

# Glutaric acid, hexyl tetrahydrofurfuryl ester

<b>Inchi:</b>	InChI=1S/C16H28O5/c1-2-3-4-5-11-20-15(17)9-6-10-16(18)21-13-14-8-7-12-19-14/h14H
<b>InchiKey:</b>	BJAANRCRIWFPRT-UHFFFAOYSA-N
<b>Formula:</b>	C16H28O5
<b>SMILES:</b>	CCCCCOC(=O)CCCC(=O)OCC1CCCO1
<b>Mol. weight [g/mol]:</b>	300.39

## Physical Properties

Property code	Value	Unit	Source
gf	-433.57	kJ/mol	Joback Method
hf	-934.69	kJ/mol	Joback Method
hfus	44.68	kJ/mol	Joback Method
hvap	74.29	kJ/mol	Joback Method
log10ws	-3.34		Crippen Method
logp	3.002		Crippen Method
mvol	246.190	ml/mol	McGowan Method
pc	1611.58	kPa	Joback Method
rinpol	2208.00		NIST Webbook
tb	760.29	K	Joback Method
tc	952.60	K	Joback Method
tf	451.87	K	Joback Method
vc	0.942	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	759.76	J/molxK	760.29	Joback Method
cpg	776.78	J/molxK	792.34	Joback Method
cpg	792.77	J/molxK	824.39	Joback Method
cpg	807.73	J/molxK	856.45	Joback Method
cpg	821.68	J/molxK	888.50	Joback Method
cpg	834.65	J/molxK	920.55	Joback Method
cpg	846.65	J/molxK	952.60	Joback Method
dvisc	0.0013349	Paxs	451.87	Joback Method
dvisc	0.0007265	Paxs	503.27	Joback Method

dvisc	0.0004426	Paxs	554.68	Joback Method
dvisc	0.0002933	Paxs	606.08	Joback Method
dvisc	0.0002073	Paxs	657.48	Joback Method
dvisc	0.0001540	Paxs	708.89	Joback Method
dvisc	0.0001192	Paxs	760.29	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=U359662&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U359662&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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