

# Glutaric acid, butyl tetrahydrofurfuryl ester

<b>Inchi:</b>	InChI=1S/C14H24O5/c1-2-3-9-18-13(15)7-4-8-14(16)19-11-12-6-5-10-17-12/h12H,2-11H
<b>InchiKey:</b>	QBHWJHBWBUUQRU-UHFFFAOYSA-N
<b>Formula:</b>	C14H24O5
<b>SMILES:</b>	CCCCOC(=O)CCCC(=O)OCC1CCCO1
<b>Mol. weight [g/mol]:</b>	272.34

## Physical Properties

Property code	Value	Unit	Source
gf	-450.41	kJ/mol	Joback Method
hf	-893.41	kJ/mol	Joback Method
hfus	39.50	kJ/mol	Joback Method
hvap	69.84	kJ/mol	Joback Method
log10ws	-2.50		Crippen Method
logp	2.222		Crippen Method
mcvol	218.010	ml/mol	McGowan Method
pc	1892.00	kPa	Joback Method
rinqol	2011.00		NIST Webbook
tb	714.53	K	Joback Method
tc	908.66	K	Joback Method
tf	429.33	K	Joback Method
vc	0.830	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	646.55	J/molxK	714.53	Joback Method
cpg	663.02	J/molxK	746.88	Joback Method
cpg	678.52	J/molxK	779.24	Joback Method
cpg	693.06	J/molxK	811.59	Joback Method
cpg	706.67	J/molxK	843.95	Joback Method
cpg	719.35	J/molxK	876.30	Joback Method
cpg	731.11	J/molxK	908.66	Joback Method
dvisc	0.0015886	Paxs	429.33	Joback Method
dvisc	0.0008884	Paxs	476.86	Joback Method

dvisc	0.0005521	Paxs	524.40	Joback Method
dvisc	0.0003713	Paxs	571.93	Joback Method
dvisc	0.0002654	Paxs	619.46	Joback Method
dvisc	0.0001990	Paxs	667.00	Joback Method
dvisc	0.0001550	Paxs	714.53	Joback Method

## Sources

<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U359659&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U359659&amp;Units=SI</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>dvisc:</b>	Dynamic viscosity
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>hvap:</b>	Enthalpy of vaporization at standard conditions
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>rinpol:</b>	Non-polar retention indices
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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