3

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340	Mixed-Integer Linear Programming (MILP) Approach for the Synthesis of Efficient Power-to-Syngas Processes. <i>Frontiers in Energy Research</i> , 2020 , 8,	3.8	4
339	Porosity and Structure of Hierarchically Porous Ni/Al2O3 Catalysts for CO2 Methanation. <i>Catalysts</i> , 2020 , 10, 1471	4	12
338	Big Data Creates New Opportunities for Materials Research: A Review on Methods and Applications of Machine Learning for Materials Design. <i>Engineering</i> , 2019 , 5, 1017-1026	9.7	78
337	Electrochemical gas phase oxidation of hydrogen chloride to chlorine: Model-based analysis of transport and reaction mechanisms. <i>Electrochimica Acta</i> , 2019 , 324, 134780	6.7	3
336	Analysis of mass transport processes in the anodic porous transport layer in PEM water electrolysers. <i>International Journal of Hydrogen Energy</i> , 2019 , 44, 28077-28087	6.7	12
335	Miniplant-Scale Evaluation of a Semibatch-Continuous Tandem Reactor System for the Hydroformylation of Long-Chain Olefins. <i>Industrial & Engineering Chemistry Research</i> , 2019 , 58, 247	1 ²⁻² 48	03
334	The FluxMax approach for simultaneous process synthesis and heat integration: Production of hydrogen cyanide. <i>AICHE Journal</i> , 2019 , 65, e16554	3.6	11
333	Energy-Efficient Gas-Phase Electrolysis of[Hydrogen Chloride. <i>Chemie-Ingenieur-Technik</i> , 2019 , 91, 795-8	3 0 88	4
332	Rational design of double salt ionic liquids as extraction solvents: Separation of thiophene/n-octane as example. <i>AICHE Journal</i> , 2019 , 65, e16625	3.6	21
331	Radiation-based model reduction for the optimization of high temperature tube bundle reactors: Synthesis of hydrogen cyanide. <i>Computers and Chemical Engineering</i> , 2019 , 127, 186-199	4	2
330	Compartments for Synthetic Cells: Osmotically Assisted Separation of Oil from Double Emulsions in a Microfluidic Chip. <i>ChemBioChem</i> , 2019 , 20, 2604-2608	3.8	11
329	Derivation of rate equations for equilibrium limited gas-solid reactions. <i>Chemical Engineering Science</i> , 2019 , 203, 76-85	4.4	1
328	Artificial Organelles for Energy Regeneration. Advanced Biology, 2019, 3, e1800323	3.5	18
327	Polymer-Based Module for NAD Regeneration with Visible Light. <i>ChemBioChem</i> , 2019 , 20, 2593-2596	3.8	18
326	Optimal Solvent Design for Extractive Distillation Processes: A Multiobjective Optimization-Based Hierarchical Framework. <i>Industrial & Engineering Chemistry Research</i> , 2019 , 58, 5777-5786	3.9	38
325	Directed Growth of Biomimetic Microcompartments. <i>Advanced Biology</i> , 2019 , 3, e1800314	3.5	14
324	Operation range extension via hot-spot control for catalytic CO2 methanation reactors. <i>Reaction Chemistry and Engineering</i> , 2019 , 4, 1019-1037	4.9	40
323	Stochastic-deterministic population balance modeling and simulation of a fluidized bed crystallizer experiment. <i>Chemical Engineering Science</i> , 2019 , 208, 115102	4.4	4

322	Systematic Selection of Green Solvents and Process Optimization for the Hydroformylation of Long-Chain Olefines. <i>Processes</i> , 2019 , 7, 882	2.9	5
321	Surrogate Modeling for Liquidliquid Equilibria Using a Parameterization of the Binodal Curve. <i>Processes</i> , 2019 , 7, 753	2.9	4
320	Global optimization of distillation columns using explicit and implicit surrogate models. <i>Chemical Engineering Science</i> , 2019 , 197, 235-245	4.4	30
319	Overview of Surrogate Modeling in Chemical Process Engineering. <i>Chemie-Ingenieur-Technik</i> , 2019 , 91, 228-239	0.8	75
318	Exergetic assessment of CO2 methanation processes for the chemical storage of renewable energies. <i>Applied Energy</i> , 2019 , 233-234, 271-282	10.7	32
317	Cyclic operation of a semi-batch reactor for the hydroformylation of long-chain olefins and integration in a continuous production process. <i>Chemical Engineering Journal</i> , 2019 , 377, 120453	14.7	4
316	Helically coiled segmented flow tubular reactor for the hydroformylation of long-chain olefins in a thermomorphic multiphase system. <i>Chemical Engineering Journal</i> , 2019 , 377, 120060	14.7	19
315	Productivity versus product quality: Exploring the limits of autothermal microchannel reactors in methane steam reforming. <i>Chemical Engineering Journal</i> , 2019 , 377, 120048	14.7	7
314	Transmembrane NADH Oxidation with Tetracyanoquinodimethane. <i>Langmuir</i> , 2018 , 34, 5435-5443	4	9
313	Techno-Ronomische Optimierung des Produktionsnetzwerkes fil die Synthese von Ameisensüre aus erneuerbaren Ressourcen. <i>Chemie-Ingenieur-Technik</i> , 2018 , 90, 256-266	0.8	2
312	Computationally efficient NMPC for batch and semi-batch processes using parsimonious input parameterization. <i>Journal of Process Control</i> , 2018 , 66, 12-22	3.9	4
311	Particle-image-velocimetry measurements in organic liquid multiphase systems for an optimal reactor design and operation. <i>Journal of Visualization</i> , 2018 , 21, 5-17	1.6	
310	Computer-aided design of ionic liquids as solvents for extractive desulfurization. <i>AICHE Journal</i> , 2018 , 64, 1013-1025	3.6	97
309	Sequential bottom-up assembly of mechanically stabilized synthetic cells by microfluidics. <i>Nature Materials</i> , 2018 , 17, 89-96	27	211
308	NMPC using Pontryagin Minimum Principle-Application to a two-phase semi-batch hydroformylation reactor under uncertainty. <i>Computers and Chemical Engineering</i> , 2018 , 108, 47-56	4	13
307	Reactor-network synthesis via flux profile analysis. <i>Chemical Engineering Journal</i> , 2018 , 335, 1018-1030	14.7	21
306	Linear Programming Approach for Structure Optimization of Renewable-to-Chemicals (R2Chem) Production Networks. <i>Industrial & Engineering Chemistry Research</i> , 2018 , 57, 9889-9902	3.9	11
305	MaxSynBio: Avenues Towards Creating Cells from the Bottom Up. <i>Angewandte Chemie - International Edition</i> , 2018 , 57, 13382-13392	16.4	155

304	Integrated Process and Ionic Liquid Design by Combining Flowsheet Simulation with Quantum-Chemical Solvent Screening. <i>Computer Aided Chemical Engineering</i> , 2018 , 44, 2167-2172	0.6	4	
303	Computer Aided Design of Green Thermomorphic Solvent Systems for Homogeneous Catalyst Recovery. <i>Computer Aided Chemical Engineering</i> , 2018 , 1783-1788	0.6	5	
302	Out-of-equilibrium microcompartments for the bottom-up integration of metabolic functions. <i>Nature Communications</i> , 2018 , 9, 2391	17.4	41	
301	Spezielle labortechnische Reaktoren: Wendelrohrreaktor. <i>Springer Reference Naturwissenschaften</i> , 2018 , 1-33	0.2		
300	Continuous production of CO from CO2 by RWGS chemical looping in fixed and fluidized bed reactors. <i>Chemical Engineering Journal</i> , 2018 , 336, 278-296	14.7	25	
299	Thermodynamic Network Flow Approach for Chemical Process Synthesis. <i>Computer Aided Chemical Engineering</i> , 2018 , 43, 881-886	0.6	2	
298	Quantitative single cell analysis uncovers the life/death decision in CD95 network. <i>PLoS Computational Biology</i> , 2018 , 14, e1006368	5	15	
297	Identification of Key Transport Phenomena in High-Temperature Reactors: Flow and Heat Transfer Characteristics. <i>Industrial & Engineering Chemistry Research</i> , 2018 , 57, 15884-15897	3.9	4	
296	Mechanisms behind overshoots in mean cluster size profiles in aggregation-breakup processes. Journal of Colloid and Interface Science, 2018 , 528, 336-348	9.3	5	
295	MaxSynBio: Wege zur Synthese einer Zelle aus nicht lebenden Komponenten. <i>Angewandte Chemie</i> , 2018 , 130, 13566-13577	3.6	25	
294	Toward Fast Dynamic Optimization: An Indirect Algorithm That Uses Parsimonious Input Parameterization. <i>Industrial & Engineering Chemistry Research</i> , 2018 , 57, 10038-10048	3.9	8	
293	A hybrid stochastic deterministic optimization approach for integrated solvent and process design. <i>Chemical Engineering Science</i> , 2017 , 159, 207-216	4.4	40	
292	Measurement and simulation of mass transfer and backmixing behavior in a gas-liquid helically coiled tubular reactor. <i>Chemical Engineering Science</i> , 2017 , 170, 410-421	4.4	23	
291	Dynamic optimization of constrained semi-batch processes using Pontryagin minimum principle An effective quasi-Newton approach. <i>Computers and Chemical Engineering</i> , 2017 , 99, 135-144	4	14	
290	Systematic Method for Screening Ionic Liquids as Extraction Solvents Exemplified by an Extractive Desulfurization Process. <i>ACS Sustainable Chemistry and Engineering</i> , 2017 , 5, 3382-3389	8.3	92	
289	Continuous Crystallization in a Helically Coiled Flow Tube: Analysis of Flow Field, Residence Time Behavior, and Crystal Growth. <i>Industrial & Engineering Chemistry Research</i> , 2017 , 56, 3699-3712	3.9	32	
288	POD-DEIM for efficient reduction of a dynamic 2D catalytic reactor model. <i>Computers and Chemical Engineering</i> , 2017 , 106, 777-784	4	9	
287	CO production from CO 2 via reverse watergas shift reaction performed in a chemical looping mode: Kinetics on modified iron oxide. <i>Journal of CO2 Utilization</i> , 2017 , 17, 60-68	7.6	35	

286	Concentration-alternating frequency response: A new method for studying polymer electrolyte membrane fuel cell dynamics. <i>Electrochimica Acta</i> , 2017 , 243, 53-64	6.7	13
285	Understanding PEM fuel cell dynamics: The reversal curve. <i>International Journal of Hydrogen Energy</i> , 2017 , 42, 15818-15827	6.7	18
284	Disorientation angle distribution of primary particles in potash alum aggregates. <i>Journal of Crystal Growth</i> , 2017 , 467, 93-106	1.6	9
283	Catalyst Layer Modeling 2017 , 259-285		1
282	Model-based Optimal Sabatier Reactor Design for Power-to-Gas Applications. <i>Energy Technology</i> , 2017 , 5, 911-921	3.5	31
281	Toward Artificial Mitochondrion: Mimicking Oxidative Phosphorylation in Polymer and Hybrid Membranes. <i>Nano Letters</i> , 2017 , 17, 6816-6821	11.5	71
2 80	Economic linear objective function approach for structure optimization of renewables-to-chemicals (R2Chem) networks. <i>Computer Aided Chemical Engineering</i> , 2017 , 40, 1975-1980	0.6	3
279	Set point tracking of a biogas plant coupled to a methanation reactor. <i>Computer Aided Chemical Engineering</i> , 2017 , 40, 1645-1650	0.6	
278	Influence of the autonomous oscillations and the CO concentration on the performance of an ECPrOx reactor. <i>Electrochimica Acta</i> , 2017 , 251, 602-612	6.7	1
277	Optimal Reactor Design via Flux Profile Analysis for an Integrated Hydroformylation Process. <i>Industrial & Design Engineering Chemistry Research</i> , 2017 , 56, 11507-11518	3.9	16
276	Carotenoid Production Process Using Green Microalgae of the Dunaliella Genus: Model-Based Analysis of Interspecies Variability. <i>Industrial & Engineering Chemistry Research</i> , 2017 , 56, 12888-12	8398	3
275	Numerical study of liquid-liquid mixing in helical pipes. <i>Chemical Engineering Science</i> , 2017 , 172, 250-267	1 4.4	46
274	Crystal Population Growth in a Continuous Helically Coiled Flow Tube Crystallizer. <i>Chemical Engineering and Technology</i> , 2017 , 40, 1584-1590	2	12
273	Thermodynamic analysis and optimization of RWGS processes for solar syngas production from CO2. <i>AICHE Journal</i> , 2017 , 63, 15-22	3.6	24
272	CO2 methanation: Optimal start-up control of a fixed-bed reactor for power-to-gas applications. <i>AICHE Journal</i> , 2017 , 63, 23-31	3.6	59
271	Recovery and Separation of Carbohydrate Derivatives from the Lipid Extracted Alga Dunaliella by Mild Liquefaction. <i>ACS Sustainable Chemistry and Engineering</i> , 2017 , 5, 588-595	8.3	6
270	Integrated reaction extraction process for the hydroformylation of long-chain alkenes with a homogeneous catalyst. <i>Computers and Chemical Engineering</i> , 2017 , 105, 212-223	4	20
269	Efficient simulation of intrinsic, extrinsic and external noise in biochemical systems. <i>Bioinformatics</i> , 2017 , 33, i319-i324	7.2	3

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268	Physics-Based Surrogate Models for Optimal Control of a CO2 Methanation Reactor. <i>Computer Aided Chemical Engineering</i> , 2017 , 40, 127-132	0.6	1
267	Process Optimization by Applying a Simultaneous Dynamic Method. <i>Computer Aided Chemical Engineering</i> , 2017 , 40, 2047-2052	0.6	
266	Dynamic Optimization of Constrained Semi-Batch Processes using Pontryagin Minimum Principle and Parsimonious Parameterization. <i>Computer Aided Chemical Engineering</i> , 2017 , 40, 2041-2046	0.6	
265	Computationally Efficient Steady-State Process Simulation by Applying a Simultaneous Dynamic Method. <i>Computer Aided Chemical Engineering</i> , 2016 , 38, 517-522	0.6	3
264	Probabilistic reactor design in the framework of elementary process functions. <i>Computers and Chemical Engineering</i> , 2016 , 94, 45-59	4	26
263	Autonomous Voltage Oscillations in a Direct Methanol Fuel Cell. <i>Electrochimica Acta</i> , 2016 , 212, 545-55.	2 6.7	17
262	Structure optimization of power-to-chemicals (P2C) networks by linear programming for the economic utilization of renewable surplus energy. <i>Computer Aided Chemical Engineering</i> , 2016 , 1551-15	5 <mark>6</mark> .6	10
261	A Short-Cut Method for the Quantification of Crystallization Kinetics. 1. Method Development. <i>Crystal Growth and Design</i> , 2016 , 16, 6743-6755	3.5	9
260	Dynamische Methode zur Berechnung thermodynamischer Gleichgewichte in reaktiven Mehrphasensystemen. <i>Chemie-Ingenieur-Technik</i> , 2016 , 88, 1617-1627	0.8	2
259	Thermomorphic solvent selection for homogeneous catalyst recovery based on COSMO-RS. <i>Chemical Engineering and Processing: Process Intensification</i> , 2016 , 99, 97-106	3.7	36
258	Flow cytometry enables dynamic tracking of algal stress response: A case study using carotenogenesis in Dunaliella salina. <i>Algal Research</i> , 2016 , 13, 227-234	5	13
257	Optimal configuration and pressure levels of electrolyzer plants in context of power-to-gas applications. <i>Applied Energy</i> , 2016 , 167, 107-124	10.7	54
256	Image-based analytical crystal shape computation exemplified for potassium dihydrogen phosphate (KDP). <i>Chemical Engineering Science</i> , 2016 , 139, 61-74	4.4	3
255	Nonlinear Model Order Reduction for Catalytic Tubular Reactors. <i>Computer Aided Chemical Engineering</i> , 2016 , 38, 2373-2378	0.6	1
254	Efficient simulation of heterogeneity and stochasticity in microbial processes. <i>Computer Aided Chemical Engineering</i> , 2016 , 1213-1218	0.6	
253	Design and Comparison of Optimal Reactor Concepts for the Hydroformylation of Olefins by Use of a Probabilistic Design Framework. <i>Computer Aided Chemical Engineering</i> , 2016 , 38, 1365-1370	0.6	
252	Optimal design of solvents for extractive reaction processes. <i>AICHE Journal</i> , 2016 , 62, 3238-3249	3.6	27
251	Hydrogen and Carbon Monoxide Production by Chemical Looping over Iron-Aluminium Oxides. <i>Energy Technology</i> , 2016 , 4, 304-313	3.5	31

250	The interaction of protein-coated bionanoparticles and surface receptors reevaluated: how important is the number of bonds?. <i>Soft Matter</i> , 2016 , 12, 6451-62	3.6	2
249	Material development and process optimization for gas-phase hydrogen chloride electrolysis with oxygen depolarized cathode. <i>Journal of Applied Electrochemistry</i> , 2016 , 46, 755-767	2.6	6
248	Optimal Control of Crystal Shapes in Batch Crystallization Experiments by Growth-Dissolution Cycles. <i>Crystal Growth and Design</i> , 2016 , 16, 3297-3306	3.5	33
247	Binding kinetics and multi-bond: Finding correlations by synthesizing interactions between ligand-coated bionanoparticles and receptor surfaces. <i>Analytical Biochemistry</i> , 2016 , 505, 8-17	3.1	1
246	Diagnostic concept for dynamically operated biogas production plants. <i>Renewable Energy</i> , 2016 , 96, 47	98489	16
245	Dynamic flux balance modeling to increase the production of high-value compounds in green microalgae. <i>Biotechnology for Biofuels</i> , 2016 , 9, 165	7.8	19
244	Valorization of the aqueous phase obtained from hydrothermally treated Dunaliella salina remnant biomass. <i>Bioresource Technology</i> , 2016 , 219, 64-71	11	22
243	Rational selection of experimental readout and intervention sites for reducing uncertainties in computational model predictions. <i>BMC Bioinformatics</i> , 2015 , 16, 13	3.6	3
242	Model-based analysis of a gas/vaporliquid microchannel membrane contactor. <i>AICHE Journal</i> , 2015 , 61, 2240-2256	3.6	10
241	Fast evaluation of univariate aggregation integrals on equidistant grids. <i>Computers and Chemical Engineering</i> , 2015 , 74, 115-127	4	12
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