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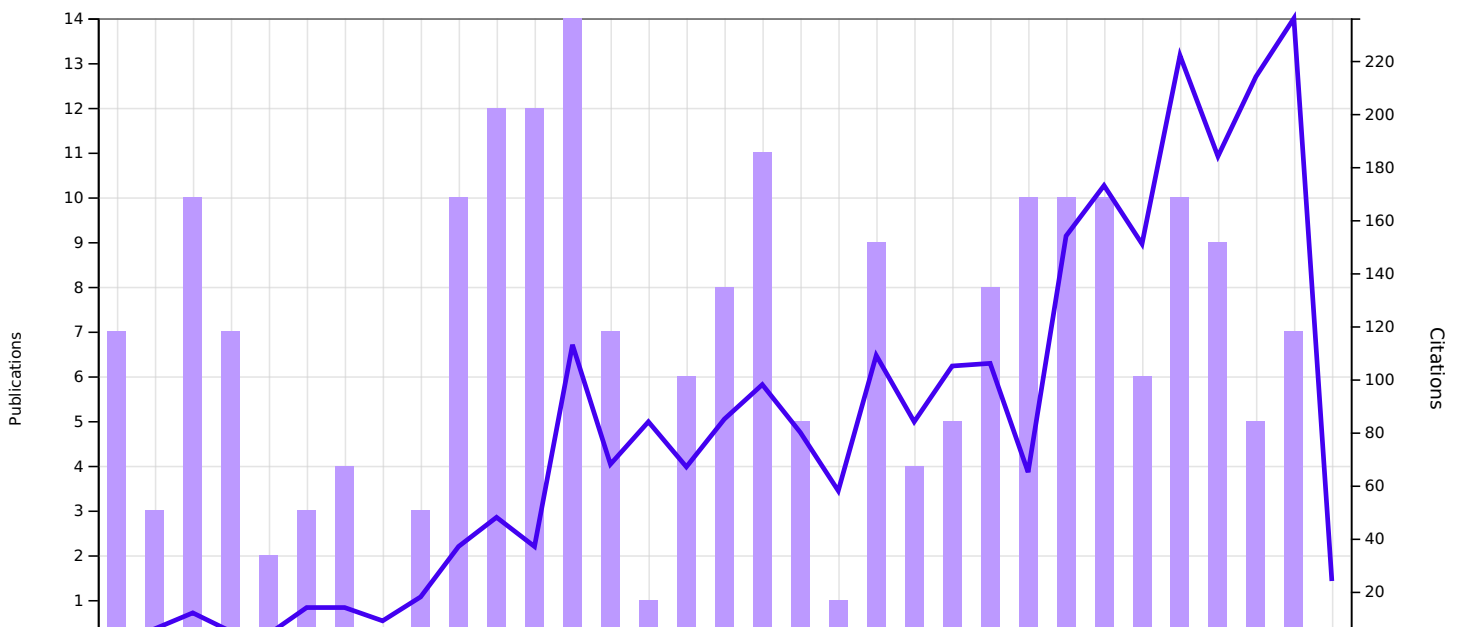
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1	Computing relative stabilities of metallofullerenes by Gibbs energy treatments Slanina, Z; Lee, SL; (...); Nagase, S 10th Electronic Computational Chemistry Conference Feb 2007 THEORETICAL CHEMISTRY ACCOUNTS 117 (2) , pp.315-322	11	6	3	4	0	5.59	95					
2	AM1 COMPUTATIONS ON 7 ISOLATED-PENTAGON-RULE ISOMERS OF C-80 SUN, ML; SLANINA, Z; (...); ADAMOWICZ, L Nov 17 1995 CHEMICAL PHYSICS LETTERS 246 (1-2) , pp.66-72	3	0	0	1	0	3.03	88					
3	Enhanced Relaxation and Intermixing in Ge Islands Grown on Pit-Patterned Si(001) Substrates Schulli, TU; Vastola, G; (...); Bauer, G Jan 16 2009 PHYSICAL REVIEW LETTERS 102 (2)	2	0	0	0	0	5.47	82					
4	Enthalpy-entropy interplay for C-36 cages: B3LYP/6-31G(*) calculations Slanina, Z; Uhlik, F; (...); Osawa, E Sep 22 2000 JOURNAL OF CHEMICAL PHYSICS 113 (12) , pp.4933-4937	2	1	0	0	0	3.29	79					
5	Multicentre bond indices from the generalized population analysis of higher order densities Ponec, R and Uhlik, F 10th Dubrovnik International Course / Conference MATH/CHEM/COMP Nov 1996 CROATICA CHEMICA ACTA 69 (3) , pp.941-954	1	0	1	0	0	2.29	64					
6	La-2@C-72 and Sc-2@C-72: Computational characterizations Slanina, Z; Chen, ZF; (...); Nagase, S Feb 16 2006 JOURNAL OF PHYSICAL CHEMISTRY A 110 (6) , pp.2231-2234	2	1	1	1	0	3.17	57					
7	Computational modeling of the elemental catalysis in the Stone-Wales fullerene rearrangements Slanina, Z; Zhao, X; (...); Osawa, E Apr 9 2000 JOURNAL OF ORGANOMETALLIC CHEMISTRY 599 (1) , pp.57-61	3	2	2	3	0	2.25	54					
8	Modeling of Ionization and Conformations of Starlike Weak Polyelectrolytes Uhlik, F; Kosovan, P; (...); Leermakers, FAM Jun 24 2014 MACROMOLECULES 47 (12) , pp.4004-4016	9	5	9	6	0	5.2	52					
9	Simulations of ionization equilibria in weak polyelectrolyte solutions and gels												

9	Landsgesell, J; Nova, L; (...); Kosovan, P Feb 14 2019 SOFT MATTER 15 (6) , pp.1155-1185	6	13	17	11	0	9.4	47
10	Local pH and effective pK(A) of weak polyelectrolytes - insights from computer simulations Nova, L; Uhlík, F and Kosovan, P Jun 14 2017 PHYSICAL CHEMISTRY CHEMICAL PHYSICS 19 (22) , pp.14376-14387	10	9	9	9	3	6.71	47
11	Electron pairing and chemical bonds. On the accuracy of the electron pair model of chemical bond Ponec, R and Uhlík, F Feb 28 1997 JOURNAL OF MOLECULAR STRUCTURE-THEOCHEM 391 (1-2) , pp.159-168	1	0	1	0	0	1.74	47
12	Polyelectrolyte behavior of polystyrene-block-poly(methacrylic acid) micelles in aqueous solutions at low ionic strength Matejíček, P; Podhájek, K; (...); Spírková, M Dec 28 2004 MACROMOLECULES 37 (26) , pp.10141-10154	1	2	2	2	0	2.3	46
13	Intrinsic Viscosity of Cyclic Polystyrene Jeong, Y; Jin, Y; (...); Roovers, J Oct 10 2017 MACROMOLECULES 50 (19) , pp.7770-7776	6	5	11	13	4	6.43	45
14	Computed structure and energetics of La@C-60 Slanina, Z; Lee, S-L; (...); Nagase, S 8th Annual Meeting of the V-A-Fock-School on Quantum and Computational Chemistry Aug 5 2005 INTERNATIONAL JOURNAL OF QUANTUM CHEMISTRY 104 (2) , pp.272-277	4	3	3	3	0	2.37	45
15	C5H3+ ISOMERIC STRUCTURES - RELATIVE STABILITIES AT HIGH-TEMPERATURES SLANINA, Z; UHLIK, F and ZERNER, MC Aug 1991 REVUE ROUMAINE DE CHIMIE 36 (8) , pp.965-974	4	3	1	2	0	1.3	43
16	Facile Synthesis of an Extensive Family of Sc2O@C-2n(n=35-47) and Chemical Insight into the Smallest Member of Sc2O@C-2(7892)-C-70 Zhang, MB; Hao, YJ; (...); Cong, HL Dec 11 2014 JOURNAL OF PHYSICAL CHEMISTRY C 118 (49) , pp.28883-28889	5	2	2	5	1	3.9	39
17	Hiding and Recovering Electrons in a Dimetallic Endohedral Fullerene: Air-Stable Products from Radical Additions Yamada, M; Kurihara, H; (...); Akasaka, T Jan 14 2015 JOURNAL OF THE AMERICAN CHEMICAL SOCIETY 137 (1) , pp.232-238	6	8	6	2	0	4.11	37
18	Experimental study of hydrophobically modified amphiphilic block copolymer micelles using light scattering and nonradiative excitation energy transfer Matejíček, P; Uhlík, F; (...); Webber, SE Dec 3 2002 MACROMOLECULES 35 (25) , pp.9487-9496	0	1	1	1	0	1.68	37
19	MPWB1K calculations of stepwise encapsulations: Li-x@C-60 Slanina, Z; Uhlík, F; (...); Nagase, S Sep 22 2008 CHEMICAL PHYSICS LETTERS 463 (1-3) , pp.121-123	1	1	0	0	0	2.13	34

20	Zeolite (In)Stability under Aqueous or Steaming Conditions Heard, C.J.; Grajciar, L.; (...); Nachtigall, P. Nov 5 2020 Aug 2020 (Early Access) ADVANCED MATERIALS 32 (44)	0	0	15	12	5	8	32
21	Computed structures of two known Yb@ C-74 isomers Slanina, Z.; Uhlik, F. and Nagase, S. Nov 30 2006 JOURNAL OF PHYSICAL CHEMISTRY A 110 (47) , pp.12860-12863	1	1	0	0	0	1.78	32
22	Mg@C(72) MNDO/d evaluation of the isomeric composition Slanina, Z.; Zhao, X.; (...); Nagase, S. 2001 JOURNAL OF MOLECULAR GRAPHICS & MODELLING 19 (2) , pp.252-255	1	0	0	1	0	1.39	32
23	Computing the relative gas-phase populations of C-60 and C-70: Beyond the traditional Delta H-f,298(o) scale Slanina, Z.; Zhao, X.N.; (...); Adamowicz, L. 2001 JOURNAL OF MOLECULAR GRAPHICS & MODELLING 19 (2) , pp.216-221	0	0	0	0	0	1.35	31
24	TEMPERATURE-DEPENDENCE OF THE GIBBS ENERGY ORDERING OF ISOMERS OF CL2O2 SLANINA, Z. and UHLIK, F. Jul 11 1991 JOURNAL OF PHYSICAL CHEMISTRY 95 (14) , pp.5432-5434	3	2	1	3	0	0.94	31
25	Isolation and Crystallographic Characterization of the Labile Isomer of Y@C-82 Cocrystallized with Ni(OEP): Unprecedented Dimerization of Pristine Metallofullerenes Bao, L.P.; Pan, C.W.; (...); Lu, X. Aug 1 2016 ANGEWANDTE CHEMIE-INTERNATIONAL EDITION 55 (32) , pp.9234-9238	11	5	4	0	0	3.75	30
26	Intermixing in heteroepitaxial islands: fast, self-consistent calculation of the concentration profile minimizing the elastic energy Gatti, R.; Uhlik, F. and Montalenti, F. Aug 28 2008 NEW JOURNAL OF PHYSICS 10	0	1	1	0	0	1.88	30
27	Polyelectrolyte shells of copolymer micelles in aqueous solutions: A Monte Carlo study Uhlik, F.; Limpouchova, Z.; (...); Prochazka, K. Aug 1 2004 JOURNAL OF CHEMICAL PHYSICS 121 (5) , pp.2367-2375	0	2	2	1	0	1.5	30
28	Ab initio SCF nonlinear pair population analysis. A new means of detection and localization of multicenter bonding Bochicchio, R.C.; Ponec, R. and Uhlik, F. Nov 5 1997 INORGANIC CHEMISTRY 36 (23) , pp.5363-5368	0	0	0	0	0	1.11	30
29	Sm@C-2V,(19138)-C-76: A Non-IPR Cage Stabilized by a Divalent Metal Ion Hao, Y.J.; Feng, L.; (...); Uhlik, F. May 4 2015 INORGANIC CHEMISTRY 54 (9) , pp.4243-4248	3	3	4	3	1	3	27
	Computed Stabilities in Metallofullerene Series: Al@C-82, Sc@C-82, Y@C-82, and La@C-82							

30	Slanina, Z; Uhlík, F; (...); Nagase, S 12th V A Fock Meeting on Quantum and Computational Chemistry Sep 2011 INTERNATIONAL JOURNAL OF QUANTUM CHEMISTRY 111 (11) , pp.2712-2718	5	2	1	3	0	2	26
31	Synthesis and photophysical properties of alpha,omega-bis(terpyridine)oligothiophenes Svoboda, J; Stenclova, P; (...); Vohlídal, J Jan 7 2011 TETRAHEDRON 67 (1) , pp.75-79	0	0	2	2	0	2	26
32	Computing enthalpy-entropy interplay for isomeric fullerenes Slanina, Z; Zhao, X; (...); Adamowicz, L 4th Congress of the International-Society-for-Theoretical-Chemical-Physics (ICTCP 4) Sep 15 2004 INTERNATIONAL JOURNAL OF QUANTUM CHEMISTRY 99 (5) , pp.640-653	1	5	2	0	0	1.3	26
33	Evidence for fullerenes in solid bitumen from pillow lavas of proterozoic age from Mitov (Bohemian Massif, Czech Republic) Jehlička, J; Svatos, A; (...); Uhlík, F Apr 2003 GEOCHIMICA ET COSMOCHIMICA ACTA 67 (8) , pp.1495-1506	2	0	0	1	0	1.24	26
34	Isomeric Sc2O@C-78 Related by a Single-Step Stone-Wales Transformation: Key Links in an Unprecedented Fullerene Formation Pathway Hao, YJ; Tang, QQ; (...); Uhlík, F Nov 7 2016 INORGANIC CHEMISTRY 55 (21) , pp.11354-11361	5	8	1	5	0	3	24
35	C-78 IPR fullerenes: Computed B3LYP/6-31G*//HF/3-21G temperature-dependent relative concentrations Uhlík, F; Slanina, Z and Osawa, E 10th International Symposium on Small Particles and Inorganic Clusters (ISSPIC 10) Sep 2001 EUROPEAN PHYSICAL JOURNAL D 16 (1-3) , pp.349-352	0	2	0	0	0	1.04	24
36	Charge-controlled nano-structuring in partially collapsed star-shaped macromolecules Uhlík, F; Kosovan, P; (...); Borisov, OV 2016 SOFT MATTER 12 (21) , pp.4846-4852	6	3	4	1	0	2.88	23
37	Silver(I) complexes with 1⁻-(diphenylphosphino)-1-cyanoferrocene: the art of improvisation in coordination Skoch, K; Uhlík, F; (...); Stepnicka, P 2016 DALTON TRANSACTIONS 45 (26) , pp.10655-10671	5	1	4	3	2	2.75	22
38	Tuning intramolecular electron and energy transfer processes in novel conjugates of La-2@C-80 and electron accepting subphthalocyanines Feng, L; Rudolf, M; (...); Akasaka, T 2015 CHEMICAL COMMUNICATIONS 51 (2) , pp.330-333	5	3	2	1	0	2.44	22
39	Enhancing Built-In Electric Field and Defect Passivation through Gradient Doping in Inverted CsPbI2Br Perovskite Solar Cells Han, DW; Yuan, Q; (...); Feng, L Jan 2021 Dec 2020 (Early Access) SOLAR RRL 5 (1)	0	0	6	13	2	5.25	21



40	La-2@C-s(17490)-C-76: A New Non-IPR Dimetallic Metallofullerene Featuring Unexpectedly Weak Metal-Pentalene Interactions Suzuki, M; Mizorogi, N; (...); Akasaka, T Dec 9 2013 CHEMISTRY-A EUROPEAN JOURNAL 19 (50) , pp.17125-17130	1	2	1	1	0	1.91	21
41	Isolation and Structural Characterization of Er@C-2v(9)-C-82 and Er@C-s(6)-C-82: Regioselective Dimerization of a Pristine Endohedral Metallofullerene Induced by Cage Symmetry Hu, SF; Liu, T; (...); Lu, X Feb 4 2019 INORGANIC CHEMISTRY 58 (3) , pp.2177-2182	4	9	5	2	0	4	20
42	The Unanticipated Dimerization of Ce@C-2v(9)-C-82 upon Co-crystallization with Ni(octaethylporphyrin) and Comparison with Monomeric M@C-2v(9)-C-82 (M = La, Sc, and Y) Suzuki, M; Yamada, M; (...); Akasaka, T Dec 12 2016 CHEMISTRY-A EUROPEAN JOURNAL 22 (50) , pp.18115-18122	8	3	3	1	0	2.38	19
43	Popular C-82 Fullerene Cage Encapsulating a Divalent Metal Ion Sm²⁺: Structure and Electrochemistry Hu, ZQ; Hao, YJ; (...); Feng, L Mar 2 2015 INORGANIC CHEMISTRY 54 (5) , pp.2103-2108	7	0	2	1	1	2.11	19
44	A Monte Carlo study of shells of hydrophobically modified amphiphilic copolymer micelles in polar solvents Uhlík, F; Limpouchova, Z; (...); Prochazka, K Jun 22 2003 JOURNAL OF CHEMICAL PHYSICS 118 (24) , pp.11258-11264	0	1	1	1	0	0.9	19
45	Nonradiative excitation energy transfer in hydrophobically modified amphiphilic block copolymer micelles. theoretical model and Monte Carlo simulations Uhlík, F; Limpouchova, Z; (...); Webber, SE Dec 3 2002 MACROMOLECULES 35 (25) , pp.9497-9505	0	1	0	0	0	0.86	19
46	Computed structures and relative stabilities of Be@C-74 Slanina, Z; Uhlík, F; (...); Nagase, S 10th Annual Meeting of the V-A-Fock-School on Quantum and Computational Chemistry Nov 5 2007 INTERNATIONAL JOURNAL OF QUANTUM CHEMISTRY 107 (13) , pp.2494-2498	0	0	0	0	0	1.06	18
47	Eu@C-72: Computed Comparable Populations of Two Non-IPR Isomers Slanina, Z; Uhlík, F; (...); Lu, X Jul 2017 MOLECULES 22 (7)	5	3	2	6	0	2.43	17
48	Calculations of Metallofullerene Yields Slanina, Z; Uhlík, F; (...); Nagase, S Nov 2011 JOURNAL OF COMPUTATIONAL AND THEORETICAL NANOSCIENCE 8 (11) , pp.2233-2239	1	0	0	1	0	1.31	17
49	Computed stabilization for a giant fullerene endohedral: Y2C2@C-1(1660)-C-108 Slanina, Z; Uhlík, F; (...); Adamowicz, L Oct 16 2018 CHEMICAL PHYSICS LETTERS 710 , pp.147-149	5	5	1	5	0	2.67	16
	Local Heat Efficiency of a Polycrystalline Silicon Thin Layer for a							



50

Local pH and Effective pK of a Polyelectrolyte Chain: Two Names for One Quantity?

[Murmiliuk, A](#); [Kosovan, P](#); (...); [Stepanek, M](#)

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3

4

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16

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Previous page

Next page

Citations

< Previous year

Next year >

Average per year

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2020

2021

2022

2023

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51	Structural and electrochemical studies of Sm@D-3h-C-74 reveal a weak metal-cage interaction and a small band gap species Xu, W; Hao, YJ; (...); Feng, L 2013 NANOSCALE 5 (21) , pp.10409-10413	1	0	1	0	1	1.45	16
52	Ho2O@C-74: Ho2O Cluster Expands within a Small Non-IPR Fullerene Cage of C-2(13333)-C-74 Liu, A; Nie, MZ; (...); Uhlik, F Apr 15 2019 INORGANIC CHEMISTRY 58 (8) , pp.4774-4781	3	5	1	6	0	3	15
53	Adamantylidene Addition to M3N@I-h-C-80 (M=Sc, Lu) and Sc3N@D-5h-C-80: Synthesis and Crystallographic Characterization of the [5,6]-Open and [6,6]-Open Adducts Yamada, M; Abe, T; (...); Akasaka, T May 11 2017 CHEMISTRY-A EUROPEAN JOURNAL 23 (27) , pp.6552-6561	3	3	2	2	0	2	14
54	Mean squared displacement from fluorecence correlation spectroscopy Kubecka, J; Uhlik, F and Kosovan, P 2016 SOFT MATTER 12 (16) , pp.3760-3769	2	0	4	3	0	1.75	14
55	Synthesis of phosphinoferrocene amides and thioamides from carbamoyl chlorides and the structural chemistry of Group 11 metal complexes with these mixed-donor ligands Fernandes, TA; Solarova, H; (...); Stepnicka, P 2015 DALTON TRANSACTIONS 44 (7) , pp.3092-3108	2	0	1	4	0	1.56	14
56	Computations on Three Isomers of La@C-74 Slanina, Z; Uhlik, F; (...); Nagase, S Nov 15 2008 INTERNATIONAL JOURNAL OF QUANTUM CHEMISTRY 108 (14) , pp.2636-2640	0	0	0	0	0	0.88	14
57	Excited electronic states and relative stabilities of C-80 Slanina, Z; Lee, SL; (...); Nagase, S 9th Annual Meeting of the V-A-Fock-School on Quantum and Computational Chemistry Aug 15 2006 INTERNATIONAL JOURNAL OF QUANTUM CHEMISTRY 106 (10) , pp.2222-2228	1	1	0	0	0	0.78	14
58	Reaction of zirconacyclopentadienes with ethynylferrocenes Dufkova, I; Matsumura, H; (...); Kotora, M 2004 COLLECTION OF CZECHOSLOVAK CHEMICAL COMMUNICATIONS 69 (2) , pp.351-364	0	0	1	0	0	0.7	14

59	Computations of model narrow nanotubes closed by fragments of smaller fullerenes and quasi-fullerenes Slanina, Z; Uhlik, F and Adamowicz, L Jun 2003 JOURNAL OF MOLECULAR GRAPHICS & MODELLING 21 (6) , pp.517-522	0	0	0	0	0	0.67	14
60	Multivalent counterions accumulate in star-like polyelectrolytes and collapse the polymer in spite of increasing its ionization Stano, R; Nova, L; (...); Kosovan, P Jan 28 2020 SOFT MATTER 16 (4) , pp.1047-1055	0	2	4	6	1	3.25	13
61	Comparing the reactivity of isomeric phosphinoferrocene nitrile and isocyanide in Pd(ii) complexes: synthesis of simple coordination compounds vs. preparation of P-chelated insertion products and Fischer-type carbenes Skoch, K; Cisarova, J; (...); Stepnicka, P Dec 7 2018 DALTON TRANSACTIONS 47 (45) , pp.16082-16101	4	4	3	2	0	2.17	13
62	Probing the Influence of Phosphine Substituents on the Donor and Catalytic Properties of Phosphinoferrocene Carboxamides: A Combined Experimental and Theoretical Study Schulz, J; Vosahlo, P; (...); Stepnicka, P May 8 2017 ORGANOMETALLICS 36 (9) , pp.1828-1841	4	0	0	3	1	1.86	13
63	Endohedrally stabilized C-70 isomer with fused pentagons characterized by crystallography Feng, L; Zhang, MR; (...); Uhlik, F 2016 DALTON TRANSACTIONS 45 (19) , pp.8142-8148	4	2	0	2	0	1.63	13
64	Calculations of the water-dimer encapsulations into C-84 Slanina, Z; Uhlik, E; (...); Adamowicz, L 2016 FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 24 (1) , pp.1-7	2	0	4	1	0	1.63	13
65	Synthesis, Crystal Structures, and Electrochemical Behavior of Fe-Ru Heterobimetallic Complexes with Bridged Metallocene Units Schulz, J; Uhlik, E; (...); Stepnicka, P Sep 22 2014 ORGANOMETALLICS 33 (18) , pp.5020-5032	1	2	2	1	1	1.3	13
66	Synthesis and characterisation of Dewar benzene-ferrocene conjugates Jankova, S; Cisarova, J; (...); Kotora, M 2009 DALTON TRANSACTIONS (17) , pp.3137-3139	1	1	1	2	0	0.87	13
67	Computations on C84O: thermodynamic, kinetic and photochemical stability Slanina, Z; Uhlik, E; (...); Osawa, E Sep 27 2004 JOURNAL OF MOLECULAR STRUCTURE-THEOCHEM 684 (1-3) , pp.129-133	0	0	0	0	0	0.65	13
68	Quantum-chemical calculations of model systems of interest in fullerene-based superconductivity Slanina, Z; Uhlik, E; (...); Adamowicz, L Jun 2003 JOURNAL OF LOW TEMPERATURE PHYSICS 131 (5-6) , pp.1259-1263	2	0	0	0	0	0.62	13
	Computing molecular complexes in Earth's and other atmospheres Slanina, Z; Uhlik, E; (...); Osawa, E							

69	25th General Assembly of the European-Geophysical-Society 2001 PHYSICS AND CHEMISTRY OF THE EARTH PART C-SOLAR-TERRESTIAL AND PLANETARY SCIENCE 26 (7) , pp.505-511	0	0	1	1	0	0.57	13
70	Evaluation of the relative stabilities of two non-IPR isomers of Sm@C-76 Slanina, Z; Uhlík, E; (...); Adamowicz, L 2016 FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 24 (5) , pp.339-344	2	2	1	2	0	1.5	12
71	Water-Dimer Stability and Its Fullerene Encapsulations Uhlík, E; Slanina, Z; (...); Nagase, S Jun 2015 JOURNAL OF COMPUTATIONAL AND THEORETICAL NANOSCIENCE 12 (6) , pp.959-964	0	0	3	1	0	1.33	12
72	A fast computational method for determining equilibrium concentration profiles in intermixed nanoislands Uhlík, E; Gatti, R and Montalenti, F Feb 25 2009 JOURNAL OF PHYSICS-CONDENSED MATTER 21 (8)	0	0	0	1	0	0.8	12
73	Electronic excited states and stabilities of fullerenes: Isomers of C-78 and Mg@C-72 Slanina, Z; Uhlík, E; (...); Nagase, S 6th Annual Meeting of the V A Fock School on Quantum and Computational Chemistry Nov 15 2004 INTERNATIONAL JOURNAL OF QUANTUM CHEMISTRY 100 (4) , pp.610-616	1	0	0	0	0	0.6	12
74	A computational characterization of CO@C-60 Slanina, Z; Uhlík, E; (...); Lu, X 2017 FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 25 (11) , pp.624-629	1	1	3	2	0	1.57	11
75	Li-x@C-60: Calculations of the encapsulation energetics and thermodynamics Slanina, Z; Uhlík, E; (...); Nagase, S Sep 2008 INTERNATIONAL JOURNAL OF MOLECULAR SCIENCES 9 (9) , pp.1841-1850	1	0	1	0	0	0.69	11
76	Hydrophobically modified amphiphilic block copolymer micelles in non-aqueous polar solvents. Fluorometric, light scattering and computer-based Monte Carlo study Matejíček, P; Uhlík, E; (...); Webber, SE 2002 COLLECTION OF CZECHOSLOVAK CHEMICAL COMMUNICATIONS 67 (5) , pp.531-556	0	1	0	0	0	0.5	11
77	Synthesis and non-conventional structure of square-planar Pd(II) and Pt(II) complexes with an N,C,N-chelated stibinidene ligand Korenkova, M; Hejda, M; (...); Dostal, L Apr 28 2018 DALTON TRANSACTIONS 47 (16) , pp.5812-5822	2	1	2	4	0	1.67	10
78	Stability Computations for Isomers of La@C-n (n=72, 74, 76) Slanina, Z; Uhlík, E; (...); Nagase, S Nov 2012 MOLECULES 17 (11) , pp.13146-13156	1	0	0	0	0	0.83	10
79	SCF study of Amphiphilic micellar shells containing polyelectrolyte and hydrophobic sequences Jelinek, K; Limpouchova, Z; (...); Prochazka, K Oct 16 2007 MACROMOLECULES 40 (21) , pp.7656-7664	0	1	0	0	0	0.59	10

80	A computational treatment of 35 IPR isomers of C(88) Slanina, Z; Uhlík, E; (...); Osawa, E 2000 FULLERENE SCIENCE AND TECHNOLOGY 8 (4-5) , pp.417-432	0	0	0	1	0	0.42	10
81	AM1 COMPUTED RELATIVE EQUILIBRIUM POPULATIONS OF THE (6/6) AND (5/6) METHANOBUCKMINSTERFULLERENES C61H2 SLANINA, Z; LEE, SL; (...); ADAMOWICZ, L Oct 7 1994 CHEMICAL PHYSICS LETTERS 228 (4-5) , pp.490-494	0	0	0	0	0	0.33	10
82	COMPUTATIONAL STUDIES OF ATMOSPHERIC CHEMISTRY SPECIES .16. A COMPUTATIONAL THERMODYNAMIC EVALUATION OF THE ALTITUDE PROFILES OF (N2)2, N2-O2 AND (O2)2 IN THE EARTHS ATMOSPHERE SLANINA, Z; UHLIK, E; (...); HINCHLIFFE, A Jan 10 1994 THERMOCHIMICA ACTA 231 , pp.55-60	0	0	0	0	0	0.33	10
83	COMPUTATIONAL STUDIES OF ATMOSPHERIC CHEMISTRY SPECIES .6. AN ABINITIO CORRELATED STUDY OF STRUCTURE, ENERGETICS AND VIBRATIONS OF (N2)2(ALPHA) UHLIK, E; SLANINA, Z and HINCHLIFFE, A Jun 1993 JOURNAL OF MOLECULAR STRUCTURE-THEOCHEM 101 (3) , pp.271-275	0	0	0	0	0	0.32	10
84	A COMPUTATIONAL EVALUATION OF THE WATER-DIMER POPULATIONS IN SATURATED STEAM RECOMMENDED FOR APPLICATIONS TO THE EARTHS, PLANETARY AND COMETARY ATMOSPHERES SLANINA, Z; UHLIK, E and CRIFO, JF Jul 1992 JOURNAL OF MOLECULAR STRUCTURE 270 , pp.1-9	0	0	1	0	0	0.31	10
85	AN ESTIMATION OF DIMERIZATION ENERGETICS OF THE CLO RADICAL SLANINA, Z and UHLIK, E Jul 19 1991 CHEMICAL PHYSICS LETTERS 182 (1) , pp.51-56	0	0	0	0	0	0.3	10
86	Antimony(i) -> Pd(ii) complexes with the (mu-Sb)Pd-2 coordination framework Korenkova, M; Hejda, M; (...); Dostal, L Aug 21 2019 DALTON TRANSACTIONS 48 (31) , pp.11912-11920	1	2	2	4	0	1.8	9
87	Ho2O@C-84: Crystallographic Evidence Showing Linear Metallic Oxide Cluster Encapsulated in IPR Fullerene Cage of D-2d(51591)-C-84 Cong, HL; Liu, A; (...); Uhlík, E Aug 19 2019 INORGANIC CHEMISTRY 58 (16) , pp.10905-10911	1	3	1	4	0	1.8	9
88	Interactions of star-like polyelectrolyte micelles with hydrophobic counterions Fernandez-Alvarez, R; Nova, L; (...); Matejcek, P Jun 15 2019 JOURNAL OF COLLOID AND INTERFACE SCIENCE 546 , pp.371-380	0	4	1	3	1	1.8	9
89	Computed Relative Populations of D-2(22)-C-84 Endohedrals with Encapsulated Monomeric and Dimeric Water Slanina, Z; Uhlík, E; (...); Adamowicz, L Apr 18 2016 CHEMPHYSICHEM 17 (8) , pp.1109-1111	0	0	4	0	0	1.13	9

90	Calculated Temperature Development of the Relative Stabilities of Yb@C-82 Isomers Slanina, Z; Uhlik, F; (...); Akasaka, T Jan 1 2014 FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 22 (1-3) , pp.147-154	3	0	1	1	0	0.9	9
91	Stability calculations for Eu@C74 isomers Uhlik, F; Slanina, Z; (...); Nagase, S Mar 5 2013 INTERNATIONAL JOURNAL OF QUANTUM CHEMISTRY 113 (5) , pp.729-733	1	3	0	1	0	0.82	9
92	Calculated relative yields for Sc2S@C-82 and Y2S@C-82 Slanina, Z; Uhlik, F; (...); Adamowicz, L Oct 2011 THEORETICAL CHEMISTRY ACCOUNTS 130 (2-3) , pp.549-554	1	0	0	0	0	0.69	9
93	Computations of endohedral fullerenes: The Gibbs energy treatment Slanina, Z; Uhlik, F; (...); Nagase, S 2006 JOURNAL OF COMPUTATIONAL METHODS IN SCIENCES AND ENGINEERING 6 (1-4) , pp.243-250	1	1	0	0	0	0.5	9
94	C94IPR isomeric set: large-scale computations of relative stabilities based on the Gibbs function Slanina, Z; Zhao, X; (...); Lee, SL 6th World Congress of Theoretically Oriented Chemists (WATOC) Jul 25 2003 JOURNAL OF MOLECULAR STRUCTURE-THEOCHEM 630 , pp.205-213	0	1	0	0	0	0.43	9
95	Energy-entropy interplay of C60F36 isomers Slanina, Z and Uhlik, F Jun 4 2003 CHEMICAL PHYSICS LETTERS 374 (1-2) , pp.100-103	0	0	0	0	0	0.43	9
96	Thermodynamic data bases and optical isomerism Slanina, Z; Uhlik, F; (...); Osawa, E 14th Dubrovnik International Course and Conference on the Interfaces between Mathematics, Chemistry and Computer Sciences Dec 2000 CROATICA CHEMICA ACTA 73 (4) , pp.1047-1055	1	0	1	0	0	0.38	9
97	Thermodynamic properties of He inside fullerenes and nanotubes Uhlik, F; Slanina, Z and Osawa, E 8th Biennial International Workshop on Fullerenes and Atomic Clusters (IWFAc 99) 2000 MOLECULAR MATERIALS 13 (1-4) , pp.231-238	1	0	0	0	0	0.38	9
98	Calculated relative populations for the Eu@C-82 isomers Slanina, Z; Uhlik, F; (...); Adamowicz, L Jul 2019 CHEMICAL PHYSICS LETTERS 726 , pp.29-33	2	2	1	3	0	1.6	8
99	Fluorescence Spectroscopy as a Tool for Investigating the Self-Organized Polyelectrolyte Systems Prochazka, K; Limpouchova, Z; (...); Hof, M 2011 SELF ORGANIZED NANOSTRUCTURES OF AMPHIPHILIC BLOCK COPOLYMERS.I 241 , pp.187-249	0	0	0	0	0	0.62	8
	Mg@C-74 isomers: Calculated relative concentrations and comparison with Ca@C-74							



100

[Uhlik, F; Slanina, Z and Nagase, S](#)

7th International Conference on Trends in Nanotechnology (TNT 2006)

Jun 2007 | [PHYSICA STATUS SOLIDIA-APPLICATIONS AND MATERIALS SCIENCE](#) 204 (6) ,

pp.1905-1910

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219 Publications		Sort by Citations: highest first	Previous page		Next page		Citations					
			< Previous year					Next year >		Average per year	Total	
			2019	2020	2021	2022	2023					
							© 2022 Clarivate					
101	Model narrow nanotubes related to C-36, C-32 and C-20: initial computational structural sampling Slanina, Z; Zhao, X and Uhlik, E 6th International Conference on Atomically Controlled Surfaces Interfaces and Nanostructures (ACSIN-6) Nov 1 2002 MATERIALS SCIENCE AND ENGINEERING B-SOLID STATE MATERIALS FOR ADVANCED TECHNOLOGY 96 (2) , pp.164-168	0	1	2	1	0	0.36	8				
102	Isomeric C60F36(g) species: Computed structures and heats of formation Slanina, Z; Uhlik, E; (...); Kolesov, VP 5th International Workshop on Fullerene and Atomic Clusters 2002 PHYSICS OF THE SOLID STATE 44 (3) , pp.534-535	0	0	0	0	0	0.36	8				
103	Non-IPR fullerenes: C-36 and C-72 Slanina, Z; Zhao, X; (...); Osawa, E 13th International Winterschool on Electronic Properties of Novel Materials 1999 ELECTRONIC PROPERTIES OF NOVEL MATERIALS - SCIENCE AND TECHNOLOGY OF MOLECULAR NANOSTRUCTURES 486 , pp.179-182	0	0	0	0	0	0.32	8				
104	ELECTRON PAIRING AND CHEMICAL-BONDS - PHYSICAL MEANING OF EFFECTIVE PAIRS PONEC, R and UHLIK, E Dec 1994 COLLECTION OF CZECHOSLOVAK CHEMICAL COMMUNICATIONS 59 (12) , pp.2567-2578	0	0	0	0	0	0.27	8				
105	Enhancing photovoltaic performance of inverted perovskite solar cells via imidazole and benzoimidazole doping of PC61BM electron transport layer Wang, Y; Yang, Y; (...); Feng, L Mar 2020 ORGANIC ELECTRONICS 78	0	0	5	2	0	1.75	7				
106	Calculated relative populations of Sm@C-82 isomers Slanina, Z; Uhlik, E; (...); Adamowicz, L 2018 FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 26 (5) , pp.233-238	4	0	1	2	0	1.17	7				
107	Sc2O@C-78: Calculations of the yield ratio for two observed isomers Slanina, Z; Uhlik, E; (...); Adamowicz, L 2017 FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 25 (2) , pp.124-127	1	0	0	0	0	1	7				
108	Thermodynamics of Small Alkali Metal Halide Cluster Ions: Comparison of Classical Molecular Simulations with Experiment and Quantum Chemistry	0	0	0	0	0	0.78	7				

	109	Vlcek, L; Uhlik, E; (...); Chialvo, AA Jan 22 2015 JOURNAL OF PHYSICAL CHEMISTRY A 119 (3) , pp.488-500	0	0	0	0	0	0.78	7
	109	Stabilities of fullerenes: Illustration on C-80 Slanina, Z; Uhlik, E; (...); Nagase, S 2008 MATCH-COMMUNICATIONS IN MATHEMATICAL AND IN COMPUTER CHEMISTRY 59 (1) , pp.225-238	0	0	1	0	0	0.44	7
	110	Computational screening of metallofullerenes for nanoscience: Sr@C-74 Slanina, Z; Uhlik, E; (...); Nagase, S 2008 MOLECULAR SIMULATION 34 (1) , pp.17-21	0	0	0	0	0	0.44	7
	111	Computations of production yields for Ba@C-74 and Yb@C-74 Slanina, Z; Uhlik, E; (...); Nagase, S Nanotech 2006 Conference 2007 MOLECULAR SIMULATION 33 (7) , pp.563-568	0	0	0	0	0	0.41	7
	112	Mean-field study of poly(methacrylic acid) shells in partly hydrophobically modified amphiphilic block copolymer micelles in polar solvents Jelinek, K; Uhlik, E; (...); Prochazka, K 4th International Symposium on Polyelectrolytes Aug 14 2003 JOURNAL OF PHYSICAL CHEMISTRY B 107 (32) , pp.8241-8247	0	1	0	0	0	0.33	7
	113	Thermodynamic properties of He@C-60 Uhlik, E; Slanina, Z and Osawa, E 2000 FULLERENE SCIENCE AND TECHNOLOGY 8 (4-5) , pp.453-460	1	0	0	1	0	0.29	7
	114	On the definitions of bond index and valence for correlated wave functions Ponec, R; Uhlik, E; (...); Jug, K 10th Dubrovnik International Course / Conference MATH/CHEM/COMP Nov 1996 CROATICA CHEMICA ACTA 69 (3) , pp.933-940	0	0	1	0	0	0.25	7
	115	POPULATION ANALYSIS OF PAIR DENSITIES - A STUDY OF BASIS SET DEPENDENCE PONEC, R; UHLIK, E and COOPER, DL Apr 1995 CROATICA CHEMICA ACTA 68 (1) , pp.149-155	0	0	0	0	0	0.24	7
	116	COMPUTATIONAL STUDIES OF ATMOSPHERIC CHEMISTRY SPECIES .13. AN AB-INITIO CORRELATED STUDY OF STRUCTURE, ENERGETICS AND VIBRATIONS OF (O2)2(ALPHA) UHLIK, E; SLANINA, Z and HINCHLIFFE, A Oct 22 1993 JOURNAL OF MOLECULAR STRUCTURE-THEOCHEM 104 (3) , pp.273-276	0	0	0	0	0	0.23	7
	117	Calculations of the relative populations of Lu@C-82 isomers Slanina, Z; Uhlik, E; (...); Adamowicz, L Sep 2 2019 FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 27 (9) , pp.710-714	1	1	1	3	0	1.2	6
	118	COMPUTATIONAL STUDIES OF ATMOSPHERIC CHEMISTRY SPECIES .12. COMPUTED GAS-PHASE THERMODYNAMICS OF THE N2-O2 COMPLEXES UHLIK, E; SLANINA, Z; (...); HINCHLIFFE, A Sep 30 1993 THERMOCHIMICA ACTA 225 (1) , pp.1-7	0	0	0	0	0	0.19	6

119	Eu@C-86 isomers: Calculated relative populations Slanina, Z; Uhlík, E; (...); Adamowicz, L Jul 2 2020 Feb 2020 (Early Access) FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 28 (7) , pp.565-570	0	2	1	2	0	1.25	5
120	Computational Modeling of the Ce@C-82 Metallofullerene Isomeric Composition Slanina, Z; Uhlík, E; (...); Adamowicz, L Dec 2 2019 ECS JOURNAL OF SOLID STATE SCIENCE AND TECHNOLOGY 8 (12) , pp.M118-M121	0	2	1	2	0	1	5
121	Computational Comparison of the Water-Dimer Encapsulations into D-2(22)-C-84 and D-2d(23)-C-84 Slanina, Z; Uhlík, E; (...); Lu, X 2017 ECS JOURNAL OF SOLID STATE SCIENCE AND TECHNOLOGY 6 (6) , pp.M3113-M3115	0	0	3	0	0	0.71	5
122	Stability issues in computational screening of carbon nanostructures: illustrations on La endohedrals Slanina, Z; Uhlík, E; (...); Lu, X 2017 MOLECULAR SIMULATION 43 (17) , pp.1472-1479	2	0	0	0	0	0.71	5
123	Water-Dimer Stability and Its Fullerene Encapsulations (vol 12, pg 959, 2015) Uhlík, E; Slanina, Z; (...); Nagase, S Sep 2015 JOURNAL OF COMPUTATIONAL AND THEORETICAL NANOSCIENCE 12 (9) , pp.2622-2622	0	0	1	0	0	0.56	5
124	Unprecedented Chemical Reactivity of a Paramagnetic Endohedral Metallofullerene La@C-5-C-82 that Leads Hydrogen Addition in the 1,3-Dipolar Cycloaddition Reaction Takano, Y; Slanina, Z; (...); Akasaka, T Dec 17 2014 JOURNAL OF THE AMERICAN CHEMICAL SOCIETY 136 (50) , pp.17537-17546	2	2	0	0	0	0.5	5
125	Synthesis, molecular structure, electrochemistry and DFT study of a ferrocenyl-substituted 4-quinazolinone and related heterocycles Tauchman, J; Hladikova, K; (...); Stepnicka, P 2013 NEW JOURNAL OF CHEMISTRY 37 (7) , pp.2019-2030	0	0	0	0	0	0.45	5
126	Predicted stabilities of endohedral metallo-fullerenes La@C76 Uhlík, E; Slanina, Z; (...); Nagase, S Dec 2012 PHYSICA STATUS SOLIDI B-BASIC SOLID STATE PHYSICS 249 (12) , pp.2585-2587	0	0	0	0	0	0.42	5
127	Double-exponential decay of orientational correlations in semiflexible polyelectrolytes Bacova, P; Kosovan, P; (...); Prochazka, K Jun 2012 EUROPEAN PHYSICAL JOURNAL E 35 (6)	0	0	1	0	0	0.42	5
128	Stimuli-responsive amphiphilic shells of kinetically frozen polymeric micelles in aqueous media: Monte Carlo simulations and comparison to self-consistent field calculations Uhlík, E; Jelinek, K; (...); Prochazka, K May 27 2008 MACROMOLECULES 41 (10) , pp.3711-3719	0	1	0	0	0	0.31	5

129	COMPUTATIONAL STUDIES OF ATMOSPHERIC CHEMISTRY SPECIES .22. A COMPUTATIONAL STUDY OF CLONO2H+ - STRUCTURE AND VIBRATIONS SLANINA, Z; UHLIK, F and HINCHLIFFE, A 1994 SPECTROSCOPY LETTERS 27 (4) , pp.563-571	0	0	1	0	0	0.17	5
130	COMPUTATIONAL STUDIES OF ATMOSPHERIC CHEMISTRY SPECIES .14. GAS-PHASE ASSOCIATION OF O-2 - A COMPUTATIONAL THERMODYNAMIC STUDY UHLIK, F; SLANINA, Z and HINCHLIFFE, A Nov 15 1993 THERMOCHIMICA ACTA 228 , pp.9-14	0	0	0	0	0	0.16	5
131	Ho2O@D-3(85)-C-92: Highly Stretched Cluster Dictated by a Giant Cage and Unexplored Isomerization Yu, YL; Slanina, Z; (...); Feng, L Aug 3 2020 INORGANIC CHEMISTRY 59 (15) , pp.11020-11027	0	0	1	3	0	1	4
132	The drag of the tails: Diffusion of sticky nanoparticles in dilute polymer solutions Kuldova, J; Uhlik, F and Kosovan, P Dec 28 2015 JOURNAL OF CHEMICAL PHYSICS 143 (24)	0	0	1	0	0	0.44	4
133	Sm@C-74: Computed Relative Isomeric Populations Slanina, Z; Uhlik, F; (...); Nagase, S Jan 1 2014 FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 22 (1-3) , pp.235-242	0	0	0	0	0	0.4	4
134	Towards Relative Populations of Non-Isomeric Metallofullerenes: La@C-76(T-d) vs. La-2@C-76(C-S, 17490) Slanina, Z; Uhlik, F; (...); Akasaka, T Jan 1 2014 FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 22 (1-3) , pp.299-306	0	0	1	0	0	0.4	4
135	MONTE CARLO SIMULATION OF FLUORESCENCE CORRELATION SPECTROSCOPY DATA Kosovan, P; Uhlik, F; (...); Hof, M 2011 COLLECTION OF CZECHOSLOVAK CHEMICAL COMMUNICATIONS 76 (3) , pp.207-222	0	0	0	1	0	0.31	4
136	Depletion profiles for dilute solutions of linear chains, stars and H-branched molecules by self-consistent field calculations and Monte Carlo simulations Preisler, Z; Kosovan, P; (...); Leermakers, FAM 2011 SOFT MATTER 7 (21) , pp.10258-10265	0	0	1	0	0	0.31	4
137	Computations of the catalytic effects in the Stone-Wales fullerene isomerizations: N and CN agents Slanina, Z; Uhlik, F; (...); Lee, SL 4th Congress of the International-Society-for-Theoretical-Chemical-Physics (ICTCP 4) Sep 15 2004 INTERNATIONAL JOURNAL OF QUANTUM CHEMISTRY 99 (5) , pp.634-639	0	0	0	0	0	0.2	4
138	Amphiphilic block copolymer micelles with hydrophobically modified shells Jelinek, K; Uhlik, F; (...); Hof, M Yangtze Conference on Fluids and Interfaces Oct-nov 2003 MOLECULAR SIMULATION 29 (10-11) , pp.655-660	0	0	1	0	0	0.19	4
	Non-central location of LiInLi@C-60							

139	Slanina, Z; Uhlík, F and Chow, TJ International Symposium on Fullerenes, Nanotubes and Carbon Nanoclusters held at the Annual Spring Meeting of the Electrochemical-Society 2003 FULLERENES AND NANOTUBES: THE BUILDING BLOCKS OF NEXT GENERATION NANODEVICES 2003 (15) , pp.569-574	0	0	0	0	0	0.19	4
140	Clustering, saturated vapors, and the atmosphere: The (H2O)(n), H2O-N-2, and H2O-O-2 cases Slanina, Z and Uhlík, F NATO Advanced Research Workshop on Weakly Interacting Molecular Pairs - Unconventional Absorbers of Radiation in the Atmospher 2003 WEAKLY INTERACTING MOLECULAR PAIRS: UNCONVENTIONAL ABSORBERS OF RADIATION IN THE ATMOSPHERE 27 , pp.101-110	0	1	0	0	0	0.19	4
141	Geometrical and thermodynamic approaches to the relative stabilities of fullerene isomers (vol 44, pg 335, 2001) Slanina, Z; Uhlík, E; (...); Osawa, E Mar 2002 MATCH-COMMUNICATIONS IN MATHEMATICAL AND IN COMPUTER CHEMISTRY (45) , pp.173-173	0	1	0	0	0	0.18	4
142	Metal-coated fullerenes C(60)M(n): Calculations for M = Be, Mg, Al and n=12, 20, 32 Slanina, Z; Miyajima, C; (...); Osawa, E 2000 FULLERENE SCIENCE AND TECHNOLOGY 8 (4-5) , pp.385-402	0	0	0	0	0	0.17	4
143	Thermodynamic properties of N@C-60 Uhlík, F; Slanina, Z and Osawa, E 14th International Winter School on Electronic Properties on Novel Materials 2000 ELECTRONIC PROPERTIES OF NOVEL MATERIALS-MOLECULAR NANOSTRUCTURES 544 , pp.183-186	0	0	0	0	0	0.17	4
144	C60O and C60O2: Thermodynamics vs kinetics of formation Slanina, Z; Lee, SL; (...); Adamovic, L 3rd Symposium of the Fullerenes-Group of the Electrochemical-Society on Fullerenes - Chemistry, Physics, and New Directions VIII 1996 PROCEEDINGS OF THE SYMPOSIUM ON RECENT ADVANCES IN THE CHEMISTRY AND PHYSICS OF FULLERENES AND RELATED MATERIALS, VOL 3 96 (10) , pp.911-918	0	0	0	0	0	0.14	4
145	AM1 COMPUTATIONS OF C60O2 SLANINA, Z; UHLIK, E; (...); ADAMOWICZ, L 1994 FULLERENE SCIENCE AND TECHNOLOGY 2 (1) , pp.73-88	0	0	0	0	0	0.13	4
146	COMPUTATIONAL STUDIES OF ATMOSPHERIC CHEMISTRY SPECIES .9. COMPUTED GAS-PHASE THERMODYNAMICS OF N2 ASSOCIATION UHLIK, F; SLANINA, Z and HINCHLIFFE, A Aug 28 1993 THERMOCHIMICA ACTA 223 , pp.1-6	0	0	0	0	0	0.13	4
147	A REFINED EVALUATION OF THE GAS-PHASE DIMERIZATION THERMODYNAMICS OF THE CLO RADICAL SLANINA, Z and UHLIK, F Mar 1992 INTERNATIONAL JOURNAL OF THERMOPHYSICS 13 (2) , pp.303-313	0	0	0	0	0	0.13	4
148	Calculated Relative Thermodynamic Stabilities of the Gd@C-82 Isomers Slanina, Z; Uhlík, E; (...); Adamowicz, L Jul 1 2021 ECS JOURNAL OF SOLID STATE SCIENCE AND TECHNOLOGY 10 (7)	0	0	0	3	0	1	3

⊖ 149 Calculated relative populations for the Eu@C-84 isomers Slanina, Z; Uhlík, E; (…); Adamowicz, L 2021 Sep 2020 (Early Access) FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 29 (2) , pp.144-148 ☰ ★ Enriched Cited References	0	0	1	2	0	0.75	3
⊖ 150 Rotameric Isomers of La-2@C-80& Dodecafluoro-Subphthalocyanine Conjugate: Computational Characterization Slanina, Z; Uhlík, E; (…); Adamowicz, L Jan 8 2020 ECS JOURNAL OF SOLID STATE SCIENCE AND TECHNOLOGY 9 (6)	0	0	0	3	0	0.75	3

Citation Report Publications Table

219 Publications		Sort by Citations: highest first		Previous page Next page		Citations							
						< Previous year					Next year >	Average per year	Total
						2019	2020	2021	2022	2023			
						© 2022 Clarivate							
151	Intramolecular micellization and nanopatterning in pH- and thermo-responsive molecular brushes Prokacheva, VM; Rud, OV; (...); Borisov, OV Jan 7 2020 SOFT MATTER 16 (1) , pp.208-218	0	1	0	2	0	0.75	3					
152	Computations on Metallofullerenes Derivatized during Extraction: La@ C-80-C6H3Cl2 and La@C-82-C6H3Cl2 Slanina, Z; Uhlik, E; (...); Akasaka, T Jan 1 2014 FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 22 (1-3) , pp.173-181	1	0	0	0	0	0.3	3					
153	Computations on Encapsulations of Lanthanides into C-74 Slanina, Z; Uhlik, E and Nagase, S Nanotech 2009 Conference 2009 NANOTECH CONFERENCE & EXPO 2009, VOL 3, TECHNICAL PROCEEDINGS , pp.312-+	0	1	0	1	0	0.2	3					
154	Computational treatment of alkaline earth encapsulations in C-74: Relative thermodynamic production abundances Uhlik, E; Slanina, Z and Nagase, S 8th Biennial International Workshop on Fullerenes and Atomic Clusters (IWFAC 2007) 2008 FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 16 (5-6) , pp.507-516	0	0	1	0	0	0.19	3					
155	Relative stabilities of C-74 isomers Slanina, Z; Uhlik, E; (...); Nagase, S May-jun 2007 FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 15 (3) , pp.195-205	0	0	0	0	0	0.18	3					
156	Copolymer micelles with polyelectrolyte shell in aqueous media. A mean-field study Jelinek, K; Uhlik, E; (...); Prochazka, K 2006 COLLECTION OF CZECHOSLOVAK CHEMICAL COMMUNICATIONS 71 (5) , pp.756-768	0	0	0	0	0	0.17	3					
157	A computational study of entropy rules for charged fullerenes Slanina, Z; Uhlik, E and Boltalina, OV 5th International Workshop on Fullerene and Atomic Clusters 2002 PHYSICS OF THE SOLID STATE 44 (3) , pp.548-550	0	0	0	0	0	0.14	3					
158	Computations of low-energy non-icosahedral structures of C-60(6-) Slanina, Z; Uhlik, E; (...); Osawa, E 2000 FULLERENE SCIENCE AND TECHNOLOGY 8 (4-5) , pp.403-415	0	0	0	0	0	0.13	3					
159	B3LYP/G-31G**//SAM1 calculations of C-36 fullerene and quasi-fullerene cages												

159	Slanina, Z; Uhlík, E; (...); Osawa, E 2000 FULLERENE SCIENCE AND TECHNOLOGY 8 (4-5) , pp.433-447	0	0	0	0	0	0.13	3
160	COMPUTATIONAL STUDIES ATMOSPHERIC CHEMISTRY .19. A COMPUTATIONAL EVALUATION OF THE OZONE DIMER ALTITUDE PROFILE UHLÍK, E; SLANINA, Z and HINCHLIFFE, A Jan 24 1994 THERMOCHIMICA ACTA 232 (1) , pp.1-6	0	0	1	0	0	0.1	3
161	MULTIMOLECULAR CLUSTERS AND THEIR ISOMERISM .69. A COMPUTATIONAL EVALUATION OF RELATIVE STABILITIES OF THE OPEN C2V AND CYCLIC D3H FORMS OF OZONE SLANINA, Z and UHLÍK, E Feb 24 1992 THERMOCHIMICA ACTA 196 (2) , pp.459-465	0	0	0	0	0	0.09	3
162	Polyelectrolyte Hydrogels as Draw Agents for Desalination of Solutions with Multivalent Ions Rud, OV; Kazakov, AD; (...); Uhlík, E Mar 8 2022 MACROMOLECULES 55 (5) , pp.1763-1770  Enriched Cited References	0	0	0	2	0	1	2
163	Phase transition in hydrophobic weak polyelectrolyte gel utilized for water desalination Prokacheva, VM; Rud, OV; (...); Borisov, OV Sep 1 2021 Apr 2021 (Early Access) DESALINATION 511	0	0	0	2	0	0.67	2
164	Computer modeling of polymer stars in variable solvent conditions: a comparison of MD simulations, self-consistent field (SCF) modeling and novel hybrid Monte Carlo SCF approach Kazakov, AD; Prokacheva, VM; (...); Leermakers, FAM Jan 21 2021 SOFT MATTER 17 (3) , pp.580-591  Enriched Cited References	0	0	0	2	0	0.67	2
165	Calculations of the Lu3N@C-80 two-isomer equilibrium Slanina, Z; Uhlík, E; (...); Adamowicz, L May 4 2019 Mar 2019 (Early Access) FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 27 (5) , pp.382-386  Enriched Cited References	0	0	1	1	0	0.4	2
166	Cyclic water-trimer encapsulation into D-2(22)-C-84 fullerene Slanina, Z; Uhlík, E; (...); Adamowicz, L Mar 2018 CHEMICAL PHYSICS LETTERS 695 , pp.245-248	0	0	1	1	0	0.33	2
167	Vibrational spectroscopic and crystallographic study of the novel guanylurea salts with sulphuric and selenic acids Matulkova, J; Fridrichova, M; (...); Nemec, J Mar 5 2017 JOURNAL OF MOLECULAR STRUCTURE 1131 , pp.294-305	0	0	0	1	0	0.29	2
168	Calculations on endohedral C-74 complexes Slanina, Z; Uhlík, E; (...); Nagase, S Apr-may 2007 JOURNAL OF NANOSCIENCE AND NANOTECHNOLOGY 7 (4-5) , pp.1339-1345	0	0	0	0	0	0.12	2

169	Two isomers of C60F48: Computed inter-isomeric equilibrium Slanina, Z; Uhlik, F and Adamowicz, L 2003 FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 11 (3) , pp.219-226	0	0	0	0	0	0.1	2
170	A model study of narrow nanotubes related to C-36, C-32, C-20, and C-16: Semiempirical and nonempirical treatments Slanina, Z and Uhlik, F 2002 FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 10 (3) , pp.207-215	0	0	0	0	0	0.09	2
171	Structure of hexa-sulfobutyl fullerenes: A computational study Slanina, Z; Uhlik, E; (...); Adamowicz, L 2002 FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 10 (4) , pp.363-372	0	0	0	1	0	0.09	2
172	C(60)(NO(2))(2), C(60)(NO(2))(4), and C(60)(NO(2))(6): Computed energetics and structures Slanina, Z; Zhao, X; (...); Osawa, E 12th International Symposium on Recent Advances in the Chemistry and Physics of Fullerenes and Related Materials at the 195th Meeting of the Electrochemical-Society 1999 RECENT ADVANCES IN THE CHEMISTRY AND PHYSICS OF FULLERENES AND RELATED MATERIAL, VOL 7 99 (12) , pp.734-744	0	0	0	1	0	0.08	2
173	AM1 COMPUTED VIBRATIONS IN 2 C60O STRUCTURES SLANINA, Z; UHLIK, E; (...); ADAMOWICZ, L 1993 SPECTROSCOPY LETTERS 26 (9) , pp.1785-1796	0	0	0	0	0	0.06	2
174	A COMPUTATIONAL EVALUATION OF THE GAS-PHASE CLO DIMERIZATION THERMODYNAMICS SLANINA, Z and UHLIK, F Sep 1991 GEOPHYSICAL RESEARCH LETTERS 18 (9) , pp.1703-1706	0	0	0	0	0	0.06	2
175	Ho@C-82 Metallofullerene: Calculated Isomeric Composition Slanina, Z; Uhlik, E; (...); Adamowicz, L May 1 2022 ECS JOURNAL OF SOLID STATE SCIENCE AND TECHNOLOGY 11 (5)  Enriched Cited References	0	0	0	1	0	0.5	1
176	Ho2C2 Cluster with Flexible Configurations inside a Large C-2(61)-C-92 Cage Dong, B; Yu, YL; (...); Feng, L Jan 10 2022 Dec 2021 (Early Access) INORGANIC CHEMISTRY 61 (1) , pp.605-612  Enriched Cited References	0	0	0	1	0	0.33	1
177	A computational characterization of H2O2@C-60 Slanina, Z; Uhlik, E; (...); Lu, X Feb 1 2022 May 2021 (Early Access) FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 30 (2) , pp.258-262  Enriched Cited References	0	0	0	1	0	0.33	1
	Quantum-Chemical Calculations of the Metallofullerene Yields in the X@ C-74, L@C-74, and Z@C-82 Series							

178	Uhlik, F; Slanina, Z and Nagase, S International Conference of Computational Methods in Sciences and Engineering (ICCMSE) 2015 PROCEEDINGS OF THE INTERNATIONAL CONFERENCE OF COMPUTATIONAL METHODS IN SCIENCES AND ENGINEERING 2010 (ICCMSE-2010) 1642 , pp.269-272	1	0	0	0	0	0.11	1
179	Predicting stabilities of endohedral metallofullerenes Yb@C-84 Uhlik, F; Slanina, Z; (...); Nagase, S Dec 2013 PHYSICA STATUS SOLIDI B-BASIC SOLID STATE PHYSICS 250 (12) , pp.2709-2712	0	0	0	0	0	0.09	1
180	Alkali-metal clusters encapsulated into fullerenes: Computations on Li-x@C-60 Slanina, Z; Uhlik, F; (...); Nagase, S 2007 JOURNAL OF COMPUTATIONAL METHODS IN SCIENCES AND ENGINEERING 7 (5-6) , pp.541-547	0	0	0	0	0	0.06	1
181	Computations of the energetics of C60F36 isomers Slanina, Z; Uhlik, F; (...); Nagase, S 2006 FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 14 (1) , pp.57-65	0	0	0	0	0	0.06	1
182	C(94)IPR isomeric set: Large-scale computations of relative stabilities Slanina, Z; Zhao, X; (...); Osawa, E International Symposium on Fullerenes, Nanotubes, and Carbon Nanoclusters held at the 199th Meeting of the Electrochemical-Society 2001 FULLERENES FOR THE NEW MILLENNIUM 2000 (11) , pp.485-497	0	0	0	0	0	0.04	1
183	Computations of model extremely narrow nanotubes related to C-36, C-32, C-20, and C-16 Uhlik, F and Slanina, Z 15th International Winter School on Electronic Properties on Novel Materials 2001 ELECTRONIC PROPERTIES OF MOLECULAR NANOSTRUCTURES 591 , pp.458-461	0	0	0	0	0	0.04	1
184	Free and hindered translation of He inside C-60 Slanina, Z; Uhlik, F and Osawa, E 12th International Symposium on Recent Advances in the Chemistry and Physics of Fullerenes and Related Materials at the 195th Meeting of the Electrochemical-Society 1999 RECENT ADVANCES IN THE CHEMISTRY AND PHYSICS OF FULLERENES AND RELATED MATERIAL, VOL 7 99 (12) , pp.816-821	0	0	0	0	0	0.04	1
185	An analytical solution of isotopomeric enumerations for C-2v cyclic odd-numbered carbon clusters C-n Slanina, Z; Uhlik, F and Adamowicz, L May 1997 JOURNAL OF RADIOANALYTICAL AND NUCLEAR CHEMISTRY 219 (1) , pp.69-71	0	0	0	0	0	0.04	1
186	C61H2 FULLEROID - AM1 COMPUTATIONAL STUDY SLANINA, Z; UHLIK, F; (...); ADAMOWICZ, L 1994 FULLERENE SCIENCE AND TECHNOLOGY 2 (1) , pp.13-24	0	0	0	0	0	0.03	1
187	COMPUTATIONAL STUDIES OF ATMOSPHERIC CHEMISTRY SPECIES .4. ON THE THERMODYNAMICS OF THE CLOO AND OCLO RADICALS SLANINA, Z and UHLIK, F Mar 22 1993 THERMOCHIMICA ACTA 216 , pp.81-85	0	0	0	0	0	0.03	1
188	MULTIMOLECULAR CLUSTERS AND THEIR ISOMERISM .63. ISOMERIC INTERPLAY IN THE TRIMERIC SYSTEMS 2HF-HCN AND HF-2HCN							

188	SLANINA, Z; UHLIK, E; (...); HINCHLIFFE, A Dec 6 1991 JOURNAL OF MOLECULAR STRUCTURE-THEOCHEM 83 , pp.49-55	0	0	0	0	0	0.03	1
189	Hairy Gels: A Computational Study Uhlik, F; Rud, OV; (...); Zhulina, EB Dec 2022 GELS 8 (12) Enriched Cited References	0	0	0	0	0	0	0
190	A Computational Characterization of 2,2 '-bis(trifluoromethyl)-[1,1 '-biphenyl]-4,4 '-diamine Iodine Dopant for Improving Power-Conversion Efficiency of Perovskite Solar Cells Slanina, Z; Uhlik, E; (...); Adamowicz, L Nov 1 2022 ECS JOURNAL OF SOLID STATE SCIENCE AND TECHNOLOGY 11 (11)	0	0	0	0	0	0	0
191	Eu@C-88 Isomers: Calculated Relative Populations Uhlik, F; Slanina, Z; (...); Adamowicz, L Oct 1 2022 ECS JOURNAL OF SOLID STATE SCIENCE AND TECHNOLOGY 11 (10) Enriched Cited References	0	0	0	0	0	0	0
192	Salt Counterion Valency Controls the Ionization and Morphology of Weak Polyelectrolyte Miktoarm Stars Nova, L and Uhlik, F Jul 26 2022 Jul 2022 (Early Access) MACROMOLECULES 55 (14) , pp.6247-6259	0	0	0	0	0	0	0
193	C-60(OH)(32) fullerenols: calculated temperature-sensitive isomeric interplay Slanina, Z; Uhlik, E; (...); Adamowicz, L Dec 2 2022 May 2022 (Early Access) FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 30 (12) , pp.1193-1198 Enriched Cited References	0	0	0	0	0	0	0
194	Relative Production Yields in Homologous Metallofullerene Series: Computations for X@C-74 and Z@C-82 Endohedrals Uhlik, F; Slanina, Z; (...); Nagase, S 7th International Conference on Computational Methods in Science and Engineering (ICCMSE) 2012 INTERNATIONAL CONFERENCE OF COMPUTATIONAL METHODS IN SCIENCES AND ENGINEERING 2009 (ICCMSE 2009) 1504 , pp.1162-1165	0	0	0	0	0	0	0
195	Monte Carlo simulation of fluorescence correlation spectroscopy data for large and rigid fluorescently labeled particles undergoing both the translational and rotational diffusion Prochazka, K; Uhlik, F; (...); Limpouchova, Z 242nd National Meeting of the American-Chemical-Society (ACS) Aug 28 2011 ABSTRACTS OF PAPERS OF THE AMERICAN CHEMICAL SOCIETY 242	0	0	0	0	0	0	0
196	Calculated Production Yields in Metallofullerene Series Slanina, Z; Uhlik, E; (...); Nagase, S NSTI Nanotechnology Conference and Expo 2011 NANOTECHNOLOGY 2011: ELECTRONICS, DEVICES, FABRICATION, MEMS, FLUIDICS AND COMPUTATIONAL, NSTI-NANOTECH 2011, VOL 2 , pp.639-642	0	0	0	0	0	0	0
	Structural and bonding features of Z@C-82 (Z = Al, Sc, Y, La) endohedrals							

197	Slanina, Z; Uhlik, E; (...); Nagase, S 2010 JOURNAL OF COMPUTATIONAL METHODS IN SCIENCES AND ENGINEERING 10 (3-6) , pp.569-574	0	0	0	0	0	0	0
198	CONCENTRATION PROFILES IN HETEROEPITAXIAL NANOISLANDS Digiuni, D; Gatti, R; (...); Montalenti, F International Conference on Physics, Chemistry and Application of Nanostructures 2009 PHYSICS, CHEMISTRY AND APPLICATION OF NANOSTRUCTURES , pp.3-+	0	0	0	0	0	0	0
199	Computer study of stimuli-responsive polyelectrolyte micelles in aqueous media. Comparison of Monte Carlo and self-consistent field approaches Jelinek, K; Uhlik, E; (...); Prochazka, K 12th IUPAC International Symposium on Macromolecule-Metal Complexes 2008 MACROMOLECULAR SYMPOSIA 270 , pp.65-73	0	0	0	0	0	0	0
200	The use of Monte Carlo simulations for the interpretation of light scattering and fluorescence data on self-assembling polymer systems in solutions Matejicek, P; Uhlik, E; (...); Prochazka, K 2008 COLLECTION OF CZECHOSLOVAK CHEMICAL COMMUNICATIONS 73 (3) , pp.293-313	0	0	0	0	0	0	0

219 Publications		Sort by Citations: highest first	Previous page		Next page		Citations						
			< Previous year					Next year >		Average per year	Total		
			2019	2020	2021	2022	2023						
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⊖ 201	Computations on Li-x@C-60 Slanina, Z ; Uhlik, F and Nagase, S Nanotechnology Conference and Trade Show (Nanotech 2008) 2008 NSTI NANOTECH 2008, VOL 3, TECHNICAL PROCEEDINGS , pp.689-+						0	0	0	0	0	0	0
⊖ 202	Gibbs energy treatment of a set of C80F36 isomers Slanina, Z ; Uhlik, F and Adamowicz, L International Symposium on Fullerenes, Nanotubes and Carbon Nanoclusters held at the Annual Spring Meeting of the Electrochemical-Society 2003 FULLERENES AND NANOTUBES: THE BUILDING BLOCKS OF NEXT GENERATION NANODEVICES 2003 (15) , pp.480-488						0	0	0	0	0	0	0
⊖ 203	Structure of hexa-sulfobutyl fullerenes: A computational study (vol 10, pg 363, 2002) Slanina, Z ; Uhlik, F ; (...); Adamowicz, L 2003 FULLERENES NANOTUBES AND CARBON NANOSTRUCTURES 11 (2) , pp.187-187						0	0	0	0	0	0	0
⊖ 204	Excited electronic states and production optimizations for promising nano-agents Slanina, Z ; Uhlik, F ; (...); Nagase, S Nanotechnology Conference and Trade Show (Nanotech 2003) 2003 NANOTECH 2003, VOL 3 , pp.504-507						0	0	0	0	0	0	0
⊖ 205	Model narrow nanotubes related to C-36, C-32 and C-20: Computational insight Slanina, Z and Uhlik, F Symposium on Making Functional Materials with Nanotubes held at the 2001 MRS Fall Meeting 2002 MAKING FUNCTIONAL MATERIALS WITH NANOTUBES 706 , pp.271-276						0	0	0	0	0	0	0
⊖ 206	Computations of nanotubes as agents for molecular electronics: Narrow nanotubes related to C-36, C-32, C-20, and C-16 Slanina, Z and Uhlik, F 2nd International Conference on Computational Nanoscience and Nanotechnology 2002 ICCN 2002: INTERNATIONAL CONFERENCE ON COMPUTATIONAL NANOSCIENCE AND NANOTECHNOLOGY , pp.470-473						0	0	0	0	0	0	0
⊖ 207	Calculations of molar fractions of the IPR C-78 fullerenes Uhlik, F and Slanina, Z 16th International Winterschool on Electronic Properties of Novel Materials 2002 STRUCTURAL AND ELECTRONIC PROPERTIES OF MOLECULAR NANOSTRUCTURES 633 , pp.409-412						0	0	0	0	0	0	0
⊖ 208	Mg@C(72) endohedrals: MNDO/d computations of the isomeric composition Slanina, Z ; Zhao, X ; (...); Adamowicz, L International Symposium on Fullerenes, Nanotubes, and Carbon Nanoclusters held at the 199th Meeting of the Electrochemical-Society 2001 FULLERENES FOR THE NEW MILLENNIUM 2000 (11) , pp.332-340						0	0	0	0	0	0	0

209	Narrow nanotubes related to C-36, C-32, C-20, and C-16: Quantum-chemical semi-empirical and non-empirical model computations Slanina, Z and Uhlik, F International Symposium on Fullerenes, Nanotubes, and Carbon Nanoclusters held at the 199th Meeting of the Electrochemical-Society 2001 FULLERENES FOR THE NEW MILLENNIUM 2000 (11) , pp.371-380	0	0	0	0	0	0	0
210	Topological, geometrical, and thermodynamic evaluations of fullerene isomers Slanina, Z; Miyajima, C; (...); Osawa, E International Symposium on Fullerenes, Nanotubes, and Carbon Nanoclusters held at the 199th Meeting of the Electrochemical-Society 2001 FULLERENES FOR THE NEW MILLENNIUM 2000 (11) , pp.462-+	0	0	0	0	0	0	0
211	Quantum-chemical computations of Delta H(f,298)(o) for selected C(60)F(30)(g) isomers Slanina, Z; Uhlik, E; (...); Kolesov, VP International Symposium on Fullerenes, Nanotubes, and Carbon Nanoclusters held at the 199th Meeting of the Electrochemical-Society 2001 FULLERENES FOR THE NEW MILLENNIUM 2000 (11) , pp.498-506	0	0	0	0	0	0	0
212	Computations of hexa-sulfobutyl fullerenes: Energetics, structures, & vibrations Slanina, Z; Uhlik, E; (...); Osawa, E International Symposium on Fullerenes, Nanotubes, and Carbon Nanoclusters held at the 199th Meeting of the Electrochemical-Society 2001 FULLERENES FOR THE NEW MILLENNIUM 2000 (11) , pp.507-517	0	0	0	0	0	0	0
213	Computations on C(84)O: Thermochemical and photochemical aspects Slanina, Z; Uhlik, E; (...); Osawa, E International Symposium on Fullerenes, Nanotubes, and Carbon Nanoclusters held at the 199th Meeting of the Electrochemical-Society 2001 FULLERENES FOR THE NEW MILLENNIUM 2000 (11) , pp.518-529	0	0	0	0	0	0	0
214	Computations of metal-covered fullerenes - C60Ben Slanina, Z; Uhlik, E; (...); Osawa, E 4th Symposium of the Fullerenes-Group of the Electrochemical-Society, at the 191st Meeting of the Electrochemical-Society - Fullerenes: Chemistry, Physics, and New Directions IX 1997 PROCEEDINGS OF THE SYMPOSIUM ON RECENT ADVANCES IN THE CHEMISTRY AND PHYSICS OF FULLERENES AND RELATED MATERIALS, VOL 4 97 (14) , pp.712-720	0	0	0	0	0	0	0
215	COMPUTATIONAL STUDIES OF ATMOSPHERIC CHEMISTRY SPECIES .8. A COMPUTATIONAL STUDY OF THE CL00 AND OCL0 RADICALS SLANINA, Z; UHLIK, F and PONEC, R Jul-aug 1993 JOURNAL DE CHIMIE PHYSIQUE ET DE PHYSICO-CHIMIE BIOLOGIQUE 90 (7-8) , pp.1459-1466	0	0	0	0	0	0	0
216	A CLASSIFICATION OF 200 ISOMERIZATIONS AMONG 51 ISOTOPOMERS OF C-8 (D2D) SLANINA, Z; UHLIK, F and ADAMOWICZ, L May 1993 JOURNAL OF RADIOANALYTICAL AND NUCLEAR CHEMISTRY-ARTICLES 170 (2) , pp.373-380	0	0	0	0	0	0	0
217	CLASSIFICATION OF 486 ISOMERIZATIONS AMONG 72 C-12/C-13 ISOTOPOMERS OF CYCLIC C-7 SLANINA, Z; UHLIK, F and ADAMOWICZ, L Apr 1993 JOURNAL OF RADIOANALYTICAL AND NUCLEAR CHEMISTRY-ARTICLES 170 (1) , pp.107-116	0	0	0	0	0	0	0

⊖ 218 DIGALLANE (4) GA2H4 AS THE IDEAL-GAS PHASE-EQUILIBRIUM MIXTURE OF ITS 6 ISOMERS SLANINA, Z and UHLIK, F Nov 22 1991 THERMOCHIMICA ACTA 191 (1) , pp.47-55	0	0	0	0	0	0	0
⊖ 219 LINEAR VERSUS CYCLIC (HCN)3 - A COMPUTATIONAL THERMODYNAMIC STUDY SLANINA, Z and UHLIK, F Oct 8 1991 THERMOCHIMICA ACTA 188 (2) , pp.273-278	0	0	0	0	0	0	0

Citation Report Publications Table