

# Carmine D'Agostino

## List of Publications by Year in descending order

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74  
papers

3,639  
citations

136885

32  
h-index

133188

59  
g-index

78  
all docs

78  
docs citations

78  
times ranked

4133  
citing authors

#	ARTICLE	IF	CITATIONS
1	Glycerol eutectics as sustainable solvent systems. <i>Green Chemistry</i> , 2011, 13, 82-90.	4.6	666
2	Molecular motion and ion diffusion in choline chloride based deep eutectic solvents studied by <sup>1</sup> H pulsed field gradient NMR spectroscopy. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 21383.	1.3	397
3	Molecular and ionic diffusion in aqueous " deep eutectic solvent mixtures: probing inter-molecular interactions using PFG NMR. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 15297-15304.	1.3	204
4	Solvent effects in the hydrogenation of 2-butanone. <i>Journal of Catalysis</i> , 2012, 289, 30-41.	3.1	140
5	Neutron diffraction, NMR and molecular dynamics study of glucose dissolved in the ionic liquid 1-ethyl-3-methylimidazolium acetate. <i>Chemical Science</i> , 2011, 2, 1594.	3.7	121
6	Structure and Dynamics of 1-Ethyl-3-methylimidazolium Acetate via Molecular Dynamics and Neutron Diffraction. <i>Journal of Physical Chemistry B</i> , 2010, 114, 7760-7768.	1.2	117
7	Mesoscopic Structuring and Dynamics of Alcohol/Water Solutions Probed by Terahertz Time-Domain Spectroscopy and Pulsed Field Gradient Nuclear Magnetic Resonance. <i>Journal of Physical Chemistry B</i> , 2014, 118, 10156-10166.	1.2	109
8	Interpretation of NMR Relaxation as a Tool for Characterising the Adsorption Strength of Liquids inside Porous Materials. <i>Chemistry - A European Journal</i> , 2014, 20, 13009-13015.	1.7	98
9	Adsorption of pyridine from aqueous solutions by polymeric adsorbents MN 200 and MN 500. Part 2: Kinetics and diffusion analysis. <i>Chemical Engineering Journal</i> , 2016, 306, 1223-1233.	6.6	85
10	Light-driven, heterogeneous organocatalysts for C-C bond formation toward valuable perfluoroalkylated intermediates. <i>Science Advances</i> , 2020, 6, .	4.7	75
11	Creation of Al-Enriched Mesoporous ZSM-5 Nanoboxes with High Catalytic Activity: Converting Tetrahedral Extra-framework Al into Framework Sites by Post Treatment. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 19478-19486.	7.2	69
12	Hydrogen Bonding Network Disruption in Mesoporous Catalyst Supports Probed by PFG-NMR Diffusometry and NMR Relaxometry. <i>Journal of Physical Chemistry C</i> , 2012, 116, 8975-8982.	1.5	65
13	Unravelling mass transport in hierarchically porous catalysts. <i>Journal of Materials Chemistry A</i> , 2019, 7, 11814-11825.	5.2	57
14	Deep eutectic solvents: alternative reaction media for organic oxidation reactions. <i>Reaction Chemistry and Engineering</i> , 2021, 6, 582-598.	1.9	57
15	Adsorption of pyridine from aqueous solutions by polymeric adsorbents MN 200 and MN 500. Part 1: Adsorption performance and PFG-NMR studies. <i>Chemical Engineering Journal</i> , 2016, 306, 67-76.	6.6	52
16	Understanding the Solvent Effect on the Catalytic Oxidation of 1,4-Butanediol in Methanol over Au/TiO <sub>2</sub> Catalyst: NMR Diffusion and Relaxation Studies. <i>Chemistry - A European Journal</i> , 2012, 18, 14426-14433.	1.7	50
17	Diffusion, Ion Pairing and Aggregation in 1-Ethyl-3-Methylimidazolium-Based Ionic Liquids Studied by <sup>1</sup> H and <sup>19</sup> F PFG NMR: Effect of Temperature, Anion and Glucose Dissolution. <i>ChemPhysChem</i> , 2018, 19, 1081-1088.	1.0	50
18	Effect of solvent on the hydrogenation of 4-phenyl-2-butanone over Pt based catalysts. <i>Journal of Catalysis</i> , 2015, 330, 344-353.	3.1	49

#	ARTICLE	IF	CITATIONS
19	Prediction of binary diffusion coefficients in non-ideal mixtures from NMR data: Hexane–nitrobenzene near its consolute point. <i>Chemical Engineering Science</i> , 2011, 66, 3898-3906.	1.9	48
20	Solvent Effect and Reactivity Trend in the Aerobic Oxidation of 1,3-Propanediols over Gold Supported on Titania: NMR Diffusion and Relaxation Studies. <i>Chemistry - A European Journal</i> , 2013, 19, 11725-11732.	1.7	46
21	Assessing the surface modifications following the mechanochemical preparation of a Ag/Al <sub>2</sub> O <sub>3</sub> selective catalytic reduction catalyst. <i>Catalysis Science and Technology</i> , 2014, 4, 531-539.	2.1	46
22	Solvent inhibition in the liquid-phase catalytic oxidation of 1,4-butanediol: understanding the catalyst behaviour from NMR relaxation time measurements. <i>Catalysis Science and Technology</i> , 2016, 6, 7896-7901.	2.1	46
23	Do group 1 metal salts form deep eutectic solvents?. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 25528-25537.	1.3	43
24	Increased Affinity of Small Gold Particles for Glycerol Oxidation over Au/TiO <sub>2</sub> Probed by NMR Relaxation Methods. <i>ACS Catalysis</i> , 2017, 7, 4235-4241.	5.5	43
25	Direct Correlation between Adsorption Energetics and Nuclear Spin Relaxation in a Liquid-Saturated Catalyst Material. <i>ChemPhysChem</i> , 2018, 19, 2472-2479.	1.0	40
26	Effect of Al content on number and location of hydroxyl acid species in zeolites: a DRIFTS quantitative protocol without the need for molar extinction coefficients. <i>RSC Advances</i> , 2017, 7, 52604-52613.	1.7	37
27	The Effect of Grafting Zirconia and Ceria onto Alumina as a Support for Silicotungstic Acid for the Catalytic Dehydration of Glycerol to Acrolein. <i>Chemistry - A European Journal</i> , 2014, 20, 1743-1752.	1.7	36
28	Probing hydrogen-bonding in binary liquid mixtures with terahertz time-domain spectroscopy: a comparison of Debye and absorption analysis. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 5999-6008.	1.3	36
29	Pulsed-Field Gradient NMR Spectroscopic Studies of Alcohols in Supported Gold Catalysts. <i>Journal of Physical Chemistry C</i> , 2011, 115, 1073-1079.	1.5	35
30	<i>In situ</i> study of reaction kinetics using compressed sensing NMR. <i>Chemical Communications</i> , 2014, 50, 14137-14140.	2.2	35
31	Deactivation studies of a carbon supported AuPt nanoparticulate catalyst in the liquid-phase aerobic oxidation of 1,2-propanediol. <i>Catalysis Science and Technology</i> , 2014, 4, 1313-1322.	2.1	34
32	A local composition model for the prediction of mutual diffusion coefficients in binary liquid mixtures from tracer diffusion coefficients. <i>Chemical Engineering Science</i> , 2015, 132, 250-258.	1.9	33
33	Prediction of mutual diffusion coefficients in non-ideal mixtures from pulsed field gradient NMR data: Triethylamine–water near its consolute point. <i>Chemical Engineering Science</i> , 2012, 74, 105-113.	1.9	32
34	Assessing the effect of reducing agents on the selective catalytic reduction of NO <sub>x</sub> over Ag/Al <sub>2</sub> O <sub>3</sub> catalysts. <i>Catalysis Science and Technology</i> , 2016, 6, 1661-1666.	2.1	32
35	The effect of coke deposition on the activity and selectivity of the HZSM-5 zeolite during ethylbenzene alkylation reaction in the presence of ethanol. <i>Catalysis Science and Technology</i> , 2014, 4, 1017.	2.1	31
36	HfN Nanoparticles: An Unexplored Catalyst for the Electrocatalytic Oxygen Evolution Reaction. <i>Angewandte Chemie - International Edition</i> , 2019, 58, 15464-15470.	7.2	31

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37	Prediction of the mutual diffusivity in acetone–chloroform liquid mixtures from the tracer diffusion coefficients. <i>Chemical Engineering Science</i> , 2013, 95, 43-47.	1.9	30
38	Structure and dynamics of aqueous 2-propanol: a THz-TDS, NMR and neutron diffraction study. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 30481-30491.	1.3	29
39	Effect of Al content on the strength of terminal silanol species in ZSM-5 zeolite catalysts: a quantitative DRIFTS study without the use of molar extinction coefficients. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 4250-4262.	1.3	28
40	Prediction of mutual diffusion coefficients in binary liquid systems with one self-associating component from viscosity data and intra-diffusion coefficients at infinite dilution. <i>Chemical Engineering Science</i> , 2016, 147, 118-127.	1.9	27
41	Nanoscale Clustering of Alcoholic Solutes in Deep Eutectic Solvents Studied by Nuclear Magnetic Resonance and Dynamic Light Scattering. <i>ACS Sustainable Chemistry and Engineering</i> , 2019, 7, 15086-15092.	3.2	26
42	Oxidative Coupling of Aldehydes with Alcohol for the Synthesis of Esters Promoted by Polystyrene-Supported N-Heterocyclic Carbene: Unraveling the Solvent Effect on the Catalyst Behavior Using NMR Relaxation. <i>Organic Letters</i> , 2020, 22, 4927-4931.	2.4	22
43	Tailoring pore structure and surface chemistry of microporous Alumina-Carbon Molecular Sieve Membranes (Al-CMSMs) by altering carbonization temperature for optimal gas separation performance: An investigation using low-field NMR relaxation measurements. <i>Chemical Engineering Journal</i> , 2021, 424, 129313.	6.6	21
44	Product Inhibition in Glycerol Oxidation over Au/TiO <sub>2</sub> Catalysts Quantified by NMR Relaxation. <i>ACS Catalysis</i> , 2018, 8, 7334-7339.	5.5	20
45	Nuclear spin relaxation as a probe of zeolite acidity: a combined NMR and TPD investigation of pyridine in HZSM-5. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 17752-17760.	1.3	19
46	Globular and Fibrous Proteins Modified with Deep Eutectic Solvents: Materials for Drug Delivery. <i>Molecules</i> , 2019, 24, 3583.	1.7	18
47	Alkaline pretreatment of walnut shells increases pore surface hydrophilicity of derived biochars. <i>Applied Surface Science</i> , 2022, 571, 151253.	3.1	17
48	Effect of paramagnetic species on T <sub>1</sub> , T <sub>2</sub> and T <sub>1</sub> /T <sub>2</sub> NMR relaxation times of liquids in porous CuSO <sub>4</sub> /Al <sub>2</sub> O <sub>3</sub> . <i>RSC Advances</i> , 2017, 7, 36163-36167.	1.7	16
49	Mutual diffusion in concentrated liquid solutions: A new model based on cluster theory. <i>Journal of Molecular Liquids</i> , 2017, 232, 516-521.	2.3	15
50	A predictive model for the diffusion of a highly non-ideal ternary system. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 18436-18446.	1.3	15
51	Maxwell–Stefan diffusion coefficient estimation for ternary systems: an ideal ternary alcohol system. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 16071-16077.	1.3	14
52	NMR relaxation time measurements of solvent effects in an organocatalysed asymmetric aldol reaction over silica SBA-15 supported proline. <i>Reaction Chemistry and Engineering</i> , 2022, 7, 269-274.	1.9	14
53	Correlating the strength of reducing agent adsorption with Ag/Al <sub>2</sub> O <sub>3</sub> catalyst performances in selective catalytic reduction (SCR) of NO <sub>x</sub> . <i>Catalysis Today</i> , 2022, 384-386, 274-278.	2.2	13
54	A kinetic analysis methodology to elucidate the roles of metal, support and solvent for the hydrogenation of 4-phenyl-2-butanone over Pt/TiO <sub>2</sub> . <i>Journal of Catalysis</i> , 2015, 330, 362-373.	3.1	12

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55	Exploring catalyst passivation with NMR relaxation. <i>Faraday Discussions</i> , 2017, 204, 439-452.	1.6	12
56	Swelling-induced structural changes and microparticle uptake of gelatin gels probed by NMR and CLSM. <i>Soft Matter</i> , 2017, 13, 2952-2961.	1.2	12
57	Creation of Al-Enriched Mesoporous ZSM-5 Nanoboxes with High Catalytic Activity: Converting Tetrahedral Extra-framework Al into Framework Sites by Post Treatment. <i>Angewandte Chemie</i> , 2020, 132, 19646-19654.	1.6	12
58	Tailoring morphology of hierarchical catalysts for tuning pore diffusion behaviour: a rational guideline exploiting bench-top pulsed-field gradient (PFG) nuclear magnetic resonance (NMR). <i>Molecular Systems Design and Engineering</i> , 2020, 5, 1193-1204.	1.7	12
59	Hole theory as a prediction tool for Brownian diffusive motion in binary mixtures of liquids. <i>RSC Advances</i> , 2017, 7, 51864-51869.	1.7	10
60	NMR Investigation into the Influence of Surface Interactions on Liquid Diffusion in a Mesoporous Catalyst Support. <i>Topics in Catalysis</i> , 2020, 63, 319-327.	1.3	10
61	HfN Nanoparticles: An Unexplored Catalyst for the Electrocatalytic Oxygen Evolution Reaction. <i>Angewandte Chemie</i> , 2019, 131, 15610-15616.	1.6	9
62	Solvent Effects in the Homogeneous Catalytic Reduction of Propionaldehyde with Aluminium Isopropoxide Catalyst: New Insights from PFG NMR and NMR Relaxation Studies. <i>ChemPhysChem</i> , 2020, 21, 1101-1106.	1.0	9
63	Inhibitory effect of oxygenated heterocyclic compounds in mesoporous catalytic materials: A pulsed-field gradient NMR diffusion study. <i>Microporous and Mesoporous Materials</i> , 2018, 269, 88-92.	2.2	7
64	In situ high-pressure $^{13}\text{C}/^1\text{H}$ NMR reaction studies of benzyl alcohol oxidation over a $\text{Pd}/\text{Al}_2\text{O}_3$ catalyst. <i>Reaction Chemistry and Engineering</i> , 2020, 5, 1053-1057.	1.9	7
65	Exploiting enhanced paramagnetic NMR relaxation for monitoring catalyst preparation using $^1\text{T}$ - $^2\text{T}$ NMR correlation maps. <i>Reaction Chemistry and Engineering</i> , 2019, 4, 268-272.	1.9	6
66	Self-diffusion of glycerol in $\text{Al}_2\text{O}_3$ -alumina nanopores. The neglected role of pore saturation in the dynamics of confined polyalcohols. <i>Applied Surface Science</i> , 2020, 516, 146089.	3.1	6
67	$^{23}\text{Na}$ NMR $^1\text{T}$ relaxation measurements as a probe for diffusion and dynamics of sodium ions in salt-glycerol mixtures. <i>Journal of Chemical Physics</i> , 2021, 154, 224501.	1.2	6
68	Insights into Substituent Effects of Benzaldehyde Derivatives in a Heterogeneous Organocatalyzed Aldol Reaction. <i>ChemCatChem</i> , 2022, 14, .	1.8	6
69	Microstructure evolution during nano-emulsification by NMR and microscopy. <i>Journal of Colloid and Interface Science</i> , 2019, 551, 138-146.	5.0	4
70	Liquid-liquid equilibrium for the ternary system ethanol/toluene/n-decane: a correction to the existing coexistence curve and NRTL parameters. <i>Physics and Chemistry of Liquids</i> , 2017, 55, 669-673.	0.4	2
71	Base adsorption mechanism over zeolite catalysts at different Al contents probed by the tapered element oscillating microbalance (TEOM). <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 25357-25364.	1.3	2
72	Assessing the use of NMR chemical shifts for prediction of VLE in non-ideal binary liquid mixtures. <i>Chemical Engineering Science</i> , 2014, 119, 331-333.	1.9	1

#	ARTICLE	IF	CITATIONS
73	Direct Correlation between Adsorption Energetics and Nuclear Spin Relaxation in a Liquid-saturated Catalyst Material. <i>ChemPhysChem</i> , 2018, 19, 2448-2448.	1.0	1
74	A Neutron Diffraction and Molecular Dynamics Investigation of Acetate-Based Ionic Liquids as Solvents for Glucose. <i>ECS Transactions</i> , 2010, 33, 611-620.	0.3	0