

# Weixing Li

## List of Publications by Year in descending order

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31  
papers

376  
citations

933447

10  
h-index

839539

18  
g-index

32  
all docs

32  
docs citations

32  
times ranked

276  
citing authors

#	ARTICLE	IF	CITATIONS
1	Theory Meets Experiment for Noncovalent Complexes: The Puzzling Case of Phicogen Interactions. <i>Angewandte Chemie - International Edition</i> , 2018, 57, 13853-13857.	13.8	60
2	The Barrier to Proton Transfer in the Dimer of Formic Acid: A Pure Rotational Study. <i>Angewandte Chemie - International Edition</i> , 2019, 58, 859-865.	13.8	48
3	Internal dynamics of cyclohexanol and the cyclohexanol-water adduct. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 3676-3682.	2.8	25
4	A General Treatment to Study Molecular Complexes Stabilized by Hydrogen, Halogen, and Carbon Bond Networks: Experiment and Theory of $(\text{CH}_2\text{F}_2)_n \cdots (\text{H}_2\text{O})_m$ . <i>Angewandte Chemie - International Edition</i> , 2019, 58, 8437-8442.	13.8	19
5	Unlocking the Water Trimer Loop: Isotopic Study of Benzophenone- $(\text{H}_2\text{O})_3$ Clusters with Rotational Spectroscopy. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 5323-5330.	13.8	18
6	A butterfly motion of formic acid and cyclobutanone in the 1:1 hydrogen bonded molecular cluster. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 204-209.	2.8	16
7	Atmospherically relevant acrolein-water complexes: spectroscopic evidence of aldehyde hydration and oxygen atom exchange. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 23559-23566.	2.8	16
8	Site-selective ionization of ethanol dimer under the tunable synchrotron VUV radiation and its subsequent fragmentation. <i>Journal of Chemical Physics</i> , 2013, 139, 024307.	3.0	15
9	The Borderline between Reactivity and Pre-reactivity of Binary Mixtures of Gaseous Carboxylic Acids and Alcohols. <i>Angewandte Chemie - International Edition</i> , 2017, 56, 3872-3875.	13.8	14
10	Shape of the Adduct Formic Acid-Dimethyl Ether: A Rotational Study. <i>Journal of Physical Chemistry A</i> , 2016, 120, 2863-2867.	2.5	12
11	Characterizing the lone pair-hole interaction in complexes of ammonia with perfluorinated arenes. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 9121-9129.	2.8	11
12	Unexpected methyl migrations of ethanol dimer under synchrotron VUV radiation. <i>Journal of Chemical Physics</i> , 2015, 142, 024306.	3.0	10
13	Conformational Equilibrium and Potential Energy Functions of the O-H Internal Rotation in the Axial and Equatorial Species of 1-Methylcyclohexanol. <i>Journal of Physical Chemistry A</i> , 2016, 120, 4338-4342.	2.5	10
14	Barrier to Proton Transfer in the Dimer of Formic Acid: A Pure Rotational Study. <i>Angewandte Chemie</i> , 2018, 131, 869.	2.0	10
15	Unlocking the Water Trimer Loop: Isotopic Study of Benzophenone- $(\text{H}_2\text{O})_3$ Clusters with Rotational Spectroscopy. <i>Angewandte Chemie</i> , 2021, 133, 5383-5390.	2.0	10
16	Characterizing hydrogen and tetrel bonds in clusters of CO <sub>2</sub> with carboxylic acids. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 16915-16922.	2.8	10
17	Double Proton Transfer Across a Table: The Formic Acid Dimer-Fluorobenzene Complex. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 25674-25679.	13.8	10
18	The microwave spectroscopy study of 1,2-dimethoxyethane. <i>Journal of Molecular Spectroscopy</i> , 2017, 337, 3-8.	1.2	9

#	ARTICLE	IF	CITATIONS
19	The Cage Structure of Indan-1,2,3-CH <sub>3</sub> is Based on the Cooperative Effects of C-H...H and C-H...F Weak Hydrogen Bonds. <i>Chemistry - A European Journal</i> , 2015, 21, 15970-15973.	3.8	7
20	Theory Meets Experiment for Noncovalent Complexes: The Puzzling Case of Pnictogen Interactions. <i>Angewandte Chemie</i> , 2018, 130, 14049-14053.	2.0	7
21	Ring puckering splitting and structure of indan. <i>Journal of Molecular Spectroscopy</i> , 2015, 316, 45-48.	1.2	6
22	The Borderline between Reactivity and Pre-reactivity of Binary Mixtures of Gaseous Carboxylic Acids and Alcohols. <i>Angewandte Chemie</i> , 2017, 129, 3930-3933.	2.0	6
23	Carboxylic Acids, Reactivity with Alcohols and Clustering with Esters: A Rotational Study of Formic Acid-Isopropylformate. <i>Journal of Physical Chemistry A</i> , 2019, 123, 1785-1789.	2.5	6
24	The rotational spectrum of methyl trifluoroacetate. <i>Molecular Physics</i> , 2018, 116, 3503-3506.	1.7	5
25	A General Treatment to Study Molecular Complexes Stabilized by Hydrogen, Halogen, and Carbon-Bond Networks: Experiment and Theory of (CH <sub>2</sub> F) <sub>2</sub> ...n...m... (H <sub>2</sub> O). <i>Angewandte Chemie</i> , 2019, 131, 8525-8530.	2.0	4
26	Interactions between azines and alcohols: a rotational study of pyridine- <i>tert</i> -butyl alcohol. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 3545-3549.	2.8	3
27	The rotational spectrum of cyclohexyl formate, chemically prepared within a supersonic expansion. <i>Journal of Molecular Structure</i> , 2020, 1209, 127952.	3.6	3
28	Rotational studies of adducts between carboxylic acids and tertiary alcohols: Formic acid- <i>tert</i> -butyl alcohol. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2021, 254, 119621.	3.9	3
29	Double proton transfer across a table: the formic acid dimer-fluorobenzene complex. <i>Angewandte Chemie</i> , 0, .	2.0	2
30	InnenrÄ4cktitelbild: Double Proton Transfer Across a Table: The Formic Acid Dimer-Fluorobenzene Complex ( <i>Angew. Chem.</i> 49/2021). <i>Angewandte Chemie</i> , 2021, 133, 26203-26203.	2.0	1
31	Site-Selective Dissociation Processes of Cationic Ethanol Conformers: The Role of Hyperconjugation. <i>Journal of Physical Chemistry A</i> , 2014, 118, 7096-7103.	2.5	0