Propanoic acid, 2-chloro-, pentyl ester

Other names: Propionic acid, 2-chloro-, pentyl ester
Pentyl 2-chloropropanoate
Inchi: InChI=1S/C8H15ClO2/c1-3-4-5-6-11-8(10)7(2)9/h7H,3-6H2,1-2H3
InchiKey: ASYLXRAVZWKEBQ-UHFFFAOYSA-N
Formula: C8H15ClO2
SMILES: CCCCCOC(=O)C(C)Cl
Mol. weight [g/mol]: 178.66
CAS: 86711-71-9

Physical Properties

Property code	Value	Unit	Source
gf	-231.81	kJ/mol	Joback Method
hf	-474.27	kJ/mol	Joback Method
hfus	19.94	kJ/mol	Joback Method
hvap	46.55	kJ/mol	Joback Method
log10ws	-2.30		Crippen Method
logp	2.347		Crippen Method
mcvol	143.260	ml/mol	McGowan Method
pc	2579.34	kPa	Joback Method
rinpol	1122.00		NIST Webbook
rinpol	1126.00		NIST Webbook
rinpol	1126.00		NIST Webbook
rinpol	1119.00		NIST Webbook
rinpol	1115.00		NIST Webbook
rinpol	1118.00		NIST Webbook
rinpol	1122.00		NIST Webbook
rinpol	1125.00		NIST Webbook
ripol	1527.00		NIST Webbook
ripol	1502.00		NIST Webbook
ripol	1487.00		NIST Webbook
ripol	1521.00		NIST Webbook
ripol	1526.00		NIST Webbook
ripol	1529.00		NIST Webbook
ripol	1500.00		NIST Webbook
ripol	1484.00		NIST Webbook
tb	495.72	K	Joback Method
tc	680.38	K	Joback Method

tf	267.00	K	Joback Method
vc	0.550	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	307.89	J/mol×K	495.72	Joback Method
cpg	363.91	J/mol×K	649.60	Joback Method
cpg	353.67	J/mol×K	618.82	Joback Method
cpg	342.95	J/mol×K	588.05	Joback Method
cpg	331.75	J/mol×K	557.27	Joback Method
cpg	320.07	J/mol×K	526.50	Joback Method
cpg	373.67	J/mol×K	680.38	Joback Method
dvisc	0.0002478	Pa×s	495.72	Joback Method
dvisc	0.0003265	Pa×s	457.60	Joback Method
dvisc	0.0004523	Pa×s	419.48	Joback Method
dvisc	0.0006687	Pa×s	381.36	Joback Method
dvisc	0.0010785	Pa×s	343.24	Joback Method
dvisc	0.0019599	Pa×s	305.12	Joback Method
dvisc	0.0042242	Pa×s	267.00	Joback Method

Sources

McGowan Method: http://link.springer.com/article/10.1007/BF02311772
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=C86711719&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method: https://en.wikipedia.org/wiki/Joback_method

Legend

cpg: Ideal gas heat capacity
dvisc: Dynamic viscosity
gf: Standard Gibbs free energy of formation
hf: Enthalpy of formation at standard conditions
hfus: Enthalpy of fusion at standard conditions

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