

# Glutaric acid, pentyl tetrahydrofurfuryl ester

<b>Inchi:</b>	InChI=1S/C15H26O5/c1-2-3-4-10-19-14(16)8-5-9-15(17)20-12-13-7-6-11-18-13/h13H,2-
<b>InchiKey:</b>	CYAMANFTXDWMFN-UHFFFAOYSA-N
<b>Formula:</b>	C15H26O5
<b>SMILES:</b>	CCCCCOC(=O)CCCC(=O)OCC1CCCO1
<b>Mol. weight [g/mol]:</b>	286.36

## Physical Properties

Property code	Value	Unit	Source
gf	-441.99	kJ/mol	Joback Method
hf	-914.05	kJ/mol	Joback Method
hfus	42.09	kJ/mol	Joback Method
hvap	72.06	kJ/mol	Joback Method
log10ws	-2.92		Crippen Method
logp	2.612		Crippen Method
mvol	232.100	ml/mol	McGowan Method
pc	1743.37	kPa	Joback Method
rinpol	2111.00		NIST Webbook
rinpol	2111.00		NIST Webbook
tb	737.41	K	Joback Method
tc	930.37	K	Joback Method
tf	440.60	K	Joback Method
vc	0.885	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	702.63	J/molxK	737.41	Joback Method
cpg	776.56	J/molxK	898.21	Joback Method
cpg	763.72	J/molxK	866.05	Joback Method
cpg	749.92	J/molxK	833.89	Joback Method
cpg	735.15	J/molxK	801.73	Joback Method
cpg	719.39	J/molxK	769.57	Joback Method
cpg	788.46	J/molxK	930.37	Joback Method
dvisc	0.0001362	Paxs	737.41	Joback Method

dvisc	0.0001754	Paxs	687.94	Joback Method
dvisc	0.0002350	Paxs	638.47	Joback Method
dvisc	0.0003307	Paxs	589.00	Joback Method
dvisc	0.0004955	Paxs	539.54	Joback Method
dvisc	0.0008054	Paxs	490.07	Joback Method
dvisc	0.0014603	Paxs	440.60	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=U359660&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U359660&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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