

Michael H Abraham

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/4588085/publications.pdf>

Version: 2024-02-01

431
papers

29,403
citations

8732

75
h-index

7333

152
g-index

435
all docs

435
docs citations

435
times ranked

11552
citing authors

#	ARTICLE	IF	CITATIONS
1	Linear solvation energy relationships. 23. A comprehensive collection of the solvatochromic parameters, π^* , α , and β , and some methods for simplifying the generalized solvatochromic equation. <i>Journal of Organic Chemistry</i> , 1983, 48, 2877-2887.	1.7	3,340
2	Scales of solute hydrogen-bonding: their construction and application to physicochemical and biochemical processes. <i>Chemical Society Reviews</i> , 1993, 22, 73.	18.7	1,969
3	Determination of sets of solute descriptors from chromatographic measurements. <i>Journal of Chromatography A</i> , 2004, 1037, 29-47.	1.8	835
4	Hydrogen Bonding. 32. An Analysis of Water-Octanol and Water-Alkane Partitioning and the $\log P$ Parameter of Seiler. <i>Journal of Pharmaceutical Sciences</i> , 1994, 83, 1085-1100.	1.6	694
5	Rate-limited steps of human oral absorption and QSAR studies. <i>Pharmaceutical Research</i> , 2002, 19, 1446-1457.	1.7	576
6	Linear solvation energy relations. <i>Journal of Solution Chemistry</i> , 1985, 14, 153-186.	0.6	515
7	Linear solvation energy relationship. 46. An improved equation for correlation and prediction of octanol/water partition coefficients of organic nonelectrolytes (including strong hydrogen bond) $T_j \text{ ETQq1 } 1 \text{ } 0.784314 \text{ rgBT } / \text{ Overlock } 1$	1.7	406
8	Evaluation of human intestinal absorption data and subsequent derivation of a quantitative structure-activity relationship (QSAR) with the Abraham descriptors. <i>Journal of Pharmaceutical Sciences</i> , 2001, 90, 749-784.	1.6	436
9	Hydrogen bonding. Part 34. The factors that influence the solubility of gases and vapours in water at 298 K, and a new method for its determination. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1994, , 1777.	0.9	425
10	Estimation of Molecular Linear Free Energy Relation Descriptors Using a Group Contribution Approach. <i>Journal of Chemical Information and Computer Sciences</i> , 1999, 39, 835-845.	2.8	410
11	Classification of stationary phases and other materials by gas chromatography. <i>Journal of Chromatography A</i> , 1999, 842, 79-114.	1.8	351
12	Hydrogen bonding. Part 10. A scale of solute hydrogen-bond basicity using log K values for complexation in tetrachloromethane. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1990, , 521.	0.9	349
13	Hydrogen bonding. 31. Construction of a scale of solute effective or summation hydrogen-bond basicity. <i>Journal of Physical Organic Chemistry</i> , 1993, 6, 660-684.	0.9	338
14	Hydrogen bonding. Part 7. A scale of solute hydrogen-bond acidity based on log K values for complexation in tetrachloromethane. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1989, , 699.	0.9	327
15	Solvent effects in organic chemistry - recent developments. <i>Canadian Journal of Chemistry</i> , 1988, 66, 2673-2686.	0.6	272
16	Hydrogen Bonding. 33. Factors That Influence the Distribution of Solutes between Blood and Brain. <i>Journal of Pharmaceutical Sciences</i> , 1994, 83, 1257-1268.	1.6	272
17	Hydrogen bonding. <i>Journal of Chromatography A</i> , 1991, 587, 213-228.	1.8	267
18	Prediction of Solubility of Drugs and Other Compounds in Organic Solvents. <i>Journal of Pharmaceutical Sciences</i> , 2010, 99, 1500-1515.	1.6	266

#	ARTICLE	IF	CITATIONS
19	The correlation and prediction of the solubility of compounds in water using an amended solvation energy relationship. <i>Journal of Pharmaceutical Sciences</i> , 1999, 88, 868-880.	1.6	262
20	Hydrogen bonding. Part 13. A new method for the characterisation of GLC stationary phases—the laffort data set. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1990, , 1451-1460.	0.9	244
21	Hydrogen Bond Structural Group Constants. <i>Journal of Organic Chemistry</i> , 2001, 66, 3484-3491.	1.7	230
22	Study of retention in reversed-phase liquid chromatography using linear solvation energy relationships I. The stationary phase. <i>Journal of Chromatography A</i> , 1996, 752, 1-18.	1.8	200
23	Correlation and prediction of a large blood–brain distribution data set—an LFER study. <i>European Journal of Medicinal Chemistry</i> , 2001, 36, 719-730.	2.6	200
24	Some Novel Liquid Partitioning Systems: Water/Ionic Liquids and Aqueous Biphasic Systems. <i>Industrial & Engineering Chemistry Research</i> , 2003, 42, 413-418.	1.8	186
25	Human Skin Permeation and Partition: General Linear Free Energy Relationship Analyses. <i>Journal of Pharmaceutical Sciences</i> , 2004, 93, 1508-1523.	1.6	182
26	Determination of olive oil–gas and hexadecane–gas partition coefficients, and calculation of the corresponding olive oil–water and hexadecane–water partition coefficients. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1987, , 797-803.	0.9	181
27	HYDROGEN BONDING. 42. CHARACTERIZATION OF REVERSED-PHASE HIGH-PERFORMANCE LIQUID CHROMATOGRAPHIC C18 STATIONARY PHASES. <i>Journal of Physical Organic Chemistry</i> , 1997, 10, 358-368.	0.9	173
28	The Factors that Influence Skin Penetration of Solutes. <i>Journal of Pharmacy and Pharmacology</i> , 2011, 47, 8-16.	1.2	166
29	Application of hydrogen bonding calculations in property based drug design. <i>Drug Discovery Today</i> , 2002, 7, 1056-1063.	3.2	163
30	Estimation of Molecular Linear Free Energy Relationship Descriptors by a Group Contribution Approach. 2. Prediction of Partition Coefficients. <i>Journal of Chemical Information and Computer Sciences</i> , 2000, 40, 71-80.	2.8	151
31	Thermodynamics of solution of homologous series of solutes in water. <i>Journal of the Chemical Society Faraday Transactions I</i> , 1984, 80, 153.	1.0	146
32	The factors that influence permeation across the blood–brain barrier. <i>European Journal of Medicinal Chemistry</i> , 2004, 39, 235-240.	2.6	143
33	Calculations on ionic solvation. Part 1.—Free energies of solvation of gaseous univalent ions using a one-layer continuum model. <i>Journal of the Chemical Society Faraday Transactions I</i> , 1978, 74, 1604.	1.0	139
34	Hydrogen bonding. 38. Effect of solute structure and mobile phase composition on reversed-phase high-performance liquid chromatographic capacity factors. <i>Journal of Physical Organic Chemistry</i> , 1994, 7, 672-684.	0.9	136
35	A data base for partition of volatile organic compounds and drugs from blood/plasma/serum to brain, and an LFER analysis of the data. <i>Journal of Pharmaceutical Sciences</i> , 2006, 95, 2091-2100.	1.6	131
36	Comparative analysis of solvation and selectivity in room temperature ionic liquids using the Abraham linear free energy relationship. <i>Green Chemistry</i> , 2006, 8, 906.	4.6	130

#	ARTICLE	IF	CITATIONS
37	On the Partition of Ampholytes: Application to Blood-Brain Distribution. <i>Journal of Pharmaceutical Sciences</i> , 1997, 86, 310-315.	1.6	128
38	Rapid Gradient RP-HPLC Method for Lipophilicity Determination: A Solvation Equation Based Comparison with Isocratic Methods. <i>Analytical Chemistry</i> , 1998, 70, 4228-4234.	3.2	126
39	A general treatment of hydrogen bond complexation constants in tetrachloromethane. <i>Journal of the American Chemical Society</i> , 1988, 110, 8534-8536.	6.6	125
40	Hydrogen bonding. <i>Journal of Chromatography A</i> , 1993, 644, 95-139.	1.8	125
41	A Comparison between the Two General Sets of Linear Free Energy Descriptors of Abraham and Klamt. <i>Journal of Chemical Information and Computer Sciences</i> , 2002, 42, 1320-1331.	2.8	122
42	Predicting Penetration Across the Blood-Brain Barrier from Simple Descriptors and Fragmentation Schemes. <i>Journal of Chemical Information and Modeling</i> , 2007, 47, 170-175.	2.5	122
43	Molecular Factors Influencing Drug Transfer across the Blood-Brain Barrier. <i>Journal of Pharmacy and Pharmacology</i> , 2011, 49, 1211-1216.	1.2	122
44	Equations for the Transfer of Neutral Molecules and Ionic Species from Water to Organic phases. <i>Journal of Organic Chemistry</i> , 2010, 75, 1006-1015.	1.7	121
45	Nasal pungency and odor of homologous aldehydes and carboxylic acids. <i>Experimental Brain Research</i> , 1998, 118, 180-188.	0.7	118
46	Calculation of Abraham descriptors from solvent-water partition coefficients in four different systems; evaluation of different methods of calculation. <i>Perkin Transactions II RSC</i> , 2002, , 470-477.	1.1	117
47	Solubility of electrolytes in 1,2-dichloroethane and 1,1-dichloroethane, and derived free energies of transfer. <i>Journal of the Chemical Society Faraday Transactions I</i> , 1976, 72, 955.	1.0	113
48	Enthalpy of Solvation Correlations for Gaseous Solutes Dissolved in Water and in 1-Octanol Based on the Abraham Model. <i>Journal of Chemical Information and Modeling</i> , 2007, 47, 115-121.	2.5	107
49	Algorithms for Skin Permeability Using Hydrogen Bond Descriptors: the Problem of Steroids. <i>Journal of Pharmacy and Pharmacology</i> , 2011, 49, 858-865.	1.2	104
50	Linear free energy relationship analysis of microemulsion electrokinetic chromatographic determination of lipophilicity. <i>Journal of Chromatography A</i> , 1996, 752, 243-249.	1.8	102
51	Solubility predictions for crystalline polycyclic aromatic hydrocarbons (PAHs) dissolved in organic solvents based upon the Abraham general solvation model. <i>Fluid Phase Equilibria</i> , 2002, 201, 245-258.	1.4	99
52	The analysis of solvation in ionic liquids and organic solvents using the Abraham linear free energy relationship. <i>Journal of Chemical Technology and Biotechnology</i> , 2006, 81, 1441-1446.	1.6	98
53	Characterization of the sorption of gaseous and organic solutes onto polydimethyl siloxane solid-phase microextraction surfaces using the Abraham model. <i>Journal of Chromatography A</i> , 2007, 1175, 162-173.	1.8	97
54	Single-ion gibbs energies, enthalpies and entropies of transfer from water to aqueous methanol based on the $(\text{Ph}_4\text{P}^+, \text{Ph}_4\text{As}^+) = \text{Ph}_4\text{B}^-$ assumption. <i>Journal of the Chemical Society Faraday Transactions I</i> , 1984, 80, 489.	1.0	94

#	ARTICLE	IF	CITATIONS
55	Correlation and estimation of gas-liquid and water-liquid partition coefficients by a linear free energy relationship method. <i>Journal of Pharmaceutical Sciences</i> , 1999, 88, 670-679.	1.6	94
56	Calculation of Abraham descriptors from experimental data from seven HPLC systems; evaluation of five different methods of calculation. Electronic supplementary information (ESI) available: Tables S1 to S5. See http://www.rsc.org/suppdata/p2/b2/b206927j/ . <i>Perkin Transactions II RSC</i> , 2002, , 2001-2010.	1.1	94
57	Chemical Toxicity Correlations for Several Fish Species Based on the Abraham Solvation Parameter Model. <i>Chemical Research in Toxicology</i> , 2005, 18, 1497-1505.	1.7	94
58	Hydrogen bonding. Part 40. Factors that influence the distribution of solutes between water and sodium dodecylsulfate micelles. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1995, , 887.	0.9	93
59	Chromatographic Estimation of Drug Disposition Properties by Means of Immobilized Artificial Membranes (IAM) and C18 Columns. <i>Journal of Medicinal Chemistry</i> , 2006, 49, 4861-4870.	2.9	92
60	An algorithm for nasal pungency thresholds in man. <i>Archives of Toxicology</i> , 1998, 72, 227-232.	1.9	89
61	Hydrogen bonding part 46: a review of the correlation and prediction of transport properties by an LFER method: physicochemical properties, brain penetration and skin permeability. <i>Pest Management Science</i> , 1999, 55, 78-88.	0.7	89
62	Evaluation of rat intestinal absorption data and correlation with human intestinal absorption. <i>European Journal of Medicinal Chemistry</i> , 2003, 38, 233-243.	2.6	89
63	The molecular properties governing solubilities of organic nonelectrolytes in water. <i>Nature</i> , 1985, 313, 384-386.	13.7	86
64	A quantitative measure of solvent solvophobic effect. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1988, , 339.	0.9	86
65	Factors that influence tadpole narcosis. An LFER analysis. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1995, , 1843.	0.9	86
66	Solute-solvent interactions in micellar electrokinetic chromatography. <i>Journal of Chromatography A</i> , 2002, 942, 237-248.	1.8	85
67	Title is missing!. <i>Journal of Solution Chemistry</i> , 2002, 31, 293-303.	0.6	83
68	Partition of solutes into wet and dry ethers; an LFER analysis. Electronic supplementary information (ESI) available: tables of solute descriptors and log P values. Values of log Kw and L. See http://www.rsc.org/suppdata/nj/b3/b303016d/ . <i>New Journal of Chemistry</i> , 2003, 27, 1041.	1.4	83
69	Determination of Solvation Descriptors for Ionic Species: Hydrogen Bond Acidity and Basicity. <i>Journal of Organic Chemistry</i> , 2004, 69, 4677-4685.	1.7	83
70	An NMR Method for the Quantitative Assessment of Intramolecular Hydrogen Bonding; Application to Physicochemical, Environmental, and Biochemical Properties. <i>Journal of Organic Chemistry</i> , 2014, 79, 11075-11083.	1.7	83
71	Solvation descriptors for ferrocene, and the estimation of some physicochemical and biochemical properties. <i>New Journal of Chemistry</i> , 2000, 24, 825-829.	1.4	81
72	Draize Rabbit Eye Test Compatibility with Eye Irritation Thresholds in Humans: A Quantitative Structure-Activity Relationship Analysis. <i>Toxicological Sciences</i> , 2003, 76, 384-391.	1.4	81

#	ARTICLE	IF	CITATIONS
73	NMR Method for the Determination of Solute Hydrogen Bond Acidity. <i>Journal of Organic Chemistry</i> , 2006, 71, 3389-3394.	1.7	81
74	Linear solvation energy relationships. Part 37. An analysis of contributions of dipolarityâ€“polarisability, nucleophilic assistance, electrophilic assistance, and cavity terms to solvent effects on t-butyl halide solvolysis rates. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1987, , 913-920.	0.9	79
75	Solvation descriptors for pesticides from the solubility of solids: diuron as an example. <i>Pest Management Science</i> , 2000, 56, 1043-1053.	1.7	79
76	Characterization of Room-Temperature Ionic Liquids by the Abraham Model with Cation-Specific and Anion-Specific Equation Coefficients. <i>Journal of Chemical Information and Modeling</i> , 2007, 47, 1123-1129.	2.5	79
77	Hydrogen bonding. <i>Journal of Chromatography A</i> , 1992, 594, 229-241.	1.8	78
78	Partition of solutes from the gas phase and from water to wet and dry di-n-butyl ether: a linear free energy relationship analysis. <i>Physical Chemistry Chemical Physics</i> , 2001, 3, 3732-3736.	1.3	78
79	Calculations on ionic solvation. III. The electrostatic free energy of solvation of ions, using a multilayered continuum model. <i>Journal of Chemical Physics</i> , 1979, 70, 2491-2496.	1.2	77
80	Descriptors for solutes from the solubility of solids: trans-stilbene as an example. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1998, , 2677-2682.	0.9	77
81	Human olfactory detection of homologous n-alcohols measured via concentrationâ€“response functions. <i>Pharmacology Biochemistry and Behavior</i> , 2008, 89, 279-291.	1.3	76
82	Partition of Volatile Organic Compounds from Air and from Water into Plant Cuticular Matrix:â€“ An LFER Analysis. <i>Environmental Science & Technology</i> , 2000, 34, 318-323.	4.6	74
83	The partition of compounds from water and from air into wet and dry ketones. <i>New Journal of Chemistry</i> , 2009, 33, 568-573.	1.4	74
84	The solubility of gases and vapours in dry octan-1-ol at 298 K. <i>Chemosphere</i> , 2001, 44, 855-863.	4.2	73
85	Air to Blood Distribution of Volatile Organic Compounds:â€“ A Linear Free Energy Analysis. <i>Chemical Research in Toxicology</i> , 2005, 18, 904-911.	1.7	73
86	Partition of compounds from gas to water and from gas to physiological saline at 310K: Linear free energy relationships. <i>Fluid Phase Equilibria</i> , 2007, 251, 93-109.	1.4	73
87	The transfer of neutral molecules, ions and ionic species from water to wet octanol. <i>Physical Chemistry Chemical Physics</i> , 2010, 12, 13182.	1.3	73
88	Method for Unknown Vapor Characterization and Classification Using a Multivariate Sorption Detector. Initial Derivation and Modeling Based on Polymer-Coated Acoustic Wave Sensor Arrays and Linear Solvation Energy Relationships. <i>Analytical Chemistry</i> , 1999, 71, 4544-4553.	3.2	72
89	The transfer of neutral molecules, ions and ionic species from water to ethylene glycol and to propylene carbonate; descriptors for pyridinium cations. <i>New Journal of Chemistry</i> , 2010, 34, 2298.	1.4	72
90	Linear solvation energy relationships. 29. Solution properties of some tetraalkylammonium halide ion pairs and dissociated ions. <i>Journal of the American Chemical Society</i> , 1985, 107, 3105-3110.	6.6	71

#	ARTICLE	IF	CITATIONS
91	On the mechanism of human intestinal absorption. <i>European Journal of Medicinal Chemistry</i> , 2002, 37, 595-605.	2.6	71
92	Linear Free Energy Relationship Correlations for Room Temperature Ionic Liquids: Revised Cation-Specific and Anion-Specific Equation Coefficients for Predictive Applications Covering a Much Larger Area of Chemical Space. <i>Industrial & Engineering Chemistry Research</i> , 2009, 48, 4145-4154.	1.8	69
93	Free energies of solution of rare gases and alkanes in water and nonaqueous solvents. A quantitative assessment of the hydrophobic effect. <i>Journal of the American Chemical Society</i> , 1979, 101, 5477-5484.	6.6	68
94	Solubility properties in polymers and biological media. 2. The correlation and prediction of the solubilities of nonelectrolytes in biological tissues and fluids. <i>Journal of Medicinal Chemistry</i> , 1985, 28, 865-870.	2.9	67
95	Chemosensory Detectability of 1-Butanol and 2-Heptanone Singly and in Binary Mixtures. <i>Physiology and Behavior</i> , 1999, 67, 269-276.	1.0	67
96	Determination of McGowan Volumes for Ions and Correlation with van der Waals Volumes. <i>Journal of Chemical Information and Computer Sciences</i> , 2003, 43, 1848-1854.	2.8	67
97	Detection of single and mixed VOCs by smell and by sensory irritation. <i>Indoor Air</i> , 2004, 14, 108-117.	2.0	67
98	Solubility of 9-fluorenone, thianthrene and xanthene in organic solvents. <i>Fluid Phase Equilibria</i> , 2005, 232, 113-121.	1.4	67
99	Diffusion coefficients in ethanol and in water at 298K: Linear free energy relationships. <i>Fluid Phase Equilibria</i> , 2011, 303, 45-55.	1.4	67
100	Lipophilicity of the Nitrophenols. <i>Journal of Organic Chemistry</i> , 2000, 65, 7114-7118.	1.7	66
101	Hydrogen bonding. Part 45. The solubility of gases and vapours in methanol at 298 K: an LFER analysis. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1998, , 1385-1390.	0.9	65
102	Characterizing the Selectivity of Stationary Phases and Organic Modifiers in Reversed-Phase High-Performance Liquid Chromatographic Systems by a General Solvation Equation Using Gradient Elution. <i>Journal of Chromatographic Science</i> , 2000, 38, 503-511.	0.7	65
103	Solubility predictions for crystalline nonelectrolyte solutes dissolved in organic solvents based upon the Abraham general solvation model. <i>Canadian Journal of Chemistry</i> , 2001, 79, 1466-1476.	0.6	65
104	Air to liver partition coefficients for volatile organic compounds and blood to liver partition coefficients for volatile organic compounds and drugs. <i>European Journal of Medicinal Chemistry</i> , 2007, 42, 743-751.	2.6	65
105	The Solvation Properties of the Aliphatic Alcohols. <i>Collection of Czechoslovak Chemical Communications</i> , 1999, 64, 1748-1760.	1.0	64
106	Correlation and prediction of the solubility of Buckminsterfullerene in organic solvents; estimation of some physicochemical properties. <i>Perkin Transactions II RSC</i> , 2000, , 281-286.	1.1	64
107	Solubility of crystalline nonelectrolyte solutes in organic solvents: mathematical correlation of 4-chloro-3-nitrobenzoic acid and 2-chloro-5-nitrobenzoic acid solubilities with the Abraham solvation parameter model. <i>Physics and Chemistry of Liquids</i> , 2005, 43, 351-360.	0.4	64
108	Correlation and prediction of partition coefficient between the gas phase and water, and the solvents dry methyl acetate, dry and wet ethyl acetate, and dry and wet butyl acetate. <i>Fluid Phase Equilibria</i> , 2008, 270, 30-44.	1.4	64

#	ARTICLE	IF	CITATIONS
109	Solute Descriptors for Phenoxide Anions and Their Use To Establish Correlations of Rates of Reaction of Anions with Iodomethane. <i>Journal of Organic Chemistry</i> , 2010, 75, 3021-3026.	1.7	64
110	Solute-solvent interactions in micellar electrokinetic chromatography. <i>Journal of Chromatography A</i> , 1999, 845, 217-226.	1.8	63
111	Air to Muscle and Blood/Plasma to Muscle Distribution of Volatile Organic Compounds and Drugs: Linear Free Energy Analyses. <i>Chemical Research in Toxicology</i> , 2006, 19, 801-808.	1.7	63
112	Hydrogen bonding. 39. The partition of solutes between water and various alcohols. <i>Journal of Physical Organic Chemistry</i> , 1994, 7, 712-716.	0.9	62
113	Solubility Properties in Biological Media. 12. Regarding the Mechanism of Nonspecific Toxicity or Narcosis by Organic Nonelectrolytes. <i>QSAR and Combinatorial Science</i> , 1988, 7, 71-78.	1.4	60
114	Partition of compounds from water and from air into amides. <i>New Journal of Chemistry</i> , 2009, 33, 2034.	1.4	60
115	An Algorithm for 353 Odor Detection Thresholds in Humans. <i>Chemical Senses</i> , 2012, 37, 207-218.	1.1	60
116	Linear solvation energy relationships. 15. Heterolytic decomposition of the tert-butyl halides. <i>Journal of Organic Chemistry</i> , 1981, 46, 3053-3056.	1.7	59
117	Hydrogen-bonding 8. Possible equivalence of solute and solvent scales of hydrogen-bond basicity of non-associated compounds. <i>Journal of Physical Organic Chemistry</i> , 1989, 2, 540-552.	0.9	59
118	Group Contribution and Machine Learning Approaches to Predict Abraham Solute Parameters, Solvation Free Energy, and Solvation Enthalpy. <i>Journal of Chemical Information and Modeling</i> , 2022, 62, 433-446.	2.5	59
119	Hydrogen bonding. 47. Characterization of the ethylene glycol-heptane partition system: Hydrogen bond acidity and basicity of peptides. <i>Journal of Pharmaceutical Sciences</i> , 1999, 88, 241-247.	1.6	58
120	Air to brain, blood to brain and plasma to brain distribution of volatile organic compounds: linear free energy analyses. <i>European Journal of Medicinal Chemistry</i> , 2006, 41, 494-502.	2.6	58
121	Descriptors for ions and ion-pairs for use in linear free energy relationships. <i>Journal of Chromatography A</i> , 2016, 1430, 2-14.	1.8	58
122	Air to fat and blood to fat distribution of volatile organic compounds and drugs: Linear free energy analyses. <i>European Journal of Medicinal Chemistry</i> , 2006, 41, 1430-1438.	2.6	57
123	Retention properties of a spacer-bonded propanediol sorbent for reversed-phase liquid chromatography and solid-phase extraction. <i>Analyst</i> , 1996, 121, 511.	1.7	56
124	Solubility of Crystalline Nonelectrolyte Solutes in Organic Solvents: Mathematical Correlation of Acetylsalicylic Acid Solubilities with the Abraham General Solvation Model. <i>Journal of Solution Chemistry</i> , 2003, 32, 1087-1102.	0.6	56
125	Octanol/Water Partition of Ionic Species, Including 544 Cations. <i>Journal of Organic Chemistry</i> , 2005, 70, 2633-2640.	1.7	56
126	Solubility of gases and vapours in propan-1-ol at 298 K. <i>Journal of Physical Organic Chemistry</i> , 1999, 12, 675-680.	0.9	55

#	ARTICLE	IF	CITATIONS
127	A Model for Odour Thresholds. <i>Chemical Senses</i> , 2002, 27, 95-104.	1.1	54
128	A Quantitative Structure Activity Analysis on the Relative Sensitivity of the Olfactory and the Nasal Trigeminal Chemosensory Systems. <i>Chemical Senses</i> , 2007, 32, 711-719.	1.1	54
129	The solubility of gases and vapours in ethanol - the connection between gaseous solubility and water-solvent partition. <i>Canadian Journal of Chemistry</i> , 1998, 76, 703-709.	0.6	53
130	Human Skin Permeation of Neutral Species and Ionic Species: Extended Linear Free Energy Relationship Analyses. <i>Journal of Pharmaceutical Sciences</i> , 2012, 101, 2034-2044.	1.6	53
131	Determinants for Nasal Trigeminal Detection of Volatile Organic Compounds. <i>Chemical Senses</i> , 2005, 30, 627-642.	1.1	52
132	Odor Detection by Humans of Lineal Aliphatic Aldehydes and Helional as Gauged by Dose-Response Functions. <i>Chemical Senses</i> , 2010, 35, 289-299.	1.1	52
133	Air to lung partition coefficients for volatile organic compounds and blood to lung partition coefficients for volatile organic compounds and drugs. <i>European Journal of Medicinal Chemistry</i> , 2008, 43, 478-485.	2.6	51
134	Development of Abraham model correlations for solvation characteristics of linear alcohols. <i>Fluid Phase Equilibria</i> , 2009, 286, 170-174.	1.4	51
135	Hydrogen Bonding 12. A New QSAR for Upper Respiratory Tract Irritation by Airborne Chemicals in Mice. <i>QSAR and Combinatorial Science</i> , 1990, 9, 6-10.	1.4	49
136	Thermochemical behavior of dissolved carboxylic acid solutes: Solubilities of 3-methylbenzoic acid and 4-chlorobenzoic acid in organic solvents. <i>Canadian Journal of Chemistry</i> , 2003, 81, 1492-1501.	0.6	49
137	Thermochemical behavior of dissolved carboxylic acid solutes: part 4 - mathematical correlation of 4-nitrobenzoic acid solubilities with the abraham solvation parameter model. <i>Physics and Chemistry of Liquids</i> , 2004, 42, 339-347.	0.4	49
138	Correlation and prediction of solute transfer to chloroalkanes from both water and the gas phase. <i>Fluid Phase Equilibria</i> , 2009, 281, 144-162.	1.4	49
139	Partition Coefficients and Solubilities of Compounds in the Water-Ethanol Solvent System. <i>Journal of Solution Chemistry</i> , 2011, 40, 1279-1290.	0.6	49
140	Abraham model correlations for describing solute transfer into diisopropyl ether. <i>Physics and Chemistry of Liquids</i> , 2015, 53, 25-37.	0.4	49
141	Hydrogen bonding. Part 25. The solvation properties of methylene iodide. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1993, , 299-304.	0.9	48
142	Inverse Least-Squares Modeling of Vapor Descriptors Using Polymer-Coated Surface Acoustic Wave Sensor Array Responses. <i>Analytical Chemistry</i> , 2001, 73, 5247-5259.	3.2	48
143	The Permeation of Neutral Molecules, Ions, and Ionic Species Through Membranes: Brain Permeation as an Example. <i>Journal of Pharmaceutical Sciences</i> , 2011, 100, 1690-1701.	1.6	48
144	Substitution at saturated carbon. Part 26. A complete analysis of solvent effects on initial states and transition states for the solvolysis of the t-butyl halides in terms of G, H, and S using the unified method. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1988, , 1717.	0.9	47

#	ARTICLE	IF	CITATIONS
145	Hydrogen-bonding. Part 6. A thermodynamically-based scale of solute hydrogen-bond basicity. <i>Tetrahedron Letters</i> , 1989, 30, 2571-2574.	0.7	47
146	Thermochemical behavior of dissolved Carboxylic Acid solutes: Part 3 – Mathematical Correlation of 2-Methylbenzoic acid solubilities with the Abraham Solvation Parameter Model. <i>Physics and Chemistry of Liquids</i> , 2004, 42, 313-322.	0.4	47
147	Correlation of the toxicity of organic compounds to tadpoles using the Abraham model. <i>Science of the Total Environment</i> , 2006, 371, 99-109.	3.9	47
148	Mathematical correlation of 1-chloroanthraquinone solubilities in organic solvents with the Abraham solvation parameter model. <i>Physics and Chemistry of Liquids</i> , 2006, 44, 377-386.	0.4	47
149	Odor detection of single chemicals and binary mixtures. <i>Behavioural Brain Research</i> , 2005, 156, 115-123.	1.2	46
150	Development of Abraham model correlations for solvation characteristics of secondary and branched alcohols. <i>Fluid Phase Equilibria</i> , 2010, 288, 121-127.	1.4	46
151	Correlation and prediction of partition coefficients between the gas phase and water, and the solvents dodecane and undecane. <i>New Journal of Chemistry</i> , 2004, 28, 1538.	1.4	45
152	Correlations for describing gas-to-ionic liquid partitioning at 323K based on ion-specific equation coefficient and group contribution versions of the Abraham model. <i>Fluid Phase Equilibria</i> , 2011, 301, 257-266.	1.4	45
153	Abraham model correlations for solute transfer into tributyl phosphate from both water and the gas phase. <i>Physics and Chemistry of Liquids</i> , 2015, 53, 10-24.	0.4	45
154	Hydrogen bonding. <i>Journal of Chromatography A</i> , 1992, 627, 294-299.	1.8	44
155	Determination of solute descriptors of tripeptide derivatives based on high-throughput gradient high-performance liquid chromatography retention data. <i>Journal of Chromatography A</i> , 1998, 803, 51-60.	1.8	44
156	Solvation parameters for the 209 PCBs: calculation of physicochemical properties. <i>Journal of Environmental Monitoring</i> , 2005, 7, 295.	2.1	44
157	Enthalpy of solvation correlations for gaseous solutes dissolved in dimethyl sulfoxide and propylene carbonate based on the Abraham model. <i>Thermochimica Acta</i> , 2007, 459, 17-25.	1.2	44
158	Correlation and Prediction of Partition Coefficients From the Gas Phase and from Water to Alkan-1-ols. <i>Industrial & Engineering Chemistry Research</i> , 2008, 47, 3990-3995.	1.8	44
159	Solubility of crystalline nonelectrolyte solutes in organic solvents – Mathematical correlation of 2-methoxybenzoic acid and 4-methoxybenzoic acid solubilities with the Abraham solvation parameter model. <i>Canadian Journal of Chemistry</i> , 2004, 82, 1353-1360.	0.6	43
160	Chemical toxicity correlations for several protozoas, bacteria, and water fleas based on the Abraham solvation parameter model. <i>Journal of Environmental Engineering and Science</i> , 2007, 6, 165-174.	0.3	43
161	The effect of ionized species on microsomal binding. <i>European Journal of Medicinal Chemistry</i> , 2012, 47, 202-205.	2.6	43
162	Psychometric functions for the olfactory and trigeminal detectability of butyl acetate and toluene. <i>Journal of Applied Toxicology</i> , 2002, 22, 25-30.	1.4	42

#	ARTICLE	IF	CITATIONS
163	Enthalpy of Solvation Correlations for Gaseous Solutes Dissolved in Toluene and Carbon Tetrachloride Based on the Abraham Model. <i>Journal of Solution Chemistry</i> , 2007, 36, 947-966.	0.6	42
164	Thermochemical investigations of solute transfer into ionic liquid solvents: updated Abraham model equation coefficients for solute activity coefficient and partition coefficient predictions. <i>Physics and Chemistry of Liquids</i> , 2014, 52, 488-518.	0.4	42
165	Ion-specific equation coefficient version of the Abraham model for ionic liquid solvents: determination of coefficients for tributylethylphosphonium, 1-butyl-1-methylmorpholinium, 1-allyl-3-methylimidazolium and octyltriethylammonium cations. <i>Physics and Chemistry of Liquids</i> , 2017, 55, 358-385.	0.4	42
166	Solvation of gaseous non-electrolytes. <i>Faraday Discussions of the Chemical Society</i> , 1988, 85, 107.	2.2	41
167	Solute effects on reversed-phase thin-layer chromatography a linear free energy relationship analysis. <i>Journal of Chromatography A</i> , 1996, 749, 201-209.	1.8	41
168	Abraham model correlations for solute transfer into 2-methoxyethanol from water and from the gas phase. <i>Journal of Molecular Liquids</i> , 2015, 209, 738-744.	2.3	40
169	Predicting Abraham model solvent coefficients. <i>Chemistry Central Journal</i> , 2015, 9, 12.	2.6	40
170	Gas-solvent and water-solvent partition of trans -stilbene at 298 K. <i>Journal of Molecular Liquids</i> , 2017, 238, 58-61.	2.3	40
171	Mathematical correlation of salicylamide solubilities in organic solvents with the Abraham solvation parameter model. <i>Physics and Chemistry of Liquids</i> , 2007, 45, 389-398.	0.4	39
172	Concentration-detection functions for the odor of homologous n-acetate esters. <i>Physiology and Behavior</i> , 2008, 95, 658-667.	1.0	39
173	Structure-activity relationships on the odor detectability of homologous carboxylic acids by humans. <i>Experimental Brain Research</i> , 2010, 207, 75-84.	0.7	39
174	Abraham Model Correlations for Transfer of Neutral Molecules to Tetrahydrofuran and to 1,4-Dioxane, and for Transfer of Ions to Tetrahydrofuran. <i>Journal of Solution Chemistry</i> , 2011, 40, 2082-2094.	0.6	39
175	Descriptors for the Prediction of Partition Coefficients and Solubilities of Organophosphorus Compounds. <i>Separation Science and Technology</i> , 2013, 48, 884-897.	1.3	39
176	Development of Abraham model correlations for solute transfer into both 2-propoxyethanol and 2-isopropoxyethanol at 298.15 K. <i>Journal of Molecular Liquids</i> , 2015, 212, 833-840.	2.3	39
177	The temperature variation of the hydrophobic effect. <i>Journal of the Chemical Society Faraday Transactions I</i> , 1988, 84, 1985.	1.0	38
178	Partition of compounds from water and from air into the wet and dry monohalobenzenes. <i>New Journal of Chemistry</i> , 2009, 33, 1685.	1.4	38
179	Mathematical correlations for describing solute transfer into functionalized alkane solvents containing hydroxyl, ether, ester or ketone solvents. <i>Fluid Phase Equilibria</i> , 2010, 298, 48-53.	1.4	38
180	Abraham model correlations for describing solute transfer into anhydrous 1,2-propylene glycol for neutral and ionic species. <i>Physics and Chemistry of Liquids</i> , 2016, 54, 1-13.	0.4	38

#	ARTICLE	IF	CITATIONS
181	Quantitative relationship between rat intestinal absorption and Abraham descriptors. <i>European Journal of Medicinal Chemistry</i> , 2003, 38, 939-947.	2.6	37
182	Determination of Abraham model solute descriptors for benzoic acid based on measured solubility ratios. <i>Physics and Chemistry of Liquids</i> , 2012, 50, 254-265.	0.4	37
183	Abraham model correlations for describing solute transfer into 2-butoxyethanol from both water and the gas phase at 298K. <i>Journal of Molecular Liquids</i> , 2015, 209, 196-202.	2.3	37
184	Abraham model correlations for solute transfer into 2-ethoxyethanol from water and from the gas phase. <i>Journal of Molecular Liquids</i> , 2015, 208, 63-70.	2.3	37
185	Characterization of Some GLC Chiral Stationary Phases: A Linear Free Energy Analysis. <i>Analytical Chemistry</i> , 1997, 69, 613-617.	3.2	36
186	Characterisation of the water-isopropyl myristate system. <i>International Journal of Pharmaceutics</i> , 2005, 294, 121-128.	2.6	36
187	Gas to Olive Oil Partition Coefficients: A Linear Free Energy Analysis. <i>Journal of Chemical Information and Modeling</i> , 2006, 46, 1735-1741.	2.5	36
188	Mathematical correlation of phenothiazine solubilities in organic solvents with the Abraham solvation parameter model. <i>Physics and Chemistry of Liquids</i> , 2006, 44, 367-376.	0.4	36
189	Linear Free Energy Relationship Correlation of the Distribution of Solutes between Water and Sodium Dodecyl Sulfate (SDS) Micelles and between Gas and SDS Micelles. <i>Journal of Chemical Information and Modeling</i> , 2007, 47, 1808-1817.	2.5	36
190	Characterization of the retention behavior of organic and pharmaceutical drug molecules on an immobilized artificial membrane column with the Abraham model. <i>Journal of Chromatography A</i> , 2007, 1160, 235-245.	1.8	36
191	Enthalpy of solvation correlations for gaseous solutes dissolved in chloroform and 1,2-dichloroethane based on the Abraham model. <i>Fluid Phase Equilibria</i> , 2007, 258, 191-198.	1.4	36
192	Chemesthesis from volatile organic compounds: Psychophysical and neural responses. <i>Physiology and Behavior</i> , 2006, 88, 317-324.	1.0	35
193	Prediction of gas to water partition coefficients from 273 to 373K using predicted enthalpies and heat capacities of hydration. <i>Fluid Phase Equilibria</i> , 2007, 262, 97-110.	1.4	35
194	The biological and toxicological activity of gases and vapors. <i>Toxicology in Vitro</i> , 2010, 24, 357-362.	1.1	35
195	Equations for the Partition of Neutral Molecules, Ions and Ionic Species from Water to Water-Ethanol Mixtures. <i>Journal of Solution Chemistry</i> , 2012, 41, 730-740.	0.6	35
196	An equation for the prediction of human skin permeability of neutral molecules, ions and ionic species. <i>International Journal of Pharmaceutics</i> , 2017, 521, 259-266.	2.6	35
197	Solubility predictions for crystalline nonelectrolyte solutes dissolved in organic solvents based upon the Abraham general solvation model. <i>Canadian Journal of Chemistry</i> , 2001, 79, 1466-1476.	0.6	35
198	Calculations on ionic solvation. Part 4. Further calculations in solvation of gaseous univalent ions using one-layer and two-layer continuum models. <i>Journal of the Chemical Society Faraday Transactions I</i> , 1980, 76, 1219.	1.0	34

#	ARTICLE	IF	CITATIONS
199	Sensory Properties of Selected Terpenes: Thresholds for Odor, Nasal Pungency, Nasal Localization, and Eye Irritation. <i>Annals of the New York Academy of Sciences</i> , 1998, 855, 648-651.	1.8	34
200	Comparison of solubility of gases and vapours in wet and dry alcohols, especially octanol. <i>Journal of Physical Organic Chemistry</i> , 2008, 21, 823-832.	0.9	34
201	Enthalpy of solvation correlations for gaseous solutes dissolved in dibutyl ether and ethyl acetate. <i>Thermochimica Acta</i> , 2008, 470, 67-76.	1.2	34
202	Experimental and predicted solubilities of 3,4-dichlorobenzoic acid in select organic solvents and in binary aqueous-ethanol mixtures. <i>Physics and Chemistry of Liquids</i> , 2012, 50, 324-335.	0.4	34
203	On the solubility of quercetin. <i>Journal of Molecular Liquids</i> , 2014, 197, 157-159.	2.3	34
204	A simple method for estimating in vitro air-tissue and in vivo blood-tissue partition coefficients. <i>Chemosphere</i> , 2015, 120, 188-191.	4.2	34
205	Estimation of molecular linear free energy relationship descriptors. 4. Correlation and prediction of cell permeation. <i>Pharmaceutical Research</i> , 2000, 17, 1013-1018.	1.7	33
206	Comparison of Two Stimulus-delivery Systems for Measurement of Nasal Pungency Thresholds. <i>Chemical Senses</i> , 2000, 25, 285-291.	1.1	33
207	RAPID METHOD FOR ESTIMATING OCTANOL-WATER PARTITION COEFFICIENT (LOG POCT) FROM ISOCRATIC RP-HPLC AND A HYDROGEN BOND ACIDITY TERM (A). <i>Journal of Liquid Chromatography and Related Technologies</i> , 2001, 24, 635-649.	0.5	33
208	Thermochemical behavior of dissolved carboxylic acid solutes: part 5 – mathematical correlation of 3,5-dinitrobenzoic acid solubilities with the abraham solvation parameter model. <i>Physics and Chemistry of Liquids</i> , 2004, 42, 457-466.	0.4	33
209	Solubility of crystalline nonelectrolyte solutes in organic solvents: mathematical correlation of 3-nitrobenzoic acid solubilities with the Abraham general solvation model. <i>Journal of Molecular Liquids</i> , 2005, 116, 19-28.	2.3	33
210	Characterisation of the water/o-nitrophenyl octyl ether system in terms of the partition of nonelectrolytes and of ions. <i>Physical Chemistry Chemical Physics</i> , 2005, 7, 2418.	1.3	33
211	Correlation and prediction of partition coefficients for solute transfer to 1,2-dichloroethane from both water and from the gas phase. <i>Fluid Phase Equilibria</i> , 2008, 273, 78-86.	1.4	33
212	Determination of Abraham model solute descriptors for 2-ethylantraquinone based on measured solubility ratios. <i>Physics and Chemistry of Liquids</i> , 2011, 49, 355-365.	0.4	33
213	Hydrogen Bond Basicity of the Chlorogroup; Hexachlorocyclohexanes as Strong Hydrogen Bond Bases. <i>Journal of Organic Chemistry</i> , 2002, 67, 4782-4786.	1.7	32
214	Mathematical correlation of naproxen solubilities in organic solvents with the abraham solvation parameter model. <i>Physics and Chemistry of Liquids</i> , 2004, 42, 481-491.	0.4	32
215	Linear Free Energy Relationship Analysis of Retention Factors in Cerasome Electrokinetic Chromatography Intended for Predicting Drug Skin Permeation. <i>Journal of Pharmaceutical Sciences</i> , 2011, 100, 3105-3113.	1.6	32
216	Examination of hydrogen-bonding interactions between dissolved solutes and alkylbenzene solvents based on Abraham model correlations derived from measured enthalpies of solvation. <i>Thermochimica Acta</i> , 2014, 594, 68-79.	1.2	32

#	ARTICLE	IF	CITATIONS
217	Equations for the Partition of Neutral Molecules, Ions and Ionic Species from Water to Water-Methanol Mixtures. <i>Journal of Solution Chemistry</i> , 2016, 45, 861-874.	0.6	32
218	Development of Abraham model correlations for describing the transfer of molecular solutes into propanenitrile and butanenitrile from water and from the gas phase. <i>Physics and Chemistry of Liquids</i> , 2018, 56, 821-833.	0.4	32
219	Physicochemical analysis of the factors governing distribution of solutes between blood and brain. <i>Bioorganic and Medicinal Chemistry Letters</i> , 1994, 4, 2511-2516.	1.0	31
220	The determination of solvation descriptors for terpenes, and the prediction of nasal pungency thresholds. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1998, , 2405-2412.	0.9	31
221	Enthalpy of solvation correlations for organic solutes and gases dissolved in N,N-dimethylformamide and tert-butanol. <i>Journal of Molecular Liquids</i> , 2009, 144, 23-31.	2.3	31
222	Enthalpy of solvation correlations for organic solutes and gases dissolved in acetonitrile and acetone. <i>Thermochimica Acta</i> , 2009, 484, 65-69.	1.2	31
223	Dose-Response Functions for the Olfactory, Nasal Trigeminal, and Ocular Trigeminal Detectability of Airborne Chemicals by Humans. <i>Chemical Senses</i> , 2016, 41, 3-14.	1.1	31
224	Linear solvation energy relationships. Part 38. An analysis of the use of solvent parameters in the correlation of rate constants, with special reference to the solvolysis of t-butyl chloride. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1987, , 1097.	0.9	30
225	Solvation properties of refrigerants, and the estimation of their water-solvent and gas-solvent partitions. <i>Fluid Phase Equilibria</i> , 2001, 180, 41-58.	1.4	30
226	Dose-addition of individual odorants in the odor detection of binary mixtures. <i>Behavioural Brain Research</i> , 2003, 138, 95-105.	1.2	30
227	Olfactory psychometric functions for homologous 2-ketones. <i>Behavioural Brain Research</i> , 2009, 201, 207-215.	1.2	30
228	Determination of the Abraham model solute descriptors for 3,5-dinitro-2-methylbenzoic acid from measured solubility data in organic solvents. <i>Physics and Chemistry of Liquids</i> , 2011, 49, 821-829.	0.4	30
229	Is there an intramolecular hydrogen bond in 2-halophenols? A theoretical and spectroscopic investigation. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 25151-25159.	1.3	30
230	Updated Abraham model correlations for correlating solute transfer into dry butanone and dry cyclohexanone solvents. <i>Physics and Chemistry of Liquids</i> , 2018, 56, 571-583.	0.4	30
231	Free energies and entropies of transfer of ions from water to methanol, ethanol and 1-propanol. <i>Journal of the Chemical Society Faraday Transactions I</i> , 1978, 74, 2101.	1.0	29
232	Effect of Anesthetic Structure on Inhalation Anesthesia: Implications for the Mechanism. <i>Journal of Pharmaceutical Sciences</i> , 2008, 97, 2373-2384.	1.6	29
233	Abraham Model Expressions for Describing Water-to-Diethylene Glycol and Gas-to-Diethylene Glycol Solute Transfer Processes at 298.15 K. <i>Journal of Solution Chemistry</i> , 2017, 46, 331-351.	0.6	29
234	Solvation descriptors for the polychloronaphthalenes: estimation of some physicochemical properties. <i>Journal of Environmental Monitoring</i> , 2001, 3, 377-381.	2.1	28

#	ARTICLE	IF	CITATIONS
235	Blood or plasma to skin distribution of drugs: A linear free energy analysis. <i>International Journal of Pharmaceutics</i> , 2007, 329, 129-134.	2.6	28
236	Prediction of milk/plasma concentration ratios of drugs and environmental pollutants. <i>European Journal of Medicinal Chemistry</i> , 2009, 44, 2452-2458.	2.6	28
237	Analysis of immobilized artificial membrane retention factors for both neutral and ionic species. <i>Journal of Chromatography A</i> , 2013, 1298, 44-49.	1.8	28
238	Compilation and analysis of types and concentrations of airborne chemicals measured in various indoor and outdoor human environments. <i>Chemosphere</i> , 2015, 127, 70-86.	4.2	28
239	Effect of halogen substitution on the enthalpies of solvation and hydrogen bonding of organic solutes in chlorobenzene and 1,2-dichlorobenzene derived using multi-parameter correlations. <i>Thermochimica Acta</i> , 2015, 617, 8-20.	1.2	28
240	The transfer of neutral molecules, ions and ionic species from water to benzonitrile; comparison with nitrobenzene. <i>Thermochimica Acta</i> , 2011, 526, 22-28.	1.2	27
241	Enthalpy of solvation correlations for organic solutes and gases dissolved in dichloromethane and 1,4-dioxane. <i>Structural Chemistry</i> , 2013, 24, 1841-1853.	1.0	27
242	Correlation of the Solubility Behavior of Crystalline 1-Nitronaphthalene in Organic Solvents With the Abraham Solvation Parameter Model. <i>Journal of Solution Chemistry</i> , 2005, 34, 1121-1133.	0.6	26
243	Solute-solvent interactions in micellar liquid chromatography. <i>Journal of Chromatography A</i> , 2006, 1117, 47-55.	1.8	26
244	Development of correlations for describing solute transfer into acyclic alcohol solvents based on the Abraham model and fragment-specific equation coefficients. <i>Fluid Phase Equilibria</i> , 2010, 288, 139-144.	1.4	26
245	Abraham model correlations for describing solute transfer into ionic liquid solvents: calculation of ion-specific equation coefficients for the 4,5-dicyano-2-(trifluoromethyl)imidazolidine anion. <i>Physics and Chemistry of Liquids</i> , 2014, 52, 777-791.	0.4	26
246	Descriptors for ferrocene and some substituted ferrocenes. <i>Journal of Molecular Liquids</i> , 2017, 232, 325-331.	2.3	26
247	Characterization of hydrophilic interaction liquid chromatography retention by a linear free energy relationship. Comparison to reversed- and normal-phase retentions. <i>Analytica Chimica Acta</i> , 2019, 1092, 132-143.	2.6	26
248	The lipophilicity and hydrogen bond strength of pyridine-N-oxides and protonated pyridine-N-oxides. <i>New Journal of Chemistry</i> , 2011, 35, 930.	1.4	25
249	Correlation of the Solubilizing Abilities of Hexyl(trimethyl)ammonium bis((Trifluoromethyl)sulfonyl)imide, 1-Propyl-1-methylpiperidinium bis((Trifluoromethyl)sulfonyl)imide, and 1-Butyl-1-methyl-pyrrolidinium Thiocyanate. <i>Journal of Solution Chemistry</i> , 2011, 40, 2000-2022.	0.6	25
250	Enthalpy of solvation correlations for organic solutes and gases dissolved in 2-propanol, 2-butanol, 2-methyl-1-propanol and ethanol. <i>Thermochimica Acta</i> , 2011, 523, 214-220.	1.2	25
251	Human Intestinal Absorption of Neutral Molecules and Ionic Species. <i>Journal of Pharmaceutical Sciences</i> , 2014, 103, 1956-1966.	1.6	25
252	Deduction of Physicochemical Properties from Solubilities: 2,4-Dihydroxybenzophenone, Biotin, and Caprolactam as Examples. <i>Journal of Chemical & Engineering Data</i> , 2015, 60, 1440-1446.	1.0	25

#	ARTICLE	IF	CITATIONS
253	Computation of Abraham model solute descriptors for 3-methyl-4-nitrobenzoic acid from measured solubility data. <i>Physics and Chemistry of Liquids</i> , 2017, 55, 482-491.	0.4	25
254	Hydrogen-bonding. Part 4. An analysis of solute hydrogen-bond basicity, in terms of complexation constants (logK), using F1 and F2 factors, the principal components of different kinds of basicity. <i>Journal of Physical Organic Chemistry</i> , 1989, 2, 243-254.	0.9	24
255	Solvatochromic analysis of di-n-butyl ether/water partition coefficients as compared to other solvent systems. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1997, , 2639-2644.	0.9	24
256	Comments regarding "predicting the equilibrium partitioning of organic compounds using just one linear solvation energy relationship (LSER)". <i>Fluid Phase Equilibria</i> , 2005, 237, 224-226.	1.4	24
257	Correlating toxicities of organic compounds to select protozoa using the Abraham model. <i>Science of the Total Environment</i> , 2006, 369, 109-118.	3.9	24
258	Enthalpy of solvation correlations for organic solutes and gases dissolved in 1-propanol and tetrahydrofuran. <i>Thermochimica Acta</i> , 2011, 519, 103-113.	1.2	24
259	Correlation of the Solubilizing Abilities of 1-Butyl-1-methylpiperidinium Bis(trifluoromethylsulfonyl)imide and 1-Butyl-1-methylpyrrolidinium Tetracyanoborate. <i>Journal of Solution Chemistry</i> , 2012, 41, 1165-1184.	0.6	24
260	Linear free-energy relationships for water/hexadec-1-ene and water/deca-1,9-diene partitions, and for permeation through lipid bilayers; comparison of permeation systems. <i>New Journal of Chemistry</i> , 2012, 36, 1798.	1.4	24
261	Determination of solvation descriptors for terpene hydrocarbons from chromatographic measurements. <i>Journal of Chromatography A</i> , 2013, 1293, 133-141.	1.8	24
262	Abraham model linear free energy relationships for describing the partitioning and solubility behavior of nonelectrolyte organic solutes dissolved in pyridine at 298.15 K. <i>Fluid Phase Equilibria</i> , 2017, 431, 66-74.	1.4	24
263	Abraham model correlations for transfer of neutral molecules and ions to sulfolane. <i>Fluid Phase Equilibria</i> , 2011, 309, 30-35.	1.4	23
264	The prediction of blood-tissue partitions, water-skin partitions and skin permeation for agrochemicals. <i>Pest Management Science</i> , 2014, 70, 1130-1137.	1.7	23
265	Solubility of the pesticide diuron in organic nonelectrolyte solvents. Comparison of observed vs. predicted values based upon Mobile Order theory. <i>Canadian Journal of Chemistry</i> , 2000, 78, 184-190.	0.6	22
266	The lipophilicity of Sudan I and its tautomeric forms. <i>Physical Chemistry Chemical Physics</i> , 2002, 4, 5748-5752.	1.3	22
267	Hydrogen bonding, steric effects and thermodynamics of partitioning. <i>Journal of Physical Organic Chemistry</i> , 2006, 19, 242-248.	0.9	22
268	Abraham model correlations for solute partitioning into o-xylene, m-xylene and p-xylene from both water and the gas phase. <i>Fluid Phase Equilibria</i> , 2011, 308, 64-71.	1.4	22
269	The solubility of liquid and solid compounds in dry octan-1-ol. <i>Chemosphere</i> , 2014, 103, 26-34.	4.2	22
270	Determination of Abraham Model Correlations for Solute Transfer into Propyl Acetate Based on Experimental Activity Coefficient and Solubility Data. <i>Journal of Solution Chemistry</i> , 2018, 47, 634-653.	0.6	22

#	ARTICLE	IF	CITATIONS
271	Abraham model correlations for describing the thermodynamic properties of solute transfer into pentyl acetate based on headspace chromatographic and solubility measurements. <i>Journal of Chemical Thermodynamics</i> , 2018, 124, 133-140.	1.0	22
272	Solubility of the Pesticide Monuron in Organic Nonelectrolyte Solvents. Comparison of Observed Versus Predicted Values Based upon Mobile Order Theory. <i>Physics and Chemistry of Liquids</i> , 2002, 40, 255-268.	0.4	21
273	Water's Solvent Partition Coefficients and $\log P$ Values as Predictors for Blood-Brain Distribution; Application of the Akaike Information Criterion. <i>Journal of Pharmaceutical Sciences</i> , 2010, 99, 2492-2501.	1.6	21
274	Correlation of the Solubilizing Abilities of 1-Butyl-1-methylpyrrolidinium Tris(pentafluoroethyl)trifluorophosphate, 1-Butyl-1-methylpyrrolidinium Triflate and 1-Methoxyethyl-1-methylmorpholinium Tris(pentafluoroethyl)trifluorophosphate. <i>Journal of Solution Chemistry</i> , 2013, 42, 772-799.	0.6	21
275	The transfer of neutral molecules from water and from the gas phase to solvents acetophenone and aniline. <i>Journal of Molecular Liquids</i> , 2015, 212, 301-306.	2.3	21
276	Determination of Abraham model solute descriptors for the monomeric and dimeric forms of trans-cinnamic acid using measured solubilities from the Open Notebook Science Challenge. <i>Chemistry Central Journal</i> , 2015, 9, 11.	2.6	21
277	Determination of Abraham model solute descriptors for monomeric 3,4,5-trimethoxybenzoic acid from experimental solubility data in organic solvents measured at 298.2 K. <i>Physics and Chemistry of Liquids</i> , 2018, 56, 381-390.	0.4	21
278	Draize Eye Scores and Eye Irritation Thresholds in Man Can Be Combined into One QSARa. <i>Annals of the New York Academy of Sciences</i> , 1998, 855, 652-656.	1.8	20
279	A Theoretical Approach to the Ferguson Principle and its Use with Non-Reactive and Reactive Airborne Chemicals. <i>Basic and Clinical Pharmacology and Toxicology</i> , 1998, 83, 270-279.	0.0	20
280	The solvation properties of nitric oxide. <i>Perkin Transactions II RSC</i> , 2000, , 2067-2070.	1.1	20
281	Prediction and Mathematical Correlation of the Solubility of Fluorene in Alcohol Solvents Based upon the Abraham General Solvation Model. <i>Physics and Chemistry of Liquids</i> , 2002, 40, 581-591.	0.4	20
282	The determination of air/water partition coefficients for alkyl carboxylic acids by a new indirect method. <i>Journal of Environmental Monitoring</i> , 2003, 5, 747.	2.1	20
283	Determination of Abraham model solute descriptors for three dichloronitrobenzenes from measured solubilities in organic solvents. <i>Physics and Chemistry of Liquids</i> , 2015, 53, 163-173.	0.4	20
284	Determination of Abraham model solute descriptors for isophthalic acid from experimental solubility data in organic solvents at 298 K. <i>Physics and Chemistry of Liquids</i> , 2016, 54, 747-757.	0.4	20
285	An assessment of air quality reflecting the chemosensory irritation impact of mixtures of volatile organic compounds. <i>Environment International</i> , 2016, 86, 84-91.	4.8	20
286	A new method for the determination of Henry's law constants (air-water-partition coefficients). <i>Fluid Phase Equilibria</i> , 2019, 502, 112300.	1.4	20
287	Hydrogen-bonding. Part 22. Characterization of soybean oil and prediction of activity coefficients in soybean oil from inverse gas chromatographic data. <i>JAOCs, Journal of the American Oil Chemists' Society</i> , 1992, 69, 1236-1238.	0.8	19
288	Chemosensory additivity in trigeminal chemoreception as reflected by detection of mixtures. <i>Experimental Brain Research</i> , 2004, 158, 196-206.	0.7	19

#	ARTICLE	IF	CITATIONS
289	Thermochemical behavior of dissolved Carboxylic Acid solutes: Part 2 – Mathematical Correlation of Ketoprofen Solubilities with the Abraham General Solvation Model. <i>Physics and Chemistry of Liquids</i> , 2004, 42, 305-312.	0.4	19
290	Mathematical correlation of 4-aminobenzoic acid solubilities in organic solvents with the abraham solvation parameter model. <i>Physics and Chemistry of Liquids</i> , 2004, 42, 633-641.	0.4	19
291	Chemical Boundaries for Detection of Eye Irritation in Humans from Homologous Vapors. <i>Toxicological Sciences</i> , 2006, 91, 600-609.	1.4	19
292	Physicochemical and biochemical properties for the dialkyl phthalates. <i>Chemosphere</i> , 2015, 119, 871-880.	4.2	19
293	Studies on the hydrogen bond acidity, and other descriptors and properties for hydroxyflavones and hydroxyisoflavones. <i>Journal of Molecular Liquids</i> , 2015, 208, 363-372.	2.3	19
294	Modeling Aquatic Toxicity through Chromatographic Systems. <i>Analytical Chemistry</i> , 2017, 89, 7996-8003.	3.2	19
295	Development of Abraham model correlations for solute transfer into 2-ethyl-1-hexanol from both water and the gas phase based on measured solubility ratios. <i>Physics and Chemistry of Liquids</i> , 2020, 58, 202-213.	0.4	19
296	Molecular restrictions for human eye irritation by chemical vapors. <i>Toxicology and Applied Pharmacology</i> , 2005, 207, 232-243.	1.3	18
297	Descriptors for the Prediction of Partition Coefficients of 8-Hydroxyquinoline and its Derivatives. <i>Separation Science and Technology</i> , 2014, 49, 2135-2141.	1.3	18
298	Development of Abraham model expressions for predicting the enthalpies of solvation of solutes dissolved in acetic acid. <i>Physics and Chemistry of Liquids</i> , 2016, 54, 141-154.	0.4	18
299	A simple and facile NMR method for the determination of hydrogen bonding by amide N-H protons in protein models and other compounds. <i>New Journal of Chemistry</i> , 2017, 41, 6064-6066.	1.4	18
300	Development of Abraham model expressions for predicting the standard molar enthalpies of vaporization of organic compounds at 298.15 K. <i>Thermochimica Acta</i> , 2019, 681, 178372.	1.2	18
301	Sensory irritation mechanisms investigated from model compounds: trifluoroethanol, hexafluoroisopropanol and methyl hexafluoroisopropyl ether. <i>Archives of Toxicology</i> , 1996, 70, 319-328.	1.9	17
302	Mathematical correlation of 1,2,4,5-tetramethylbenzene solubilities in organic solvents with the Abraham solvation parameter model. <i>Physics and Chemistry of Liquids</i> , 2006, 44, 173-182.	0.4	17
303	Equations for water-triolein partition coefficients for neutral species; comparison with other water-solvent partitions, and environmental and toxicological processes. <i>Chemosphere</i> , 2016, 154, 48-54.	4.2	17
304	Determination of the solubilising character of 2-methoxyethyl-(dimethyl)ethylammonium tris(pentafluoroethyl)trifluorophosphate based on the Abraham solvation parameter model. <i>Physics and Chemistry of Liquids</i> , 2016, 54, 110-126.	0.4	17
305	Abraham model expressions for describing water-to-organic solvent and gas-to-organic solvent partition coefficients for solute transfer into anhydrous poly(ethylene glycol) dialkyl ether solvents at 298.15 K. <i>Physics and Chemistry of Liquids</i> , 2017, 55, 347-357.	0.4	17
306	Determination of Abraham model solute descriptors for o-acetoacetanilide based on experimental solubility data in organic mono-solvents. <i>Physics and Chemistry of Liquids</i> , 2019, 57, 528-535.	0.4	17

#	ARTICLE	IF	CITATIONS
307	Estimation of enthalpies of sublimation of organic, organometallic and inorganic compounds. Fluid Phase Equilibria, 2020, 515, 112575.	1.4	17
308	Solvation descriptors for N-nitrosodialkylamines; calculation of some of their properties of environmental significance. Journal of Environmental Monitoring, 2002, 4, 743-746.	2.1	16
309	Henry's Law constants or air to water partition coefficients for 1,3,5-triazines by an LFER method. Journal of Environmental Monitoring, 2007, 9, 234-239.	2.1	16
310	The hydrogen bond acidity and other descriptors for oximes. New Journal of Chemistry, 2009, 33, 76-81.	1.4	16
311	The hydrogen bond properties of water from 273 K to 573 K; equations for the prediction of gas-water partition coefficients. Physical Chemistry Chemical Physics, 2012, 14, 7433.	1.3	16
312	Abraham model enthalpy of solvation correlations for solutes dissolved in dimethyl carbonate and diethyl carbonate. Physics and Chemistry of Liquids, 2015, 53, 732-747.	0.4	16
313	Updated Abraham model correlations for enthalpies of solvation of organic solutes dissolved in benzene and acetonitrile. Physics and Chemistry of Liquids, 2019, 57, 84-99.	0.4	16
314	Estimation of vapor pressures of liquid and solid organic and organometallic compounds at 298.15 K. Fluid Phase Equilibria, 2020, 519, 112595.	1.4	16
315	100 years of chromatography or is it 171?. Journal of Chromatography A, 2004, 1061, 113-114.	1.8	15
316	SOLUBILITY BEHAVIOR OF CRYSTALLINE POLYCYCLIC AROMATIC HYDROCARBONS (PAHs): PREDICTION OF FLUORENE SOLUBILITIES IN ORGANIC SOLVENTS WITH THE ABRAHAM SOLVATION PARAMETER MODEL. Polycyclic Aromatic Compounds, 2005, 25, 313-326.	1.4	15
317	Comment on "Systematic Investigation of the Sorption Properties of Polyurethane Foams for Organic Vapors". Analytical Chemistry, 2007, 79, 6891-6893.	3.2	15
318	LFER correlations for the solubilising characterisation of room temperature ionic liquids containing trifluoromethanesulfonate and trifluoroacetate anions. Physics and Chemistry of Liquids, 2008, 46, 631-642.	0.4	15
319	Linear free energy relationship (LFER) correlations for the solubilising characterisation of room temperature ionic liquids containing triethylsulphonium and 1-butyl-1-methylpyrrolidinium cations. Physics and Chemistry of Liquids, 2010, 48, 385-393.	0.4	15
320	Abraham model enthalpy of solvation correlations for solutes dissolved in 1-alkanol solvents (C_4 to C_6). Physics and Chemistry of Liquids, 2015, 53, 638-659.	0.4	15
321	Quantifying solvent effects through QSPR: A new look over different model equations. Journal of Molecular Liquids, 2019, 291, 111244.	2.3	15
322	Abraham model correlations for describing solute transfer into 4-methyl-2-pentanol from both water and the gas phase. Journal of Molecular Liquids, 2019, 278, 335-341.	2.3	15
323	Solvation Descriptors for Zwitterionic $\hat{\pm}$ -Aminoacids; Estimation of Water's Solvent Partition Coefficients, Solubilities, and Hydrogen-Bond Acidity and Hydrogen-Bond Basicity. ACS Omega, 2019, 4, 2883-2892.	1.6	15
324	Descriptors for terpene esters from chromatographic and partition measurements: Estimation of human odor detection thresholds. Journal of Chromatography A, 2020, 1609, 460428.	1.8	15

#	ARTICLE	IF	CITATIONS
325	Linear free energy relationship models for the retention of partially ionized acid-base compounds in reversed-phase liquid chromatography. <i>Journal of Chromatography A</i> , 2021, 1635, 461720.	1.8	15
326	Hydrogen bonding part 46: a review of the correlation and prediction of transport properties by an lfer method: physicochemical properties, brain penetration and skin permeability. , 1999, 55, 78.		15
327	Concentration-detection functions for eye irritation evoked by homologous n-alcohols and acetates approaching a cut-off point. <i>Experimental Brain Research</i> , 2007, 182, 71-79.	0.7	14
328	Solvation parameters for mercury and mercury(ii) compounds: calculation of properties of environmental interest. <i>Journal of Environmental Monitoring</i> , 2008, 10, 435.	2.1	14
329	Comparison of lipid membrane water partitioning with various organic solvent water partitions of neutral species and ionic species: Uniqueness of cerasome as a model for the stratum corneum in partition processes. <i>International Journal of Pharmaceutics</i> , 2015, 494, 1-8.	2.6	14
330	Solubility of sorbic acid in organic mono-solvents: calculation of Abraham model solute descriptors from measured solubility data. <i>Physics and Chemistry of Liquids</i> , 2017, 55, 650-658.	0.4	14
331	Structural properties governing drug-plasma protein binding determined by high-performance liquid chromatography method. <i>Journal of Pharmaceutical and Biomedical Analysis</i> , 2018, 149, 16-21.	1.4	14
332	Abraham model solute descriptors reveal strong intramolecular hydrogen bonding in 1,4-dihydroxyanthraquinone and 1,8-dihydroxyanthraquinone. <i>Physics and Chemistry of Liquids</i> , 2018, 56, 416-420.	0.4	14
333	A cut-off in ocular chemesthesis from vapors of homologous alkylbenzenes and 2-ketones as revealed by concentration-detection functions. <i>Toxicology and Applied Pharmacology</i> , 2008, 230, 298-303.	1.3	13
334	Abraham model correlations for estimating solute transfer of neutral molecules into anhydrous acetic acid from water and from the gas phase. <i>Journal of Molecular Liquids</i> , 2015, 212, 16-22.	2.3	13
335	Development of Abraham model correlations for predicting enthalpies of solvation of nonionic solutes dissolved in formamide. <i>Physics and Chemistry of Liquids</i> , 2016, 54, 313-324.	0.4	13
336	Abraham Model Correlations for Triethylene Glycol Solvent Derived from Infinite Dilution Activity Coefficient, Partition Coefficient and Solubility Data Measured at 298.15ÅK. <i>Journal of Solution Chemistry</i> , 2017, 46, 2249-2267.	0.6	13
337	Determination of Abraham model solute descriptors for 4- <i>tert</i> -butylbenzoic acid from experimental solubility data in organic mono-solvents. <i>Physics and Chemistry of Liquids</i> , 2019, 57, 445-452.	0.4	13
338	Descriptors for vitamin K3 (menadione); calculation of biological and physicochemical properties. <i>Journal of Molecular Liquids</i> , 2021, 330, 115707.	2.3	13
339	Application of reaction field theory to the calculation of solvent effects on the t-butyl chloride solvolysis, and on the Me ₄ N ⁺ Cl ⁻ ion pair. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1974, , 47.	0.9	12
340	Hydrogen Bonding between Solutes in Solvents Octan-1-ol and Water. <i>Journal of Organic Chemistry</i> , 2010, 75, 7651-7658.	1.7	12
341	The correlation and prediction of infinite dilution activity coefficients of compounds in water at 298.15ÅK. <i>Fluid Phase Equilibria</i> , 2017, 449, 117-129.	1.4	12
342	Abraham model correlations for describing the solubilising character of 3-Methoxy-1-butanol and 1- <i>tert</i> -Butoxy-2-propanol solvents. <i>Physics and Chemistry of Liquids</i> , 2019, 57, 163-173.	0.4	12

#	ARTICLE	IF	CITATIONS
343	Abraham model correlations for solute transfer into 2-methyl-2-butanol based on measured activity coefficient and solubility data at 298.15 K. <i>Journal of Molecular Liquids</i> , 2019, 293, 111454.	2.3	12
344	Development of Abraham model correlations for describing solute transfer into 2-methyl-1-butanol from both water and the gas phase from experimental solubility data of crystalline organic compounds. <i>Physics and Chemistry of Liquids</i> , 2020, 58, 623-635.	0.4	12
345	Abraham model correlations for describing dissolution of organic solutes and inorganic gases in dimethyl carbonate. <i>Physics and Chemistry of Liquids</i> , 2021, 59, 181-195.	0.4	12
346	Hydrogen bonding part 46: A review of the correlation and prediction of transport properties by an LFER method: physicochemical properties, brain penetration and skin permeability??. <i>Pest Management Science</i> , 1999, 55, 78-88.	0.7	12
347	Heats of solution of 1:1 electrolytes in 1,2- and 1,1-dichloroethane and derived enthalpies and entropies of transfer of electrolytes from water to these solvents. <i>Journal of Solution Chemistry</i> , 1976, 5, 529-542.	0.6	11
348	Hydrogen bond descriptors and other properties of ion pairs. <i>New Journal of Chemistry</i> , 2011, 35, 1740.	1.4	11
349	On the solubility of nicotinic acid and isonicotinic acid in water and organic solvents. <i>Journal of Chemical Thermodynamics</i> , 2013, 61, 74-78.	1.0	11
350	Abraham model linear free energy relationships as a means of extending solubility studies to include the estimation of solute solubilities in additional organic solvents. <i>Journal of Chemical Thermodynamics</i> , 2016, 102, 392-397.	1.0	11
351	Abraham model ion-specific equation coefficients for the 1-butyl-2,3-dimethylimidazolium and 4-cyano-1-butylpyridinium cations calculated from measured gas-to-liquid partition coefficient data. <i>Physics and Chemistry of Liquids</i> , 2017, 55, 218-237.	0.4	11
352	Development of Abraham model IL-specific correlations for N-triethyl(octyl)ammonium bis(fluorosulfonyl)imide and 1-butyl-3-methylpyrrolidinium bis(fluorosulfonyl)imide. <i>Physics and Chemistry of Liquids</i> , 2019, 57, 733-745.	0.4	11
353	Abraham model correlations for enthalpies of solvation of organic solutes dissolved in methyl acetate and octane. <i>Physics and Chemistry of Liquids</i> , 2020, 58, 18-30.	0.4	11
354	Abraham model correlations for enthalpies of solvation of organic solutes dissolved in N,N-Dimethylacetamide, 2-butanone and tetrahydrofuran (UPDATED) at 298.15 K. <i>Physics and Chemistry of Liquids</i> , 2020, 58, 675-692.	0.4	11
355	Thermodynamic parameters for the transfer of ions from water to propylene carbonate. <i>Monatshefte für Chemie</i> , 1979, 110, 517-524.	0.9	10
356	Characterisation of room temperature ionic liquid chromatographic stationary phases by combining experimental retention factor and partition coefficient data into a single model. <i>Physics and Chemistry of Liquids</i> , 2009, 47, 74-83.	0.4	10
357	Prediction of convulsant activity of gases and vapors. <i>European Journal of Medicinal Chemistry</i> , 2009, 44, 885-890.	2.6	10
358	The factors that influence solubility in perfluoroalkane solvents. <i>Fluid Phase Equilibria</i> , 2016, 421, 59-66.	1.4	10
359	Determination of molar refractions and Abraham descriptors for tris(acetylacetonato)chromium(III), tris(acetylacetonato)iron(III) and tris(acetylacetonato)cobalt(III). <i>New Journal of Chemistry</i> , 2017, 41, 14259-14265.	1.4	10
360	Estimation of heat capacities of gases, liquids and solids, and heat capacities of vaporization and of sublimation of organic chemicals at 298.15 K. <i>Journal of Molecular Liquids</i> , 2020, 317, 113969.	2.3	10

#	ARTICLE	IF	CITATIONS
361	Solute-solvent interactions in chemical and biological systems. IV. Correlations of ΔG , ΔH and ΔS of transfer of aliphatic and aromatic solutes from 2,2,4-trimethylpentane to aqueous solution. <i>Journal of Physical Organic Chemistry</i> , 1989, 2, 559-564.	0.9	9
362	Self-Assembly Does Not Account for the Hydrophobic Effect. <i>Journal of the American Chemical Society</i> , 2002, 124, 7853-7856.	6.6	9
363	Comments on "Solvation Parameters. 2. A Simplified Molecular Topology To Generate Easily Optimized Values". <i>Journal of Chemical Information and Modeling</i> , 2006, 46, 1879-1881.	2.5	9
364	A linear free energy analysis of PAMPA models for biological systems. <i>International Journal of Pharmaceutics</i> , 2015, 496, 717-722.	2.6	9
365	Development of Abraham model correlations for enthalpies of solvation of organic solutes dissolved in 1,3-dioxolane. <i>Physics and Chemistry of Liquids</i> , 2016, 54, 786-796.	0.4	9
366	Determination of Abraham model solute descriptors for 2-methyl-3-nitrobenzoic acid from measured solubility data in alcohol, alkyl ether, alkyl acetate and 2-alkoxyalcohol mono-solvents. <i>Physics and Chemistry of Liquids</i> , 0, , 1-9.	0.4	9
367	Revisiting blood-brain barrier: A chromatographic approach. <i>Journal of Pharmaceutical and Biomedical Analysis</i> , 2017, 145, 98-109.	1.4	9
368	Abraham model correlations for solute transfer into benzyl alcohol from both water and the gas phase. <i>Physics and Chemistry of Liquids</i> , 2020, 58, 116-126.	0.4	9
369	Abraham model correlation for direct water-to-2,2,5,5-tetramethyloxolane solute transfer partitioning process revisited. <i>Physics and Chemistry of Liquids</i> , 2020, 58, 833-838.	0.4	9
370	Equations for the Correlation and Prediction of Partition Coefficients of Neutral Molecules and Ionic Species in the Water-Isopropanol Solvent System. <i>Journal of Solution Chemistry</i> , 2021, 50, 458-472.	0.6	9
371	Substitution at saturated carbon. Part XIX. The effect of alcohols and water on the free energy of solutes and on the free energy of transition states in S _N and S _E reactions. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1975, , 1856.	0.9	8
372	The Correlation and Prediction of Butane/Water and Gas/Butane Partition Coefficients. <i>Canadian Journal of Chemical Engineering</i> , 2008, 83, 362-365.	0.9	8
373	Descriptors for Pentane-2,4-dione and Its Derivatives. <i>Journal of Solution Chemistry</i> , 2017, 46, 1625-1638.	0.6	8
374	Descriptors for the α,ω -dicarboxylic acids from oxalic acid to sebacic acid. <i>Fluid Phase Equilibria</i> , 2018, 467, 17-24.	1.4	8
375	Descriptors for Cyclooctasulfur: Estimation of Water-Solvent Partition Coefficients, Solubilities in Solvents, and Physicochemical Properties. <i>ACS Omega</i> , 2018, 3, 5516-5521.	1.6	8
376	Linear free energy relationship analysis of permeability across polydimethylsiloxane (PDMS) membranes and comparison with human skin permeation in vitro. <i>European Journal of Pharmaceutical Sciences</i> , 2018, 123, 524-530.	1.9	8
377	Limiting Diffusion Coefficients for Ions and Nonelectrolytes in Solvents Water, Methanol, Ethanol, Propan-1-ol, Butan-1-ol, Octan-1-ol, Propanone and Acetonitrile at 298 K, Analyzed Using Abraham Descriptors. <i>Journal of Solution Chemistry</i> , 2019, 48, 748-757.	0.6	8
378	Applications of Abraham solvation parameter model: estimation of the lethal median molar concentration of the antiepileptic drug levetiracetam towards aquatic organisms from measured solubility data. <i>Physics and Chemistry of Liquids</i> , 2020, 58, 302-308.	0.4	8

#	ARTICLE	IF	CITATIONS
379	Abraham model correlations for describing solute transfer processes into diethyl carbonate. <i>Physics and Chemistry of Liquids</i> , 2021, 59, 26-39.	0.4	8
380	Solvent and gas-phase effects on the equilibrium between configurational isomers of some 4-t-butylcyclohexanes. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1982, , 1503.	0.9	7
381	Partition between phases of a solute that exists as two interconverting species. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1995, , 1839.	0.9	7
382	Hydrogen bond and other descriptors for thalidomide and its N-alkyl analogs; prediction of physicochemical and biological properties. <i>European Journal of Pharmaceutical Sciences</i> , 2004, 21, 465-469.	1.9	7
383	Linear free energy relationship correlations for the solubilising characterisation of room temperature ionic liquids containing 1-hexyloxymethyl-3-methylimidazolium and 1,3-dihexyloxymethylimidazolium cations. <i>Physics and Chemistry of Liquids</i> , 2010, 48, 394-402.	0.4	7
384	Abraham model correlations describing the solubilising ability of peanut oil. <i>Physics and Chemistry of Liquids</i> , 2014, 52, 792-803.	0.4	7
385	Analysis of the solubility of betaine: calculation of descriptors and physicochemical properties. <i>Fluid Phase Equilibria</i> , 2015, 387, 1-4.	1.4	7
386	Solvation descriptors for porphyrins (porphines). <i>New Journal of Chemistry</i> , 2016, 40, 9945-9950.	1.4	7
387	Partition of Neutral Molecules and Ions from Water to o-Nitrophenyl Octyl Ether and of Neutral Molecules from the Gas Phase to o-Nitrophenyl Octyl Ether. <i>Journal of Solution Chemistry</i> , 2018, 47, 293-307.	0.6	7
388	Determination of Abraham model correlations for describing solute transfer into the methyl butyrate mono-solvent at 298 K. <i>Physics and Chemistry of Liquids</i> , 2020, 58, 792-802.	0.4	7
389	Abraham solvation parameter model: calculation of ion-specific equation coefficients for the N-Ethyl-N-methylmorpholinium and N-Octyl-N-methylmorpholinium cations. <i>Physics and Chemistry of Liquids</i> , 2021, 59, 575-584.	0.4	7
390	The estimation of physicochemical properties of methyl and other alkyl naphthalenes. <i>Journal of Environmental Monitoring</i> , 2005, 7, 445.	2.1	6
391	Gas-liquid and water-liquid partition coefficients of the tetraphenyl compounds of group (IV). <i>New Journal of Chemistry</i> , 2012, 36, 626-631.	1.4	6
392	Comment on "Measurement and Correlation of the Solubility of <i>p</i> -Coumaric Acid in Nine Pure and Water + Ethanol Mixed Solvents at Temperatures from 293.15 to 333.15 K". <i>Journal of Chemical & Engineering Data</i> , 2017, 62, 578-583.	1.0	6
393	Calculation of the Abraham model solute descriptors for the pharmaceutical compound acipimox based on experimental solubility data. <i>Physics and Chemistry of Liquids</i> , 2019, 57, 382-387.	0.4	6
394	Determination of Abraham model solute descriptors for xanthone based on experimental solubility measurements at 298.2 K. <i>Physics and Chemistry of Liquids</i> , 2020, 58, 214-221.	0.4	6
395	Heterolytic cleavage of main group metal-carbon bonds. , 0, , 25-149.		6
396	Design and Information Content of Arrays of Sorption-Based Vapor Sensors Using Solubility Interactions and Linear Solvation Energy Relationships. , 2009, , 193-218.		6

#	ARTICLE	IF	CITATIONS
397	Descriptors for some compounds with pharmacological activity; calculation of properties. International Journal of Pharmaceutics, 2022, 617, 121597.	2.6	6
398	Response to "A critique of Abraham and Acree's correlation for deca-1,9-diene" water partition coefficients. New Journal of Chemistry, 2013, 37, 882.	1.4	5
399	The correlation and prediction of the temperature variation of infinite dilution activity coefficients of compounds in water. Fluid Phase Equilibria, 2018, 455, 1-5.	1.4	5
400	Descriptors for the hydrogen halides, their solution properties and hydrogen-bonding acidity and basicity: Comparison of the latter with gas phase data. Journal of Molecular Liquids, 2019, 275, 667-673.	2.3	5
401	Solubility of 4-methyl-3-nitrobenzoic acid in organic mono-solvents: calculation of Abraham model solute descriptors. Physics and Chemistry of Liquids, 2020, 58, 782-791.	0.4	5
402	The partition of organic compounds from water into the methyl isobutyl ketone extraction solvent with updated Abraham model equation. Physics and Chemistry of Liquids, 2021, 59, 431-441.	0.4	5
403	Descriptors for High-Energy Nitro Compounds; Estimation of Thermodynamic, Physicochemical and Environmental Properties. Propellants, Explosives, Pyrotechnics, 2021, 46, 267-279.	1.0	5
404	Descriptors for Edaravone; studies on its structure, and prediction of properties. Journal of Molecular Liquids, 2021, 332, 115821.	2.3	5
405	Summation solute hydrogen bonding acidity values for hydroxyl substituted flavones determined by NMR spectroscopy. Natural Product Communications, 2013, 8, 85-98.	0.2	5
406	Updated Abraham model correlations to describe enthalpies of solvation of solutes dissolved in heptane, cyclohexane and N,N-dimethylformamide. Physics and Chemistry of Liquids, 2021, 59, 442-453.	0.4	4
407	Properties of the <i>tert</i> -butyl halide solvolysis transition states. Physical Chemistry Chemical Physics, 2021, 23, 3311-3320.	1.3	4
408	Abraham model correlations for describing the partition of organic compounds from water into the methyl ethyl ketone extraction solvent. Physics and Chemistry of Liquids, 2022, 60, 47-58.	0.4	4
409	Professor Robert W. Taft. Journal of Physical Organic Chemistry, 1994, 7, 655-656.	0.9	3
410	Determination of partition coefficients of refrigerants by gas liquid chromatographic headspace analysis. Journal of Chromatography A, 2012, 1265, 144-148.	1.8	3
411	Commentary on "Measurement and Correlation of the Solubility of Telmisartan (Form A) in Nine Different Solvents from 277.85 to 338.35 K". Journal of Solution Chemistry, 2016, 45, 1902-1905.	0.6	3
412	Illustration of the calculation of solute descriptors for maltol from published solubility data. Physics and Chemistry of Liquids, 2018, 56, 403-409.	0.4	3
413	The assessment of intramolecular hydrogen bonding in ortho-substituted anilines by an NMR method. Journal of Molecular Liquids, 2020, 315, 113730.	2.3	3
414	Descriptors for fluorotelomere alcohols. Calculation of physicochemical properties. Physics and Chemistry of Liquids, 0, , 1-6.	0.4	3

#	ARTICLE	IF	CITATIONS
415	Abraham Model Descriptors for Melatonin; Prediction of Solution, Biological and Thermodynamic Properties. <i>Journal of Solution Chemistry</i> , 0, , 1.	0.6	3
416	Solvent effects on the cis- and trans-4-t-butylcyclohexanol equilibrium by a new method, using gas-chromatographic head-space analysis. <i>Journal of the Chemical Society Chemical Communications</i> , 1980, , 312.	2.0	2
417	A critique of a thermodynamic description of hydrophobic aggregation in aqueous solution. <i>Thermochimica Acta</i> , 2003, 403, 219-222.	1.2	2
418	Comments concerning "characterizations for vinylimidazolium based ionic liquid polymer stationary phases for capillary gas chromatography". <i>Chromatographia</i> , 2008, 68, 1075-1078.	0.7	2
419	Summation Solute Hydrogen Bonding Acidity Values for Hydroxyl Substituted Flavones Determined by NMR Spectroscopy. <i>Natural Product Communications</i> , 2013, 8, 1934578X1300800.	0.2	2
420	Abraham solvation parameter model: updated correlations for describing solute partitioning into plant cuticles from water and from air. <i>Physics and Chemistry of Liquids</i> , 2020, , 1-17.	0.4	2
421	Comments on "Classification of biphasic solvent systems according to Abraham descriptors for countercurrent chromatography". <i>Journal of Chromatography A</i> , 2020, 1618, 460889.	1.8	2
422	Descriptors for adamantane and some of its derivatives. <i>Journal of Molecular Liquids</i> , 2021, 325, 114894.	2.3	2
423	Abraham model solute descriptors: effect of structural features on the calculated numerical descriptor values for vanillin and select derivatives. <i>Physics and Chemistry of Liquids</i> , 2023, 61, 1-13.	0.4	2
424	Substitution at saturated carbon. Part 25. The iododemallation of benzyltrialkyltins in methanol. <i>Bulletin Des Sociétés Chimiques Belges</i> , 1980, 89, 819-829.	0.0	1
425	Comment on "Structural Determinants of Drug Partitioning in Surrogates of Phosphatidylcholine Bilayer Strata". <i>Molecular Pharmaceutics</i> , 2015, 12, 1328-1329.	2.3	1
426	Comments concerning "A possible simplification of the Goss-modified Abraham solvation equation". <i>Chemosphere</i> , 2015, 138, 1058-1061.	4.2	1
427	Comment on "Solubility Measurement and Thermodynamic Modeling of		