

# Glutaric acid, heptyl nonyl ester

<b>Inchi:</b>	InChI=1S/C21H40O4/c1-3-5-7-9-10-12-14-19-25-21(23)17-15-16-20(22)24-18-13-11-8-6
<b>InchiKey:</b>	YLZNKOSNJSNTGC-UHFFFAOYSA-N
<b>Formula:</b>	C21H40O4
<b>SMILES:</b>	CCCCCCCCCOC(=O)CCCC(=O)OCCCCCCC
<b>Mol. weight [g/mol]:</b>	356.54

## Physical Properties

Property code	Value	Unit	Source
gf	-341.90	kJ/mol	Joback Method
hf	-966.37	kJ/mol	Joback Method
hfus	55.72	kJ/mol	Joback Method
hvap	80.65	kJ/mol	Joback Method
log10ws	-6.34		Crippen Method
logp	5.964		Crippen Method
mcvol	321.630	ml/mol	McGowan Method
pc	1012.95	kPa	Joback Method
rinpol	2490.00		NIST Webbook
tb	832.46	K	Joback Method
tc	1019.85	K	Joback Method
tf	470.75	K	Joback Method
vc	1.260	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1030.26	J/mol×K	832.46	Joback Method
cpg	1048.96	J/mol×K	863.69	Joback Method
cpg	1066.54	J/mol×K	894.92	Joback Method
cpg	1083.03	J/mol×K	926.15	Joback Method
cpg	1098.44	J/mol×K	957.39	Joback Method
cpg	1112.80	J/mol×K	988.62	Joback Method
cpg	1126.13	J/mol×K	1019.85	Joback Method
dvisc	0.0007511	Paxs	470.75	Joback Method
dvisc	0.0003622	Paxs	531.03	Joback Method

dvisc	0.0002027	Paxs	591.32	Joback Method
dvisc	0.0001263	Paxs	651.61	Joback Method
dvisc	0.0000852	Paxs	711.89	Joback Method
dvisc	0.0000612	Paxs	772.17	Joback Method
dvisc	0.0000461	Paxs	832.46	Joback Method

## Sources

<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=U358282&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U358282&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</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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