

Paweł, Oracz

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Predicted Mutual Solubilities in Water + C5-C12 Hydrocarbon Systems. Results at 298 K. ChemEngineering, 2021, 5, 89.	2.4	1
2	IUPAC-NIST Solubility Data Series. 96. Amines with Water Part 1. C4-C6 Aliphatic Amines. Journal of Physical and Chemical Reference Data, 2012, 41, .	4.2	10
3	IUPAC-NIST Solubility Data Series. 96. Amines with Water Part 2. C7-C24 Aliphatic Amines. Journal of Physical and Chemical Reference Data, 2012, 41, .	4.2	10
4	IUPAC-NIST Solubility Data Series. 96. Amines with Water Part 3. Non-Aliphatic Amines. Journal of Physical and Chemical Reference Data, 2012, 41, .	4.2	4
5	Vapor Pressures and Vaporization Enthalpy of 2-Ethoxyphenol. Journal of Chemical & Engineering Data, 2012, 57, 3176-3179.	1.9	2
6	Application of the Unified Functional Activity Coefficient (UNIFAC) and Analytical Solution of Groups (ASOG) for the Calculation of Mutual Solubilities in Water Systems of Alkanes, Arenes, and Alkanols. Journal of Chemical & Engineering Data, 2011, 56, 4853-4861.	1.9	10
7	Vapor-Liquid Equilibria in the 2-Methylcyclohexanol + 2-Methylcyclohexyl Ethanoate System at 101.325 kPa. Journal of Chemical & Engineering Data, 2007, 52, 645-649.	1.9	2
8	Recommended Liquid-Liquid Equilibrium Data. Part 5. Ether-Water Systems. Journal of Physical and Chemical Reference Data, 2007, 36, 1399-1415.	4.2	10
9	Recommended Vapor-Liquid Equilibrium Data. Part 4. Binary Alkanol-Alkene/Alkyne Systems. Journal of Physical and Chemical Reference Data, 2006, 35, 1577-1596.	4.2	2
10	Recommended Vapor-Liquid Equilibrium Data. Part 3. Binary Alkanol-Aromatic Hydrocarbon Systems. Journal of Physical and Chemical Reference Data, 2004, 33, 959-997.	4.2	16
11	True reactivity ratios for styrene-methyl methacrylate copolymerization in bulk. Macromolecular Theory and Simulations, 1998, 7, 97-103.	1.4	3
12	Non-uniqueness in determination of terminal and penultimate model reactivity ratios in the styrene-methyl methacrylate free-radical copolymerization system. Macromolecular Theory and Simulations, 1997, 6, 565-576.	1.4	21