

Tibor Nagy

List of Publications by Year in descending order

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64
papers

1,402
citations

394421

19
h-index

345221

36
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66
all docs

66
docs citations

66
times ranked

1477
citing authors

#	ARTICLE	IF	CITATIONS
1	Design of combustion experiments using differential entropy. <i>Combustion Theory and Modelling</i> , 2022, 26, 67-90.	1.9	4
2	Comparison of methane combustion mechanisms using laminar burning velocity measurements. <i>Combustion and Flame</i> , 2022, 238, 111867.	5.2	17
3	A Short-Cut Data Mining Method for the Mass Spectrometric Characterization of Block Copolymers. <i>Processes</i> , 2022, 10, 42.	2.8	1
4	Quantification of Polyethylene Glycol 400 Excreted in the Urine by MALDI-TOF Mass Spectrometry. <i>Pharmaceutics</i> , 2022, 14, 1341.	4.5	1
5	Isocyanonaphthol Derivatives: Excited-State Proton Transfer and Solvatochromic Properties. <i>International Journal of Molecular Sciences</i> , 2022, 23, 7250.	4.1	3
6	Mass Spectral Filtering by Mass-Remainder Analysis (MARA) at High Resolution and Its Application to Metabolite Profiling of Flavonoids. <i>International Journal of Molecular Sciences</i> , 2021, 22, 864.	4.1	4
7	Structural Characterization of Daunomycin-Peptide Conjugates by Various Tandem Mass Spectrometric Techniques. <i>International Journal of Molecular Sciences</i> , 2021, 22, 1648.	4.1	1
8	Correlation between Heavy Metal-Induced Histopathological Changes and Trophic Interactions between Different Fish Species. <i>Applied Sciences (Switzerland)</i> , 2021, 11, 3760.	2.5	15
9	Comparison of Methane Combustion Mechanisms Using Shock Tube and Rapid Compression Machine Ignition Delay Time Measurements. <i>Energy & Fuels</i> , 2021, 35, 12329-12351.	5.1	23
10	Polydispersity Ratio and Its Application for the Characterization of Poloxamers. <i>Macromolecules</i> , 2021, 54, 9984-9991.	4.8	5
11	Study on the bZIP-Type Transcription Factors NapA and RsmA in the Regulation of Intracellular Reactive Species Levels and Sterigmatocystin Production of <i>Aspergillus nidulans</i> . <i>International Journal of Molecular Sciences</i> , 2021, 22, 11577.	4.1	4
12	Synthesis and characterization of isophorondiamine-based oligoamides: catalytic effect of amides during the curing of epoxy resins. <i>Polymer Bulletin</i> , 2020, 77, 4655-4678.	3.3	7
13	Biochemical characterization of Ty1 retrotransposon protease. <i>PLoS ONE</i> , 2020, 15, e0227062.	2.5	9
14	Automatic kinetic model generation and selection based on concentration versus time curves. <i>International Journal of Chemical Kinetics</i> , 2020, 52, 109-123.	1.6	3
15	Biochemical Characterization of Human Retroviral-Like Aspartic Protease 1 (ASPRV1). <i>Biomolecules</i> , 2020, 10, 1004.	4.0	4
16	Identification of Host Cellular Protein Substrates of SARS-COV-2 Main Protease. <i>International Journal of Molecular Sciences</i> , 2020, 21, 9523.	4.1	22
17	Tandem Mass-Remainder Analysis of Industrially Important Polyether Polyols. <i>Polymers</i> , 2020, 12, 2768.	4.5	3
18	Encoding Information into Polyethylene Glycol Using an Alcohol-Isocyanate "Click" Reaction. <i>International Journal of Molecular Sciences</i> , 2020, 21, 1318.	4.1	6

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19	Multistep Mass-Remainder Analysis and its Application in Copolymer Blends. <i>Macromolecules</i> , 2020, 53, 1199-1204.	4.8	8
20	Sequential Oscillations in the Ferrocene-Catalyzed Belousov-Zhabotinsky Reaction: The Case of Oxalacetic Acid Substrate. <i>ChemistrySelect</i> , 2019, 4, 451-456.	1.5	0
21	Spilanthol Inhibits Inflammatory Transcription Factors and iNOS Expression in Macrophages and Exerts Anti-inflammatory Effects in Dermatitis and Pancreatitis. <i>International Journal of Molecular Sciences</i> , 2019, 20, 4308.	4.1	20
22	Improved Modeling of Peptidic Foldamers Using a Quantum Chemical Parametrization Based on Torsional Minimum Energy Path Matching. <i>ChemPlusChem</i> , 2019, 84, 927-941.	2.8	5
23	Mass Spectrometric Characterization of Epoxidized Vegetable Oils. <i>Polymers</i> , 2019, 11, 394.	4.5	6
24	Mass-Remainder Analysis (MARA): An Improved Method for Elemental Composition Assignment in Petroleomics. <i>Analytical Chemistry</i> , 2019, 91, 6479-6486.	6.5	10
25	Isocyanonaphthalenes as extremely low molecular weight, selective, ratiometric fluorescent probes for Mercury(II). <i>Talanta</i> , 2019, 201, 165-173.	5.5	21
26	Synthesis and supramolecular assembly of fluorinated biogenic amine recognition host polymers. <i>Polymer Chemistry</i> , 2019, 10, 5626-5634.	3.9	8
27	Rapid mapping of various chemicals in personal care and healthcare products by direct analysis in real time mass spectrometry. <i>Talanta</i> , 2019, 192, 241-247.	5.5	11
28	Mass-Remainder Analysis (MARA): a New Data Mining Tool for Copolymer Characterization. <i>Analytical Chemistry</i> , 2018, 90, 3892-3897.	6.5	21
29	A general formulation of the quasiclassical trajectory method for reduced-dimensionality reaction dynamics calculations. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 13224-13240.	2.8	4
30	Rapid qualitative analysis of 2 flavonoids, rutin and silybin, in medical pills by direct analysis in real-time mass spectrometry (<sc>DART-MS</sc>) combined with <i>in situ</i> derivatization. <i>Journal of Mass Spectrometry</i> , 2018, 53, 240-246.	1.6	8
31	One-pot Synthesis of 1,3-Butadiene and 1,6-Hexanediol Derivatives from Cyclopentadiene (CPD) via Tandem Olefin Metathesis Reactions. <i>ChemCatChem</i> , 2018, 10, 4870-4877.	3.7	1
32	Reply to Comment on "Mass-Remainder Analysis (MARA): A New Data Mining Tool for Copolymer Characterization" (An Example of Multiple Discovery). <i>Analytical Chemistry</i> , 2018, 90, 8719-8720.	6.5	2
33	Screening of additives and other chemicals in polyurethanes by direct analysis in real time mass spectrometry (DART-MS). <i>Analytical and Bioanalytical Chemistry</i> , 2017, 409, 6149-6162.	3.7	14
34	Adiabatic Switching Extended To Prepare Semiclassically Quantized Rotational "Vibrational Initial States for Quasiclassical Trajectory Calculations. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 4621-4626.	4.6	29
35	Kinetics of Uncatalyzed Reactions of 2,4- and 4,4-Diphenylmethane-Diisocyanate with Primary and Secondary Alcohols. <i>International Journal of Chemical Kinetics</i> , 2017, 49, 643-655.	1.6	13
36	Reactions of 2,6-Toluene Diisocyanate with Alcohols: Kinetic Studies in the Absence and Presence of Catalysts. <i>ChemistrySelect</i> , 2017, 2, 11302-11306.	1.5	1

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37	WACCM-rSIC. Geoscientific Model Development, 2016, 9, 3123-3136.	3.6	16
38	Oscillatory reaction cross sections caused by normal mode sampling in quasiclassical trajectory calculations. Journal of Chemical Physics, 2016, 144, 014104.	3.0	12
39	Uncatalyzed reactions of 4,4-diphenylmethane-diisocyanate with polymer polyols as revealed by matrix-assisted laser desorption/ionization mass spectrometry. RSC Advances, 2016, 6, 47023-47032.	3.6	7
40	Rapid detection of hazardous chemicals in textiles by direct analysis in real-time mass spectrometry (DART-MS). Analytical and Bioanalytical Chemistry, 2016, 408, 5189-5198.	3.7	30
41	Development of a Joint Hydrogen and Syngas Combustion Mechanism Based on an Optimization Approach. International Journal of Chemical Kinetics, 2016, 48, 407-422.	1.6	122
42	Can Nonpolar Polyisobutylenes be Measured by Electrospray Ionization Mass Spectrometry? Anion-Attachment Proved to be an Appropriate Method. Journal of the American Society for Mass Spectrometry, 2016, 27, 432-442.	2.8	4
43	Testing the Palma-Clary Reduced Dimensionality Model Using Classical Mechanics on the $\text{CH}_4 + \text{H} \rightarrow \text{CH}_3 + \text{H}_2$ Reaction. Journal of Physical Chemistry A, 2016, 120, 5083-5093.	2.5	7
44	HSO_3Cl : a prototype molecule for studying OH-stretching overtone induced photodissociation. Physical Chemistry Chemical Physics, 2016, 18, 6780-6788.	2.8	15
45	Following the molecular motion of near-resonant excited CO on Pt(111): A simulated x-ray photoelectron diffraction study based on molecular dynamics calculations. Structural Dynamics, 2015, 2, 035102.	2.3	6
46	Electrospray ionization tandem mass spectrometry of the star-shaped propoxylated diethylenetriamine polyols. Journal of Mass Spectrometry, 2015, 50, 914-917.	1.6	4
47	New insight into the kinetics of diisocyanate-alcohol reactions by high-performance liquid chromatography and mass spectrometry. Journal of Applied Polymer Science, 2015, 132, .	2.6	21
48	Direct analysis in real time mass spectrometry (DART-MS) of highly nonpolar low molecular weight polyisobutylenes. Journal of Mass Spectrometry, 2015, 50, 1071-1078.	1.6	7
49	Chiral differentiation of the noscapine and hydrastine stereoisomers by electrospray ionization tandem mass spectrometry. Journal of Mass Spectrometry, 2015, 50, 240-246.	1.6	9
50	Uncertainty of the rate parameters of several important elementary reactions of the H_2 and syngas combustion systems. Combustion and Flame, 2015, 162, 2059-2076.	5.2	55
51	Histopathological alterations and oxidative stress in liver and kidney of <i>Leuciscus cephalus</i> following exposure to heavy metals in the Tur River, North Western Romania. Ecotoxicology and Environmental Safety, 2015, 119, 198-205.	6.0	71
52	Computational study of collisions between $\text{O}(^3\text{P})$ and $\text{NO}(^2\text{I})$ at temperatures relevant to the hypersonic flight regime. Journal of Chemical Physics, 2014, 141, 164319.	3.0	34
53	Multisurface Adiabatic Reactive Molecular Dynamics. Journal of Chemical Theory and Computation, 2014, 10, 1366-1375.	5.3	60
54	Kinetic isotope effect in malonaldehyde determined from path integral Monte Carlo simulations. Physical Chemistry Chemical Physics, 2014, 16, 204-211.	2.8	35

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55	Competitive reaction pathways in vibrationally induced photodissociation of H_2SO_4 . <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 18533.	2.8	17
56	Comparison of the performance of several recent hydrogen combustion mechanisms. <i>Combustion and Flame</i> , 2014, 161, 2219-2234.	5.2	144
57	Arm length distribution in four arm star shaped propoxylated ethylenediamine polyol by tandem mass spectrometry. <i>Journal of Mass Spectrometry</i> , 2013, 48, 1125-1127.	1.6	2
58	Determination of the uncertainty domain of the Arrhenius parameters needed for the investigation of combustion kinetic models. <i>Reliability Engineering and System Safety</i> , 2012, 107, 29-34.	8.9	50
59	State-selected ion-molecule reactions with Coulomb-crystallized molecular ions in traps. <i>Chemical Physics Letters</i> , 2012, 547, 1-8.	2.6	39
60	A shock tube and chemical kinetic modeling study of the pyrolysis and oxidation of butanols. <i>Combustion and Flame</i> , 2012, 159, 2009-2027.	5.2	87
61	Uncertainty of Arrhenius parameters. <i>International Journal of Chemical Kinetics</i> , 2011, 43, 359-378.	1.6	96
62	Uncertainty analysis of varying temperature chemical kinetic systems. <i>Procedia, Social and Behavioral Sciences</i> , 2010, 2, 7757-7758.	0.5	1
63	Relaxation of concentration perturbation in chemical kinetic systems. <i>Reaction Kinetics and Catalysis Letters</i> , 2009, 96, 269-278.	0.6	4
64	Reduction of very large reaction mechanisms using methods based on simulation error minimization. <i>Combustion and Flame</i> , 2009, 156, 417-428.	5.2	131