

Sule Atahan-Evrenk

List of Publications by Year in descending order

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18
papers

2,613
citations

567281

15
h-index

839539

18
g-index

19
all docs

19
docs citations

19
times ranked

4200
citing authors

#	ARTICLE	IF	CITATIONS
1	Tuning charge transport in solution-sheared organic semiconductors using lattice strain. <i>Nature</i> , 2011, 480, 504-508.	27.8	981
2	The Harvard Clean Energy Project: Large-Scale Computational Screening and Design of Organic Photovoltaics on the World Community Grid. <i>Journal of Physical Chemistry Letters</i> , 2011, 2, 2241-2251.	4.6	470
3	From computational discovery to experimental characterization of a high hole mobility organic crystal. <i>Nature Communications</i> , 2011, 2, 437.	12.8	321
4	Lead candidates for high-performance organic photovoltaics from high-throughput quantum chemistry – the Harvard Clean Energy Project. <i>Energy and Environmental Science</i> , 2014, 7, 698-704.	30.8	189
5	Accelerated computational discovery of high-performance materials for organic photovoltaics by means of cheminformatics. <i>Energy and Environmental Science</i> , 2011, 4, 4849.	30.8	169
6	Multiple coherent states for first-principles semiclassical initial value representation molecular dynamics. <i>Journal of Chemical Physics</i> , 2009, 130, 234113.	3.0	82
7	Effects of Odd – Even Side Chain Length of Alkyl-Substituted Diphenylbithiophenes on First Monolayer Thin Film Packing Structure. <i>Journal of the American Chemical Society</i> , 2013, 135, 11006-11014.	13.7	81
8	Theoretical Characterization of the Air-Stable, High-Mobility Dinaphtho[2,3- <i>b</i> :2'- <i>f</i>]thieno[3,2- <i>b</i>]-thiophene Organic Semiconductor. <i>Journal of Physical Chemistry C</i> , 2010, 114, 2334-2340.	3.1	73
9	First-principles semiclassical initial value representation molecular dynamics. <i>Physical Chemistry Chemical Physics</i> , 2009, 11, 3861.	2.8	70
10	Prediction of Intramolecular Reorganization Energy Using Machine Learning. <i>Journal of Physical Chemistry A</i> , 2019, 123, 7855-7863.	2.5	36
11	Prediction and Theoretical Characterization of p-Type Organic Semiconductor Crystals for Field-Effect Transistor Applications. <i>Topics in Current Chemistry</i> , 2014, 345, 95-138.	4.0	30
12	Coherent Dynamics of Mixed Frenkel and Charge-Transfer Excitons in Dinaphtho[2,3- <i>b</i> :2'- <i>f</i>]thieno[3,2- <i>b</i>]-thiophene Thin Films: The Importance of Hole Delocalization. <i>Journal of Physical Chemistry Letters</i> , 2016, 7, 1374-1380.	4.6	24
13	Coupled-States Statistical Investigation of Vibrational and Rotational Relaxation of OH(2 $\hat{1}$) by Collisions with Atomic Hydrogen. <i>Journal of Physical Chemistry A</i> , 2006, 110, 5436-5445.	2.5	17
14	An ab initio investigation of the O(3P) – H ₂ (1 $\hat{1}$ g) van der Waals well. <i>Physical Chemistry Chemical Physics</i> , 2006, 8, 4420-4426.	2.8	17
15	Confined organization of fullerene units along high polymer chains. <i>Journal of Materials Chemistry C</i> , 2013, 1, 5747.	5.5	16
16	Cross sections and thermal rate constants for the isotope exchange reaction: D(S ₂) + OH(1 $\hat{2}$) – OD(1 $\hat{2}$) + H(S ₂). <i>Journal of Chemical Physics</i> , 2005, 123, 204306.	3.0	15
17	A quantitative structure – property study of reorganization energy for known p-type organic semiconductors. <i>RSC Advances</i> , 2018, 8, 40330-40337.	3.6	13
18	Laser spectroscopic study of the SiAr van der Waals complex. <i>Journal of Chemical Physics</i> , 2002, 116, 9239-9248.	3.0	9