

Chang Yachao

List of Publications by Year in descending order

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71
papers

3,617
citations

126907

33
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155660

55
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all docs

71
docs citations

71
times ranked

1528
citing authors

#	ARTICLE	IF	CITATIONS
1	Construction and derivation of a series of skeletal chemical mechanisms for n-alkanes with uniform and decoupling structure based on reaction rate rules. <i>Combustion and Flame</i> , 2022, 236, 111785.	5.2	18
2	Collaborative optimization of fuel composition and operating parameters of gasoline compression ignition (GCI) engine in a wide load range. <i>Fuel</i> , 2022, 310, 122366.	6.4	11
3	Construction of reduced chemical mechanisms orientated toward specific applications: a case study of primary reference fuel. <i>Combustion Theory and Modelling</i> , 2022, 26, 560-589.	1.9	4
4	Development of a 5-component gasoline surrogate model using recent advancements in the detailed H ₂ /O ₂ /CO/C ₁ -C ₃ mechanism for decoupling methodology. <i>Fuel</i> , 2021, 283, 118793.	6.4	12
5	Feasibility study of the combustion strategy of n-butanol/diesel dual direct injection (DI2) in a compression-ignition engine. <i>Fuel</i> , 2021, 289, 119865.	6.4	21
6	Novel Method for Quantitative Assessment of Reduced Chemical Mechanisms Based on the Inherent Similarity Evaluation: Case Study of n-Heptane. <i>Energy & Fuels</i> , 2021, 35, 4478-4492.	5.1	3
7	Towards a comprehensive optimization of engine efficiency and emissions by coupling artificial neural network (ANN) with genetic algorithm (GA). <i>Energy</i> , 2021, 225, 120331.	8.8	82
8	Spray-turbulence chemistry interactions under engine-like conditions. <i>Progress in Energy and Combustion Science</i> , 2021, 86, 100939.	31.2	22
9	Construction of reduced oxidation mechanisms of polyoxymethylene dimethyl ethers (PODE ₁₋₆) with consistent structure using decoupling methodology and reaction rate rule. <i>Combustion and Flame</i> , 2021, 232, 111534.	5.2	20
10	Computational optimization of the piston bowl geometry for the different combustion regimes of the dual-mode dual-fuel (DMDF) concept through an improved genetic algorithm. <i>Energy Conversion and Management</i> , 2021, 246, 114658.	9.2	13
11	Mechanism of micro-explosion of water-in-oil emulsified fuel droplet and its effect on soot generation. <i>Energy</i> , 2020, 191, 116488.	8.8	41
12	Reduction of large-scale chemical mechanisms using global sensitivity analysis on reaction class/sub-mechanism. <i>Combustion and Flame</i> , 2020, 212, 355-366.	5.2	29
13	Multiple-objective optimization of methanol/diesel dual-fuel engine at low loads: A comparison of reactivity controlled compression ignition (RCCI) and direct dual fuel stratification (DDFS) strategies. <i>Fuel</i> , 2020, 262, 116673.	6.4	80
14	Construction of a Skeletal Oxidation Mechanism for 2,5-Dimethylfuran Using Decoupling Methodology and Reaction Class-Based Global Sensitivity Analysis. <i>Energy & Fuels</i> , 2020, 34, 16654-16665.	5.1	4
15	Kinetic Modeling Study of Polycyclic Aromatic Hydrocarbon Formation and Oxidation for Oxygenated Fuels including Methanol, n-Butanol, Methyl Butanoate, and Dimethyl Ether. <i>Energy & Fuels</i> , 2020, 34, 4882-4898.	5.1	16
16	Computational optimization of the dual-mode dual-fuel concept through genetic algorithm at different engine loads. <i>Energy Conversion and Management</i> , 2020, 208, 112577.	9.2	20
17	Development of a practical reaction model of polycyclic aromatic hydrocarbon (PAH) formation and oxidation for diesel surrogate fuel. <i>Fuel</i> , 2020, 267, 117159.	6.4	25
18	Influence of the functional group of fuels on the construction of skeletal chemical mechanisms: A case study of 1-hexane, 1-hexene, and 1-hexanol. <i>Combustion and Flame</i> , 2020, 221, 120-135.	5.2	17

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19	Construction and assessment of reduced oxidation mechanisms using global sensitivity analysis and uncertainty analysis. Proceedings of the Combustion Institute, 2019, 37, 751-761.	3.9	16
20	Reduction of Detailed Chemical Mechanisms Using Reaction Class-Based Global Sensitivity and Path Sensitivity Analyses. Energy & Fuels, 2019, 33, 9289-9301.	5.1	4
21	Experimental study on the influence of low-temperature combustion (LTC) mode and fuel properties on cyclic variations in a compression-ignition engine. Fuel, 2019, 256, 115907.	6.4	26
22	Flash boiling explosion model of dimethyl ether (DME)/n-pentane bi-component droplet under gasoline engine conditions. Fuel, 2019, 243, 240-250.	6.4	8
23	Establishment of an improved heat transfer model based on an enhanced thermal wall function for internal combustion engines operated under different combustion modes. Energy Conversion and Management, 2019, 195, 748-759.	9.2	26
24	Numerical analysis of jet breakup based on a modified compressible two-fluid-LES model. Fuel, 2019, 254, 115608.	6.4	9
25	Evaluation of variable compression ratio (VCR) and variable valve timing (VVT) strategies in a heavy-duty diesel engine with reactivity controlled compression ignition (RCCI) combustion under a wide load range. Fuel, 2019, 253, 114-128.	6.4	40
26	Numerical investigation into primary breakup of diesel spray with residual bubbles in the nozzle. Fuel, 2019, 250, 265-276.	6.4	14
27	Understanding of the Ignition Behavior of n-Heptane Spray in Constant Volume Combustion Bombs Focusing on Chemical Kinetics. Energy & Fuels, 2019, 33, 12830-12838.	5.1	2
28	Construction of a decoupling physical-chemical surrogate (DPCS) for practical diesel fuel. Applied Thermal Engineering, 2019, 149, 536-547.	6.0	14
29	Development of a decoupling physical-chemical surrogate (DPCS) model for simulation of the spray and combustion of multi-component biodiesel fuels. Fuel, 2019, 240, 16-30.	6.4	15
30	Comprehensive analysis of exergy destruction sources in different engine combustion regimes. Energy, 2018, 149, 697-708.	8.8	47
31	Principle of determining the optimal operating parameters based on fuel properties and initial conditions for RCCI engines. Fuel, 2018, 216, 284-295.	6.4	41
32	Efficient Approach for the Optimization of Skeletal Chemical Mechanisms with Multiobjective Genetic Algorithm. Energy & Fuels, 2018, 32, 7086-7102.	5.1	32
33	Construction of a skeletal oxidation mechanism of n-pentanol by integrating decoupling methodology, genetic algorithm, and uncertainty quantification. Combustion and Flame, 2018, 194, 15-27.	5.2	45
34	Numerical optimization and comparative study of n-butanol concentration stratification combustion and n-butanol/diesel reactivity stratification combustion for advanced compression ignition (CI) engine. Fuel, 2018, 213, 83-97.	6.4	24
35	An improved method for coupling the in-nozzle cavitation with Multi-fluid-quasi-VOF model for diesel spray. Computers and Fluids, 2018, 177, 20-32.	2.5	29
36	Experimental and modeling study of liquid fuel injection and combustion in diesel engines with a common rail injection system. Applied Energy, 2018, 230, 287-304.	10.1	94

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37	Potential of reactivity controlled compression ignition (RCCI) combustion coupled with variable valve timing (VVT) strategy for meeting Euro 6 emission regulations and high fuel efficiency in a heavy-duty diesel engine. <i>Energy Conversion and Management</i> , 2018, 171, 683-698.	9.2	42
38	Computational optimization of fuel supply, syngas composition, and intake conditions for a syngas/diesel RCCI engine. <i>Fuel</i> , 2018, 234, 120-134.	6.4	49
39	Multi-objective optimization of the combustion of a heavy-duty diesel engine with low temperature combustion under a wide load range: (I) Computational method and optimization results. <i>Energy</i> , 2017, 126, 707-719.	8.8	31
40	Multi-objective optimization of the combustion of a heavy-duty diesel engine with low temperature combustion (LTC) under a wide load range: (II) Detailed parametric, energy, and exergy analysis. <i>Energy</i> , 2017, 139, 247-261.	8.8	23
41	An efficient liquid film vaporization model for multi-component fuels considering thermal and mass diffusions. <i>Applied Thermal Engineering</i> , 2017, 112, 534-548.	6.0	48
42	Simulations of gasoline engine combustion and emissions using a chemical-kinetics-based turbulent premixed combustion modeling approach. <i>Proceedings of the Institution of Mechanical Engineers, Part D: Journal of Automobile Engineering</i> , 2017, 231, 743-765.	1.9	11
43	Construction of a skeletal mechanism for butanol isomers based on the decoupling methodology. <i>Energy Conversion and Management</i> , 2016, 128, 250-260.	9.2	33
44	Thermodynamic energy and exergy analysis of three different engine combustion regimes. <i>Applied Energy</i> , 2016, 180, 849-858.	10.1	117
45	Towards a comprehensive understanding of the influence of fuel properties on the combustion characteristics of a RCCI (reactivity controlled compression ignition) engine. <i>Energy</i> , 2016, 99, 69-82.	8.8	86
46	Development of a quasi-dimensional vaporization model for multi-component fuels focusing on forced convection and high temperature conditions. <i>International Journal of Heat and Mass Transfer</i> , 2016, 97, 130-145.	4.8	51
47	Development of an improved liquid film model for spray/wall interaction under engine-relevant conditions. <i>International Journal of Multiphase Flow</i> , 2016, 79, 74-87.	3.4	59
48	Effect of combustion regime on in-cylinder heat transfer in internal combustion engines. <i>International Journal of Engine Research</i> , 2016, 17, 331-346.	2.3	51
49	Application of the Optimized Decoupling Methodology for the Construction of a Skeletal Primary Reference Fuel Mechanism Focusing on Engine-Relevant Conditions. <i>Frontiers in Mechanical Engineering</i> , 2015, 1, .	1.8	53
50	A reduced toluene reference fuel chemical kinetic mechanism for combustion and polycyclic-aromatic hydrocarbon predictions. <i>Combustion and Flame</i> , 2015, 162, 2390-2404.	5.2	171
51	Construction of Skeletal Oxidation Mechanisms for the Saturated Fatty Acid Methyl Esters from Methyl Butanoate to Methyl Palmitate. <i>Energy & Fuels</i> , 2015, 29, 1076-1089.	5.1	15
52	Development of a skeletal mechanism for diesel surrogate fuel by using a decoupling methodology. <i>Combustion and Flame</i> , 2015, 162, 3785-3802.	5.2	162
53	Understanding the Relationship between Cetane Number and the Ignition Delay in Shock Tubes for Different Fuels Based on a Skeletal Primary Reference Fuel (<i>n</i> -Hexadecane/Iso-cetane) Mechanism. <i>Energy & Fuels</i> , 2015, 29, 3413-3427.	5.1	23
54	Evaluation of the necessity of exhaust gas recirculation employment for a methanol/diesel reactivity controlled compression ignition engine operated at medium loads. <i>Energy Conversion and Management</i> , 2015, 101, 40-51.	9.2	57

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55	Development of a skeletal oxidation mechanism for biodiesel surrogate. Proceedings of the Combustion Institute, 2015, 35, 3037-3044.	3.9	69
56	Numerical simulation of cyclic variability in reactivity-controlled compression ignition combustion with a focus on the initial temperature at intake valve closing. International Journal of Engine Research, 2015, 16, 441-460.	2.3	43
57	Parametric study and optimization of a RCCI (reactivity controlled compression ignition) engine fueled with methanol and diesel. Energy, 2014, 65, 319-332.	8.8	173
58	Development of a reduced n-dodecane-PAH mechanism and its application for n-dodecane soot predictions. Fuel, 2014, 136, 25-36.	6.4	111
59	DEVELOPMENT OF A NEW SPRAY/WALL INTERACTION MODEL FOR DIESEL SPRAY UNDER PCCI-ENGINE RELEVANT CONDITIONS. Atomization and Sprays, 2014, 24, 41-80.	0.8	74
60	Numerical evaluation of the potential of late intake valve closing strategy for diesel PCCI (premixed) Tj ETQq0 0 0 rgBT /Overlock 10 Tf 5	8.8	61
61	Numerical study on the combustion and emission characteristics of a methanol/diesel reactivity controlled compression ignition (RCCI) engine. Applied Energy, 2013, 106, 184-197.	10.1	237
62	Development of a new skeletal mechanism for n-decane oxidation under engine-relevant conditions based on a decoupling methodology. Combustion and Flame, 2013, 160, 1315-1332.	5.2	154
63	Application of a Decoupling Methodology for Development of Skeletal Oxidation Mechanisms for Heavy <i>n</i> -Alkanes from <i>n</i> -Octane to <i>n</i> -Hexadecane. Energy & Fuels, 2013, 27, 3467-3479.	5.1	62
64	Improvement on a skeletal chemical kinetic model of iso-octane for internal combustion engine by using a practical methodology. Fuel, 2013, 103, 884-891.	6.4	67
65	Development of a Phenomenological Soot Model Coupled with a Skeletal PAH Mechanism for Practical Engine Simulation. Energy & Fuels, 2013, 27, 1699-1711.	5.1	49
66	Development of a New Skeletal Chemical Kinetic Model of Toluene Reference Fuel with Application to Gasoline Surrogate Fuels for Computational Fluid Dynamics Engine Simulation. Energy & Fuels, 2013, 27, 4899-4909.	5.1	105
67	Enhancement on a Skeletal Kinetic Model for Primary Reference Fuel Oxidation by Using a Semidecoupling Methodology. Energy & Fuels, 2012, 26, 7069-7083.	5.1	260
68	The effect of injection timing and intake valve close timing on performance and emissions of diesel PCCI engine with a full engine cycle CFD simulation. Applied Energy, 2011, 88, 2967-2975.	10.1	126
69	Development of a Reduced Chemical Mechanism for Dimethyl Ether (DME) Using a Decoupling Methodology. , 0, , .		6
70	Comparing the Exergy Destruction of Methanol and Gasoline in Reactivity Controlled Compression Ignition (RCCI) Engine. , 0, , .		13
71	Numerical Optimization of the Piston Bowl Geometry and Investigation of the Key Geometric Parameters for the Dual-Mode Dual-Fuel (DMDF) Concept under a Wide Load Range. , 0, , .		1