2- Chloropropionic acid, decyl ester

Other names: Propanoic acid, 2-chloro, decyl ester
Decyl 2-chloropropanoate
Inchi: InChI=1S/C13H25ClO2/c1-3-4-5-6-7-8-9-10-11-16-13(15)12(2)14/h12H,3-11H2,1-2H3
InchiKey: ZULQXFOAUSXCIJ-UHFFFAOYSA-N
Formula: C13H25ClO2
SMILES: CCCCCCCCCCOC(=O)C(C)Cl
Mol. weight [g/mol]: 248.79
CAS: 86711-78-6

Physical Properties

Property code	Value	Unit	Source
gf	-189.71	kJ/mol	Joback Method
hf	-577.47	kJ/mol	Joback Method
hfus	32.89	kJ/mol	Joback Method
hvap	57.69	kJ/mol	Joback Method
log10ws	-4.39		Crippen Method
logp	4.298		Crippen Method
mcvol	213.710	ml/mol	McGowan Method
pc	1667.33	kPa	Joback Method
rinpol	1635.00		NIST Webbook
rinpol	1633.00		NIST Webbook
rinpol	1644.00		NIST Webbook
rinpol	1612.00		NIST Webbook
rinpol	1620.00		NIST Webbook
rinpol	1623.00		NIST Webbook
rinpol	1642.60		NIST Webbook
rinpol	1629.00		NIST Webbook
rinpol	1625.00		NIST Webbook
rinpol	1627.00		NIST Webbook
ripol	2014.00		NIST Webbook
ripol	2022.00		NIST Webbook
ripol	2030.00		NIST Webbook
ripol	1981.00		NIST Webbook
ripol	1966.00		NIST Webbook
ripol	2013.00		NIST Webbook
ripol	2037.00		NIST Webbook
tb	610.12	K	Joback Method

tc	786.44	K	Joback Method
tf	323.35	K	Joback Method
vc	0.831	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	548.21	J/mol×K	610.12	Joback Method
cpg	564.14	J/mol×K	639.51	Joback Method
cpg	579.37	J/mol×K	668.89	Joback Method
cpg	593.91	J/mol×K	698.28	Joback Method
cpg	607.77	J/mol×K	727.67	Joback Method
cpg	620.96	J/mol×K	757.05	Joback Method
cpg	633.51	J/mol×K	786.44	Joback Method
dvisc	0.0031957	Pa×s	323.35	Joback Method
dvisc	0.0013781	Pa×s	371.14	Joback Method
dvisc	0.0007200	Pa×s	418.94	Joback Method
dvisc	0.0004297	Pa×s	466.74	Joback Method
dvisc	0.0002822	Pa×s	514.53	Joback Method
dvisc	0.0001991	Pa×s	562.33	Joback Method
dvisc	0.0001484	Pa×s	610.12	Joback Method

Sources

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
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=C86711786&Units=SI

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