

Sebacic acid, 3-methylbutyl octyl ester

Inchi:	InChI=1S/C23H44O4/c1-4-5-6-7-12-15-19-26-22(24)16-13-10-8-9-11-14-17-23(25)27-20
InchiKey:	JAXXJFOVSJVBQO-UHFFFAOYSA-N
Formula:	C23H44O4
SMILES:	CCCCCCCCOC(=O)CCCCCCCCC(=O)OCCC(C)C
Mol. weight [g/mol]:	384.59

Physical Properties

Property code	Value	Unit	Source
gf	-327.50	kJ/mol	Joback Method
hf	-1012.93	kJ/mol	Joback Method
hfus	57.38	kJ/mol	Joback Method
hvap	84.72	kJ/mol	Joback Method
log10ws	-6.93		Crippen Method
logp	6.600		Crippen Method
mcvol	349.810	ml/mol	McGowan Method
pc	903.97	kPa	Joback Method
rinpol	2640.00		NIST Webbook
rinpol	2640.00		NIST Webbook
tb	877.78	K	Joback Method
tc	1074.72	K	Joback Method
tf	478.29	K	Joback Method
vc	1.365	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1154.96	J/mol×K	877.78	Joback Method
cpg	1174.58	J/mol×K	910.60	Joback Method
cpg	1192.90	J/mol×K	943.43	Joback Method
cpg	1209.94	J/mol×K	976.25	Joback Method
cpg	1225.74	J/mol×K	1009.07	Joback Method
cpg	1240.32	J/mol×K	1041.89	Joback Method
cpg	1253.72	J/mol×K	1074.72	Joback Method
dvisc	0.0006901	Pa×s	478.29	Joback Method

dvisc	0.0003008	Pa×s	544.87	Joback Method
dvisc	0.0001571	Pa×s	611.45	Joback Method
dvisc	0.0000932	Pa×s	678.03	Joback Method
dvisc	0.0000607	Pa×s	744.62	Joback Method
dvisc	0.0000424	Pa×s	811.20	Joback Method
dvisc	0.0000313	Pa×s	877.78	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U355372&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

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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