

# Glutaric acid, butyl 4-methyl-3-nitrobenzyl ester

<b>Inchi:</b>	InChI=1S/C17H23NO6/c1-3-4-10-23-16(19)6-5-7-17(20)24-12-14-9-8-13(2)15(11-14)18(
<b>InchiKey:</b>	LVRFZEXMUUTLBY-UHFFFAOYSA-N
<b>Formula:</b>	C17H23NO6
<b>SMILES:</b>	CCCCOC(=O)CCCC(=O)OCc1ccc(C)c([N+](=O)[O-])c1
<b>Mol. weight [g/mol]:</b>	337.37

## Physical Properties

Property code	Value	Unit	Source
gf	-246.88	kJ/mol	Joback Method
hf	-680.98	kJ/mol	Joback Method
hfus	49.98	kJ/mol	Joback Method
hvap	91.94	kJ/mol	Joback Method
log10ws	-4.97		Crippen Method
logp	3.460		Crippen Method
mvol	258.930	ml/mol	McGowan Method
pc	1672.79	kPa	Joback Method
rmpol	2569.00		NIST Webbook
tb	929.42	K	Joback Method
tc	1152.32	K	Joback Method
tf	620.74	K	Joback Method
vc	1.010	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	811.03	J/molxK	929.42	Joback Method
cpg	823.17	J/molxK	966.57	Joback Method
cpg	834.07	J/molxK	1003.72	Joback Method
cpg	843.77	J/molxK	1040.87	Joback Method
cpg	852.26	J/molxK	1078.02	Joback Method
cpg	859.58	J/molxK	1115.17	Joback Method
cpg	865.74	J/molxK	1152.32	Joback Method

# Sources

<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307I">http://pubs.acs.org/doi/abs/10.1021/ci990307I</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=U376780&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U376780&amp;Units=SI</a>

# Legend

<b>cpg:</b>	Ideal gas heat capacity
<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>mccvol:</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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