

# Feng Luan

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/7302340/publications.pdf>

Version: 2024-02-01

133  
papers

3,680  
citations

109321

35  
h-index

161849

54  
g-index

136  
all docs

136  
docs citations

136  
times ranked

4011  
citing authors

#	ARTICLE	IF	CITATIONS
1	Selective detection of enrofloxacin in biological and environmental samples using a molecularly imprinted electrochemiluminescence sensor based on functionalized copper nanoclusters. <i>Talanta</i> , 2022, 236, 122835.	5.5	38
2	Encapsulating Ru(bpy) <sub>3</sub> <sup>2+</sup> in an infinite coordination polymer network: Towards a solid-state electrochemiluminescence sensing platform for histamine to evaluate fish product quality. <i>Food Chemistry</i> , 2022, 368, 130852.	8.2	20
3	An ultrasensitive electrochemiluminescence biosensor for the detection of total bacterial count in environmental and biological samples based on a novel sulfur quantum dot luminophore. <i>Analyst, The</i> , 2022, 147, 1716-1721.	3.5	4
4	A portable visual coffee ring based on carbon dot sensitized lanthanide complex coordination to detect bisphenol A in water. <i>RSC Advances</i> , 2022, 12, 7306-7312.	3.6	8
5	Electrochemiluminescent determination of CYFRA21-1 serum levels using Ti-Fe@O nanotubes immunoassay. <i>Mikrochimica Acta</i> , 2022, 189, 136.	5.0	1
6	Selective Detection of Alkaline Phosphatase Activity in Environmental Water Samples by Copper Nanoclusters Doped Lanthanide Coordination Polymer Nanocomposites as the Ratiometric Fluorescent Probe. <i>Biosensors</i> , 2022, 12, 372.	4.7	5
7	An Electrochemical Sensor Based on Gold and Bismuth Bimetallic Nanoparticles Decorated L-Cysteine Functionalized Graphene Oxide Nanocomposites for Sensitive Detection of Iron Ions in Water Samples. <i>Nanomaterials</i> , 2021, 11, 2386.	4.1	8
8	Ru(bpy) <sub>3</sub> <sup>2+</sup> encapsulated cyclodextrin based metal organic framework with improved biocompatibility for sensitive electrochemiluminescence detection of CYFRA21-1 in cell. <i>Biosensors and Bioelectronics</i> , 2021, 190, 113371.	10.1	38
9	Binding specificities of estrogen receptor with perfluorinated compounds: A cross species comparison. <i>Environment International</i> , 2020, 134, 105284.	10.0	31
10	Functionalized Copper Nanoclusters-Based Fluorescent Probe with Aggregation-Induced Emission Property for Selective Detection of Sulfide Ions in Food Additives. <i>Journal of Agricultural and Food Chemistry</i> , 2020, 68, 11301-11308.	5.2	46
11	Interaction of Coumarin Phytoestrogens with ER $\alpha$ and ER $\beta$ : A Molecular Dynamics Simulation Study. <i>Molecules</i> , 2020, 25, 1165.	3.8	11
12	Facile synthesis of a cyclodextrin-metal organic framework decorated with Ketjen Black and platinum nanoparticles and its application in the electrochemical detection of ofloxacin. <i>Analyst, The</i> , 2020, 145, 1943-1949.	3.5	32
13	The Interaction of Isoflavone Phytoestrogens with ER $\alpha$ and ER $\beta$ by Molecular Docking and Molecular Dynamics Simulations. <i>Current Computer-Aided Drug Design</i> , 2020, 16, 655-665.	1.2	5
14	An electrochemiluminescence sensor for the detection of prostate protein antigen based on the graphene quantum dots infilled TiO <sub>2</sub> nanotube arrays. <i>Talanta</i> , 2019, 191, 103-108.	5.5	60
15	Simultaneous voltammetric determination of guanine and adenine using MnO <sub>2</sub> nanosheets and ionic liquid-functionalized graphene combined with a permeation-selective polydopamine membrane. <i>Mikrochimica Acta</i> , 2019, 186, 450.	5.0	51
16	Understanding the interaction of single-walled carbon nanotube (SWCNT) on estrogen receptor: A combined molecular dynamics and experimental study. <i>Ecotoxicology and Environmental Safety</i> , 2019, 172, 373-379.	6.0	10
17	A novel electrochemiluminescent emitter of europium hydroxide nanorods and its application in bioanalysis. <i>Chemical Communications</i> , 2019, 55, 12479-12482.	4.1	20
18	CoS <sub>2</sub> -decorated ionic liquid-functionalized graphene as a novel hydrazine electrochemical sensor. <i>Talanta</i> , 2018, 182, 529-535.	5.5	59

#	ARTICLE	IF	CITATIONS
19	Reduced graphene oxide functionalized with a CoS <sub>2</sub> /ionic liquid composite and decorated with gold nanoparticles for voltammetric sensing of dopamine. <i>Mikrochimica Acta</i> , 2018, 185, 166.	5.0	48
20	A further development of the QNAR model to predict the cellular uptake of nanoparticles by pancreatic cancer cells. <i>Food and Chemical Toxicology</i> , 2018, 112, 571-580.	3.6	13
21	Prediction of the Toxicity of Binary Mixtures by QSAR Approach Using the Hypothetical Descriptors. <i>International Journal of Molecular Sciences</i> , 2018, 19, 3423.	4.1	28
22	Estimation of the Toxicity of Different Substituted Aromatic Compounds to the Aquatic Ciliate <i>Tetrahymena pyriformis</i> by QSAR Approach. <i>Molecules</i> , 2018, 23, 1002.	3.8	5
23	Preparation of gold nanoparticles supported on graphene oxide with flagella as the template for nonenzymatic hydrogen peroxide sensing. <i>Analytical and Bioanalytical Chemistry</i> , 2018, 410, 5915-5921.	3.7	10
24	What's the Key Factor to Ensure the Photoactivity Enhancement of Fe <sub>2</sub> O <sub>3</sub> Films with Ni(OH) <sub>2</sub> Loading: Clues from a Structural Modification with Flagella Nanowires. <i>Journal of Physical Chemistry C</i> , 2017, 121, 25364-25371.	3.1	4
25	Rational Design of Multi-Target Estrogen Receptors ER <sup>1</sup> and ER <sup>2</sup> by QSAR Approaches. <i>Current Drug Targets</i> , 2017, 18, 576-591.	2.1	8
26	QSAR-Based Studies of Nanomaterials in the Environment. , 2017, , 1504-1532.		2
27	QSAR-Based Studies of Nanomaterials in the Environment. , 2017, , 1339-1366.		1
28	Determination of 6-Benzylaminopurine in Bean Sprouts by Capillary Electrophoresis Compared with HPLC. <i>Food Analytical Methods</i> , 2016, 9, 3025-3031.	2.6	14
29	One-step electrochemical fabrication of a nickel oxide nanoparticle/polyaniline nanowire/graphene oxide hybrid on a glassy carbon electrode for use as a non-enzymatic glucose biosensor. <i>RSC Advances</i> , 2016, 6, 92541-92546.	3.6	57
30	Determination of Vanillin in Milk Powder by Capillary Electrophoresis Combined with Dispersive Liquid-Liquid Microextraction. <i>Food Analytical Methods</i> , 2016, 9, 1706-1712.	2.6	36
31	Detection of sibutramine and phenolphthalein in functional foods using capillary electrophoresis. <i>Analytical Methods</i> , 2016, 8, 621-626.	2.7	10
32	Review of Structures Containing Fullerene-C60 for Delivery of Antibacterial Agents. Multitasking model for Computational Assessment of Safety Profiles. <i>Current Bioinformatics</i> , 2015, 10, 565-578.	1.5	9
33	Dispersive liquid-liquid microextraction combined with non-aqueous capillary electrophoresis for the determination of imazalil, prochloraz and thiabendazole in apples, cherry tomatoes and grape juice. <i>Journal of the Science of Food and Agriculture</i> , 2015, 95, 745-751.	3.5	27
34	Computational modeling in nanomedicine: prediction of multiple antibacterial profiles of nanoparticles using a quantitative structure-activity relationship perturbation model. <i>Nanomedicine</i> , 2015, 10, 193-204.	3.3	55
35	QSAR-Based Studies of Nanomaterials in the Environment. <i>Advances in Chemical and Materials Engineering Book Series</i> , 2015, , 506-534.	0.3	1
36	Classification of natural estrogen-like isoflavonoids and diphenolics by QSAR tools. <i>Combinatorial Chemistry and High Throughput Screening</i> , 2015, 18, 712-722.	1.1	0

#	ARTICLE	IF	CITATIONS
37	Development of a Capillary Zone Electrophoresis Method for Determination of Mebendazole and Levamisole Hydrochloride in a Combined Tablet and a Comparison with a LC Method. <i>Journal of AOAC INTERNATIONAL</i> , 2014, 97, 128-132.	1.5	9
38	Computational Tool for Risk Assessment of Nanomaterials: Novel QSTR-Perturbation Model for Simultaneous Prediction of Ecotoxicity and Cytotoxicity of Uncoated and Coated Nanoparticles under Multiple Experimental Conditions. <i>Environmental Science &amp; Technology</i> , 2014, 48, 14686-14694.	10.0	124
39	Application of a non-aqueous capillary electrophoresis method to the analysis of triclosan in personal care products. <i>Analytical Methods</i> , 2014, 6, 4723-4728.	2.7	6
40	Computer-aided nanotoxicology: assessing cytotoxicity of nanoparticles under diverse experimental conditions by using a novel QSTR-perturbation approach. <i>Nanoscale</i> , 2014, 6, 10623.	5.6	118
41	Computational ecotoxicology: Simultaneous prediction of ecotoxic effects of nanoparticles under different experimental conditions. <i>Environment International</i> , 2014, 73, 288-294.	10.0	102
42	QSPR and Flow Cytometry Analysis (QSPR-FCA): Review and New Findings on Parallel Study of Multiple Interactions of Chemical Compounds with Immune Cellular and Molecular Targets. <i>Current Drug Metabolism</i> , 2014, 15, 414-428.	1.2	24
43	Matrix Trace Operators: From Spectral Moments of Molecular Graphs and Complex Networks to Perturbations in Synthetic Reactions, Micelle Nanoparticles, and Drug ADME Processes. <i>Current Drug Metabolism</i> , 2014, 15, 470-488.	1.2	26
44	Prediction of the Estrogen Receptor Binding Affinity for both hER<sub>α</sub> and hER<sub>β</sub> by QSAR Approaches. <i>Letters in Drug Design and Discovery</i> , 2014, 11, 265-278.	0.7	4
45	Use of Experimental Design and Artificial Neural Network in Optimization of Capillary Electrophoresis for the Determination of Nicotinic Acid and Nicotinamide in Food Compared with High-Performance Liquid Chromatography. <i>Food Analytical Methods</i> , 2013, 6, 191-200.	2.6	11
46	Model for High-Throughput Screening of Multitarget Drugs in Chemical Neurosciences: Synthesis, Assay, and Theoretic Study of Rasagiline Carbamates. <i>ACS Chemical Neuroscience</i> , 2013, 4, 1393-1403.	3.5	50
47	SEPARATION AND DETERMINATION OF FIVE ACTIVE COMPONENTS IN EYE DROPS BY CAPILLARY ELECTROPHORESIS IN COMPARISON WITH HPLC. <i>Journal of Liquid Chromatography and Related Technologies</i> , 2013, 36, 549-560.	1.0	4
48	DETERMINATION OF AMILORIDE HYDROCHLORIDE AND FUROSEMIDE IN COMPOUND FUROSEMIDE TABLETS BY CAPILLARY ELECTROPHORESIS COMBINED WITH RESPONSE SURFACE METHODOLOGY AND ARTIFICIAL NEURAL NETWORK. <i>Journal of Liquid Chromatography and Related Technologies</i> , 2013, 36, 2905-2918.	1.0	5
49	Prediction of the baseline toxicity of non-polar narcotic chemical mixtures by QSAR approach. <i>Chemosphere</i> , 2013, 90, 1980-1986.	8.2	22
50	TOPS-MODE model of multiplexing neuroprotective effects of drugs and experimental-theoretic study of new 1,3-rasagiline derivatives potentially useful in neurodegenerative diseases. <i>Bioorganic and Medicinal Chemistry</i> , 2013, 21, 1870-1879.	3.0	48
51	Development and validation of a non-aqueous capillary electrophoresis method for simultaneous estimation of mebendazole and levamisole hydrochloride in compound mebendazole tablets. <i>Analytical Methods</i> , 2013, 5, 762-766.	2.7	8
52	Review of quantitative structure-activity/property relationship studies of dyes: recent advances and perspectives. <i>Coloration Technology</i> , 2013, 129, 173-186.	1.5	18
53	QSAR Modeling for the Antimalarial Activity of 1,4-Naphthoquinonyl Derivatives as Potential Antimalarial Agents. <i>Current Computer-Aided Drug Design</i> , 2013, 9, 95-107.	1.2	0
54	Determination of neotame in non-alcoholic beverage by capillary zone electrophoresis. <i>Journal of the Science of Food and Agriculture</i> , 2013, 93, 3334-3338.	3.5	13

#	ARTICLE	IF	CITATIONS
55	Multi-Target Inhibitors for Proteins Associated with Alzheimer: In Silico Discovery using Fragment-Based Descriptors. <i>Current Alzheimer Research</i> , 2013, 10, 117-124.	1.4	41
56	Abelson Tyrosine-Protein Kinase 1 as Principal Target for Drug Discovery Against Leukemias. Role of the Current Computer-Aided Drug Design Methodologies. <i>Current Topics in Medicinal Chemistry</i> , 2013, 12, 2745-2762.	2.1	15
57	Recent Advances on QSAR-Based Profiling of Agonist and Antagonist A3 Adenosine Receptor Ligands. <i>Current Topics in Medicinal Chemistry</i> , 2013, 13, 1048-1068.	2.1	4
58	Unified Multi-target Approach for the Rational in silico Design of Anti-bladder Cancer Agents. <i>Anti-Cancer Agents in Medicinal Chemistry</i> , 2013, 13, 791-800.	1.7	41
59	Photophysical Behavior and the Binding to Human Serum Albumin of a Novel Triazole Compound. <i>Acta Chimica Sinica</i> , 2013, 71, 837.	1.4	0
60	QSAR modeling for the antimalarial activity of 1,4-naphthoquinonyl derivatives as potential antimalarial agents. <i>Current Computer-Aided Drug Design</i> , 2013, 9, 95-107.	1.2	1
61	Cheminformatics in Multi-target Drug Discovery for Anti-cancer Therapy: In Silico Design of Potent and Versatile Anti-brain Tumor Agents. <i>Anti-Cancer Agents in Medicinal Chemistry</i> , 2012, 12, 678-685.	1.7	41
62	In Silico Discovery and Virtual Screening of Multi-Target Inhibitors for Proteins in Mycobacterium tuberculosis. <i>Combinatorial Chemistry and High Throughput Screening</i> , 2012, 15, 666-673.	1.1	41
63	Enhancing the function, non-additivity, and substitution position effect of the Li atom in the cation- $\pi$ interaction and its mechanism: an ab initio study of Li- $\pi$ -Li-substituted benzene complexes. <i>Molecular Physics</i> , 2012, 110, 65-74.	1.7	5
64	QSAR Studies of PTP1B Inhibitors: Recent Advances and Perspectives. <i>Current Medicinal Chemistry</i> , 2012, 19, 4208-4217.	2.4	10
65	Separation and Determination of Honokiol and Magnolol in Chinese Traditional Medicines by Capillary Electrophoresis with the Application of Response Surface Methodology and Radial Basis Function Neural Network. <i>Journal of Chromatographic Science</i> , 2012, 50, 71-75.	1.4	15
66	Determination of formaldehyde in aquatic products by micellar electrokinetic capillary chromatography with 2,4-dinitrophenylhydrazine derivatization. <i>Acta Chromatographica</i> , 2012, 24, 519-528.	1.3	9
67	Theoretical study on polyaniline gas sensors: Examinations of response mechanism for alcohol. <i>Synthetic Metals</i> , 2012, 162, 862-867.	3.9	27
68	Rational drug design for anti-cancer chemotherapy: Multi-target QSAR models for the in silico discovery of anti-colorectal cancer agents. <i>Bioorganic and Medicinal Chemistry</i> , 2012, 20, 4848-4855.	3.0	87
69	Predicting multiple ecotoxicological profiles in agrochemical fungicides: A multi-species cheminformatic approach. <i>Ecotoxicology and Environmental Safety</i> , 2012, 80, 308-313.	6.0	58
70	Cheminformatics in anti-cancer chemotherapy: Multi-target QSAR model for the in silico discovery of anti-breast cancer agents. <i>European Journal of Pharmaceutical Sciences</i> , 2012, 47, 273-279.	4.0	74
71	A ligand-based approach for the in silico discovery of multi-target inhibitors for proteins associated with HIV infection. <i>Molecular BioSystems</i> , 2012, 8, 2188.	2.9	38
72	Determination of purines in soybean milk by capillary electrophoresis in comparison with high performance liquid chromatography. <i>Analytical Methods</i> , 2012, 4, 3386.	2.7	15

#	ARTICLE	IF	CITATIONS
73	Recent Advances on A3 Adenosine Receptor Antagonists by QSAR Tools. <i>Current Topics in Medicinal Chemistry</i> , 2012, 12, 878-894.	2.1	7
74	Self-doped polyaniline on functionalized carbon cloth as electroactive materials for supercapacitor. <i>Electrochimica Acta</i> , 2012, 64, 17-22.	5.2	111
75	Determination of benzoyl peroxide, as benzoic acid, in wheat flour by capillary electrophoresis compared with HPLC. <i>Journal of the Science of Food and Agriculture</i> , 2012, 92, 960-964.	3.5	33
76	Matrix Solid-Phase Dispersion Extraction and Capillary Electrophoresis Determination of Tetracycline Residues in Milk. <i>Food Analytical Methods</i> , 2012, 5, 148-153.	2.6	68
77	QSAR Studies of PTP1B Inhibitors: 1, 2-Naphthoquinone Derivatives. <i>Letters in Drug Design and Discovery</i> , 2012, 9, 915-925.	0.7	0
78	Some measures for making halogen bonds stronger than hydrogen bonds in H <sub>2</sub> CSâ€“HOX (X) Tj ETQq0,0 0 rgBTj/Overlock	2.8	66
79	Influence of Hybridization and Cooperativity on the Properties of Au-Bonding Interaction: Comparison with Hydrogen Bonds. <i>Journal of Physical Chemistry A</i> , 2011, 115, 2853-2858.	2.5	26
80	Affinity prediction on A3 adenosine receptor antagonists: The chemometric approach. <i>Bioorganic and Medicinal Chemistry</i> , 2011, 19, 6853-6859.	3.0	12
81	Quantitative Structure-Wavelength Relationship Modeling of Porphin -Derivative Photosensitizers. <i>Combinatorial Chemistry and High Throughput Screening</i> , 2011, 14, 601-615.	1.1	1
82	Prediction of the maximum absorption wavelength of azobenzene dyes by QSPR tools. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2011, 83, 353-361.	3.9	16
83	Fragment-based QSAR model toward the selection of versatile anti-sarcoma leads. <i>European Journal of Medicinal Chemistry</i> , 2011, 46, 5910-5916.	5.5	43
84	Multi-target drug discovery in anti-cancer therapy: Fragment-based approach toward the design of potent and versatile anti-prostate cancer agents. <i>Bioorganic and Medicinal Chemistry</i> , 2011, 19, 6239-6244.	3.0	60
85	Application of quantitative structureâ€“activity relationship to the determination of binding constant based on fluorescence quenching. <i>Journal of Luminescence</i> , 2011, 131, 126-133.	3.1	8
86	Quantitative structureâ€“property relationship study for the prediction of characteristic infrared absorption of carbonyl group of commonly used carbonyl compounds. <i>Vibrational Spectroscopy</i> , 2011, 55, 49-57.	2.2	3
87	QSPR Study on the prediction of ionization potential of various organic compounds by heuristic method and radial basis function neural network. , 2011, , .		3
88	A 3D Nitrate-bridging Heterometal-azido Inorganic Polymer Containing Cd <sub>8</sub> Cube: Synthesis, Structure, Topology and Fluorescence. <i>Bulletin of the Korean Chemical Society</i> , 2011, 32, 1769-1772.	1.9	1
89	Determination of melamine in milk powder, milk and fish feed by capillary electrophoresis: a good alternative to HPLC. <i>Journal of the Science of Food and Agriculture</i> , 2010, 90, 2178-2182.	3.5	39
90	Partially covalent nature and substitution non-additivity of Au-bonding in H <sub>2</sub> Oâ€“AuCH <sub>3</sub> complex. <i>Chemical Physics Letters</i> , 2010, 498, 259-262.	2.6	11

#	ARTICLE	IF	CITATIONS
91	Quantitative structure–electrochemistry relationship for variously-substituted 9, 10-anthraquinones using both an heuristic method and a radial basis function neural network. <i>Dyes and Pigments</i> , 2010, 84, 148-152.	3.7	12
92	QSPR Study for Estimation of Density of Some Aromatic Explosives by Multiple Linear Regression Approach. <i>Propellants, Explosives, Pyrotechnics</i> , 2010, 35, 169-174.	1.6	6
93	QSPR study for the prediction of UV maximum absorption wavelength of coumarins by heuristic method and radial basis function neural network. , 2010, , .		0
94	Analysis of alkaloids in pharmaceutical preparations containing Kushen by capillary electrophoresis with application of experimental design and a quantitative structure-property relationship approach. <i>Acta Chromatographica</i> , 2010, 22, 445-457.	1.3	1
95	QSPR model to predict the thermal stabilities of second-order nonlinear optical (NLO) chromophore molecules. <i>Molecular Simulation</i> , 2009, 35, 248-257.	2.0	9
96	Studies on the quantitative relationship between the olfactory thresholds of pyrazine derivatives and their molecular structures. <i>Flavour and Fragrance Journal</i> , 2009, 24, 62-68.	2.6	6
97	A Quantitative Structure–Activity Relationship Study of Some Commercially Available Cephalosporins. <i>QSAR and Combinatorial Science</i> , 2009, 28, 1003-1009.	1.4	2
98	pH-controlled morphological structure of polyaniline during electrochemical deposition. <i>Electrochimica Acta</i> , 2009, 54, 6172-6177.	5.2	59
99	Prediction of hydrophile–lipophile balance values of anionic surfactants using a quantitative structure–property relationship. <i>Journal of Colloid and Interface Science</i> , 2009, 336, 773-779.	9.4	13
100	Application of experimental design and radial basis function neural network to the separation and determination of active components in traditional Chinese medicines by capillary electrophoresis. <i>Analytica Chimica Acta</i> , 2009, 638, 88-93.	5.4	33
101	Quantitative structure–max relationship study on flavones by heuristic method and radial basis function neural network. <i>Analytica Chimica Acta</i> , 2009, 649, 52-61.	5.4	6
102	Quantitative structure-property relationship study on the determination of binding constant by fluorescence quenching. <i>Open Chemistry</i> , 2009, 7, 59-65.	1.9	1
103	QSPR study for the prediction of half-wave potentials of benzoxazines by heuristic method and radial basis function neural network. <i>Open Chemistry</i> , 2009, 7, 439-445.	1.9	5
104	Quantitative structure–property relationship study for estimation of quantitative calibration factors of some organic compounds in gas chromatography. <i>Analytica Chimica Acta</i> , 2008, 612, 126-135.	5.4	14
105	The effect of methyl group on the cooperativity between three types of hydrogen bond: $O_i \cdots H_j \cdots O$ , $C_i \cdots H_j \cdots O$ , and $O_i \cdots H_j \cdots i$ . <i>International Journal of Quantum Chemistry</i> , 2008, 108, 558-566.	2.0	20
106	Classification of the fragrance properties of chemical compounds based on support vector machine and linear discriminant analysis. <i>Flavour and Fragrance Journal</i> , 2008, 23, 232-238.	2.6	12
107	Classification of estrogen receptor- $\beta$ ligands on the basis of their binding affinities using support vector machine and linear discriminant analysis. <i>European Journal of Medicinal Chemistry</i> , 2008, 43, 43-52.	5.5	14
108	Regulating Function of Methyl Group in Strength of $CH \cdots O$ Hydrogen Bond: A High-Level Ab Initio Study. <i>Journal of Physical Chemistry A</i> , 2008, 112, 3985-3990.	2.5	45

#	ARTICLE	IF	CITATIONS
109	Analysis of food additives by capillary electrophoresis. <i>Acta Chromatographica</i> , 2008, 20, 239-246.	1.3	8
110	QSPR analysis of air-to-blood distribution of volatile organic compounds. <i>Ecotoxicology and Environmental Safety</i> , 2008, 71, 731-739.	6.0	10
111	Prediction of quantitative calibration factors of some organic compounds in gas chromatography. <i>Analyst, The</i> , 2008, 133, 881.	3.5	10
112	Cooperativity between two types of hydrogen bond in H <sub>3</sub> Câ€“HCNâ€“HCN and H <sub>3</sub> Câ€“HNCâ€“HNC complexes. <i>Journal of Chemical Physics</i> , 2008, 128, 154102.	3.0	56
113	Prediction of atmospheric degradation data for POPs by gene expression programming. <i>SAR and QSAR in Environmental Research</i> , 2008, 19, 465-479.	2.2	6
114	Prediction of retention times for a large set of pesticides or toxicants based on support vector machine and the heuristic method. <i>Toxicology Letters</i> , 2007, 175, 136-144.	0.8	23
115	QSAR method for prediction of protein-peptide binding affinity: Application to MHC class I molecule HLA-A*0201. <i>Journal of Molecular Graphics and Modelling</i> , 2007, 26, 246-254.	2.4	16
116	Accurate quantitative structureâ€“property relationship model of mobilities of peptides in capillary zone electrophoresis. <i>Analyst, The</i> , 2006, 131, 1254.	3.5	24
117	Quantitative structure-activity relationship models for prediction of sensory irritants (logRD50) of volatile organic chemicals. <i>Chemosphere</i> , 2006, 63, 1142-1153.	8.2	30
118	QSPR study of permeability coefficients through low-density polyethylene based on radial basis function neural networks and the heuristic method. <i>Computational Materials Science</i> , 2006, 37, 454-461.	3.0	21
119	Comparison of the characterization on binding of alpinetin and cardamonin to lysozyme by spectroscopic methods. <i>International Journal of Biological Macromolecules</i> , 2006, 39, 165-173.	7.5	54
120	QSPR Study of Fluorescence Wavelengths ( $\lambda_{ex}/\lambda_{em}$ ) Based on the Heuristic Method and Radial Basis Function Neural Networks. <i>QSAR and Combinatorial Science</i> , 2006, 25, 147-155.	1.4	24
121	QSAR Prediction of the Penetration of Drugs Across a Polydimethylsiloxane Membrane. <i>QSAR and Combinatorial Science</i> , 2006, 25, 895-904.	1.4	8
122	QSAR Study of Polychlorinated Dibenzodioxins, Dibenzofurans, and Biphenyls using the Heuristic Method and Support Vector Machine. <i>QSAR and Combinatorial Science</i> , 2006, 25, 46-55.	1.4	22
123	Binding analysis of glycyrrhetic acid to human serum albumin: Fluorescence spectroscopy, FTIR, and molecular modeling. <i>Bioorganic and Medicinal Chemistry</i> , 2006, 14, 3210-3217.	3.0	341
124	Quantitative structureâ€“property relationships for pesticides in biopartitioning micellar chromatography. <i>Journal of Chromatography A</i> , 2006, 1113, 140-147.	3.7	28
125	Prediction of Milk/Plasma Drug Concentration (M/P) Ratio Using Support Vector Machine (SVM) Method. <i>Pharmaceutical Research</i> , 2006, 23, 41-48.	3.5	25
126	Prediction of retention time of a variety of volatile organic compounds based on the heuristic method and support vector machine. <i>Analytica Chimica Acta</i> , 2005, 537, 101-110.	5.4	51



#	ARTICLE	IF	CITATIONS
127	Prediction of pKa for Neutral and Basic Drugs Based on Radial Basis Function Neural Networks and the Heuristic Method. <i>Pharmaceutical Research</i> , 2005, 22, 1454-1460.	3.5	48
128	Support Vector Machine-based QSPR for the Prediction of Van der Waals' Constants. <i>QSAR and Combinatorial Science</i> , 2005, 24, 227-239.	1.4	13
129	Classification of the Carcinogenicity of N-Nitroso Compounds Based on Support Vector Machines and Linear Discriminant Analysis. <i>Chemical Research in Toxicology</i> , 2005, 18, 198-203.	3.3	63
130	Support Vector Machine and the Heuristic Method to Predict the Solubility of Hydrocarbons in Electrolyte. <i>Journal of Physical Chemistry A</i> , 2005, 109, 3485-3492.	2.5	23
131	Binding of the bioactive compound 5,7,4-trihydroxy-6,3,5-trimethoxyflavone to human serum albumin. <i>International Journal of Biological Macromolecules</i> , 2005, 37, 85-91.	7.5	9
132	Diagnosing Breast Cancer Based on Support Vector Machines. <i>Journal of Chemical Information and Computer Sciences</i> , 2003, 43, 900-907.	2.8	109
133	Iridium-Complex-Functionalized Magnetic Nanoparticles for Fluorescent Detection of Mercapto Drugs. <i>ACS Applied Nano Materials</i> , 0, , .	5.0	3