

Yorck-Michael Neuhold

List of Publications by Year in descending order

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30
papers

1,241
citations

471509

17
h-index

526287

27
g-index

31
all docs

31
docs citations

31
times ranked

1330
citing authors

#	ARTICLE	IF	CITATIONS
1	Activity-based analysis of potentiometric pH titrations. <i>Analytica Chimica Acta</i> , 2019, 1075, 49-56.	5.4	4
2	HERFD XAS/ATR-FTIR batch reactor cell. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 2164-2170.	2.8	29
3	Combined in Situ Monitoring Method for Analysis and Optimization of the Lithiation-Fluoroacetylation of N-(4-Chlorophenyl)-Pivalamide. <i>Industrial & Engineering Chemistry Research</i> , 2011, 50, 5982-5991.	3.7	4
4	In situ infrared monitoring of the solid/liquid catalyst interface during the three-phase hydrogenation of nitrobenzene over nanosized Au on TiO ₂ . <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 12463.	2.8	77
5	On-line Reaction Monitoring of Lithiation of Halogen Substituted Acetanilides via in situ Calorimetry, ATR Spectroscopy, and Endoscopy. <i>Chimia</i> , 2011, 65, 253.	0.6	0
6	Conjoint Analysis of Kinetic and Equilibrium Data for Mechanistic Elucidation in Polynuclear Complexation Reactions, Exemplified by Metal(II) Helicate Complex Formation. <i>Australian Journal of Chemistry</i> , 2010, 63, 141.	0.9	0
7	Nonisothermal Calorimetry for Fast Thermokinetic Reaction Analysis: Solvent-Free Esterification of <i>n</i> -Butanol by Acetic Anhydride. <i>Organic Process Research and Development</i> , 2010, 14, 524-536.	2.7	11
8	Photoreactions of Mercury in Surface Ocean Water: Gross Reaction Kinetics and Possible Pathways. <i>Environmental Science & Technology</i> , 2010, 44, 644-649.	10.0	106
9	Systematic prediction of linear dependencies in the concentration profiles and implications on the kinetic hard-modelling of spectroscopic data. <i>Chemometrics and Intelligent Laboratory Systems</i> , 2009, 95, 170-187.	3.5	6
10	Kinetic hard-modelling and spectral validation of rank-deficient spectroscopic data: A case study. <i>Chemometrics and Intelligent Laboratory Systems</i> , 2009, 98, 213-226.	3.5	4
11	Temperature oscillation calorimetry for the determination of the heat capacity in a small-scale reactor. <i>Chemical Engineering Science</i> , 2008, 63, 3755-3765.	3.8	10
12	Multivariate kinetic hard-modelling of spectroscopic data: A comparison of the esterification of butanol by acetic anhydride on different scales and with different instruments. <i>Chemical Engineering Science</i> , 2008, 63, 4800-4809.	3.8	17
13	Uncertainties and error propagation in kinetic hard-modelling of spectroscopic data. <i>Chemometrics and Intelligent Laboratory Systems</i> , 2008, 93, 120-131.	3.5	21
14	Complexation Kinetics of Copper(II) and Nickel(II) with Macrocycles: Identification of an Outer-Sphere Chelate Effect. <i>Inorganic Chemistry</i> , 2007, 46, 4002-4009.	4.0	21
15	Advances in the modelling and analysis of complex and industrial processes. <i>Chemometrics and Intelligent Laboratory Systems</i> , 2006, 82, 75-82.	3.5	15
16	Kinetic Modeling of Multivariate Measurements with Nonlinear Regression. , 2006, , 217-262.		0
17	Application of a genetic algorithm: near optimal estimation of the rate and equilibrium constants of complex reaction mechanisms. <i>Chemometrics and Intelligent Laboratory Systems</i> , 2004, 70, 193-203.	3.5	32
18	Analyses of three-way data from equilibrium and kinetic investigations. <i>Analytica Chimica Acta</i> , 2003, 490, 99-108.	5.4	23

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19	Analysis of reactions in aqueous solution at non-constant pH: no more buffers?. <i>Physical Chemistry Chemical Physics</i> , 2003, 5, 2836.	2.8	52
20	Copper(I) Dioxygen Reactivity of [(L)CuI]+ (L = Tris(2-pyridylmethyl)amine): Kinetic/Thermodynamic and Spectroscopic Studies Concerning the Formation of Cu ^{II} O ₂ and Cu ²⁺ O ₂ Adducts as a Function of Solvent Medium and 4-Pyridyl Ligand Substituent Variations. <i>Inorganic Chemistry</i> , 2003, 42, 1807-1824.	4.0	179
21	Simulation of Complex Chemical Kinetics. <i>Bioinorganic Reaction Mechanisms</i> , 2003, 5, .	0.4	5
22	Simulation of Complex Chemical Kinetics. <i>Bioinorganic Reaction Mechanisms</i> , 2003, 5, 39-46.	0.4	4
23	An electrospray ionization mass spectrometry study of the nitroprusside cation-thiolate system. <i>Dalton Transactions RSC</i> , 2002, , 3649-3655.	2.3	14
24	Hard-modelled trilinear decomposition (HTD) for an enhanced kinetic multicomponent analysis. <i>Journal of Chemometrics</i> , 2002, 16, 218-227.	1.3	22
25	Rank annihilation correction for the amendment of instrumental inconsistencies. <i>Analytica Chimica Acta</i> , 2002, 464, 249-259.	5.4	25
26	(F8TPP)FeII/O ₂ Reactivity Studies {F8TPP = Tetrakis(2,6-difluorophenyl)porphyrinate(2-)}: Spectroscopic (UV-Visible and NMR) and Kinetic Study of Solvent-Dependent (Fe/O ₂ = 1:1 or 2:1) Reversible O ₂ -Reduction and Ferryl Formation. <i>Inorganic Chemistry</i> , 2001, 40, 5754-5767.	4.0	121
27	Formation and Characterization of a High-Spin Heme-Copper Dioxygen (Peroxo) Complex. <i>Journal of the American Chemical Society</i> , 1999, 121, 9885-9886.	13.7	78
28	Oxo- and Hydroxo-Bridged Heme-Copper Assemblies Formed from Acid-Base or Metal Dioxygen Chemistry. <i>Inorganic Chemistry</i> , 1999, 38, 3093-3102.	4.0	43
29	Peroxo-, Oxo-, and Hydroxo-Bridged Dicopper Complexes: Observation of Exogenous Hydrocarbon Substrate Oxidation. <i>Journal of the American Chemical Society</i> , 1998, 120, 12960-12961.	13.7	180
30	Binucleating Ligand Structural Effects on (1/4-Peroxo)- and Bis(1/4-oxo)dicopper Complex Formation and Decay: Competition between Arene Hydroxylation and Aliphatic C-H Bond Activation. <i>Inorganic Chemistry</i> , 1997, 36, 6343-6356.	4.0	138