Butanedioic acid, octyl phenylmethyl ester

Other names:	Octyl benzyl succinate
Inchi:	InChI=1S/C19H28O4/c1-2-3-4-5-6-10-15-22-18(20)13-14-19(21)23-16-17-11-8-7-9-12-17
InchiKey:	JKDPVSKXVPWGCQ-UHFFFAOYSA-N
Formula:	C19H28O4
SMILES:	CCCCCCCCOC(=O)CCC(=O)OCc1ccccc1
Mol. weight [g/mol]:	320.42
CAS:	119450-16-7

Physical Properties

Property code	Value	Unit	Source
gf	-246.33	kJ/mol	Joback Method
hf	-688.56	kJ/mol	Joback Method
hfus	44.58	kJ/mol	Joback Method
hvap	78.48	kJ/mol	Joback Method
log10ws	-5.10		Crippen Method
logp	4.414		Crippen Method
mcvol	269.690	ml/mol	McGowan Method
pc	1457.90	kPa	Joback Method
rinpol	2344.00		NIST Webbook
tb	813.38	K	Joback Method
tc	1012.05	K	Joback Method
tf	474.63	K	Joback Method
vc	1.040	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	819.71	J/mol×K	813.38	Joback Method
cpg	835.71	J/mol×K	846.49	Joback Method
cpg	850.62	J/mol×K	879.60	Joback Method
cpg	864.47	J/mol×K	912.71	Joback Method
cpg	877.28	J/mol×K	945.83	Joback Method
cpg	889.08	J/mol×K	978.94	Joback Method
cpg	899.89	J/mol×K	1012.05	Joback Method

dvisc	0.0007727	Pa×s	474.63	Joback Method
dvisc	0.0004063	Pa×s	531.09	Joback Method
dvisc	0.0002418	Pa×s	587.55	Joback Method
dvisc	0.0001576	Pa×s	644.00	Joback Method
dvisc	0.0001100	Pa×s	700.46	Joback Method
dvisc	0.0000811	Pa×s	756.92	Joback Method
dvisc	0.0000623	Pa×s	813.38	Joback Method

Sources

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=C119450167&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

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
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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