

# Glutaric acid, 2,2-dichloroethyl neopentyl ester

Inchi:	InChI=1S/C12H20Cl2O4/c1-12(2,3)8-18-11(16)6-4-5-10(15)17-7-9(13)14/h9H,4-8H2,1-3H
InchiKey:	QLEGELAPPQLNPY-UHFFFAOYSA-N
Formula:	C12H20Cl2O4
SMILES:	CC(C)(C)COC(=O)CCCC(=O)OCC(Cl)Cl
Mol. weight [g/mol]:	299.19

## Physical Properties

Property code	Value	Unit	Source
gf	-441.14	kJ/mol	Joback Method
hf	-826.12	kJ/mol	Joback Method
hfus	29.87	kJ/mol	Joback Method
hvap	67.70	kJ/mol	Joback Method
log10ws	-3.24		Crippen Method
logp	3.093		Crippen Method
mvol	219.300	ml/mol	McGowan Method
pc	1849.92	kPa	Joback Method
rinpol	1786.00		NIST Webbook
rinpol	1786.00		NIST Webbook
tb	697.73	K	Joback Method
tc	894.59	K	Joback Method
tf	416.58	K	Joback Method
vc	0.837	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	576.25	J/molxK	697.73	Joback Method
cpg	589.82	J/molxK	730.54	Joback Method
cpg	602.57	J/molxK	763.35	Joback Method
cpg	614.51	J/molxK	796.16	Joback Method
cpg	625.68	J/molxK	828.97	Joback Method
cpg	636.08	J/molxK	861.78	Joback Method
cpg	645.74	J/molxK	894.59	Joback Method
dvisc	0.0014409	Paxs	416.58	Joback Method

dvisc	0.0007306	Paxs	463.44	Joback Method
dvisc	0.0004197	Paxs	510.30	Joback Method
dvisc	0.0002646	Paxs	557.15	Joback Method
dvisc	0.0001793	Paxs	604.01	Joback Method
dvisc	0.0001284	Paxs	650.87	Joback Method
dvisc	0.0000962	Paxs	697.73	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=U391610&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U391610&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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