

# Acetic acid, chloro-, ethyl ester

Other names:	2-Chloroacetic acid ethyl ester
	Acetic acid, 2-chloro-, ethyl ester
	Chloroacetic acid ethyl ester
	Ethyl chloracetate
	Ethyl chloroacetate
	Ethyl chloroethanoate
	Ethyl ester of chloroacetic acid
	Ethyl monochloracetate
	Ethyl monochloroacetate
	Ethyl «alpha»-chloroacetate
	Ethyl «alpha»-chloroacetate
	Ethylester kyseliny chloroctove
	NSC 8833
	UN 1181
	chloroacetic acid, ethyl ester
	ethanoic acid, chloro-, ethyl ester
Inchi:	InChI=1S/C4H7ClO2/c1-2-7-4(6)3-5/h2-3H2,1H3
InchiKey:	VEUUMBGHMNQHGGO-UHFFFAOYSA-N
Formula:	C4H7ClO2
SMILES:	CCOC(=O)CCl
Mol. weight [g/mol]:	122.55
CAS:	105-39-5

## Physical Properties

Property code	Value	Unit	Source
chl	-2052.00	kJ/mol	NIST Webbook
gf	-263.05	kJ/mol	Joback Method
hf	-386.43	kJ/mol	Joback Method
hfus	13.10	kJ/mol	Joback Method
hvap	49.48	kJ/mol	NIST Webbook
hvap	49.47 ± 0.08	kJ/mol	NIST Webbook
hvap	49.50 ± 0.10	kJ/mol	NIST Webbook
log10ws	-0.51		Crippen Method
logp	0.788		Crippen Method
mcvol	86.900	ml/mol	McGowan Method
pc	3940.66	kPa	Joback Method
rinpol	791.00		NIST Webbook

rinpol	806.40		NIST Webbook
rinpol	791.00		NIST Webbook
rinpol	776.00		NIST Webbook
rinpol	774.00		NIST Webbook
rinpol	802.00		NIST Webbook
rinpol	839.00		NIST Webbook
rinpol	806.40		NIST Webbook
rinpol	839.00		NIST Webbook
rinpol	810.00		NIST Webbook
rinpol	810.00		NIST Webbook
rinpol	832.00		NIST Webbook
ripol	1336.00		NIST Webbook
ripol	1380.00		NIST Webbook
ripol	1302.00		NIST Webbook
ripol	1292.00		NIST Webbook
ripol	1337.00		NIST Webbook
ripol	1355.00		NIST Webbook
ripol	1302.00		NIST Webbook
ripol	1380.00		NIST Webbook
tb	404.64	K	Joback Method
tc	592.13	K	Joback Method
tf	236.92	K	Joback Method
vc	0.333	m3/kmol	Joback Method

# Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	188.84	J/molxK	592.13	Joback Method
cpg	157.93	J/molxK	435.89	Joback Method
cpg	164.55	J/molxK	467.14	Joback Method
cpg	170.96	J/molxK	498.39	Joback Method
cpg	177.14	J/molxK	529.63	Joback Method
cpg	183.10	J/molxK	560.88	Joback Method
cpg	151.10	J/molxK	404.64	Joback Method

dvisc	0.0009692	Paxs	303.15	Volumetric and Transport Properties of Binary Liquid Mixtures of Phenylacetoneitrile with Aliphatic Esters at Temperatures of (303.15 to 313.15) K
dvisc	0.0008979	Paxs	308.15	Volumetric and Transport Properties of Binary Liquid Mixtures of Phenylacetoneitrile with Aliphatic Esters at Temperatures of (303.15 to 313.15) K
dvisc	0.0008341	Paxs	313.15	Volumetric and Transport Properties of Binary Liquid Mixtures of Phenylacetoneitrile with Aliphatic Esters at Temperatures of (303.15 to 313.15) K
hvapt	48.50	kJ/mol	358.00	NIST Webbook
hvapt	45.00	kJ/mol	346.00	NIST Webbook
hvapt	40.43	kJ/mol	417.40	NIST Webbook
rfi	1.42000		298.15	Density, Viscosity, Refractive Index, and Speed of Sound in the Binary Mixtures of Ethyl Chloroacetate + Cyclohexanone, + Chlorobenzene, + Bromobenzene, or + Benzyl Alcohol at (298.15, 303.15, and 308.15) K

rfi	1.41730		303.15	Density, Viscosity, Refractive Index, and Speed of Sound in the Binary Mixtures of Ethyl Chloroacetate + Cyclohexanone, + Chlorobenzene, + Bromobenzene, or + Benzyl Alcohol at (298.15, 303.15, and 308.15) K
rfi	1.41480		308.15	Density, Viscosity, Refractive Index, and Speed of Sound in the Binary Mixtures of Ethyl Chloroacetate + Cyclohexanone, + Chlorobenzene, + Bromobenzene, or + Benzyl Alcohol at (298.15, 303.15, and 308.15) K
rhoI	1144.11	kg/m3	298.15	Excess molar volumes, viscosity deviations and excess thermal expansion coefficients for binary and ternary mixtures consist of diethylketone + 2-butanol + ethylchloroacetate at (298.15, 308.15 and 318.15) K

rhoI	1131.46	kg/m3	308.15	Excess molar volumes, viscosity deviations and excess thermal expansion coefficients for binary and ternary mixtures consist of diethylketone + 2-butanol + ethylchloroacetate at (298.15, 308.15 and 318.15) K
rhoI	1118.74	kg/m3	318.15	Excess molar volumes, viscosity deviations and excess thermal expansion coefficients for binary and ternary mixtures consist of diethylketone + 2-butanol + ethylchloroacetate at (298.15, 308.15 and 318.15) K
speedsl	1229.00	m/s	303.15	Density and Speed of Sound of Binary Mixtures of N-Methylacetamide with Ethyl Acetate, Ethyl Chloroacetate, and Ethyl Cyanoacetate in the Temperature Interval (303.15 to 318.15) K
speedsl	1212.00	m/s	308.15	Density and Speed of Sound of Binary Mixtures of N-Methylacetamide with Ethyl Acetate, Ethyl Chloroacetate, and Ethyl Cyanoacetate in the Temperature Interval (303.15 to 318.15) K

speedsl	1198.00	m/s	313.15	Density and Speed of Sound of Binary Mixtures of N-Methylacetamide with Ethyl Acetate, Ethyl Chloroacetate, and Ethyl Cyanoacetate in the Temperature Interval (303.15 to 318.15) K
speedsl	1177.00	m/s	318.15	Density and Speed of Sound of Binary Mixtures of N-Methylacetamide with Ethyl Acetate, Ethyl Chloroacetate, and Ethyl Cyanoacetate in the Temperature Interval (303.15 to 318.15) K

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.48407e+01
Coeff. B	-3.71707e+03
Coeff. C	-5.37270e+01
Temperature range (K), min.	309.14
Temperature range (K), max.	443.80

Datasets

Viscosity, Pa\*s

Temperature, K - Liquid	Pressure, kPa - Liquid	Viscosity, Pa*s - Liquid
303.15	101.30	0.0012750

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

**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci990307l>

**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)

**Density and Speed of Sound of Binary Mixtures of N-Methylacetamide with Polyethylene Glycols: Properties and Binary Critical Densities at Temperature Ranges from 115 to 310 K and Aromatic Esters at Temperatures of (303.15 to 313.15) K: Density, Viscosity, Refractive Index, and Speed of Sound in the Binary Mixtures of Ethyl Chloroacetate + Carbon Tetrachloride, + Carbon Disulfide, + Carbon Monoxide, + Benzene, + Bromobenzene, + Nitrobenzene, + Toluene, + Diethylketone + 2-butanol + ethylchloroacetate at (298.15, 308.15 and 318.15) K: Density, Viscosity, Compressibilities, viscosities, and excess molar volumes of binary mixtures of alkanooates with tetra- and trichloromethanes at 303.15 K:** <https://www.doi.org/10.1021/je060343y>, <https://www.doi.org/10.1021/je900525f>, <https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>, <https://www.doi.org/10.1021/je0201828>, <https://www.doi.org/10.1016/j.tca.2013.03.008>, <http://webbook.nist.gov/cgi/cbook.cgi?ID=C105395&Units=SI>, <http://link.springer.com/article/10.1007/BF02311772>, <https://www.doi.org/10.1016/j.tca.2004.07.014>

<b>chl:</b>	Standard liquid enthalpy of combustion
<b>cpg:</b>	Ideal gas heat capacity
<b>dvisc:</b>	Dynamic viscosity
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>hvap:</b>	Enthalpy of vaporization at standard conditions
<b>hvapt:</b>	Enthalpy of vaporization at a given temperature
<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>pc:</b>	Critical Pressure
<b>pvap:</b>	Vapor pressure
<b>rfi:</b>	Refractive Index
<b>rho:</b>	Liquid Density
<b>rinpol:</b>	Non-polar retention indices
<b>ripol:</b>	Polar retention indices
<b>speedsl:</b>	Speed of sound in fluid
<b>tb:</b>	Normal Boiling Point Temperature

**tc:** Critical Temperature  
**tf:** Normal melting (fusion) point  
**vc:** Critical Volume

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