

Juan Teodomiro LÃ³pez Navarrete

List of Publications by Year in descending order

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263
papers

9,452
citations

36303

51
h-index

54911

84
g-index

263
all docs

263
docs citations

263
times ranked

9421
citing authors

#	ARTICLE	IF	CITATIONS
1	Backbone Configuration and Electronic Property Tuning of Imide-Functionalized Ladder-Type Heteroarenes-Based Polymer Acceptors for Efficient All-Polymer Solar Cells. <i>Advanced Functional Materials</i> , 2022, 32, .	14.9	12
2	Poly(3-hexylthiophene-2,5-diyl): Evidence of different polymer chain conformations in the solid state from a combined study of regioregularity control and Raman spectroscopy. <i>Journal of Molecular Structure</i> , 2020, 1221, 128882.	3.6	4
3	Fluorene-Based Donor-Acceptor Copolymers Containing Functionalized Benzotriazole Units: Tunable Emission and their Electrical Properties. <i>Polymers</i> , 2020, 12, 256.	4.5	6
4	Effect of the Linkage Position on the Conjugation Length of Truxene-Based Porous Polymers: Implications for Their Sensing Performance of Nitroaromatics. <i>Chemistry of Materials</i> , 2019, 31, 6971-6978.	6.7	21
5	(Semi)ladder-Type Bithiophene Imide-Based All-Acceptor Semiconductors: Synthesis, Structure-Property Correlations, and Unipolar n-Type Transistor Performance. <i>Journal of the American Chemical Society</i> , 2018, 140, 6095-6108.	13.7	178
6	Infrared and multi-wavelength Raman spectroscopy of regio-regular P3HT and its deuterio derivatives. <i>Journal of Raman Spectroscopy</i> , 2018, 49, 569-580.	2.5	16
7	Ladder-Type Heteroarenes: Up to 15 Rings with Five Imide Groups. <i>Angewandte Chemie - International Edition</i> , 2017, 56, 9924-9929.	13.8	105
8	Functionalized branched EDOT-terthiophene copolymer films by electropolymerization and post-polymerization -click-reactions. <i>Beilstein Journal of Organic Chemistry</i> , 2015, 11, 335-347.	2.2	15
9	Robust Ethylenedioxythiophene-Vinylene Oligomers from Fragile Thiophene-Vinylene Cores: Synthesis and Optical, Chemical and Electrochemical Properties of Multicharged Shapes. <i>Chemistry - A European Journal</i> , 2015, 21, 1713-1725.	3.3	13
10	High Yield Ultrafast Intramolecular Singlet Exciton Fission in a Quinoidal Bithiophene. <i>Journal of Physical Chemistry Letters</i> , 2015, 6, 1375-1384.	4.6	106
11	On the handedness of helical aggregates of C ₃ tricarboxamides: a multichiroptical characterization. <i>Chemical Communications</i> , 2015, 51, 9781-9784.	4.1	26
12	Triindole-Bridge-Triindole Dimers as Models for Two Dimensional Microporous Polymers. <i>Organic Letters</i> , 2015, 17, 2258-2261.	4.6	18
13	Planarization, Fusion, and Strain of Carbon-Bridged Phenylenevinylene Oligomers Enhance π -Electron and Charge Conjugation: A Dissectional Vibrational Raman Study. <i>Journal of the American Chemical Society</i> , 2015, 137, 3834-3843.	13.7	44
14	Understanding the Origin of the VCD Signals on the Basis of a Nonredundant Coordinate Definition. <i>Journal of Chemical Theory and Computation</i> , 2015, 11, 2633-2641.	5.3	2
15	Polarization, second-order nonlinear optical properties and electrochromism in 4H-pyranylidene chromophores with a quinoid/aromatic thiophene ring bridge. <i>RSC Advances</i> , 2015, 5, 231-242.	3.6	35
16	Combined Raman spectroscopic and Rietveld analyses as a useful and nondestructive approach to studying flint raw materials at prehistoric archaeological sites. <i>Archaeological and Anthropological Sciences</i> , 2015, 7, 235-243.	1.8	11
17	A combined MD/QM and experimental exploration of conformational richness in branched oligothiophenes. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 24841-24852.	2.8	13
18	Branched polythiophenes by Ni-catalyzed Kumada coupling. <i>Polymer Chemistry</i> , 2014, 5, 6824-6833.	3.9	10

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19	Mode Robustness in Raman Optical Activity. <i>Journal of Chemical Theory and Computation</i> , 2014, 10, 5520-5527.	5.3	23
20	Multistep π -Dimerization of Tetrakis(<i>n</i> -decyl)heptathienoacene Radical Cations: A Combined Experimental and Theoretical Study. <i>Chemistry - A European Journal</i> , 2014, 20, 10351-10359.	3.3	12
21	Raman Spectroscopic Characterization of Polyselenophenes and Poly(3,4-ethylenedioxy-selenophene)s. <i>Israel Journal of Chemistry</i> , 2014, 54, 759-766.	2.3	8
22	Alkoxy-Functionalized Thienyl-Vinylene Polymers for Field-Effect Transistors and All-Polymer Solar Cells. <i>Advanced Functional Materials</i> , 2014, 24, 2782-2793.	14.9	83
23	Properties of Sizeable [<i>n</i>]Cycloparaphenylenes as Molecular Models of Single-Wall Carbon Nanotubes Elucidated by Raman Spectroscopy: Structural and Electron-Transfer Responses under Mechanical Stress. <i>Angewandte Chemie - International Edition</i> , 2014, 53, 7033-7037.	13.8	77
24	EDOT-Based Copolymers with Pendant Anthraquinone Units: Analysis of Their Optoelectronic Properties within the Double-Cable Context. <i>Journal of Physical Chemistry C</i> , 2014, 118, 9899-9910.	3.1	2
25	Diradicals acting through diamagnetic phenylene vinylene bridges: Raman spectroscopy as a probe to characterize spin delocalization. <i>Journal of Chemical Physics</i> , 2014, 140, 164903.	3.0	6
26	Antiaromatic bisindeno-[<i>n</i>]thienoacenes with small singlet biradical characters: syntheses, structures and chain length dependent physical properties. <i>Chemical Science</i> , 2014, 5, 4490-4503.	7.4	62
27	Unfolding Pathway of a Globular Protein by Surfactants Monitored with Raman Optical Activity. <i>Journal of Physical Chemistry Letters</i> , 2014, 5, 8-13.	4.6	9
28	Zethrene biradicals: How pro-aromaticity is expressed in the ground electronic state and in the lowest energy singlet, triplet, and ionic states. <i>Journal of Chemical Physics</i> , 2014, 140, 054706.	3.0	28
29	Phenyl- and Thienyl-Ended Symmetric Azomethines and Azines as Model Compounds for n-Channel Organic Field-Effect Transistors: An Electrochemical and Computational Study. <i>Journal of Physical Chemistry C</i> , 2014, 118, 3984-3993.	3.1	30
30	Carbon dots obtained using hydrothermal treatment of formaldehyde. Cell imaging in vitro. <i>Nanoscale</i> , 2014, 6, 9071-9077.	5.6	79
31	Symmetry Lowering in Triindoles: Impact on the Electronic and Photophysical Properties. <i>Journal of Physical Chemistry C</i> , 2014, 118, 5470-5477.	3.1	27
32	Turning on the biradical state of tetracyano-perylene and quaterrylenequinodimethanes by incorporation of additional thiophene rings. <i>Chemical Science</i> , 2014, 5, 3072-3080.	7.4	48
33	Inversion of Supramolecular Helicity in Oligo(<i>p</i> -phenylene)-Based Supramolecular Polymers: Influence of Molecular Atropisomerism. <i>Angewandte Chemie - International Edition</i> , 2014, 53, 1373-1377.	13.8	96
34	Chameleon-like behaviour of cyclo[<i>n</i>]paraphenylenes in complexes with C ₇₀ : on their impressive electronic and structural adaptability as probed by Raman spectroscopy. <i>Faraday Discussions</i> , 2014, 173, 157-171.	3.2	30
35	The unusual electronic structure of ambipolar dicyanovinyl-substituted diketopyrrolopyrrole derivatives. <i>Journal of Materials Chemistry C</i> , 2014, 2, 6376.	5.5	55
36	Tetracyanoquaterrylene and Tetracyanohexarylenequinodimethanes with Tunable Ground States and Strong Near-Infrared Absorption. <i>Angewandte Chemie - International Edition</i> , 2013, 52, 8561-8565.	13.8	94

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37	Molecular and Electronic Structure Basis of the Ambipolar Behavior of Naphthalimide-Terthiophene Derivatives: Implementation in Organic Field-Effect Transistors. <i>Chemistry - A European Journal</i> , 2013, 19, 12458-12467.	3.3	37
38	Push-pull systems bearing a quinoid/aromatic thieno[3,2-b]thiophene moiety: synthesis, ground state polarization and second-order nonlinear properties. <i>Organic and Biomolecular Chemistry</i> , 2013, 11, 6338.	2.8	25
39	Polymer solar cells with enhanced fill factors. <i>Nature Photonics</i> , 2013, 7, 825-833.	31.4	887
40	Interpretation of the infrared and Raman spectra of zwitterionic push-pull dyes based on quinoidal thiazole. <i>Journal of Molecular Structure</i> , 2013, 1044, 55-60.	3.6	2
41	Novel Thiophene-Phenylene-Thiophene Fused Bis lactam-Based Donor-Acceptor Type Conjugate Polymers: Synthesis by Direct Arylation and Properties. <i>Macromolecules</i> , 2013, 46, 9220-9230.	4.8	41
42	Radical cations of end-capped tetrathienoacenes and their π -dimerization controlled by the nature of λ -substituents and counterion concentration. <i>RSC Advances</i> , 2013, 3, 25644.	3.6	9
43	The first chiral Raman spectrum report of a protein: a perspective of 20 years. <i>Chemical Communications</i> , 2013, 49, 8893.	4.1	8
44	Designing new symmetrical facial oligothiophene amphiphiles. <i>Organic and Biomolecular Chemistry</i> , 2013, 11, 8435.	2.8	7
45	Impact of the Synergistic Collaboration of Oligothiophene Bridges and Ruthenium Complexes on the Optical Properties of Dumbbell-Shaped Compounds. <i>Chemistry - A European Journal</i> , 2013, 19, 1476-1488.	3.3	9
46	Evidence for Multicenter Bonding in Dianionic Tetracyanoethylene Dimers by Raman Spectroscopy. <i>Angewandte Chemie - International Edition</i> , 2013, 52, 6421-6425.	13.8	33
47	Influence of Processing Solvents on Optical Properties and Morphology of a Semicrystalline Low Bandgap Polymer in the Neutral and Charged States. <i>Macromolecules</i> , 2013, 46, 4924-4931.	4.8	36
48	Pushing Extended π -Quinodimethanes to the Limit: Stable Tetracyano-oligo(π -annulated) Tj ETQqO O O rgBT /Overlock 10 2013, 135, 6363-6371.	13.7	170
49	Amplified Spontaneous Emission in Pentathienoacene Dioxides by Direct Optical Pump and by Energy Transfer: Correlation with Photophysical Parameters. <i>Advanced Optical Materials</i> , 2013, 1, 588-599.	7.3	11
50	Thermomagnetic Molecular System Based on TTF-PTM Radical: Switching the Spin and Charge Delocalization. <i>Journal of Physical Chemistry Letters</i> , 2013, 4, 2721-2726.	4.6	32
51	Electropolymerized Three-Dimensional Randomly Branched EDOT-Containing Copolymers. <i>Langmuir</i> , 2013, 29, 15463-15473.	3.5	21
52	Linear and Nonlinear Optical Properties of Ramified Hexaazatriphenylenes: Charge Transfer Contributions to the Octupolar Response. <i>Journal of Physical Chemistry C</i> , 2013, 117, 626-632.	3.1	18
53	Interplay of λ , λ versus λ , λ Conjugation in the Excited States and Charged Defects of Branched Oligothiophenes as Models for Dendrimeric Materials. <i>Chemistry - A European Journal</i> , 2013, 19, 17165-17171.	3.3	8
54	Carbon-Bridged Oligo(phenylenevinylene)s: Stable π -Systems with High Responsiveness to Doping and Excitation. <i>Journal of the American Chemical Society</i> , 2012, 134, 19254-19259.	13.7	87

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55	Carbonyl-Functionalized Quaterthiophenes: A Study of the Vibrational Raman and Electronic Absorption/Emission Properties Guided by Theoretical Calculations. <i>ChemPhysChem</i> , 2012, 13, 168-176.	2.1	8
56	Molecular tuning in highly fluorescent dithieno[3,2-b:2'-3'-d]pyrrole-based oligomers: effects of N-functionalization and terminal aryl unit. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 6101.	2.8	36
57	Kinetically Blocked Stable Heptazethrene and Octazethrene: Closed-Shell or Open-Shell in the Ground State?. <i>Journal of the American Chemical Society</i> , 2012, 134, 14913-14922.	13.7	256
58	Vibrational Circular Dichroism Shows Reversible Helical Handedness Switching in Peptidomimetic l-Valine Fibrils. <i>Journal of Physical Chemistry Letters</i> , 2012, 3, 2120-2124.	4.6	21
59	Delocalization-to-Localization Charge Transition in Diferrocenyl-Oligothiophene-Vinylene Molecular Wires as a Function of the Size by Raman Spectroscopy. <i>Journal of the American Chemical Society</i> , 2012, 134, 5675-5681.	13.7	33
60	Stable Tetrabenzo-Chichibabin-TMs Hydrocarbons: Tunable Ground State and Unusual Transition between Their Closed-Shell and Open-Shell Resonance Forms. <i>Journal of the American Chemical Society</i> , 2012, 134, 14513-14525.	13.7	218
61	Organic Materials in the Undergraduate Laboratory: Microscale Synthesis and Investigation of a Donor-Acceptor Molecule. <i>Journal of Chemical Education</i> , 2012, 89, 1461-1465.	2.3	12
62	Self-Assembly Studies of a Chiral Bisurea-Based Superhydrogelator. <i>Chemistry - A European Journal</i> , 2012, 18, 14725-14731.	3.3	40
63	Conformational Control of the Electronic Properties of an $\hat{\pi}$ -Terthiophene: Lessons from a Precursor Towards Dendritic Hyperbranched Oligo- and Poly-Thiophenes. <i>ChemPhysChem</i> , 2012, 13, 3893-3900.	2.1	11
64	Bithiopheneimide-Dithienosilole/Dithienogermole Copolymers for Efficient Solar Cells: Information from Structure-Property-Device Performance Correlations and Comparison to Thieno[3,4- <i>c</i>]pyrrole-4,6-dione Analogues. <i>Journal of the American Chemical Society</i> , 2012, 134, 18427-18439.	13.7	257
65	$\hat{\pi}$ -Oligofurans show a sizeable extent of $\hat{\pi}$ -conjugation as probed by Raman spectroscopy. <i>Chemical Communications</i> , 2012, 48, 6732.	4.1	37
66	Electronic and vibrational circular dichroism spectroscopies for the understanding of chiral organization in porphyrin aggregates. <i>Chemical Communications</i> , 2012, 48, 9147.	4.1	16
67	Controlling the Macroscopic Chirality of Organic Materials Based on 1,3,5-Trialkynylbenzenes. <i>European Journal of Organic Chemistry</i> , 2012, 2012, 1577-1582.	2.4	5
68	Synthesis of the Smallest Axially Chiral Molecule by Asymmetric Carbon-Fluorine Bond Activation. <i>Angewandte Chemie - International Edition</i> , 2012, 51, 2218-2220.	13.8	43
69	Substituent and counterion effects on the formation of $\hat{\pi}$ -dimer dications of end-capped heptathienoacenes. <i>Chemical Communications</i> , 2011, 47, 12622.	4.1	14
70	Oligothiophene Tetracyanobutadienes: Alternative Donor-Acceptor Architectures for Molecular and Polymeric Materials. <i>Chemistry of Materials</i> , 2011, 23, 823-831.	6.7	42
71	Raman Optical Activity Spectra and Conformational Elucidation of Chiral Drugs. The Case of the Antiangiogenic Aerophysinin-1. <i>Journal of Physical Chemistry A</i> , 2011, 115, 2752-2755.	2.5	22
72	Two-Photon Mediated Three-Photon Fluorescence: Lessons from a Quinoidal Oligothiophene Dimer. <i>Journal of Physical Chemistry Letters</i> , 2011, 2, 2179-2183.	4.6	13

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73	The Frontiers of Quinoidal Stability in Long Oligothiophenes: Raman Spectra of Dicationic Polaron Pairs. <i>Journal of the American Chemical Society</i> , 2011, 133, 16350-16353.	13.7	55
74	Understanding Optoelectronic Properties of Cyano-Terminated Oligothiophenes in the Context of Intramolecular Charge Transfer. <i>Journal of Physical Chemistry B</i> , 2011, 115, 10573-10585.	2.6	23
75	π-conjugation and charge polarization in fluorene-dibenzothiophene-S,S-dioxide co-oligomers by Raman spectroscopy and quantum chemistry. <i>Journal of Chemical Physics</i> , 2011, 134, 044520.	3.0	13
76	On the Origin of the Chiro-Optical Activity in Supramolecular Assemblies: A Quantum Chemical Study of C ₃ Octopolar Systems. <i>Journal of Chemical Theory and Computation</i> , 2011, 7, 3314-3322.	5.3	5
77	Functionalized pentacenes: a combined theoretical, Raman and UV-Vis spectroscopic study. <i>Theoretical Chemistry Accounts</i> , 2011, 128, 521-530.	1.4	22
78	Theoretical evaluation of the nature and strength of the F ^{δ+} ⋯F ^{δ-} intermolecular interactions present in fluorinated hydrocarbons. <i>Theoretical Chemistry Accounts</i> , 2011, 128, 541-553.	1.4	58
79	The longest quinoidal oligothiophene: A Raman story. <i>Chemical Record</i> , 2011, 11, 45-53.	5.8	20
80	Enhanced Functionality for Donor-Acceptor Oligothiophenes by means of Inclusion of BODIPY: Synthesis, Electrochemistry, Photophysics, and Model Chemistry. <i>Chemistry - A European Journal</i> , 2011, 17, 498-507.	3.3	63
81	Aromatic/Proaromatic Donors in Dicyanomethylenethiazole Merocyanines: From Neutral to Strongly Zwitterionic Nonlinear Optical Chromophores. <i>Chemistry - A European Journal</i> , 2011, 17, 826-838.	3.3	64
82	Enantiopure, Monodisperse Allenyl-acetylenic Cyclooligomers: Effect of Symmetry and Conformational Flexibility on the Chiroptical Properties of Carbon-Rich Compounds. <i>Chemistry - A European Journal</i> , 2011, 17, 3876-3885.	3.3	25
83	Hexaazatriphenylene (HAT) versus tri-HAT: The Bigger the Better?. <i>Chemistry - A European Journal</i> , 2011, 17, 10312-10322.	3.3	40
84	Diferrocenyl oligothiophene wires: Raman and quantum chemical study of valence-trapped cations. <i>Journal of Chemical Physics</i> , 2011, 135, 234705.	3.0	2
85	Quinoidal Oligothiophenes: Towards Biradical Ground-State Species. <i>Chemistry - A European Journal</i> , 2010, 16, 470-484.	3.3	74
86	Neutral and Oxidized Triisopropylsilyl End-Capped Oligothienoacenes: A Combined Electrochemical, Spectroscopic, and Theoretical Study. <i>Chemistry - A European Journal</i> , 2010, 16, 5481-5491.	3.3	25
87	Comparison of Thiophene-Pyrrole Oligomers with Oligothiophenes: A Joint Experimental and Theoretical Investigation of Their Structural and Spectroscopic Properties. <i>Chemistry - A European Journal</i> , 2010, 16, 6866-6876.	3.3	27
88	Optical absorption and emission properties of end-capped oligothienoacenes: A joint theoretical and experimental study. <i>Organic Electronics</i> , 2010, 11, 1701-1712.	2.6	19
89	Mesomeric betaine chemistry in solution: Solvent effect on the structure and spectra of uracilyl-pyridinium betaine. <i>Chemical Physics</i> , 2010, 371, 1-9.	1.9	3
90	Raman Spectra and Quantum Chemistry Calculations of Fluorene-Dibenzothiophene-S,S-dioxide Oligomers. , 2010, , .		0

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91	Ultrafast and High-Contrast Electrochromism on Bendable Transparent Carbon Nanotube Electrodes. <i>Journal of Physical Chemistry Letters</i> , 2010, 1, 1367-1371.	4.6	26
92	Do [all]-S,Sâ€²-Dioxide Oligothiophenes Show Electronic and Optical Properties of Oligoenes and/or of Oligothiophenes?. <i>Journal of the American Chemical Society</i> , 2010, 132, 6231-6242.	13.7	51
93	SEIRA and SERS Effects in Cyclopentabithiophenethiol-Capped Gold Nanoparticles. <i>Journal of Physical Chemistry C</i> , 2010, 114, 12900-12904.	3.1	11
94	Tuning the Supramolecular Chirality of One- and Two-Dimensional Aggregates with the Number of Stereogenic Centers in the Component Porphyrins. <i>Journal of the American Chemical Society</i> , 2010, 132, 9350-9362.	13.7	98
95	Aggregation Behavior of a Conjugated C₃-Symmetric Molecule: A Description Based on Chiro-Optical Experimental and Theoretical Spectroscopies. <i>Journal of Physical Chemistry B</i> , 2010, 114, 5710-5717.	2.6	7
96	Ambipolar Organic Fieldâ€¢Effect Transistors from Crossâ€¢Conjugated Aromatic Quaterthiophenes; Comparisons with Quinoidal Parent Materials. <i>Advanced Functional Materials</i> , 2009, 19, 386-394.	14.9	71
97	Ferrocenylâ€¢Ended Thienoâ€¢Vinylene Oligomers: Donorâ€¢Acceptor Polarization and Mixedâ€¢Valence Properties with Emphasis on the Raman Mapping of Localizedâ€¢Delocalized Transitions. <i>Chemistry - A European Journal</i> , 2009, 15, 2548-2559.	3.3	19
98	Thiopheneâ€¢Diazine Molecular Semiconductors: Synthesis, Structural, Electrochemical, Optical, and Electronic Structural Properties; Implementation in Organic Fieldâ€¢Effect Transistors. <i>Chemistry - A European Journal</i> , 2009, 15, 5023-5039.	3.3	82
99	Oxidation of Endâ€¢Capped Pentathienoacenes and Characterization of Their Radical Cations. <i>Chemistry - A European Journal</i> , 2009, 15, 12346-12361.	3.3	17
100	Synthesis, Spectroscopy, Nonlinear Optics, and Theoretical Investigations of Thienylethynyl Octopoles with a Tunable Core. <i>Chemistry - A European Journal</i> , 2009, 15, 8223-8234.	3.3	14
101	Electronic Studies on Oligothiophenevinylenes: Understanding the Nature of Their Ground and Excited Electronic States. <i>ChemPhysChem</i> , 2009, 10, 1901-1910.	2.1	6
102	FT Raman and DFT Study on a Series of Allâ€¢anti</i> Oligothienoacenes Endâ€¢Capped with Triisopropylsilyl Groups. <i>ChemPhysChem</i> , 2009, 10, 3069-3076.	2.1	11
103	A Raman approach to pseudo-cross-conjugation in mesomeric betaines. <i>Journal of Raman Spectroscopy</i> , 2009, 40, 238-239.	2.5	5
104	Sensing properties of organised films based on a bithiophene derivative. <i>Sensors and Actuators B: Chemical</i> , 2009, 141, 625-633.	7.8	11
105	Impact of Perfluorination on the Charge-Transport Parameters of Oligoacene Crystals. <i>Journal of the American Chemical Society</i> , 2009, 131, 1502-1512.	13.7	174
106	Quantum mechanical study and vibrational spectra of indazolium-3-carboxylate and its decarboxylation product, the N-heterocyclic carbene indazol-3-ylidene. <i>Physical Chemistry Chemical Physics</i> , 2009, 11, 341-348.	2.8	5
107	Effect of ring fusion on the amplified spontaneous emission properties of oligothiophenes. <i>Journal of Materials Chemistry</i> , 2009, 19, 6556.	6.7	17
108	Raman Detection of â€¢Ambiguousâ€¢Conjugated Biradicals: Rapid Thermal Singletâ€¢Triplet Intersystem Crossing in an Extended Viologen. <i>Angewandte Chemie - International Edition</i> , 2008, 47, 1443-1446.	13.8	53

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109	A $\hat{\Gamma}^2$ -Naphthaleneimide-Modified Terthiophene Exhibiting Charge Transfer and Polarization Through the Short Molecular Axis. Joint Spectroscopic and Theoretical Study. <i>Journal of Physical Chemistry A</i> , 2008, 112, 6732-6740.	2.5	27
110	Raman Spectroscopy Shows Interchain through Space Charge Delocalization in a Mixed Valence Oligothiophene Cation and in Its $\hat{\Gamma}^2$ -Dimeric Biradicaloid Dication. <i>Journal of the American Chemical Society</i> , 2008, 130, 14028-14029.	13.7	36
111	Electronic, Optical, and Vibrational Properties of Bridged Dithienylethylene-Based NLO Chromophores. <i>Journal of Physical Chemistry C</i> , 2008, 112, 3109-3120.	3.1	48
112	Electrochemical, Magnetic, and Electrical Properties of $\hat{\Gamma}^2$ -Capped Sexithiophene Films. Part 3. Conduction in Poly(bis-terthienyl-B)s (B = Ethane, Disulfide, Diacetylene, Acetylene, Ethylene). <i>Chemistry of Materials</i> , 2008, 20, 6847-6856.	6.7	12
113	Vibrational fingerprint of the structural tuning in push-pull organic chromophores with quinoid or proaromatic spacers. <i>Journal of Chemical Physics</i> , 2007, 126, 074701.	3.0	7
114	Theoretical understanding of the increment of $\hat{\Gamma}^2$ upon protonation of pyridine peripheral octupolar molecules: Toward nonlinear optical sensors. <i>Journal of Chemical Physics</i> , 2007, 127, 164704.	3.0	11
115	NLO properties of dithienothiophene-based chromophores: a comparison study between the donor/donor and donor/acceptor substitution patterns. , 2007, , .		1
116	Tetrathiafulvalene-Based Materials for Organic Field Effect Transistors. Inspection of Their Semiconductor Properties by Means of Molecular Spectroscopy and Quantum Chemistry. <i>Journal of Physical Chemistry C</i> , 2007, 111, 10110-10118.	3.1	20
117	Linear and Nonlinear Optical Properties of Pyridine-Based Octopolar Chromophores Designed for Chemical Sensing. Joint Spectroscopic and Theoretical Study. <i>Journal of Physical Chemistry C</i> , 2007, 111, 18778-18784.	3.1	25
118	Thiophene- and Selenophene-Based Heteroacenes: $\hat{\Gamma}^2$ Combined Quantum Chemical DFT and Spectroscopic Raman and UV-Vis-NIR Study. <i>Journal of Physical Chemistry B</i> , 2007, 111, 7488-7496.	2.6	32
119	Synthesis and Doping of a Multifunctional Tetrathiafulvalene- Substituted Poly(isocyanide). <i>Macromolecules</i> , 2007, 40, 7521-7531.	4.8	54
120	Push-Pull Bithienyl Chromophore with an Unusual Transverse Path of Conjugation. <i>Journal of Physical Chemistry A</i> , 2007, 111, 841-851.	2.5	5
121	Electronic and Molecular Structures of Trigonal Truxene-Core Systems Conjugated to Peripheral Fluorene Branches. Spectroscopic and Theoretical Study. <i>Journal of Physical Chemistry B</i> , 2007, 111, 4026-4035.	2.6	36
122	Helically Annulated and Cross-Conjugated $\hat{\Gamma}^2$ -Oligothiophenes: A Fourier Transform Raman Spectroscopic and Quantum Chemical Density Functional Theory Study. <i>Journal of Physical Chemistry C</i> , 2007, 111, 4854-4860.	3.1	14
123	On the Biradicaloid Nature of Long Quinoidal Oligothiophenes: Experimental Evidence Guided by Theoretical Studies. <i>Angewandte Chemie - International Edition</i> , 2007, 46, 9057-9061.	13.8	143
124	Fourier Transform Raman and DFT Study of Three Annulated Oligothiophenes with Different Molecular Shapes. <i>ChemPhysChem</i> , 2007, 8, 745-750.	2.1	6
125	The first synthesis of a conjugated hybrid of C60 fullerene and a single-wall carbon nanotube. <i>Carbon</i> , 2007, 45, 2250-2252.	10.3	60
126	Vibrational spectra of oligothiophenyl-vinylenes with donor-acceptor and donor-acceptor substitution patterns. <i>Journal of Molecular Structure</i> , 2007, 834-836, 374-379.	3.6	1

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127	Vibrational spectra of nonlinear optical chromophores based on octopolar C ₃ -symmetric 1,3,5 trisalkynylbenzenes. <i>Journal of Molecular Structure</i> , 2007, 834-836, 369-373.	3.6	2
128	Electronic spectroscopy study and molecular docking simulation of the interaction of terthiophene with DNA. <i>Journal of Molecular Structure</i> , 2007, 834-836, 176-181.	3.6	4
129	Oligothiophene- and Oligopyrrole-Mediated Aggregation of Gold Nanoparticles. <i>Journal of Physical Chemistry C</i> , 2007, 111, 5886-5892.	3.1	18
130	Structure-Property Relationships in Push-Pull Amino/Cyanovinyl End-Capped Oligothiophenes: A Quantum Chemical and Experimental Studies. <i>Journal of Organic Chemistry</i> , 2006, 71, 7509-7520.	3.2	81
131	Octopolar Chromophores Based on Donor- and Acceptor-Substituted 1,3,5-Tris(phenylethynyl)benzenes: Impact of meta-Conjugation on the Molecular and Electronic Structure by Means of Spectroscopy and Theory. <i>Journal of Physical Chemistry B</i> , 2006, 110, 19198-19206.	2.6	31
132	Combined Quantum Chemical Density Functional Theory and Spectroscopic Raman and UV-vis-NIR Study of Oligothienoacenes with Five and Seven Rings. <i>Journal of Physical Chemistry A</i> , 2006, 110, 5058-5065.	2.5	39
133	Exploration of Ground and Excited Electronic States of Aromatic and Quinoid S,S-Dioxide Terthiophenes. <i>Complementary Systems for Enhanced Electronic Organic Materials</i> . <i>Journal of the American Chemical Society</i> , 2006, 128, 10134-10144.	13.7	55
134	Magnetic and Conductive Properties of Quinoidal Oligothiophenes. <i>Chemistry of Materials</i> , 2006, 18, 1539-1545.	6.7	32
135	Hybrid Organic Semiconductors Including Chalcogen Atoms in π -Conjugated Skeletons. Tuning of Optical, Redox, and Vibrational Properties by Heavy Atom Conjugation. <i>Journal of Physical Chemistry A</i> , 2006, 110, 7422-7430.	2.5	25
136	Perfluorination of tetracene: effects on the optical gap and electron-acceptor properties. An electrochemical, theoretical DFT, and Raman spectroscopic study. , 2006, , .		3
137	Regioselective hydroxylation of phenols by simultaneous photochemical generation of phenol cation-radical and hydroxyl radical. <i>Tetrahedron</i> , 2006, 62, 2927-2935.	1.9	15
138	Tuning of Electronic Properties in Thienyl-Phosphole π -Conjugated Systems through P-Functionalization Monitored by Raman Spectroscopy. <i>Chemistry - A European Journal</i> , 2006, 12, 3759-3767.	3.3	26
139	Optical, Redox, and NLO Properties of Tricyanovinyl Oligothiophenes: Comparisons between Symmetric and Asymmetric Substitution Patterns. <i>Chemistry - A European Journal</i> , 2006, 12, 5458-5470.	3.3	37
140	Magnetic Properties of Quinoidal Oligothiophenes: More Than Good Candidates for Ambipolar Organic Semiconductors?. <i>Advanced Functional Materials</i> , 2006, 16, 531-536.	14.9	42
141	Structural implications of ring shape, dimension, and metal atom insertion in nanosized cyclic oligothiophenes: Joint Raman and density functional theory study. <i>Journal of Chemical Physics</i> , 2006, 125, 044518.	3.0	12
142	Vibrational dynamics study of the effect of the substituents on the π -conjugation of different bithiophene molecules. <i>Journal of Molecular Structure</i> , 2005, 744-747, 393-401.	3.6	4
143	FT-Raman spectroscopic study, aided by quantum chemical DFT calculations, of a series of oligothiophenes end-capped by nitriles. <i>Journal of Molecular Structure</i> , 2005, 744-747, 403-409.	3.6	6
144	Combined theoretical and spectroscopic Raman study of 3,4-ethylenedioxy and S,S-dioxide substituted terthiophenes and their parent polymers. <i>Journal of Molecular Structure</i> , 2005, 744-747, 551-556.	3.6	5

#	ARTICLE	IF	CITATIONS
145	Multidisciplinary Physicochemical Analysis of Oligothiophenes End-Capped by Nitriles:  Electrochemistry, UV-Vis-Near-IR, IR, and Raman Spectroscopies and Quantum Chemistry. <i>Journal of Physical Chemistry B</i> , 2005, 109, 10115-10125.	2.6	40
146	Spectroscopic and DFT studies of donor-acceptor molecules containing phenylquinoline and phenothiazine moieties in various redox states. <i>International Journal of Quantum Chemistry</i> , 2005, 104, 635-644.	2.0	7
147	Synthesis and Characterization of a Novel Terthiophene-Based Quinodimethane Bearing a 3,4-Ethylenedioxythiophene Central Unit. <i>Journal of Physical Chemistry B</i> , 2005, 109, 22308-22318.	2.6	18
148	Tuning First Molecular Hyperpolarizabilities through the Use of Proaromatic Spacers. <i>Journal of the American Chemical Society</i> , 2005, 127, 8835-8845.	13.7	95
149	Synthesis and Characterization of Three Novel Perfluoro-oligothiophenes Ranging in Length from the Trimer to the Pentamer. <i>Journal of Physical Chemistry B</i> , 2005, 109, 20737-20745.	2.6	16
150	Alternated Quinoid/Aromatic Units in Terthiophenes Building Blocks for Electroactive Narrow Band Gap Polymers. Extended Spectroscopic, Solid State, Electrochemical, and Theoretical Study. <i>Journal of Physical Chemistry B</i> , 2005, 109, 16616-16627.	2.6	48
151	Combined Raman, electrochemical and DFT studies on a series of $\hat{1}$, $\hat{1}$ -thiophene-phosphole oligomers and their corresponding polymers. <i>Synthetic Metals</i> , 2005, 153, 249-252.	3.9	12
152	Synthesis, spectroscopy and quantum chemical DFT studies on new pleiadene-based materials. <i>Synthetic Metals</i> , 2005, 153, 245-248.	3.9	3
153	Raman and Theoretical Study of the Solvent Effects on the Sizable Intramolecular Charge Transfer in the Push-Pull 5-(Dimethylamino)-5-nitro-2,2-bithiophene. <i>Journal of Physical Chemistry A</i> , 2005, 109, 8724-8731.	2.5	28
154	Incisive Structure-Spectroscopic Correlation in Oligothiophenes Functionalized with (\hat{A}) Inductive/Mesomeric Fluorine Groups: A Joint Raman and DFT Study. <i>Journal of the American Chemical Society</i> , 2005, 127, 13364-13372.	13.7	29
155	Exploration of the electronic structure of dendrimerlike acetylene-bridged oligothiophenes by correlating Raman spectroscopy, electrochemistry, and theory. <i>Journal of Chemical Physics</i> , 2004, 120, 11874-11881.	3.0	10
156	Microwave-assisted sidewall functionalization of single-wall carbon nanotubes by Diels-Alder cycloaddition. <i>Chemical Communications</i> , 2004, , 1734-1735.	4.1	149
157	Spectroscopic and Theoretical Study of the Molecular and Electronic Structures of a Terthiophene-Based Quinodimethane. <i>ChemPhysChem</i> , 2004, 5, 529-539.	2.1	46
158	Application of Raman spectroscopy and quantum chemistry for featuring the structure of positively charged species in macrocyclic $\hat{1}$ -conjugated diacetylene-bridged oligothiophenes. <i>Journal of Raman Spectroscopy</i> , 2004, 35, 592-599.	2.5	25
159	Vibrational and Quantum-Chemical Study of Nonlinear Optical Chromophores Containing Dithienothiophene as the Electron Relay. <i>Chemistry - A European Journal</i> , 2004, 10, 3805-3816.	3.3	44
160	Quantum chemical DFT and spectroscopic study of a push-pull chromophore for second-order nonlinear optics containing bithiophene as the electron relay. <i>Computational and Theoretical Chemistry</i> , 2004, 709, 187-193.	1.5	27
161	A Practical Spectroscopic and Theoretical Approach To Study the Electrochromism in Molecular-Based Materials: The Case of a Family of Dendrimerlike Poly(6-azulenylethenyl)benzenes. <i>Journal of Physical Chemistry B</i> , 2004, 108, 18463-18471.	2.6	6
162	A Raman and Computational Study of Two Dithienyl Naphthodithiophenes: A Synthesis and Characterization of New Polymers Showing Low Band Gap Optical and Electroactive Features. <i>Journal of Physical Chemistry B</i> , 2004, 108, 7611-7619.	2.6	4

#	ARTICLE	IF	CITATIONS
163	Combined Raman and Computational Study of a Novel Series of Macrocyclic π -Conjugated Diacetylene-Bridged π -Linked Oligothiophenes. <i>Journal of Physical Chemistry B</i> , 2004, 108, 3158-3167.	2.6	24
164	Electronic Modulation of Dithienothiophene (DTT) as π -Center of D- π -D Chromophores on Optical and Redox Properties: A Analysis by UV-Vis-NIR and Raman Spectroscopies Combined with Electrochemistry and Quantum Chemical DFT Calculations. <i>Journal of the American Chemical Society</i> , 2004, 126, 13363-13376.	13.7	52
165	Study of the ac conductivity of π - π -dimethyl sexithiophene in pristine and doped states. <i>Journal of Non-Crystalline Solids</i> , 2004, 342, 146-151.	3.1	2
166	Combined Spectroscopic and Theoretical Study of Narrow Band Gap Heterocyclic Co-oligomers Containing Alternating Aromatic Donor and Quinoid Acceptor Units. <i>Journal of Physical Chemistry B</i> , 2004, 108, 2516-2526.	2.6	66
167	Vibrational and Quantum-Chemical Study of Push-Pull Chromophores for Second-Order Nonlinear Optics from Rigidified Thiophene-Based π -Conjugating Spacers. <i>Chemistry - A European Journal</i> , 2003, 9, 3670-3682.	3.3	57
168	Vibrational study of push-pull chromophores for second-order non-linear optics derived from rigidified thiophene π -conjugating spacers. <i>Journal of Molecular Structure</i> , 2003, 651-653, 151-158.	3.6	34
169	Theoretical description of the Raman spectrum of a vinylene-bridged quaterthiophene oligomer. <i>Journal of Molecular Structure</i> , 2003, 651-653, 657-664.	3.6	9
170	UV-Vis, IR, Raman and theoretical characterization of a novel quinoid oligothiophene molecular material. <i>Journal of Molecular Structure</i> , 2003, 651-653, 665-673.	3.6	10
171	Computation and Spectroelectrochemistry as Complementary Tools for the Study of Electrochemically Induced Charged Defects in 4-[Bis(4-methylphenyl)amino]phenyl Oligothiophenes as Model Systems for Hole-Transporting Materials. <i>Journal of Physical Chemistry B</i> , 2003, 107, 2637-2644.	2.6	42
172	Spectroscopic and Theoretical Study of Push-Pull Chromophores Containing Thiophene-Based Quinonoid Structures as Electron Spacers. <i>Journal of Physical Chemistry B</i> , 2003, 107, 12175-12183.	2.6	40
173	Nitro-Functionalized Oligothiophenes as a Novel Type of Electroactive Molecular Material: A Spectroscopic, Electrochemical, and Computational Study. <i>Journal of the American Chemical Society</i> , 2003, 125, 2524-2534.	13.7	106
174	Infrared and Raman features of a series of π -bis(arylthio)oligothiophenes as molecular wires. A π -electron delocalization efficiency study. <i>Journal of Chemical Physics</i> , 2003, 118, 1912-1920.	3.0	21
175	Efficiency of the π conjugation in a novel family of π -bisphenyl end-capped oligothiophenes by means of Raman spectroscopy. <i>Journal of Chemical Physics</i> , 2002, 116, 10419-10427.	3.0	63
176	Experimental and Theoretical Study of the Infrared and Raman Spectra of a Substituted Sexithiophene in Five Oxidation States. <i>Journal of Physical Chemistry B</i> , 2002, 106, 3597-3605.	2.6	63
177	Vibrational Spectroscopic Features of a Novel Family of Amorphous Molecular Materials Containing an Oligothiophene Moiety as Color-Tunable Emitting Materials. <i>Journal of Physical Chemistry B</i> , 2002, 106, 7163-7170.	2.6	41
178	Combined Spectroelectrochemical and Theoretical Study of a Vinylene-Bridged Sexithiophene Cooligomer: A Analysis of the π -Electron Delocalization and of the Electronic Defects Generated upon Doping. <i>Journal of Physical Chemistry B</i> , 2002, 106, 3872-3881.	2.6	63
179	Quinonoid Oligothiophenes as Electron-Donor and Electron-Acceptor Materials. A Spectroelectrochemical and Theoretical Study. <i>Journal of the American Chemical Society</i> , 2002, 124, 12380-12388.	13.7	109
180	Infrared spectra of two sexithiophenes in neutral and doped states: a theoretical and spectroscopic study. <i>Vibrational Spectroscopy</i> , 2002, 30, 175-189.	2.2	7

#	ARTICLE	IF	CITATIONS
181	Tetrathiafulvalene Derivatives as NLO-phores:Â Synthesis, Electrochemistry, Raman Spectroscopy, Theoretical Calculations, and NLO Properties of Novel TTF-Derived Donor-Î€-Acceptor Dyads. Journal of Organic Chemistry, 2001, 66, 8872-8882.	3.2	127
182	Intramolecular charge transfer in push-pull oligothiophenes from their vibrational spectra. Synthetic Metals, 2001, 119, 551-552.	3.9	4
183	Theoretical and vibrational study of electron-acceptor oligothienoquinonoids with well defined substitution patterns. Synthetic Metals, 2001, 119, 553-554.	3.9	2
184	Spectroelectrochemical Raman study of a new series of thiophene/phenylene co-oligomers. Synthetic Metals, 2001, 119, 305-306.	3.9	8
185	Vibrational and theoretical DFT study of two regioregular methyl-disubstituted bithiophenes. Journal of Molecular Structure, 2001, 563-564, 539-544.	3.6	3
186	Spectroelectrochemical Raman Study of two Î±,Î±'-End Capped Sexithiophenes: The Effect of the Introduction of a Polarisable Sulfur Atom in the Side Chain. Materials Research Society Symposia Proceedings, 2000, 660, .	0.1	0
187	Infrared and Raman spectra of a new radical cation charged defect created on a well-barrier-well thiophene-based oligomer. Journal of Raman Spectroscopy, 2000, 31, 565-570.	2.5	23
188	Vibrational spectra of charged defects in a series of Î±,Î±'-bis(aminomethyl) end-capped oligothiophenes induced by chemical doping with iodine. Journal of Molecular Structure, 2000, 521, 239-247.	3.6	3
189	Density functional study on the structures and vibrational spectra of the radical cation and dication of Î±,Î±'-bis(aminomethyl)quaterthiophene. Journal of Molecular Structure, 2000, 521, 249-260.	3.6	6
190	Spectroelectrochemical Raman Study of two Î±,Î±'-End Capped Sexithiophenes: The Effect of the Introduction of a Polarisable Sulfur Atom in the Side Chain. Materials Research Society Symposia Proceedings, 2000, 660, 1.	0.1	0
191	Raman spectra and conformational properties of hexyl- and methylsulphanyl-substituted oligothiophenes. Synthetic Metals, 2000, 108, 27-31.	3.9	9
192	A Combined Spectroscopic and Theoretical Study of a Series of Aminomethyl End-Capped Oligothiophenes with Potential Applications in Thin Film Devices. Journal of Physical Chemistry A, 2000, 104, 735-740.	2.5	24
193	Spectroelectrochemical Raman Study of a Novel Well-Barrier-Well Vinylene-Bridged-Octithiophene Oligomer:Â An Analysis of the Conjugation Length and of the Electronic Defects Created upon Doping. Journal of Physical Chemistry A, 2000, 104, 10656-10661.	2.5	14
194	Combined Theoretical and Vibrational Study of Dihexylbithienoquinonoid Derivatives with Regioregular Head-to-Head, Head-to-Tail, and Tail-to-Tail Orientations. Journal of Physical Chemistry A, 2000, 104, 661-672.	2.5	30
195	FT-IR and FT-Raman spectra of a series of oxidized Î±,Î±'-diethyl end-capped oligothiophenyls: a spectroscopic study of conjugational model defects. Optical Materials, 1999, 12, 321-325.	3.6	3
196	Infrared and Raman Spectra of a Well-Barrier-Well 1,2-Di(Î±,Î±'-Bithienyl)Vinylene. Synthetic Metals, 1999, 101, 548.	3.9	1
197	Vibrational and electronic spectroscopic study of two oligothiophene materials bearing a heteroquinonoid structure.. Synthetic Metals, 1999, 101, 549-550.	3.9	1
198	Vibrational spectra and ab initio dft calculations of 3,3'-and 4,4'-dimethyl substituted 2,2'-bithiophene.. Synthetic Metals, 1999, 101, 590-591.	3.9	4

#	ARTICLE	IF	CITATIONS
199	Vibrational Spectroscopic Study of a Series of $\hat{I}_{\pm}, \hat{I}_{\pm}^{\sim}$ -Diethyl End-Capped Oligothiophenes with Different Chain Lengths in the Neutral State. <i>Journal of Physical Chemistry A</i> , 1999, 103, 816-822.	2.5	34
200	FT-Raman Studies of Charged Defects Created on Methyl End-Capped Oligothiophenes by Doping with NOBF ₄ . <i>Advanced Materials</i> , 1998, 10, 1458-1461.	21.0	68
201	Electrochemical doping in a series of $\hat{I}_{\pm}, \hat{I}_{\pm}^{\sim}$ -dimethyl end-capped oligothiophenyls An FT-Raman confirmation of a radical cation generation. <i>Optical Materials</i> , 1998, 9, 82-87.	3.6	14
202	$\hat{I}_{\pm}, \hat{I}_{\pm}^{\sim}$ in situ TM spectroelectrochemical study of a series of $\hat{I}_{\pm}, \hat{I}_{\pm}^{\sim}$ -dimethyl end-capped oligothiophene films. <i>Synthetic Metals</i> , 1998, 95, 93-100.	3.9	18
203	Vibrational spectroscopic study of 5,5- \hat{I}_{\pm}^{\sim} -bis(dicyanomethylene)-5,5- \hat{I}_{\pm}^{\sim} -dihydro- \hat{I}_{\pm}^{\sim} -2,2- \hat{I}_{\pm}^{\sim} ;5- \hat{I}_{\pm}^{\sim} ,2- \hat{I}_{\pm}^{\sim} -terthiophene bearing a heteroquinonoid structure as a model of doped polythiophene. <i>Journal of Chemical Physics</i> , 1998, 109, 2543-2548.	3.0	26
204	Vibrational spectra of charged defects in a series of $\hat{I}_{\pm}, \hat{I}_{\pm}^{\sim}$ -dimethyl end-capped oligothiophenes induced by chemical doping with iodine. <i>Journal of Chemical Physics</i> , 1998, 109, 10419-10429.	3.0	107
205	Vibrational Spectroscopy study of doping induced charged defects in a series of $\hat{I}_{\pm}, \hat{I}_{\pm}^{\sim}$ -Dimethyl end-capped Oligothiophenes. <i>Synthetic Metals</i> , 1997, 84, 571-572.	3.9	8
206	A theoretical investigation of $\hat{I}_{\pm}, \hat{I}_{\pm}^{\sim}$ -dimethyl end-capped oligothiophenes: Structures, vibrational spectra and conjugational defects. <i>Synthetic Metals</i> , 1997, 85, 1157-1158.	3.9	2
207	Vibrational analysis as a function of the chain-length of well-barrier-well $\hat{I}_{\pm}, \hat{I}_{\pm}^{\sim}$ -conjugated thiophene-based oligomers. <i>Synthetic Metals</i> , 1997, 84, 591-592.	3.9	4
208	A theoretical investigation of $\hat{I}_{\pm}, \hat{I}_{\pm}^{\sim}$ -dimethyl end-capped oligothiophenes: structures, vibrational spectra and conjugation defects. <i>Synthetic Metals</i> , 1997, 89, 159-160.	3.9	2
209	Ab initio self-consistent reaction field calculations on amino acids: asparagine zwitterions in polar medium and gas phase. <i>Theoretical Chemistry Accounts</i> , 1997, 98, 5-15.	1.4	9
210	Ab initio theoretical study of thiophene derivatives: 2-methylthiophene and 3-methylthiophene. <i>Journal of Molecular Structure</i> , 1997, 410-411, 311-314.	3.6	7
211	Solvent effects on the structure and spectra of glutamine studied by the SCRF theory. <i>Journal of Molecular Structure</i> , 1997, 410-411, 353-356.	3.6	0
212	Infrared and Raman spectra of two well-barrier-well 1,2-di($\hat{I}_{\pm}, \hat{I}_{\pm}^{\sim}$ -oligothienyl)ethanes. <i>Journal of Raman Spectroscopy</i> , 1997, 28, 855-865.	2.5	3
213	Solvent effects on electronic properties, geometries and internal rotation barriers of bithiophenes. An ab initio self-consistent reaction field theoretical study. <i>Synthetic Metals</i> , 1996, 76, 221-224.	3.9	16
214	Delocalization length, electronic properties and vibrational spectra of neutral $\hat{I}_{\pm}, \hat{I}_{\pm}^{\sim}$ -dimethyl end-capped oligothiophenes. <i>Synthetic Metals</i> , 1996, 76, 277-280.	3.9	15
215	$\hat{I}_{\pm}, \hat{I}_{\pm}^{\sim}$ Electron delocalization in pristine polyfuran: from the oligomers to the polymer. <i>Acta Polymerica</i> , 1996, 47, 62-65.	0.9	18
216	Scaled Quantum-Mechanical Force Field and Vibrational Spectra of 3-Methylthiophene. <i>The Journal of Physical Chemistry</i> , 1996, 100, 2907-2914.	2.9	14

#	ARTICLE	IF	CITATIONS
217	Conformational Disorder and Mean Conjugation of Neutral β,β -Dimethyl End-Capped Oligothiophenes in Solution: A FT-Raman and FT-Infrared Spectroscopic Study. <i>The Journal of Physical Chemistry</i> , 1996, 100, 289-293.	2.9	27
218	Infrared and Raman spectra of L-asparagine and L-asparagine-d5 in the solid state. <i>Journal of Raman Spectroscopy</i> , 1995, 26, 1003-1008.	2.5	31
219	Vibrational study of aspartic acid and glutamic acid dipeptides. <i>Journal of Molecular Structure</i> , 1995, 348, 249-252.	3.6	57
220	Vibrational spectra and assignments of amino acid L-asparagine. <i>Journal of Molecular Structure</i> , 1995, 349, 57-60.	3.6	9
221	Force field and normal coordinate calculations for glutamic acid. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 1995, 51, 293-302.	3.9	35
222	Force field and normal coordinate calculations of the amino acid l-asparagine. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 1995, 51, 2347-2356.	3.9	8
223	A semiempirical approach for the calculation of the vibrational spectra of conducting polymers: the case of polyselenophene. <i>Journal of Molecular Structure</i> , 1995, 348, 91-94.	3.6	10
224	Lattice dynamics and vibrational spectra of polyfuran: effective conjugation coordinate and photoexcited spectrum. <i>Synthetic Metals</i> , 1995, 69, 391-392.	3.9	10
225	Structural and spectroscopical study of glutamic acid in the nonzwitterionic form. <i>Computational and Theoretical Chemistry</i> , 1995, 330, 261-266.	1.5	28
226	Ab initio study of torsional potentials in 2,2'-bithiophene and 3,4'- and 3,3'-dimethyl-2,2'-bithiophene as models of the backbone flexibility in polythiophene and poly(3-methylthiophene). <i>Journal of Chemical Physics</i> , 1994, 101, 1369-1377.	3.0	78
227	Vibrational spectra of $[^{15}\text{N}]$ glutamic acid and $[2\text{H}_4]$ glutamic acid. <i>Journal of Raman Spectroscopy</i> , 1994, 25, 861-867.	2.5	12
228	Transferable semiempirical quadratic force fields: The case of polythiophene and shorter oligomers. <i>Journal of Computational Chemistry</i> , 1994, 15, 405-423.	3.3	26
229	Ir and Raman spectra of L-aspartic acid and isotopic derivatives. <i>Biopolymers</i> , 1994, 34, 1065-1077.	2.4	67
230	An interpretation of the vibrational spectra of insulating and electrically conducting poly(3-methylthiophene) aided by a theoretical dynamical model. <i>Journal of Chemical Physics</i> , 1994, 100, 114-129.	3.0	66
231	Conformational and vibrational study on 1,2-dihydroxybenzene. <i>Journal of Molecular Structure</i> , 1993, 293, 59-62.	3.6	5
232	Comparison between semiempirical and experimental force fields of oligothiophenes as an approach for the calculations of the vibrational spectrum of the polymer. <i>Journal of Molecular Structure</i> , 1993, 294, 37-40.	3.6	3
233	Harmonic force field for amino acid L-glutamine by MNDO semiempirical method. <i>Journal of Molecular Structure</i> , 1993, 294, 49-52.	3.6	8
234	Normal coordinate and rotational barrier calculations on 1,2-dihydroxybenzene. <i>Vibrational Spectroscopy</i> , 1993, 4, 321-334.	2.2	35

#	ARTICLE	IF	CITATIONS
235	A study by Raman spectroscopy and the semiempirical AM1 method on several 1,2-dihydroxybenzene solutions. <i>Spectrochimica Acta Part A: Molecular Spectroscopy</i> , 1993, 49, 1759-1767.	0.1	12
236	Resonance raman spectra and lattice dynamics calculations of pristine and doped polyconjugated polyfuran. <i>Synthetic Metals</i> , 1993, 57, 4467-4472.	3.9	12
237	Vibrational Spectra and lattice dynamics calculations of poly (p-phenylene): oligomers and polymer. <i>Synthetic Metals</i> , 1993, 57, 4461-4466.	3.9	10
238	Lattice dynamics and vibrational spectra of pristine and doped polyconjugated polyfuran. <i>Journal of Chemical Physics</i> , 1993, 98, 769-783.	3.0	50
239	Lattice dynamics and vibrational spectra of pristine, doped, and photoexcited poly(3-methylthiophene). <i>Synthetic Metals</i> , 1992, 51, 211-218.	3.9	9

240

#	ARTICLE	IF	CITATIONS
253	A simple interpretation of the vibrational spectra of undoped, doped and photoexcited polyacetylene: Amplitude mode theory in the GF formalism. Solid State Communications, 1988, 65, 625-630.	1.9	221
254	FTIR spectra (frequency and intensity) of poly-(para-phenylenes). Mikrochimica Acta, 1988, 94, 247-249.	5.0	8
255	Lattice dynamics and infrared spectra of doping induced and photoexcited poly (acetylene). Solid State Communications, 1988, 65, 409-414.	1.9	3
256	Interpretation of vibrational spectra of pristine, doped and photoinduced polyacetylene.. Journal of Molecular Structure, 1988, 174, 375-382.	3.6	1
257	Vibrational spectrum and internal rotation in 2-methylpyrazine. Journal of the Chemical Society, Faraday Transactions 2, 1988, 84, 53-65.	1.1	43
258	On the problem of π electron delocalisation across sp^3 carbon atoms introduced as defects in polyacetylene. Solid State Communications, 1987, 64, 1183-1186.	1.9	14
259	Vibrational spectrum and internal rotation in 2,5-dimethylpyrazine. Journal of Molecular Structure, 1987, 162, 263-272.	3.6	15
260	Force field for in-plane vibrations of pyrazine. Spectrochimica Acta Part A: Molecular Spectroscopy, 1986, 42, 1343-1348.	0.1	27
261	A MINDO/3 harmonic force field for pyrazine. In-plane A_g and B_{3u} vibrations. Journal of Molecular Structure, 1986, 142, 295-298.	3.6	7
262	Conformations and vibrational spectra of methyl-pyrazines. Journal of Molecular Structure, 1986, 142, 423-426.	3.6	5
263	Vibrational spectra of [1H4]pyrazine and [2H4]pyrazine. Journal of the Chemical Society, Faraday Transactions 2, 1985, 81, 405.	1.1	51