

Tetrahydrofuran

Other names:	1,4-epoxybutane Butane «alpha», «delta»-oxide Butane Â«alphaÂ», Â«deltaÂ»-oxide Butane, 1,4-epoxy- Cyclotetramethylene oxide Diethylene oxide Dynasolve 150 Furan, tetrahydro- Furanidine Hydrofuran NCI-C60560 NSC 57858 Oxacyclopentane Oxolane Rcra waste number U213 THF THF (tetrahydrofuran) Tetrahydrofuraan Tetrahydrofurane Tetrahydrofuranne Tetraidrofurano Tetramethylene oxide UN 2056
Inchi:	InChI=1S/C4H8O/c1-2-4-5-3-1/h1-4H2
InchiKey:	WYURNTSHIVDZCO-UHFFFAOYSA-N
Formula:	C4H8O
SMILES:	C1CCOC1
Mol. weight [g/mol]:	72.11
CAS:	109-99-9

Physical Properties

Property code	Value	Unit	Source
affp	822.10	kJ/mol	NIST Webbook
basg	794.70	kJ/mol	NIST Webbook
chg	-2533.20 ± 0.67	kJ/mol	NIST Webbook
chl	-2505.80 ± 2.10	kJ/mol	NIST Webbook

chl	-2501.20 ± 0.84	kJ/mol	NIST Webbook
dvisc	0.0004750	Paxs	A volumetric and viscosity study for the binary mixtures of 1-hexyl-3-methylimidazolium tetrafluoroborate with some molecular solvents
gf	-59.06	kJ/mol	
hf	-184.20 ± 0.71	kJ/mol	NIST Webbook
hfus	6.96	kJ/mol	Joback Method
hvap	32.00	kJ/mol	NIST Webbook
hvap	32.90	kJ/mol	NIST Webbook
hvap	32.00	kJ/mol	NIST Webbook
hvap	32.16	kJ/mol	NIST Webbook
ie	9.38 ± 0.05	eV	NIST Webbook
ie	9.38	eV	NIST Webbook
ie	9.41	eV	NIST Webbook
ie	9.42 ± 0.01	eV	NIST Webbook
ie	9.40 ± 0.02	eV	NIST Webbook
ie	9.74	eV	NIST Webbook
ie	9.71	eV	NIST Webbook
ie	9.65	eV	NIST Webbook
ie	9.53	eV	NIST Webbook
ie	9.57 ± 0.02	eV	NIST Webbook
ie	9.54	eV	NIST Webbook
log10ws	0.49		Estimated Solubility Method
logp	0.797		Crippen Method
mcvol	62.230	ml/mol	McGowan Method
pc	5190.00 ± 50.66	kPa	NIST Webbook
rinpol	626.00		NIST Webbook
rinpol	640.00		NIST Webbook
rinpol	618.00		NIST Webbook
rinpol	612.90		NIST Webbook
rinpol	618.00		NIST Webbook
rinpol	631.00		NIST Webbook
rinpol	620.00		NIST Webbook
rinpol	626.00		NIST Webbook
rinpol	631.00		NIST Webbook
rinpol	630.00		NIST Webbook
rinpol	611.00		NIST Webbook
rinpol	609.80		NIST Webbook
rinpol	609.90		NIST Webbook
rinpol	617.00		NIST Webbook
rinpol	621.00		NIST Webbook
rinpol	609.00		NIST Webbook

rinpol	632.00	NIST Webbook
rinpol	619.00	NIST Webbook
rinpol	610.00	NIST Webbook
rinpol	632.00	NIST Webbook
rinpol	629.00	NIST Webbook
rinpol	617.00	NIST Webbook
rinpol	624.00	NIST Webbook
rinpol	622.00	NIST Webbook
rinpol	623.00	NIST Webbook
rinpol	638.00	NIST Webbook
rinpol	627.00	NIST Webbook
rinpol	629.00	NIST Webbook
rinpol	612.00	NIST Webbook
rinpol	622.00	NIST Webbook
rinpol	608.00	NIST Webbook
rinpol	638.00	NIST Webbook
rinpol	615.00	NIST Webbook
rinpol	624.00	NIST Webbook
rinpol	615.00	NIST Webbook
rinpol	636.00	NIST Webbook
rinpol	619.00	NIST Webbook
rinpol	618.00	NIST Webbook
rinpol	615.00	NIST Webbook
rinpol	633.00	NIST Webbook
rinpol	620.00	NIST Webbook
rinpol	618.00	NIST Webbook
rinpol	623.20	NIST Webbook
rinpol	614.00	NIST Webbook
rinpol	630.00	NIST Webbook
rinpol	624.00	NIST Webbook
rinpol	620.00	NIST Webbook
rinpol	620.00	NIST Webbook
rinpol	629.00	NIST Webbook
rinpol	638.00	NIST Webbook
rinpol	640.00	NIST Webbook
ripol	898.00	NIST Webbook
ripol	888.00	NIST Webbook
ripol	907.00	NIST Webbook
ripol	915.00	NIST Webbook
ripol	887.00	NIST Webbook
ripol	829.00	NIST Webbook
ripol	900.00	NIST Webbook
ripol	868.00	NIST Webbook
ripol	872.00	NIST Webbook

ripol	895.00		NIST Webbook
ripol	854.00		NIST Webbook
ripol	829.00		NIST Webbook
ripol	857.00		NIST Webbook
ripol	854.00		NIST Webbook
ripol	867.00		NIST Webbook
ripol	866.00		NIST Webbook
ripol	861.00		NIST Webbook
ripol	868.00		NIST Webbook
ripol	903.00		NIST Webbook
ripol	868.00		NIST Webbook
ripol	900.00		NIST Webbook
sg	301.70 ± 1.70	J/molxK	NIST Webbook
sl	203.90	J/molxK	NIST Webbook
sl	203.80	J/molxK	NIST Webbook
sl	203.90	J/molxK	NIST Webbook
tb	337.65 ± 2.00	K	NIST Webbook
tb	337.65 ± 2.00	K	NIST Webbook
tb	339.65 ± 0.30	K	NIST Webbook
tb	339.10 ± 0.20	K	NIST Webbook
tb	340.00 ± 2.00	K	NIST Webbook
tb	339.00 ± 2.00	K	NIST Webbook
tb	337.90 ± 0.50	K	NIST Webbook
tb	340.15 ± 2.00	K	NIST Webbook
tb	338.90 ± 0.15	K	NIST Webbook
tb	339.10	K	NIST Webbook
tb	340.20	K	NIST Webbook
tb	339.20 ± 0.50	K	NIST Webbook
tb	339.15 ± 0.10	K	NIST Webbook
tb	338.90	K	NIST Webbook
tb	339.20	K	Experimental Isobaric Vapor-Liquid Equilibrium Data for Binary Mixtures of Cyclic Ethers with (1-Methylethyl)benzene
tb	339.12	K	Isobaric Vapor-Liquid Equilibrium for Binary System of Tetrahydrofuran + 1,4-Butanediol and gamma-Butyrolactone at 50.0 and 70.0 kPa
tb	338.65	K	Measurement and correlation of isobaric vapor-liquid equilibria of methanol + tetrahydrofuran + 1-alkyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide at 101.3 kPa

tb	339.12	K	Study of isobaric vapour liquid equilibrium of some cyclic ethers with 1-chloropropane: Experimental results and SAFT-VR modelling
tb	339.35 ± 0.50	K	NIST Webbook
tb	330.15 ± 2.00	K	NIST Webbook
tb	339.05 ± 0.30	K	NIST Webbook
tc	541.00 ± 1.50	K	NIST Webbook
tc	540.10 ± 0.30	K	NIST Webbook
tc	540.20	K	NIST Webbook
tf	164.05 ± 0.40	K	NIST Webbook
tf	164.63 ± 0.20	K	NIST Webbook
tf	164.15 ± 1.00	K	NIST Webbook
tf	165.10 ± 0.50	K	NIST Webbook
tt	164.76 ± 0.02	K	NIST Webbook
tt	164.76 ± 0.02	K	NIST Webbook
tt	164.76 ± 0.05	K	NIST Webbook
vc	0.225 ± 0.003	m3/kmol	NIST Webbook
volm	8.17e-05	m3/mol	Excess Gibbs Energies of the Ternary System 2-Methoxyethanol + Tetrahydrofuran + Cyclohexane and Other Relevant Binaries at 298.15 K
volm	8.17e-05	m3/mol	Excess Gibbs energies and volumes of the ternary system chloroform + tetrahydrofuran + cyclohexane at 298.15 K

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	133.68 ± 0.27	J/mol×K	500.15	NIST Webbook
cpg	120.39 ± 0.24	J/mol×K	449.15	NIST Webbook
cpg	106.12 ± 0.21	J/mol×K	399.15	NIST Webbook
cpg	91.36 ± 0.18	J/mol×K	349.15	NIST Webbook
cpg	85.13 ± 0.17	J/mol×K	328.15	NIST Webbook
cpl	106.70	J/mol×K	210.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry

cpl	124.50	J/mol×K	305.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry
cpl	123.10	J/mol×K	300.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry
cpl	121.70	J/mol×K	295.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry
cpl	120.30	J/mol×K	290.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry
cpl	119.20	J/mol×K	285.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry
cpl	120.50	J/mol×K	298.00	NIST Webbook
cpl	120.00	J/mol×K	298.15	NIST Webbook
cpl	123.90	J/mol×K	298.15	NIST Webbook
cpl	123.56	J/mol×K	298.15	NIST Webbook
cpl	122.92	J/mol×K	298.15	NIST Webbook
cpl	124.10	J/mol×K	298.15	NIST Webbook
cpl	124.10	J/mol×K	298.15	NIST Webbook
cpl	126.00	J/mol×K	310.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry
cpl	135.20	J/mol×K	318.15	Thermophysical properties of dimethyl sulfoxide + cyclic and linear ethers at 308.15K Application of an extended cell model
cpl	128.80	J/mol×K	320.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry

cpl	133.20	J/molxK	313.15	Thermophysical properties of dimethyl sulfoxide + cyclic and linear ethers at 308.15K Application of an extended cell model
cpl	131.40	J/molxK	308.15	Thermophysical properties of dimethyl sulfoxide + cyclic and linear ethers at 308.15K Application of an extended cell model
cpl	128.50	J/molxK	303.15	Thermophysical properties of dimethyl sulfoxide + cyclic and linear ethers at 308.15K Application of an extended cell model
cpl	136.60	J/molxK	323.15	Thermophysical properties of dimethyl sulfoxide + cyclic and linear ethers at 308.15K Application of an extended cell model
cpl	126.40	J/molxK	298.15	Thermophysical properties of dimethyl sulfoxide + cyclic and linear ethers at 308.15K Application of an extended cell model
cpl	123.80	J/molxK	293.15	Thermophysical properties of dimethyl sulfoxide + cyclic and linear ethers at 308.15K Application of an extended cell model
cpl	121.10	J/molxK	288.15	Thermophysical properties of dimethyl sulfoxide + cyclic and linear ethers at 308.15K Application of an extended cell model

cpl	115.50	J/mol×K	270.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry
cpl	130.50	J/mol×K	325.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry
cpl	105.50	J/mol×K	180.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry
cpl	105.60	J/mol×K	185.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry
cpl	105.80	J/mol×K	190.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry
cpl	106.00	J/mol×K	195.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry
cpl	106.30	J/mol×K	200.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry
cpl	117.90	J/mol×K	280.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry
cpl	116.60	J/mol×K	275.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry
cpl	107.90	J/mol×K	225.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry

cpl	114.40	J/mol×K	265.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry	
cpl	113.40	J/mol×K	260.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry	
cpl	112.40	J/mol×K	255.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry	
cpl	111.50	J/mol×K	250.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry	
cpl	110.60	J/mol×K	245.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry	
cpl	109.80	J/mol×K	240.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry	
cpl	109.20	J/mol×K	235.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry	
cpl	108.50	J/mol×K	230.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry	
cpl	106.50	J/mol×K	205.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry	
cpl	107.60	J/mol×K	220.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry	
cpl	123.90	J/mol×K	298.15	NIST Webbook	

cpl	127.30	J/molxK	315.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry
cpl	107.20	J/molxK	215.15	Measurement of heat capacities of ionic liquids by differential scanning calorimetry
dvisc	0.0004540	Paxs	298.15	Excess molar volume, viscosity, and refractive index study for the ternary mixture {2-methyl-2-butanol (1) + tetrahydrofuran (2) + propylamine (3)} at different temperatures. Application of the ERAS-model and Peng Robinson Stryjek Vera equation of state
dvisc	0.0004276	Paxs	308.15	Studies on Thermodynamic and Transport Properties of Binary Mixtures of Acetonitrile with Some Cyclic Ethers at Different Temperatures by Volumetric, Viscometric, and Interferometric Techniques
dvisc	0.0004631	Paxs	298.15	Studies on Thermodynamic and Transport Properties of Binary Mixtures of Acetonitrile with Some Cyclic Ethers at Different Temperatures by Volumetric, Viscometric, and Interferometric Techniques

dvisc	0.0003902	Paxs	318.15	Densities, Viscosities, and Sound Speeds of Some Acetate Salts in Binary Mixtures of Tetrahydrofuran and Methanol at (303.15, 313.15, and 323.15) K
dvisc	0.0004277	Paxs	308.15	Densities, Viscosities, and Sound Speeds of Some Acetate Salts in Binary Mixtures of Tetrahydrofuran and Methanol at (303.15, 313.15, and 323.15) K
dvisc	0.0004630	Paxs	298.15	Densities, Viscosities, and Sound Speeds of Some Acetate Salts in Binary Mixtures of Tetrahydrofuran and Methanol at (303.15, 313.15, and 323.15) K
dvisc	0.0004631	Paxs	298.15	Densities and Viscosities of the Binary Mixtures of Tetrahydrofuran with Isomeric Chlorobutanes at 298.15 K and 313.15 K
dvisc	0.0003992	Paxs	313.15	Excess Enthalpy, Excess Volume, Viscosity Deviation, and Speed of Sound Deviation for the Mixture Tetrahydrofuran + 2,2,2-Trifluoroethanol at (283.15, 298.15, and 313.15) K

dvisc	0.0004645	Paxs	298.15	Excess Enthalpy, Excess Volume, Viscosity Deviation, and Speed of Sound Deviation for the Mixture Tetrahydrofuran + 2,2,2-Trifluoroethanol at (283.15, 298.15, and 313.15) K
dvisc	0.0005438	Paxs	283.15	Excess Enthalpy, Excess Volume, Viscosity Deviation, and Speed of Sound Deviation for the Mixture Tetrahydrofuran + 2,2,2-Trifluoroethanol at (283.15, 298.15, and 313.15) K
dvisc	0.0004172	Paxs	308.15	Excess molar volume, viscosity, and refractive index study for the ternary mixture {2-methyl-2-butanol (1) + tetrahydrofuran (2) + propylamine (3)} at different temperatures. Application of the ERAS-model and Peng Robinson Stryjek Vera equation of state
dvisc	0.0003903	Paxs	318.15	Studies on Thermodynamic and Transport Properties of Binary Mixtures of Acetonitrile with Some Cyclic Ethers at Different Temperatures by Volumetric, Viscometric, and Interferometric Techniques

dvisc	0.0004991	Paxs	288.15	Excess molar volume, viscosity, and refractive index study for the ternary mixture {2-methyl-2-butanol (1) + tetrahydrofuran (2) + propylamine (3)} at different temperatures. Application of the ERAS-model and Peng Robinson Stryjek Vera equation of state
dvisc	0.0004125	Paxs	308.15	Ultrasonic and viscometric study of molecular interactions in binary mixtures of aniline with 1-propanol, 2-propanol, 2-methyl-1-propanol, and 2-methyl-2-propanol at different temperatures
dvisc	0.0004580	Paxs	298.15	Ultrasonic and viscometric study of molecular interactions in binary mixtures of aniline with 1-propanol, 2-propanol, 2-methyl-1-propanol, and 2-methyl-2-propanol at different temperatures
dvisc	0.0003900	Paxs	313.00	Ultrasonic velocity, viscosity and excess properties of binary mixture of tetrahydrofuran with 1-propanol and 2-propanol
dvisc	0.0004200	Paxs	303.00	Ultrasonic velocity, viscosity and excess properties of binary mixture of tetrahydrofuran with 1-propanol and 2-propanol

dvisc	0.0004700	Paxs	293.00	Ultrasonic velocity, viscosity and excess properties of binary mixture of tetrahydrofuran with 1-propanol and 2-propanol
dvisc	0.0004900	Paxs	293.00	Ultrasonic velocity, viscosity and excess properties of binary mixture of tetrahydrofuran with 1-propanol and 2-propanol
dvisc	0.0004260	Paxs	308.15	Viscosity Behavior of Some Oxygen Containing Compounds
dvisc	0.0004620	Paxs	298.15	Viscosity Behavior of Some Oxygen Containing Compounds
dvisc	0.0005080	Paxs	288.15	Viscosity Behavior of Some Oxygen Containing Compounds
dvisc	0.0003979	Paxs	313.15	Densities and Viscosities of the Binary Mixtures of Tetrahydrofuran with Isomeric Chlorobutanes at 298.15 K and 313.15 K
dvisc	0.0003762	Paxs	318.15	Ultrasonic and viscometric study of molecular interactions in binary mixtures of aniline with 1-propanol, 2-propanol, 2-methyl-1-propanol, and 2-methyl-2-propanol at different temperatures
hfust	8.54	kJ/mol	164.80	NIST Webbook
hfust	8.54	kJ/mol	164.76	NIST Webbook
hfust	8.54	kJ/mol	164.80	NIST Webbook
hfust	8.54	kJ/mol	164.76	NIST Webbook
hvapt	33.10	kJ/mol	306.00	NIST Webbook
hvapt	30.80	kJ/mol	320.50	NIST Webbook
hvapt	32.50 ± 0.20	kJ/mol	287.50	NIST Webbook

hvapt	29.60	kJ/mol	504.00	NIST Webbook
hvapt	29.00	kJ/mol	439.00	NIST Webbook
hvapt	31.90	kJ/mol	334.50	NIST Webbook
hvapt	32.30	kJ/mol	314.50	NIST Webbook
hvapt	29.81	kJ/mol	339.10	NIST Webbook
hvapt	32.80	kJ/mol	290.50	NIST Webbook
hvapt	33.00	kJ/mol	293.00	NIST Webbook
hvapt	31.80	kJ/mol	303.00	NIST Webbook
kvisc	0.0000005	m2/s	313.15	Experimental and predicted viscosities of binary mixtures of cyclic ethers with 1-chloropentane or 1-chlorohexane at 283.15, 298.15, and 313.15K
kvisc	0.0000005	m2/s	298.15	Experimental and predicted viscosities of binary mixtures of cyclic ethers with 1-chloropentane or 1-chlorohexane at 283.15, 298.15, and 313.15K
kvisc	0.0000006	m2/s	283.15	Experimental and predicted viscosities of binary mixtures of cyclic ethers with 1-chloropentane or 1-chlorohexane at 283.15, 298.15, and 313.15K
pvap	11.68	kPa	284.88	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran

pvap	21.60	kPa	298.15	Isothermal (vapour + liquid) equilibrium for binary mixtures of (tetrahydrofuran + 1,1,2,2-tetrachloroethane or tetrachloroethene) at nine temperatures
pvap	17.25	kPa	293.15	Isothermal (vapour + liquid) equilibrium for binary mixtures of (tetrahydrofuran + 1,1,2,2-tetrachloroethane or tetrachloroethene) at nine temperatures
pvap	13.67	kPa	288.15	Isothermal (vapour + liquid) equilibrium for binary mixtures of (tetrahydrofuran + 1,1,2,2-tetrachloroethane or tetrachloroethene) at nine temperatures
pvap	10.73	kPa	283.15	Isothermal (vapour + liquid) equilibrium for binary mixtures of (tetrahydrofuran + 1,1,2,2-tetrachloroethane or tetrachloroethene) at nine temperatures
pvap	101.30	kPa	338.65	Measurement and correlation of isobaric vapor-liquid equilibria of methanol + tetrahydrofuran + 1-alkyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide at 101.3 kPa
pvap	70.61	kPa	328.15	Isothermal vapour-liquid equilibrium for cyclic ethers with 1-chloropentane

pvap	40.62	kPa	313.15	Isothermal vapour-liquid equilibrium for cyclic ethers with 1-chloropentane
pvap	21.61	kPa	298.15	Isothermal vapour-liquid equilibrium for cyclic ethers with 1-chloropentane
pvap	8.61	kPa	279.41	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	26.81	kPa	303.15	Isothermal (vapour + liquid) equilibrium for binary mixtures of (tetrahydrofuran + 1,1,2,2-tetrachloroethane or tetrachloroethene) at nine temperatures
pvap	32.94	kPa	308.15	Isothermal (vapour + liquid) equilibrium for binary mixtures of (tetrahydrofuran + 1,1,2,2-tetrachloroethane or tetrachloroethene) at nine temperatures
pvap	40.21	kPa	313.15	Isothermal (vapour + liquid) equilibrium for binary mixtures of (tetrahydrofuran + 1,1,2,2-tetrachloroethane or tetrachloroethene) at nine temperatures

pvap	48.68	kPa	318.15	Isothermal (vapour + liquid) equilibrium for binary mixtures of (tetrahydrofuran + 1,1,2,2-tetrachloroethane or tetrachloroethene) at nine temperatures
pvap	58.54	kPa	323.15	Isothermal (vapour + liquid) equilibrium for binary mixtures of (tetrahydrofuran + 1,1,2,2-tetrachloroethane or tetrachloroethene) at nine temperatures
pvap	13.79	kPa	288.50	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	21.61	kPa	298.15	Isothermal (vapour + liquid) equilibrium of (cyclic ethers + chlorohexane) mixtures: Experimental results and SAFT modelling
pvap	14.00	kPa	289.00	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	14.48	kPa	289.66	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran

pvap	15.65	kPa	291.04	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	16.65	kPa	292.31	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	17.43	kPa	293.30	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	19.72	kPa	296.05	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	20.37	kPa	296.84	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	20.65	kPa	297.09	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran

pvap	21.50	kPa	298.15	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	22.24	kPa	298.72	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	70.61	kPa	328.15	Isothermal (vapour + liquid) equilibrium of (cyclic ethers + chlorohexane) mixtures: Experimental results and SAFT modelling
pvap	25.28	kPa	301.73	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	26.87	kPa	303.23	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	27.79	kPa	304.04	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran

pvap	7.10	kPa	275.99	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	7.56	kPa	276.53	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	7.84	kPa	277.76	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	8.33	kPa	278.19	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	12.44	kPa	286.55	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	9.04	kPa	279.86	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran

pvap	8.93	kPa	280.16	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	9.96	kPa	281.95	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	10.55	kPa	283.07	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	10.32	kPa	283.15	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	10.90	kPa	283.36	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	29.43	kPa	305.35	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran

pvap	30.77	kPa	306.54	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	32.49	kPa	308.07	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	35.55	kPa	310.20	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	36.65	kPa	310.86	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	38.54	kPa	311.90	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	40.13	kPa	313.12	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran

pvap	40.09	kPa	313.18	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	44.65	kPa	315.86	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	47.86	kPa	317.71	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	51.31	kPa	319.58	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	55.34	kPa	321.57	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	59.46	kPa	323.59	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran

pvap	6.59	kPa	273.60	Vapor Pressure and Its Temperature Dependence of 28 Organic Compounds: Cyclic Amines, Cyclic Ethers, and Cyclic and Open Chain Secondary Alcohols
pvap	8.38	kPa	278.30	Vapor Pressure and Its Temperature Dependence of 28 Organic Compounds: Cyclic Amines, Cyclic Ethers, and Cyclic and Open Chain Secondary Alcohols
pvap	11.81	kPa	285.02	Vapor Pressure and Its Temperature Dependence of 28 Organic Compounds: Cyclic Amines, Cyclic Ethers, and Cyclic and Open Chain Secondary Alcohols
pvap	13.02	kPa	286.97	Vapor Pressure and Its Temperature Dependence of 28 Organic Compounds: Cyclic Amines, Cyclic Ethers, and Cyclic and Open Chain Secondary Alcohols
pvap	14.89	kPa	289.75	Vapor Pressure and Its Temperature Dependence of 28 Organic Compounds: Cyclic Amines, Cyclic Ethers, and Cyclic and Open Chain Secondary Alcohols

pvap	17.00	kPa	292.65	Vapor Pressure and Its Temperature Dependence of 28 Organic Compounds: Cyclic Amines, Cyclic Ethers, and Cyclic and Open Chain Secondary Alcohols
pvap	21.27	kPa	297.57	Vapor Pressure and Its Temperature Dependence of 28 Organic Compounds: Cyclic Amines, Cyclic Ethers, and Cyclic and Open Chain Secondary Alcohols
pvap	26.67	kPa	302.75	Vapor Pressure and Its Temperature Dependence of 28 Organic Compounds: Cyclic Amines, Cyclic Ethers, and Cyclic and Open Chain Secondary Alcohols
pvap	32.88	kPa	307.83	Vapor Pressure and Its Temperature Dependence of 28 Organic Compounds: Cyclic Amines, Cyclic Ethers, and Cyclic and Open Chain Secondary Alcohols
pvap	26.67	kPa	303.04	Double Azeotropy in Binary Mixtures 1,1,1,2,3,4,4,5,5,5-Decafluoropentane High-Boiling Diastereomer + Tetrahydrofuran
pvap	26.68	kPa	303.05	Double Azeotropy in Binary Mixtures 1,1,1,2,3,4,4,5,5,5-Decafluoropentane High-Boiling Diastereomer + Tetrahydrofuran

pvap	29.09	kPa	305.11	Double Azeotropy in Binary Mixtures 1,1,1,2,3,4,4,5,5,5-Decafluoropentane High-Boiling Diastereomer + Tetrahydrofuran
pvap	34.50	kPa	309.28	Double Azeotropy in Binary Mixtures 1,1,1,2,3,4,4,5,5,5-Decafluoropentane High-Boiling Diastereomer + Tetrahydrofuran
pvap	40.15	kPa	313.11	Double Azeotropy in Binary Mixtures 1,1,1,2,3,4,4,5,5,5-Decafluoropentane High-Boiling Diastereomer + Tetrahydrofuran
pvap	45.51	kPa	316.36	Double Azeotropy in Binary Mixtures 1,1,1,2,3,4,4,5,5,5-Decafluoropentane High-Boiling Diastereomer + Tetrahydrofuran
pvap	48.62	kPa	318.10	Double Azeotropy in Binary Mixtures 1,1,1,2,3,4,4,5,5,5-Decafluoropentane High-Boiling Diastereomer + Tetrahydrofuran
pvap	23.84	kPa	300.57	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	26.62	kPa	302.96	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran

rfi	1.40720	298.15	Bubble temperature measurements on the binary mixtures formed by decane with a variety of compounds at 95.8 kPa
rfi	1.39510	318.15	Refractive Indices of Binary Mixtures of Tetrahydrofuran with Aromatic Hydrocarbon at Temperatures from (288.15 to 318.15) K
rfi	1.39750	313.15	Refractive Indices of Binary Mixtures of Tetrahydrofuran with Aromatic Hydrocarbon at Temperatures from (288.15 to 318.15) K
rfi	1.39990	308.15	Refractive Indices of Binary Mixtures of Tetrahydrofuran with Aromatic Hydrocarbon at Temperatures from (288.15 to 318.15) K
rfi	1.40230	303.15	Refractive Indices of Binary Mixtures of Tetrahydrofuran with Aromatic Hydrocarbon at Temperatures from (288.15 to 318.15) K
rfi	1.40460	298.15	Refractive Indices of Binary Mixtures of Tetrahydrofuran with Aromatic Hydrocarbon at Temperatures from (288.15 to 318.15) K
rfi	1.40700	293.15	Refractive Indices of Binary Mixtures of Tetrahydrofuran with Aromatic Hydrocarbon at Temperatures from (288.15 to 318.15) K

rfi	1.40940	288.15	Refractive Indices of Binary Mixtures of Tetrahydrofuran with Aromatic Hydrocarbon at Temperatures from (288.15 to 318.15) K
rfi	1.40500	298.15	Viscosity, Density, Speed of Sound, and Refractive Index of Binary Mixtures of Organic Solvent + Ionic Liquid, 1-Butyl-3-methylimidazolium Hexafluorophosphate at 298.15 K
rfi	1.40520	298.15	Acoustic and Volumetric Properties of Binary Mixtures of Ionic Liquid 1-Butyl-3-methylimidazolium Bis(trifluoromethylsulfonyl)imide with Acetonitrile and Tetrahydrofuran
rfi	1.40610	298.15	Vapor Liquid Equilibrium for Ternary and Binary Mixtures of Tetrahydrofuran, Cyclohexane, and 1,2-Propanediol at 101.3 kPa
rfi	1.40500	293.15	Solubilities of Phosphorus-Containing Compounds in Selected Solvents
rfi	1.40470	298.15	Solubilities of Bis (2,2,6,6-Tetramethyl-4-Piperidiny) Maleate in Hexane, Heptane, Octane, m-Xylene and Tetrahydrofuran from (253.15 to 310.15) K
rfi	1.40480	303.15	Thermodynamic Properties of Water + Tetrahydrofuran and Water + 1,4-Dioxane Mixtures at (303.15, 313.15, and 323.15) K

rfi	1.40260	303.15	Density, Viscosity, Refractive Index, and Speed of Sound in the Binary Mixtures of Tri-n-butylamine + Triethylamine, + Tetrahydrofuran, + Tetradecane, + Tetrachloroethylene, + Pyridine, or + Trichloroethylene at (298.15, 303.15, and 308.15) K
rfi	1.39980	308.15	Density, Viscosity, Refractive Index, and Speed of Sound in the Binary Mixtures of Tri-n-butylamine + Triethylamine, + Tetrahydrofuran, + Tetradecane, + Tetrachloroethylene, + Pyridine, or + Trichloroethylene at (298.15, 303.15, and 308.15) K
rfi	1.40490	298.15	Density, Viscosity, Refractive Index, and Speed of Sound in the Binary Mixtures of Tri-n-butylamine + Triethylamine, + Tetrahydrofuran, + Tetradecane, + Tetrachloroethylene, + Pyridine, or + Trichloroethylene at (298.15, 303.15, and 308.15) K

rfi	1.40720		293.15	Bubble-Temperature Measurements on Some Binary Mixtures Formed by Tetrahydrofuran or Amyl Alcohol with Hydrocarbons, Chlorohydrocarbons, or Butanols at (94.6 or 95.8) kPa
rfi	1.40280		303.15	Densities, speeds of sound, isentropic compressibilities, refractive indexes, and viscosities of tetrahydrofuran with haloalkane or alkyl ethanoate at T = 303.15 K
rfi	1.40480		298.15	To the issue of temperature-dependent behavior of standard molar volumes of components in the binary system (water + tetrahydrofuran) at ambient pressure
rhoI	882.00	kg/m3	298.15	Vapour liquid equilibrium of cyclic ethers with 1-chlorohexane: Experimental results and UNIFAC predictions
rhoI	878.00	kg/m3	303.00	Densities, ultrasonic speeds and refractive indices of phenetole with N-methyl-2-pyrrolidone, N,N-dimethylformamide and tetrahydrofuran binary mixtures at different temperatures

rhoI	882.00	kg/m3	298.00	Densities, ultrasonic speeds and refractive indices of phenetole with N-methyl-2-pyrrolidone, N,N-dimethylformamide and tetrahydrofuran binary mixtures at different temperatures
rhoI	855.46	kg/m3	323.15	Investigation on thermophysical and excess properties of binary mixtures of imidazolium based ionic liquids at temperatures (293.15 to 323.15) K: III [Cnmim][PF6] (n = 4, 6, 8) + THF
rhoI	874.00	kg/m3	308.00	Densities, ultrasonic speeds and refractive indices of phenetole with N-methyl-2-pyrrolidone, N,N-dimethylformamide and tetrahydrofuran binary mixtures at different temperatures
rhoI	887.01	kg/m3	293.15	Temperature and concentration dependence of volumetric properties of (tetrahydrofuran + methyl acrylate, or + ethyl acrylate, or + n-butyl acrylate, or + tert-butyl acrylate) binary mixtures
rhoI	882.05	kg/m3	298.15	Temperature and concentration dependence of volumetric properties of (tetrahydrofuran + methyl acrylate, or + ethyl acrylate, or + n-butyl acrylate, or + tert-butyl acrylate) binary mixtures

rhoI	872.14	kg/m3	308.15	Temperature and concentration dependence of volumetric properties of (tetrahydrofuran + methyl acrylate, or + ethyl acrylate, or + n-butyl acrylate, or + tert-butyl acrylate) binary mixtures
rhoI	867.19	kg/m3	313.15	Temperature and concentration dependence of volumetric properties of (tetrahydrofuran + methyl acrylate, or + ethyl acrylate, or + n-butyl acrylate, or + tert-butyl acrylate) binary mixtures
rhoI	877.10	kg/m3	303.15	Temperature and concentration dependence of volumetric properties of (tetrahydrofuran + methyl acrylate, or + ethyl acrylate, or + n-butyl acrylate, or + tert-butyl acrylate) binary mixtures
rhoI	882.00	kg/m3	298.15	Volumetric Properties of the Ionic Liquid, 1-Butyl-3-methylimidazolium Tetrafluoroborate, in Organic Solvents at T = 298.15 K
rhoI	882.09	kg/m3	298.15	Excess Enthalpies of the Ternary Mixtures: Tetrahydrofuran + (Hexane or Cyclohexane) + Decane at 298.15K

rhoI	882.06	kg/m3	298.15	Excess Molar volumes and Surface Tensions of Trimethylbenzene with Tetrahydrofuran Tetrachloromethane and Dimethylsulfoxide at 298.15 K	
rhoI	882.10	kg/m3	298.15	Study of the Surface Tension of Chlorocyclohexane or Bromocyclohexane with Some Cyclic Ethers	
rhoI	887.41	kg/m3	293.15	Volumetric Properties of Binary Mixtures Containing Ionic Liquids and Some Aprotic Solvents	
rhoI	881.98	kg/m3	298.15	Volumetric Properties of Binary Mixtures Containing Ionic Liquids and Some Aprotic Solvents	
rhoI	876.47	kg/m3	303.13	Volumetric Properties of Binary Mixtures Containing Ionic Liquids and Some Aprotic Solvents	
rhoI	870.96	kg/m3	308.15	Volumetric Properties of Binary Mixtures Containing Ionic Liquids and Some Aprotic Solvents	
rhoI	882.08	kg/m3	298.15	Excess Molar Enthalpies of Dipropyl Ether + Dibutyl Ether + (1-Hexene or Tetrahydrofuran) at 298.15 K	

rhoI	881.50	kg/m3	298.15	Densities, Viscosities, and Sound Speed of Binary Mixtures of Hexyl Acetate with Tetrahydrofuran, 1,4-Dioxane, Anisole, and Butyl Vinyl Ether
rhoI	878.70	kg/m3	303.15	Densities, Viscosities, and Sound Speed of Binary Mixtures of Hexyl Acetate with Tetrahydrofuran, 1,4-Dioxane, Anisole, and Butyl Vinyl Ether
rhoI	873.00	kg/m3	308.15	Densities, Viscosities, and Sound Speed of Binary Mixtures of Hexyl Acetate with Tetrahydrofuran, 1,4-Dioxane, Anisole, and Butyl Vinyl Ether
rhoI	866.90	kg/m3	313.15	Densities, Viscosities, and Sound Speed of Binary Mixtures of Hexyl Acetate with Tetrahydrofuran, 1,4-Dioxane, Anisole, and Butyl Vinyl Ether
rhoI	887.52	kg/m3	293.15	Volumetric properties of binary mixtures of ethers and acetonitrile: Experimental results and application of the Prigogine Flory Patterson theory
rhoI	861.11	kg/m3	318.15	Investigation on thermophysical and excess properties of binary mixtures of imidazolium based ionic liquids at temperatures (293.15 to 323.15) K: III [Cnmim][PF6] (n = 4, 6, 8) + THF

rhoI	866.71	kg/m3	313.15	Investigation on thermophysical and excess properties of binary mixtures of imidazolium based ionic liquids at temperatures (293.15 to 323.15) K: III [Cnmim][PF6] (n = 4, 6, 8) + THF
rhoI	872.28	kg/m3	308.15	Investigation on thermophysical and excess properties of binary mixtures of imidazolium based ionic liquids at temperatures (293.15 to 323.15) K: III [Cnmim][PF6] (n = 4, 6, 8) + THF
rhoI	877.81	kg/m3	303.15	Investigation on thermophysical and excess properties of binary mixtures of imidazolium based ionic liquids at temperatures (293.15 to 323.15) K: III [Cnmim][PF6] (n = 4, 6, 8) + THF
rhoI	883.30	kg/m3	298.15	Investigation on thermophysical and excess properties of binary mixtures of imidazolium based ionic liquids at temperatures (293.15 to 323.15) K: III [Cnmim][PF6] (n = 4, 6, 8) + THF
rhoI	888.77	kg/m3	293.15	Investigation on thermophysical and excess properties of binary mixtures of imidazolium based ionic liquids at temperatures (293.15 to 323.15) K: III [Cnmim][PF6] (n = 4, 6, 8) + THF

rhoI	859.96	kg/m3	318.15	Experimental and theoretical excess molar properties of imidazolium based ionic liquids with molecular organic solvents I. 1-Hexyl-3-methylimidazolium tetrafluoroborate and 1-octyl-3-methylimidazolium tetrafluoroborate with cyclic ethers
rhoI	871.16	kg/m3	308.15	Experimental and theoretical excess molar properties of imidazolium based ionic liquids with molecular organic solvents I. 1-Hexyl-3-methylimidazolium tetrafluoroborate and 1-octyl-3-methylimidazolium tetrafluoroborate with cyclic ethers
rhoI	882.21	kg/m3	298.15	Experimental and theoretical excess molar properties of imidazolium based ionic liquids with molecular organic solvents I. 1-Hexyl-3-methylimidazolium tetrafluoroborate and 1-octyl-3-methylimidazolium tetrafluoroborate with cyclic ethers
rhoI	855.29	kg/m3	323.15	Experimental and theoretical excess molar properties of imidazolium based ionic liquids with molecular organic solvents I. 1-Hexyl-3-methylimidazolium tetrafluoroborate and 1-octyl-3-methylimidazolium tetrafluoroborate with cyclic ethers

rhoI	860.96	kg/m3	318.15	Experimental and theoretical excess molar properties of imidazolium based ionic liquids with molecular organic solvents I. 1-Hexyl-3-methylimidazolium tetrafluoroborate and 1-octyl-3-methylimidazolium tetrafluoroborate with cyclic ethers
rhoI	866.58	kg/m3	313.15	Experimental and theoretical excess molar properties of imidazolium based ionic liquids with molecular organic solvents I. 1-Hexyl-3-methylimidazolium tetrafluoroborate and 1-octyl-3-methylimidazolium tetrafluoroborate with cyclic ethers
rhoI	872.16	kg/m3	308.15	Experimental and theoretical excess molar properties of imidazolium based ionic liquids with molecular organic solvents I. 1-Hexyl-3-methylimidazolium tetrafluoroborate and 1-octyl-3-methylimidazolium tetrafluoroborate with cyclic ethers
rhoI	877.70	kg/m3	303.15	Experimental and theoretical excess molar properties of imidazolium based ionic liquids with molecular organic solvents I. 1-Hexyl-3-methylimidazolium tetrafluoroborate and 1-octyl-3-methylimidazolium tetrafluoroborate with cyclic ethers

rhoI	882.21	kg/m3	298.15	Experimental and theoretical excess molar properties of imidazolium based ionic liquids with molecular organic solvents I. 1-Hexyl-3-methylimidazolium tetrafluoroborate and 1-octyl-3-methylimidazolium tetrafluoroborate with cyclic ethers
rhoI	888.69	kg/m3	293.15	Experimental and theoretical excess molar properties of imidazolium based ionic liquids with molecular organic solvents I. 1-Hexyl-3-methylimidazolium tetrafluoroborate and 1-octyl-3-methylimidazolium tetrafluoroborate with cyclic ethers
rhoI	882.15	kg/m3	298.15	Densities, speeds of sound and viscosities of binary mixtures of tetrahydrofuran with 1-hexanol, 1-octanol and 1-decanol at T = (298.15 to 313.15) K
rhoI	866.23	kg/m3	313.15	Volumetric properties of binary mixtures of N-ethylformamide with tetrahydrofuran, 2-butanone and ethylacetate from (293.15 to 313.15) K
rhoI	872.02	kg/m3	308.15	Volumetric properties of binary mixtures of N-ethylformamide with tetrahydrofuran, 2-butanone and ethylacetate from (293.15 to 313.15) K

rhoI	877.60	kg/m3	303.15	Volumetric properties of binary mixtures of N-ethylformamide with tetrahydrofuran, 2-butanone and ethylacetate from (293.15 to 313.15) K
rhoI	883.24	kg/m3	298.15	Volumetric properties of binary mixtures of N-ethylformamide with tetrahydrofuran, 2-butanone and ethylacetate from (293.15 to 313.15) K
rhoI	888.73	kg/m3	293.15	Volumetric properties of binary mixtures of N-ethylformamide with tetrahydrofuran, 2-butanone and ethylacetate from (293.15 to 313.15) K
rhoI	859.84	kg/m3	318.15	Phenomenon of "negative partial molar expansibility" of water in tetrahydrofuran: How plausible is it? Discussion on the paper "Volumetric properties on the (tetrahydrofuran + water) and (tetra-n-butylammonium bromide + water) systems: Experimental measurements and correlations" by Veronica Belandria, Ammir H. Mohammadi and Dominique Richon [J. Chem. Thermodyn. 41 (2009) 1382 1386]

rhoI	871.01	kg/m3	308.15	Phenomenon of "negative partial molar expansibility" of water in tetrahydrofuran: How plausible is it? Discussion on the paper "Volumetric properties on the (tetrahydrofuran + water) and (tetra-n-butylammonium bromide + water) systems: Experimental measurements and correlations" by Veronica Belandria, Ammir H. Mohammadi and Dominique Richon [J. Chem. Thermodyn. 41 (2009) 1382 1386]
------	--------	-------	--------	--

rhoI	882.04	kg/m3	298.15	Phenomenon of "negative partial molar expansibility" of water in tetrahydrofuran: How plausible is it? Discussion on the paper "Volumetric properties on the (tetrahydrofuran + water) and (tetra-n-butylammonium bromide + water) systems: Experimental measurements and correlations" by Veronica Belandria, Ammir H. Mohammadi and Dominique Richon [J. Chem. Thermodyn. 41 (2009) 1382 1386]
------	--------	-------	--------	--

rhoI	892.94	kg/m3	288.15	Phenomenon of "negative partial molar expansibility" of water in tetrahydrofuran: How plausible is it? Discussion on the paper "Volumetric properties on the (tetrahydrofuran + water) and (tetra-n-butylammonium bromide + water) systems: Experimental measurements and correlations" by Veronica Belandria, Ammir H. Mohammadi and Dominique Richon [J. Chem. Thermodyn. 41 (2009) 1382 1386]
------	--------	-------	--------	--

rhoI	903.71	kg/m3	278.15	Phenomenon of "negative partial molar expansibility" of water in tetrahydrofuran: How plausible is it? Discussion on the paper "Volumetric properties on the (tetrahydrofuran + water) and (tetra-n-butylammonium bromide + water) systems: Experimental measurements and correlations" by Veronica Belandria, Ammir H. Mohammadi and Dominique Richon [J. Chem. Thermodyn. 41 (2009) 1382 1386]
------	--------	-------	--------	--

rhoI	876.54	kg/m3	303.15	Volumetric properties of binary mixtures of ethers and acetonitrile: Experimental results and application of the Prigogine Flory Patterson theory
------	--------	-------	--------	---

rhoI	882.04	kg/m3	298.15	Volumetric properties of binary mixtures of ethers and acetonitrile: Experimental results and application of the Prigogine Flory Patterson theory
rhoI	881.95	kg/m3	298.15	Surface Tension of Mixtures of Tetrahydrofuran or Tetrahydropyran with Isomeric Chlorobutanes
rhoI	892.86	kg/m3	288.15	Volumetric properties of binary mixtures of ethers and acetonitrile: Experimental results and application of the Prigogine Flory Patterson theory
rhoI	882.09	kg/m3	298.15	Excess molar enthalpies of the ternary mixtures (1-hexene + tetrahydrofuran or 2-methyltetrahydrofuran + methyl tert-butyl ether) at the temperature 298.15K.
rhoI	882.00	kg/m3	298.15	Volumetric and compressibility behaviour of ionic liquid, 1-n-butyl-3-methylimidazolium hexafluorophosphate and tetrabutylammonium hexafluorophosphate in organic solvents at T = 298.15 K
rhoI	882.09	kg/m3	298.15	Excess molar enthalpies of the ternary mixtures: (tetrahydrofuran or 2-methyltetrahydrofuran + methyl tert-butyl ether + n-octane) at the temperature 298.15 K

rhoI	882.09	kg/m3	298.15	Excess molar enthalpies of the ternary mixtures: {tetrahydrofuran or 2-methyltetrahydrofuran + methyl tert-butyl ether + n-dodecane} at the temperature 298.15 K
rhoI	882.01	kg/m3	298.15	Standard partial molar volumes of some electrolytes in ethylene carbonate based mixtures
rhoI	876.50	kg/m3	303.15	Unravelling various types of non-covalent interactions of benzyl amine with ethers in n-hexane at 303.15 K by ultrasonic and DFT methods
rhoI	884.99	kg/m3	298.15	Quaternary isothermal vapor-liquid equilibrium of the model biofuel 2-butanone + n-heptane + tetrahydrofuran + cyclohexane using Raman spectroscopic characterization
rhoI	854.56	kg/m3	323.15	Volumetric and viscometric behavior of the binary systems ethyl lactate + 1,2- propanediol, + 1,3-propanediol, + tetrahydrofuran and + tetraethylene glycol dimethyl ether. New UNIFAC-VISCO and ASOG-VISCO parameters determination

rhoI	860.21	kg/m3	318.15	Volumetric and viscometric behavior of the binary systems ethyl lactate + 1,2- propanediol, + 1,3-propanediol, + tetrahydrofuran and + tetraethylene glycol dimethyl ether. New UNIFAC-VISCO and ASOG-VISCO parameters determination
rhoI	865.82	kg/m3	313.15	Volumetric and viscometric behavior of the binary systems ethyl lactate + 1,2- propanediol, + 1,3-propanediol, + tetrahydrofuran and + tetraethylene glycol dimethyl ether. New UNIFAC-VISCO and ASOG-VISCO parameters determination
rhoI	871.38	kg/m3	308.15	Volumetric and viscometric behavior of the binary systems ethyl lactate + 1,2- propanediol, + 1,3-propanediol, + tetrahydrofuran and + tetraethylene glycol dimethyl ether. New UNIFAC-VISCO and ASOG-VISCO parameters determination

rhoI	876.91	kg/m3	303.15	Volumetric and viscometric behavior of the binary systems ethyl lactate + 1,2- propanediol, + 1,3-propanediol, + tetrahydrofuran and + tetraethylene glycol dimethyl ether. New UNIFAC-VISCO and ASOG-VISCO parameters determination
rhoI	882.40	kg/m3	298.15	Volumetric and viscometric behavior of the binary systems ethyl lactate + 1,2- propanediol, + 1,3-propanediol, + tetrahydrofuran and + tetraethylene glycol dimethyl ether. New UNIFAC-VISCO and ASOG-VISCO parameters determination
rhoI	887.87	kg/m3	293.15	Volumetric and viscometric behavior of the binary systems ethyl lactate + 1,2- propanediol, + 1,3-propanediol, + tetrahydrofuran and + tetraethylene glycol dimethyl ether. New UNIFAC-VISCO and ASOG-VISCO parameters determination

rhoI	893.29	kg/m3	288.15	Volumetric and viscometric behavior of the binary systems ethyl lactate + 1,2- propanediol, + 1,3-propanediol, + tetrahydrofuran and + tetraethylene glycol dimethyl ether. New UNIFAC-VISCO and ASOG-VISCO parameters determination
rhoI	862.23	kg/m3	318.15	Temperature and concentration dependence of volumetric properties of (tetrahydrofuran + methyl acrylate, or + ethyl acrylate, or + n-butyl acrylate, or + tert-butyl acrylate) binary mixtures
sfust	51.80	J/molxK	164.76	NIST Webbook
sfust	51.83	J/molxK	164.76	NIST Webbook
speedsl	1207.00	m/s	313.15	Thermodynamic and acoustic properties of binary mixtures of ethers. V. Diisopropyl ether or oxolane with 2- or 3-chloroanilines at 303.15, 313.15 and 323.15 K
speedsl	1229.61	m/s	308.15	Ultrasonic and Volumetric Properties of 1-Ethyl-3-methylimidazolium Trifluoromethanesulfonate Ionic Liquid with 2-Propanol or Tetrahydrofuran at Several Temperatures
speedsl	1277.37	m/s	298.15	Ultrasonic and Volumetric Properties of 1-Ethyl-3-methylimidazolium Trifluoromethanesulfonate Ionic Liquid with 2-Propanol or Tetrahydrofuran at Several Temperatures

speedsl	1325.78	m/s	288.15	Ultrasonic and Volumetric Properties of 1-Ethyl-3-methylimidazolium Trifluoromethanesulfonate Ionic Liquid with 2-Propanol or Tetrahydrofuran at Several Temperatures
speedsl	1374.98	m/s	278.15	Ultrasonic and Volumetric Properties of 1-Ethyl-3-methylimidazolium Trifluoromethanesulfonate Ionic Liquid with 2-Propanol or Tetrahydrofuran at Several Temperatures
speedsl	1280.20	m/s	298.15	Compressibility Studies of Binary Solutions Involving Water as a Solute in Nonaqueous Solvents at T) 298.15 K
speedsl	1165.00	m/s	323.15	Thermodynamic and acoustic properties of binary mixtures of ethers. V. Diisopropyl ether or oxolane with 2- or 3-chloroanilines at 303.15, 313.15 and 323.15 K
speedsl	1182.37	m/s	318.15	Ultrasonic and Volumetric Properties of 1-Ethyl-3-methylimidazolium Trifluoromethanesulfonate Ionic Liquid with 2-Propanol or Tetrahydrofuran at Several Temperatures
speedsl	1254.00	m/s	303.15	Thermodynamic and acoustic properties of binary mixtures of ethers. V. Diisopropyl ether or oxolane with 2- or 3-chloroanilines at 303.15, 313.15 and 323.15 K

speedsl	1165.00	m/s	323.15	Thermodynamic and acoustic properties of binary mixtures of ethers. IV. Diisopropyl ether or oxolane with N,N-dimethylaniline or N,Ndiethylaniline at 303.15, 313.15 and 323.15 K
speedsl	1207.00	m/s	313.15	Thermodynamic and acoustic properties of binary mixtures of ethers. IV. Diisopropyl ether or oxolane with N,N-dimethylaniline or N,Ndiethylaniline at 303.15, 313.15 and 323.15 K
speedsl	1254.00	m/s	303.15	Thermodynamic and acoustic properties of binary mixtures of ethers. IV. Diisopropyl ether or oxolane with N,N-dimethylaniline or N,Ndiethylaniline at 303.15, 313.15 and 323.15 K
speedsl	1165.00	m/s	323.15	Thermodynamic and acoustic properties of binary mixtures of ethers. III. Diisopropyl ether or oxolane with o- or m-toluidines at 303.15, 313.15 and 323.15 K
speedsl	1256.00	m/s	303.15	Thermodynamic and acoustic properties of binary mixtures of oxolane with aniline and substituted anilines at 303.15, 313.15 and 323.15 K

speedsl	1254.00	m/s	303.15	Thermodynamic and acoustic properties of binary mixtures of ethers. III. Diisopropyl ether or oxolane with o- or m-toluidines at 303.15, 313.15 and 323.15 K
speedsl	1171.00	m/s	323.15	Thermodynamic and acoustic properties of binary mixtures of oxolane with aniline and substituted anilines at 303.15, 313.15 and 323.15 K
speedsl	1212.00	m/s	313.15	Thermodynamic and acoustic properties of binary mixtures of oxolane with aniline and substituted anilines at 303.15, 313.15 and 323.15 K
speedsl	1207.00	m/s	313.15	Thermodynamic and acoustic properties of binary mixtures of ethers. III. Diisopropyl ether or oxolane with o- or m-toluidines at 303.15, 313.15 and 323.15 K
srf	0.03	N/m	293.15	Thermodynamic surface properties of [BMIm][NTf2] or [EMIm][NTf2] binary mixtures with tetrahydrofuran, acetonitrile or dimethylsulfoxide
srf	0.03	N/m	308.15	Densities, Viscosities, Refractive Indexes, and Surface Tensions for Binary and Ternary Mixtures of Tetrahydrofuran, 2-Propanol, and 2,2,4-Trimethylpentane

srf	0.03	N/m	298.15	Densities, Viscosities, Refractive Indexes, and Surface Tensions for Binary and Ternary Mixtures of Tetrahydrofuran, 2-Propanol, and 2,2,4-Trimethylpentane
srf	0.03	N/m	288.15	Densities, Viscosities, Refractive Indexes, and Surface Tensions for Binary and Ternary Mixtures of Tetrahydrofuran, 2-Propanol, and 2,2,4-Trimethylpentane
srf	0.03	N/m	308.15	Surface Tensions and the Gibbs Excess Surface Concentration of Binary Mixtures of the Ionic Liquid 1-Ethyl-3-methylimidazolium Bis[(trifluoromethyl)sulfonyl]imide with Tetrahydrofuran and Acetonitrile
srf	0.03	N/m	298.15	Surface Tensions and the Gibbs Excess Surface Concentration of Binary Mixtures of the Ionic Liquid 1-Ethyl-3-methylimidazolium Bis[(trifluoromethyl)sulfonyl]imide with Tetrahydrofuran and Acetonitrile
srf	0.03	N/m	293.15	Surface Tensions and the Gibbs Excess Surface Concentration of Binary Mixtures of the Ionic Liquid 1-Ethyl-3-methylimidazolium Bis[(trifluoromethyl)sulfonyl]imide with Tetrahydrofuran and Acetonitrile

srf	0.03	N/m	308.15	Thermodynamic surface properties of [BMIm][NTf2] or [EMIm][NTf2] binary mixtures with tetrahydrofuran, acetonitrile or dimethylsulfoxide
srf	0.03	N/m	303.15	Thermodynamic surface properties of [BMIm][NTf2] or [EMIm][NTf2] binary mixtures with tetrahydrofuran, acetonitrile or dimethylsulfoxide
srf	0.03	N/m	298.15	Thermodynamic surface properties of [BMIm][NTf2] or [EMIm][NTf2] binary mixtures with tetrahydrofuran, acetonitrile or dimethylsulfoxide
srf	0.03	N/m	303.13	Surface Tensions and the Gibbs Excess Surface Concentration of Binary Mixtures of the Ionic Liquid 1-Ethyl-3-methylimidazolium Bis[(trifluoromethyl)sulfonyl]imide with Tetrahydrofuran and Acetonitrile

Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbp	339.15	K	96.60	Low cost apparatus for rapid boiling point determination of small air sensitive samples under inert atmosphere

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/T + C \cdot \ln(T) + D \cdot T^2$
Coeff. A	7.83577e+01
Coeff. B	-6.33998e+03
Coeff. C	-9.60880e+00
Coeff. D	8.18348e-06
Temperature range (K), min.	164.65
Temperature range (K), max.	540.15

Datasets

Mass density, kg/m3

Pressure, kPa - Liquid	Temperature, K - Liquid	Mass density, kg/m3 - Liquid
85.90	298.15	882.02
Reference		https://www.doi.org/10.1016/j.fluid.2013.05.001

Viscosity, Pa*s

Temperature, K - Liquid	Pressure, kPa - Liquid	Viscosity, Pa*s - Liquid
333.20	10200.00	0.0003650
Reference		https://www.doi.org/10.1007/s10765-008-0542-6

Temperature, K	Pressure, kPa	Viscosity, Pa*s
303.15	101.33	0.0003260
Reference		https://www.doi.org/10.1016/j.tca.2009.07.008

Sources

[illegible]

<https://www.doi.org/10.1016/j.jct.2017.07.012>
<https://www.doi.org/10.1016/j.fluid.2007.01.003>
<https://www.doi.org/10.1016/j.jct.2012.06.005>
<https://www.doi.org/10.1016/j.fluid.2018.11.011>
<https://www.doi.org/10.1021/acs.jced.6b00576>
<https://www.doi.org/10.1016/j.jct.2015.11.005>
<https://www.doi.org/10.1016/j.jct.2013.05.023>
<https://www.doi.org/10.1016/j.jct.2008.03.017>
<https://www.doi.org/10.1021/je050406x>
<https://www.doi.org/10.1016/j.fluid.2017.06.001>
https://en.wikipedia.org/wiki/Joback_method
<https://www.doi.org/10.1016/j.jct.2013.10.038>
<https://www.doi.org/10.1016/j.jct.2016.09.011>
<https://www.doi.org/10.1021/je1008284>
<https://www.doi.org/10.1021/je200609g>
<https://www.doi.org/10.1016/j.jct.2018.01.003>
<https://www.doi.org/10.1016/j.jct.2003.09.005>
<https://www.doi.org/10.1016/j.fluid.2011.07.018>
<https://www.doi.org/10.1021/je400781b>
<https://www.doi.org/10.1016/j.jct.2005.06.003>
<https://www.doi.org/10.1021/je8007099>
<https://www.doi.org/10.1016/j.jct.2018.05.003>
<https://www.doi.org/10.1016/j.tca.2011.01.019>
<https://www.doi.org/10.1016/j.jct.2013.11.027>
<https://www.doi.org/10.1016/j.tca.2011.07.005>
<https://www.doi.org/10.1021/je700497h>
<https://www.doi.org/10.1016/j.jct.2009.06.014>
<https://www.doi.org/10.1021/acs.jced.5b00007>
<https://www.doi.org/10.1021/je800428r>
<https://www.doi.org/10.1021/je060311a>
<https://www.doi.org/10.1016/j.fluid.2014.03.027>
<https://www.doi.org/10.1016/j.jct.2016.07.017>
<https://www.doi.org/10.1016/j.jct.2016.12.002>
<https://www.doi.org/10.1016/j.fluid.2017.12.029>
<https://www.doi.org/10.1016/j.fluid.2018.06.013>
<https://www.doi.org/10.1016/j.fluid.2004.11.006>
<https://www.doi.org/10.1016/j.jct.2011.11.021>
<https://www.doi.org/10.1021/acs.jced.8b01256>

Solubility of 5-Amino Salicylic Acid in Different Solvents at Various Temperatures and activity coefficients at infinite dilution for organic compounds in the ionic liquid 1-hexyl-3-methylimidazolium chloride:

[illegible]

<https://www.doi.org/10.1016/j.ijct.2018.09.003>

<https://www.doi.org/10.1016/j.ijct.2007.09.007>

<https://www.doi.org/10.1016/j.ijct.2013.02.004>

<https://www.doi.org/10.1021/ie034076x>

<https://www.doi.org/10.1016/j.tca.2011.02.004>

<https://www.doi.org/10.1021/acs.iced.7b00585>

<http://www.ddbst.com/en/EED/VLE/VLE%20Acetonitrile%3BTetrahydrofuran.php>

<https://www.doi.org/10.1021/ie101008v>

<https://www.doi.org/10.1016/j.fluid.2007.07.016>

<https://www.doi.org/10.1016/j.ijct.2013.05.008>

<https://www.doi.org/10.1016/j.fluid.2016.02.004>

<https://www.doi.org/10.1021/ie700560s>

<https://www.doi.org/10.1016/j.ijct.2012.03.005>

<https://www.doi.org/10.1016/j.ijct.2006.08.007>

<https://www.doi.org/10.1021/ie900838a>

<https://www.doi.org/10.1016/j.ijct.2010.10.026>

<https://www.doi.org/10.1016/j.ijct.2014.04.024>

<https://www.doi.org/10.1021/ie700426k>

<https://www.doi.org/10.1021/je700338t>

<https://www.doi.org/10.1016/j.fluid.2007.08.013>

<https://www.doi.org/10.1021/ie800218a>

<https://www.doi.org/10.1016/j.ijct.2013.01.005>

<https://www.doi.org/10.1016/j.fluid.2014.04.003>

<https://www.doi.org/10.1016/j.ijct.2014.04.030>

<https://www.doi.org/10.1016/j.tsc.2006.05.010>

<https://www.doi.org/10.1016/j.fluid.2016.11.001>

<https://www.doi.org/10.1016/j.ijct.2013.08.016>

<https://www.doi.org/10.1031/acs.iced.9b00381>

<https://www.doi.org/10.1016/j.fluid.2018.04.004>

<https://www.doi.org/10.1031/acs.jced.9b00564>

<https://www.doi.org/10.1031/acs.joc.3b00551>

<https://www.doi.org/10.1031/acs.joc.3b00340>

3.3] <https://www.doi.org/10.1016/j.fluid.2014.10.041>

<https://www.doi.org/10.1016/j.tsc.2019.03.007>

<https://www.doi.org/10.1016/j.jst.2020.03.000>

<https://www.doi.org/10.1001/ja.10000010>

[illegible]

imide

[illegible]

11. "10-1010" 51-10010-00-050

11. "1994-1995" (1994-1995) (1994-1995) (1994-1995)

[https://doi.org/10.1080/09670262.2021.2001331](#)

http://www.elsevier.com/locate/jbiotec

http://dx.doi.org/10.1016/j.jmbs.2014.07.001

[illegible]

<https://www.doi.org/10.1016/j.fluid.2018.04.009>

<https://www.doi.org/10.1016/j.fluid.2013.07.034>

<https://www.doi.org/10.1016/j.ijct.2005.07.010>

<https://www.doi.org/10.1016/j.ijct.2009.06.011>

<https://www.doi.org/10.1016/j.fluid.2004.10.027>

<https://www.doi.org/10.1016/j.fluid.2012.02.013>

<https://www.doi.org/10.1021/je100315a>

<https://www.doi.org/10.1021/ie034183r>

<https://www.doi.org/10.1016/j.ijct.2016.01.022>

<https://www.doi.org/10.1021/acs.iced.9b00341>

<https://www.doi.org/10.1021/acs.jced.8b00425>

<https://www.doi.org/10.1021/acs.jced.7b00244>

<https://www.doi.org/10.1016/j.fluid.2014.04.006>

<https://www.doi.org/10.1021/acs.jced.7b00584>

<https://www.doi.org/10.1016/j.ijct.2007.10.006>

<https://www.doi.org/10.1021/acs.iced.5b00857>

<https://www.doi.org/10.1021/je100556t>

<https://www.doi.org/10.1021/ie9003178>

<https://www.doi.org/10.1016/j.ijct.2008.12.005>

<https://www.doi.org/10.1021/ie030141r>

<https://www.doi.org/10.1021/acs.iced.9b00696>

<https://www.doi.org/10.1016/j.ijct.2015.05.019>

<https://www.doi.org/10.1016/j.ijct.2013.08.033>

<https://www.doi.org/10.1016/j.ijct.2019.05.011>

<https://www.doi.org/10.1031/je5010565>

<http://webbook.nist.gov/cgi/cheek.cgi?ID=C100000&Units=SI>

<https://www.doi.org/10.1016/j.ijct.2013.03.031>

<https://www.doi.org/10.1016/j.fluid.2015.03.036>

<https://www.doi.org/10.1016/j.ijat.2015.04.025>

<https://www.doi.org/10.1016/j.ijat.2010.05.009>

<https://www.doi.org/10.1031/acs.joc.3b01011>

<https://www.doi.org/10.1016/j.jast.2016.01.017>

<https://www.doi.org/10.1016/j.fluid.2014.01.030>

<https://www.doi.org/10.1001/jco.2024.1534>

<https://www.doi.org/10.1016/j.jst.2019.05.017>

<https://www.doi.org/10.1001/jco.2024.1002>

[illegible]

[View all posts by Dr. David M. Williams](#)

[View all posts by J. L. Brainerd](#)

[illegible]

11. // 11.10.10 // 10.10.10.10

DOI: 10.1002/for

doi:10.1371/journal.pone.0142037.g002

[illegible]

<https://www.doi.org/10.1016/j.jct.2016.07.035>
<https://www.doi.org/10.1016/j.jct.2013.10.025>
<https://www.doi.org/10.1021/acs.jced.7b00851>
<https://www.doi.org/10.1016/j.jct.2005.04.013>
<https://www.doi.org/10.1016/j.jct.2005.07.018>
<https://www.doi.org/10.1021/je800354p>
<https://www.doi.org/10.1021/je700640r>
<https://www.doi.org/10.1016/j.jct.2013.01.007>
<https://www.doi.org/10.1016/j.jct.2017.03.015>
<https://www.doi.org/10.1016/j.fluid.2019.03.023>
<https://www.doi.org/10.1016/j.jct.2011.04.018>
<https://www.doi.org/10.1021/je900629v>
<https://www.doi.org/10.1007/s10765-010-0860-3>
<https://www.doi.org/10.1016/j.jct.2018.07.024>
<https://www.doi.org/10.1016/j.fluid.2012.06.012>
<https://www.doi.org/10.1016/j.jct.2010.10.002>
<https://www.doi.org/10.1016/j.jct.2011.09.022>
<https://www.doi.org/10.1016/j.jct.2013.10.026>
<https://www.doi.org/10.1016/j.jct.2015.05.022>
<https://www.doi.org/10.1016/j.tca.2009.07.008>
<https://www.doi.org/10.1021/je0500577>
<https://www.doi.org/10.1016/j.fluid.2010.10.008>
<https://www.doi.org/10.1021/je1002016>
<https://www.doi.org/10.1016/j.jct.2016.06.028>
<https://www.doi.org/10.1021/je200531k>
<https://www.doi.org/10.1016/j.jct.2017.10.003>
<https://www.doi.org/10.1021/je200822w>
<https://www.doi.org/10.1016/j.jct.2016.10.042>
<https://www.doi.org/10.1021/je0498762>
<https://www.doi.org/10.1021/je9001976>
<https://www.doi.org/10.1021/acs.jced.8b00566>
<https://www.doi.org/10.1016/j.jct.2011.11.025>
<https://www.doi.org/10.1021/je700297c>
<https://www.doi.org/10.1016/j.fluid.2016.10.009>
<https://www.doi.org/10.1016/j.fluid.2016.08.012>
<https://www.doi.org/10.1021/acs.jced.5b00783>
<https://www.doi.org/10.1021/je900547w>
<https://www.doi.org/10.1021/je1005517>
<https://www.doi.org/10.1016/j.jct.2016.09.025>
<https://www.doi.org/10.1016/j.jct.2010.01.004>
<https://www.doi.org/10.1021/acs.jced.8b00080>
<https://www.doi.org/10.1021/acs.jced.7b00005>
<https://www.doi.org/10.1021/je500396b>

[illegible]

phosphoramidic

A volumetric and viscosity study for
the binary mixtures of
KDB-Pure-(Ketone)-Thermophysics
Fenaric Acid Database) some molecular
Measurement of heat capacities of
solvents.
ionic liquids by differential scanning
calorimetry Derived Platform Chemicals:
Thermodynamic Studies on the
Separation of Phthaloxymethane
hexamethylurea;
Equilibrium of Equilibrium of
Dinitrophenylhydrazine based on melting
acid; P. R. Z. Thermophysical Ester in 10
Pure Solvents:

Solubility and Thermodynamic Behavior of Veramoss in Different Pure Solvents and Ethanol-water and propylene carbonate-organic solvents. Measurement and correlation of solubility and mixing properties of dimethyl sulfoxide in binary mixtures with tetrahydrofuran + 1-alkyl-3-methylimidazolium bis(trifluoromethylsulfonylethyl)phosphonium salts. Isentropic compressibilities, refractive indexes, and viscosities of binary mixtures with solutes in water. VII. Cyclic ethers at higher pressures (2000-6500 kPa) of pressures based on compatibility coefficients and the addition of isobaric supercritical fluids of methylcyclopentane, tetrahydrofuran + 1-alkyl-3-methylimidazolium bis(trifluoromethylsulfonylethyl)phosphonium salts

<https://www.doi.org/10.1021/acs.jced.6b00646>
<https://www.doi.org/10.1016/j.jct.2013.01.024>
<https://www.doi.org/10.1016/j.fluid.2013.06.037>
<https://www.doi.org/10.1021/je0497471>
<https://www.doi.org/10.1016/j.tca.2004.08.001>
<https://www.doi.org/10.1016/j.jct.2009.11.005>
<https://www.doi.org/10.1016/j.jct.2018.05.017>
<https://www.doi.org/10.1016/j.fluid.2016.04.012>

Legend

affp:	Proton affinity
basg:	Gas basicity
chg:	Standard gas enthalpy of combustion
chl:	Standard liquid enthalpy of combustion
cpg:	Ideal gas heat capacity
cpl:	Liquid phase heat capacity
dvisc:	Dynamic viscosity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hfust:	Enthalpy of fusion at a given temperature
hvap:	Enthalpy of vaporization at standard conditions
hvapt:	Enthalpy of vaporization at a given temperature
ie:	Ionization energy
kvisc:	Kinematic viscosity
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
pvap:	Vapor pressure
rfi:	Refractive Index
rho:	Liquid Density
rinpol:	Non-polar retention indices
ripol:	Polar retention indices
sfust:	Entropy of fusion at a given temperature
sg:	Molar entropy at standard conditions
sl:	Liquid phase molar entropy at standard conditions
speedsl:	Speed of sound in fluid
srf:	Surface Tension
tb:	Normal Boiling Point Temperature

tbp:	Boiling point at given pressure
tc:	Critical Temperature
tf:	Normal melting (fusion) point
tt:	Triple Point Temperature
vc:	Critical Volume
volm:	Molar Volume

Latest version available from:

<https://www.chemeo.com/cid/66-029-7/Tetrahydrofuran.pdf>

Generated by Cheméo on 2025-12-24 00:47:49.901221305 +0000 UTC m=+6285467.431261967.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.