

# Takao Okazaki

## List of Publications by Year in descending order

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38  
papers

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citations

840776

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times ranked

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#	ARTICLE	IF	CITATIONS
1	Brønsted acid-catalyzed aldol cyclotrimerization of indanones in ionic liquid: An experimental and DFT study of substituent effect. <i>Journal of Physical Organic Chemistry</i> , 2021, 34, .	1.9	0
2	Phenyl(triazolyl)carbene revisited: Unique role of triazolyl group on carbene chemistry. <i>Journal of Physical Organic Chemistry</i> , 2021, 34, e4187.	1.9	0
3	Charge delocalization mode and change in aromaticity of protonated 7-phenylbenzo[ <i>k</i> ]fluoranthenes studied by experimental observation and DFT calculations. <i>Journal of Physical Organic Chemistry</i> , 2019, 32, e3998.	1.9	1
4	Electron-Transfer Properties of Phenyleneethynylene Linkers Bound to Gold via a Self-Assembled Monolayer of Molecular Tripod. <i>Molecules</i> , 2018, 23, 2893.	3.8	5
5	Tuning the coverage of self-assembled monolayer by introducing bulky substituents onto rigid adamantane tripod. <i>Arkivoc</i> , 2018, 2018, 131-144.	0.5	4
6	Effect of solvation of ionic liquid on Brønsted acid-catalyzed aldol cyclotrimerization of indanones and related cyclic ketones. <i>Journal of Physical Organic Chemistry</i> , 2018, 31, e3887.	1.9	1
7	NMR and DFT studies on persistent carbocations derived from benzo[ <i>k</i> ]xanthene, dibenzo[ <i>b</i> , <i>d</i> ]benzo[1,2- <i>b</i> :4,3- <i>b</i> ]difuran, and dibenzo[ <i>d</i> , <i>d</i> ]benzo[1,2- <i>b</i> :4,5- <i>b</i> ]difuran in superacidic media. <i>Journal of Physical Organic Chemistry</i> , 2016, 29, 107-111.	1.9	8
8	Cationic Intermediates for Electrophilic Reactions from 9,9-Dimethyl-9H-9-silafluorene. <i>Current Organic Chemistry</i> , 2016, 20, 3014-3021.	1.6	2
9	Solvolysis of 2-adamantyl <i>p</i> -toluenesulfonate in ionic liquid 1-butyl-3-methylimidazolium tetrafluoroborate. <i>Tetrahedron Letters</i> , 2015, 56, 6066-6068.	1.4	4
10	Electrochemistry of the Self-Assembled Monolayers of Dyads Consisting of Tripod-Shaped Trithiol and Bithiophene on Gold. <i>Molecules</i> , 2014, 19, 15298-15313.	3.8	12
11	4-(Pentafluorosulfonyl)benzenediazonium Tetrafluoroborate: A Versatile Launch Pad for the Synthesis of Aromatic SF <sub>5</sub> Compounds via Cross Coupling, Azo Coupling, Homocoupling, Dediazonation, and Click Chemistry. <i>European Journal of Organic Chemistry</i> , 2014, 2014, 1630-1644.	2.4	31
12	Experimental NMR and DFT Studies of Persistent Carbocations Derived from Hetero-Polycyclic Aromatic Hydrocarbons Containing Oxygen Atom: Dibenzo[ <i>b</i> , <i>d</i> ]furan, Benzo[ <i>b</i> ]naphtho[1,2- <i>d</i> ]furan, Benzo[ <i>b</i> ]naphtho[2,3- <i>d</i> ]furan, Benzo[ <i>b</i> ]naphtho[2,1- <i>d</i> ]furan, and Dinaphtho[2,1- <i>b</i> :1,2- <i>d</i> ]furan. <i>Bulletin of the Chemical Society of Japan</i> , 2014, 87, 1235-1244.	3.2	4
13	Ideal Redox Behavior of the High-Density Self-Assembled Monolayer of a Molecular Tripod on a Au(111) Surface with a Terminal Ferrocene Group. <i>Langmuir</i> , 2013, 29, 4275-4282.	3.5	49
14	NMR and DFT Study on Onium Ions Derived from Substituted Fluoranthenes and Benzo[ <i>k</i> ]fluoranthenes. <i>Bulletin of the Chemical Society of Japan</i> , 2013, 86, 464-471.	3.2	5
15	A Computational (DFT, MP2) and GIAO NMR Study of Substituent Effects in Benzenediazonium Mono- and Dications. <i>European Journal of Organic Chemistry</i> , 2011, 2011, 1771-1775.	2.4	6
16	Self-Sensitized Photooxygenation of a C <sub>60</sub> -Cycloheptatriene Dyad to Form Norcaradiene-Derived Endoperoxides. <i>European Journal of Organic Chemistry</i> , 2010, 2010, 3257-3264.	2.4	5
17	Stable NMR Spectroscopy and GIAO-DFT Study of Carbocations Derived from Multibridged [3< <i>n</i> >-Cyclophanes. <i>European Journal of Organic Chemistry</i> , 2009, 2009, 4451-4457.	2.4	4
18	Stable NMR and GIAO-DFT Study of the Carbocations from Benzofluorenes and Dibenzofluorenes; Synthesis of Nitro Derivatives; Mutagenicity Assay and X-ray Analysis. <i>European Journal of Organic Chemistry</i> , 2008, 2008, 1740-1752.	2.4	9

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19	Superacid-Catalyzed Dimerization/Cyclization of Isopropenyl-PAHs: Novel Pathways to PAH Dimers, Phenalenes and Their Stable Carbocations. <i>European Journal of Organic Chemistry</i> , 2008, 2008, 3700-3708.	2.4	4
20	Synthesis and Stability Studies of Regioisomeric Acetylnitropyrenes and Nitropyrenyl Carbinols and GIAO-DFT Study of Nitro Substituent Effects on Pyrenyl Carbocations. <i>European Journal of Organic Chemistry</i> , 2008, 2008, 6093-6105.	2.4	6
21	R(Ar)O <sup>+</sup> N <sub>2</sub> vs. R(Ar)N <sub>2</sub> O <sup>+</sup> : Are Alkoxy-(Aryloxy)-diazonium Ions or Alkyl-(Aryl)-N-nitroso-onium Ions Formed in the Gas-Phase Reactions of N <sub>2</sub> O with H <sup>+</sup> , Me <sup>+</sup> , Ph <sup>+</sup> , PhCH <sub>2</sub> <sup>+</sup> , Tr <sup>+</sup> and PhCO <sup>+</sup> ? <i>European Journal of Organic Chemistry</i> , 2007, 2007, 70-77.	2.4	4
22	Stable Ion and Electrophilic Substitution (Nitration and Bromination) Study of A-Ring Substituted Phenanthrenes: Novel Carbocations and Substituted Derivatives; NMR, X-ray Analysis, and Comparative DNA Binding. <i>European Journal of Organic Chemistry</i> , 2007, 2007, 487-497.	2.4	8
23	Rigid Molecular Tripod with an Adamantane Framework and Thiol Legs. Synthesis and Observation of an Ordered Monolayer on Au(111). <i>Journal of Organic Chemistry</i> , 2006, 71, 1362-1369.	3.2	90
24	Transannular $\pi$ - $\pi$ interactions in janusenes and in related rigid systems with cofacial aromatic rings; gauging aromaticity in the hydrocarbons and in model carbocations; a DFT study. <i>Organic and Biomolecular Chemistry</i> , 2006, 4, 3085-3095.	2.8	18
25	Intermediates of Halogen Addition to Phenylethyne and Protonation of Phenylethynyl Halides. Open Halovinyl Cations, Bridged Halonium, or Phenyl-Bridged Ions: A Substituent Effect Study by DFT and GIAO-DFT. <i>Journal of Organic Chemistry</i> , 2006, 71, 9643-9650.	3.2	12
26	Probing the Intermediates of Halogen Addition to Alkynes: A Bridged Halonium versus Open Halovinyl Cation; A Theoretical Study. <i>Journal of Organic Chemistry</i> , 2005, 70, 9139-9146.	3.2	11
27	Theoretical (DFT, GIAO-NMR, NICS) study of carbocations (M+H) <sup>+</sup> , dications (M <sub>2</sub> <sup>+</sup> ) and dianions (M <sub>2</sub> <sup>2-</sup> ) from dihydro-dicyclopenta[ef,kl]heptalene (dihydro-azupyrene), dihydro-dicyclohepta[ed,gh]pentalene, and related bridged [14]annulenes. <i>Organic and Biomolecular Chemistry</i> , 2005, 3, 286-294.	2.8	7
28	Electrophilic and oxidative chemistry of pyrene and its non-alternant isomers: theoretical (DFT, Tj ETQq0 0 0 rgBT /Overlock 10 Tf 50 39 (dicyclopenta[ef,kl]heptalene) and dicyclohepta[ed,gh]pentalene. <i>Organic and Biomolecular Chemistry</i> , 2004, 2, 2214-2219.	2.8	15
29	Pyrene (BaP) derivatives: 7,8-dihydro-BaP, 9,10-dihydro-BaP and its 6-halo derivatives, 1- and 3-methoxy-9,10-dihydro-BaP-7(8H)-one, as well as the proximate carcinogen BaP 7,8-dihydrodiol and its dibenzoate, combined with a comparative DNA binding study of regioisomeric (1-, 4-, 2-) pyrenylcarbinols. Electronic supplementary information (ESI) available: Selected NMR spectra (Fig. S1 and Charts S1-S10) and DFT computed energies for carbocations (Table S1). See <a href="http://www.rsc.org/s/organicandbiomolecularchemistry">http://www.rsc.org/s/organicandbiomolecularchemistry</a> , 2003, 1, 1509-1516.	2.8	15
30	$\beta$ -Silyl-Substituted Siladamantyl, Silabicyclo[2.2.2]octyl, Silanorbornyl, and 1-Silacyclohexyl Cations. A Theoretical (DFT and GIAO NMR) Study. <i>Journal of Organic Chemistry</i> , 2003, 68, 1827-1833.	3.2	7
31	Novel Examples of Three-Dimensional Aromaticity: 1,3-Dehydro-siladamantane Dications. A Theoretical (DFT, GIAO NMR, NICS) Study. <i>Journal of Organic Chemistry</i> , 2002, 67, 8721-8725.	3.2	11
32	Protonation studies on epimeric homoallylic adamantylideneadamantyl alcohols, 4-methyleneadamantylideneadamantane, adamantylideneadamantane (Ad <sup>+</sup> Ad) and sesquihomoadamantene, and reaction of Ad <sup>+</sup> Ad and sesquihomoadamantene with NO <sub>2</sub> +BF <sub>4</sub> <sup>-</sup> and PhI(OH)OTs: a stable-ion NMR and theoretical (GIAO-NMR) study. Electronic supplementary information (ESI) available: representative 1D-NMR spectra and tables of cartesian coordinates. See <a href="http://www.rsc.org/suppdata/p2/b2/b201660e/">http://www.rsc.org/suppdata/p2/b2/b201660e/</a> . <i>Perkin Transactions II RSC</i> , 2002, 1105-1111.	1.1	6
33	Stable ion and electrophilic chemistry of fluoranthene-PAHs. Electronic supplementary information (ESI) available: Table S1, NMR spectra and results of calculations. See <a href="http://www.rsc.org/suppdata/p2/b1/b108025n/">http://www.rsc.org/suppdata/p2/b1/b108025n/</a> . <i>Perkin Transactions II RSC</i> , 2002, 621-629.	1.1	10
34	Generation and NMR studies of stable cations derived from monothia[3.2]- and dithia[3.3]metacyclophanes. <i>Perkin Transactions II RSC</i> , 2001, 745-748.	1.1	4
35	First Examples of Stable Arenium Ions from Large Methylene-Bridged Polycyclic Aromatic Hydrocarbons (PAHs). Directive Effects and Charge Delocalization Mode. <i>Journal of Organic Chemistry</i> , 2001, 66, 3977-3983.	3.2	19
36	Substituent Effects and Charge Delocalization Mode in Chrysenium, Benzo[c]phenanthrenium, and Benzo[g]chrysenium Cations: A Stable Ion and Electrophilic Substitution Study. <i>Journal of Organic Chemistry</i> , 2001, 66, 780-788.	3.2	28

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37	Persistent Oxidation Dications from Twisted Fluoranthenes, Benzo[k]fluoranthene and Dimethyldibenzo[j,l]fluoranthene: A Charge Delocalization Mode, Tropicity, and Formation of Novel 8,8- $\pi$ -Bifluoranthenyls. An NMR and Theoretical Study. <i>Journal of Organic Chemistry</i> , 2001, 66, 8701-8708.	3.2	6
38	Electrophilic reactivity and $\pi$ -complexation studies in 1,8-naphthylene-bridged [3.2]paracyclophane with a cyclobutane calliper. <i>Perkin Transactions II RSC</i> , 2000, , 2347-2350.	1.1	5