

Succinic acid, ethyl 3-methoxybenzyl ester

Inchi:	InChI=1S/C14H18O5/c1-3-18-13(15)7-8-14(16)19-10-11-5-4-6-12(9-11)17-2/h4-6,9H,3,7
InchiKey:	XOUJEOKVDGEISQ-UHFFFAOYSA-N
Formula:	C14H18O5
SMILES:	CCOC(=O)CCC(=O)OCc1cccc(OC)c1
Mol. weight [g/mol]:	266.29

Physical Properties

Property code	Value	Unit	Source
gf	-403.06	kJ/mol	Joback Method
hf	-729.05	kJ/mol	Joback Method
hfus	32.43	kJ/mol	Joback Method
hvap	70.42	kJ/mol	Joback Method
log10ws	-2.71		Crippen Method
logp	2.082		Crippen Method
mcvol	205.110	ml/mol	McGowan Method
pc	2121.68	kPa	Joback Method
rinpol	1983.00		NIST Webbook
rinpol	1983.00		NIST Webbook
tb	726.38	K	Joback Method
tc	930.82	K	Joback Method
tf	453.03	K	Joback Method
vc	0.777	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	566.63	J/molxK	726.38	Joback Method
cpg	580.70	J/molxK	760.45	Joback Method
cpg	593.86	J/molxK	794.53	Joback Method
cpg	606.11	J/molxK	828.60	Joback Method
cpg	617.43	J/molxK	862.68	Joback Method
cpg	627.82	J/molxK	896.75	Joback Method
cpg	637.28	J/molxK	930.82	Joback Method
dvisc	0.0007156	Paxs	453.03	Joback Method

dvisc	0.0004349	Paxs	498.59	Joback Method
dvisc	0.0002873	Paxs	544.15	Joback Method
dvisc	0.0002023	Paxs	589.71	Joback Method
dvisc	0.0001498	Paxs	635.26	Joback Method
dvisc	0.0001155	Paxs	680.82	Joback Method
dvisc	0.0000920	Paxs	726.38	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=U381250&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
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

Latest version available from:

<https://www.chemeo.com/cid/89-819-5/Succinic-acid-ethyl-3-methoxybenzyl-ester.pdf>

Generated by Cheméo on 2024-04-28 13:06:36.200121174 +0000 UTC m=+16598845.120698502.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.