

Dong Tian

List of Publications by Year in descending order

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Version: 2024-02-01

39
papers

1,907
citations

331670

21
h-index

289244

40
g-index

41
all docs

41
docs citations

41
times ranked

2006
citing authors

#	ARTICLE	IF	CITATIONS
19	Optimization of Ni-Based Catalysts for Dry Reforming of Methane via Alloy Design: A Review. <i>Energy & Fuels</i> , 2022, 36, 5102-5151.	5.1	29
20	Noises- and delay-enhanced stability in a bistable dynamical system describing chemical reaction. <i>European Physical Journal B</i> , 2014, 87, 1.	1.5	24
21	Enhanced performance of chemical looping combustion of methane by combining oxygen carriers via optimizing the stacking sequences. <i>Applied Energy</i> , 2018, 230, 696-711.	10.1	22
22	The mechanism of photocatalyst and the effects of co-doping CeO ₂ on refractive index and reflectivity from DFT calculation. <i>Computational Materials Science</i> , 2019, 158, 197-208.	3.0	21
23	Tailoring of crystalline structure of carbon nitride for superior photocatalytic hydrogen evolution. <i>Journal of Colloid and Interface Science</i> , 2019, 556, 324-334.	9.4	20
24	Enhanced resistance to carbon deposition in chemical-looping combustion of methane: Synergistic effect of different oxygen carriers via sequence filling. <i>Chemical Engineering Journal</i> , 2021, 421, 129776.	12.7	20
25	Stochastic delayed monomer-dimer surface reaction model with various dimer adsorption. <i>European Physical Journal B</i> , 2014, 87, 1.	1.5	19
26	Surface amorphous carbon doping of carbon nitride for efficient acceleration of electron transfer to boost photocatalytic activities. <i>Applied Surface Science</i> , 2020, 507, 145145.	6.1	19
27	Delays-based protein switches in a stochastic single-gene network. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2015, 434, 68-83.	2.6	17
28	Exploring electrocatalytic stability and activity of unmodified and platinum-modified tungsten and niobium nitrides. <i>International Journal of Hydrogen Energy</i> , 2020, 45, 22883-22892.	7.1	17
29	Suppressing byproduct formation for high selective CO ₂ reduction over optimized Ni/TiO ₂ based catalysts. <i>Journal of Energy Chemistry</i> , 2022, 72, 465-478.	12.9	17
30	Modification of KNO ₃ on the reducibility and reactivity of Fe ₂ O ₃ -based oxygen carriers for chemical-looping combustion of methane. <i>Canadian Journal of Chemical Engineering</i> , 2017, 95, 1569-1578.	1.7	15
31	Noises-induced toggle switch and stability in a gene regulation network. <i>International Journal of Modern Physics B</i> , 2014, 28, 1450223.	2.0	13
32	First-principles study the behavior of oxygen vacancy on the surface of ZrO ₂ and Zr _{0.97} M _{0.03} O ₂ . <i>Computational Condensed Matter</i> , 2017, 11, 1-10.	2.1	13
33	Transition Metal Nitrides as Promising Catalyst Supports for Tuning CO/H ₂ Syngas Production from Electrochemical CO ₂ Reduction. <i>Angewandte Chemie</i> , 2020, 132, 11441-11444.	2.0	11
34	Water splitting for hydrogen generation over lanthanum-calcium-iron perovskite-type membrane driven by reducing atmosphere. <i>International Journal of Hydrogen Energy</i> , 2017, 42, 19776-19787.	7.1	10
35	Effect of Fe doping concentration on photocatalytic performance of CeO ₂ from DFT insight into analysis. <i>AIP Advances</i> , 2019, 9, .	1.3	10
36	Theoretical investigation the growth of Fe ₃ Si on GaAs: Stability and electronic properties of Fe ₃ Si/GaAs(0 0 1), (1 1 0) via DFT. <i>Applied Surface Science</i> , 2020, 506, 144691.	6.1	6

#	ARTICLE	IF	CITATIONS
37	Insights into the influence of functional groups on the properties of graphene from firstâ€principles calculations. Journal of Physical Organic Chemistry, 2022, 35, .	1.9	2
38	Hydrostatic pressures effect on structure stability, electronic, optical and elastic properties of rutile VO ₂ doped TiO ₂ by density functional theory investigation. Materials Research Express, 2019, 6, 0965c2.	1.6	1
39	Modeling and simulation of dual-three-phase induction machine with two opened phases. , 2008, , .		0