

# Acetonitrile

Other names:	Acetonitril
	CH3CN
	Cyanomethane
	Cyanure de methyl
	ETHANENITIRILE
	Ethanenitrile
	Ethanonitrile
	Ethyl nitrile
	Methane, cyano-
	Methanecarbonitrile
	Methyl cyanide
	Methylkyanid
	NA 1648
	NCI-C60822
	Rcra waste number U003
	UN 1648
	USAF EK-488
Inchi:	InChI=1S/C2H3N/c1-2-3/h1H3
InchiKey:	WEVYAHXRMPXWCK-UHFFFAOYSA-N
Formula:	C2H3N
SMILES:	CC#N
Mol. weight [g/mol]:	41.05
CAS:	75-05-8

## Physical Properties

Property code	Value	Unit	Source
af	0.3270		KDB
affp	787.40 ± 5.90	kJ/mol	NIST Webbook
affp	779.20	kJ/mol	NIST Webbook
aigt	797.04	K	KDB
basg	748.00	kJ/mol	NIST Webbook
chl	-1256.33 ± 0.30	kJ/mol	NIST Webbook
chl	-1247.20 ± 7.20	kJ/mol	NIST Webbook
chl	-1270.00	kJ/mol	NIST Webbook

cpl	92.36	J/molxK	Volumes, Heat Capacities, and Compressibilities of the Mixtures of Acetonitrile with N,N-Dimethylacetamide and Propylene Carbonate
cpl	90.93	J/molxK	Heat capacities of the mixtures of ionic liquids with acetonitrile
dm	3.50	debye	KDB
ea	0.00 ± 0.01	eV	NIST Webbook
ea	0.01	eV	NIST Webbook
ea	0.01	eV	NIST Webbook
fill	4.40	% in Air	KDB
flu	16.00	% in Air	KDB
fpc	278.71	K	KDB
gf	105.70	kJ/mol	KDB
gyrad	1.8210		KDB
hf	87.92	kJ/mol	KDB
hf	65.86	kJ/mol	NIST Webbook
hf	74.04 ± 0.37	kJ/mol	NIST Webbook
hfl	31.40	kJ/mol	NIST Webbook
hfl	40.56 ± 0.40	kJ/mol	NIST Webbook
hfus	2.44	kJ/mol	Joback Method
hvap	33.00	kJ/mol	NIST Webbook
hvap	33.40	kJ/mol	NIST Webbook
hvap	33.45 ± 0.21	kJ/mol	NIST Webbook
hvap	33.00	kJ/mol	NIST Webbook
hvap	32.94 ± 0.06	kJ/mol	NIST Webbook
ie	12.21	eV	NIST Webbook
ie	15.11	eV	NIST Webbook
ie	12.20 ± 0.01	eV	NIST Webbook
ie	12.20 ± 0.00	eV	NIST Webbook
ie	12.12	eV	NIST Webbook
ie	12.46	eV	NIST Webbook
ie	13.11	eV	NIST Webbook
ie	15.12	eV	NIST Webbook
ie	16.98	eV	NIST Webbook
ie	12.19 ± 0.01	eV	NIST Webbook
ie	12.23 ± 0.05	eV	NIST Webbook
ie	12.38 ± 0.04	eV	NIST Webbook
ie	12.22 ± 0.01	eV	NIST Webbook
ie	12.30 ± 0.25	eV	NIST Webbook
ie	12.33 ± 0.08	eV	NIST Webbook
ie	12.19 ± 0.01	eV	NIST Webbook
ie	12.20	eV	NIST Webbook
ie	12.21 ± 0.00	eV	NIST Webbook

ie	13.14	eV	NIST Webbook
ie	12.20 ± 0.01	eV	NIST Webbook
log10ws	0.26		Aqueous Solubility Prediction Method
log10ws	0.26		Estimated Solubility Method
logp	0.530		Crippen Method
mcvol	40.420	ml/mol	McGowan Method
nfpaf	%!d(float64=3)		KDB
nfpah	%!d(float64=2)		KDB
nfpas	%!d(float64=1)		KDB
pc	4830.00 ± 20.00	kPa	NIST Webbook
pc	4850.00	kPa	KDB
pc	4890.00 ± 10.00	kPa	NIST Webbook
pc	4830.00 ± 81.06	kPa	NIST Webbook
pc	4934.00 ± 3.00	kPa	NIST Webbook
pc	4934.00 ± 3.00	kPa	NIST Webbook
pc	4833.20 ± 81.06	kPa	NIST Webbook
pc	4833.20 ± 81.06	kPa	NIST Webbook
pc	4830.00 ± 20.00	kPa	NIST Webbook
pc	4934.00 ± 20.00	kPa	NIST Webbook
rinpol	457.67		NIST Webbook
rinpol	455.00		NIST Webbook
rinpol	447.00		NIST Webbook
rinpol	455.00		NIST Webbook
rinpol	464.00		NIST Webbook
rinpol	490.00		NIST Webbook
rinpol	455.00		NIST Webbook
rinpol	443.00		NIST Webbook
rinpol	452.00		NIST Webbook
rinpol	443.00		NIST Webbook
rinpol	447.00		NIST Webbook
rinpol	470.00		NIST Webbook
rinpol	456.00		NIST Webbook
rinpol	500.00		NIST Webbook
rinpol	467.00		NIST Webbook
rinpol	446.00		NIST Webbook
rinpol	439.00		NIST Webbook
rinpol	439.00		NIST Webbook
rinpol	447.00		NIST Webbook
rinpol	444.00		NIST Webbook
rinpol	460.00		NIST Webbook
rinpol	440.00		NIST Webbook
rinpol	425.00		NIST Webbook
rinpol	464.00		NIST Webbook

rinpol	460.00		NIST Webbook
rinpol	450.00		NIST Webbook
rinpol	442.00		NIST Webbook
rinpol	432.00		NIST Webbook
rinpol	456.90		NIST Webbook
rinpol	452.35		NIST Webbook
rinpol	452.50		NIST Webbook
rinpol	452.53		NIST Webbook
rinpol	452.92		NIST Webbook
rinpol	453.32		NIST Webbook
rinpol	453.90		NIST Webbook
rinpol	454.52		NIST Webbook
rinpol	455.45		NIST Webbook
rinpol	452.72		NIST Webbook
rinpol	452.90		NIST Webbook
rinpol	452.71		NIST Webbook
rinpol	453.18		NIST Webbook
rinpol	453.70		NIST Webbook
rinpol	455.25		NIST Webbook
rinpol	455.74		NIST Webbook
rinpol	456.69		NIST Webbook
rinpol	439.00		NIST Webbook
rinpol	454.45		NIST Webbook
ripol	1003.00		NIST Webbook
ripol	1002.00		NIST Webbook
ripol	1011.00		NIST Webbook
ripol	1045.00		NIST Webbook
ripol	1010.00		NIST Webbook
ripol	1012.00		NIST Webbook
ripol	1026.00		NIST Webbook
ripol	1045.00		NIST Webbook
ripol	988.00		NIST Webbook
ripol	1030.00		NIST Webbook
ripol	1026.00		NIST Webbook
ripol	1002.00		NIST Webbook
ripol	1026.00		NIST Webbook
ripol	1025.00		NIST Webbook
ripol	1045.00		NIST Webbook
ripol	1013.00		NIST Webbook
ripol	1013.00		NIST Webbook
ripol	1012.00		NIST Webbook
sl	149.62	J/molxK	NIST Webbook

tb	354.25	K	Vapor Liquid Equilibria Measurement of (Methanol + Ethanenitrile + Bis(trifluoromethylsulfonyl) Imide)-Based Ionic Liquids at 101.3 kPa
tb	354.74	K	Isobaric Vapor-Liquid Equilibrium of the Acetonitrile + 1-Propanol + Ionic Liquids at an Atmospheric Pressure
tb	354.65	K	Vapor-Liquid Equilibria for the Ternary System Acetonitrile + 1-Propanol + Dimethyl Sulfoxide and the Corresponding Binary Systems at 101.3 kPa
tb	354.75	K	Isobaric Vapor Liquid Equilibrium for the Acetonitrile + Water System Containing Different Ionic Liquids at Atmospheric Pressure
tb	354.75 ± 0.20	K	NIST Webbook
tb	354.80	K	NIST Webbook
tb	354.25 ± 0.30	K	NIST Webbook
tb	355.00	K	NIST Webbook
tb	354.60 ± 0.30	K	NIST Webbook
tb	354.25 ± 0.50	K	NIST Webbook
tb	370.55 ± 0.10	K	NIST Webbook
tb	354.70	K	NIST Webbook
tb	354.75 ± 0.50	K	NIST Webbook
tb	354.70	K	NIST Webbook
tb	354.80 ± 0.40	K	NIST Webbook
tb	354.80 ± 0.50	K	NIST Webbook
tb	354.75 ± 0.20	K	NIST Webbook
tb	354.80 ± 0.30	K	NIST Webbook
tb	354.75	K	NIST Webbook
tb	354.45 ± 0.50	K	NIST Webbook
tb	354.65 ± 0.50	K	NIST Webbook
tb	354.80 ± 0.30	K	NIST Webbook
tb	355.00 ± 0.50	K	NIST Webbook
tb	351.00 ± 3.00	K	NIST Webbook
tb	354.71 ± 0.20	K	NIST Webbook
tb	354.35 ± 0.50	K	NIST Webbook
tb	354.70 ± 2.00	K	NIST Webbook
tb	355.00 ± 1.50	K	NIST Webbook
tb	354.72 ± 0.10	K	NIST Webbook
tb	355.00 ± 2.00	K	NIST Webbook
tb	355.00 ± 0.50	K	NIST Webbook
tb	354.90 ± 0.30	K	NIST Webbook

tb	354.70 ± 2.00	K	NIST Webbook
tb	355.15 ± 1.50	K	NIST Webbook
tb	354.80 ± 0.50	K	NIST Webbook
tb	354.00 ± 2.00	K	NIST Webbook
tb	355.00 ± 2.00	K	NIST Webbook
tb	354.90 ± 0.70	K	NIST Webbook
tb	353.00 ± 2.00	K	NIST Webbook
tb	354.70	K	Acetonitrile Dehydration via Extractive Distillation Using Low Transition Temperature Mixtures as Entrainers
tb	353.90 ± 1.50	K	NIST Webbook
tb	354.80 ± 0.30	K	NIST Webbook
tb	354.80 ± 0.30	K	NIST Webbook
tb	354.90 ± 0.20	K	NIST Webbook
tb	354.80 ± 0.30	K	NIST Webbook
tb	354.80 ± 0.60	K	NIST Webbook
tb	354.80 ± 0.40	K	NIST Webbook
tb	354.75 ± 0.30	K	NIST Webbook
tb	354.95 ± 1.50	K	NIST Webbook
tb	354.75 ± 0.30	K	NIST Webbook
tb	354.70 ± 0.40	K	NIST Webbook
tb	354.75 ± 0.20	K	NIST Webbook
tb	354.68	K	Measurement and correlation of isobaric vapour-liquid equilibrium for the (acetonitrile + water) system containing different ionic liquids at atmospheric pressure
tb	354.69 ± 0.30	K	NIST Webbook
tb	354.90 ± 0.60	K	NIST Webbook
tb	354.69 ± 0.50	K	NIST Webbook
tb	355.00 ± 1.50	K	NIST Webbook
tb	354.65	K	(Vapour + liquid) equilibria in the ternary system (acetonitrile + n-propanol + ethylene glycol) and corresponding binary systems at 101.3 kPa
tb	355.00 ± 2.00	K	NIST Webbook
tb	354.80	K	KDB
tb	354.65	K	The isobaric vapor liquid equilibria of ethyl acetate p acetonitrile p bis(trifluoromethylsulfonyl)imide-based ionic liquids at 101.3 kPa
tc	543.45 ± 1.00	K	NIST Webbook
tc	545.50	K	KDB
tc	544.99 ± 0.20	K	NIST Webbook

tc	544.99 ± 0.20	K	NIST Webbook
tc	544.99 ± 0.20	K	NIST Webbook
tc	543.20 ± 2.00	K	NIST Webbook
tc	543.15 ± 1.00	K	NIST Webbook
tc	545.50 ± 0.20	K	NIST Webbook
tc	545.50 ± 0.30	K	NIST Webbook
tc	545.50 ± 0.30	K	NIST Webbook
tf	227.45	K	Aqueous Solubility Prediction Method
tf	229.32	K	KDB
tf	229.07	K	Efficient determination of crystallisation and melting points at low cooling and heating rates with novel computer controlled equipment
tf	230.42	K	Experimental (Solid + Liquid) and (Liquid + Liquid) Equilibria and Excess Molar Volume of Alkanol + Acetonitrile, Propanenitrile, and Butanenitrile Mixtures
tt	229.32 ± 0.02	K	NIST Webbook
vc	0.173	m3/kmol	KDB
zc	0.1849930		KDB
zra	0.20		KDB

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	59.46	J/molxK	379.59	Joback Method
cpg	62.33	J/molxK	411.95	Joback Method
cpg	65.11	J/molxK	444.30	Joback Method
cpg	67.79	J/molxK	476.66	Joback Method
cpg	70.38	J/molxK	509.01	Joback Method
cpg	72.87	J/molxK	541.37	Joback Method
cpg	56.48	J/molxK	347.24	Joback Method
cpl	92.21	J/molxK	303.15	Volumetric Properties, Viscosities, and Isobaric Heat Capacities of Imidazolium Octanoate Protic Ionic Liquid in Molecular Solvents

cpl	92.40	J/molxK	308.15	Volumetric Properties, Viscosities, and Isobaric Heat Capacities of Imidazolium Octanoate Protic Ionic Liquid in Molecular Solvents
cpl	82.50	J/molxK	297.00	NIST Webbook
cpl	91.69	J/molxK	298.15	Volumetric Properties, Viscosities, and Isobaric Heat Capacities of Imidazolium Octanoate Protic Ionic Liquid in Molecular Solvents
cpl	91.70	J/molxK	298.15	NIST Webbook
cpl	91.70	J/molxK	298.15	NIST Webbook
cpl	91.70	J/molxK	298.15	NIST Webbook
cpl	81.80	J/molxK	303.15	NIST Webbook
cpl	77.40	J/molxK	298.15	NIST Webbook
cpl	92.82	J/molxK	313.15	Volumetric Properties, Viscosities, and Isobaric Heat Capacities of Imidazolium Octanoate Protic Ionic Liquid in Molecular Solvents
cpl	91.69	J/molxK	298.15	NIST Webbook
cpl	91.46	J/molxK	298.15	NIST Webbook
dvisc	0.0003270	Paxs	313.15	Properties of Pure 1-Butyl-2,3-dimethylimidazolium Tetrafluoroborate Ionic Liquid and Its Binary Mixtures with Dimethyl Sulfoxide and Acetonitrile
dvisc	0.0003900	Paxs	293.15	Viscosities of 1-Hexyl-3-methylimidazolium Tetrafluoroborate and Its Binary Mixtures with Dimethyl Sulfoxide and Acetonitrile



dvisc	0.0003500	Paxs	303.15	Viscosities of 1-Hexyl-3-methylimidazolium Tetrafluoroborate and Its Binary Mixtures with Dimethyl Sulfoxide and Acetonitrile
dvisc	0.0003300	Paxs	313.15	Viscosities of 1-Hexyl-3-methylimidazolium Tetrafluoroborate and Its Binary Mixtures with Dimethyl Sulfoxide and Acetonitrile
dvisc	0.0003000	Paxs	323.15	Viscosities of 1-Hexyl-3-methylimidazolium Tetrafluoroborate and Its Binary Mixtures with Dimethyl Sulfoxide and Acetonitrile
dvisc	0.0002800	Paxs	333.15	Viscosities of 1-Hexyl-3-methylimidazolium Tetrafluoroborate and Its Binary Mixtures with Dimethyl Sulfoxide and Acetonitrile
dvisc	0.0002600	Paxs	343.15	Viscosities of 1-Hexyl-3-methylimidazolium Tetrafluoroborate and Its Binary Mixtures with Dimethyl Sulfoxide and Acetonitrile
dvisc	0.0003443	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	0.0003124	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	0.0002890	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	0.0002891	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	0.0003125	Paxs	308.15	Electrical Conductances of 1-Butyl-3-propylimidazolium Bromide and 1-Butyl-3-propylbenzimidazolium Bromide in Water, Methanol, and Acetonitrile at (308, 313, and 318) K at 0.1 MPa	
dvisc	0.0003042	Paxs	313.15	Electrical Conductances of 1-Butyl-3-propylimidazolium Bromide and 1-Butyl-3-propylbenzimidazolium Bromide in Water, Methanol, and Acetonitrile at (308, 313, and 318) K at 0.1 MPa	

dvisc	0.0002903	Paxs	318.15	Electrical Conductances of 1-Butyl-3-propylimidazolium Bromide and 1-Butyl-3-propylbenzimidazolium Bromide in Water, Methanol, and Acetonitrile at (308, 313, and 318) K at 0.1 MPa
dvisc	0.0003700	Paxs	298.15	Properties of Pure 1-Butyl-2,3-dimethylimidazolium Tetrafluoroborate Ionic Liquid and Its Binary Mixtures with Dimethyl Sulfoxide and Acetonitrile
dvisc	0.0003540	Paxs	303.15	Properties of Pure 1-Butyl-2,3-dimethylimidazolium Tetrafluoroborate Ionic Liquid and Its Binary Mixtures with Dimethyl Sulfoxide and Acetonitrile
dvisc	0.0003414	Paxs	308.15	Molecular interaction studies and theoretical estimation of ultrasonic speeds using scaled particle theory in binary mixtures of toluene with homologous nitriles at different temperatures
dvisc	0.0003020	Paxs	323.15	Properties of Pure 1-Butyl-2,3-dimethylimidazolium Tetrafluoroborate Ionic Liquid and Its Binary Mixtures with Dimethyl Sulfoxide and Acetonitrile

dvisc	0.0002810	Paxs	333.15	Properties of Pure 1-Butyl-2,3-dimethylimidazolium Tetrafluoroborate Ionic Liquid and Its Binary Mixtures with Dimethyl Sulfoxide and Acetonitrile	
dvisc	0.0002620	Paxs	343.15	Properties of Pure 1-Butyl-2,3-dimethylimidazolium Tetrafluoroborate Ionic Liquid and Its Binary Mixtures with Dimethyl Sulfoxide and Acetonitrile	
dvisc	0.0003440	Paxs	298.15	Viscosity, Density, and Speed of Sound for the Binary Mixtures of Formamide with 2-Methoxyethanol, Acetophenone, Acetonitrile, 1,2-Dimethoxyethane, and Dimethylsulfoxide at Different Temperatures	
dvisc	0.0003554	Paxs	303.15	Molecular interaction studies and theoretical estimation of ultrasonic speeds using scaled particle theory in binary mixtures of toluene with homologous nitriles at different temperatures	
dvisc	0.0002890	Paxs	318.15	Viscosity, Density, and Speed of Sound for the Binary Mixtures of Formamide with 2-Methoxyethanol, Acetophenone, Acetonitrile, 1,2-Dimethoxyethane, and Dimethylsulfoxide at Different Temperatures	

dvisc	0.0003417	Paxs	298.15	Density and Viscosity of Anhydrous Mixtures of Dimethylsulfoxide with Acetonitrile in the Range (298.15 to 318.15) K
dvisc	0.0003280	Paxs	303.15	Density and Viscosity of Anhydrous Mixtures of Dimethylsulfoxide with Acetonitrile in the Range (298.15 to 318.15) K
dvisc	0.0003129	Paxs	308.15	Density and Viscosity of Anhydrous Mixtures of Dimethylsulfoxide with Acetonitrile in the Range (298.15 to 318.15) K
dvisc	0.0003009	Paxs	313.15	Density and Viscosity of Anhydrous Mixtures of Dimethylsulfoxide with Acetonitrile in the Range (298.15 to 318.15) K
dvisc	0.0002899	Paxs	318.15	Density and Viscosity of Anhydrous Mixtures of Dimethylsulfoxide with Acetonitrile in the Range (298.15 to 318.15) K
dvisc	0.0003081	Paxs	293.15	Thermodynamic Study of Phenyl Salicylate Solutions in Aprotic Solvents at Different Temperatures
dvisc	0.0002857	Paxs	298.15	Thermodynamic Study of Phenyl Salicylate Solutions in Aprotic Solvents at Different Temperatures

dvisc	0.0002642	Paxs	303.15	Thermodynamic Study of Phenyl Salicylate Solutions in Aprotic Solvents at Different Temperatures	
dvisc	0.0002444	Paxs	308.15	Thermodynamic Study of Phenyl Salicylate Solutions in Aprotic Solvents at Different Temperatures	
dvisc	0.0002374	Paxs	310.15	Thermodynamic Study of Phenyl Salicylate Solutions in Aprotic Solvents at Different Temperatures	
dvisc	0.0002267	Paxs	313.15	Thermodynamic Study of Phenyl Salicylate Solutions in Aprotic Solvents at Different Temperatures	
dvisc	0.0003696	Paxs	298.15	Molecular interaction studies and theoretical estimation of ultrasonic speeds using scaled particle theory in binary mixtures of toluene with homologous nitriles at different temperatures	
dvisc	0.0002620	Paxs	343.15	Viscosity of binary mixtures of 1-ethyl-3-methylimidazolium tetrafluoroborate ionic liquid with four organic solvents	
dvisc	0.0002810	Paxs	333.15	Viscosity of binary mixtures of 1-ethyl-3-methylimidazolium tetrafluoroborate ionic liquid with four organic solvents	

dvisc	0.0003020	Paxs	323.15	Viscosity of binary mixtures of 1-ethyl-3-methylimidazolium tetrafluoroborate ionic liquid with four organic solvents
dvisc	0.0003270	Paxs	313.15	Viscosity of binary mixtures of 1-ethyl-3-methylimidazolium tetrafluoroborate ionic liquid with four organic solvents
dvisc	0.0003540	Paxs	303.15	Viscosity of binary mixtures of 1-ethyl-3-methylimidazolium tetrafluoroborate ionic liquid with four organic solvents
dvisc	0.0003700	Paxs	298.15	Viscosity of binary mixtures of 1-ethyl-3-methylimidazolium tetrafluoroborate ionic liquid with four organic solvents
dvisc	0.0003860	Paxs	293.15	Viscosity of binary mixtures of 1-ethyl-3-methylimidazolium tetrafluoroborate ionic liquid with four organic solvents
dvisc	0.0003009	Paxs	313.15	The (water + acetonitrile) mixture revisited: A new approach for calculating partial molar volumes
dvisc	0.0003129	Paxs	308.15	The (water + acetonitrile) mixture revisited: A new approach for calculating partial molar volumes
dvisc	0.0003280	Paxs	303.15	The (water + acetonitrile) mixture revisited: A new approach for calculating partial molar volumes

dvisc	0.0003417	Paxs	298.15	The (water + acetonitrile) mixture revisited: A new approach for calculating partial molar volumes
dvisc	0.0002680	Paxs	328.15	Volumetric properties of ionic liquid 1,3-dimethylimidazolium methyl sulfate + molecular solvents at T = (298.15 - 328.15) K
dvisc	0.0002918	Paxs	318.15	Volumetric properties of ionic liquid 1,3-dimethylimidazolium methyl sulfate + molecular solvents at T = (298.15 - 328.15) K
dvisc	0.0003171	Paxs	308.15	Volumetric properties of ionic liquid 1,3-dimethylimidazolium methyl sulfate + molecular solvents at T = (298.15 - 328.15) K
dvisc	0.0003431	Paxs	298.15	Volumetric properties of ionic liquid 1,3-dimethylimidazolium methyl sulfate + molecular solvents at T = (298.15 - 328.15) K
dvisc	0.0002899	Paxs	318.15	The (water + acetonitrile) mixture revisited: A new approach for calculating partial molar volumes
dvisc	0.0003130	Paxs	308.15	Viscosity, Density, and Speed of Sound for the Binary Mixtures of Formamide with 2-Methoxyethanol, Acetophenone, Acetonitrile, 1,2-Dimethoxyethane, and Dimethylsulfoxide at Different Temperatures



dvisc	0.0003700	Paxs	298.15	Viscosities of 1-Hexyl-3-methylimidazolium Tetrafluoroborate and Its Binary Mixtures with Dimethyl Sulfoxide and Acetonitrile
hfust	8.17	kJ/mol	229.30	NIST Webbook
hvapt	34.20	kJ/mol	298.00	NIST Webbook
hvapt	34.80	kJ/mol	321.00	NIST Webbook
hvapt	33.30	kJ/mol	334.50	NIST Webbook
hvapt	29.75	kJ/mol	354.70	NIST Webbook
hvapt	33.23	kJ/mol	298.15	NIST Webbook
hvapt	31.38	kJ/mol	352.80	KDB
hvapt	33.90	kJ/mol	290.00	NIST Webbook
hvapt	33.80	kJ/mol	325.00	NIST Webbook
pvap	4.54	kPa	278.88	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	55.46	kPa	336.56	Vapor Liquid Equilibrium, Excess Molar Enthalpies, and Excess Molar Volumes of Binary Mixtures Containing 2-Ethoxy-2-methylpropane or 2-Ethoxy-2-methylbutane and Acetonitrile or Propanenitrile
pvap	60.76	kPa	339.20	Vapor Liquid Equilibrium, Excess Molar Enthalpies, and Excess Molar Volumes of Binary Mixtures Containing 2-Ethoxy-2-methylpropane or 2-Ethoxy-2-methylbutane and Acetonitrile or Propanenitrile

pvap	64.81	kPa	341.07	Vapor Liquid Equilibrium, Excess Molar Enthalpies, and Excess Molar Volumes of Binary Mixtures Containing 2-Ethoxy-2-methylpropane or 2-Ethoxy-2-methylbutane and Acetonitrile or Propanenitrile
pvap	71.24	kPa	343.86	Vapor Liquid Equilibrium, Excess Molar Enthalpies, and Excess Molar Volumes of Binary Mixtures Containing 2-Ethoxy-2-methylpropane or 2-Ethoxy-2-methylbutane and Acetonitrile or Propanenitrile
pvap	75.40	kPa	345.58	Vapor Liquid Equilibrium, Excess Molar Enthalpies, and Excess Molar Volumes of Binary Mixtures Containing 2-Ethoxy-2-methylpropane or 2-Ethoxy-2-methylbutane and Acetonitrile or Propanenitrile
pvap	80.49	kPa	347.55	Vapor Liquid Equilibrium, Excess Molar Enthalpies, and Excess Molar Volumes of Binary Mixtures Containing 2-Ethoxy-2-methylpropane or 2-Ethoxy-2-methylbutane and Acetonitrile or Propanenitrile
pvap	86.34	kPa	349.69	Vapor Liquid Equilibrium, Excess Molar Enthalpies, and Excess Molar Volumes of Binary Mixtures Containing 2-Ethoxy-2-methylpropane or 2-Ethoxy-2-methylbutane and Acetonitrile or Propanenitrile

pvap	92.16	kPa	351.76	Vapor Liquid Equilibrium, Excess Molar Enthalpies, and Excess Molar Volumes of Binary Mixtures Containing 2-Ethoxy-2-methylpropane or 2-Ethoxy-2-methylbutane and Acetonitrile or Propanenitrile
pvap	99.40	kPa	354.12	Vapor Liquid Equilibrium, Excess Molar Enthalpies, and Excess Molar Volumes of Binary Mixtures Containing 2-Ethoxy-2-methylpropane or 2-Ethoxy-2-methylbutane and Acetonitrile or Propanenitrile
pvap	101.30	kPa	354.75	Isobaric Vapor Liquid Equilibrium for the Acetonitrile + Water System Containing Different Ionic Liquids at Atmospheric Pressure
pvap	1096.90	kPa	453.15	Vapor-Liquid Equilibria on Seven Binary Systems: Ethylene Oxide + 2-Methylpropane; Acetophenone + Phenol; cis-1,3-Dichloropropene + 1,2-Dichloropropane; 1,5-Hexadiene + Allyl Chloride; Isopropyl Acetate + Acetonitrile; Vinyl Chloride + Methyl Chloride; and 1,4-Butanediol + c-Butyrolactone

pvap	297.37	kPa	393.15	Vapor-Liquid Equilibria on Seven Binary Systems: Ethylene Oxide + 2-Methylpropane; Acetophenone + Phenol; cis-1,3-Dichloropropene + 1,2-Dichloropropane; 1,5-Hexadiene + Allyl Chloride; Isopropyl Acetate + Acetonitrile; Vinyl Chloride + Methyl Chloride; and 1,4-Butanediol + c-Butyrolactone
pvap	49.16	kPa	333.15	Vapor-Liquid Equilibria on Seven Binary Systems: Ethylene Oxide + 2-Methylpropane; Acetophenone + Phenol; cis-1,3-Dichloropropene + 1,2-Dichloropropane; 1,5-Hexadiene + Allyl Chloride; Isopropyl Acetate + Acetonitrile; Vinyl Chloride + Methyl Chloride; and 1,4-Butanediol + c-Butyrolactone
pvap	4174.61	kPa	535.03	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	4001.46	kPa	532.08	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	3760.37	kPa	527.79	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	3512.89	kPa	523.15	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry

pvap	3254.08	kPa	518.01	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	2999.22	kPa	512.61	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	2747.95	kPa	506.92	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	2523.66	kPa	501.48	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	2303.51	kPa	495.75	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	2110.77	kPa	490.37	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	1923.88	kPa	484.77	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	1749.84	kPa	479.15	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	49.76	kPa	333.51	Vapor Liquid Equilibrium, Excess Molar Enthalpies, and Excess Molar Volumes of Binary Mixtures Containing 2-Ethoxy-2-methylpropane or 2-Ethoxy-2-methylbutane and Acetonitrile or Propanenitrile

pvap	43.71	kPa	329.94	Vapor Liquid Equilibrium, Excess Molar Enthalpies, and Excess Molar Volumes of Binary Mixtures Containing 2-Ethoxy-2-methylpropane or 2-Ethoxy-2-methylbutane and Acetonitrile or Propanenitrile
pvap	37.80	kPa	326.03	Vapor Liquid Equilibrium, Excess Molar Enthalpies, and Excess Molar Volumes of Binary Mixtures Containing 2-Ethoxy-2-methylpropane or 2-Ethoxy-2-methylbutane and Acetonitrile or Propanenitrile
pvap	27.19	kPa	317.53	Vapor Liquid Equilibrium, Excess Molar Enthalpies, and Excess Molar Volumes of Binary Mixtures Containing 2-Ethoxy-2-methylpropane or 2-Ethoxy-2-methylbutane and Acetonitrile or Propanenitrile
pvap	18.59	kPa	308.32	Vapor Liquid Equilibrium, Excess Molar Enthalpies, and Excess Molar Volumes of Binary Mixtures Containing 2-Ethoxy-2-methylpropane or 2-Ethoxy-2-methylbutane and Acetonitrile or Propanenitrile
pvap	49.76	kPa	333.57	Vapor Liquid Equilibrium, Excess Molar Enthalpies, and Excess Molar Volumes of Binary Mixtures Containing 2-Ethoxy-2-methylpropane or 2-Ethoxy-2-methylbutane and Acetonitrile or Propanenitrile

pvap	71.24	kPa	343.94	Vapor Liquid Equilibrium, Excess Molar Enthalpies, and Excess Molar Volumes of Binary Mixtures Containing 2-Ethoxy-2-methylpropane or 2-Ethoxy-2-methylbutane and Acetonitrile or Propanenitrile
pvap	33.86	kPa	323.15	Vapor Pressures for the Acetonitrile + Tetrabutylammonium Bromide, Water + Tetrabutylammonium Bromide, and Acetonitrile + Water + Tetrabutylammonium Bromide Systems
pvap	101.30	kPa	354.65	The isobaric vapor liquid equilibria of ethyl acetate p acetonitrile p bis(trifluoromethylsulfonyl)imide-based ionic liquids at 101.3 kPa
pvap	16.47	kPa	305.51	Experimental determination and prediction of gas solubility data for oxygen in acetonitrile
pvap	33.71	kPa	323.12	Experimental determination and prediction of gas solubility data for oxygen in acetonitrile
pvap	96.17	kPa	353.13	Experimental determination and prediction of gas solubility data for oxygen in acetonitrile
pvap	4.44	kPa	278.72	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran

pvap	1602.63	kPa	474.05	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	4.51	kPa	278.98	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	4.73	kPa	279.81	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	4.86	kPa	280.27	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	5.34	kPa	281.91	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	5.64	kPa	283.17	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran



pvap	5.84	kPa	283.56	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	5.92	kPa	283.81	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	6.10	kPa	284.56	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	6.47	kPa	285.70	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	6.59	kPa	286.17	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	6.75	kPa	286.61	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran

pvap	7.50	kPa	288.72	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	7.56	kPa	288.80	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	7.59	kPa	288.83	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	7.80	kPa	289.39	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	8.03	kPa	290.09	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	8.30	kPa	290.60	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran

pvap	8.60	kPa	291.36	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	9.16	kPa	292.48	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	9.41	kPa	293.21	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	9.82	kPa	293.92	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	10.03	kPa	294.40	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	10.43	kPa	295.33	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran

pvap	10.77	kPa	295.94	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	11.02	kPa	296.44	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	11.60	kPa	297.48	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	12.05	kPa	298.30	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	12.17	kPa	298.67	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	12.61	kPa	299.42	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran

pvap	13.61	kPa	301.19	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	14.18	kPa	302.14	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	15.57	kPa	304.16	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	16.10	kPa	305.01	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	16.96	kPa	306.04	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	17.53	kPa	306.90	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran

pvap	18.22	kPa	307.95	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	19.85	kPa	309.92	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	20.49	kPa	310.71	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	21.38	kPa	311.79	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	21.55	kPa	311.88	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	22.50	kPa	312.81	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran

pvap	23.42	kPa	313.77	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	24.17	kPa	314.79	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	25.10	kPa	315.71	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	25.15	kPa	315.79	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	26.11	kPa	316.63	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	27.21	kPa	317.72	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran

pvap	28.27	kPa	318.66	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	29.39	kPa	319.69	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	30.52	kPa	320.67	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	31.73	kPa	321.65	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	32.95	kPa	322.54	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	32.94	kPa	322.56	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran



pvap	33.19	kPa	322.64	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	34.49	kPa	323.64	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	35.65	kPa	324.59	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	37.00	kPa	325.58	Vapor pressures and activity coefficients of binary mixtures of 1-ethyl-3-methylimidazolium bis(trifluoromethylsulfonyl)imide with acetonitrile and tetrahydrofuran
pvap	101.32	kPa	354.25	Vapor Liquid Equilibria Measurement of (Methanol + Ethanenitrile + Bis(trifluoromethylsulfonyl) Imide)-Based Ionic Liquids at 101.3 kPa
pvap	10.00	kPa	295.25	Phase Behavior of Binary Mixtures Containing Succinic Acid or Its Esters
pvap	10.00	kPa	295.40	Phase Behavior of Binary Mixtures Containing Succinic Acid or Its Esters

pvap	101.32	kPa	354.70	Acetonitrile Dehydration via Extractive Distillation Using Low Transition Temperature Mixtures as Entrainers
pvap	1450.28	kPa	468.37	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	131.21	kPa	363.15	Vapor-Liquid Equilibria for Four Binary Systems at 363.15 K: N-Methylformamide + Hexane, + Benzene, + Chlorobenzene, and + Acetonitrile
pvap	14.15	kPa	302.25	Vapor-Liquid Equilibrium for Acetonitrile + Propanenitrile and 1-Pentanamine + 1-Methoxy-2-propanol
pvap	17.52	kPa	307.09	Vapor-Liquid Equilibrium for Acetonitrile + Propanenitrile and 1-Pentanamine + 1-Methoxy-2-propanol
pvap	21.13	kPa	311.49	Vapor-Liquid Equilibrium for Acetonitrile + Propanenitrile and 1-Pentanamine + 1-Methoxy-2-propanol
pvap	22.65	kPa	313.15	Vapor-Liquid Equilibrium for Acetonitrile + Propanenitrile and 1-Pentanamine + 1-Methoxy-2-propanol
pvap	25.72	kPa	316.27	Vapor-Liquid Equilibrium for Acetonitrile + Propanenitrile and 1-Pentanamine + 1-Methoxy-2-propanol

pvap	27.78	kPa	318.18	Vapor-Liquid Equilibrium for Acetonitrile + Propanenitrile and 1-Pentanamine + 1-Methoxy-2-propanol
pvap	30.64	kPa	320.65	Vapor-Liquid Equilibrium for Acetonitrile + Propanenitrile and 1-Pentanamine + 1-Methoxy-2-propanol
pvap	33.80	kPa	323.16	Vapor-Liquid Equilibrium for Acetonitrile + Propanenitrile and 1-Pentanamine + 1-Methoxy-2-propanol
pvap	37.90	kPa	326.16	Vapor-Liquid Equilibrium for Acetonitrile + Propanenitrile and 1-Pentanamine + 1-Methoxy-2-propanol
pvap	41.38	kPa	328.50	Vapor-Liquid Equilibrium for Acetonitrile + Propanenitrile and 1-Pentanamine + 1-Methoxy-2-propanol
pvap	45.62	kPa	331.15	Vapor-Liquid Equilibrium for Acetonitrile + Propanenitrile and 1-Pentanamine + 1-Methoxy-2-propanol
pvap	49.06	kPa	333.15	Vapor-Liquid Equilibrium for Acetonitrile + Propanenitrile and 1-Pentanamine + 1-Methoxy-2-propanol
pvap	54.01	kPa	335.84	Vapor-Liquid Equilibrium for Acetonitrile + Propanenitrile and 1-Pentanamine + 1-Methoxy-2-propanol
pvap	58.58	kPa	338.15	Vapor-Liquid Equilibrium for Acetonitrile + Propanenitrile and 1-Pentanamine + 1-Methoxy-2-propanol

pvap	69.62	kPa	343.18	Vapor-Liquid Equilibrium for Acetonitrile + Propanenitrile and 1-Pentanamine + 1-Methoxy-2-propanol	
pvap	82.37	kPa	348.25	Vapor-Liquid Equilibrium for Acetonitrile + Propanenitrile and 1-Pentanamine + 1-Methoxy-2-propanol	
pvap	96.34	kPa	353.10	Vapor-Liquid Equilibrium for Acetonitrile + Propanenitrile and 1-Pentanamine + 1-Methoxy-2-propanol	
pvap	4.32	kPa	277.92	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry	
pvap	4.49	kPa	278.62	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry	
pvap	5.25	kPa	281.57	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry	
pvap	6.38	kPa	285.38	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry	
pvap	7.17	kPa	287.67	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry	
pvap	8.30	kPa	290.65	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry	
pvap	9.18	kPa	292.75	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry	
pvap	10.60	kPa	295.78	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry	

pvap	13.37	kPa	300.82	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	15.27	kPa	303.81	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	19.64	kPa	309.64	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	24.97	kPa	315.43	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	31.31	kPa	321.12	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	36.39	kPa	325.02	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	45.91	kPa	331.27	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	55.17	kPa	336.41	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	65.07	kPa	341.18	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	75.44	kPa	345.57	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	85.31	kPa	349.33	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	95.59	kPa	352.88	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry

pvap	100.75	kPa	354.55	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	101.12	kPa	354.67	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	110.61	kPa	357.56	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	122.63	kPa	360.94	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	125.13	kPa	361.61	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	155.33	kPa	368.97	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	168.12	kPa	371.74	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	175.04	kPa	373.17	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	204.59	kPa	378.81	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	235.79	kPa	384.11	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	306.28	kPa	394.29	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	399.50	kPa	405.24	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry

pvap	502.67	kPa	415.21	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	605.60	kPa	423.68	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	708.99	kPa	431.12	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	804.86	kPa	437.30	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	910.82	kPa	443.50	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	1025.47	kPa	449.60	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	1151.97	kPa	455.74	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	1290.20	kPa	461.87	Vapor Pressures of Acetonitrile Determined by Comparative Ebulliometry
pvap	101.30	kPa	354.74	Isobaric Vapor-Liquid Equilibrium of the Acetonitrile + 1-Propanol + Ionic Liquids at an Atmospheric Pressure
rfi	1.34390		293.15	A novel static analytical apparatus for phase equilibrium measurements
rfi	1.34139		298.15	Isothermal Vapor-Liquid Equilibrium of Binary Mixtures Containing 1-Chlorobutane, Ethanol, or Acetonitrile

rfi	1.34410	295.10	Ternary Liquid-Liquid Equilibria of Acetonitrile and Water with Heptanoic Acid and Nonanol at 323.15 K and 1 atm
rfi	1.34190	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	1.34386	293.15	Ternary Liquid Liquid Equilibrium Data for the Water + Acetonitrile + {Butan-1-ol or 2-Methylpropan-1-ol} Systems at (303.2, 323.2, 343.2) K and 1 atm
rfi	1.34160	298.15	Acoustic and Volumetric Properties of Binary Mixtures of Ionic Liquid 1-Butyl-3-methylimidazolium Bis(trifluoromethylsulfonyl)imide with Acetonitrile and Tetrahydrofuran
rfi	1.34250	293.15	Isobaric Vapor Liquid Equilibrium for the Extractive Distillation of Acetonitrile + Water Mixtures Using Dimethyl Sulfoxide at 101.3 kPa
rfi	1.34300	293.15	Solubilities of Phosphorus-Containing Compounds in Selected Solvents
rfi	1.34300	293.15	Solubilities of Some Phosphaspirocyclic Compounds in Selected Solvents



rfi	1.34180	298.15	Vapor-Liquid Equilibrium Data for the Binary Methyl Esters (Butyrate, Pentanoate, and Hexanoate) (1) + Acetonitrile (2) Systems at 93.32 kPa
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rfi	1.34180	298.15	Physics and Chemistry of Lithium Halides in 1,3-Dioxolane and Its Binary Mixtures with Acetonitrile probed by Conductometric, Volumetric, Viscometric, Refractometric and Acoustic Study
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rfi	1.33710	308.15	Densities, viscosities, excess molar volumes, and refractive indices of acetonitrile and 2-alkanols binary mixtures at different temperatures: Experimental results and application of the Prigogine Flory Patterson theory
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rfi	1.33900	303.15	Densities, viscosities, excess molar volumes, and refractive indices of acetonitrile and 2-alkanols binary mixtures at different temperatures: Experimental results and application of the Prigogine Flory Patterson theory
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rfi	1.34110	298.15	Densities, viscosities, excess molar volumes, and refractive indices of acetonitrile and 2-alkanols binary mixtures at different temperatures: Experimental results and application of the Prigogine Flory Patterson theory
rfi	1.34360	293.15	Densities, viscosities, excess molar volumes, and refractive indices of acetonitrile and 2-alkanols binary mixtures at different temperatures: Experimental results and application of the Prigogine Flory Patterson theory
rfi	1.34420	293.15	ERAS modeling of the excess molar enthalpies of binary liquid mixtures of 1-pentanol and 1-hexanol with acetonitrile at atmospheric pressure and 288, 298, 313 and 323K
rfi	1.34390	293.15	Experimental solubility for betulin and estrone in various solvents within the temperature range $T = (293.2 \text{ to } 328.2) \text{ K}$
rfi	1.35900	298.15	Volumetric properties, viscosity and refractive index of the protic ionic liquid, pyrrolidinium octanoate, in molecular solvents

rfi	1.34410		293.15	Experimental study and ERAS modeling of the excess molar enthalpy of (acetonitrile + 1-heptanol or 1-octanol) mixtures at (298.15, 313.15, and 323.15) K and atmospheric pressure
rfi	1.34160		298.15	(Vapor + liquid) equilibrium of the binary mixtures formed by acetonitrile with selected compounds at 95.5 kPa
rfi	1.34160		298.15	Excess Gibbs free energies of the binary mixtures of acetonitrile with butanols at 94.83 kPa
rfi	1.33960		303.15	Liquid liquid equilibria measurements of ternary systems (acetonitrile + a carboxylic acid + dodecane) at 303.15 K
rhoI	776.85	kg/m3	298.15	Compressibility Studies of Some Copper(I), Silver(I), and Tetrabutylammonium Salts in Acetonitrile + Adiponitrile Binary Mixtures
rhoI	761.45	kg/m3	308.15	Densities and Viscosities of Ternary N,N'-Bis(2-pyridylmethylidene)-1,2-diiminoethane Schiff Base + Imidazolium Based Ionic Liquids + Acetonitrile Solutions at T = (298.15 to 313.15) K

rhoI	771.43	kg/m3	303.15	Densities and Viscosities of Ternary N,N'-Bis(2-pyridylmethylidene)-1,2-diiminoethane Schiff Base + Imidazolium Based Ionic Liquids + Acetonitrile Solutions at T = (298.15 to 313.15) K
rhoI	776.65	kg/m3	298.15	Densities and Viscosities of Ternary N,N'-Bis(2-pyridylmethylidene)-1,2-diiminoethane Schiff Base + Imidazolium Based Ionic Liquids + Acetonitrile Solutions at T = (298.15 to 313.15) K
rhoI	754.85	kg/m3	318.15	Density, Viscosity, Speed of Sound, and Refractive Index of a Ternary Solution of Aspirin, 1-Butyl-3-methylimidazolium Bromide, and Acetonitrile at Different Temperatures T = (288.15 to 318.15) K
rhoI	765.81	kg/m3	308.15	Density, Viscosity, Speed of Sound, and Refractive Index of a Ternary Solution of Aspirin, 1-Butyl-3-methylimidazolium Bromide, and Acetonitrile at Different Temperatures T = (288.15 to 318.15) K

rhoI	776.62	kg/m3	298.15	Density, Viscosity, Speed of Sound, and Refractive Index of a Ternary Solution of Aspirin, 1-Butyl-3-methylimidazolium Bromide, and Acetonitrile at Different Temperatures T = (288.15 to 318.15) K
rhoI	787.39	kg/m3	288.15	Density, Viscosity, Speed of Sound, and Refractive Index of a Ternary Solution of Aspirin, 1-Butyl-3-methylimidazolium Bromide, and Acetonitrile at Different Temperatures T = (288.15 to 318.15) K
rhoI	726.43	kg/m3	343.15	Volumetric Properties of Binary Mixtures of Two 1-Alkyl-3-Methylimidazolium Tetrafluoroborate Ionic Liquids with Molecular Solvents
rhoI	737.86	kg/m3	333.15	Volumetric Properties of Binary Mixtures of Two 1-Alkyl-3-Methylimidazolium Tetrafluoroborate Ionic Liquids with Molecular Solvents
rhoI	749.11	kg/m3	323.15	Volumetric Properties of Binary Mixtures of Two 1-Alkyl-3-Methylimidazolium Tetrafluoroborate Ionic Liquids with Molecular Solvents
rhoI	760.18	kg/m3	313.15	Volumetric Properties of Binary Mixtures of Two 1-Alkyl-3-Methylimidazolium Tetrafluoroborate Ionic Liquids with Molecular Solvents

rhoI	755.36	kg/m3	313.15	Densities and Viscosities of Ternary N,N'-Bis(2-pyridylmethylidene)-1,2-diiminoethane Schiff Base + Imidazolium Based Ionic Liquids + Acetonitrile Solutions at T = (298.15 to 313.15) K
rhoI	768.14	kg/m3	303.15	Densities and Viscosities of Ternary N,N'-Bis(2-pyridylmethylidene)-1,2-diiminoethane Schiff Base + Imidazolium Based Ionic Liquids + Acetonitrile Solutions at T = (298.15 to 313.15) K
rhoI	760.00	kg/m3	308.15	Densities and Viscosities of Ternary N,N'-Bis(2-pyridylmethylidene)-1,2-diiminoethane Schiff Base + Imidazolium Based Ionic Liquids + Acetonitrile Solutions at T = (298.15 to 313.15) K
rhoI	752.55	kg/m3	313.15	Densities and Viscosities of Ternary N,N'-Bis(2-pyridylmethylidene)-1,2-diiminoethane Schiff Base + Imidazolium Based Ionic Liquids + Acetonitrile Solutions at T = (298.15 to 313.15) K
rhoI	761.40	kg/m3	308.15	Densities and Viscosities of Ternary N,N'-Bis(2-pyridylmethylidene)-1,2-diiminoethane Schiff Base + Imidazolium Based Ionic Liquids + Acetonitrile Solutions at T = (298.15 to 313.15) K

rhoI	752.52	kg/m3	313.15	Densities and Viscosities of Ternary N,N'-Bis(2-pyridylmethylidene)-1,2-diiminoethane Schiff Base + Imidazolium Based Ionic Liquids + Acetonitrile Solutions at T = (298.15 to 313.15) K	
rhoI	763.20	kg/m3	308.15	Densities and Viscosities of Ternary N,N'-Bis(2-pyridylmethylidene)-1,2-diiminoethane Schiff Base + Imidazolium Based Ionic Liquids + Acetonitrile Solutions at T = (298.15 to 313.15) K	
rhoI	755.34	kg/m3	313.15	Densities and Viscosities of Ternary N,N'-Bis(2-pyridylmethylidene)-1,2-diiminoethane Schiff Base + Imidazolium Based Ionic Liquids + Acetonitrile Solutions at T = (298.15 to 313.15) K	
rhoI	782.30	kg/m3	293.15	Densities of Ionic Liquids, 1-Butyl-3-methylimidazolium Hexafluorophosphate and 1-Butyl-3-methylimidazolium Tetrafluoroborate, with Benzene, Acetonitrile, and 1-Propanol at T = (293.15 to 343.15) K	
rhoI	776.93	kg/m3	298.15	Densities of Ionic Liquids, 1-Butyl-3-methylimidazolium Hexafluorophosphate and 1-Butyl-3-methylimidazolium Tetrafluoroborate, with Benzene, Acetonitrile, and 1-Propanol at T = (293.15 to 343.15) K	

rhoI	771.44	kg/m3	303.15	Densities of Ionic Liquids, 1-Butyl-3-methylimidazolium Hexafluorophosphate and 1-Butyl-3-methylimidazolium Tetrafluoroborate, with Benzene, Acetonitrile, and 1-Propanol at T = (293.15 to 343.15) K
rhoI	765.97	kg/m3	308.15	Densities of Ionic Liquids, 1-Butyl-3-methylimidazolium Hexafluorophosphate and 1-Butyl-3-methylimidazolium Tetrafluoroborate, with Benzene, Acetonitrile, and 1-Propanol at T = (293.15 to 343.15) K
rhoI	760.68	kg/m3	313.15	Densities of Ionic Liquids, 1-Butyl-3-methylimidazolium Hexafluorophosphate and 1-Butyl-3-methylimidazolium Tetrafluoroborate, with Benzene, Acetonitrile, and 1-Propanol at T = (293.15 to 343.15) K
rhoI	749.73	kg/m3	323.15	Densities of Ionic Liquids, 1-Butyl-3-methylimidazolium Hexafluorophosphate and 1-Butyl-3-methylimidazolium Tetrafluoroborate, with Benzene, Acetonitrile, and 1-Propanol at T = (293.15 to 343.15) K
rhoI	738.19	kg/m3	333.15	Densities of Ionic Liquids, 1-Butyl-3-methylimidazolium Hexafluorophosphate and 1-Butyl-3-methylimidazolium Tetrafluoroborate, with Benzene, Acetonitrile, and 1-Propanol at T = (293.15 to 343.15) K



rhoI	726.94	kg/m3	343.15	Densities of Ionic Liquids, 1-Butyl-3-methylimidazolium Hexafluorophosphate and 1-Butyl-3-methylimidazolium Tetrafluoroborate, with Benzene, Acetonitrile, and 1-Propanol at T = (293.15 to 343.15) K
rhoI	782.04	kg/m3	293.15	Volumetric Properties of Binary Mixtures of 1-Butyl-3-methylimidazolium Chloride + Water or Hydrophilic Solvents at Different Temperatures
rhoI	776.64	kg/m3	298.15	Volumetric Properties of Binary Mixtures of 1-Butyl-3-methylimidazolium Chloride + Water or Hydrophilic Solvents at Different Temperatures
rhoI	771.21	kg/m3	303.15	Volumetric Properties of Binary Mixtures of 1-Butyl-3-methylimidazolium Chloride + Water or Hydrophilic Solvents at Different Temperatures
rhoI	765.75	kg/m3	308.15	Volumetric Properties of Binary Mixtures of 1-Butyl-3-methylimidazolium Chloride + Water or Hydrophilic Solvents at Different Temperatures
rhoI	760.25	kg/m3	313.15	Volumetric Properties of Binary Mixtures of 1-Butyl-3-methylimidazolium Chloride + Water or Hydrophilic Solvents at Different Temperatures

rhoI	754.73	kg/m3	318.15	Volumetric Properties of Binary Mixtures of 1-Butyl-3-methylimidazolium Chloride + Water or Hydrophilic Solvents at Different Temperatures
rhoI	771.13	kg/m3	303.15	Volumetric Properties of Binary Mixtures of Two 1-Alkyl-3-Methylimidazolium Tetrafluoroborate Ionic Liquids with Molecular Solvents
rhoI	776.55	kg/m3	298.15	Volumetric Properties of Binary Mixtures of Two 1-Alkyl-3-Methylimidazolium Tetrafluoroborate Ionic Liquids with Molecular Solvents
rhoI	781.95	kg/m3	293.15	Volumetric Properties of Binary Mixtures of Two 1-Alkyl-3-Methylimidazolium Tetrafluoroborate Ionic Liquids with Molecular Solvents
rhoI	776.62	kg/m3	298.15	Calorimetric Study of Nitrile Group-Solvent Interactions and Comparison with Dispersive Quasi-Chemical (DISQUAC) Predictions
rhoI	760.20	kg/m3	313.15	Volumetric Properties of Binary Mixtures Containing Ionic Liquids and Some Aprotic Solvents
rhoI	765.68	kg/m3	308.15	Volumetric Properties of Binary Mixtures Containing Ionic Liquids and Some Aprotic Solvents

rhoI	771.12	kg/m3	303.13	Volumetric Properties of Binary Mixtures Containing Ionic Liquids and Some Aprotic Solvents
rhoI	776.52	kg/m3	298.15	Volumetric Properties of Binary Mixtures Containing Ionic Liquids and Some Aprotic Solvents
rhoI	781.89	kg/m3	293.15	Volumetric Properties of Binary Mixtures Containing Ionic Liquids and Some Aprotic Solvents
rhoI	749.07	kg/m3	323.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	749.07	kg/m3	323.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	749.11	kg/m3	323.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile

rhoI	749.07	kg/m3	323.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	760.16	kg/m3	313.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	760.15	kg/m3	313.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	782.00	kg/m3	293.00	KDB
rhoI	776.53	kg/m3	298.15	Volumetric Properties of the Ionic Liquid, 1-Butyl-3-methylimidazolium Tetrafluoroborate, in Organic Solvents at T = 298.15 K
rhoI	765.98	kg/m3	308.10	Synthesis of 1,3-Dimethylimidazolium Chloride and Volumetric Property Investigations of Its Aqueous Solution
rhoI	786.50	kg/m3	293.15	Interpretation of Association Behavior and Molecular Interactions in Binary Mixtures from Thermoacoustics and Molecular Compression Data

rho1	781.10	kg/m3	298.15	Interpretation of Association Behavior and Molecular Interactions in Binary Mixtures from Thermoacoustics and Molecular Compression Data
rho1	773.30	kg/m3	303.15	Interpretation of Association Behavior and Molecular Interactions in Binary Mixtures from Thermoacoustics and Molecular Compression Data
rho1	766.50	kg/m3	308.15	Interpretation of Association Behavior and Molecular Interactions in Binary Mixtures from Thermoacoustics and Molecular Compression Data
rho1	760.50	kg/m3	313.15	Interpretation of Association Behavior and Molecular Interactions in Binary Mixtures from Thermoacoustics and Molecular Compression Data
rho1	776.50	kg/m3	298.15	Bubble point measurements of binary mixtures formed by 1-hexanol with selected nitro-compounds and substituted benzenes at 95.6 kPa
rho1	783.38	kg/m3	293.15	Isobaric vapor liquid equilibrium for binary system of methanol and acetonitrile

rhoI	776.55	kg/m3	298.15	Liquid-liquid equilibrium data for ternary mixtures composed of n-hexane, benzene and acetonitrile at (298.15, 308.15, and 318.15) K
rhoI	776.64	kg/m3	298.15	Liquid liquid equilibria in the ternary systems {hexadecane + BTX aromatics + 2-methoxyethanol or acetonitrile} at 298.15 K
rhoI	782.10	kg/m3	293.15	Effect of imidazolium-based ionic liquid on vapor-liquid equilibria of 2-propanol + acetonitrile binary system at 101.3 kPa
rhoI	776.69	kg/m3	298.15	Standard partial molar volumes of some electrolytes in ethylene carbonate based mixtures
rhoI	776.53	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	782.10	kg/m3	293.15	Densities and volumetric properties of (acetonitrile + an amide) binary mixtures at temperatures between 293.15 K and 318.15 K
rhoI	776.80	kg/m3	298.15	Densities and volumetric properties of (acetonitrile + an amide) binary mixtures at temperatures between 293.15 K and 318.15 K

rhoI	771.50	kg/m3	303.15	Densities and volumetric properties of (acetonitrile + an amide) binary mixtures at temperatures between 293.15 K and 318.15 K
rhoI	766.30	kg/m3	308.15	Densities and volumetric properties of (acetonitrile + an amide) binary mixtures at temperatures between 293.15 K and 318.15 K
rhoI	761.00	kg/m3	313.15	Densities and volumetric properties of (acetonitrile + an amide) binary mixtures at temperatures between 293.15 K and 318.15 K
rhoI	755.70	kg/m3	318.15	Densities and volumetric properties of (acetonitrile + an amide) binary mixtures at temperatures between 293.15 K and 318.15 K
rhoI	776.60	kg/m3	298.15	Density and speed of sound of lithium bromide with organic solvents: Measurement and correlation
rhoI	787.38	kg/m3	288.15	Volumetric properties of binary mixtures of ethers and acetonitrile: Experimental results and application of the Prigogine Flory Patterson theory
rhoI	782.01	kg/m3	293.15	Volumetric properties of binary mixtures of ethers and acetonitrile: Experimental results and application of the Prigogine Flory Patterson theory

rhoI	776.62	kg/m3	298.15	Volumetric properties of binary mixtures of ethers and acetonitrile: Experimental results and application of the Prigogine Flory Patterson theory
rhoI	771.19	kg/m3	303.15	Volumetric properties of binary mixtures of ethers and acetonitrile: Experimental results and application of the Prigogine Flory Patterson theory
rhoI	776.74	kg/m3	298.15	Probing subsistence of ion-pair and triple-ion of an ionic salt in liquid environments by means of conductometric contrivance
rhoI	776.47	kg/m3	298.15	Solution thermodynamics of iron(III)-N,N'-ethylene-bis(salicylideneiminato)-chloride in binary mixtures of N,N-dimethylformamide and acetonitrile at T = (298.15, 303.15, 308.15 and 313.15) K
rhoI	771.56	kg/m3	303.15	Solution thermodynamics of iron(III)-N,N'-ethylene-bis(salicylideneiminato)-chloride in binary mixtures of N,N-dimethylformamide and acetonitrile at T = (298.15, 303.15, 308.15 and 313.15) K
rhoI	765.26	kg/m3	308.15	Solution thermodynamics of iron(III)-N,N'-ethylene-bis(salicylideneiminato)-chloride in binary mixtures of N,N-dimethylformamide and acetonitrile at T = (298.15, 303.15, 308.15 and 313.15) K



rhoI	759.55	kg/m3	313.15	Solution thermodynamics of iron(III)-N,N'-ethylene-bis(salicylideneiminato)-chloride in binary mixtures of N,N-dimethylformamide and acetonitrile at T = (298.15, 303.15, 308.15 and 313.15) K
rhoI	776.62	kg/m3	298.15	Ionic molar volumes in methanol mixtures with acetonitrile, N,N-dimethylformamide and propylene carbonate at T = 298.15 K
rhoI	776.30	kg/m3	298.15	Solubility and solution thermodynamics of sorbic acid in eight pure organic solvents
rhoI	782.01	kg/m3	293.15	Volumetric properties of binary mixtures of (acetonitrile + amines) at several temperatures with application of the ERAS model
rhoI	776.61	kg/m3	298.15	Volumetric properties of binary mixtures of (acetonitrile + amines) at several temperatures with application of the ERAS model
rhoI	771.20	kg/m3	303.15	Volumetric properties of binary mixtures of (acetonitrile + amines) at several temperatures with application of the ERAS model

rhoI	765.75	kg/m3	308.15	Volumetric properties of binary mixtures of (acetonitrile + amines) at several temperatures with application of the ERAS model
rhoI	777.10	kg/m3	298.15	Solubility and solution thermodynamics of thymol in six pure organic solvents
rhoI	777.00	kg/m3	298.15	Measurement and correlation of solubility and solution thermodynamics of 1,3-dimethylurea in different solvents from T = (288.15 to 328.15) K
rhoI	782.10	kg/m3	293.15	Measurement and correlation of the vapor-liquid equilibrium for methanol + acetonitrile + imidazolium-based ionic liquids at 101.3 kPa
rhoI	782.12	kg/m3	293.15	Densities and volumetric properties of (acetonitrile + alkyl acrylate monomer) binary mixtures at temperatures from 293.15 K to 318.15 K
rhoI	776.82	kg/m3	298.15	Densities and volumetric properties of (acetonitrile + alkyl acrylate monomer) binary mixtures at temperatures from 293.15 K to 318.15 K

rhoI	771.52	kg/m3	303.15	Densities and volumetric properties of (acetonitrile + alkyl acrylate monomer) binary mixtures at temperatures from 293.15 K to 318.15 K
rhoI	766.21	kg/m3	308.15	Densities and volumetric properties of (acetonitrile + alkyl acrylate monomer) binary mixtures at temperatures from 293.15 K to 318.15 K
rhoI	760.90	kg/m3	313.15	Densities and volumetric properties of (acetonitrile + alkyl acrylate monomer) binary mixtures at temperatures from 293.15 K to 318.15 K
rhoI	755.59	kg/m3	318.15	Densities and volumetric properties of (acetonitrile + alkyl acrylate monomer) binary mixtures at temperatures from 293.15 K to 318.15 K
rhoI	782.00	kg/m3	293.15	Thermodynamic and transport properties of acetonitrile + alkanediols liquid mixtures at different temperatures, experimental measurements and modeling
rhoI	776.00	kg/m3	298.15	Thermodynamic and transport properties of acetonitrile + alkanediols liquid mixtures at different temperatures, experimental measurements and modeling

rhoI	771.00	kg/m3	303.15	Thermodynamic and transport properties of acetonitrile + alkanediols liquid mixtures at different temperatures, experimental measurements and modeling
rhoI	782.00	kg/m3	293.15	Thermophysical approach to understand the nature of molecular interactions and structural factor between methyl isobutyl ketone and organic solvents mixtures
rhoI	760.16	kg/m3	313.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	760.00	kg/m3	313.15	Thermophysical approach to understand the nature of molecular interactions and structural factor between methyl isobutyl ketone and organic solvents mixtures
rhoI	782.12	kg/m3	293.15	Densities, ultrasonic speeds, excess and partial molar properties of binary mixtures of acetonitrile with some alkyl methacrylates at temperatures from 293.15 K to 318.15 K

rhoI	776.82	kg/m3	298.15	Densities, ultrasonic speeds, excess and partial molar properties of binary mixtures of acetonitrile with some alkyl methacrylates at temperatures from 293.15 K to 318.15 K
rhoI	771.52	kg/m3	303.15	Densities, ultrasonic speeds, excess and partial molar properties of binary mixtures of acetonitrile with some alkyl methacrylates at temperatures from 293.15 K to 318.15 K
rhoI	766.21	kg/m3	308.15	Densities, ultrasonic speeds, excess and partial molar properties of binary mixtures of acetonitrile with some alkyl methacrylates at temperatures from 293.15 K to 318.15 K
rhoI	760.90	kg/m3	313.15	Densities, ultrasonic speeds, excess and partial molar properties of binary mixtures of acetonitrile with some alkyl methacrylates at temperatures from 293.15 K to 318.15 K
rhoI	755.59	kg/m3	318.15	Densities, ultrasonic speeds, excess and partial molar properties of binary mixtures of acetonitrile with some alkyl methacrylates at temperatures from 293.15 K to 318.15 K

rhoI	798.12	kg/m3	278.15	Thermodynamic evidence for nano-heterogeneity in solutions of the macrocycle C-butylresorcin[4]arene in non-aqueous solvents
rhoI	792.79	kg/m3	283.15	Thermodynamic evidence for nano-heterogeneity in solutions of the macrocycle C-butylresorcin[4]arene in non-aqueous solvents
rhoI	787.44	kg/m3	288.15	Thermodynamic evidence for nano-heterogeneity in solutions of the macrocycle C-butylresorcin[4]arene in non-aqueous solvents
rhoI	782.06	kg/m3	293.15	Thermodynamic evidence for nano-heterogeneity in solutions of the macrocycle C-butylresorcin[4]arene in non-aqueous solvents
rhoI	776.66	kg/m3	298.15	Thermodynamic evidence for nano-heterogeneity in solutions of the macrocycle C-butylresorcin[4]arene in non-aqueous solvents
rhoI	771.23	kg/m3	303.15	Thermodynamic evidence for nano-heterogeneity in solutions of the macrocycle C-butylresorcin[4]arene in non-aqueous solvents
rhoI	765.77	kg/m3	308.15	Thermodynamic evidence for nano-heterogeneity in solutions of the macrocycle C-butylresorcin[4]arene in non-aqueous solvents

rhoI	776.68	kg/m3	298.15	Exploration of Solvation Consequence of Ionic Liquid [Bu4PCH3SO3] in Various Solvent Systems by Conductance and FTIR Study
rhoI	776.51	kg/m3	298.15	Excess Enthalpies in Binary Systems of Isomeric C8 Aliphatic Monoethers with Acetonitrile and Their Description by the COSMO-SAC Model
rhoI	771.06	kg/m3	303.15	Excess Enthalpies in Binary Systems of Isomeric C8 Aliphatic Monoethers with Acetonitrile and Their Description by the COSMO-SAC Model
rhoI	765.59	kg/m3	308.15	Excess Enthalpies in Binary Systems of Isomeric C8 Aliphatic Monoethers with Acetonitrile and Their Description by the COSMO-SAC Model
rhoI	760.10	kg/m3	313.15	Excess Enthalpies in Binary Systems of Isomeric C8 Aliphatic Monoethers with Acetonitrile and Their Description by the COSMO-SAC Model
rhoI	754.56	kg/m3	318.15	Excess Enthalpies in Binary Systems of Isomeric C8 Aliphatic Monoethers with Acetonitrile and Their Description by the COSMO-SAC Model

rhoI	748.99	kg/m3	323.15	Excess Enthalpies in Binary Systems of Isomeric C8 Aliphatic Monoethers with Acetonitrile and Their Description by the COSMO-SAC Model
rhoI	776.71	kg/m3	298.15	Excess Molar Volumes, Excess Molar Isentropic Compressibilities, Viscosity Deviations, and Activation Parameters for 1-Ethyl-3-methyl-imidazolium Trifluoro-methanesulfonate + Dimethyl Sulfoxide and/or Acetonitrile at T = 298.15 to 323.15 K and P = 0.1 MPa
rhoI	771.29	kg/m3	303.15	Excess Molar Volumes, Excess Molar Isentropic Compressibilities, Viscosity Deviations, and Activation Parameters for 1-Ethyl-3-methyl-imidazolium Trifluoro-methanesulfonate + Dimethyl Sulfoxide and/or Acetonitrile at T = 298.15 to 323.15 K and P = 0.1 MPa
rhoI	765.83	kg/m3	308.15	Excess Molar Volumes, Excess Molar Isentropic Compressibilities, Viscosity Deviations, and Activation Parameters for 1-Ethyl-3-methyl-imidazolium Trifluoro-methanesulfonate + Dimethyl Sulfoxide and/or Acetonitrile at T = 298.15 to 323.15 K and P = 0.1 MPa



rhoI	760.34	kg/m3	313.15	Excess Molar Volumes, Excess Molar Isentropic Compressibilities, Viscosity Deviations, and Activation Parameters for 1-Ethyl-3-methyl-imidazolium Trifluoro-methanesulfonate + Dimethyl Sulfoxide and/or Acetonitrile at T = 298.15 to 323.15 K and P = 0.1 MPa
rhoI	754.82	kg/m3	318.15	Excess Molar Volumes, Excess Molar Isentropic Compressibilities, Viscosity Deviations, and Activation Parameters for 1-Ethyl-3-methyl-imidazolium Trifluoro-methanesulfonate + Dimethyl Sulfoxide and/or Acetonitrile at T = 298.15 to 323.15 K and P = 0.1 MPa
rhoI	749.25	kg/m3	323.15	Excess Molar Volumes, Excess Molar Isentropic Compressibilities, Viscosity Deviations, and Activation Parameters for 1-Ethyl-3-methyl-imidazolium Trifluoro-methanesulfonate + Dimethyl Sulfoxide and/or Acetonitrile at T = 298.15 to 323.15 K and P = 0.1 MPa
rhoI	787.33	kg/m3	288.15	Density, Electrical Conductivity, Dynamic Viscosity, Excess Properties, and Molecular Interactions of Ionic Liquid 1-Cyanopropyl-3-methylimidazolium Tetrafluoroborate and Binary System with Acetonitrile

rhoI	781.96	kg/m3	293.15	Density, Electrical Conductivity, Dynamic Viscosity, Excess Properties, and Molecular Interactions of Ionic Liquid 1-Cyanopropyl-3-methylimidazolium Tetrafluoroborate and Binary System with Acetonitrile
rhoI	776.61	kg/m3	298.15	Density, Electrical Conductivity, Dynamic Viscosity, Excess Properties, and Molecular Interactions of Ionic Liquid 1-Cyanopropyl-3-methylimidazolium Tetrafluoroborate and Binary System with Acetonitrile
rhoI	771.18	kg/m3	303.15	Density, Electrical Conductivity, Dynamic Viscosity, Excess Properties, and Molecular Interactions of Ionic Liquid 1-Cyanopropyl-3-methylimidazolium Tetrafluoroborate and Binary System with Acetonitrile
rhoI	765.73	kg/m3	308.15	Density, Electrical Conductivity, Dynamic Viscosity, Excess Properties, and Molecular Interactions of Ionic Liquid 1-Cyanopropyl-3-methylimidazolium Tetrafluoroborate and Binary System with Acetonitrile

rhoI	760.25	kg/m3	313.15	Density, Electrical Conductivity, Dynamic Viscosity, Excess Properties, and Molecular Interactions of Ionic Liquid 1-Cyanopropyl-3-methylimidazolium Tetrafluoroborate and Binary System with Acetonitrile
rhoI	754.72	kg/m3	318.15	Density, Electrical Conductivity, Dynamic Viscosity, Excess Properties, and Molecular Interactions of Ionic Liquid 1-Cyanopropyl-3-methylimidazolium Tetrafluoroborate and Binary System with Acetonitrile
rhoI	749.16	kg/m3	323.15	Density, Electrical Conductivity, Dynamic Viscosity, Excess Properties, and Molecular Interactions of Ionic Liquid 1-Cyanopropyl-3-methylimidazolium Tetrafluoroborate and Binary System with Acetonitrile
rhoI	787.58	kg/m3	288.15	The Thermodynamic and Excess Properties of Trialkyl-Substituted Imidazolium-Based Ionic Liquids with Thiocyanate and Its Binary Systems with Acetonitrile
rhoI	781.89	kg/m3	293.15	The Thermodynamic and Excess Properties of Trialkyl-Substituted Imidazolium-Based Ionic Liquids with Thiocyanate and Its Binary Systems with Acetonitrile

rhoI	776.52	kg/m3	298.15	The Thermodynamic and Excess Properties of Trialkyl-Substituted Imidazolium-Based Ionic Liquids with Thiocyanate and Its Binary Systems with Acetonitrile
rhoI	771.12	kg/m3	303.15	The Thermodynamic and Excess Properties of Trialkyl-Substituted Imidazolium-Based Ionic Liquids with Thiocyanate and Its Binary Systems with Acetonitrile
rhoI	765.83	kg/m3	308.15	The Thermodynamic and Excess Properties of Trialkyl-Substituted Imidazolium-Based Ionic Liquids with Thiocyanate and Its Binary Systems with Acetonitrile
rhoI	760.36	kg/m3	313.15	The Thermodynamic and Excess Properties of Trialkyl-Substituted Imidazolium-Based Ionic Liquids with Thiocyanate and Its Binary Systems with Acetonitrile
rhoI	755.14	kg/m3	318.15	The Thermodynamic and Excess Properties of Trialkyl-Substituted Imidazolium-Based Ionic Liquids with Thiocyanate and Its Binary Systems with Acetonitrile

rhoI	749.44	kg/m3	323.15	The Thermodynamic and Excess Properties of Trialkyl-Substituted Imidazolium-Based Ionic Liquids with Thiocyanate and Its Binary Systems with Acetonitrile
rhoI	785.71	kg/m3	298.15	Liquid-Liquid Equilibrium for Ternary Systems, Water + 5-Hydroxymethylfurfural + (1-Butanol, Isobutanol, Methyl Isobutyl Ketone), at 313.15, 323.15, and 333.15 K
rhoI	776.71	kg/m3	298.15	Effect of Solvents and Temperature on Interactions in Binary and Ternary Mixtures of 1-Butyl-3-methylimidazolium Trifluoromethanesulfonate with Acetonitrile or/and N,N-Dimethylformamide
rhoI	771.29	kg/m3	303.15	Effect of Solvents and Temperature on Interactions in Binary and Ternary Mixtures of 1-Butyl-3-methylimidazolium Trifluoromethanesulfonate with Acetonitrile or/and N,N-Dimethylformamide
rhoI	765.83	kg/m3	308.15	Effect of Solvents and Temperature on Interactions in Binary and Ternary Mixtures of 1-Butyl-3-methylimidazolium Trifluoromethanesulfonate with Acetonitrile or/and N,N-Dimethylformamide

rhoI	760.34	kg/m3	313.15	Effect of Solvents and Temperature on Interactions in Binary and Ternary Mixtures of 1-Butyl-3-methylimidazolium Trifluoromethanesulfonate with Acetonitrile or/and N,N-Dimethylformamide
rhoI	754.82	kg/m3	318.15	Effect of Solvents and Temperature on Interactions in Binary and Ternary Mixtures of 1-Butyl-3-methylimidazolium Trifluoromethanesulfonate with Acetonitrile or/and N,N-Dimethylformamide
rhoI	749.25	kg/m3	323.15	Effect of Solvents and Temperature on Interactions in Binary and Ternary Mixtures of 1-Butyl-3-methylimidazolium Trifluoromethanesulfonate with Acetonitrile or/and N,N-Dimethylformamide
rhoI	787.31	kg/m3	288.13	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	787.31	kg/m3	288.13	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile

rhoI	787.31	kg/m3	288.13	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	787.31	kg/m3	288.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	781.93	kg/m3	293.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	781.93	kg/m3	293.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	781.93	kg/m3	293.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile

rhoI	781.93	kg/m3	293.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	776.53	kg/m3	298.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	776.53	kg/m3	298.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	776.53	kg/m3	298.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	776.53	kg/m3	298.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	776.53	kg/m3	298.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile



rhoI	771.10	kg/m3	303.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	771.11	kg/m3	303.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	771.10	kg/m3	303.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	771.11	kg/m3	303.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	765.64	kg/m3	308.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile

rhoI	765.66	kg/m3	308.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	765.65	kg/m3	308.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
rhoI	771.00	kg/m3	303.15	Thermophysical approach to understand the nature of molecular interactions and structural factor between methyl isobutyl ketone and organic solvents mixtures
rhoI	760.18	kg/m3	313.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. II. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in Acetonitrile
speedsl	1198.32	m/s	318.15	Volumetric and Speed of Sound of Ionic Liquid, 1-Butyl-3-methylimidazolium Hexafluorophosphate with Acetonitrile and Methanol at T ) (298.15 to 318.15) K

speedsl	1258.73	m/s	303.15	Application of Prigogine Flory Patterson theory to excess molar volume and speed of sound of 1-n-butyl-3-methylimidazolium hexafluorophosphate or 1-n-butyl-3-methylimidazolium tetrafluoroborate in methanol and acetonitrile
speedsl	1218.52	m/s	313.15	Application of Prigogine Flory Patterson theory to excess molar volume and speed of sound of 1-n-butyl-3-methylimidazolium hexafluorophosphate or 1-n-butyl-3-methylimidazolium tetrafluoroborate in methanol and acetonitrile
speedsl	1319.14	m/s	288.15	Volumetric and Isentropic Compressibility Behavior of Ionic Liquid, 1-Propyl-3-Methylimidazolium Bromide in Acetonitrile, Dimethylformamide, and Dimethylsulfoxide at T = (288.15 to 308.15) K
speedsl	1299.01	m/s	293.15	Volumetric and Isentropic Compressibility Behavior of Ionic Liquid, 1-Propyl-3-Methylimidazolium Bromide in Acetonitrile, Dimethylformamide, and Dimethylsulfoxide at T = (288.15 to 308.15) K

speedsl	1278.77	m/s	298.15	Volumetric and Isentropic Compressibility Behavior of Ionic Liquid, 1-Propyl-3-Methylimidazolium Bromide in Acetonitrile, Dimethylformamide, and Dimethylsulfoxide at T = (288.15 to 308.15) K
speedsl	1258.51	m/s	303.15	Volumetric and Isentropic Compressibility Behavior of Ionic Liquid, 1-Propyl-3-Methylimidazolium Bromide in Acetonitrile, Dimethylformamide, and Dimethylsulfoxide at T = (288.15 to 308.15) K
speedsl	1238.40	m/s	308.15	Volumetric and Isentropic Compressibility Behavior of Ionic Liquid, 1-Propyl-3-Methylimidazolium Bromide in Acetonitrile, Dimethylformamide, and Dimethylsulfoxide at T = (288.15 to 308.15) K
speedsl	1239.00	m/s	308.15	Ultrasonic studies on binary mixtures of some aromatic ketones with acetonitrile at T = 308.15 K
speedsl	1278.62	m/s	298.15	Application of Prigogine Flory Patterson theory to excess molar volume and speed of sound of 1-n-butyl-3-methylimidazolium hexafluorophosphate or 1-n-butyl-3-methylimidazolium tetrafluoroborate in methanol and acetonitrile

speedsl	1278.28	m/s	298.15	Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. IV. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Bromides and Chlorides in Acetonitrile
speedsl	1238.66	m/s	308.15	Application of Prigogine Flory Patterson theory to excess molar volume and speed of sound of 1-n-butyl-3-methylimidazolium hexafluorophosphate or 1-n-butyl-3-methylimidazolium tetrafluoroborate in methanol and acetonitrile
speedsl	1281.30	m/s	298.15	Compressibility Studies of Binary Solutions Involving Water as a Solute in Nonaqueous Solvents at T ) 298.15 K
speedsl	1198.32	m/s	318.15	Application of Prigogine Flory Patterson theory to excess molar volume and speed of sound of 1-n-butyl-3-methylimidazolium hexafluorophosphate or 1-n-butyl-3-methylimidazolium tetrafluoroborate in methanol and acetonitrile
speedsl	1278.62	m/s	298.15	Volumetric and Speed of Sound of Ionic Liquid, 1-Butyl-3-methylimidazolium Hexafluorophosphate with Acetonitrile and Methanol at T ) (298.15 to 318.15) K

speedsl	1258.73	m/s	303.15	Volumetric and Speed of Sound of Ionic Liquid, 1-Butyl-3-methylimidazolium Hexafluorophosphate with Acetonitrile and Methanol at T ) (298.15 to 318.15) K
speedsl	1238.66	m/s	308.15	Volumetric and Speed of Sound of Ionic Liquid, 1-Butyl-3-methylimidazolium Hexafluorophosphate with Acetonitrile and Methanol at T ) (298.15 to 318.15) K
speedsl	1218.52	m/s	313.15	Volumetric and Speed of Sound of Ionic Liquid, 1-Butyl-3-methylimidazolium Hexafluorophosphate with Acetonitrile and Methanol at T ) (298.15 to 318.15) K
srf	0.03	N/m	283.15	Application of the Extended Langmuir Model for the Determination of Lyophobicity of 1-Propanol in 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	313.15	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	293.20	KDB
srf	0.03	N/m	278.15	Application of the Extended Langmuir Model for the Determination of Lyophobicity of 1-Propanol in Acetonitrile
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
srf	0.03	N/m	288.15	Application of the Extended Langmuir Model for the Determination of Lyophobicity of 1-Propanol in Acetonitrile
srf	0.03	N/m	293.15	Application of the Extended Langmuir Model for the Determination of Lyophobicity of 1-Propanol in Acetonitrile
srf	0.03	N/m	298.15	Application of the Extended Langmuir Model for the Determination of Lyophobicity of 1-Propanol in Acetonitrile
srf	0.03	N/m	303.15	Application of the Extended Langmuir Model for the Determination of Lyophobicity of 1-Propanol in Acetonitrile

srf	0.03	N/m	308.15	Application of the Extended Langmuir Model for the Determination of Lyophobicity of 1-Propanol in Acetonitrile
srf	0.03	N/m	313.15	Application of the Extended Langmuir Model for the Determination of Lyophobicity of 1-Propanol in Acetonitrile
srf	0.03	N/m	283.15	Surface tension of non-ideal binary and ternary liquid mixtures at various temperatures and p = 81.5 kPa
srf	0.03	N/m	298.15	Surface tension of non-ideal binary and ternary liquid mixtures at various temperatures and p = 81.5 kPa
srf	0.03	N/m	308.15	Surface tension of non-ideal binary and ternary liquid mixtures at various temperatures and p = 81.5 kPa
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	298.15	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	308.15	Thermodynamic surface properties of [BMIm][NTf2] or [EMIm][NTf2] binary mixtures with tetrahydrofuran, acetonitrile or dimethylsulfoxide
srf	0.03	N/m	313.15	Thermodynamic surface properties of [BMIm][NTf2] or [EMIm][NTf2] binary mixtures with tetrahydrofuran, acetonitrile or dimethylsulfoxide
srf	0.03	N/m	298.15	Physicochemical properties of two 1-alkyl-1-methylpyrrolidinium bis[(trifluoromethyl)sulfonyl]imide ionic liquids and of binary mixtures of 1-butyl-1-methylpyrrolidinium bis[(trifluoromethyl)sulfonyl]imide with methanol or acetonitrile
srf	0.03	N/m	288.15	Physicochemical properties of two 1-alkyl-1-methylpyrrolidinium bis[(trifluoromethyl)sulfonyl]imide ionic liquids and of binary mixtures of 1-butyl-1-methylpyrrolidinium bis[(trifluoromethyl)sulfonyl]imide with methanol or acetonitrile
srf	0.03	N/m	308.15	Physicochemical properties of two 1-alkyl-1-methylpyrrolidinium bis[(trifluoromethyl)sulfonyl]imide ionic liquids and of binary mixtures of 1-butyl-1-methylpyrrolidinium bis[(trifluoromethyl)sulfonyl]imide with methanol or acetonitrile

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	Experimental Data and Correlation of Surface Tensions of the Binary and Ternary Systems of Water + Acetonitrile + 2-Propanol at 298.15 K and Atmospheric Pressure
srf	0.03	N/m	293.15	Density and Surface Tension of Binary Mixtures of Acetonitrile + 1-Alkanol at 293.15 K
svapt	111.44	J/mol×K	298.15	NIST Webbook
tcondl	0.18	W/m×K	313.15	Liquid Thermal Conductivities of Acetonitrile, Diethyl Sulfide, Hexamethyleneimine, Tetrahydrothiophene, and Tetramethylethylenediamine
tcondl	0.21	W/m×K	273.15	Liquid Thermal Conductivities of Acetonitrile, Diethyl Sulfide, Hexamethyleneimine, Tetrahydrothiophene, and Tetramethylethylenediamine
tcondl	0.20	W/m×K	293.15	Liquid Thermal Conductivities of Acetonitrile, Diethyl Sulfide, Hexamethyleneimine, Tetrahydrothiophene, and Tetramethylethylenediamine

## Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.53562e+01
Coeff. B	-3.64591e+03
Coeff. C	-1.52630e+01
Temperature range (K), min.	229.32
Temperature range (K), max.	545.50

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/T + C \cdot \ln(T) + D \cdot T^2$
Coeff. A	4.10590e+01
Coeff. B	-4.99962e+03
Coeff. C	-3.88171e+00
Coeff. D	3.51596e-06
Temperature range (K), min.	229.32
Temperature range (K), max.	545.50

## Datasets

### Refractive index (Na D-line)

Pressure, kPa - Liquid	Temperature, K - Liquid	Refractive index (Na D-line) - Liquid
81.50	298.15	1.3417

Reference

<https://www.doi.org/10.1021/je700645p>

### Mass density, kg/m3

Pressure, kPa - Liquid	Temperature, K - Liquid	Mass density, kg/m3 - Liquid
200.00	313.15	755.0
200.00	353.15	710.0
600.00	313.15	755.0



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<https://www.doi.org/10.1016/j.jct.2006.08.007>  
<https://www.doi.org/10.1021/je060178m>  
<https://www.doi.org/10.1021/je050317k>  
<https://www.doi.org/10.1016/j.jct.2012.09.017>  
<https://www.doi.org/10.1016/j.jct.2016.10.020>  
<https://www.doi.org/10.1016/j.fluid.2005.02.016>  
<https://www.doi.org/10.1021/je300339q>  
<https://www.doi.org/10.1016/j.tca.2012.08.009>  
<https://www.doi.org/10.1016/j.fluid.2012.12.003>  
<https://www.doi.org/10.1021/acs.jced.5b01053>  
<https://www.doi.org/10.1016/j.jct.2006.01.016>  
<https://www.doi.org/10.1021/je060305e>  
<https://www.doi.org/10.1016/j.jct.2013.08.011>  
<https://www.doi.org/10.1016/j.jct.2016.02.006>  
<https://www.doi.org/10.1021/je1007235>  
<https://www.doi.org/10.1016/j.fluid.2014.01.029>  
<https://www.doi.org/10.1016/j.fluid.2010.10.008>  
<https://www.doi.org/10.1021/acs.jced.9b00350>  
<https://www.doi.org/10.1021/acs.jced.7b01085>  
<https://www.doi.org/10.1016/j.fluid.2013.09.023>  
<https://www.doi.org/10.1016/j.jct.2016.05.011>  
<https://www.doi.org/10.1021/acs.jced.9b00460>  
<https://www.doi.org/10.1016/j.fluid.2016.12.012>  
<https://www.doi.org/10.1021/acs.jced.8b00684>  
<https://www.doi.org/10.1016/j.fluid.2015.10.035>  
<https://www.doi.org/10.1016/j.jct.2018.09.017>  
<https://www.doi.org/10.1021/acs.jced.5b00075>  
<https://www.doi.org/10.1021/je900733j>  
<https://www.doi.org/10.1021/je100671v>  
<https://www.doi.org/10.1016/j.jct.2015.09.002>  
<https://www.doi.org/10.1021/acs.jced.8b00292>  
<https://www.doi.org/10.1021/je020226c>  
<https://www.doi.org/10.1021/je4001894>  
<https://www.doi.org/10.1021/acs.jced.6b00911>  
<https://www.doi.org/10.1016/j.fluid.2012.05.006>  
<https://www.doi.org/10.1016/j.fluid.2018.04.004>  
<https://www.doi.org/10.1021/acs.jced.6b00349>  
<https://www.doi.org/10.1016/j.fluid.2018.05.005>  
<https://www.doi.org/10.1016/j.jct.2016.09.011>  
<https://www.doi.org/10.1021/je900401z>  
<https://www.doi.org/10.1016/j.jct.2016.08.034>  
<https://www.doi.org/10.1021/je500867u>  
<https://www.doi.org/10.1021/acs.jced.8b00888>

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<https://www.doi.org/10.1016/j.jct.2014.12.025>  
<https://www.doi.org/10.1021/je4003114>  
**a-5,5-tetramethyl**  
<https://www.doi.org/10.1021/acs.jced.9b00220>  
<https://www.doi.org/10.1021/je100125x>  
<https://www.doi.org/10.1021/acs.jced.8b01051>  
<https://www.doi.org/10.1021/acs.jced.8b00931>  
<https://www.doi.org/10.1021/acs.jced.6b00163>  
<https://www.doi.org/10.1021/acs.jced.8b00536>  
<https://www.doi.org/10.1016/j.jct.2013.10.026>  
<https://www.doi.org/10.1021/je400790d>  
<https://www.doi.org/10.1021/acs.jced.8b00080>  
<https://www.doi.org/10.1021/je8007099>  
<https://www.doi.org/10.1016/j.jct.2015.05.022>  
<https://www.doi.org/10.1021/je1005483>  
**um**  
<https://www.doi.org/10.1016/j.jct.2013.05.008>  
<https://www.doi.org/10.1016/j.jct.2015.12.032>  
<https://www.doi.org/10.1021/je5002158>  
<https://www.doi.org/10.1021/je3009792>  
<https://www.doi.org/10.1021/je400531a>  
<https://www.doi.org/10.1021/acs.jced.8b00288>  
<https://www.doi.org/10.1021/acs.jced.6b00664>  
**axazaisowurtziane**  
<https://www.doi.org/10.1021/je8001909>  
<https://www.doi.org/10.1016/j.fluid.2016.08.019>  
<https://www.doi.org/10.1016/j.jct.2016.07.001>  
<https://www.doi.org/10.1021/acs.jced.6b00721>  
<https://www.doi.org/10.1021/acs.jced.9b00278>  
<https://www.doi.org/10.1016/j.jct.2016.08.021>  
<https://www.doi.org/10.1021/je900554r>  
<https://www.doi.org/10.1016/j.jct.2015.09.010>  
<https://www.doi.org/10.1016/j.jct.2016.07.050>  
<https://www.doi.org/10.1021/je0498661>  
<https://www.doi.org/10.1016/j.jct.2011.09.028>  
<https://www.doi.org/10.1021/acs.jced.9b00308>  
<https://www.doi.org/10.1021/je800754w>  
<https://www.doi.org/10.1016/j.jct.2015.07.010>  
<https://www.doi.org/10.1016/j.jct.2014.05.018>  
<https://www.doi.org/10.1021/acs.jced.7b01108>  
<https://www.doi.org/10.1021/acs.jced.8b01181>  
<https://www.doi.org/10.1021/acs.jced.9b00844>  
<https://www.doi.org/10.1021/je1002237>  
<https://www.doi.org/10.1016/j.jct.2010.04.019>  
<https://www.doi.org/10.1016/j.jct.2018.06.006>  
<https://www.doi.org/10.1016/j.jct.2011.11.005>



[illegible]

<https://www.doi.org/10.1016/j.jct.2016.11.029>  
<https://www.doi.org/10.1016/j.fluid.2018.06.017>  
<https://www.doi.org/10.1016/j.fluid.2016.03.008>  
<https://www.doi.org/10.1016/j.jct.2011.11.021>  
<https://www.doi.org/10.1016/j.jct.2016.06.015>  
<https://www.doi.org/10.1021/je2000292>  
<https://www.doi.org/10.1021/je3010535>  
<https://www.doi.org/10.1016/j.jct.2017.03.004>  
<https://www.doi.org/10.1021/acs.jced.5b00714>  
<https://www.doi.org/10.1021/acs.jced.7b00178>  
<https://www.doi.org/10.1016/j.jct.2012.01.019>  
<http://link.springer.com/article/10.1007/BF02311772>  
<https://www.doi.org/10.1021/acs.jced.5b00175>  
<https://www.doi.org/10.1016/j.jct.2015.02.024>  
<https://www.doi.org/10.1016/j.fluid.2016.06.022>  
<https://www.doi.org/10.1021/acs.jced.5b00995>  
<https://www.doi.org/10.1021/acs.jced.9b00190>  
<https://www.doi.org/10.1016/j.fluid.2013.09.028>  
<https://www.doi.org/10.1021/acs.jced.8b01265>  
<https://www.doi.org/10.1016/j.jct.2006.01.015>  
<https://www.doi.org/10.1016/j.jct.2016.11.019>  
<https://www.doi.org/10.1016/j.jct.2014.02.021>  
<https://www.doi.org/10.1021/acs.jced.8b01101>  
<https://www.doi.org/10.1016/j.jct.2017.05.013>  
<https://www.doi.org/10.1021/je900609f>  
<https://www.doi.org/10.1021/acs.jced.9b00620>  
<https://www.doi.org/10.1016/j.jct.2007.09.007>  
<https://www.doi.org/10.1016/j.jct.2017.06.008>  
<https://www.doi.org/10.1021/je020216b>  
<https://www.doi.org/10.1021/je0498560>  
<https://www.doi.org/10.1021/je3007474>  
<https://www.doi.org/10.1021/je400781b>  
<https://www.doi.org/10.1016/j.fluid.2014.01.043>  
<https://www.doi.org/10.1016/j.jct.2016.12.014>  
<https://www.doi.org/10.1016/j.jct.2016.07.021>  
<https://www.doi.org/10.1016/j.fluid.2015.03.036>  
<https://www.doi.org/10.1016/j.fluid.2013.06.037>  
<https://www.doi.org/10.1021/je200195q>  
<https://www.doi.org/10.1021/acs.jced.8b00785>  
<https://www.doi.org/10.1021/acs.jced.9b00661>  
<https://www.doi.org/10.1007/s10765-016-2096-3>  
<https://www.doi.org/10.1021/acs.jced.8b00717>  
<https://www.doi.org/10.1021/je5008372>

[illegible]

<https://www.doi.org/10.1016/j.jct.2017.12.012>

<https://www.doi.org/10.1021/acs.iced.7b00755>

<https://www.doi.org/10.1021/acs.jced.7b00615>

<https://www.doi.org/10.1021/je050165t>

<https://www.doi.org/10.1016/j.jct.2016.07.009>

<https://www.doi.org/10.1016/j.tca.2012.03.007>

<https://www.doi.org/10.1016/j.fluid.2019.01.005>

<https://www.doi.org/10.1016/j.ijct.2016.07.017>

<https://www.doi.org/10.1016/j.fluid.2010.10.013>

<https://www.doi.org/10.1021/je300692s>

<https://www.doi.org/10.1021/acs.iced.5b00670>

<https://www.doi.org/10.1016/j.ijct.2013.02.021>

<https://www.doi.org/10.1021/acs.jced.7b00743>

<https://www.doi.org/10.1021/ie301238f>

<https://www.doi.org/10.1021/acs.iced.6b01044>

<https://www.doi.org/10.1016/j.ijct.2016.12.028>

<https://www.doi.org/10.1021/ie034144f>

<https://www.doi.org/10.1016/j.ijct.2016.12.002>

<https://www.doi.org/10.1021/acs.iced.9b00562>

<https://www.doi.org/10.1021/acs.iced.7b01134>

<https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>

<https://www.doi.org/10.1021/acs.iced.8b01214>

<https://www.doi.org/10.1016/j.ijct.2004.03.007>

<https://www.doi.org/10.1021/jo020150k>

<https://www.doi.org/10.1016/j.fluid.2017.08.007>

<https://www.doi.org/10.1016/j.ijct.2017.01.006>

<https://www.doi.org/10.1016/j.fluid.2016.03.004>

<https://www.doi.org/10.1016/j.ijet.2016.11.033>

<https://www.doi.org/10.1016/j.fluid.2010.01.003>

<https://www.doi.org/10.1031/acs.joc.4b00635>

<https://www.doi.org/10.1016/j.ijet.2016.08.025>

<https://www.doi.org/10.1021/acs.joc.5b00400>

<https://www.doi.org/10.1021/acs.joc.5b00659>

<https://www.doi.org/10.1001/ajph.2019.0352>

<https://www.doi.org/10.1016/j.fluid.2019.09.027>

<https://www.doi.org/10.1001/psap.2021.5b00100>

**10**        "I'm not going to let you go," he said.

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<http://www.fishbase.org/10.1016/j.jct.2010.03.003>

<http://dx.doi.org/10.1016/j.joule.2019.05.005>

<http://pubs.acs.org, doi, suppl, 10.1021/acs.chem.5b00000>

<https://www.eia.gov/tools/batchquery/batchquery.cfm>

<https://www.ashm.org/forfe=7/assjccat/02000000>

14 Pure Solvents from 278.15 to 323.15 K:



[illegible]

https://www.doi.org/10.1016/j.jct.2012.03.015  
https://www.doi.org/10.1021/je050440b  
https://www.doi.org/10.1016/j.fluid.2013.06.038  
https://www.doi.org/10.1021/acs.jced.9b00353  
https://www.doi.org/10.1021/acs.jced.5b00796  
https://www.doi.org/10.1016/j.fluid.2014.04.006  
https://www.doi.org/10.1021/je900888e  
https://www.doi.org/10.1021/acs.jced.6b00552  
https://www.doi.org/10.1021/je500392g  
https://www.doi.org/10.1021/acs.jced.8b01256  
https://www.doi.org/10.1021/acs.jced.8b00176  
https://www.doi.org/10.1021/je020130i  
https://www.doi.org/10.1021/je4010917  
https://www.doi.org/10.1016/j.jct.2011.09.022  
https://www.doi.org/10.1016/j.jct.2005.08.009  
https://www.doi.org/10.1016/j.fluid.2013.05.038  
https://www.doi.org/10.1021/acs.jced.6b00230  
https://www.doi.org/10.1021/acs.jced.8b00196  
https://www.doi.org/10.1021/acs.jced.8b00271  
https://www.doi.org/10.1021/je500437m  
https://www.doi.org/10.1021/acs.jced.8b00551  
https://www.doi.org/10.1021/je200637v  
https://www.doi.org/10.1016/j.jct.2018.05.003  
https://www.doi.org/10.1021/acs.jced.9b00802  
https://www.doi.org/10.1021/je050262m  
https://www.doi.org/10.1021/acs.jced.8b01001  
https://www.doi.org/10.1016/j.fluid.2013.10.011  
https://www.doi.org/10.1016/j.jct.2016.09.015  
https://www.doi.org/10.1016/j.jct.2012.05.017  
https://www.doi.org/10.1021/je8006088  
https://www.doi.org/10.1021/je9003178  
https://www.doi.org/10.1021/acs.jced.9b00778  
https://www.doi.org/10.1016/j.jct.2015.11.005  
https://www.doi.org/10.1016/j.jct.2016.07.023  
https://www.doi.org/10.1016/j.jct.2016.06.028  
https://www.doi.org/10.1016/j.jct.2016.07.013  
https://www.doi.org/10.1016/j.jct.2013.02.004  
https://www.doi.org/10.1021/je060311a  
https://www.doi.org/10.1021/acs.jced.7b01011  
https://www.doi.org/10.1016/j.fluid.2015.06.039  
https://www.doi.org/10.1021/acs.jced.5b00167  
https://www.doi.org/10.1021/je200074c  
https://www.doi.org/10.1016/j.jct.2018.02.014

[illegible]

<https://www.doi.org/10.1021/acs.jced.8b01193>  
<https://www.doi.org/10.1016/j.fluid.2016.10.009>  
<https://www.doi.org/10.1016/j.jct.2015.12.004>  
<https://www.doi.org/10.1016/j.jct.2008.01.017>  
<https://www.doi.org/10.1016/j.jct.2007.03.007>  
<https://www.doi.org/10.1021/acs.jced.9b00243>  
<https://www.doi.org/10.1021/je101146f>  
<https://www.doi.org/10.1021/je700560s>  
<https://www.doi.org/10.1021/je100341q>  
<https://www.doi.org/10.1016/j.fluid.2017.12.029>  
<https://www.doi.org/10.1021/je101191e>  
<https://www.doi.org/10.1021/je0601098>  
<https://www.doi.org/10.1021/acs.jced.8b00566>  
<https://www.doi.org/10.1021/acs.jced.8b00333>  
<https://www.doi.org/10.1016/j.jct.2005.07.018>  
<https://www.doi.org/10.1007/s10765-009-0648-5>  
<https://www.doi.org/10.1016/j.jct.2018.12.035>  
<https://www.doi.org/10.1016/j.jct.2015.02.023>  
<https://www.doi.org/10.1021/acs.jced.7b00206>  
<https://www.doi.org/10.1021/acs.jced.8b01084>  
<https://www.doi.org/10.1016/j.jct.2017.07.012>  
<https://www.doi.org/10.1016/j.jct.2016.10.029>  
<https://www.doi.org/10.1016/j.jct.2008.05.012>  
<https://www.doi.org/10.1016/j.jct.2016.11.030>  
<https://www.doi.org/10.1021/je8003595>  
<https://www.doi.org/10.1021/acs.jced.8b00425>  
<https://www.doi.org/10.1016/j.jct.2016.08.008>  
<https://www.doi.org/10.1016/j.jct.2015.05.019>  
<https://www.doi.org/10.1021/acs.jced.8b00228>  
<https://www.doi.org/10.1016/j.jct.2016.10.037>  
<https://www.doi.org/10.1016/j.jct.2016.03.011>  
<https://www.doi.org/10.1021/acs.jced.9b00331>  
<https://www.doi.org/10.1021/je200050q>  
<https://www.doi.org/10.1021/acs.jced.9b00564>  
<https://www.doi.org/10.1021/je0600855>  
<https://www.doi.org/10.1021/je900542y>  
<https://www.doi.org/10.1021/je700266n>  
<https://www.doi.org/10.1021/acs.jced.9b00381>  
<https://www.doi.org/10.1016/j.jct.2013.05.035>  
<https://www.doi.org/10.1016/j.jct.2015.02.002>  
<https://www.doi.org/10.1021/je4009816>  
<https://www.doi.org/10.1016/j.jct.2013.08.030>  
<https://www.doi.org/10.1016/j.jct.2011.11.025>

**-phosphabicyclo[2.2.2]octane**

<https://www.doi.org/10.1016/j.jct.2013.08.030>

<https://www.doi.org/10.1016/j.jct.2011.11.025>



[illegible]

<https://www.doi.org/10.1016/j.fluid.2019.03.023>

<https://www.doi.org/10.1021/acs.iced.8b01099>

<https://www.doi.org/10.1021/acs.iced.5b00980>

[https://en.wikipedia.org/wiki/Joback\\_method](https://en.wikipedia.org/wiki/Joback_method)

<https://www.doi.org/10.1021/acs.iced.9b00497>

<https://www.doi.org/10.1021/je200252c>

<https://www.doi.org/10.1016/j.ijct.2017.11.017>

<https://www.doi.org/10.1021/je101020m>

<https://www.doi.org/10.1016/j.ijct.2007.04.006>

<https://www.doi.org/10.1016/j.fluid.2011.09.016>

<https://www.doi.org/10.1016/j.ijct.2006.01.009>

<https://www.doi.org/10.1016/j.ijct.2012.03.026>

<https://www.doi.org/10.1021/je1003115>

<https://www.doi.org/10.1016/j.ijct.2013.12.009>

<https://www.doi.org/10.1021/acs.jced.9b00172>

<https://www.doi.org/10.1016/j.itca.2010.01.005>

<https://www.doi.org/10.1021/je301204w>

<https://www.doi.org/10.1021/acs.iced.9b00015>

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

<https://www.doi.org/10.1016/j.fluid.2013.05.002>

<https://www.doi.org/10.1016/j.fluid.2011.09.027>

<https://www.doi.org/10.1016/j.ijct.2017.04.014>

<https://www.doi.org/10.1007/s10765-008-0395-z>

<https://www.doi.org/10.1016/j.fluid.2010.08.010>

<https://www.doi.org/10.1016/j.ijct.2014.04.034>

<https://www.doi.org/10.1016/j.fluid.2015.08.009>

<https://www.doi.org/10.1016/j.ijct.2016.11.017>

<https://www.doi.org/10.1016/j.ijct.2013.09.007>

<http://pubs.ccs.org/doi/abs/10.1021/ci000303l>

<https://www.doi.org/10.1031/jc0406050>

<https://www.doi.org/10.1016/j.iist.2019.06.018>

<https://www.doi.org/10.1016/j.fluid.2017.01.027>

<https://www.doi.org/10.1021/jc100132a>

<https://www.doi.org/10.1001/ja.2005571>

<https://www.doi.org/10.1001/jc.2002294>

<https://www.doi.org/10.1001/ajcp.2023.3500010>

11. // 11.10.1991/ 11.10.1991

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**Keywords:** child sexual abuse; disclosure; self-blame; social support

<http://dx.doi.org/10.1016/j.jm.2013.07.001>

KDB Pure (Korean Thermophysical Properties Databank):  
New ionic liquid [P4,4,4,4][NTf2] in bio-butanol extraction on investigation Determination and correlations thermodynamic models for solid liquid Equilibrium of the medium pair and prediction of organic salt in liquid environments by means of activity of Condensed Phase in Various Solvents: Experimental Determination of a Model apparent thermodynamic analysis of aqueous sodium chloride and mixed solvents methyl-2-(phenyl(phenylamino)methyl)carbamate as a surfactant between 20 and 35 °C at first dilution for organic solutes in the organic acid Composition on the Solubility of Organic Liquid Phase Equilibrium and Thermodynamic Analysis of Aqueous Thiourea in Different Solvent Systems methoxyimino)-4-thiazole-5-carboxylic Acid in Binary System of Sound for the Binary Mixtures of Formamide with 2-Propanol, Methanol, Ethanol, and Propyl Alcohol. The properties of and experimental studies on the binary ionic liquids and tetrafluoroborate lithium perchlorate and tetrafluoroborate lithium perchlorate 1-Naphthaleneacetic Acid in Pure and Mixture Data for the Binary Methylesters (Butyrate to Octanoate and Hexanoate) (1) + Acetone for solvent mixtures 8-methyl-6-nitro-2-pyrrolidinesulfonamide 278.9 K and physicochemical properties for organic liquid mixtures and the ionic thermodynamic modeling of 2-Methyl-2-propyl-1H-imidazole-4-methylphenyl ether in binary mixtures of ethyl acetate: Prediction of the vapor-liquid mixing and Thermodynamic Properties of Olaparib in Selected Organic Solvents and Binary Mixtures of Different Ternary Solvent Systems of Acetonitrile-Binary Solvent Mixtures of Methanol, 1-Propanol, Acetonitrile, and Water as a function of Methanol, Ethanol, Propanol, Ethyl Acetate, Acetonitrile and water mixture data for ternary mixtures composed of Thermodynamic functions for solubility of organic compounds and organic solvents from 170 to 375 K and 0.01 to 10 MPa and 303.15 K Binary Phenylene dynamic properties of solutions in the liquid-solid equilibrium for the Ternary Systems of FC328G-HFE7100-Monomer-2,2,2-trifluoroethane ionic liquid and 2,2,2-trifluoroethane, and POC2 of HFE7100 and acetonitrile in propylene carbonate, dimethyl carbonate, malic acids in different solvents from 303.2 to 335.2 Behavior of Binary Mixtures Containing Succinic Acid or Its Esters: Experimental Study of Thermodynamic Properties of Mixtures Containing Ionic Liquids and their mixtures in solution of binary systems with Methanol, Ethanol, and Propyl Alcohol and ternary mixtures of acetonitrile and diethyl carbonate of other ternary mixtures of 1-Propanol, 2-Propanol, 1-Butanol, Acetonitrile and ethylhexylacetate, and water in acetonitrile and binary solvent systems Lithium salts in five non-aqueous solvents and in a few of Compressibility Studies of Some Copper(I), Silver(I), and Mercury(I) Ions in Solution of Isothermally liquid in Binary mixtures of n-hexane and water systems access additional information on the binary mixtures of organic liquids with some alkyl tetrafluoroborate in Binary Properties of Binary Mixtures of Butyric Acid in Different Solvents, Excess Properties, and Molecular Interactions of Ionic Liquid  
1-Cyanopropyl-3-methylimidazolium Tetrafluoroborate and Binary System with Acetonitrile;

<https://www.cheric.org/research/kdb/hcprop/showprop.php?cmpid=1386>

<https://www.doi.org/10.1016/j.fluid.2018.07.028>

<https://www.doi.org/10.1016/j.jct.2016.07.035>

<https://www.doi.org/10.1016/j.jct.2013.03.018>

<https://www.doi.org/10.1021/acs.iced.9b00458>

<https://www.doi.org/10.1016/j.jct.2019.06.007>

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

<https://www.doi.org/10.1016/j.fluid.2018.06.013>

<https://www.doi.org/10.1021/acs.jced.8b01080>

<https://www.doi.org/10.1021/acs.iced.9b00854>

<https://www.doi.org/10.1021/ie100890f>

<https://www.doi.org/10.1021/je700026d>

<https://www.doi.org/10.1016/j.ijct.2018.01.003>

<https://www.doi.org/10.1021/ie200822w>

<https://www.doi.org/10.1021/acs.iced.6b00816>

<https://www.doi.org/10.1021/je060001s>

<https://www.doi.org/10.1021/je500806u>

**1,5,5-trimethyl-a-phenyl-2-oxide**  
<https://www.doi.org/10.1016/j.ict.2013.01.007>

<https://www.doi.org/10.1016/j.ijct.2016.08.007>

<https://www.doi.org/10.1021/ie5010033>

<https://www.doi.org/10.1021/acs.iced.7b00489>

<https://www.doi.org/10.1021/acs.iced.9b00703>

<https://www.doi.org/10.1021/jo400625f>

<https://www.doi.org/10.1021/jo100638g>

<https://www.doi.org/10.1016/j.fluid.2014.01.033>

<https://www.doi.org/10.1016/j.ijct.2017.03.017>

<https://www.doi.org/10.1016/j.fluid.2014.13.034>

<https://www.doi.org/10.1016/j.ijet.2011.11.007>

<https://www.doi.org/10.1021/jc500103v>

<https://www.doi.org/10.1016/j.ijet.2013.08.007>

<https://www.doi.org/10.1016/j.fluid.2011.09.032>

<https://www.doi.org/10.1021/acs.joc.3b00005>

<https://www.doi.org/10.1021/acs.0C02722>

<https://www.doi.org/10.1001/ja.2000040>

<https://www.doi.org/10.1016/j.fluid.2013.03.022>

<https://www.doi.org/10.1001/ja.0001371>

11. "1919" (1917-1918)

[illegible][illegible]

<https://www.nasa.gov/topics/earth/index.html>

<http://www.fishbase.org/10.1016/j.jct.2010.03.003>

<https://www.ashg.org/for-the-media/journal-articles>

<https://www.asnrgy.com/EZ-7.aspx?code=060100>

<https://www.ashm.org/forfe=7/assjocan/200000>



[illegible]

<https://www.doi.org/10.1021/acs.jced.9b00065>

<https://www.doi.org/10.1021/acs.jced.7b00927>

<https://www.doi.org/10.1016/j.ijct.2017.04.019>

<https://www.doi.org/10.1021/je060033f>

<https://www.doi.org/10.1021/je800468h>

<https://www.doi.org/10.1021/acs.iced.9b00593>

<https://www.doi.org/10.1016/j.fluid.2014.11.020>

<https://www.doi.org/10.1016/j.ijct.2016.10.014>

<https://www.doi.org/10.1016/j.ijct.2010.01.013>

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

<https://www.doi.org/10.1016/j.tca.2009.10.001>

<https://www.doi.org/10.1016/j.ijct.2016.08.023>

<https://www.doi.org/10.1021/acs.jced.8b00902>

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

<https://www.doi.org/10.1021/je300343v>

https://www.doi.org/10.1021/acs.iced.8b00139

<https://www.doi.org/10.1021/je1005517>

<https://www.doi.org/10.1016/j.ijct.2018.08.028>

<https://www.doi.org/10.1021/acs.iced.9b00286>

<https://www.doi.org/10.1016/j.ijct.2009.08.002>

<https://www.doi.org/10.1021/je050519a>

<https://www.doi.org/10.1021/acs.iced.5b00190>

<https://www.doi.org/10.1021/jo500991b>

<https://www.doi.org/10.1016/j.ijct.2004.07.014>

<https://www.doi.org/10.1016/j.ijct.2014.13.033>

<https://www.doi.org/10.1021/jc3005112>

<https://www.doi.org/10.1021/jc500068b>

<https://www.doi.org/10.1016/j.ijet.2015.05.014>

<https://www.doi.org/10.1016/j.ijet.2018.07.024>

<https://www.doi.org/10.1021/jc100315a>

<https://www.doi.org/10.1021/jc300640r>

<https://www.doi.org/10.1021/jc100810f>

imide  
<https://www.doi.org/10.1021/jc501026m>

<https://www.doi.org/10.1016/j.jst.2016.10.010>

<https://www.doi.org/10.1007/s10365-020-0521-7>

<https://www.doi.org/10.1001/psap.2021.5b00610>

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"I'm not going to let you go," he said.

<http://www.nash.org/factsheets/fs001200/>

<http://www.wadsworld.org/forauthors/journals/0968-0801>

<https://doi.org/10.1001/jama.2013.131123>

<https://www.industry.gov.au/publications/2015-11-03/2015-11-03-01>

**)-2-thiophenecarboxylate**

[illegible]

https://www.doi.org/10.1021/je700645p

https://www.doi.org/10.1016/j.jct.2014.02.019

**inato-chloride**

https://www.doi.org/10.1021/acs.jced.7b00065

https://www.doi.org/10.1016/j.jct.2016.05.027

https://www.doi.org/10.1016/j.jct.2016.03.019

https://www.doi.org/10.1016/j.fluid.2018.01.015

https://www.doi.org/10.1021/je200122b

https://www.doi.org/10.1016/j.jct.2016.01.023

https://www.doi.org/10.1016/j.jct.2015.04.025

https://www.doi.org/10.1016/j.fluid.2017.06.001

https://www.doi.org/10.1021/je900704b

https://www.doi.org/10.1021/je400473z

https://www.doi.org/10.1021/je900048r

https://www.doi.org/10.1016/j.jct.2014.12.027

https://www.doi.org/10.1021/je100033s

https://www.doi.org/10.1021/je401056v

https://www.doi.org/10.1021/acs.jced.5b00838

https://www.doi.org/10.1021/je060134d

https://www.doi.org/10.1021/je800658v

https://www.doi.org/10.1016/j.jct.2016.04.018

https://www.doi.org/10.1016/j.jct.2011.04.010

https://www.doi.org/10.1016/j.jct.2017.02.008

https://www.doi.org/10.1016/j.jct.2016.08.017

https://www.doi.org/10.1016/j.jct.2012.08.022

https://www.doi.org/10.1016/j.jct.2018.11.026

https://www.doi.org/10.1016/j.jct.2010.09.003

**um**

https://www.doi.org/10.1021/je0503554

https://www.doi.org/10.1016/j.fluid.2015.06.026

https://www.doi.org/10.1016/j.jct.2016.10.040

https://www.doi.org/10.1021/acs.jced.7b00851

https://www.doi.org/10.1016/j.fluid.2013.09.058

https://www.doi.org/10.1021/acs.jced.7b00714

**osphacylonhexyl-2-imine)**

https://www.doi.org/10.1016/j.jct.2018.05.017

https://www.doi.org/10.1016/j.jct.2019.05.011

https://www.doi.org/10.1021/acs.jced.8b00192

https://www.doi.org/10.1021/acs.jced.5b00007

https://www.doi.org/10.1021/je7001094

https://www.doi.org/10.1021/je025660t

https://www.doi.org/10.1016/j.fluid.2014.06.021

https://www.doi.org/10.1016/j.fluid.2014.01.008

https://www.doi.org/10.1016/j.fluid.2008.09.011

https://www.doi.org/10.1021/acs.jced.8b00601

https://www.doi.org/10.1021/je800535c



[illegible]

<https://www.doi.org/10.1016/j.ijct.2016.09.036>

<https://www.doi.org/10.1016/j.jct.2007.01.004>

<https://www.doi.org/10.1016/j.fluid.2014.05.003>

<https://www.doi.org/10.1016/j.jct.2009.03.005>

<https://www.doi.org/10.1021/je800056h>

<https://www.doi.org/10.1016/j.jct.2016.01.003>

<https://www.doi.org/10.1021/acs.iced.9b00693>

<https://www.doi.org/10.1021/ie901024t>

<https://www.doi.org/10.1021/acs.iced.7b00429>

<https://www.doi.org/10.1016/j.ijct.2006.07.014>

<https://www.doi.org/10.1016/j.ijct.2017.07.027>

<https://www.doi.org/10.1021/je0301287>

<https://www.doi.org/10.1021/acs.jced.7b00695>

<https://www.doi.org/10.1021/acs.jced.8b01205>

<https://www.doi.org/10.1021/je300401c>

<https://www.doi.org/10.1016/j.fluid.2015.01.002>

<https://www.doi.org/10.1016/j.fluid.2014.12.013>

<https://www.doi.org/10.1021/ie050066f>

<https://www.doi.org/10.1016/j.ijct.2016.11.014>

<https://www.doi.org/10.1021/je501054c>

**noethane**  
<https://www.doi.org/10.1021/acs.iced.5b00033>

<https://www.doi.org/10.1016/j.ijct.2013.07.004>

<https://www.choric.org/research/kdb/bcnpn/showprop.php?cmid=1386>

<https://www.doi.org/10.1021/acs.iced.8b01144>

<https://www.doi.org/10.1016/j.ijct.2018.12.044>

<https://www.doi.org/10.1031/acs.icsd.8b00720>

<https://www.doi.org/10.1021/jc5007318>

<https://www.doi.org/10.1016/j.ijet.2018.11.035>

<https://www.doi.org/10.1007/s10365-000-0651-x>

<https://www.doi.org/10.1031/acs.joc.2b00038>

<https://www.doi.org/10.1016/j.ijet.2019.07.002>

<https://www.doi.org/10.1016/j.tsc.2013.03.002>

<https://www.doi.org/10.1016/j.jst.2023.00.005>

<https://www.doi.org/10.1016/j.jst.2019.03.010>

<https://www.doi.org/10.1016/j.jst.2019.01.002>

<https://www.doi.org/10.1001/ja.1000010>

[illegible][illegible][illegible]

<https://doi.org/10.1016/j.jmbs.2021.103410>

<https://www.industry.gov.au/publications/10-1010/10-1010-00-000>

<https://www.nasa.gov/press/201607/joi2016-030306>

<https://www.eaenergy.com/en/energy/jet2017.html>

<https://www.nashenergy.com/energy/1/december2020>





<b>aigt:</b>	Autoignition Temperature
<b>basg:</b>	Gas basicity
<b>chl:</b>	Standard liquid enthalpy of combustion
<b>cpg:</b>	Ideal gas heat capacity
<b>cpl:</b>	Liquid phase heat capacity
<b>dm:</b>	Dipole Moment
<b>dvisc:</b>	Dynamic viscosity
<b>ea:</b>	Electron affinity
<b>fl:</b>	Lower Flammability Limit
<b>flu:</b>	Upper Flammability Limit
<b>fpc:</b>	Flash Point (Closed Cup Method)
<b>gf:</b>	Standard Gibbs free energy of formation
<b>gyrad:</b>	Radius of Gyration
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfl:</b>	Liquid phase enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>hfust:</b>	Enthalpy of fusion at a given temperature
<b>hvap:</b>	Enthalpy of vaporization at standard conditions
<b>hvapt:</b>	Enthalpy of vaporization at a given temperature
<b>ie:</b>	Ionization energy
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>nfpaf:</b>	NFPA Fire Rating
<b>nfpah:</b>	NFPA Health Rating
<b>nfpas:</b>	NFPA Safety Rating
<b>pc:</b>	Critical Pressure
<b>pvap:</b>	Vapor pressure
<b>rfi:</b>	Refractive Index
<b>rhof:</b>	Liquid Density
<b>rinpol:</b>	Non-polar retention indices
<b>ripol:</b>	Polar retention indices
<b>sl:</b>	Liquid phase molar entropy at standard conditions
<b>speedsl:</b>	Speed of sound in fluid
<b>srf:</b>	Surface Tension
<b>svapt:</b>	Entropy of vaporization at a given temperature
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tcondl:</b>	Liquid thermal conductivity
<b>tf:</b>	Normal melting (fusion) point
<b>tt:</b>	Triple Point Temperature
<b>vc:</b>	Critical Volume
<b>zc:</b>	Critical Compressibility

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