Anna Spalletti

List of Publications by Year
in descending order

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121 papers	2,630 citations	172457 29 h-index	276875 41 g-index
127 all docs	127 docs citations	127 times ranked	1583 citing authors

#	Article	IF	Citations
1	Exploring a new class of singlet fission fluorene derivatives with high-energy triplets. Chemical Science, 2022, 13, 2071-2078.	7.4	3
2	Acidâ€“base strength and acido(fluoro)chromism of three pushâ€“pull derivatives of 2,6-distyrylpyridine. Photochemical and Photobiological Sciences, 2022, 21, 935-947.	2.9	5
3	Amphiphilicity-Controlled Localization of Red Emitting Bicationic Fluorophores in Tumor Cells Acting as Bio-Probes and Anticancer Drugs. Molecules, 2022, 27, 3713.	3.8	3
4	The role of twisting in driving excited-state symmetry breaking and enhanced two-photon absorption in quadrupolar cationic pyridinium derivatives. Physical Chemistry Chemical Physics, 2021, 23, 16739-16753.	2.8	9
5	Metal complexes with sterically demanding phenanthroline ligands: A combined spectroscopic study. Dyes and Pigments, 2021, 187, 109150.	3.7	3
6	Nonlinear optical properties of a new panchromatic series of water-soluble bicationic push-pull fluorophores. Dyes and Pigments, 2021, 194, 109620.	3.7	15
7	Competition between fluorescence and triplet production ruled by nitro groups in one-arm and two-arm styrylbenzene heteroanalogues. Photochemical and Photobiological Sciences, 2020, 19, 1665-1676.	2.9	23
8	Uncovering Structureâ€“Property Relationships in Pushâ€“Pull Chromophores: A Promising Route to Large Hyperpolarizability and Two-Photon Absorption. Journal of Physical Chemistry C, 2020, 124, 15739-15748.	3.1	27
9	Energy-Transfer and Charge-Transfer Dynamics in Highly Fluorescent Naphthalimideâ€“BODIPY Dyads: Effect of BODIPY Orientation. Journal of Physical Chemistry C, 2019, 123, 24362-24374.	3.1	25
10	In memory of Professor Ugo Mazzucato (1929â€“2017). Photochemical and Photobiological Sciences, 2019, 18, 2092-2093.	2.9	0
11	Effect of the size of polycyclic aryl groups on the competition between adiabatic/diabatic photoisomerization mechanisms of cis-styrylarenes. Photochemical and Photobiological Sciences, 2019, 18, 2125-2135.	2.9	0
12	Four styryl phenanthroline derivatives as excellent acidochromic probes. Dyes and Pigments, 2019, 162, 440-450.	3.7	11
13	Fluorosolvatochromism and hyperpolarizability of one-arm and two-arms nitro-compounds bearing heterocyclic rings. Journal of Photochemistry and Photobiology A: Chemistry, 2019, 368, 190-199.	3.9	23
14	Photoinduced ICT vs. excited rotamer intercoversion in two quadrupolar polyaromatic N-methylpyridinium cations. Physical Chemistry Chemical Physics, 2018, 20, 2851-2864.	2.8	18
15	Evaluation of Hyperpolarizability from the Solvatochromic Method: Thiophene Containing Pushâ€“Pull Cationic Dyes as a Case Study. Journal of Physical Chemistry C, 2018, 122, 2285-2296.	3.1	32
16	New Styryl Phenanthroline Derivatives as Model Dâ^'Ï€â^'Aâ^'Ï€â^'D Materials for Nonâ€Linear Optics. ChemPhysChem, 2018, 19, 1917-1929.	2.1	20
17	Spectral properties and photoreactivity of sydnonyl-stilbenes. Journal of Photochemistry and Photobiology A: Chemistry, 2018, 351, 124-130.	3.9	3
18	Photoinduced Intramolecular Charge Transfer and Hyperpolarizability Coefficient in Pushâ€Pull Pyridinium Salts with Increasing Strength of the Acceptor Group. ChemPlusChem, 2018, 83, 1021-1031.	2.8	18

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19	A cationic naphthyl derivative defies the non-equilibrated excited rotamers principle. Physical Chemistry Chemical Physics, 2017, 19, 5262-5272.	2.8	8
20	Enhancement of Two-Photon Absorption Parallels Intramolecular Charge-Transfer Efficiency in Quadrupolar versus Dipolar Cationic Chromophores. Journal of Physical Chemistry C, 2017, 121, 3987-4001.	3.1	30
21	Optical Communication among Oscillatory Reactions and Photoâ€Excitable Systems: UV and Visible Radiation Can Synchronize Artificial Neuron Models. Angewandte Chemie - International Edition, 2017, 56, 7535-7540.	13.8	43
22	Optical Communication among Oscillatory Reactions and Photoâ€Excitable Systems: UV and Visible Radiation Can Synchronize Artificial Neuron Models. Angewandte Chemie, 2017, 129, 7643-7648.	2.0	3
23	Proton transfer in the ground and excited state and photobehaviour of the positional isomers of E -5-[2-(pyrid- n -yl)ethenyl]oxazoleâ€™s (n = 2, 3 and 4). Journal of Photochemistry and Photobiology A: Chemistry, 2017, 333, 33-39.	3.9	2
24	Effect of the positional isomerism on the photoreactivity of styryloxazoles. Journal of Photochemistry and Photobiology A: Chemistry, 2016, 316, 95-103.	3.9	8
25	Efficient Excitedâ€State Symmetry Breaking in a Cationic Quadrupolar System Bearing Diphenylamino Donors. ChemPhysChem, 2016, 17, 136-146.	2.1	42
26	Photobehavior and Nonlinear Optical Properties of Pushâ€“Pull, Symmetrical, and Highly Fluorescent Benzothiadiazole Derivatives. Journal of Physical Chemistry C, 2016, 120, 23726-23739.	3.1	52
27	A two excited state model to explain the peculiar photobehaviour of a flexible quadrupolar Dâ€“Ïâ€“D anthracene derivative. Physical Chemistry Chemical Physics, 2016, 18, 23389-23399.	2.8	16
28	Deactivating effect of the pyridine n,Ï€* states on the photoreactivity of 5-[2-(pyrid-n-yl)ethenyl]oxazole (n= 2, 3 and 4). Journal of Photochemistry and Photobiology A: Chemistry, 2016, 329, 262-272.	3.9	7
29	Twisting in the excited state of an N-methylpyridinium fluorescent dye modulated by nano-heterogeneous micellar systems. Photochemical and Photobiological Sciences, 2016, 15, 525-535.	2.9	11
30	Unexpected multiple activated steps in the excited state decay of some bis(phenylethynyl)-fluorenes and -anthracenes. Physical Chemistry Chemical Physics, 2016, 18, 285-294.	2.8	3
31	The role of acrylic acid impurity as a sensitizing component in electrocardiogram electrodes. Contact Dermatitis, 2015, 73, 44-48.	1.4	34
32	Spectroscopic and Photophysical Characterization of Acetylenic Fluorophores: The Role of the Proximity Effect on Increasing Internal Conversion. ChemPlusChem, 2015, 80, 1045-1051.	2.8	4
33	Presence of Two Emissive Minima in the Lowest Excited State of a Pushâ€“Pull Cationic Dye Unequivocally Proved by Femtosecond Up-Conversion Spectroscopy and Vibronic Quantum-Mechanical Computations. Journal of Physical Chemistry B, 2015, 119, 6035-6040.	2.6	37
34	Inclusion of Two Pushâ€“Pull N-Methylpyridinium Salts in Anionic Surfactant Solutions: A Comprehensive Photophysical Investigation. Journal of Physical Chemistry B, 2015, 119, 6658-6667.	2.6	16
35	Spectroscopic Investigation of Interactions of New Potential Anticancer Drugs with DNA and Non-Ionic Micelles. Journal of Physical Chemistry B, 2015, 119, 1483-1495.	2.6	27
36	Acidâ€“Base Strength and Acidochromism of Some Dimethylaminoâ€“Azinium Iodides. An Integrated Experimental and Theoretical Study. Journal of Physical Chemistry A, 2015, 119, 323-333.	2.5	23

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37	Effect of the Ï€ Bridge and Acceptor on Intramolecular Charge Transfer in Pushâ€“Pull Cationic Chromophores: An Ultrafast Spectroscopic and TDâ€DFT Computational Study. ChemPhysChem, 2015, 16, 1440-1450.	2.1	40
38	An ultrafast spectroscopic and quantum mechanical investigation of multiple emissions in pushâ€“pull pyridinium derivatives bearing different electron donors. Physical Chemistry Chemical Physics, 2015, 17, 20981-20989.	2.8	30
39	Inclusion of pushâ€“pull N-methylpyridinium salts within surfactant hydrogels: is their excited state intramolecular charge transfer mediated by twisting?. Physical Chemistry Chemical Physics, 2015, 17, 17214-17220.	2.8	18
40	Unusual high fluorescence of two nitro-distyrylbenzene-like compounds induced by CT processes affecting the fluorescence/intersystem-crossing competition. Physical Chemistry Chemical Physics, 2015, 17, 14740-14749.	2.8	37
41	Experimental evidence of dual emission in a negatively solvatochromic pushâ€“pull pyridinium derivative. Physical Chemistry Chemical Physics, 2015, 17, 1877-1882.	2.8	48
42	Photoisomerization and Photocyclization of 5-Styryloxazole. Croatica Chemica Acta, 2014, 87, 327-333.	0.4	8
43	Quantitative cascade energy transfer in semiconductor thin films. Photochemical and Photobiological Sciences, 2014, 13, 1031-1038.	2.9	6
44	Photobehaviour of methyl-pyridinium and quinolinium iodide derivatives, free and complexed with DNA. A case of bisintercalation. Photochemical and Photobiological Sciences, 2014, 13, 939-950.	2.9	19
45	Photoinduced symmetry-breaking intramolecular charge transfer in a quadrupolar pyridinium derivative. Physical Chemistry Chemical Physics, 2014, 16, 13984-13994.	2.8	62
46	Intramolecular Charge Transfer of Pushâ€“Pull Pyridinium Salts in the Triplet Manifold. Journal of Physical Chemistry A, 2014, 118, 7782-7787.	2.5	32
47	Intramolecular Charge Transfer of Pushâ€“Pull Pyridinium Salts in the Singlet Manifold. Journal of Physical Chemistry A, 2014, 118, 3580-3592.	2.5	77
48	Photochemical and photobiological studies on furoquinazolines as new psoralen analogs. Journal of Photochemistry and Photobiology B: Biology, 2014, 138, 43-54.	3.8	9
49	Spectral properties and photophysics of arylacetylenes in thin films. Organic Photonics and Photovoltaics, 2014, 2, .	1.3	4
50	Photophysics of Pushâ€“Pull Distyrylfurans, Thiophenes and Pyridines by Fast and Ultrafast Techniques. ChemPhysChem, 2013, 14, 970-981.	2.1	32
51	Intramolecular charge transfer, solvatochromism and hyperpolarizability of compounds bearing ethenylene or ethynylene bridges. Chemical Physics, 2012, 407, 9-19.	1.9	104
52	Synthesis, spectral properties and photobehaviour of pushâ€“pull distyrylbenzene nitro-derivatives. Journal of Photochemistry and Photobiology A: Chemistry, 2012, 244, 38-46.	3.9	27
53	Comprehensive Photophysical Behaviour of Ethynyl Fluorenes and Ethynyl Anthracenes Investigated by Fast and Ultrafast Timeâ€Resolved Spectroscopy. ChemPhysChem, 2012, 13, 724-735.	2.1	40
54	A peculiar dependence of intersystem crossing of p-nitro-2,5-distyrylfuran on the dielectric properties of the solvent. Physical Chemistry Chemical Physics, 2011, 13, 20787.	2.8	16

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55	Ultrafast photoinduced intramolecular charge transfer in pushâ€“pull distyryl furan and benzofuran: solvent and molecular structure effect. Physical Chemistry Chemical Physics, 2011, 13, 4519.	2.8	52
56	Photobehaviour and DNA interaction of styrylquinolinium salts bearing thiophene substituents. Photochemical and Photobiological Sciences, 2011, 10, 973-979.	2.9	17
57	Photochemistry and DNA-affinity of some pyrimidine-substituted styryl-azinium iodides. Photochemical and Photobiological Sciences, 2011, 10, 1830-1836.	2.9	22
58	Synthesis and photobehaviour of donor-Ï€-acceptor conjugated arylacetylenes. Journal of Photochemistry and Photobiology A: Chemistry, 2011, 223, 140-148.	3.9	29
59	Spectral properties and photobehaviour of 2,5-distyrylfuran derivatives. Journal of Photochemistry and Photobiology A: Chemistry, 2011, 219, 1-9.	3.9	21
60	Induced phosphorescence of some aza- and thio-stilbenes embedded in thallium-exchanged zeolites. Journal of Luminescence, 2011, 131, 1193-1197.	3.1	6
61	Synthesis and photophysical properties of conjugated anthracene-based compounds. Journal of Photochemistry and Photobiology A: Chemistry, 2010, 211, 162-169.	3.9	19
62	Photochemistry and DNA-affinity of some stilbene and distyrylbenzene analogues containing pyridinium and imidazolium iodides. Journal of Photochemistry and Photobiology A: Chemistry, 2010, 216, 66-72.	3.9	33
63	Photobehavior of the Geometrical Isomers of Two 1,4-Distyrylbenzene Analogues with Side Groups of Different Electron Donor/Acceptor Character. Journal of Physical Chemistry A, 2010, 114, 10761-10768.	2.5	26
64	Fluorescence/photoisomerization competition in trans-aza-1,2-diarylethenes. Journal of Fluorescence, 2009, 19, 759-768.	2.5	10
65	Novel conjugated Ï‰,Ï‰â€²-diaryl hexatriene derivatives with the central double bond in the benzofuran ring and their photochemical and photophysical properties. Journal of Photochemistry and Photobiology A: Chemistry, 2009, 202, 136-141.	3.9	19
66	Spectra and photophysics of new organic fluorophores: 2,3-Di(phenylethenyl)benzofuran derivatives. Chemical Physics, 2009, 361, 61-67.	1.9	42
67	Competition between Photoisomerization and Photocyclization of the Cis Isomers of n-Styrylnaphthalenes and -Phenanthrenes. Journal of Physical Chemistry A, 2009, 113, 14521-14529.	2.5	17
68	Adiabatic Pathways in the Conformational and Geometrical Photoisomerizations of the 1,2-Distyrylbenzene Isomers. Journal of Physical Chemistry A, 2009, 113, 8557-8568.	2.5	7
69	cis peak as probe to investigate the molecular structure. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2008, 71, 543-549.	3.9	11
70	Excited state behaviour of some thio-analogues of 1,3-distyrylbenzene. Journal of Photochemistry and Photobiology A: Chemistry, 2008, 196, 233-238.	3.9	9
71	Photoisomerization mechanism of the cis isomers of 1,2-distyrylbenzene and two hetero-analogues. Journal of Photochemistry and Photobiology A: Chemistry, 2008, 195, 301-306.	3.9	15
72	Photobehaviour of thio-analogues of stilbene and 1,4-distyrylbenzene studied by time-resolved absorption techniques. Chemical Physics, 2008, 352, 28-34.	1.9	19

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73	Structures, spectra and photophysics of new organic fluorophores: 2,3- and 2,5-di(phenylethenyl)furan. Chemical Physics, 2008, 353, 163-169.	1.9	16
74	S0â†'Sn and S1â†'Sn absorption spectra of thio-distyrylbenzenes. Chemical Physics, 2007, 337, 168-176.	1.9	16
75	Role of adiabatic pathways in the photoisomerization of aromatic olefins. Inorganica Chimica Acta, 2007, 360, 961-969.	2.4	17
76	Heteroatom effect on the radiative and reactive photobehaviour of E,E-1,2-distyrylbenzene. Journal of Photochemistry and Photobiology A: Chemistry, 2007, 187, 325-331.	3.9	10
77	Conformational equilibria in EE-1,3-di-(3â€²-thienylethenyl)benzene: One-way adiabatic interconversion of rotamers in S1 and their excited state properties. Chemical Physics, 2006, 328, 275-283.	1.9	7
78	Effect of stereoisomerism on the radiative and reactive relaxation channels of two thio-analogues of distyrylbenzene. Chemical Physics, 2006, 331, 164-172.	1.9	11
79	Triplet-sensitized photobehaviour of the three stereoisomers of 1,4-distyrylbenzene and some aza-analogues. Journal of Photochemistry and Photobiology A: Chemistry, 2006, 177, 307-313.	3.9	23
80	Photobehaviour of some 1-heteroaryl-2-(1-methylpyridinium-2-yl)ethene iodides (free and complexed) Tj ETQq0 0 0 rgBT /Overlock 10 Tf	3.9	14
81	Effect of the chain length on the excited state properties of Î±-naphthyl,Ï‰-phenyl-polyenes. Journal of Photochemistry and Photobiology A: Chemistry, 2005, 174, 181-186.	3.9	4
82	Excited state properties of cross-conjugated 1,2- and 1,3-distyrylbenzene and some aza-analogues. Chemical Physics, 2005, 312, 205-211.	1.9	30
83	Competitive radiative and reactive relaxation channels in the excited state decay of some thio-analogues of EE-distyrylbenzene. Photochemical and Photobiological Sciences, 2005, 4, 547.	2.9	30
84	Protonation effect on the excited state behaviour of some aza-analogues ofEE-distyrylbenzene. International Journal of Photoenergy, 2004, 6, 241-250.	2.5	8
85	Effect of thienyl groups on the photoisomerization and rotamerism of symmetric and asymmetric diaryl-ethenes and diaryl-butadienesElectronic supplementary information (ESI) available: (1) Calculated electronic spectra (transition energy and oscillator strength) and ground state total energy of the rotamers of the trans isomers; (2) Absorption and emission spectra. See http://www.rsc.org/suppdata/pp/b3/b309267d/. Photochemical and Photobiological Sciences, 2004, 3,	2.9	22
86	Photobehaviour of di(phenylbutadienyl)arenes containing nitrogen and sulfur heteroatomsElectronic supplementary information (ESI) available: calculated electronic spectra (transition energy and oscillator strength) and heats of formation in the ground state of the elongated and compressed conformations of the investigated compounds. See http://www.rsc.org/suppdata/pp/b3/b309267d/. Photochemical and Photobiological Sciences, 2004, 3, 205.	2.9	9
87	Conformer-specific and two-fold adiabatic photoisomerization of ZZ-1,4-di-(2-quinolylethenyl)benzene. Photochemical and Photobiological Sciences, 2004, 3, 695.	2.9	15
88	Rotamerism and electronic spectra of aza-derivatives of stilbene and diphenylbutadiene. A combined experimental and theoretical study. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2003, 59, 75-86.	3.9	29
89	Protonation effect on the excited state behaviour of EE-1-(n-pyridyl)-4-phenylbutadienes (n = 2, 3 and) Tj ETQq1 1 0.784314 rgBT /Ov and Photobiological Sciences, 2003, 2, 282.	2.9	9
90	Effect of the Nitrogen Heteroatom on the Excited State Properties of 1,4-Distyrylbenzene. Journal of Physical Chemistry A, 2003, 107, 11231-11238.	2.5	34

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91	Photoisomerization mechanisms and photoselectivity of the stereoisomers of 1-(pyrid-n-yl),4-phenylbuta-1,3-diene. Physical Chemistry Chemical Physics, 2002, 4, 2911-2916.	2.8	14
92	Photophysical and Photochemical Behavior of the Three Conformational Isomers oftrans-1,2-di(2-Naphthyl)ethene in Nonpolar Solvent. Journal of Physical Chemistry A, 2002, 106, 7068-7074.	2.5	16
93	Excited-State Properties of the Four Stereoisomers of 1-(9-Anthryl)-4-phenyl-1,3-butadiene:Â Evidence of Adiabatic and Diabatic Deactivation Pathways. Journal of Physical Chemistry A, 2002, 106, 11400-11407.	2.5	3
94	Conformational equilibria in EE-2,6-di-[2-(furan-2-yl)vinyl]pyridine controlled by intramolecular hydrogen-type bonds. Journal of Molecular Structure, 2002, 612, 339-347.	3.6	13
95	Photophysics and photochemistry of the EE and ZE isomers of 1-(n-pyridyl)-4-phenyl-1,3-butadiene (n = 2,)	2.8	19
96	Effect of the nature of aryl and heteroaryl groups on the excited state properties of asymmetric 1,4-diarylbutadienes. Chemical Physics, 2001, 272, 213-225.	1.9	18
97	Effect of solvent polarizability on dual fluorescence of EE-1-phenyl,4-(1â€²-pyrenyl)-1,3-butadiene. Chemical Physics, 2000, 260, 383-390.	1.9	14
98	Photophysics and photochemistry of 2,6-distyrylpyridine and some heteroanalogues. Physical Chemistry Chemical Physics, 2000, 2, 4005-4012.	2.8	34
99	Temperature and solvent effects on rotamer-specific photobehaviour of the cis and trans isomers of 2-styrylanthracene. Physical Chemistry Chemical Physics, 1999, 1, 5623-5632.	2.8	16
100	Spectral Characterization, Photophysics, and Photochemistry of the Four Stereoisomers of 1-(2-anthryl)-4-phenyl-1,3-butadiene. Journal of Physical Chemistry A, 1999, 103, 8994-9002.	2.5	11
101	Excited-State Behavior of Someall-trans-Î±,ï‰-Dithienylpolyenes. Journal of the American Chemical Society, 1999, 121, 1065-1075.	13.7	46
102	Synthesis and basicity of 2,6-di-[2-(heteroaryl)vinyl]pyridines. Tetrahedron, 1998, 54, 9721-9730.	1.9	10
103	Effect of pyridyl and thienyl groups on the excited state properties of stilbene-like molecules. Journal of Chemical Sciences, 1998, 110, 297-310.	1.5	12
104	Temperature effects on the photoreactivity and rotamerism of (Z)-1-styrylanthracene in non-polar and polar solvents. Journal of the Chemical Society, Faraday Transactions, 1997, 93, 211-219.	1.7	8
105	A theoretical and experimental study of the excited state relaxation properties of mono-aza- and di-aza-trans-stilbenes. AIP Conference Proceedings, 1996, , .	0.4	5
106	Spectral and photophysical properties of trans-2-styrylanthracene rotamers, derived by kinetic fluorescence analysis. A comparison with the results obtained by statistical procedures. Chemical Physics, 1996, 202, 367-376.	1.9	13
107	Excited state behaviour of some trans-stilbene analogues bearing thiophene rings. Journal of Photochemistry and Photobiology A: Chemistry, 1996, 100, 57-64.	3.9	36
108	Solvent and temperature effects on the fluorescence and competitive photoreactions of cis-9-styrylanthracene. Research on Chemical Intermediates, 1995, 21, 735-747.	2.7	16

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109	Effect of the nitrogen heteroatom on the photophysics and photochemistry of transâ€1â€styrylnaphthalene and transâ€9â€styrylphenanthrene in different solvents. Recueil Des Travaux Chimiques Des Pays-Bas, 1995, 114, 459-464.	0.0	24
110	Role of internal conversion on the excited state properties of trans-styrylpyridines. Chemical Physics, 1995, 196, 383-393.	1.9	48
111	Evidence of adiabatic channels in the singlet photoisomerization of cis-1,2-diarylethenes: a fluorimetric study. Coordination Chemistry Reviews, 1993, 125, 251-260.	18.8	34
112	Conformational equilibria in trans-diarylethylenes: spectral and photophysical properties of rotamers of 1-(2-naphthyl)-2-(6â€²-quinolyl)ethylene, derived from kinetic and statistical fluorescence analysis. Journal of Molecular Structure, 1993, 298, 165-175.	3.6	6
113	Effect of the nature of the aromatic groups on the lowest excited states of trans-1,2-diarylethenes. Journal of the Chemical Society, Faraday Transactions, 1992, 88, 3139.	1.7	50
114	Principal-component self-modeling analysis of fluorescence for some trans-diarylethylenes. A comparison with kinetic analysis. Chemical Physics, 1992, 160, 131-144.	1.9	30
115	Decay pathways of the first excited singlet state of cis-1-styrylpyrene. Chemical Physics Letters, 1991, 186, 297-302.	2.6	25
116	The three-component fluorescence emission of trans-2-styrylanthracene in fluid solution. The implication of an upper excited singlet state. Spectrochimica Acta Part A: Molecular Spectroscopy, 1990, 46, 413-418.	0.1	22
117	A photophysical and theoretical study of styrylanthracenes. Journal of the Chemical Society, Faraday Transactions 2, 1988, 84, 385.	1.1	44
118	Photophysical and theoretical studies of photoisomerism and rotamerism of trans-styrylphenanthrenes. The Journal of Physical Chemistry, 1987, 91, 4733-4743.	2.9	109
119	Rotamerism and transâ€“cis photoisomerization of 1-(2-naphthyl)-2-(nâ€²-pyridyl)ethylenes studied by stationary and pulsed fluorescence techniques. Journal of the Chemical Society, Faraday Transactions 2, 1986, 82, 775-788.	1.1	19
120	Excited State Reactivity of Aza-Aromatics. Zeitschrift Fur Physikalische Chemie, 1983, 138, 199-206.	2.8	6
121	Photochemical and Photophysical Behaviour of 9-Styrylphenanthrene and its Aza-Analogues. Zeitschrift Fur Physikalische Chemie, 1982, 133, 107-118.	2.8	17