

RÃ©mi PÃ©tuya

List of Publications by Year in descending order

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Version: 2024-02-01

21
papers

487
citations

759233

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23
times ranked

700
citing authors

#	ARTICLE	IF	CITATIONS
1	Machine Learning Prediction of Metal-Organic Framework Guest Accessibility from Linker and Metal Chemistry. <i>Angewandte Chemie - International Edition</i> , 2022, 61, .	13.8	24
2	Machine Learning Prediction of Metal-Organic Framework Guest Accessibility from Linker and Metal Chemistry. <i>Angewandte Chemie</i> , 2022, 134, .	2.0	2
3	Size dependence of the dissociation process of spherical hydrate particles <i>via</i> microsecond molecular dynamics simulations. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 11180-11185.	2.8	10
4	Amino Acid Residues Determine the Response of Flexible Metal-Organic Frameworks to Guests. <i>Journal of the American Chemical Society</i> , 2020, 142, 14903-14913.	13.7	29
5	A new fluorescent hemicyptophane for acetylcholine recognition with an unusual recognition mode. <i>New Journal of Chemistry</i> , 2020, 44, 11853-11860.	2.8	11
6	Controlling Single Molecule Conductance by a Locally Induced Chemical Reaction on Individual Thiophene Units. <i>Angewandte Chemie</i> , 2020, 132, 6266-6271.	2.0	2
7	Controlling Single Molecule Conductance by a Locally Induced Chemical Reaction on Individual Thiophene Units. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 6207-6212.	13.8	9
8	The Anisotropic Responses of a Flexible Metal-Organic Framework Constructed from Asymmetric Flexible Linkers and Heptanuclear Zinc Carboxylate Secondary Building Units. <i>Crystal Growth and Design</i> , 2019, 19, 5604-5618.	3.0	6
9	Selective recognition of acetylcholine over choline by a fluorescent cage. <i>Organic and Biomolecular Chemistry</i> , 2019, 17, 5253-5257.	2.8	7
10	Magnetic coupling between 3d transition metal adatoms on graphene supported by metallic substrates. <i>Carbon</i> , 2017, 116, 599-605.	10.3	14
11	Electric-Field-Driven Direct Desulfurization. <i>ACS Nano</i> , 2017, 11, 4703-4709.	14.6	43
12	Ultrafast electronic response of graphene to a strong and localized electric field. <i>Nature Communications</i> , 2016, 7, 13948.	12.8	125
13	CO ₂ Binding and Induced Structural Collapse of a Surface-Supported Metal-Organic Network. <i>Journal of Physical Chemistry C</i> , 2016, 120, 18622-18630.	3.1	12
14	Hydrogen abstraction from metal surfaces: when electron-hole pair excitations strongly affect hot-atom recombination. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 31378-31383.	2.8	30
15	Scattering of Atomic Hydrogen Off a H-Covered W(110) Surface: Hot-Atom versus Eley-Rideal Abstraction Dynamics. <i>Journal of Physical Chemistry C</i> , 2015, 119, 3171-3179.	3.1	28
16	Isotope Effects in Eley-Rideal and Hot-Atom Abstraction Dynamics of Hydrogen from Tungsten (100) and (110) Surfaces. <i>Journal of Physical Chemistry C</i> , 2015, 119, 15325-15332.	3.1	14
17	Energy Dissipation to Tungsten Surfaces upon Eley-Rideal Recombination of N ₂ and H ₂ . <i>Journal of Physical Chemistry C</i> , 2015, 119, 15434-15442.	3.1	40
18	Bipolar Conductance Switching of Single Anthradithiophene Molecules. <i>ACS Nano</i> , 2015, 9, 12506-12512.	14.6	37

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19	Dynamics of H ₂ Eley-Rideal abstraction from W(110): Sensitivity to the representation of the molecule-surface potential. <i>Journal of Chemical Physics</i> , 2014, 141, 024701.	3.0	15
20	Comparative Theoretical Study of H ₂ Eley-Rideal Recombination Dynamics on W(100) and W(110). <i>Journal of Physical Chemistry C</i> , 2014, 118, 11704-11710.	3.1	19
21	Revisiting the Nonreactive Scattering of N ₂ off W(100): On the Influence of the Scattering Azimuth on In-Plane Angular Distributions. <i>Journal of Physical Chemistry C</i> , 2014, 118, 21904-21910.	3.1	10