

# Margarida S Miranda

## List of Publications by Year in descending order

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76  
papers

1,460  
citations

361413

20  
h-index

377865

34  
g-index

76  
all docs

76  
docs citations

76  
times ranked

1481  
citing authors

#	ARTICLE	IF	CITATIONS
1	Epitope-Imprinted Nanoparticles as Transforming Growth Factor- $\beta$ 3 Sequestering Ligands to Modulate Stem Cell Fate. <i>Advanced Functional Materials</i> , 2021, 31, 2003934.	14.9	21
2	Magnetic triggers in biomedical applications – prospects for contact free cell sensing and guidance. <i>Journal of Materials Chemistry B</i> , 2021, 9, 1259-1271.	5.8	7
3	Toward Spinning Greener Advanced Silk Fibers by Feeding Silkworms with Nanomaterials. <i>ACS Sustainable Chemistry and Engineering</i> , 2020, 8, 11872-11887.	6.7	20
4	Remote triggering of TGF- $\beta$ 2/Smad2/3 signaling in human adipose stem cells laden on magnetic scaffolds synergistically promotes tenogenic commitment. <i>Acta Biomaterialia</i> , 2020, 113, 488-500.	8.3	12
5	Exploring inhalable polymeric dry powders for anti-tuberculosis drug delivery. <i>Materials Science and Engineering C</i> , 2018, 93, 1090-1103.	7.3	23
6	Magnetic responsive cell-based strategies for diagnostics and therapeutics. <i>Biomedical Materials (Bristol)</i> , 2018, 13, 054001.	3.3	24
7	Development of Inhalable Superparamagnetic Iron Oxide Nanoparticles (SPIONs) in Microparticulate System for Antituberculosis Drug Delivery. <i>Advanced Healthcare Materials</i> , 2018, 7, e1800124.	7.6	34
8	A theoretical study of the strong interactions between carbon dioxide and OH+ and NH2 + products resulting from protonation of 1,2-dioxirane-3-one and 1,2-oxaziridine-3-one, respectively. <i>Structural Chemistry</i> , 2016, 27, 1743-1751.	2.0	6
9	What is the enthalpy of formation of pyrazine-2-carboxylic acid?. <i>Journal of Chemical Thermodynamics</i> , 2016, 97, 261-263.	2.0	3
10	Theoretical characterization of the chemical bonds of some three-membered ring compounds through QTAIM theory. <i>Structural Chemistry</i> , 2016, 27, 663-670.	2.0	10
11	Three-membered ring amides – a calculational and conceptual study of the structure and energetics of 1,2-oxaziridine-3-one and aziridine-2,3-dione. <i>Canadian Journal of Chemistry</i> , 2015, 93, 406-413.	1.1	6
12	Reprint of Indenone and cyclopentadienone energetics via mass spectrometry and computations: Are these species antiaromatic or nonaromatic?. <i>International Journal of Mass Spectrometry</i> , 2015, 378, 175-179.	1.5	2
13	Protonated heterocyclic derivatives of cyclopropane and cyclopropanone: classical species, alternate sites, and ring fragmentation. <i>Canadian Journal of Chemistry</i> , 2015, 93, 708-714.	1.1	5
14	Degradation in chlorinated water of the UV filter 4- <i>tert</i> -butyl-4-methoxydibenzoylmethane present in commercial sunscreens. <i>Environmental Technology (United Kingdom)</i> , 2015, 36, 1319-1326.	2.2	25
15	A theoretical study of the UV absorption of 4-methylbenzylidene camphor: from the UVB to the UVA region. <i>Photochemical and Photobiological Sciences</i> , 2015, 14, 465-472.	2.9	9
16	Gas-phase thermochemical properties of some tri-substituted phenols: A density functional theory study. <i>Journal of Chemical Thermodynamics</i> , 2015, 80, 65-72.	2.0	5
17	Theoretical characterization of molecular complexes formed between triplet vinyl nitrene and Lewis acids. <i>Structural Chemistry</i> , 2015, 26, 565-571.	2.0	2
18	Indenone and cyclopentadienone energetics via mass spectrometry and computations: Are these species antiaromatic or nonaromatic?. <i>International Journal of Mass Spectrometry</i> , 2014, 369, 87-91.	1.5	4

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19	Theoretical study of the effect of resonance on $\pi$ - $\pi$ stacked firefly oxyluciferin dimers. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 2014, 278, 9-13.	3.9	1
20	A computational study of the structure, aromaticity and enthalpy of formation of UVA filter 4-tert-butyl-4'-methoxydibenzoylmethane. <i>Computational and Theoretical Chemistry</i> , 2014, 1038, 6-16.	2.5	1
21	UV filter 2-ethylhexyl 4'-methoxycinnamate: a structure, energetic and UV-vis spectral analysis based on density functional theory. <i>Journal of Physical Organic Chemistry</i> , 2014, 27, 47-56.	1.9	24
22	Photodegradation of avobenzone: Stabilization effect of antioxidants. <i>Journal of Photochemistry and Photobiology B: Biology</i> , 2014, 140, 36-40.	3.8	131
23	Study of the transformation of two salicylates used in personal care products in chlorinated water. <i>Water Research</i> , 2014, 65, 32-39.	11.3	4
24	Structural, Energetic, and UV-Vis Spectral Analysis of UVA Filter 4-tert-Butyl-4'-methoxydibenzoylmethane. <i>Journal of Physical Chemistry A</i> , 2014, 118, 1511-1518.	2.5	26
25	Feeling and investigating blue: On the enthalpy of formation of indigo. <i>Journal of Chemical Thermodynamics</i> , 2014, 73, 69-75.	2.0	3
26	Occurrence of Personal Care Products and Transformation Processes in Chlorinated Waters. <i>Handbook of Environmental Chemistry</i> , 2014, , 123-136.	0.4	1
27	Gas-phase molecular structure and energetics of UVB filter 4-methylbenzylidene camphor: A computational study. <i>Computational and Theoretical Chemistry</i> , 2014, 1033, 67-73.	2.5	3
28	Enthalpy of formation of 5-fluoro-1,3-dimethyluracil: 5-Fluorouracil revisited. <i>Journal of Chemical Thermodynamics</i> , 2014, 75, 106-115.	2.0	13
29	Computational Study on the Vinyl Azide Decomposition. <i>Journal of Physical Chemistry A</i> , 2014, 118, 5038-5045.	2.5	14
30	Solid luminescent CdSe-thiolated porous phosphate heterostructures. Application in fingerprint detection in different surfaces. <i>Surface and Interface Analysis</i> , 2013, 45, 612-618.	1.8	16
31	Luminescent behavior of CdTe quantum dots: Neodymium(III) complex-capped nanoparticles. <i>Journal of Luminescence</i> , 2013, 134, 408-413.	3.1	2
32	The structure and energetics of pyrrolidinones, tetrahydrofuranones, piperidinones, and tetrahydropyranones: a computational study. <i>Structural Chemistry</i> , 2013, 24, 1829-1839.	2.0	8
33	Structure and energetics correlations in some chlorohydroxypyridines. <i>Journal of Chemical Thermodynamics</i> , 2013, 62, 170-177.	2.0	3
34	Amino, ammonio and aminioethenes: a theoretical study of their structure and energetics. <i>Journal of Physical Organic Chemistry</i> , 2013, 26, 613-625.	1.9	4
35	Degradation of UV filters 2-ethylhexyl-4-methoxycinnamate and 4-tert-butyl-4'-methoxydibenzoylmethane in chlorinated water. <i>Environmental Chemistry</i> , 2013, 10, 127.	1.5	32
36	Thiolated DAB dendrimers and CdSe quantum dots nanocomposites for Cd(II) or Pb(II) sensing. <i>Talanta</i> , 2012, 88, 403-407.	5.5	48

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37	The degradation products of UV filters in aqueous and chlorinated aqueous solutions. <i>Water Research</i> , 2012, 46, 3167-3176.	11.3	133
38	Aza-Diels-Alder reaction between cyclopentadiene and protonated N-phenylethyliminoacetates of 8-phenylmenthol and 8-phenylneomenthol: a density functional theory study. <i>Journal of Physical Organic Chemistry</i> , 2012, 25, 515-522.	1.9	4
39	Paradigms and paradoxes: en route to the understanding of the aromaticity of the $\sigma$ -isobenzofuran, anthranil, benzofurazan and 2,1,3-benzothiadiazole. <i>Structural Chemistry</i> , 2012, 23, 1241-1243.	2.0	7
40	2,1,3-Benzothiadiazole: Study of its structure, energetics and aromaticity. <i>Journal of Chemical Thermodynamics</i> , 2012, 50, 30-36.	2.0	19
41	Study of Energetics and Structure of 1,2,3-Benzotriazin-4(3H)-one and its 1H and Enol Tautomers. <i>Journal of Physical Chemistry B</i> , 2011, 115, 6616-6622.	2.6	14
42	Energetics of Quinazoline-2,4(1H,3H)-dione: An Experimental and Computational Study. <i>Journal of Chemical &amp; Engineering Data</i> , 2011, 56, 4516-4523.	1.9	12
43	CdSe quantum dots capped PAMAM dendrimer nanocomposites for sensing nitroaromatic compounds. <i>Talanta</i> , 2011, 83, 1335-1340.	5.5	56
44	Paradigms and paradoxes: the aromaticity of 6:6 fused carbocycles and heterocycles as an extension of a study of indane and indene derivatives. <i>Structural Chemistry</i> , 2011, 22, 1221-1224.	2.0	8
45	Combined experimental and computational study on the energetics of 1,2-benzisothiazol-3(2H)-one and 1,4-benzothiazin-3(2H,4H)-one. <i>Journal of Chemical Thermodynamics</i> , 2011, 43, 635-644.	2.0	13
46	Experimental and computational study on the energetics of 10,11-dihydro-5H-dibenzo[a,d]cycloheptene (dibenzosuberane). <i>Journal of Chemical Thermodynamics</i> , 2011, 43, 364-370.	2.0	8
47	Experimental and computational thermochemical study of oxindole. <i>Journal of Chemical Thermodynamics</i> , 2010, 42, 1101-1106.	2.0	13
48	Standard molar enthalpy of formation of 1-benzosuberone: An experimental and computational study. <i>Journal of Chemical Thermodynamics</i> , 2010, 42, 1094-1100.	2.0	8
49	Energetics of Coumarin and Chromone. <i>Journal of Physical Chemistry B</i> , 2009, 113, 11216-11221.	2.6	38
50	3,4,5-Trimethoxyphenol: A combined experimental and theoretical thermochemical investigation of its antioxidant capacity. <i>Journal of Chemical Thermodynamics</i> , 2008, 40, 625-631.	2.0	20
51	Calorimetric and Computational Thermochemical Study of 3,3-Tetramethyleneglutaric Acid, 3,3-Tetramethyleneglutaric Anhydride, and 3,3-Tetramethyleneglutarimide. <i>Journal of Physical Chemistry A</i> , 2008, 112, 10053-10058.	2.5	5
52	Thermochemical Parameters of the Chloronitrophenol Isomers: A Combined Experimental and Theoretical Investigation. <i>Journal of Chemical &amp; Engineering Data</i> , 2007, 52, 627-634.	1.9	20
53	The Experimental and Calculational Thermochemistry of 1,2,4,5-Benzenetetracarboxylic Dianhydride: Is This 10 $\pi$ Multiring Species Aromatic?. <i>Journal of Physical Chemistry A</i> , 2007, 111, 7181-7188.	2.5	15
54	Calorimetric and Computational Study of Indanones. <i>Journal of Physical Chemistry A</i> , 2007, 111, 11153-11159.	2.5	20

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55	Experimental and computational thermochemistry of three nitrogen-containing heterocycles: 2-benzimidazolinone, 2-benzoxazolinone and 3-indazolinone. <i>Molecular Physics</i> , 2006, 104, 325-334.	1.7	38
56	Benzanilide: on the crossroads of calorimetry, computations and concepts. <i>Molecular Physics</i> , 2006, 104, 2855-2860.	1.7	10
57	Calorimetric and computational study of 2H-1, 4-benzoxazin-3(4H)-one and of related species. <i>Molecular Physics</i> , 2006, 104, 1833-1841.	1.7	7
58	Thermochemical study of cyanopyrazines: Experimental and theoretical approaches. <i>Journal of Chemical Thermodynamics</i> , 2006, 38, 559-564.	2.0	8
59	Experimental and computational thermochemistry of the dihydroxypyridine isomers. <i>Journal of Chemical Thermodynamics</i> , 2006, 38, 450-454.	2.0	11
60	Experimental thermochemical study of three monosubstituted pyrazines. <i>Journal of Chemical Thermodynamics</i> , 2005, 37, 49-53.	2.0	19
61	Thermochemistry of diphenic anhydride. A combined experimental and theoretical study. <i>Molecular Physics</i> , 2005, 103, 1885-1894.	1.7	9
62	Saccharin: a combined experimental and computational thermochemical investigation of a sweetener and sulfonamide. <i>Molecular Physics</i> , 2005, 103, 221-228.	1.7	32
63	The Enthalpies of Formation of o-, m-, and p-Benzoquinone: Gas-Phase Ion Energetics, Combustion Calorimetry, and Quantum Chemical Computations Combined. <i>Journal of the American Chemical Society</i> , 2005, 127, 6116-6122.	13.7	72
64	Thermochemical Study of the Cyanophenol Isomers. <i>Structural Chemistry</i> , 2004, 15, 103-116.	2.0	24
65	Standard molar enthalpies of formation of the methoxynitrophenol isomers: a combined experimental and theoretical investigation. <i>Journal of Chemical Thermodynamics</i> , 2004, 36, 431-436.	2.0	23
66	Aspects of the Aromaticity of Anthranil. <i>European Journal of Organic Chemistry</i> , 2004, 2004, 3340-3345.	2.4	28
67	Experimental and theoretical study of the dissociation enthalpy of the N=O bond on 2-hydroxypyridine N-oxide: theoretical analysis of the energetics of the N=O bond for hydroxypyridine N-oxide isomers. <i>Journal of Chemical Thermodynamics</i> , 2004, 36, 107-113.	2.0	14
68	Thermochemical study of chloropyrazines and chloroquinoxalines. <i>Journal of Chemical Thermodynamics</i> , 2004, 36, 377-383.	2.0	13
69	Surprises with strain energy and sulpholane (tetrahydrothiophene 1,1-dioxide): a combined experimental and theoretical investigation. <i>Molecular Physics</i> , 2004, 102, 525-530.	1.7	9
70	Thermochemistry of biphenylcarboxylic and dicarboxylic acids. A combined experimental and theoretical study. This paper is dedicated to Kurt Mislow, a pioneering biphenyl stereochemist, on the occasion of his 80th birthday. <i>Organic and Biomolecular Chemistry</i> , 2004, 2, 1353.	2.8	15
71	The energetics of isomeric benzoxazine diones: isatoic anhydride revisited. <i>Organic and Biomolecular Chemistry</i> , 2004, 2, 1647.	2.8	16
72	Thermochemical Study of the Methoxy- and Dimethoxyphenol Isomers. <i>Journal of Chemical &amp; Engineering Data</i> , 2003, 48, 669-679.	1.9	50

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73	Are isatin and isatoic anhydride antiaromatic and aromatic respectively? A combined experimental and theoretical investigation. <i>Organic and Biomolecular Chemistry</i> , 2003, 1, 2566.	2.8	40
74	Thermochemical study of the ethylpyridine and ethylpyrazine isomers. <i>Organic and Biomolecular Chemistry</i> , 2003, 1, 4329.	2.8	19
75	Thermochemistry of (E)- and (Z)-disubstituted alkene species: a combined experimental and theoretical investigation of isomeric dimethyl fumarate and dimethyl maleate. <i>Organic and Biomolecular Chemistry</i> , 2003, 1, 2930.	2.8	17
76	Title is missing!. <i>Structural Chemistry</i> , 2001, 12, 171-181.	2.0	21