

# Nonanoic acid, octadecyl ester

<b>Inchi:</b>	InChI=1S/C27H54O2/c1-3-5-7-9-11-12-13-14-15-16-17-18-19-20-22-24-26-29-27(28)25-
<b>InchiKey:</b>	IUITWWQQPHHQUA-UHFFFAOYSA-N
<b>Formula:</b>	C27H54O2
<b>SMILES:</b>	CCCCCCCCCCCCCCCCCOC(=O)CCCCCCCC
<b>Mol. weight [g/mol]:</b>	410.72

## Physical Properties

Property code	Value	Unit	Source
gf	-57.46	kJ/mol	Joback Method
hf	-845.41	kJ/mol	Joback Method
hfus	68.47	kJ/mol	Joback Method
hvap	84.85	kJ/mol	Joback Method
log10ws	-9.99		Crippen Method
logp	9.542		Crippen Method
mvol	398.730	ml/mol	McGowan Method
pc	705.83	kPa	Joback Method
rinpol	2861.00		NIST Webbook
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tb	893.45	K	Joback Method
tc	1097.77	K	Joback Method
tf	466.21	K	Joback Method
vc	1.571	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1350.54	J/molxK	893.45	Joback Method
cpg	1374.21	J/molxK	927.50	Joback Method
cpg	1396.42	J/molxK	961.56	Joback Method
cpg	1417.22	J/molxK	995.61	Joback Method
cpg	1436.68	J/molxK	1029.66	Joback Method
cpg	1454.85	J/molxK	1063.71	Joback Method
cpg	1471.80	J/molxK	1097.77	Joback Method
dvisc	0.0007281	Paxs	466.21	Joback Method

dvisc	0.0002915	Paxs	537.42	Joback Method
dvisc	0.0001446	Paxs	608.62	Joback Method
dvisc	0.0000831	Paxs	679.83	Joback Method
dvisc	0.0000530	Paxs	751.04	Joback Method
dvisc	0.0000366	Paxs	822.24	Joback Method
dvisc	0.0000268	Paxs	893.45	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=U340285&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U340285&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>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>h<sub>vap</sub>:</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>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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