

# Jean-François Gal

## List of Publications by Year in descending order

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109  
papers

2,961  
citations

186265

28  
h-index

206112

48  
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115  
all docs

115  
docs citations

115  
times ranked

2553  
citing authors

#	ARTICLE	IF	CITATIONS
1	Free energies of cation-molecule complex formation and cation-solvent transfers. <i>Pure and Applied Chemistry</i> , 1990, 62, 17-23.	1.9	150
2	Revised and Expanded Scale of Gas-Phase Lithium Cation Basicities. An Experimental and Theoretical Study. <i>Journal of Physical Chemistry A</i> , 2000, 104, 2824-2833.	2.5	131
3	Consequences of proton transfer in guanidine. <i>Journal of Physical Organic Chemistry</i> , 2003, 16, 91-106.	1.9	109
4	Dinuclear Copper Intermediates in Copper(I)-Catalyzed Azide-Alkyne Cycloaddition Directly Observed by Electrospray Ionization Mass Spectrometry. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 3065-3068.	13.8	98
5	Thermochemical aspects of proton transfer in the gas phase. <i>Journal of Mass Spectrometry</i> , 2001, 36, 699-716.	1.6	94
6	Superbases in the gas phase. Part II. Further extension of the basicity scale using acyclic and cyclic guanidines. <i>Journal of Physical Organic Chemistry</i> , 1994, 7, 725-733.	1.9	89
7	Enhanced Basicity of Push-Pull Nitrogen Bases in the Gas Phase. <i>Chemical Reviews</i> , 2016, 116, 13454-13511.	47.7	85
8	Lithium-Cation/π Complexes of Aromatic Systems. The Effect of Increasing the Number of Fused Rings. <i>Journal of the American Chemical Society</i> , 2003, 125, 10394-10401.	13.7	82
9	Site of Protonation of Nicotine and Nornicotine in the Gas Phase: Pyridine or Pyrrolidine Nitrogen?. <i>Journal of the American Chemical Society</i> , 2002, 124, 10552-10562.	13.7	77
10	Characterization of oligomers from methylglyoxal under dark conditions: a pathway to produce secondary organic aerosol through cloud processing during nighttime. <i>Atmospheric Chemistry and Physics</i> , 2010, 10, 3803-3812.	4.9	74
11	Gas-phase structural (internal) effects in strong organic nitrogen bases. <i>Journal of Physical Organic Chemistry</i> , 2001, 14, 25-34.	1.9	72
12	Elemental characterization and source identification of PM2.5 using Positive Matrix Factorization: The Malraux road tunnel, Nice, France. <i>Atmospheric Research</i> , 2009, 94, 320-329.	4.1	71
13	Superbases in the gas phase: Amidine and guanidine derivatives with proton affinities larger than 1000 kJ mol <sup>-1</sup> . <i>Rapid Communications in Mass Spectrometry</i> , 1993, 7, 599-602.	1.5	68
14	An Overview of Lewis Basicity and Affinity Scales. <i>Journal of Chemical Education</i> , 2011, 88, 1651-1657.	2.3	67
15	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.	1.9	59
16	Determination of benzene, toluene, ethylbenzene and xylenes in air by solid phase micro-extraction/gas chromatography/mass spectrometry. <i>Analytical and Bioanalytical Chemistry</i> , 2004, 380, 824-830.	3.7	55
17	Gas-Phase Basicity and Acidity Trends in $\hat{1},\hat{1}^2$ -Unsaturated Amines, Phosphines, and Arsines. <i>Journal of the American Chemical Society</i> , 1999, 121, 4653-4663.	13.7	47
18	Gas-phase lithium-cation basicities of some benzene derivatives. <i>International Journal of Mass Spectrometry</i> , 2002, 219, 445-456.	1.5	47

#	ARTICLE	IF	CITATIONS
19	Linear solvation energy relationships. Part 32. A co-ordinate covalency parameter, $\hat{\nu}_4$ , which, in combination with the hydrogen bond acceptor basicity parameter, $\hat{\nu}_2$ , permits correlation of many properties of neutral oxygen and nitrogen bases (including aqueous pKa). Journal of the Chemical Society Perkin Transactions II, 1985, , 1583-1589.	0.9	46
20	Cesium cation affinities and basicities. International Journal of Mass Spectrometry, 2007, 267, 7-23.	1.5	45
21	Acidity Trends in $\hat{\nu}_2$ -Unsaturated Alkanes, Silanes, Germanes, and Stannanes. Journal of the American Chemical Society, 2001, 123, 6353-6359.	13.7	43
22	The P4 $\hat{\nu}_2$ -Li <sup>+</sup> Ion in the Gas Phase: A Planetary System. Journal of the American Chemical Society, 2000, 122, 4451-4454.	13.7	42
23	The Nicotinic Pharmacophore: Thermodynamics of the Hydrogen-Bonding Complexation of Nicotine, Nornicotine, and Models. Journal of Organic Chemistry, 2003, 68, 8208-8221.	3.2	42
24	Insight into the Mechanisms of the Multicomponent Ugi and Ugi $\hat{\nu}_2$ -Smiles Reactions by ESI $\hat{\nu}_2$ MS(/MS). European Journal of Organic Chemistry, 2014, 2014, 7087-7090.	2.4	39
25	Comparison of Br $\hat{\nu}_2$ sted acidities of neutral NH-acids in gas phase, dimethyl sulfoxide and water. International Journal of Mass Spectrometry and Ion Processes, 1998, 175, 61-69.	1.8	33
26	Thermogravimetric calibration of permeation tubes used for the preparation of gas standards for air pollution analysis. Analyst, The, 2005, 130, 1369.	3.5	33
27	The heats of formation, gas-phase acidities, and related thermochemical properties of the third-row hydrides GeH <sub>4</sub> , AsH <sub>3</sub> , SeH <sub>2</sub> and HBr from G2 ab initio calculations. International Journal of Mass Spectrometry and Ion Processes, 1997, 167-168, 689-696.	1.8	32
28	Complexes between Lithium Cation and Diphenylalkanes in the Gas Phase: The Pincer Effect. Chemistry - A European Journal, 2006, 12, 7676-7683.	3.3	32
29	Ab initio study of tautomerism and of basicity center preference in histamine, from gas phase to solution-comparison with experimental data (gas phase, solution, solid state). Journal of Physical Organic Chemistry, 2003, 16, 783-796.	1.9	30
30	Using Thermogravimetry for Weight Loss Monitoring of Permeation Tubes Used for Generation of Trace Concentration Gas Standards. Analytical Chemistry, 2002, 74, 305-307.	6.5	29
31	The Gas-Phase Acidity of HCP, CH <sub>3</sub> CP, HCAs, and CH <sub>3</sub> CAs: An Unexpected Enhanced Acidity of the Methyl Group. Chemistry - A European Journal, 2002, 8, 4919-4924.	3.3	29
32	On the Use of the Kinetic Method for the Determination of Proton Affinities by Fourier-transform Ion Cyclotron Resonance Mass Spectrometry. Rapid Communications in Mass Spectrometry, 1996, 10, 242-245.	1.5	28
33	Gas-Phase Basicities and Acidities of Ethyl-, Vinyl-, and Ethynylarsine. An Experimental and Theoretical Study. Journal of Physical Chemistry A, 1997, 101, 9525-9530.	2.5	28
34	Enhanced Li <sup>+</sup> Binding Energies in Alkylbenzene Derivatives: The Scorpion Effect. Chemistry - A European Journal, 2003, 9, 4330-4338.	3.3	28
35	Experimental and Theoretical Evidence of Basic Site Preference in Polyfunctional Superbasic Amidinazine: N <sup>1</sup> ,N <sup>1</sup> -Dimethyl-N <sup>2</sup> - $\hat{\nu}_2$ -(2-pyridylethyl)formamidine. Journal of Organic Chemistry, 2004, 69, 4023-4030.	3.2	28
36	Acidity Trends in $\hat{\nu}_2$ -Unsaturated Sulfur, Selenium, and Tellurium Derivatives: Comparison with C-, Si-, Ge-, Sn-, N-, P-, As-, and Sb-Containing Analogues. Chemistry - A European Journal, 2005, 11, 2145-2153.	3.3	28

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37	Lithium-cation and proton affinities of sulfoxides and sulfones: A fourier transform ion cyclotron resonance study. <i>Journal of the American Society for Mass Spectrometry</i> , 1997, 8, 262-269.	2.8	27
38	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.	1.9	24
39	Gas-Phase Basicity of Polyfunctional Amidinazines: Experimental Evidence of Preferred Site(s) of Protonation. <i>Journal of Organic Chemistry</i> , 2000, 65, 4635-4640.	3.2	24
40	Meta versus para substituent effect in the gas phase: Separation of inductive and resonance components. <i>Journal of Physical Organic Chemistry</i> , 1994, 7, 615-624.	1.9	23
41	Metal triflates and triflimides as Lewis superacids: preparation, synthetic application and affinity tests by mass spectrometry. <i>Journal of Physical Organic Chemistry</i> , 2013, 26, 87-97.	1.9	23
42	Application of FT-ICR-MS for the study of proton-transfer reactions involving biomolecules. <i>Analytical and Bioanalytical Chemistry</i> , 2007, 389, 1365-1380.	3.7	22
43	Experimental and theoretical study of carbon suboxide C <sub>3</sub> O <sub>2</sub> , protonated carbon suboxide C <sub>3</sub> HO <sub>2</sub> <sup>+</sup> and C <sub>3</sub> HO <sub>2</sub> <sup>•</sup> radical in the gas phase. <i>International Journal of Mass Spectrometry and Ion Processes</i> , 1995, 141, 241-255.	1.8	21
44	Fourier transform ion cyclotron resonance determination of lithium cation basicities by the kinetic method: upward extension of the scale to phosphoryl compounds. <i>Journal of Mass Spectrometry</i> , 1998, 33, 757-765.	1.6	21
45	Exceptionally High Proton and Lithium Cation Gas-Phase Basicity of the Anti-Diabetic Drug Metformin. <i>Journal of Physical Chemistry A</i> , 2017, 121, 8706-8718.	2.5	21
46	The gas-phase acidity and bond dissociation energies of hydrogen telluride. <i>International Journal of Mass Spectrometry and Ion Processes</i> , 1989, 93, 87-94.	1.8	20
47	Sterically Hindered Resonance in Methyl-Substituted Anilines in the Gas Phase. <i>Journal of Organic Chemistry</i> , 1994, 59, 8127-8131.	3.2	20
48	Gas-phase basicity of simple amides toward proton and lithium cation: an experimental and theoretical study. <i>European Journal of Mass Spectrometry</i> , 1999, 5, 259.	0.7	20
49	Computational Study of Cesium Cation Interactions with Neutral and Anionic Compounds Related to Soil Organic Matter. <i>Journal of Physical Chemistry A</i> , 2009, 113, 10734-10744.	2.5	19
50	Cyano substituent effects on enol and enethiol acidity and basicity: The protonation and deprotonation of 3-hydroxy-2-propenenitrile and its thio analogue. <i>International Journal of Mass Spectrometry</i> , 2007, 267, 125-133.	1.5	18
51	Gas-Phase Lithium Cation Basicity: Revisiting the High Basicity Range by Experiment and Theory. <i>Journal of the American Society for Mass Spectrometry</i> , 2014, 25, 1962-1973.	2.8	18
52	Gas-Phase Cation Basicities for Sulfuryl Species from Calculation and Experiment. <i>Journal of Organic Chemistry</i> , 1997, 62, 9203-9209.	3.2	17
53	Gas-Phase Basicities of Acid Anhydrides. <i>Journal of Physical Chemistry A</i> , 1998, 102, 9183-9192.	2.5	17
54	Are nicotinoids protonated on the pyridine or the amino nitrogen in the gas phase?. <i>Journal of Physical Organic Chemistry</i> , 2006, 19, 104-114.	1.9	17

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55	Gas-Phase Protonation and Deprotonation of Acrylonitrile Derivatives $Ni\frac{1}{2}C\frac{1}{2}CH\frac{3}{4}CH\frac{1}{4}X$ (X=CH <sub>3</sub> , NH <sub>2</sub> ) <i>J. Phys. Chem. A</i> , 2001, 105, 10784-10791.	3.31	17
56	Mass spectrometric characterization of metal triflates and triflimides (Lewis superacid catalysts) by electrospray ionization and tandem mass spectrometry. <i>Rapid Communications in Mass Spectrometry</i> , 2010, 24, 2611-2619.	1.5	17
57	Glossary of terms used in physical organic chemistry (IUPAC Recommendations 2021). <i>Pure and Applied Chemistry</i> , 2022, 94, 353-534.	1.9	17
58	Investigations of cluster ions formed between cesium cations and benzoic, salicylic and phthalic acids by electrospray mass spectrometry and density-functional theory calculations. Toward a modeling of the interaction of Cs <sup>+</sup> with humic substances. <i>Rapid Communications in Mass Spectrometry</i> , 2005, 19, 568-573.	1.5	16
59	Acidity and basicity of thiocarboxamides in the gas phase. <i>Journal of Physical Organic Chemistry</i> , 1994, 7, 511-517.	1.9	15
60	Interaction of the cesium cation with mono-, di-, and tricarboxylic acids in the gas phase. A Cs <sup>+</sup> affinity scale for cesium carboxylates ion pairs. <i>Journal of the American Society for Mass Spectrometry</i> , 2009, 20, 1912-1924.	2.8	15
61	Can Nitriles Be Stronger Bases Than Proton Sponges in the Gas Phase? A Computational Analysis. <i>Journal of Physical Chemistry A</i> , 2015, 119, 8225-8236.	2.5	15
62	A microcalorimetric method for the measurement of enthalpies of solution of gases in liquids. <i>Thermochimica Acta</i> , 1987, 115, 67-81.	2.7	14
63	Adduct formation between phthalate esters and Li <sup>+</sup> in the gas phase: a thermochemical study by FT-ICR mass spectrometry. <i>International Journal of Mass Spectrometry</i> , 2002, 217, 75-79.	1.5	14
64	SPME Sampling of BTEX before GC/MS Analysis: Examples of outdoor and indoor Air Quality Measurements in Public and Private Sites. <i>Annali Di Chimica</i> , 2005, 95, 757-766.	0.6	14
65	Exceptional proton affinities of push-pull nitriles substituted by the guanidino and phosphazeno groups. <i>RSC Advances</i> , 2015, 5, 25513-25517.	3.6	14
66	The guanylated bioamine agmatine: A theoretical investigation of its structure and exceptional high basicity in the gas phase. <i>Computational and Theoretical Chemistry</i> , 2017, 1109, 10-18.	2.5	14
67	Gas-phase basicity of aromatic azines: A short review on structural effects. <i>International Journal of Mass Spectrometry</i> , 2017, 418, 130-139.	1.5	14
68	Catalytic intramolecular carbonyl-ene reaction with ketones: evidence for a retro-ene process. <i>New Journal of Chemistry</i> , 2015, 39, 7453-7458.	2.8	13
69	Biguanide Antidiabetic Drugs: Ipeglimin Exhibits Higher Proton Basicity but Smaller Lithium-Cation Basicity than Metformin in Vacuo. <i>ACS Omega</i> , 2018, 3, 17842-17852.	3.5	13
70	Gas-phase ion-molecule reactions in organophosphorus esters. <i>J. Phys. Chem. A</i> , 1999, 34, 1296-1302.		12
71	Vinyl- and ethynylsilanes, -germanes and -stannanes. A new case of dissociative proton attachment. <i>Journal of Physical Organic Chemistry</i> , 2002, 15, 509-513.	1.9	12
72	Steric Effects in Isolated Molecules: Gas-Phase Basicity of Methyl-Substituted Acetophenones. <i>European Journal of Organic Chemistry</i> , 2005, 2005, 2580-2588.	2.4	12

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73	Experimental (FT-ICR) and theoretical (DFT) estimation of the basic site preference for the bidentate molecule 2-( $\beta$ -aminoethyl)-pyridine: similarity with histamine. <i>Journal of Physical Organic Chemistry</i> , 2005, 18, 856-863.	1.9	12
74	A study of the cesium cation bonding to carboxylate anions by the kinetic method and quantum chemical calculations. <i>Journal of Mass Spectrometry</i> , 2010, 45, 520-527.	1.6	12
75	Bond Strength and Reactivity Scales for Lewis Superacid Adducts: A Comparative Study with $\text{In}(\text{OTf})_3$ and $\text{Al}(\text{OTf})_3$ . <i>ChemPhysChem</i> , 2017, 18, 683-691.	2.1	12
76	Bonding energetics in clusters formed by cesium salts: a study by collision-induced dissociation and density functional theory. <i>Rapid Communications in Mass Spectrometry</i> , 2006, 20, 2057-2062.	1.5	11
77	Theoretical modeling of sensitivity factors of Bayard-Alpert ionization gauges. <i>International Journal of Mass Spectrometry</i> , 2013, 341-342, 52-58.	1.5	11
78	Ab Initio and Experimental Thermodynamic and Kinetic Study of Proton-Assisted Bond Activation in Gaseous Hydrocarbons: Deconvolution of Reaction Efficiencies in the Case of Adamantane. <i>Chemistry - A European Journal</i> , 2006, 12, 5505-5513.	3.3	10
79	Effect of the Number of Methyl Groups on the Cation Affinity of Oxygen, Nitrogen, and Phosphorus Sites of Lewis Bases. <i>Journal of Physical Chemistry A</i> , 2016, 120, 9109-9116.	2.5	10
80	Push-Pull Effect on the Gas-Phase Basicity of Nitriles: Transmission of the Resonance Effects by Methylene-cyclopropene and Cyclopropenimine $\pi$ -Systems Substituted by Two Identical Strong Electron Donors. <i>Symmetry</i> , 2021, 13, 1554.	2.2	10
81	Relative basicities toward metal triflates Lewis acids by electrospray mass spectrometry. <i>Chemical Communications</i> , 2010, 46, 8472.	4.1	9
82	A Quantitative Approach of the Interaction between Metal Triflates and Organic Ligands Using Electrospray Mass Spectrometry. <i>Journal of the American Society for Mass Spectrometry</i> , 2012, 23, 2059-2062.	2.8	9
83	On the Significance of Lone Pair/Lone Pair and Lone Pair/Bond Pair Repulsions in the Cation Affinity and Lewis Acid/Lewis Base Interactions. <i>ACS Omega</i> , 2018, 3, 11331-11339.	3.5	9
84	Gas-phase lithium cation basicity of histamine and its agonist 2-( $\beta$ -aminoethyl)-pyridine. <i>International Journal of Mass Spectrometry</i> , 2007, 267, 315-323.	1.5	8
85	Bonding between the cesium cation and substituted benzoic acids or benzoate anions in the gas phase: A density functional theory and mass spectrometric study. <i>Collection of Czechoslovak Chemical Communications</i> , 2009, 74, 167-188.	1.0	8
86	Measuring Gas-Phase Basicities Relative to the Lithium Cation by Mass Spectrometry: A Physical Chemistry Experiment. <i>Journal of Chemical Education</i> , 2012, 89, 1476-1478.	2.3	8
87	Aluminum Monocation Basicity and Affinity Scales. <i>European Journal of Mass Spectrometry</i> , 2015, 21, 517-532.	1.0	8
88	Directionality of Cation/Molecule Bonding in Lewis Bases Containing the Carbonyl Group. <i>Journal of Physical Chemistry A</i> , 2017, 121, 6810-6822.	2.5	8
89	The gas-phase basicity of ethyl-, ethenyl- and ethynylphosphines and arsines. <i>International Journal of Mass Spectrometry and Ion Processes</i> , 1998, 175, 27-33.	1.8	7
90	Interaction between the Cesium Cation and Cesium Carboxylates: An Extended $\text{Cs}^+$ Basicity Scale. <i>ChemPlusChem</i> , 2013, 78, 1195-1204.	2.8	7

#	ARTICLE	IF	CITATIONS
91	Comment on the Article "Gutmann Donor and Acceptor Numbers for Ionic Liquids" by M. Schmeisser, P. Illner, R. Puchta, A. Zahl, and R. vanâ€¦Eldik (<i>Chem. Eur. J.</i> 2012, <i>18</i>, 10969â€“10982). Chemistry 3.3 A European Journal, 2013, 19, 16832-16834.		7
92	Enthalpies of Adduct Formation between Boron Trifluoride and Selected Organic Bases in Solution: Toward an Accurate Theoretical Entry to Lewis Basicity. <i>Molecules</i> , 2021, 26, 6659.	3.8	7
93	On the Lewis Basicity of Phosphoramides: A Critical Examination of Their Donor Number through Comparison of Enthalpies of Adduct Formation with SbCl 5 and BF 3. <i>ChemPhysChem</i> , 2019, 20, 2566-2576.	2.1	6
94	Quality Control of a Functionalized Polymer Catalyst by Energy Dispersive X-ray Spectrometry (EDX or Tj ETQq0 0 0 rgBT /Overlock 10 T	6.5	6
95	Is Allylphosphine a Carbon or a Phosphorus Base in the Gas Phase?. <i>European Journal of Mass Spectrometry</i> , 2003, 9, 245-255.	1.0	5
96	Catalytic Effect of Cesium Cation Adduct Formation on the Decarboxylation of Carboxylate Ions in the Gas Phase. <i>Chemistry - A European Journal</i> , 2014, 20, 815-823.	3.3	5
97	Quantitative Ligand Affinity Scales for Metal Triflate Salts: Application to Isomer Differentiation. <i>ChemPlusChem</i> , 2017, 82, 498-506.	2.8	5
98	Assessment of the nature of interactions of cations with cycloheptatriene derivatives using change in the aromaticity: Comparison with electron density and NBO results. <i>Molecular Physics</i> , 2020, 118, e1662507.	1.7	5
99	Purine tautomeric preferences and bond-length alternation in relation with protonation-deprotonation and alkali metal cationization. <i>Journal of Molecular Modeling</i> , 2020, 26, 93.	1.8	5
100	Can the Lewis Basicity of an Isolated Solvent Molecule be Used for Characterizing Solvent Effects?. <i>Current Analytical Chemistry</i> , 2021, 17, 328-338.	1.2	4
101	Nitriles with High Gas-Phase Basicityâ€”Part II Transmission of the Pushâ€“Pull Effect through Methylene-cyclopropene and Cyclopropanimine Scaffolds Intercalated between Different Electron Donor(s) and the Cyano N-Protonation Site. <i>Molecules</i> , 2022, 27, 4370.	3.8	4
102	Metal Triflates as Catalysts in Organic Synthesis: Determination of Their Lewis Acidity by Mass Spectrometry. <i>ChemPlusChem</i> , 2022, 87, e202200037.	2.8	3
103	A Short Note on the History of Chromatography at the University of Tartu, Estonia. <i>Chromatographia</i> , 2010, 72, 203-204.	1.3	2
104	An Excursion into the History of Chromatography: Mikhail Tswett, From Asti, Italy, to Tartu, Estonia. <i>Chromatographia</i> , 2019, 82, 519-521.	1.3	2
105	Alkali Metal Cations Bonding to Carboxylate Anions: Studies using Mass Spectrometry and Quantum Chemical Calculations. <i>Journal of Physical Chemistry A</i> , 2020, 124, 4390-4399.	2.5	2
106	Organometallic superacids and hyperacids: Acidity enhancement by internal bonding with a strong electron-pair acceptor group BX2. <i>Chemical Physics Letters</i> , 2021, 763, 138207.	2.6	2
107	Energetics and Structures of Adducts of JohnPhos(Au+), PPh3(Au+), and IPr(Au+) with Organic Substrates: A Mass Spectrometry and DFT Study. <i>Organometallics</i> , 2021, 40, 1642-1653.	2.3	0
108	Evolution of Chemical Research in Nice, CÃ“te dÃ“Azur: From Early Laboratories to the â€“Institut de Chimie de Niceâ€™. <i>ChemPlusChem</i> , 2022, , e202100532.	2.8	0

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109	Front Cover: Evolution of Chemical Research in Nice, C�te d'Azur: From Early Laboratories to the Institut de Chimie de Nice (ChemPlusChem 6/2022). ChemPlusChem, 2022, 87, .	2.8	0