

Succinic acid, isobutyl 2-phenoxyethyl ester

Inchi:	InChI=1S/C16H22O5/c1-13(2)12-21-16(18)9-8-15(17)20-11-10-19-14-6-4-3-5-7-14/h3-7,
InchiKey:	IWGSNKDKLBAGQH-UHFFFAOYSA-N
Formula:	C16H22O5
SMILES:	CC(C)COC(=O)CCC(=O)OCCOc1ccccc1
Mol. weight [g/mol]:	294.34

Physical Properties

Property code	Value	Unit	Source
gf	-379.03	kJ/mol	Joback Method
hf	-764.14	kJ/mol	Joback Method
hfus	34.48	kJ/mol	Joback Method
hvap	73.82	kJ/mol	Joback Method
log10ws	-2.84		Crippen Method
logp	2.588		Crippen Method
mvol	233.290	ml/mol	McGowan Method
pc	1826.28	kPa	Joback Method
rinpol	2119.00		NIST Webbook
rinpol	2119.00		NIST Webbook
tb	766.72	K	Joback Method
tc	969.72	K	Joback Method
tf	448.05	K	Joback Method
vc	0.883	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	677.69	J/molxK	766.72	Joback Method
cpg	742.52	J/molxK	935.89	Joback Method
cpg	731.63	J/molxK	902.05	Joback Method
cpg	719.71	J/molxK	868.22	Joback Method
cpg	706.75	J/molxK	834.39	Joback Method
cpg	692.74	J/molxK	800.55	Joback Method
cpg	752.38	J/molxK	969.72	Joback Method
dvisc	0.0000653	Paxs	766.72	Joback Method

dvisc	0.0000852	Paxs	713.61	Joback Method
dvisc	0.0001162	Paxs	660.50	Joback Method
dvisc	0.0001671	Paxs	607.38	Joback Method
dvisc	0.0002578	Paxs	554.27	Joback Method
dvisc	0.0004359	Paxs	501.16	Joback Method
dvisc	0.0008347	Paxs	448.05	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
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=U381193&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
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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