

Florence Vermeire

List of Publications by Year in descending order

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34
papers

1,603
citations

516215

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395343

33
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docs citations

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times ranked

1077
citing authors

#	ARTICLE	IF	CITATIONS
1	Group Contribution and Machine Learning Approaches to Predict Abraham Solute Parameters, Solvation Free Energy, and Solvation Enthalpy. <i>Journal of Chemical Information and Modeling</i> , 2022, 62, 433-446.	2.5	59
2	Detailed Kinetic Modeling for the Pyrolysis of a Jet A Surrogate. <i>Energy & Fuels</i> , 2022, 36, 1304-1315.	2.5	11
3	Variable pressure JSR study of low temperature oxidation chemistry of n-heptane by synchrotron photoionization mass spectrometry. <i>Combustion and Flame</i> , 2022, 240, 111946.	2.8	7
4	Analytics Driving Kinetics: Advanced Mass Spectrometric Characterization of Petroleum Products. <i>Energy & Fuels</i> , 2022, 36, 6-59.	2.5	10
5	Machine Learning for Physicochemical Property Prediction of Complex Hydrocarbon Mixtures. <i>Industrial & Engineering Chemistry Research</i> , 2022, 61, 8581-8594.	1.8	17
6	Predicting Solubility Limits of Organic Solutes for a Wide Range of Solvents and Temperatures. <i>Journal of the American Chemical Society</i> , 2022, 144, 10785-10797.	6.6	31
7	Combustion of ethylamine, dimethylamine and diethylamine: Theoretical and kinetic modeling study. <i>Proceedings of the Combustion Institute</i> , 2021, 38, 585-592.	2.4	12
8	Thermal decomposition of furans with oxygenated substituents: A combined experimental and quantum chemical study. <i>Proceedings of the Combustion Institute</i> , 2021, 38, 699-707.	2.4	9
9	Bond additivity corrections for CBSâ€³ calculated standard enthalpies of formation of H, C, O, N, and S containing species. <i>International Journal of Chemical Kinetics</i> , 2021, 53, 345-355.	1.0	5
10	Transfer learning for solvation free energies: From quantum chemistry to experiments. <i>Chemical Engineering Journal</i> , 2021, 418, 129307.	6.6	77
11	Detailed experimental and kinetic modeling study of 3â€arene pyrolysis. <i>International Journal of Chemical Kinetics</i> , 2020, 52, 785-795.	1.0	4
12	Experimental and kinetic modeling study of the pyrolysis and oxidation of diethylamine. <i>Fuel</i> , 2020, 275, 117744.	3.4	11
13	The thermal decomposition of furfural: molecular chemistry unraveled. <i>Proceedings of the Combustion Institute</i> , 2019, 37, 445-452.	2.4	16
14	QUANTIS: Data quality assessment tool by clustering analysis. <i>International Journal of Chemical Kinetics</i> , 2019, 51, 872-885.	1.0	7
15	A first evaluation of butanoic and pentanoic acid oxidation kinetics. <i>Chemical Engineering Journal</i> , 2019, 373, 973-984.	6.6	27
16	Exploring hydroperoxides in combustion: History, recent advances and perspectives. <i>Progress in Energy and Combustion Science</i> , 2019, 73, 132-181.	15.8	119
17	Experimental and modeling study of the pyrolysis and combustion of dimethoxymethane. <i>Combustion and Flame</i> , 2018, 190, 270-283.	2.8	78
18	A study of thermal decomposition of bromoethane. <i>Journal of Analytical and Applied Pyrolysis</i> , 2018, 136, 199-207.	2.6	7

#	ARTICLE	IF	CITATIONS
19	Experimental and Kinetic Modeling Study of Cyclohexane Pyrolysis. <i>Energy & Fuels</i> , 2018, 32, 7153-7168.	2.5	12
20	An evaluation of the impact of SG1 disproportionation and the addition of styrene in NMP of methyl methacrylate. <i>AIChE Journal</i> , 2018, 64, 2545-2559.	1.8	15
21	Group additive modeling of cyclopentane pyrolysis. <i>Journal of Analytical and Applied Pyrolysis</i> , 2017, 128, 437-450.	2.6	12
22	Experimental and kinetic modeling study of the pyrolysis and oxidation of 1,5-hexadiene: The reactivity of allylic radicals and their role in the formation of aromatics. <i>Fuel</i> , 2017, 208, 779-790.	3.4	17
23	Steam cracking of bio-derived normal and branched alkanes: Influence of branching on product distribution and formation of aromatics. <i>Journal of Analytical and Applied Pyrolysis</i> , 2016, 122, 468-478.	2.6	8
24	Experimental and Kinetic Modeling Study of 2-Methyl-2-Butene: Allylic Hydrocarbon Kinetics. <i>Journal of Physical Chemistry A</i> , 2015, 119, 7462-7480.	1.1	62
25	Experimental and Modeling Investigation of the Low-Temperature Oxidation of Dimethyl Ether. <i>Journal of Physical Chemistry A</i> , 2015, 119, 7905-7923.	1.1	85
26	Progress in Understanding Low-Temperature Organic Compound Oxidation Using a Jet-Stirred Reactor. <i>International Journal of Chemical Kinetics</i> , 2014, 46, 619-639.	1.0	80
27	An experimental and modeling study of propene oxidation. Part 1: Speciation measurements in jet-stirred and flow reactors. <i>Combustion and Flame</i> , 2014, 161, 2765-2784.	2.8	251
28	An experimental and modeling study of the low- and high-temperature oxidation of cyclohexane. <i>Combustion and Flame</i> , 2013, 160, 2319-2332.	2.8	71
29	Low temperature oxidation of benzene and toluene in mixture with n-decane. <i>Proceedings of the Combustion Institute</i> , 2013, 34, 297-305.	2.4	42
30	Towards cleaner combustion engines through groundbreaking detailed chemical kinetic models. <i>Chemical Society Reviews</i> , 2011, 40, 4762.	18.7	111
31	Experimental Confirmation of the Low-Temperature Oxidation Scheme of Alkanes. <i>Angewandte Chemie - International Edition</i> , 2010, 49, 3169-3172.	7.2	180
32	Thermal Decomposition of Norbornane (bicyclo[2.2.1]heptane) Dissolved in Benzene: An Experimental Study and Mechanism Investigation. <i>Energy & Fuels</i> , 2007, 21, 1406-1414.	2.5	7
33	Thermal decomposition of n-dodecane: Experiments and kinetic modeling. <i>Journal of Analytical and Applied Pyrolysis</i> , 2007, 78, 419-429.	2.6	138
34	The merit of pressure dependent kinetic modelling in steam cracking. <i>Faraday Discussions</i> , 0, 238, 491-511.	1.6	5