

decyl lactate

Other names:	decyl 2-hydroxypropanoate
Inchi:	InChI=1S/C13H26O3/c1-3-4-5-6-7-8-9-10-11-16-13(15)12(2)14/h12,14H,3-11H2,1-2H3
InchiKey:	FUGCGCXGFWNOSY-UHFFFAOYSA-N
Formula:	C13H26O3
SMILES:	CCCCCCCCCOC(=O)C(C)O
Mol. weight [g/mol]:	230.34
CAS:	42175-34-8

Physical Properties

Property code	Value	Unit	Source
gf	-314.60	kJ/mol	Joback Method
hf	-713.96	kJ/mol	Joback Method
hfus	32.78	kJ/mol	Joback Method
hvap	69.98	kJ/mol	Joback Method
log10ws	-3.06		Aqueous Solubility Prediction Method
logp	3.051		Crippen Method
mvol	207.340	ml/mol	McGowan Method
pc	1857.91	kPa	Joback Method
tb	664.87	K	Joback Method
tc	833.68	K	Joback Method
tf	354.25	K	Joback Method
vc	0.800	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	583.63	J/mol×K	664.87	Joback Method
cpg	649.56	J/mol×K	805.54	Joback Method
cpg	637.60	J/mol×K	777.41	Joback Method
cpg	625.04	J/mol×K	749.27	Joback Method
cpg	611.86	J/mol×K	721.14	Joback Method
cpg	598.06	J/mol×K	693.00	Joback Method
cpg	660.93	J/mol×K	833.68	Joback Method

dvisc	0.0000414	Paxs	664.87	Joback Method
dvisc	0.0000659	Paxs	613.10	Joback Method
dvisc	0.0001143	Paxs	561.33	Joback Method
dvisc	0.0002217	Paxs	509.56	Joback Method
dvisc	0.0004994	Paxs	457.79	Joback Method
dvisc	0.0013843	Paxs	406.02	Joback Method
dvisc	0.0051687	Paxs	354.25	Joback Method

Sources

Joback Method:	https://en.wikipedia.org/wiki/Joback_method
Aqueous Solubility Prediction Method:	http://onschallenge.wikispaces.com/file/view/AqueousDataset002.xlsx/351826032/AqueousDa
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C42175348&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

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
mvol:	McGowan's characteristic volume
pc:	Critical Pressure
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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