

Succinic acid, hexyl 3-methylbutyl ester

Inchi:	InChI=1S/C15H28O4/c1-4-5-6-7-11-18-14(16)8-9-15(17)19-12-10-13(2)3/h13H,4-12H2,1
InchiKey:	VIUUJFDIZCUSEM-UHFFFAOYSA-N
Formula:	C15H28O4
SMILES:	CCCCCOC(=O)CCC(=O)OCCC(C)C
Mol. weight [g/mol]:	272.38

Physical Properties

Property code	Value	Unit	Source
gf	-394.86	kJ/mol	Joback Method
hf	-847.81	kJ/mol	Joback Method
hfus	36.66	kJ/mol	Joback Method
hvap	66.91	kJ/mol	Joback Method
log10ws	-3.58		Crippen Method
logp	3.479		Crippen Method
mcvol	237.090	ml/mol	McGowan Method
pc	1528.27	kPa	Joback Method
rinpol	1807.00		NIST Webbook
rinpol	1807.00		NIST Webbook
tb	694.74	K	Joback Method
tc	872.74	K	Joback Method
tf	388.13	K	Joback Method
vc	0.917	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	677.76	J/mol×K	694.74	Joback Method
cpg	694.10	J/mol×K	724.41	Joback Method
cpg	709.65	J/mol×K	754.07	Joback Method
cpg	724.41	J/mol×K	783.74	Joback Method
cpg	738.39	J/mol×K	813.41	Joback Method
cpg	751.59	J/mol×K	843.08	Joback Method
cpg	764.02	J/mol×K	872.74	Joback Method
dvisc	0.0016719	Pa×s	388.13	Joback Method

dvisc	0.0007909	Pa×s	439.23	Joback Method
dvisc	0.0004373	Pa×s	490.33	Joback Method
dvisc	0.0002704	Pa×s	541.43	Joback Method
dvisc	0.0001817	Pa×s	592.54	Joback Method
dvisc	0.0001300	Pa×s	643.64	Joback Method
dvisc	0.0000977	Pa×s	694.74	Joback Method

Sources

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=U349246&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

cp_g:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
g_f:	Standard Gibbs free energy of formation
h_f:	Enthalpy of formation at standard conditions
h_{fus}:	Enthalpy of fusion at standard conditions
h_{vap}:	Enthalpy of vaporization at standard conditions
log_{10ws}:	Log10 of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mc_{vol}:	McGowan's characteristic volume
p_c:	Critical Pressure
rin_{pol}:	Non-polar retention indices
t_b:	Normal Boiling Point Temperature
t_c:	Critical Temperature
t_f:	Normal melting (fusion) point
v_c:	Critical Volume

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