

Succinic acid, 3-methylbut-2-yl tetrahydrofurfuryl ester

Inchi:	InChI=1S/C14H24O5/c1-10(2)11(3)19-14(16)7-6-13(15)18-9-12-5-4-8-17-12/h10-12H,4-9
InchiKey:	UOUYPQGZNCGMBE-UHFFFAOYSA-N
Formula:	C14H24O5
SMILES:	CC(C)C(C)OC(=O)CCC(=O)OCC1CCCO1
Mol. weight [g/mol]:	272.34

Physical Properties

Property code	Value	Unit	Source
gf	-455.29	kJ/mol	Joback Method
hf	-903.97	kJ/mol	Joback Method
hfus	32.46	kJ/mol	Joback Method
hvap	69.06	kJ/mol	Joback Method
log10ws	-2.37		Crippen Method
logp	2.077		Crippen Method
mvol	218.010	ml/mol	McGowan Method
pc	1918.62	kPa	Joback Method
rinpol	1865.00		NIST Webbook
rinpol	1865.00		NIST Webbook
tb	713.65	K	Joback Method
tc	913.71	K	Joback Method
tf	399.33	K	Joback Method
vc	0.818	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	647.55	J/molxK	713.65	Joback Method
cpg	664.56	J/molxK	746.99	Joback Method
cpg	680.53	J/molxK	780.34	Joback Method
cpg	695.46	J/molxK	813.68	Joback Method
cpg	709.38	J/molxK	847.02	Joback Method
cpg	722.29	J/molxK	880.37	Joback Method
cpg	734.21	J/molxK	913.71	Joback Method
dvisc	0.0022756	Paxs	399.33	Joback Method

dvisc	0.0010761	Paxs	451.72	Joback Method
dvisc	0.0005946	Paxs	504.10	Joback Method
dvisc	0.0003674	Paxs	556.49	Joback Method
dvisc	0.0002466	Paxs	608.88	Joback Method
dvisc	0.0001763	Paxs	661.26	Joback Method
dvisc	0.0001324	Paxs	713.65	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=U390714&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

Latest version available from:

<https://www.chemeo.com/cid/92-704-8/Succinic-acid-3-methylbut-2-yl-tetrahydrofurfuryl-ester.pdf>

Generated by Cheméo on 2024-04-19 18:46:37.560553959 +0000 UTC m=+15841646.481131276.

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