

# Mark J Dellostritto

## List of Publications by Year in descending order

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Version: 2024-02-01

13

papers

255

citations

933447

10

h-index

1125743

13

g-index

13

all docs

13

docs citations

13

times ranked

406

citing authors

#	ARTICLE	IF	CITATIONS
1	Importance of nuclear quantum effects on the hydration of chloride ion. <i>Physical Review Materials</i> , 2021, 5, .	2.4	11
2	Modeling the Ultrafast Electron Attachment Dynamics of Solvated Uracil. <i>Journal of Physical Chemistry A</i> , 2021, 125, 6995-7003.	2.5	8
3	Probing Heterogeneous Charge Distributions at the $\text{Al}_2\text{O}_3(0001)/\text{H}_2\text{O}$ Interface. <i>Journal of the American Chemical Society</i> , 2020, 142, 12096-12105.	13.7	21
4	Aqueous solvation of the chloride ion revisited with density functional theory: impact of correlation and exchange approximations. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 10666-10675.	2.8	20
5	Bond-Dependent Thole Model for Polarizability and Spectroscopy. <i>Journal of Physical Chemistry A</i> , 2019, 123, 5378-5387.	2.5	6
6	Effect of Interlayer $\text{Co}^{2+}$ on Structure and Charge Transfer in NiFe Layered Double Hydroxides. <i>Journal of Physical Chemistry C</i> , 2019, 123, 13593-13599.	3.1	11
7	Sodium Halide Adsorption and Water Structure at the $\text{Al}_2\text{O}_3(0001)/\text{Water}$ Interface. <i>Journal of Physical Chemistry C</i> , 2019, 123, 15618-15628.	3.1	19
8	Effect of Functional and Electron Correlation on the Structure and Spectroscopy of the $\text{Al}_2\text{O}_3(001)\text{-H}_2\text{O}$ Interface. <i>Journal of Physical Chemistry Letters</i> , 2019, 10, 2031-2036.	4.6	22
9	Relating Interfacial Order to Sum Frequency Generation with Ab Initio Simulations of the Aqueous $\text{Al}_2\text{O}_3(0001)$ and $(112\bar{1}0)$ Interfaces. <i>Journal of Physical Chemistry C</i> , 2018, 122, 21284-21294.	3.1	30
10	Bond Polarizability Model for Sum Frequency Generation at the $\text{Al}_2\text{O}_3(0001)\text{-H}_2\text{O}$ Interface. <i>Journal of Physical Chemistry A</i> , 2017, 121, 3045-3055.	2.5	16
11	Effect of Ions on H-Bond Structure and Dynamics at the Quartz(101)-Water Interface. <i>Langmuir</i> , 2016, 32, 11353-11365.	3.5	41
12	Vibrational Density of States of Strongly H-Bonded Interfacial Water: Insights from Inelastic Neutron Scattering and Theory. <i>Journal of Physical Chemistry C</i> , 2014, 118, 10805-10813.	3.1	48
13	Tan $\frac{1}{4}s$ universal contact and collective oscillations of strongly interacting Fermi gases. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2012, 376, 2298-2305.	2.1	2