3-Chloropropionic acid, heptyl ester

Other names:	Propanoic acid, 3-chloro, heptyl ester
	Heptyl 3-chloropropanoate
Inchi:	InChI=1S/C10H19ClO2/c1-2-3-4-5-6-9-13-10(12)7-8-11/h2-9H2,1H3
InchiKey:	BQZQUWKYBJWQNL-UHFFFAOYSA-N
Formula:	C10H19ClO2
SMILES:	CCCCCCCOC(=O)CCCl
Mol. weight [g/mol]:	206.71
CAS:	74306-04-0

Physical Properties

Property code	Value	Unit	Source
gf	-212.53	kJ/mol	Joback Method
hf	-510.27	kJ/mol	Joback Method
hfus	28.64	kJ/mol	Joback Method
hvap	51.39	kJ/mol	Joback Method
log10ws	-3.02		Crippen Method
logp	3.129		Crippen Method
mcvol	171.440	ml/mol	McGowan Method
pc	2125.60	kPa	Joback Method
rinpol	1387.00		NIST Webbook
rinpol	1385.00		NIST Webbook
rinpol	1399.00		NIST Webbook
rinpol	1394.00		NIST Webbook
rinpol	1393.00		NIST Webbook
rinpol	1391.00		NIST Webbook
rinpol	1399.00		NIST Webbook
rinpol	1386.00		NIST Webbook
rinpol	1387.00		NIST Webbook
rinpol	1405.00		NIST Webbook
rinpol	1402.00		NIST Webbook
ripol	1828.00		NIST Webbook
ripol	1874.00		NIST Webbook
ripol	1883.00		NIST Webbook
ripol	1895.00		NIST Webbook
ripol	1912.00		NIST Webbook
ripol	1848.00		NIST Webbook
ripol	1871.00		NIST Webbook

ripol	1895.00		NIST Webbook
ripol	1848.00		NIST Webbook
tb	541.92	K	Joback Method
tc	719.34	K	Joback Method
tf	304.54	K	Joback Method
vc	0.668	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	398.46	J/mol×K	541.92	Joback Method
cpg	461.13	J/mol×K	689.77	Joback Method
cpg	449.68	J/mol×K	660.20	Joback Method
cpg	437.70	J/mol×K	630.63	Joback Method
cpg	425.18	J/mol×K	601.06	Joback Method
cpg	412.10	J/mol×K	571.49	Joback Method
cpg	472.05	J/mol×K	719.34	Joback Method
dvisc	0.0002192	Pa×s	541.92	Joback Method
dvisc	0.0002844	Pa×s	502.36	Joback Method
dvisc	0.0003858	Pa×s	462.79	Joback Method
dvisc	0.0005540	Pa×s	423.23	Joback Method
dvisc	0.0008572	Pa×s	383.67	Joback Method
dvisc	0.0014664	Pa×s	344.10	Joback Method
dvisc	0.0028840	Pa×s	304.54	Joback Method

Sources

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=C74306040&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

McGowan Method: http://link.springer.com/article/10.1007/BF02311772

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