Propanoic acid, 2-chloro, 2-methylpropyl ester

Other names:	Isobutyl 2-chloropropanoate
	isobutyl 2-chloropropionate
Inchi:	InChI=1S/C7H13ClO2/c1-5(2)4-10-7(9)6(3)8/h5-6H,4H2,1-3H3
InchiKey:	AEEGJNJYJCGJFZ-UHFFFAOYSA-N
Formula:	C7H13ClO2
SMILES:	CC(C)COC(=O)C(C)Cl
Mol. weight [g/mol]:	164.63
CAS:	62108-67-2

Physical Properties

Property code	Value	Unit	Source
chl	-4033.00 ± 8.40	kJ/mol	NIST Webbook
gf	-242.67	kJ/mol	Joback Method
hf	-549.80 ± 9.60	kJ/mol	NIST Webbook
hfl	-603.30 ± 8.40	kJ/mol	NIST Webbook
hfus	13.82	kJ/mol	Joback Method
hvap	53.60 ± 4.20	kJ/mol	NIST Webbook
log10ws	-1.64		Crippen Method
logp	1.813		Crippen Method
mcvol	129.170	ml/mol	McGowan Method
pc	2875.03	kPa	Joback Method
rinpol	976.00		NIST Webbook
rinpol	994.00		NIST Webbook
rinpol	990.00		NIST Webbook
rinpol	988.00		NIST Webbook
rinpol	999.00		NIST Webbook
rinpol	989.00		NIST Webbook
rinpol	983.00		NIST Webbook
ripol	1340.00		NIST Webbook
ripol	1344.00		NIST Webbook
ripol	1336.00		NIST Webbook
ripol	1342.00		NIST Webbook
ripol	1363.00		NIST Webbook
ripol	1344.00		NIST Webbook
tb	472.40	K	Joback Method
tc	663.16	K	Joback Method
tf	240.73	K	Joback Method

vc	0.488	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	265.59	J/mol×K	472.40	Joback Method
cpg	277.03	J/mol×K	504.19	Joback Method
cpg	288.00	J/mol×K	535.99	Joback Method
cpg	298.52	J/mol×K	567.78	Joback Method
cpg	308.57	J/mol×K	599.57	Joback Method
cpg	318.17	J/mol×K	631.37	Joback Method
cpg	327.32	J/mol×K	663.16	Joback Method
dvisc	0.0064287	Pa×s	240.73	Joback Method
dvisc	0.0025878	Pa×s	279.34	Joback Method
dvisc	0.0012993	Pa×s	317.95	Joback Method
dvisc	0.0007574	Pa×s	356.56	Joback Method
dvisc	0.0004906	Pa×s	395.18	Joback Method
dvisc	0.0003433	Pa×s	433.79	Joback Method
dvisc	0.0002547	Pa×s	472.40	Joback Method

Sources

Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C62108672&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

Legend

chl:	Standard liquid enthalpy of combustion
cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions

hfl:	Liquid phase enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	Non-polar retention indices
ripol:	Polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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