

Xin Xu

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/5710284/publications.pdf>

Version: 2024-02-01

9
papers

159
citations

1478505
6
h-index

1588992
8
g-index

9
all docs

9
docs citations

9
times ranked

217
citing authors

#	ARTICLE	IF	CITATIONS
1	Quantum interference in $H + HD \rightarrow H_2 + D$ between direct abstraction and roaming insertion pathways. <i>Science</i> , 2020, 368, 767-771.	12.6	52
2	Direct observation of forward-scattering oscillations in the $H+HD \rightarrow H_2+D$ reaction. <i>Nature Chemistry</i> , 2018, 10, 653-658.	13.6	46
3	A full-dimensional time-dependent wave packet study of the $OH + CO \rightarrow H + CO_2$ reaction. <i>Theoretical Chemistry Accounts</i> , 2012, 131, 1.	1.4	21
4	A neural network potential energy surface for the $F + CH_4$ reaction including multiple channels based on coupled cluster theory. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 9090-9100.	2.8	21
5	An <i>ab initio</i> -based global potential energy surface for the SH_3 system and full-dimensional state-to-state quantum dynamics study for the $H_2 + HS \rightarrow H_2S + H$ reaction. <i>Journal of Computational Chemistry</i> , 2019, 40, 1151-1160.		10
6	State-to-state differential cross-sections for the reactive scattering of $H^*(n)$ with <i>o</i> -D ₂ . <i>Chemical Science</i> , 2012, 3, 2839.	7.4	6
7	Strong non-Arrhenius behavior at low temperatures in the $OH + HCl \rightarrow H_2O + Cl$ reaction due to resonance induced quantum tunneling. <i>Chemical Science</i> , 2022, 13, 7955-7961.	7.4	2
8	Quantum Wave Packet Study of the $H + Br_2 \rightarrow HBr + Br$ Reaction on a New <i>Ab Initio</i> Potential Energy Surface. <i>Journal of Physical Chemistry A</i> , 2021, 125, 7289-7296.	2.5	1
9	$D_2 + C(CH_3)_4 \rightarrow HD + C(CH_3)_3CH_2$: possible concerted flow of vibration energy. <i>Molecular Physics</i> , 2012, 110, 1713-1720.	1.7	0