

# Radha Dilip Banhatti

## List of Publications by Year in descending order

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53  
papers

1,204  
citations

331670

21  
h-index

377865

34  
g-index

55  
all docs

55  
docs citations

55  
times ranked

900  
citing authors

#	ARTICLE	IF	CITATIONS
1	Ionic motion in materials with disordered structures: conductivity spectra and the concept of mismatch and relaxation. <i>Physical Chemistry Chemical Physics</i> , 2002, 4, 3155-3167.	2.8	110
2	Ionic motion in materials with disordered structures. <i>Solid State Ionics</i> , 2006, 177, 1551-1557.	2.7	87
3	Anion reorientation in an ion conducting plastic crystal $\hat{\alpha}$ coherent quasielastic neutron scattering from sodium ortho-phosphate. <i>Physica B: Condensed Matter</i> , 1999, 266, 60-68.	2.7	68
4	Modelling frequency-dependent conductivities and permittivities in the framework of the MIGRATION concept. <i>Solid State Ionics</i> , 2004, 169, 1-8.	2.7	68
5	Characterization of the complex ion dynamics in lithium silicate glasses via computer simulations. <i>Physical Chemistry Chemical Physics</i> , 2002, 4, 3185-3192.	2.8	60
6	Correlated ionic hopping processes in crystalline and glassy electrolytes resulting in MIGRATION-type and nearly-constant-loss-type conductivities. <i>Physical Chemistry Chemical Physics</i> , 2005, 7, 157.	2.8	45
7	Frequency-dependent fluidity and conductivity of an ionic liquid. <i>Physical Chemistry Chemical Physics</i> , 2009, 11, 5930.	2.8	44
8	Anion Rotation and Cation Transport in the Rotor Phase $\hat{\Gamma}_2$ -Sodium Orthophosphate: Paddle-Wheel Mechanism Redefined in View of New Experimental Results. <i>Zeitschrift Fur Physikalische Chemie</i> , 2000, 214, .	2.8	40
9	Structure and dynamics of lithium silicate melts: molecular dynamics simulations. <i>Physical Chemistry Chemical Physics</i> , 2001, 3, 5104-5108.	2.8	40
10	Synthesis and Modeling of Polysiloxane-Based Salt-in-Polymer Electrolytes with Various Additives. <i>Journal of Physical Chemistry B</i> , 2009, 113, 15473-15484.	2.6	39
11	Low-Temperature Phases of Rubidium Silver Iodide: Crystal Structures and Dynamics of the Mobile Silver Ions. <i>Journal of Physical Chemistry A</i> , 2006, 110, 3010-3016.	2.5	36
12	Conductivity spectra and ion dynamics of a salt-in-polymer electrolyte. <i>Solid State Ionics</i> , 2006, 177, 3135-3139.	2.7	36
13	First and Second Universalities: Expeditions Towards and Beyond. <i>Zeitschrift Fur Physikalische Chemie</i> , 2010, 224, 1891-1950.	2.8	35
14	Nearly constant loss effects in borate glasses. <i>Physical Chemistry Chemical Physics</i> , 2009, 11, 3158.	2.8	31
15	New nearly constant loss feature detected in glass at low temperatures. <i>Physical Chemistry Chemical Physics</i> , 2010, 12, 14102.	2.8	30
16	Conductivity dispersion in supercooled calcium potassium nitrate: caged ionic motion viewed as part of standard behaviour. <i>Physical Chemistry Chemical Physics</i> , 2007, 9, 5582.	2.8	26
17	Translational and localised ionic motion in materials with disordered structures. <i>Solid State Sciences</i> , 2008, 10, 790-803.	3.2	26
18	Nearly constant loss effect in sodium borate and silver meta-phosphate glasses: New insights. <i>Solid State Ionics</i> , 2011, 192, 70-75.	2.7	25

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19	Conductivity spectroscopy covering 17 decades on the frequency scale. <i>Solid State Ionics</i> , 2005, 176, 1971-1978.	2.7	24
20	Insights from molecular dynamics simulations on structural organization and diffusive dynamics of an ionic liquid at solid and vacuum interfaces. <i>Journal of Colloid and Interface Science</i> , 2019, 553, 350-363.	9.4	23
21	Non-Arrhenius viscosity related to short-time ion dynamics in a fragile molten salt. <i>Physical Chemistry Chemical Physics</i> , 2005, 7, 1096.	2.8	21
22	Dielectric function and localized diffusion in ion conducting glasses. <i>Solid State Ionics</i> , 2004, 175, 661-663.	2.7	20
23	Coupling model and MIGRATION concept – Equivalence and mutual mapping. <i>Journal of Non-Crystalline Solids</i> , 2007, 353, 3845-3852.	3.1	20
24	Low-temperature $\hat{\Gamma}$ -AgI confined in glass: Structure and dynamics. <i>Solid State Ionics</i> , 2015, 271, 2-9.	2.7	20
25	Polaronic transport in iron phosphate glasses containing $\text{HfO}_2$ and $\text{CeO}_2$ . <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 3999-4009.	2.8	20
26	Insights into Ion-Network Interactions and Ion Transport in Glass. <i>Zeitschrift Fur Physikalische Chemie</i> , 2009, 223, 1201-1215.	2.8	19
27	Insights from Local Network Structures and Localized Diffusion on the Ease of Lithium Ion Transport in Two Mixed Glass-Former Systems. <i>Journal of Physical Chemistry C</i> , 2017, 121, 17641-17657.	3.1	18
28	Scaling features of conductivity spectra reveal complexities in ionic, polaronic and mixed ionic-polaronic conduction in phosphate glasses. <i>Acta Materialia</i> , 2019, 175, 46-54.	7.9	18
29	Backward correlations and dynamic heterogeneities: a computer study of ion dynamics. <i>Physical Review B</i> , 2002, 66, .	3.2	17
30	Dynamics of mobile ions in crystals, glasses and melts, described by the concept of mismatch and relaxation. <i>Solid State Ionics</i> , 2002, 154-155, 65-74.	2.7	16
31	From Ostwald's Times to Solid State Ionics: Migration and Localised Hopping of Silver Ions in Crystalline Rubidium Silver Iodide. <i>Zeitschrift Fur Physikalische Chemie</i> , 2003, 217, 1245-1264.	2.8	15
32	Using pressure, temperature and frequency as variables to study the dynamics of mobile ions in materials with disordered structures. <i>European Physical Journal: Special Topics</i> , 2008, 161, 65-78.	2.6	13
33	Anion reorientation in $\text{Na}_3\text{PO}_4$ . <i>Physica B: Condensed Matter</i> , 1997, 241-243, 338-340.	2.7	11
34	Toward understanding the second universality – A journey inspired by Arthur Stanley Nowick. <i>Journal of Electroceramics</i> , 2015, 34, 4-14.	2.0	10
35	Nearly constant loss behavior in $\text{Li}^{1/2}\text{-RbAg}_4\text{I}_5$ : microwave conductivity plateau identified. <i>Solid State Ionics</i> , 2004, 175, 819-822.	2.7	9
36	A Mechanistic Approach to Conductivity Relaxation in Ionic Glasses. <i>Zeitschrift Fur Physikalische Chemie</i> , 2004, 218, 1401-1412.	2.8	9

#	ARTICLE	IF	CITATIONS
37	Ionic transport and localized ionic motion in Na <sup>+</sup> -alumina, Na <sub>1.70</sub> Li <sub>0.32</sub> Al <sub>10.66</sub> O <sub>17</sub> . Journal of Materials Science, 2007, 42, 1942-1947.	3.7	8
38	Computational Study of Glycerol Binding within the Active Site of Coenzyme B <sub>12</sub> -Dependent Diol Dehydratase. Journal of Physical Chemistry B, 2019, 123, 6178-6187.	2.6	6
39	A theoretical calculation on Schottky defects in AgCl. Physica Status Solidi (B): Basic Research, 1991, 164, 357-367.	1.5	5
40	Role of induced quadrupoles in the simulation of intrinsic point defects in AgCl and NaCl. Physical Review B, 1993, 48, 6839-6853.	3.2	5
41	The cationic energy landscape in alkali silicate glasses: Properties and relevance. Journal of Chemical Physics, 2009, 131, 224708.	3.0	5
42	Broadband Conductivities and Fluidities of Fragile Ionic Liquids. Electrochemistry, 2009, 77, 573-581.	1.4	4
43	Exploring Reactive Conformations of Coenzyme A during Binding and Unbinding to Pyruvate Formate Lyase. Journal of Physical Chemistry A, 2019, 123, 9345-9356.	2.5	4
44	Ionic conductivity of a fragile glass-forming molten salt: Modelling its dependence on frequency, temperature, and pressure. International Journal of Materials Research, 2022, 95, 921-927.	0.3	4
45	Defect models in silver halides. Bulletin of Materials Science, 1997, 20, 435-440.	1.7	3
46	The Influence of Chemical Change on Protein Dynamics: A Case Study with Pyruvate Formate Lyase. Chemistry - A European Journal, 2019, 25, 8741-8753.	3.3	3
47	Structural characterization of an ionic liquid in bulk and in nano-confined environment using data from MD simulations. Data in Brief, 2020, 28, 104794.	1.0	3
48	Point defect modelling and transport processes in AgBr. Radiation Effects and Defects in Solids, 1995, 134, 157-159.	1.2	2
49	An Extended Polarizable Point Ion Model for Schottky Defects. Physica Status Solidi (B): Basic Research, 1991, 166, 15-23.	1.5	1
50	Majority and minority intrinsic defects in lithium and sodium halides. Bulletin of Materials Science, 1997, 20, 451-454.	1.7	1
51	A Schematic Model for Multi-particle Dynamics in Ion Transport: From Mean Field to Non-mean Field Effects. Zeitschrift Fur Physikalische Chemie, 2009, 223, 1259-1272.	2.8	0
52	The Influence of Chemical Change on Protein Dynamics: A Case Study with Pyruvate Formate Lyase. Chemistry - A European Journal, 2019, 25, 8653.	3.3	0
53	Ion Transport in Glass-Forming Calcium Potassium Nitrate: From Complex Behaviours to Unexpected Simplicities. , 2019, 22, 140-159.		0