

Yongfa Zhu

List of Publications by Year in descending order

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Version: 2024-02-01

11
papers

105
citations

1684188

5
h-index

1372567

10
g-index

11
all docs

11
docs citations

11
times ranked

64
citing authors

#	ARTICLE	IF	CITATIONS
1	Dynamics and kinetics of the reaction $\text{OH} + \text{H}_2\text{S} \rightarrow \text{H}_2\text{O} + \text{SH}$ on an accurate potential energy surface. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 26315-26324.	2.8	25
2	Theoretical study of the $\text{F}(\text{P}) + \text{NH}_3 \rightarrow \text{HF} + \text{NH}_2$ reaction on an accurate potential energy surface: dynamics and kinetics. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 11385-11394.	2.8	21
3	Tracking the energy flow in the hydrogen exchange reaction $\text{OH} + \text{H}_2\text{O} \rightarrow \text{H}_2\text{O} + \text{OH}$. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 12543-12556.	2.8	19
4	Kinetic and dynamic studies of the $\text{H}_3^+ + \text{CO} \rightarrow \text{H}_2 + \text{HCO}^+/\text{HOC}^+$ reaction on a high-level <i>ab initio</i> potential energy surface. <i>Journal of Chemical Physics</i> , 2019, 151, .	3.0	12
5	Mode specific dynamics in bond selective reaction $\text{O}(\text{3P}) + \text{HOD} \rightarrow \text{O}(\text{H}) + \text{OD}/\text{O}(\text{D}) + \text{OH}$. <i>Journal of Chemical Physics</i> , 2018, 149, 054304.	3.0	8
6	Mode-specific quantum dynamics and kinetics of the hydrogen abstraction reaction $\text{OH} + \text{H}_2\text{O} \rightarrow \text{H}_2\text{O} + \text{OH}$. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 24054-24060.	2.8	6
7	Final-State-Resolved Dynamics of the $\text{H}_3^+ + \text{CO} \rightarrow \text{H}_2^+ + \text{HCO}^+/\text{HOC}^+$ Reaction: A Quasi-Classical Trajectory Study. <i>Journal of Physical Chemistry A</i> , 2020, 124, 6794-6800.	2.5	5
8	Thermal rate coefficients and kinetic isotope effects of the reaction $\text{HO} + \text{H}_2\text{O} \rightarrow \text{H}_2\text{O} + \text{OH}$. <i>Theoretical Chemistry Accounts</i> , 2019, 138, 1.	1.4	4
9	Dissociative photodetachment of H_3O^+ : a full-dimensional quantum dynamics study. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 22298-22304.	2.8	2
10	Product vibrational state distributions of $\text{F} + \text{CH}_3\text{OH}$ reaction on full-dimensional accurate potential energy surface. <i>Chinese Journal of Chemical Physics</i> , 2022, 35, 153-166.	1.3	2
11	Assessing the applicability of the MBE approach for constructing potential energy surfaces of nitrogen clusters. <i>Chemical Physics</i> , 2021, 549, 111272.	1.9	1