

# Sung-Woo Park

## List of Publications by Year in descending order

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34  
papers

1,399  
citations

331670

21  
h-index

345221

36  
g-index

40  
all docs

40  
docs citations

40  
times ranked

1605  
citing authors

#	ARTICLE	IF	CITATIONS
1	Generation of N-Centered Radicals via a Photocatalytic Energy Transfer: Remote Double Functionalization of Arenes Facilitated by Singlet Oxygen. <i>Journal of the American Chemical Society</i> , 2019, 141, 10538-10545.	13.7	75
2	Conjugate Addition of Perfluoroarenes to $\hat{I}^{\pm}, \hat{I}^2$ -Unsaturated Carbonyls Enabled by an Alkoxide-Hydrosilane System: Implication of a Radical Pathway. <i>Journal of the American Chemical Society</i> , 2018, 140, 9659-9668.	13.7	15
3	Visible light sensitization of benzoyl azides: cascade cyclization toward oxindoles via a non-nitrene pathway. <i>Chemical Communications</i> , 2017, 53, 8798-8801.	4.1	34
4	High-temperature in situ crystallographic observation of reversible gas sorption in impermeable organic cages. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2015, 112, 14156-14161.	7.1	27
5	Cp*Ir(III)-Catalyzed Mild and Broad C-H Arylation of Arenes and Alkenes with Aryldiazonium Salts Leading to the External Oxidant-Free Approach. <i>Journal of the American Chemical Society</i> , 2015, 137, 8584-8592.	13.7	125
6	Chemoselective Silylative Reduction of Conjugated Nitriles under Metal-Free Catalytic Conditions: $\hat{I}^2$ -Silyl Amines and Enamines. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 6832-6836.	13.8	67
7	Complete Switch of Selectivity in the C-H Alkenylation and Hydroarylation Catalyzed by Iridium: The Role of Directing Groups. <i>Journal of the American Chemical Society</i> , 2015, 137, 13448-13451.	13.7	127
8	Novel Ionophores with 2-Crown-1 Topology: Anion Sensing via Pure Aliphatic C-H...Anion Hydrogen Bonding. <i>Organic Letters</i> , 2014, 16, 334-337.	4.6	21
9	Boron-Catalyzed Silylative Reduction of Quinolines: Selective $sp^3$ C-Si Bond Formation. <i>Journal of the American Chemical Society</i> , 2014, 136, 16780-16783.	13.7	113
10	Gas phase hydration of amino acids and dipeptides: effects on the relative stability of zwitterion vs. canonical conformers. <i>RSC Advances</i> , 2014, 4, 16352-16361.	3.6	26
11	Design of Carbene-Based Organocatalysts for Nitrogen Fixation: Theoretical Study. <i>Journal of Chemical Theory and Computation</i> , 2012, 8, 1983-1988.	5.3	20
12	An unprecedented $\hat{I}^2$ -linear-bent isomerism in tri-nuclear Cu <sub>2</sub> ZnII complexes with a salen type di-Schiff base ligand. <i>Dalton Transactions</i> , 2012, 41, 11009.	3.3	69
13	Very Efficient Nucleophilic Aromatic Fluorination Reaction in Molten Salts: A Mechanistic Study. <i>Bulletin of the Korean Chemical Society</i> , 2012, 33, 881-884.	1.9	5
14	SN <sub>2</sub> Fluorination reactions in ionic liquids: a mechanistic study towards solvent engineering. <i>Organic and Biomolecular Chemistry</i> , 2011, 9, 418-422.	2.8	37
15	Counterion-Mediated Hydrogen-Bonding Effects: Mechanistic Study of Gold(I)-Catalyzed Enantioselective Hydroamination of Allenes. <i>Chemistry - an Asian Journal</i> , 2011, 6, 1982-1986.	3.3	45
16	Effects of Ion and Protic Solvent on Nucleophilic Aromatic Substitution ( $S_NAr$ ) Reactions. <i>Bulletin of the Korean Chemical Society</i> , 2010, 31, 2571-2574.	1.9	15
17	A Mechanistic Study of SN <sub>2</sub> Reaction in a Diol Solvent. <i>Journal of Physical Chemistry A</i> , 2009, 113, 3685-3689.	2.5	15
18	Bis-Terminal Hydroxy Polyethers as All-Purpose, Multifunctional Organic Promoters: A Mechanistic Investigation and Applications. <i>Angewandte Chemie - International Edition</i> , 2009, 48, 7683-7686.	13.8	103

#	ARTICLE	IF	CITATIONS
19	Ab Initio and DFT Studies of the Thermal Rearrangement of Trimethylsilylsilylene. <i>Organometallics</i> , 2008, 27, 2123-2127.	2.3	4
20	Efficiency of Bulky Protic Solvent for S <sub>N</sub> 2 Reaction. <i>Organic Letters</i> , 2008, 10, 61-64.	4.6	23
21	Facile SN2 Reaction in Protic Solvent: A Quantum Chemical Analysis. <i>Journal of Physical Chemistry A</i> , 2007, 111, 10152-10161.	2.5	52
22	Structure and stability of glycine(H <sub>2</sub> O) <sub>3</sub> cluster and anion: Zwitterion vs. canonical glycine. <i>International Journal of Quantum Chemistry</i> , 2007, 107, 1316-1327.	2.0	27
23	Geometry, Chemical Bonding, and Electronic Spectra of Sinand Sin <sup>n</sup> Glycine (n= 3~5) Complexes. <i>Journal of Physical Chemistry A</i> , 2006, 110, 7173-7177.	2.5	5
24	Structures and isomerization of serine in aqueous solution: Computational study. <i>Chemical Physics Letters</i> , 2005, 403, 72-76.	2.6	12
25	Structures and isomerization of neutral and zwitterion serine-water clusters: Computational study. <i>International Journal of Quantum Chemistry</i> , 2005, 101, 55-66.	2.0	40
26	Intermediate Complexes in S <sub>N</sub> 2 Reaction: [Na <sup>+</sup> , F <sup>-</sup> , H <sub>2</sub> O, CH <sub>3</sub> Cl] System. <i>Bulletin of the Korean Chemical Society</i> , 2005, 26, 2081-2083.	1.9	2
27	Computational Study of Proline - Water Cluster. <i>Bulletin of the Korean Chemical Society</i> , 2005, 26, 909-912.	1.9	19
28	Dynamic paths between neutral alanine-water and zwitterionic alanine-water clusters: single, double and triple proton transfer. <i>Chemical Physics Letters</i> , 2003, 371, 74-79.	2.6	58
29	Effects of Substituting Group on the Hydrogen Bonding in Phenol-H <sub>2</sub> O Complexes: Ab Initio Study. <i>Journal of Physical Chemistry A</i> , 2003, 107, 131-139.	2.5	47
30	Effects of Microsolvation on the Structures and Reactions of Neutral and Zwitterion Alanine: Computational Study. <i>Journal of Physical Chemistry B</i> , 2003, 107, 14109-14118.	2.6	76
31	Hydrogen Bonding in Aromatic Alcohol-Water Clusters: A Brief Review. <i>Bulletin of the Korean Chemical Society</i> , 2003, 24, 695-702.	1.9	32
32	Structure, Spectroscopic Properties and Reactions of Interstellar Molecule HC <sub>2</sub> N and Isomers: Ab initio Study. <i>Bulletin of the Korean Chemical Society</i> , 2002, 23, 1553-1559.	1.9	5
33	Computational Study of Catechol-(H <sub>2</sub> O) <sub>n</sub> (n=1-3) Clusters. <i>Bulletin of the Korean Chemical Society</i> , 2002, 23, 1297-1303.	1.9	6
34	Computational study of medium-sized cumulenethiones H <sub>2</sub> C <sub>n</sub> S (n=3~9). <i>Chemical Physics Letters</i> , 2000, 326, 530-536.	2.6	7