

# Nicolae Viorel Pavel

## List of Publications by Year in descending order

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111  
papers

3,606  
citations

109137

35  
h-index

155451

55  
g-index

113  
all docs

113  
docs citations

113  
times ranked

2780  
citing authors

#	ARTICLE	IF	CITATIONS
1	Effect of temperature on the association behavior in aqueous mixtures of an oppositely charged amphiphilic block copolymer and bile salt. <i>Polymer</i> , 2020, 206, 122871.	1.8	8
2	Self-Assembly of Model Amphiphilic Peptides in Nonaqueous Solvents: Changing the Driving Force for Aggregation Does Not Change the Fibril Structure. <i>Langmuir</i> , 2020, 36, 8451-8460.	1.6	7
3	Polymorphic Self-Organization of Lauroyl Peptide in Response to pH and Concentration. <i>Langmuir</i> , 2020, 36, 3941-3951.	1.6	7
4	Block copolymers as bile salt sequestrants: intriguing structures formed in a mixture of an oppositely charged amphiphilic block copolymer and bile salt. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 12518-12529.	1.3	18
5	<i>Arabidopsis</i> and <i>Chlamydomonas</i> phosphoribulokinase crystal structures complete the redox structural proteome of the Calvin-Benson cycle. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2019, 116, 8048-8053.	3.3	25
6	Bile Salts: Natural Surfactants and Precursors of a Broad Family of Complex Amphiphiles. <i>Langmuir</i> , 2019, 35, 6803-6821.	1.6	64
7	A Stereochemically Driven Supramolecular Polymerisation. <i>Chemistry - A European Journal</i> , 2018, 24, 8195-8204.	1.7	11
8	The effect of fatty acid binding in the acid isomerizations of albumin investigated with a continuous acidification method. <i>Colloids and Surfaces B: Biointerfaces</i> , 2018, 168, 109-116.	2.5	3
9	Bile acid derivative-based catanionic mixtures: versatile tools for superficial charge modulation of supramolecular lamellae and nanotubes. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 18957-18968.	1.3	17
10	Time-Dependent pH Scanning of the Acid-Induced Unfolding of Human Serum Albumin Reveals Stabilization of the Native Form by Palmitic Acid Binding. <i>Journal of Physical Chemistry B</i> , 2017, 121, 4388-4399.	1.2	20
11	Wormlike reverse micelles in lecithin/bile salt/water mixtures in oil. <i>Colloids and Surfaces A: Physicochemical and Engineering Aspects</i> , 2017, 532, 411-419.	2.3	19
12	Twisted nanoribbons from a RGD-bearing cholic acid derivative. <i>Colloids and Surfaces B: Biointerfaces</i> , 2017, 159, 183-190.	2.5	11
13	Supramolecular assembly of a thermoresponsive steroidal surfactant with an oppositely charged thermoresponsive block copolymer. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 1504-1515.	1.3	19
14	On the stability of lithocholate derivative supramolecular tubules. <i>RSC Advances</i> , 2017, 7, 512-517.	1.7	13
15	Crystal structure of a lithium salt of a glucosyl derivative of lithocholic acid. <i>Steroids</i> , 2016, 113, 87-94.	0.8	4
16	Structural Response of Human Serum Albumin to Oxidation: Biological Buffer to Local Formation of Hypochlorite. <i>Journal of Physical Chemistry B</i> , 2016, 120, 12261-12271.	1.2	13
17	Interaction between bile salt sodium glycodeoxycholate and PEO-PPO-PEO triblock copolymers in aqueous solution. <i>RSC Advances</i> , 2016, 6, 69313-69325.	1.7	19
18	Tailoring Supramolecular Nanotubes by Bile Salt Based Surfactant Mixtures. <i>Angewandte Chemie</i> , 2015, 127, 7124-7127.	1.6	7

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19	Tailoring Supramolecular Nanotubes by Bile Salt Based Surfactant Mixtures. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 7018-7021.	7.2	37
20	Multi stimuli response of a single surfactant presenting a rich self-assembly behavior. <i>RSC Advances</i> , 2015, 5, 37800-37806.	1.7	27
21	A tryptophan-substituted cholic acid: Expanding the family of labelled biomolecules. <i>Colloids and Surfaces A: Physicochemical and Engineering Aspects</i> , 2015, 483, 142-149.	2.3	9
22	Bile salts and derivatives: Rigid unconventional amphiphiles as dispersants, carriers and superstructure building blocks. <i>Current Opinion in Colloid and Interface Science</i> , 2015, 20, 170-182.	3.4	87
23	Unravelling the shape and structural assembly of the photosynthetic GAPDH-CP12-PRK complex from <i>Arabidopsis thaliana</i> by small-angle X-ray scattering analysis. <i>Acta Crystallographica Section D: Biological Crystallography</i> , 2015, 71, 2372-2385.	2.5	13
24	Characterization of Carbon Nanotube Dispersions in Solutions of Bile Salts and Derivatives Containing Aromatic Substituents. <i>Journal of Physical Chemistry B</i> , 2014, 118, 1012-1021.	1.2	35
25	On the self-assembly of a tryptophan labeled deoxycholic acid. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 19492.	1.3	19
26	Sugar-Bile Acid-Based Bolaamphiphiles: From Scrolls to Monodisperse Single-Walled Tubules. <i>Langmuir</i> , 2014, 30, 6358-6366.	1.6	27
27	Ibuprofen and Propofol Cobinding Effect on Human Serum Albumin Unfolding in Urea. <i>Journal of Physical Chemistry B</i> , 2014, 118, 10043-10051.	1.2	11
28	Drug-loaded nanoparticles and supramolecular nanotubes formed from a volatile microemulsion with bile salt derivatives. <i>Physical Chemistry Chemical Physics</i> , 2013, 15, 6016.	1.3	18
29	Between Peptides and Bile Acids: Self-Assembly of Phenylalanine Substituted Cholic Acids. <i>Journal of Physical Chemistry B</i> , 2013, 117, 9248-9257.	1.2	33
30	Crystal structure of head-to-head dimers of cholic and deoxycholic acid derivatives with different symmetric bridges. <i>Steroids</i> , 2013, 78, 247-254.	0.8	8
31	pH sensitive tubules of a bile acid derivative: a tubule opening by release of wall leaves. <i>Physical Chemistry Chemical Physics</i> , 2013, 15, 7560.	1.3	37
32	Cationic Gels Based on Cholic Acid Derivatives. <i>Langmuir</i> , 2013, 29, 12342-12351.	1.6	33
33	A Single Amino-Acid Substitution Allows Endo-Polygalacturonase of <i>Fusarium verticillioides</i> to Acquire Recognition by PGIP2 from <i>Phaseolus vulgaris</i> . <i>PLoS ONE</i> , 2013, 8, e80610.	1.1	23
34	Amino acid-bile acid based molecules: extremely narrow surfactant nanotubes formed by a phenylalanine-substituted cholic acid. <i>Chemical Communications</i> , 2012, 48, 12011.	2.2	34
35	Nanoparticles with a Bicontinuous Cubic Internal Structure Formed by Cationic and Non-ionic Surfactants and an Anionic Polyelectrolyte. <i>Langmuir</i> , 2012, 28, 16536-16546.	1.6	34
36	Formation of tubules by p-tert-butylphenylamide derivatives of chenodeoxycholic and ursodeoxycholic acids in aqueous solution. <i>Steroids</i> , 2012, 77, 1205-1211.	0.8	23

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37	Formation of host-guest and sandwich complexes by a $\beta$ -cyclodextrin derivative. <i>Journal of Inclusion Phenomena and Macrocyclic Chemistry</i> , 2011, 69, 245-253.	1.6	5
38	Structural Resolution of the Complex between a Fungal Polygalacturonase and a Plant Polygalacturonase-Inhibiting Protein by Small-Angle X-Ray Scattering Å. <i>Plant Physiology</i> , 2011, 157, 599-607.	2.3	38
39	Human serum albumin binding ibuprofen: A 3D description of the unfolding pathway in urea. <i>Biophysical Chemistry</i> , 2010, 147, 111-122.	1.5	40
40	Catanionic Tubules with Tunable Charge. <i>Angewandte Chemie - International Edition</i> , 2010, 49, 6604-6607.	7.2	55
41	Supramolecular Structures Generated by a <i>p</i> - <i>tert</i> -Butylphenylamide Derivative of Deoxycholic Acid. From Planar Sheets to Tubular Structures through Helical Ribbons. <i>Langmuir</i> , 2010, 26, 7768-7773.	1.6	20
42	Urea-Induced Denaturation Process on Defatted Human Serum Albumin and in the Presence of Palmitic Acid. <i>Journal of Physical Chemistry B</i> , 2009, 113, 12590-12602.	1.2	46
43	Aggregation Behavior of Tetracarboxylic Surfactants Derived from Cholic and Deoxycholic Acids and Ethylenediaminetetraacetic Acid. <i>Langmuir</i> , 2009, 25, 9037-9044.	1.6	13
44	Sodium Glycodeoxycholate and Glycocholate Mixed Aggregates in Gas and Solution Phases. <i>Journal of Physical Chemistry B</i> , 2009, 113, 7162-7169.	1.2	10
45	Kinetics of formation of supramolecular tubules of a sodium cholate derivative. <i>Soft Matter</i> , 2009, 5, 3018.	1.2	46
46	About the albumin structure in solution: cigar Expanded form versus heart Normal shape. <i>Physical Chemistry Chemical Physics</i> , 2008, 10, 6741.	1.3	70
47	Measurement of x-ray multielectron photoexcitations at the $\alpha$ K-edge of $\text{Hg}^{2+}$ . <i>Physical Review B</i> , 2008, 78, .	1.1	8
48	Human Serum Albumin Unfolding: A Small-Angle X-ray Scattering and Light Scattering Study. <i>Journal of Physical Chemistry B</i> , 2008, 112, 15460-15469.	1.2	32
49	Synthesis and Characterization of a New Gemini Surfactant Derived from 3 $\beta$ ,12 $\beta$ -Dihydroxy-5 $\alpha$ -cholan-24-amine (Steroid Residue) and Ethylenediaminetetraacetic Acid (Spacer). <i>Langmuir</i> , 2008, 24, 6060-6066.	1.6	47
50	Early Stages of Formation of Branched Host-Guest Supramolecular Polymers. <i>Journal of Physical Chemistry B</i> , 2008, 112, 8536-8541.	1.2	14
51	Evidence for Sevenfold Coordination in the First Solvation Shell of Hg(II) Aqua Ion. <i>Journal of the American Chemical Society</i> , 2007, 129, 5430-5436.	6.6	78
52	Study on the Structure of Host-Guest Supramolecular Polymers. <i>Macromolecules</i> , 2007, 40, 5899-5906.	2.2	22
53	Supramolecular Structures Generated by <i>p</i> - <i>tert</i> -Butylphenyl-amide Derivative of Cholic Acid: From Vesicles to Molecular Tubes. <i>Advanced Materials</i> , 2007, 19, 1752-1756.	11.1	78
54	Structure of Sodium Glycodeoxycholate Micellar Aggregates from Small-Angle X-ray Scattering and Light-Scattering Techniques. <i>Journal of Physical Chemistry B</i> , 2006, 110, 12351-12359.	1.2	20

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55	New Lamellar Structure Formed by an Adamantyl Derivative of Cholic Acid. <i>Journal of Physical Chemistry B</i> , 2006, 110, 13679-13681.	1.2	43
56	Comment on "Two-Dimensional NMR Study on the Structures of Micelles of Sodium Taurocholate". <i>Journal of Physical Chemistry B</i> , 2005, 109, 9849-9850.	1.2	6
57	Small-Angle X-ray Scattering and Light Scattering on Lysozyme and Sodium Glycocholate Micelles. <i>Journal of Physical Chemistry B</i> , 2005, 109, 23857-23869.	1.2	32
58	Correlation between Small-Angle X-ray Scattering Spectra and Apparent Diffusion Coefficients in the Study of Structure and Interaction of Sodium Taurodeoxycholate Micelles. <i>Journal of Physical Chemistry B</i> , 2005, 109, 6111-6120.	1.2	13
59	QELS and X-ray study of two dihydroxy bile salt aqueous solutions. <i>Colloids and Surfaces A: Physicochemical and Engineering Aspects</i> , 2004, 248, 79-84.	2.3	10
60	An Integrated Study of Small-Angle X-ray Scattering and Dynamic Light Scattering on Cylindrical Micelles of Sodium Glycodeoxycholate. <i>Journal of Physical Chemistry B</i> , 2004, 108, 3078-3085.	1.2	41
61	Study of Intermicellar Interactions and Micellar Sizes in Ionic Micelle Solutions by Comparing Collective Diffusion and Self-Diffusion Coefficients. <i>Journal of Physical Chemistry B</i> , 2004, 108, 4799-4805.	1.2	37
62	Double-Electron Excitation Channels at the Ca <sup>2+</sup> K-Edge of Hydrated Calcium Ion. <i>Journal of Physical Chemistry B</i> , 2004, 108, 11857-11865.	1.2	28
63	Diffusivity Study of Dihydroxy and Trihydroxy Bile Salt Systems. <i>Langmuir</i> , 2003, 19, 1319-1323.	1.6	12
64	Collective diffusion and self-diffusion coefficients comparison to separate interactions and micellar size effects on ionic micelle diffusivities: Cylindrical micelles of sodium taurodeoxycholate. <i>Journal of Chemical Physics</i> , 2003, 118, 2865.	1.2	22
65	Sodium Taurodeoxycholate Structure from Solid to Liquid Phase. <i>Langmuir</i> , 2002, 18, 2812-2816.	1.6	16
66	Hydrogen and Higher Shell Contributions in Zn <sup>2+</sup> , Ni <sup>2+</sup> , and Co <sup>2+</sup> Aqueous Solutions: An X-ray Absorption Fine Structure and Molecular Dynamics Study. <i>Journal of the American Chemical Society</i> , 2002, 124, 1958-1967.	6.6	175
67	Development and Validation of an Integrated Computational Approach for the Study of Ionic Species in Solution by Means of Effective Two-Body Potentials. The Case of Zn <sup>2+</sup> , Ni <sup>2+</sup> , and Co <sup>2+</sup> in Aqueous Solutions. <i>Journal of the American Chemical Society</i> , 2002, 124, 1968-1976.	6.6	92
68	Combined XANES and EXAFS analysis of Co <sup>2+</sup> , Ni <sup>2+</sup> , and Zn <sup>2+</sup> aqueous solutions. <i>Physical Review B</i> , 2002, 66, .	1.1	88
69	Evidence of distorted fivefold coordination of the Cu <sup>2+</sup> aqua ion from an x-ray-absorption spectroscopy quantitative analysis. <i>Physical Review B</i> , 2002, 65, .	1.1	131
70	EXAFS and molecular dynamics studies of ionic solutions. <i>Journal of Synchrotron Radiation</i> , 2001, 8, 173-177.	1.0	10
71	K- and L-edge XAFS determination of the local structure of aqueous Nd(III) and Eu(III). <i>Journal of Synchrotron Radiation</i> , 2001, 8, 666-668.	1.0	17
72	Complete spectrum of multielectron excitations at the Br K edge x-ray absorption spectra. <i>Physical Review B</i> , 2001, 64, .	1.1	10

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73	Determination of two- and three-body correlation functions in ionic solutions by means of MD and EXAFS investigations. <i>Journal of Synchrotron Radiation</i> , 1999, 6, 281-283.	1.0	5
74	Evidence of three-body correlation functions in Rb <sup>+</sup> and Sr <sup>2+</sup> acetonitrile solutions. <i>Journal of Chemical Physics</i> , 1999, 111, 5107-5115.	1.2	24
75	X-ray Absorption Study of Copper(II) <sup>2+</sup> Glycinate Complexes in Aqueous Solution. <i>Journal of Physical Chemistry B</i> , 1998, 102, 3114-3122.	1.2	49
76	Structural investigation of copper(II) chloride solutions using x-ray absorption spectroscopy. <i>Journal of Chemical Physics</i> , 1997, 107, 2807-2812.	1.2	59
77	Evidence for multielectron resonances at the SrKedge. <i>Physical Review A</i> , 1996, 53, 798-805.	1.0	57
78	Multielectron excitations at theLedges of barium in aqueous solution. <i>Physical Review B</i> , 1996, 54, 12129-12138.	1.1	48
79	An extended x-ray absorption fine structure study by employing molecular dynamics simulations: Bromide ion in methanolic solution. <i>Journal of Chemical Physics</i> , 1996, 104, 1779-1790.	1.2	48
80	Differential anomalous scattering study of probe molecules. <i>Nuclear Instruments &amp; Methods in Physics Research B</i> , 1995, 97, 539-542.	0.6	0
81	EXAFS Study of Micellar Aggregates of Bile Acid Rubidium Salts. <i>The Journal of Physical Chemistry</i> , 1995, 99, 5471-5480.	2.9	26
82	XAS Study of Solubilization Loci of Brominated Molecules in Aqueous Micellar Solutions. <i>The Journal of Physical Chemistry</i> , 1994, 98, 2982-2990.	2.9	10
83	Triplet correlations in the hydration shell of aquaions. <i>Chemical Physics Letters</i> , 1994, 225, 150-155.	1.2	59
84	An extended x-ray absorption fine structure study of aqueous solutions by employing molecular dynamics simulations. <i>Journal of Chemical Physics</i> , 1994, 100, 985-994.	1.2	133
85	Double-electron excitation channels at the BrKedge of HBr andBr <sub>2</sub> . <i>Physical Review A</i> , 1993, 47, 2055-2063.	1.0	101
86	Multiple scattering x-ray absorption analysis of simple brominated hydrocarbon molecules. <i>The Journal of Physical Chemistry</i> , 1993, 97, 5486-5494.	2.9	44
87	EXAFS study of probe molecules in micellar solutions. <i>The Journal of Physical Chemistry</i> , 1991, 95, 7880-7886.	2.9	32
88	From crystal to micelle: A new approach to the micellar structure. <i>Journal of Inclusion Phenomena and Macrocyclic Chemistry</i> , 1989, 7, 391-400.	1.6	48
89	EXAFS: a new approach to the structure of micellar aggregates. <i>The Journal of Physical Chemistry</i> , 1988, 92, 2858-2862.	2.9	70
90	Size and shape of sodium deoxycholate micellar aggregates. <i>The Journal of Physical Chemistry</i> , 1987, 91, 356-362.	2.9	141

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91	Intermolecular interactions in sodium deoxycholate micelles: an NMR study involving a spin-labeled cholestane. <i>The Journal of Physical Chemistry</i> , 1987, 91, 83-89.	2.9	39
92	Multiple-scattering effects in the K-edge x-ray-absorption near-edge structure of crystalline and amorphous silicon. <i>Physical Review B</i> , 1987, 36, 6426-6433.	1.1	73
93	Spherical wave exafs analysis of the silicon K-edge X-ray absorption spectrum. <i>Solid State Communications</i> , 1987, 61, 635-639.	0.9	15
94	SPHERICAL WAVES EXAFS AND MULTIPLE SCATTERING EFFECTS IN XANES OF THE K-EDGE SPECTRUM OF SILICON. <i>Journal De Physique Colloque</i> , 1986, 47, C8-71-C8-74.	0.2	3
95	Study of the 4:1 inclusion compound between deoxycholic acid and (E)-p-dimethylaminoazobenzene by vapour pressure measurements. <i>Thermochimica Acta</i> , 1985, 87, 231-238.	1.2	2
96	The structure of the 2/1 'channel' inclusion compound between deoxycholic acid and pinacolone, 2C <sub>24</sub> H <sub>40</sub> O <sub>4</sub> .C <sub>6</sub> H <sub>12</sub> O. <i>Acta Crystallographica Section C: Crystal Structure Communications</i> , 1985, 41, 229-232.	0.4	4
97	Nuclear magnetic resonance and x-ray studies on micellar aggregates of sodium deoxycholate. <i>The Journal of Physical Chemistry</i> , 1984, 88, 5720-5724.	2.9	100
98	Crystal structure of the 2:1 inclusion compound between deoxycholic acid and quadricyclane. <i>Journal of Inclusion Phenomena</i> , 1984, 1, 329-337.	0.6	8
99	Chain folding in single crystals of polytetrafluoroethylene. <i>Journal of Polymer Science, Polymer Physics Edition</i> , 1983, 21, 321-328.	1.0	4
100	Structure of the 4:1 inclusion compound between deoxycholic acid and (E)-p-dimethylaminoazobenzene. <i>Acta Crystallographica Section B: Structural Crystallography and Crystal Chemistry</i> , 1982, 38, 2615-2620.	0.4	14
101	Crystal structure and van der Waals energy study of the 2:1 inclusion compound between deoxycholic acid and norbornadiene. <i>Acta Crystallographica Section B: Structural Crystallography and Crystal Chemistry</i> , 1981, 37, 368-372.	0.4	12
102	Possible models for the polyethylene hexagonal phase. <i>Polymer</i> , 1980, 21, 973-974.	1.8	3
103	Chain folding in crystalline syndiotactic poly(vinyl chloride). <i>Journal of Polymer Science, Polymer Physics Edition</i> , 1979, 17, 753-762.	1.0	10
104	The crystal structure of the inclusion compound between cyclohexatriene, benzene and water. <i>Acta Crystallographica Section B: Structural Crystallography and Crystal Chemistry</i> , 1979, 35, 2605-2609.	0.4	34
105	Physicochemical study of crystals of ethylene-butadiene copolymers. <i>Journal of Polymer Science: Polymer Chemistry Edition</i> , 1978, 16, 115-127.	0.8	4
106	Minimization programs for potential energy calculations in crystals and isolated molecules and macromolecules. <i>Zeitschrift für Kristallographie</i> , 1976, 144, 64-75.	1.1	14
107	The molecular and crystal structure of an allylpalladium(II) triazenido complex: [(1-3- $\hat{I}$ -C <sub>3</sub> H <sub>5</sub> )Pd(II)(p-CH <sub>3</sub> C <sub>6</sub> H <sub>4</sub> NNNC <sub>6</sub> H <sub>4</sub> CH <sub>3</sub> -p)] <sub>2</sub> . <i>Journal of Organometallic Chemistry</i> , 1976, 108, 409-421.	0.8	32
108	An x-ray and conformational study of Kapton H. <i>Journal of Polymer Science, Polymer Physics Edition</i> , 1976, 14, 1553-1560.	1.0	38

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109	Potential energy calculations about the chain folding of $\hat{\pm}(\hat{\alpha}^{\sim})$ poly(L-alanine). <i>Polymer</i> , 1976, 17, 257-259.	1.8	5
110	Potential energy calculations of ethylene-butadiene copolymers. <i>Journal of Polymer Science: Polymer Chemistry Edition</i> , 1975, 13, 125-131.	0.8	12
111	Generation and best fitting of molecular models. <i>Acta Crystallographica Section B: Structural Crystallography and Crystal Chemistry</i> , 1972, 28, 1968-1969.	0.4	8