

John C Mackie

List of Publications by Year in descending order

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papers

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126907

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136
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2594
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#	ARTICLE	IF	CITATIONS
1	Modeling and Experimental Study on the Thermal Decomposition of Perfluorooctanesulfonic Acid (PFOS) in an γ -Alumina Reactor. <i>Industrial & Engineering Chemistry Research</i> , 2022, 61, 5453-5463.	3.7	7
2	Kinetics of Decomposition of PFOS Relevant to Thermal Desorption Remediation of Soils. <i>Industrial & Engineering Chemistry Research</i> , 2021, 60, 9080-9087.	3.7	11
3	The Catalyzed Conversion of Methane to Value-Added Products. <i>Energy Technology</i> , 2020, 8, 1900665.	3.8	13
4	Study on Catalyst Deactivation During the Hydrodeoxygenation of Model Compounds. <i>Topics in Catalysis</i> , 2020, 63, 778-792.	2.8	16
5	Products and mechanism of thermal decomposition of chlorpyrifos under inert and oxidative conditions. <i>Environmental Sciences: Processes and Impacts</i> , 2020, 22, 2084-2094.	3.5	8
6	Mechanisms of thermal decomposition of cyclodiene pesticides, identification and possible mitigation of their toxic products. <i>Proceedings of the Combustion Institute</i> , 2019, 37, 1143-1150.	3.9	5
7	Formation of polychlorinated dibenzo-p-dioxins and dibenzofurans (PCDD/F) from oxidation of 4,4'-dichlorobiphenyl (4,4'-DCB). <i>Proceedings of the Combustion Institute</i> , 2019, 37, 1075-1082.	3.9	12
8	Thermal oxidation of dieldrin and concomitant formation of toxic products including polychlorinated dibenzo-p-dioxin and dibenzofuran (PCDD/F). <i>Chemosphere</i> , 2019, 225, 209-216.	8.2	5
9	Pyrolysis of Glyphosate and Its Toxic Products. <i>Environmental Science & Technology</i> , 2019, 53, 13742-13747.	10.0	13
10	Mechanism of the Thermal Decomposition of Chlorpyrifos and Formation of the Dioxin Analog, 2,3,7,8-Tetrachloro-1,4-dioxino-dipyridine (TCDDpy). <i>Environmental Science & Technology</i> , 2018, 52, 7327-7333.	10.0	17
11	Process for Chloroform Decomposition: Nonthermal Plasma Polymerization with Methane and Hydrogen. <i>Industrial & Engineering Chemistry Research</i> , 2018, 57, 9075-9082.	3.7	1
12	Mechanism and Rate of Thermal Decomposition of Hexachlorocyclopentadiene and Its Importance in PCDD/F Formation from the Combustion of Cyclodiene Pesticides. <i>Journal of Physical Chemistry A</i> , 2017, 121, 5871-5883.	2.5	8
13	Gas phase pyrolysis of endosulfan and formation of dioxin precursors of polychlorinated dibenzo-p-dioxins and dibenzofurans (PCDD/F). <i>Proceedings of the Combustion Institute</i> , 2017, 36, 1119-1127.	3.9	5
14	Oxidation of 4-bromo-4'-chlorobiphenyl, model species for forming mixed halogenated aromatic compounds. <i>International Journal of Environment and Pollution</i> , 2017, 61, 243.	0.2	4
15	Experimental investigation of the reaction of HCFC-22 and methane in a dielectric barrier discharge non-equilibrium plasma. <i>Chemical Engineering Journal</i> , 2016, 301, 73-82.	12.7	4
16	Gas Phase Thermal Oxidation of Endosulfan and Formation of Polychlorinated Dibenzodioxins and Dibenzofurans. <i>Environmental Science & Technology</i> , 2016, 50, 10106-10113.	10.0	6
17	Reaction of dichloromethane under non-oxidative conditions in a dielectric barrier discharge reactor and characterisation of the resultant polymer. <i>Chemical Engineering Journal</i> , 2016, 290, 499-506.	12.7	4
18	Experimental Study on the Reaction of CCl_3F and CH_4 in a Dielectric Barrier Discharge Nonequilibrium Plasma Reactor. <i>Industrial & Engineering Chemistry Research</i> , 2016, 55, 463-471.	3.7	2

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19	Effect of methane on the conversion of HFC-134a in a dielectric barrier discharge non-equilibrium plasma reactor. <i>Chemical Engineering Journal</i> , 2016, 284, 412-421.	12.7	8
20	Thermodynamic stability and structure of cuprous chloride surfaces: a DFT investigation. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 7038-7045.	2.8	8
21	<i>S</i> -Nitrosation of Aminothiones. <i>Journal of Organic Chemistry</i> , 2015, 80, 6951-6958.	3.2	1
22	Characterization of Polymer Synthesized from the Nonequilibrium Plasma Conversion of CFC-12 and Methane in a Dielectric Barrier Discharge Reactor. <i>Industrial & Engineering Chemistry Research</i> , 2014, 53, 19380-19386.	3.7	4
23	Water formation via HCl oxidation on Cu(100). <i>Applied Surface Science</i> , 2014, 299, 156-161.	6.1	5
24	Nonequilibrium Plasma Polymerization of HFC-134a in a Dielectric Barrier Discharge Reactor: Polymer Characterization and a Proposed Mechanism for Polymer Formation. <i>IEEE Transactions on Plasma Science</i> , 2014, 42, 3095-3100.	1.3	7
25	Reaction of carbon tetrachloride with methane in a non-equilibrium plasma at atmospheric pressure, and characterisation of the polymer thus formed. <i>Journal of Hazardous Materials</i> , 2014, 280, 38-45.	12.4	3
26	Comparative Study on the Formation of Toxic Species from 4-chlorobiphenyl in Fires: Effect of Catalytic Surfaces. <i>Procedia Engineering</i> , 2013, 62, 350-358.	1.2	6
27	Study on the Reaction of CCl_2F_2 with CH_4 in a Dielectric Barrier Discharge Nonequilibrium Plasma. <i>Plasma Processes and Polymers</i> , 2013, 10, n/a-n/a.	3.0	2
28	Reaction of chloroform in a non-oxidative atmosphere using dielectric barrier discharge. , 2013, , .		0
29	Roles of peroxides and unsaturation in spontaneous heating of linseed oil. <i>Fire Safety Journal</i> , 2013, 61, 108-115.	3.1	9
30	Formation of polychlorinated dibenzo-p-dioxins and polychlorinated dibenzofurans (PCDD/F) by precursor pathways in oxidation of pesticide alpha-cypermethrin. <i>Proceedings of the Combustion Institute</i> , 2013, 34, 3499-3507.	3.9	7
31	Comparative Study of the Physicochemical Properties of Ortho-Substituted Aromatic Nitroso Compounds. <i>Journal of Chemical & Engineering Data</i> , 2013, 58, 1005-1010.	1.9	5
32	Rate constants for reactions of ethylbenzene with hydroperoxyl radical. <i>Combustion and Flame</i> , 2013, 160, 9-16.	5.2	30
33	A Melamine-Modified Zeolite with Enhanced CO_2 Capture Properties. <i>Energy Technology</i> , 2013, 1, 345-349.	3.8	18
34	Trapping of Nitric Oxide, Generated during Sensitization of Ammonium Nitrate Emulsion Explosive, by Aromatic Nitroso Sulfonates. <i>Industrial & Engineering Chemistry Research</i> , 2013, 52, 10561-10568.	3.7	2
35	Non-thermal plasma polymerization of HFC-134A in a dielectric barrier discharge reactor; Polymer characterization and a proposed mechanism for polymer formation. , 2013, , .		0
36	Non-thermal plasma polymerization of HFC-134a in a dielectric barrier discharge reactor; Polymer characterization and understanding the mechanism of polymer formation. , 2013, , .		1

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37	Non-Oxidative Conversion of 1,2-Dichloroethane in a Non-Thermal Plasma and Characterisation of the Polymer Formed. <i>Plasma Processes and Polymers</i> , 2013, 10, 141-149.	3.0	7
38	Accurate Rate Constants for Decomposition of Aqueous Nitrous Acid. <i>Inorganic Chemistry</i> , 2012, 51, 2178-2185.	4.0	40
39	Mechanistic Study of Trapping of NO by 3,5-Dibromo-4-Nitrosobenzene Sulfonate. <i>Industrial & Engineering Chemistry Research</i> , 2012, 51, 14325-14336.	3.7	3
40	Conversion of Fluorine-Containing Ozone-Depleting and Greenhouse Gases to Valuable Polymers in a Nonthermal Plasma. <i>Industrial & Engineering Chemistry Research</i> , 2012, 51, 11279-11283.	3.7	20
41	Mechanism of Formation of Volatile Organic Compounds from Oxidation of Linseed Oil. <i>Industrial & Engineering Chemistry Research</i> , 2012, 51, 5653-5661.	3.7	22
42	Identification and Quantitation of Volatile Organic Compounds from Oxidation of Linseed Oil. <i>Industrial & Engineering Chemistry Research</i> , 2012, 51, 5645-5652.	3.7	32
43	Oxidation of dibenzo-p-dioxin: Formation of initial products, 2-methylbenzofuran and 3-hydro-2-methylenebenzofuran. <i>Combustion and Flame</i> , 2012, 159, 3056-3065.	5.2	7
44	Low temperature oxidation of linseed oil: a review. <i>Fire Science Reviews</i> , 2012, 1, .	0.9	80
45	Determination of toxic products released in combustion of pesticides. <i>Progress in Energy and Combustion Science</i> , 2012, 38, 400-418.	31.2	24
46	An equilibrium ab initio atomistic thermodynamics study of chlorine adsorption on the Cu(001) surface. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 10306.	2.8	25
47	Air Pollutants Formed in Thermal Decomposition of Folpet Fungicide under Oxidative Conditions. <i>Environmental Science & Technology</i> , 2011, 45, 554-560.	10.0	11
48	Chlorination of the Cu(110) Surface and Copper Nanoparticles: A Density Functional Theory Study. <i>Journal of Physical Chemistry C</i> , 2011, 115, 13412-13419.	3.1	25
49	Experimental Study of Decomposition of Aqueous Nitrosyl Thiocyanate. <i>Inorganic Chemistry</i> , 2011, 50, 7440-7452.	4.0	10
50	Formation of toxic species and precursors of PCDD/F in thermal decomposition of alpha-cypermethrin. <i>Chemosphere</i> , 2011, 85, 143-150.	8.2	9
51	Oxidation reactions and spontaneous ignition of linseed oil. <i>Proceedings of the Combustion Institute</i> , 2011, 33, 2625-2632.	3.9	18
52	Rate constants for hydrogen abstraction reactions by the hydroperoxyl radical from methanol, ethenol, acetaldehyde, toluene, and phenol. <i>Journal of Computational Chemistry</i> , 2011, 32, 1725-1733.	3.3	43
53	Mechanistic study of the reaction of CHF ₃ with CH ₄ . <i>Chemical Engineering Journal</i> , 2011, 166, 822-831.	12.7	18
54	Quantum chemical study of copper (II) chloride and the Deacon reaction. <i>Chemical Physics Letters</i> , 2011, 501, 215-220.	2.6	12

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55	Toxic pollutants emitted from thermal decomposition of phthalimide compounds. Journal of Hazardous Materials, 2011, 187, 407-412.	12.4	7
56	Formation of polychlorinated dibenzo-p-dioxins and dibenzofurans (PCDD/F) in oxidation of captan pesticide. Proceedings of the Combustion Institute, 2011, 33, 701-708.	3.9	24
57	Catalytic pyrolysis of CHF ₃ over activated carbon and activated carbon supported potassium catalyst. Journal of Fluorine Chemistry, 2010, 131, 698-703.	1.7	16
58	Conversion of a CFCs, HFCs and HCFCs waste mixture via reaction with methane. Journal of Hazardous Materials, 2010, 184, 696-703.	12.4	7
59	Experimental and chemical kinetic study of the pyrolysis of trifluoroethane and the reaction of trifluoromethane with methane. Journal of Fluorine Chemistry, 2010, 131, 751-760.	1.7	38
60	Conversion of CHF ₃ to CH ₂ CF ₂ via reaction with CH ₄ in the presence of CBrF ₃ : An experimental and kinetic modelling study. Journal of Hazardous Materials, 2010, 180, 181-187.	12.4	19
61	A DFT study on the self-coupling reactions of the three isomeric semiquinone radicals. Computational and Theoretical Chemistry, 2010, 958, 106-115.	1.5	9
62	Adsorption of 2-chlorophenol on Cu ₂ O(111)â€“CuCUS: A first-principles density functional study. Applied Surface Science, 2010, 256, 4764-4770.	6.1	8
63	Theoretical study of reactions of HO ₂ in low-temperature oxidation of benzene. Combustion and Flame, 2010, 157, 1325-1330.	5.2	17
64	Theoretical Study of Unimolecular Decomposition of Catechol. Journal of Physical Chemistry A, 2010, 114, 1060-1067.	2.5	44
65	Effect of Methanol on the Gas-Phase Reaction of Trifluoromethane with Methane. Industrial & Engineering Chemistry Research, 2010, 49, 8406-8414.	3.7	11
66	Thermal Decomposition of Captan and Formation Pathways of Toxic Air Pollutants. Environmental Science & Technology, 2010, 44, 4149-4154.	10.0	17
67	Synthesis of Vinylidene Fluoride via Reaction of Chlorodifluoromethane (HCFC-22) with Methane. Industrial & Engineering Chemistry Research, 2010, 49, 6010-6019.	3.7	5
68	Theoretical Study of the Ammoniaâˆ“Hypochlorous Acid Reaction Mechanism. Journal of Physical Chemistry A, 2010, 114, 2597-2606.	2.5	39
69	Theoretical Study on the Thermodynamic Properties and Self-Decomposition of Methylbenzenediol Isomers. Journal of Physical Chemistry A, 2010, 114, 11751-11760.	2.5	4
70	Thermochemical Properties and Decomposition Pathways of Three Isomeric Semiquinone Radicals. Journal of Physical Chemistry A, 2010, 114, 1098-1108.	2.5	36
71	Interaction of Chlorine and Oxygen with the Cu(100) Surface. Journal of Physical Chemistry C, 2010, 114, 19048-19054.	3.1	19
72	A first-principles density functional study of chlorophenol adsorption on Cu ₂ O(110):CuO. Journal of Chemical Physics, 2009, 130, 184505.	3.0	30

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73	Mechanisms for formation, chlorination, dechlorination and destruction of polychlorinated dibenzo-p-dioxins and dibenzofurans (PCDD/Fs). <i>Progress in Energy and Combustion Science</i> , 2009, 35, 245-274.	31.2	401
74	Pyrolysis of permethrin and formation of precursors of polychlorinated dibenzo-p-dioxins and dibenzofurans (PCDD/F) under non-oxidative conditions. <i>Chemosphere</i> , 2009, 74, 1435-1443.	8.2	32
75	Adsorption of chlorophenol on the Cu(111) surface: A first-principles density functional theory study. <i>Applied Surface Science</i> , 2008, 254, 4218-4224.	6.1	27
76	2-Chlorophenol adsorption on Cu(100): First-principles density functional study. <i>Surface Science</i> , 2008, 602, 1554-1562.	1.9	14
77	Experimental and Kinetic Studies of Gas-phase Pyrolysis of $C_{10}H_8Cl_4$. <i>Industrial & Engineering Chemistry Research</i> , 2008, 47, 2579-2584.	3.7	16
78	Conversion of CHF_3 to CH_2F_2 via Reaction with CH_4 and $CaBr_2$. <i>Environmental Science & Technology</i> , 2008, 42, 5795-5799.	10.0	23
79	Computational Study of the Oxidation and Decomposition of Dibenzofuran under Atmospheric Conditions. <i>Journal of Physical Chemistry A</i> , 2008, 112, 6960-6967.	2.5	30
80	Quantum Chemical and Kinetic Study of Formation of 2-Chlorophenoxy Radical from 2-Chlorophenol: Unimolecular Decomposition and Bimolecular Reactions with H, OH, Cl, and O_2 . <i>Journal of Physical Chemistry A</i> , 2008, 112, 3680-3692.	2.5	34
81	Simultaneous conversion of $CHClF_2$ and CH_3Br to CH_2CF_2 . <i>Chemosphere</i> , 2007, 68, 2003-2006.	8.2	10
82	Catalytic Effect of CuO and Other Transition Metal Oxides in Formation of Dioxins: Theoretical Investigation of Reaction Between 2,4,5-Trichlorophenol and CuO. <i>Environmental Science & Technology</i> , 2007, 41, 5708-5715.	10.0	36
83	Quantum Chemical Investigation of Formation of Polychlorodibenzo-p-Dioxins and Dibenzofurans from Oxidation and Pyrolysis of 2-Chlorophenol. <i>Journal of Physical Chemistry A</i> , 2007, 111, 2563-2573.	2.5	73
84	Theoretical Study of Reaction Pathways of Dibenzofuran and Dibenzo-p-Dioxin under Reducing Conditions. <i>Journal of Physical Chemistry A</i> , 2007, 111, 7133-7140.	2.5	30
85	An Experimental and Kinetic Modeling Study of the Reaction of CHF_3 with Methane. <i>Environmental Science & Technology</i> , 2006, 40, 5778-5785.	10.0	27
86	Quantum Chemical Study of Low Temperature Oxidation Mechanism of Dibenzofuran. <i>Journal of Physical Chemistry A</i> , 2006, 110, 13560-13567.	2.5	29
87	Experimental and Quantum Chemical Study of the Reaction $CF_2 + CH_3 \rightarrow CF_2CH_3 + H$: A Key Mechanism in the Reaction between Methane and Fluorocarbons. <i>Industrial & Engineering Chemistry Research</i> , 2006, 45, 3758-3762.	3.7	15
88	Oxidation of CO by SO_2 : A Theoretical Study. <i>ChemInform</i> , 2005, 36, no.	0.0	1
89	Oxidation of CO by SO_2 : A Theoretical Study. <i>Journal of Physical Chemistry A</i> , 2005, 109, 2019-2025.	2.5	19
90	Quantum Chemical Study of the Mechanism of Reaction between $NH(X^3\Sigma^-)$ and H_2 , H_2O , and CO_2 under Combustion Conditions. <i>Journal of Physical Chemistry A</i> , 2005, 109, 11967-11974.	2.5	35

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91	An ab initio Quantum Chemical and Kinetic Study of the NNH + O Reaction Potential Energy Surface: How Important Is this Route to NO in Combustion?. ChemInform, 2003, 34, no.	0.0	0
92	An Ab Initio Quantum Chemical and Kinetic Study of the NNH + O Reaction Potential Energy Surface: How Important Is This Route to NO in Combustion?. Journal of Physical Chemistry A, 2003, 107, 6792-6803.	2.5	40
93	Reactions of Phosphorus-Containing Species of Importance in the Catalytic Recombination of H + OH: A Quantum Chemical and Kinetic Study. Journal of Physical Chemistry A, 2002, 106, 10825-10830.	2.5	16
94	The Role of Phosphorus Dioxide in the H + OH Recombination Reaction: Ab Initio Quantum Chemical Computation of Thermochemical and Rate Parameters. Journal of Physical Chemistry A, 2002, 106, 1533-1541.	2.5	21
95	The pyrolysis of cyclopentadiene: quantum chemical and kinetic modelling studies of the acetylene plus propyne/allene decomposition channels. Physical Chemistry Chemical Physics, 2001, 3, 2467-2473.	2.8	39
96	A Study of Furan as a Model Oxygenated Reburn Fuel for Nitric Oxide Reduction. Energy & Fuels, 2001, 15, 743-750.	5.1	9
97	A kinetic study of the oxidation of pyridine. Proceedings of the Combustion Institute, 2000, 28, 1709-1716.	3.9	36
98	An experimental and kinetic modeling study of the reduction of NO by coal volatiles in a flow reactor. Proceedings of the Combustion Institute, 2000, 28, 2345-2351.	3.9	7
99	Heats of Formation of Hydrofluorocarbons Obtained by Gaussian-3 and Related Quantum Chemical Computations. Journal of Physical Chemistry A, 2000, 104, 7600-7611.	2.5	60
100	Pyrolysis of Furan: Ab Initio Quantum Chemical and Kinetic Modeling Studies. Journal of Physical Chemistry A, 2000, 104, 1861-1875.	2.5	102
101	The thermal decomposition of pyrrole: an ab initio quantum chemical study of the potential energy surface associated with the hydrogen cyanide plus propyne channel. Chemical Physics Letters, 1999, 300, 321-330.	2.6	46
102	Shock-Tube Study of the Pyrolysis of the Halon Replacement Molecule CF ₃ CHF ₂ CF ₃ . Journal of Physical Chemistry A, 1999, 103, 54-61.	2.5	52
103	Shock Tube Study of the Oxidation of C ₃ F ₆ by N ₂ O. Journal of Physical Chemistry A, 1999, 103, 5967-5977.	2.5	13
104	Ab Initio Quantum Chemical and Experimental (Shock Tube) Studies of the Pyrolysis Kinetics of Acetonitrile. Journal of Physical Chemistry A, 1999, 103, 1054-1072.	2.5	31
105	Observation of Cyclopenta-Fused and Ethynyl-Substituted PAH during the Fuel-Rich Combustion of Primary Tar from a Bituminous Coal. Energy & Fuels, 1999, 13, 1167-1172.	5.1	37
106	Ab Initio Quantum Chemical and Kinetic Modeling Study of the Pyrolysis Kinetics of Pyrrole. Journal of Physical Chemistry A, 1999, 103, 3923-3934.	2.5	69
107	An ab initio quantum chemical study of the electronic structure and stability of the pyrrolyl radical: Comparison with the isoelectronic cyclopentadienyl radical. Chemical Physics Letters, 1998, 290, 391-398.	2.6	42
108	The formation of nitrogen species and oxygenated PAH during the combustion of coal volatiles. Proceedings of the Combustion Institute, 1998, 27, 1687-1693.	0.3	8

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109	Release of HCN, NH ₃ , and HNCO from the Thermal Gas-Phase Cracking of Coal Pyrolysis Tars. <i>Energy & Fuels</i> , 1998, 12, 536-541.	5.1	86
110	Decomposition of the Benzyl Radical: Quantum Chemical and Experimental (Shock Tube) Investigations of Reaction Pathways. <i>Journal of Physical Chemistry A</i> , 1997, 101, 7105-7113.	2.5	59
111	The Pyrolysis of 3-Picoline: Ab Initio Quantum Chemical and Experimental (Shock Tube) Kinetic Studies. <i>Israel Journal of Chemistry</i> , 1996, 36, 239-248.	2.3	6
112	A kinetic study of the oxidation of acetonitrile: A model for NO formation from fuel-bound nitrogen. <i>Proceedings of the Combustion Institute</i> , 1996, 26, 597-604.	0.3	7
113	Thermal decomposition of two coal model compounds – pyridine and 2-picoline. Kinetics and product distributions. <i>Journal of Analytical and Applied Pyrolysis</i> , 1995, 34, 47-63.	5.5	45
114	Ab initio studies of the thermal decomposition of azaaromatics: free radical versus intramolecular mechanism. <i>Journal of the Chemical Society, Faraday Transactions</i> , 1995, 91, 1587.	1.7	22
115	Experimental and ab Initio Theoretical Study of the Kinetics of Rearrangement of Ketene Imine to Acetonitrile. <i>The Journal of Physical Chemistry</i> , 1994, 98, 13546-13555.	2.9	33
116	Experimental and theoretical study of the isomerisation of N-methylpyrrole. <i>Chemical Physics Letters</i> , 1994, 221, 267-273.	2.6	9
117	Kinetics of the thermal decomposition and isomerisation of pyrazine (1,4 diazine). <i>Proceedings of the Combustion Institute</i> , 1994, 25, 893-900.	0.3	6
118	Kinetics of thermal decomposition of the diazines: shock-tube pyrolysis of pyrimidine. <i>Journal of the Chemical Society, Faraday Transactions</i> , 1994, 90, 541.	1.7	25
119	Modelling studies of gas-phase processes in OXCO reactors. <i>Studies in Surface Science and Catalysis</i> , 1994, , 137-142.	1.5	0
120	Kinetics of pyrolysis of the isomeric butenenitriles and kinetic modeling. <i>The Journal of Physical Chemistry</i> , 1992, 96, 272-281.	2.9	24
121	Kinetics of pyrolysis of a coal model compound, 2-picoline, the nitrogen heteroaromatic analogue of toluene. 2. The 2-picoly radical and kinetic modeling. <i>The Journal of Physical Chemistry</i> , 1992, 96, 10339-10348.	2.9	25
122	Kinetics of pyrolysis of a coal model compound, 2-picoline, the nitrogen heteroaromatic analog of toluene. 1. Product distributions. <i>The Journal of Physical Chemistry</i> , 1992, 96, 10334-10339.	2.9	39
123	Kinetics of pyrolysis of furan. <i>Journal of the Chemical Society, Faraday Transactions</i> , 1991, 87, 815.	1.7	63
124	Partial Oxidation of Methane: The Role of the Gas Phase Reactions. <i>Catalysis Reviews - Science and Engineering</i> , 1991, 33, 169-240.	12.9	108
125	Shock tube pyrolysis of pyrrole and kinetic modeling. <i>International Journal of Chemical Kinetics</i> , 1991, 23, 733-760.	1.6	115
126	Partial oxidation of methane by nitrous oxide. <i>Energy & Fuels</i> , 1990, 4, 285-290.	5.1	6

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127	Shock tube pyrolysis of pyridine. The Journal of Physical Chemistry, 1990, 94, 4099-4106.	2.9	151
128	Inhibition of C2 oxidation by methane under oxidative coupling conditions. Energy & Fuels, 1990, 4, 277-285.	5.1	24
129	Pyrolysis of coal at high temperatures. Energy & Fuels, 1988, 2, 391-400.	5.1	56
130	Products from the rapid pyrolysis of a brown coal in inert and reducing atmospheres. Fuel, 1985, 64, 400-405.	6.4	15
131	Products from rapid heating of a brown coal in the temperature range 400–2300 Å°C. Fuel, 1984, 63, 394-400.	6.4	51
132	Homogeneous nucleation of cyclohexane: A double-diaphragm shock tube study. Journal of Colloid and Interface Science, 1981, 83, 547-557.	9.4	2
133	Double-diaphragm shock tube - Comparison between theory and experiment. AIAA Journal, 1981, 19, 405-406.	2.6	2
134	The isomerization of cycloheptatriene at high temperatures. International Journal of Chemical Kinetics, 1976, 8, 695-707.	1.6	15
135	Search for a Charge-Transfer State in Crystalline Anthracene. Journal of Chemical Physics, 1965, 42, 1535-1540.	3.0	61