

Richard J Sadus

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140
papers

2,595
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26
h-index

42
g-index

151
ext. papers

2,806
ext. citations

3.2
avg, IF

5.58
L-index

#	Paper	IF	Citations
140	Equations of state for the calculation of fluid-phase equilibria. <i>AIChE Journal</i> , 2000 , 46, 169-196	3.6	298
139	Molecular simulation of the phase behavior of noble gases using accurate two-body and three-body intermolecular potentials. <i>Journal of Chemical Physics</i> , 1999 , 111, 1533-1540	3.9	90
138	High-pressure phase equilibria and supercritical pVT data of the binary water + methane mixture to 723 K and 200 MPa. <i>The Journal of Physical Chemistry</i> , 1993 , 97, 9054-9059		73
137	Calculating critical transitions of fluid mixtures: Theory vs. experiment. <i>AIChE Journal</i> , 1994 , 40, 1376-1403	3.6	72
136	Molecular dynamics simulation of the dielectric constant of water: the effect of bond flexibility. <i>Journal of Chemical Physics</i> , 2011 , 134, 234501	3.9	70
135	Solid-liquid equilibria and triple points of n-6 Lennard-Jones fluids. <i>Journal of Chemical Physics</i> , 2009 , 131, 174504	3.9	69
134	Viscoelastic properties of dendrimers in the melt from nonequilibrium molecular dynamics. <i>Journal of Chemical Physics</i> , 2004 , 121, 12050-9	3.9	51
133	Atomistic water models: Aqueous thermodynamic properties from ambient to supercritical conditions. <i>Fluid Phase Equilibria</i> , 2016 , 407, 7-30	2.5	48
132	Molecular dynamics simulation of the effect of bond flexibility on the transport properties of water. <i>Journal of Chemical Physics</i> , 2012 , 137, 104512	3.9	47
131	Hard sphere compressibility factors for equation of state development. <i>AIChE Journal</i> , 2005 , 51, 309-313	3.6	45
130	Thermodynamics of Fluid Polyamorphism. <i>Physical Review X</i> , 2018 , 8,	9.1	43
129	Internal structure of dendrimers in the melt under shear: a molecular dynamics study. <i>Journal of Chemical Physics</i> , 2004 , 121, 1091-6	3.9	41
128	Molecular simulation of the vapor-liquid coexistence of mercury. <i>Journal of Chemical Physics</i> , 2003 , 119, 6691-6697	3.9	41
127	Phase diagram of the Weeks-Chandler-Andersen potential from very low to high temperatures and pressures. <i>Physical Review E</i> , 2009 , 80, 061101	2.4	40
126	Influence of bond flexibility on the vapor-liquid phase equilibria of water. <i>Journal of Chemical Physics</i> , 2007 , 126, 044701	3.9	40
125	A link between the two-body and three-body interaction energies of fluids from molecular simulation. <i>Journal of Chemical Physics</i> , 2000 , 112, 6382-6385	3.9	38
124	Molecular simulation of the vapour-liquid phase coexistence of neon and argon using ab initio potentials. <i>Physical Chemistry Chemical Physics</i> , 2001 , 3, 1297-1302	3.6	38

123	A molecular dynamics study of nitric oxide in water: diffusion and structure. <i>Journal of Chemical Physics</i> , 2005 , 123, 054505	3.9	35
122	Relationships between three-body and two-body interactions in fluids and solids. <i>Journal of Chemical Physics</i> , 2006 , 125, 144509	3.9	35
121	Equations of state for fluids: The Dieterici approach revisited. <i>Journal of Chemical Physics</i> , 2001 , 115, 1460-1462	3.9	34
120	Three-body interactions in fluids from molecular simulation: Vapor-liquid phase coexistence of argon. <i>Journal of Chemical Physics</i> , 1996 , 104, 4784-4787	3.9	34
119	Exact calculation of the effect of three-body Axilrod-Teller interactions on vapour-liquid phase coexistence. <i>Fluid Phase Equilibria</i> , 1998 , 144, 351-359	2.5	33
118	Molecular simulation of dendrimers and their mixtures under shear: comparison of isothermal-isobaric (NpT) and isothermal-isochoric (NVT) ensemble systems. <i>Journal of Chemical Physics</i> , 2005 , 123, 34905	3.9	33
117	Role of nonadditive forces on the structure and properties of liquid water. <i>Journal of Chemical Physics</i> , 2007 , 127, 154509	3.9	31
116	Analytic dependence of the pressure and energy of an atomic fluid under shear. <i>Physical Review E</i> , 2001 , 63, 021204	2.4	31
115	Molecular simulation of the vapour-liquid equilibria of pure fluids and binary mixtures containing dipolar components: the effect of Keesom interactions. <i>Molecular Physics</i> , 1996 , 87, 979-990	1.7	27
114	Effect of potential truncations and shifts on the solid-liquid phase coexistence of Lennard-Jones fluids. <i>Journal of Chemical Physics</i> , 2010 , 133, 124515	3.9	26
113	High-Pressure Phase Equilibria and Critical Curve of the Water + Helium System to 200 MPa and 723 K. <i>The Journal of Physical Chemistry</i> , 1995 , 99, 4273-4277		25
112	Equations of State for Hard-Sphere Chains. <i>The Journal of Physical Chemistry</i> , 1995 , 99, 12363-12366		25
111	Analysis of the shape of dendrimers under shear. <i>Journal of Chemical Physics</i> , 2006 , 124, 044910	3.9	24
110	Kinetics and chemomechanical properties of the F1-ATPase molecular motor. <i>Journal of Chemical Physics</i> , 2003 , 118, 9890-9898	3.9	24
109	Intermolecular interactions and the thermodynamic properties of supercritical fluids. <i>Journal of Chemical Physics</i> , 2013 , 138, 194502	3.9	23
108	The Dieterici alternative to the van der Waals approach for equations of state: second virial coefficients. <i>Physical Chemistry Chemical Physics</i> , 2002 , 4, 919-921	3.6	23
107	Closed-loop liquid-liquid equilibria and the global phase behaviour of binary mixtures involving hard-sphere + van der Waals interactions. <i>Molecular Physics</i> , 2000 , 98, 715-723	1.7	22
106	On the relationship between two-body and three-body interactions from nonequilibrium molecular dynamics simulation. <i>Journal of Chemical Physics</i> , 2001 , 115, 9410-9413	3.9	22

105	Second virial coefficient properties of the - Lennard-Jones/Mie potential. <i>Journal of Chemical Physics</i> , 2018 , 149, 074504	3.9	22
104	Thermodynamic properties of supercritical n-m Lennard-Jones fluids and isochoric and isobaric heat capacity maxima and minima. <i>Journal of Chemical Physics</i> , 2013 , 139, 154503	3.9	21
103	Thermodynamic properties in the molecular dynamics ensemble applied to the gaussian core model fluid. <i>Journal of Chemical Physics</i> , 2011 , 134, 114515	3.9	21
102	Equilibrium and nonequilibrium molecular dynamics methods for determining solid-liquid phase coexistence at equilibrium. <i>Journal of Chemical Physics</i> , 2003 , 119, 11017-11023	3.9	21
101	Scaling behavior for the pressure and energy of shearing fluids. <i>Physical Review E</i> , 2003 , 67, 061201	2.4	20
100	Molecular simulation of the thermodynamic, structural, and vapor-liquid equilibrium properties of neon. <i>Journal of Chemical Physics</i> , 2016 , 145, 104501	3.9	20
99	Effect of three-body interactions on the vapor-liquid phase equilibria of binary fluid mixtures. <i>Journal of Chemical Physics</i> , 2006 , 125, 074503	3.9	19
98	Vapour-liquid and liquid-liquid phase equilibria of binary mixtures containing helium: comparison of experiment with predictions using equations of state. <i>Fluid Phase Equilibria</i> , 1996 , 122, 1-15	2.5	19
97	Application of hard convex body and hard sphere equations of state to the critical properties of binary mixtures. <i>Fluid Phase Equilibria</i> , 1988 , 39, 89-99	2.5	19
96	Phase behaviour of fluorocarbon and hydrocarbon mixtures: interpretation of type II and type III behaviour in terms of a hard-sphere + attractive force equation of state. <i>Fluid Phase Equilibria</i> , 1986 , 25, 263-272	2.5	18
95	Critical properties of ternary mixtures: hydrocarbon, acetone and alkanenitrile mixtures. <i>Chemical Engineering Science</i> , 1987 , 42, 1717-1722	4.4	18
94	Phase behaviour of binary fluid mixtures: a global phase diagram solely in terms of pure component properties. <i>Fluid Phase Equilibria</i> , 2003 , 214, 67-78	2.5	17
93	Two-body interatomic potentials for He, Ne, Ar, Kr, and Xe from ab initio data. <i>Journal of Chemical Physics</i> , 2019 , 150, 134504	3.9	16
92	Thermodynamic properties and diffusion of water + methane binary mixtures. <i>Journal of Chemical Physics</i> , 2014 , 140, 104505	3.9	16
91	Molecular Simulation of Henry's Constant at Vapor-Liquid and Liquid-Liquid Phase Boundaries. <i>Journal of Physical Chemistry B</i> , 1997 , 101, 3834-3838	3.4	16
90	Thermophysical properties of supercritical water and bond flexibility. <i>Physical Review E</i> , 2015 , 92, 012124	4.4	15
89	Thermodynamic properties of liquid water from a polarizable intermolecular potential. <i>Journal of Chemical Physics</i> , 2013 , 138, 044503	3.9	15
88	Solid-liquid phase equilibria of the Gaussian core model fluid. <i>Journal of Chemical Physics</i> , 2009 , 131, 184507	3.9	14

87	Three-body interactions and solid-liquid phase equilibria: application of a molecular dynamics algorithm. <i>Physical Review E</i> , 2006 , 74, 031203	2.4	14
86	Energy and pressure of shearing fluids at different state points. <i>Physical Review E</i> , 2001 , 64, 021201	2.4	14
85	Influence of combining rules and molecular shape on the high pressure phase equilibria of binary fluid mixtures. <i>The Journal of Physical Chemistry</i> , 1993 , 97, 1985-1992		14
84	Deiters' equation of state and critical phenomena. <i>Chemical Engineering Science</i> , 1988 , 43, 459-466	4.4	14
83	Molecular simulation of intermolecular attraction and repulsion in coexisting liquid and vapour phases. <i>Fluid Phase Equilibria</i> , 1997 , 134, 77-85	2.5	13
82	Molecular simulation of fluids with non-identical intermolecular potentials: thermodynamic properties of 10-5 + 12-6 Mie potential binary mixtures. <i>Journal of Chemical Physics</i> , 2015 , 142, 084504	3.9	12
81	Intermolecular potentials and the accurate prediction of the thermodynamic properties of water. <i>Journal of Chemical Physics</i> , 2013 , 139, 194505	3.9	12
80	Allosteric Conformational Transition in Adenylate Kinase: Dynamic Correlations and Implication for Allostery. <i>Australian Journal of Chemistry</i> , 2010 , 63, 405	1.2	12
79	Structure and polarization properties of water: molecular dynamics with a nonadditive intermolecular potential. <i>Physical Review E</i> , 2012 , 85, 051509	2.4	12
78	New Dieterici-type equations of state for fluid phase equilibria. <i>Fluid Phase Equilibria</i> , 2003 , 212, 31-39	2.5	12
77	Simple equation of state for hard-sphere chains. <i>AIChE Journal</i> , 1999 , 45, 2454-2457	3.6	12
76	RESEARCH NOTE Molecular simulation of the liquid-liquid equilibria of binary mixtures containing dipolar and non-polar components interacting via the Keesom potential. <i>Molecular Physics</i> , 1996 , 89, 1187-1194	1.7	12
75	Influence of quantum effects on the high-pressure phase behavior of binary mixtures containing hydrogen. <i>The Journal of Physical Chemistry</i> , 1992 , 96, 3855-3860		12
74	Ab initio interatomic potentials and the thermodynamic properties of fluids. <i>Journal of Chemical Physics</i> , 2017 , 147, 024505	3.9	11
73	Predicting vapor-liquid phase equilibria with augmented ab initio interatomic potentials. <i>Journal of Chemical Physics</i> , 2017 , 146, 244504	3.9	11
72	The effect of three-body interactions on the liquid-liquid phase coexistence of binary fluid mixtures. <i>Fluid Phase Equilibria</i> , 1998 , 150-151, 63-72	2.5	11
71	Coarse-grained dynamics of the receiver domain of NtrC: fluctuations, correlations and implications for allosteric cooperativity. <i>Proteins: Structure, Function and Bioinformatics</i> , 2008 , 73, 218-27	4.2	11
70	Molecular simulation of the high-pressure phase equilibria of binary atomic fluid mixtures using the exponential-6 intermolecular potential. <i>Fluid Phase Equilibria</i> , 2000 , 170, 269-284	2.5	11

69	Predicting the Gas-Liquid Critical Properties of Binary Mixtures: An Alternative to Conventional Mixing Rules. <i>Zeitschrift Fur Elektrotechnik Und Elektrochemie</i> , 1992 , 96, 1454-1459		11
68	SURVEY OF THE HIGH PRESSURE PHASE BEHAVIOUR OF FLUIDS 1992 , 1-12		11
67	Attractive forces in binary mixtures: a priori prediction of gas-liquid critical properties. <i>The Journal of Physical Chemistry</i> , 1989 , 93, 3787-3790		11
66	Critical properties of ternary mixtures: Siloxane and perfluoromethylcyclohexane mixtures. <i>Chemical Engineering Science</i> , 1988 , 43, 883-885	4.4	11
65	The Widom Line and the Lennard-Jones Potential. <i>Journal of Physical Chemistry B</i> , 2019 , 123, 8268-8273	3.4	10
64	Flow of water through carbon nanotubes predicted by different atomistic water models. <i>Journal of Chemical Physics</i> , 2019 , 150, 194501	3.9	10
63	Strain-rate dependent shear viscosity of the Gaussian core model fluid. <i>Journal of Chemical Physics</i> , 2009 , 131, 224511	3.9	10
62	Dynamic and coordinating domain motions in the active subunits of the F1-ATPase molecular motor. <i>Biochimica Et Biophysica Acta - Proteins and Proteomics</i> , 2006 , 1764, 1553-60	4	10
61	The strain rate dependence of shear viscosity, pressure and energy from two-body and three-body interactions. <i>Fluid Phase Equilibria</i> , 2001 , 183-184, 371-379	2.5	10
60	Molecular simulation of the phase behaviour of ternary fluid mixtures: the effect of a third component on vapour-liquid and liquid-liquid coexistence. <i>Fluid Phase Equilibria</i> , 1999 , 157, 169-180	2.5	10
59	Molecular simulation of the shear viscosity and the self-diffusion coefficient of mercury along the vapor-liquid coexistence curve. <i>Journal of Chemical Physics</i> , 2005 , 123, 34511	3.9	9
58	Novel critical transitions in ternary fluid mixtures. <i>The Journal of Physical Chemistry</i> , 1992 , 96, 5197-5202		9
57	Intermolecular Potential-Based Equations of State from Molecular Simulation and Second Virial Coefficient Properties. <i>Journal of Physical Chemistry B</i> , 2018 , 122, 7757-7763	3.4	8
56	A suite of domain-specific visual languages for scientific software application modelling 2013 ,		8
55	Complex cooperativity of ATP hydrolysis in the F(1)-ATPase molecular motor. <i>Biochimica Et Biophysica Acta - Proteins and Proteomics</i> , 2004 , 1698, 197-202	4	8
54	Cooperativity in the motor activities of the ATP-fueled molecular motors. <i>Biochimica Et Biophysica Acta - Proteins and Proteomics</i> , 2005 , 1752, 111-23	4	8
53	Two-body intermolecular potentials from second virial coefficient properties. <i>Journal of Chemical Physics</i> , 2019 , 150, 024503	3.9	7
52	Fully a priori prediction of the vapor-liquid equilibria of Ar, Kr, and Xe from ab initio two-body plus three-body interatomic potentials. <i>Journal of Chemical Physics</i> , 2019 , 151, 034509	3.9	7

51	Dielectric and structural properties of aqueous nonpolar solute mixtures. <i>Journal of Chemical Physics</i> , 2012 , 137, 124501	3.9	7
50	Pressure and energy behavior of the Gaussian core model fluid under shear. <i>Physical Review E</i> , 2010 , 82, 011201	2.4	7
49	Effect of Three-Body Interactions between Dissimilar Molecules on the Phase Behavior of Binary Mixtures: The Transition from Vapor-Liquid Equilibria to Type III Behavior. <i>Industrial & Engineering Chemistry Research</i> , 1998 , 37, 2977-2982	3.9	7
48	A Simplified Thermodynamic Perturbation Theory-Dimer Equation of State for Mixtures of Hard-Sphere Chains. <i>Macromolecules</i> , 1996 , 29, 7212-7216	5.5	7
47	Phase behaviour of 1-alkanol + alkane mixtures: gas-liquid critical temperatures. <i>Fluid Phase Equilibria</i> , 1991 , 67, 259-271	2.5	7
46	Novel high pressure critical phase transitions in multicomponent fluid mixtures. <i>Fluid Phase Equilibria</i> , 1993 , 83, 101-108	2.5	7
45	Gas-liquid critical properties of binary mixtures of n-alkanes and 2,2,4-trimethylpentane with the weakly polar halocarbons 1,2-dichloroethane, cis-1,2-dichloroethene trans-1,2-dichloroethene and tetrachloromethane. <i>Industrial & Engineering Chemistry Research</i> , 1989 , 28, 481-484	3.9	7
44	Ab Initio Interatomic Potentials and the Classical Molecular Simulation Prediction of the Thermophysical Properties of Helium. <i>Journal of Physical Chemistry B</i> , 2020 , 124, 2268-2276	3.4	6
43	Mg ²⁺ coordinating dynamics in Mg:ATP fueled motor proteins. <i>Journal of Chemical Physics</i> , 2014 , 140, 115102	3.9	6
42	Modified force decomposition algorithms for calculating three-body interactions via molecular dynamics. <i>Computer Physics Communications</i> , 2006 , 175, 683-691	4.2	6
41	Liquid-crystal behavior of hard ellipsoid dimers. <i>Journal of Chemical Physics</i> , 2001 , 114, 5432-5434	3.9	6
40	Phase behaviour of carbon dioxide-benzene-water ternary mixtures at high pressures and temperatures up to 300 MPa and 600 K. <i>Physical Chemistry Chemical Physics</i> , 2000 , 2, 4157-4164	3.6	6
39	Phase behaviour of ternary mixtures: a theoretical investigation of the critical properties of mixtures with equal size components. <i>Physical Chemistry Chemical Physics</i> , 1999 , 1, 4329-4336	3.6	6
38	Conformational dynamics of ATP/Mg:ATP in motor proteins via data mining and molecular simulation. <i>Journal of Chemical Physics</i> , 2012 , 137, 075101	3.9	5
37	Parallel algorithms for molecular dynamics with induction forces. <i>Computer Physics Communications</i> , 2008 , 178, 384-392	4.2	5
36	Influence of two-body and three-body interatomic forces on gas, liquid, and solid phases. <i>Physical Review E</i> , 2006 , 74, 021202	2.4	5
35	Response to [Comment on Equations of state for fluids: The Dieterici approach revisited]. <i>J. Chem. Phys.</i> 115, 1460 (2001)]. <i>Journal of Chemical Physics</i> , 2004 , 120, 8870-8870	3.9	5
34	New phase for one-component hard spheres. <i>Journal of Chemical Physics</i> , 2004 , 120, 11686-91	3.9	5

33	Three-body interactions and the phase equilibria of mixtures. <i>High Temperatures - High Pressures</i> , 2001 , 33, 111-118	1.3	5
32	Molecular simulation of the vapour-liquid equilibria of pure fluids and binary mixtures containing dipolar components: the effect of Keesom interactions		5
31	Structural behavior of fluids from the vapor and liquid region to the supercritical phase. <i>Physical Review E</i> , 2019 , 100, 052132	2.4	5
30	Nonequilibrium equation of state for Lennard-Jones fluids and the calculation of strain-rate dependent shear viscosity. <i>AIChE Journal</i> , 2011 , 57, 250-258	3.6	4
29	Molecular simulation of the thermophysical properties of fluids: phase behaviour and transport properties. <i>Molecular Simulation</i> , 2006 , 32, 185-189	2	4
28	Prediction of one-component vapour-liquid equilibria from the triple point to the critical point using a simplified perturbed hard-chain theory equation of state. <i>Fluid Phase Equilibria</i> , 1995 , 109, 171-182	2.5	4
27	Monte Carlo simulation of vapour-liquid equilibria in Lennard-Jones + three-body potential binary fluid mixtures. <i>Fluid Phase Equilibria</i> , 1996 , 116, 289-295	2.5	4
26	Calculation of the critical high pressure liquid-liquid phase equilibria of binary mixtures containing ammonia Unlike interaction parameters for the HeiligBranck equation of state. <i>Fluid Phase Equilibria</i> , 1994 , 101, 89-99	2.5	4
25	Prediction of binary and ternary critical properties using deiters' equation, hard sphere and hard convex body equations of state. <i>Fluid Phase Equilibria</i> , 1988 , 42, 85-103	2.5	4
24	Molecular simulation of orthobaric isochoric heat capacities near the critical point. <i>Physical Review E</i> , 2019 , 99, 012139	2.4	3
23	Thermodynamic properties and anomalous behavior of double-Gaussian core model potential fluids. <i>Physical Review E</i> , 2019 , 100, 012112	2.4	3
22	Molecular simulation of the phase behavior of fluids and fluid mixtures using the synthetic method. <i>Journal of Chemical Physics</i> , 2012 , 137, 054507	3.9	3
21	Molecular simulation and theory for nanosystems: Insights for molecular motors. <i>Molecular Simulation</i> , 2008 , 34, 23-27	2	3
20	Influence of molecular shape on vapour-liquid equilibria. <i>Fluid Phase Equilibria</i> , 1994 , 95, 371-381	2.5	3
19	Combining intermolecular potentials for the prediction of fluid properties: Two-body and three-body interactions. <i>Journal of Chemical Physics</i> , 2020 , 153, 214509	3.9	3
18	Beyond Traditional Effective Intermolecular Potentials and Pairwise Interactions in Molecular Simulation. <i>Lecture Notes in Computer Science</i> , 2002 , 932-941	0.9	3
17	Global phase diagram for anisotropic binary fluid mixtures: reverse type IV behaviour. <i>Molecular Physics</i> , 2003 , 101, 2211-2217	1.7	2
16	Molecular simulation of liquid-crystal transitions in hard prolate ellipsoid monomers and dimers. <i>Fluid Phase Equilibria</i> , 2002 , 194-197, 227-231	2.5	2

15	An equation of state for hard convex body chains. <i>Molecular Physics</i> , 1999 , 97, 1279-1284	1.7	2
14	Calculating one-component vapour-liquid equilibria using equations of state: a new temperature-dependence for the attractive term in the Christoforakos-Franck equation. <i>Fluid Phase Equilibria</i> , 1995 , 103, 41-50	2.5	2
13	MODELS FOR HIGH PRESSURE MULTICOMPONENT EQUILIBRIA 1992 , 25-68		2
12	Unusual critical temperature phenomena in fluid mixtures containing perfluorotributylamine. <i>Fluid Phase Equilibria</i> , 1992 , 77, 269-283	2.5	2
11	Parallelization Algorithms for Three-Body Interactions in Molecular Dynamics Simulation. <i>Lecture Notes in Computer Science</i> , 2006 , 374-382	0.9	1
10	Binary mixtures of water + five noble gases: comparison of binodal and critical curves at high pressures. <i>Fluid Phase Equilibria</i> , 1996 , 123, 1-15	2.5	1
9	Vapor-liquid equilibria and cohesive r interactions. <i>Journal of Chemical Physics</i> , 2020 , 153, 204504	3.9	1
8	Effect of the range of particle cohesion on the phase behavior and thermodynamic properties of fluids. <i>Journal of Chemical Physics</i> , 2020 , 153, 244502	3.9	1
7	Interatomic Interactions Responsible for the Solid-Liquid and Vapor-Liquid Phase Equilibria of Neon. <i>Journal of Physical Chemistry B</i> , 2021 , 125, 8522-8531	3.4	1
6	A mechanochemical theory for the ATP-fuelled biomolecular motors. <i>International Journal of Nanotechnology</i> , 2009 , 6, 1121	1.5	
5	Response to [Comment on: New phase for one-component hard spheres][J. Chem. Phys. 120, 11686 (2004)]. <i>Journal of Chemical Physics</i> , 2004 , 121, 12117-12118	3.9	
4	PROCEDURE FOR DETERMINING THE PHASE BEHAVIOUR OF TERNARY MIXTURES: COMPARISON OF EXPERIMENT WITH THEORY FOR GAS-LIQUID CRITICAL TEMPERATURES 1992 , 69-85		
3	NOVEL HIGHER ORDER CRITICAL PHENOMENA IN TERNARY MIXTURES 1992 , 87-102		
2	TERNARY TETRAFLUOROMETHANE AND n-ALKANE MIXTURES 1992 , 177-204		
1	TERNARY MIXTURES CONTAINING WATER AND n-ALKANES 1992 , 229-255		