Dylan M Anstine

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	A molecular dynamics study of water-soluble polymers: analysis of force fields from atomistic simulations. Molecular Simulation, 2019, 45, 310-321.	2.0	20
2	Incorporating Flexibility Effects into Metal–Organic Framework Adsorption Simulations Using Different Models. ACS Applied Materials & Interfaces, 2021, 13, 61305-61315.	8.0	17
3	Adsorption space for microporous polymers with diverse adsorbate species. Npj Computational Materials, 2021, 7, .	8.7	16
4	Screening PIM-1 performance as a membrane for binary mixture separation of gaseous organic compounds. Journal of Membrane Science, 2020, 599, 117798.	8.2	13
5	Sorptionâ€induced polymer rearrangement: approaches from molecular modeling. Polymer International, 2020, 70, 984.	3.1	13
6	Effects of an atomistic modeling approach on predicted mechanical properties of glassy polymers via molecular dynamics. Modelling and Simulation in Materials Science and Engineering, 2020, 28, 025006.	2.0	9
7	PECDA hydrogel structure from semi-dilute concentrations: insights from experiments and molecular simulations. Soft Matter, 2022, 18, 3565-3574.	2.7	9
8	An Insight into Structural and Mechanical Properties of Idealâ€Networked Poly(Ethylene) Tj ETQq0 0 0 rgBT /Ove Physics, 2020, 221, 1900326.	rlock 10 T 2.2	50 467 Td 7
9	In silico design of microporous polymers for chemical separations and storage. Current Opinion in Chemical Engineering, 2022, 36, 100795.	7.8	6
10	Sulfonyl <scp>PIM</scp> â€l: A diverse separation membrane with dilation resistance. AICHE Journal, 2021, 67, e17006.	3.6	4
11	Attosecond time delay in the valence photoionization of C ₂₄₀ versus C ₆₀ . Journal of Physics: Conference Series, 2015, 635, 112025.	0.4	0