

Glutaric acid, tetrahydrofurfuryl tridecyl ester

Inchi:	InChI=1S/C23H42O5/c1-2-3-4-5-6-7-8-9-10-11-12-18-27-22(24)16-13-17-23(25)28-20-21
InchiKey:	VBICSTPSVCFUSY-UHFFFAOYSA-N
Formula:	C23H42O5
SMILES:	CCCCCCCCCCCCOC(=O)CCCC(=O)OCC1CCCO1
Mol. weight [g/mol]:	398.58

Physical Properties

Property code	Value	Unit	Source
gf	-374.63	kJ/mol	Joback Method
hf	-1079.17	kJ/mol	Joback Method
hfus	62.81	kJ/mol	Joback Method
hvap	89.87	kJ/mol	Joback Method
log10ws	-6.27		Crippen Method
logp	5.733		Crippen Method
mvol	344.820	ml/mol	McGowan Method
pc	999.54	kPa	Joback Method
rinpol	2946.00		NIST Webbook
rinpol	2946.00		NIST Webbook
tb	920.45	K	Joback Method
tc	1126.92	K	Joback Method
tf	530.76	K	Joback Method
vc	1.333	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1182.69	J/molxK	920.45	Joback Method
cpg	1201.40	J/molxK	954.86	Joback Method
cpg	1218.68	J/molxK	989.27	Joback Method
cpg	1234.56	J/molxK	1023.68	Joback Method
cpg	1249.08	J/molxK	1058.09	Joback Method
cpg	1262.29	J/molxK	1092.50	Joback Method
cpg	1274.22	J/molxK	1126.92	Joback Method
dvisc	0.0006283	Paxs	530.76	Joback Method

dvisc	0.0003157	Paxs	595.71	Joback Method
dvisc	0.0001816	Paxs	660.66	Joback Method
dvisc	0.0001153	Paxs	725.61	Joback Method
dvisc	0.0000789	Paxs	790.55	Joback Method
dvisc	0.0000572	Paxs	855.50	Joback Method
dvisc	0.0000434	Paxs	920.45	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=U359669&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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