

# Glutaric acid, di(5-methylhex-2-yl) ester

<b>Inchi:</b>	InChI=1S/C19H36O4/c1-14(2)10-12-16(5)22-18(20)8-7-9-19(21)23-17(6)13-11-15(3)4/h1
<b>InchiKey:</b>	SWPFRISFZMPKCU-UHFFFAOYSA-N
<b>Formula:</b>	C19H36O4
<b>SMILES:</b>	CC(C)CCC(C)OC(=O)CCCC(=O)OC(C)CCC(C)C
<b>Mol. weight [g/mol]:</b>	328.49

## Physical Properties

Property code	Value	Unit	Source
gf	-368.50	kJ/mol	Joback Method
hf	-946.21	kJ/mol	Joback Method
hfus	36.45	kJ/mol	Joback Method
hvap	74.65	kJ/mol	Joback Method
log10ws	-5.24		Crippen Method
logp	4.893		Crippen Method
mcvol	293.450	ml/mol	McGowan Method
pc	1174.44	kPa	Joback Method
rinqol	2239.00		NIST Webbook
tb	784.94	K	Joback Method
tc	970.70	K	Joback Method
tf	388.21	K	Joback Method
vc	1.123	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	910.84	J/molxK	784.94	Joback Method
cpg	929.17	J/molxK	815.90	Joback Method
cpg	946.44	J/molxK	846.86	Joback Method
cpg	962.66	J/molxK	877.82	Joback Method
cpg	977.85	J/molxK	908.78	Joback Method
cpg	992.02	J/molxK	939.74	Joback Method
cpg	1005.19	J/molxK	970.70	Joback Method
dvisc	0.0021477	Paxs	388.21	Joback Method
dvisc	0.0007008	Paxs	454.33	Joback Method

dvisc	0.0003040	Paxs	520.45	Joback Method
dvisc	0.0001591	Paxs	586.58	Joback Method
dvisc	0.0000950	Paxs	652.70	Joback Method
dvisc	0.0000624	Paxs	718.82	Joback Method
dvisc	0.0000439	Paxs	784.94	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=U377666&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U377666&amp;Units=SI</a>

## Legend

<b>cp<sub>g</sub>:</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>log<sub>10</sub>ws:</b>	Log <sub>10</sub> of Water solubility in mol/l
<b>log<sub>p</sub>:</b>	Octanol/Water partition coefficient
<b>m<sub>cvol</sub>:</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

Latest version available from:

<https://www.chemeo.com/cid/46-321-4/Glutaric-acid-di-5-methylhex-2-yl-ester.pdf>

Generated by Cheméo on 2024-04-24 15:52:55.4018934 +0000 UTC m=+16263224.322470719.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.