

# Paweł, Oracz

## List of Publications by Year in descending order

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12  
papers

91  
citations

1478505

6  
h-index

1372567

10  
g-index

12  
all docs

12  
docs citations

12  
times ranked

91  
citing authors

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