

# 11,19,23-trimethyl-heptatriacontane

<b>Inchi:</b>	InChI=1S/C40H82/c1-6-8-10-12-14-16-17-18-19-21-24-28-34-39(4)36-31-37-40(5)35-30
<b>InchiKey:</b>	XFBXDFLXACOLRR-UHFFFAOYSA-N
<b>Formula:</b>	C40H82
<b>SMILES:</b>	CCCCCCCCCCCCCCC(C)CCCC(C)CCCCCCCC(C)CCCCCCCCC
<b>Mol. weight [g/mol]:</b>	563.08

## Physical Properties

Property code	Value	Unit	Source
gf	278.60	kJ/mol	Joback Method
hf	-884.77	kJ/mol	Joback Method
hfus	88.79	kJ/mol	Joback Method
hvap	103.47	kJ/mol	Joback Method
log10ws	-15.84		Crippen Method
logp	15.418		Crippen Method
mcvol	574.460	ml/mol	McGowan Method
pc	394.61	kPa	Joback Method
rinpol	3780.00		NIST Webbook
rinpol	3780.00		NIST Webbook
rinpol	3780.00		NIST Webbook
tb	1113.28	K	Joback Method
tc	1468.24	K	Joback Method
tf	495.56	K	Joback Method
vc	2.257	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	2175.93	J/molxK	1113.28	Joback Method
cpg	2217.11	J/molxK	1172.44	Joback Method
cpg	2254.57	J/molxK	1231.60	Joback Method
cpg	2288.88	J/molxK	1290.76	Joback Method
cpg	2320.61	J/molxK	1349.92	Joback Method
cpg	2350.35	J/molxK	1409.08	Joback Method
cpg	2378.66	J/molxK	1468.24	Joback Method

dvisc	0.0004037	Paxs	495.56	Joback Method
dvisc	0.0000943	Paxs	598.51	Joback Method
dvisc	0.0000337	Paxs	701.47	Joback Method
dvisc	0.0000157	Paxs	804.42	Joback Method
dvisc	0.0000087	Paxs	907.37	Joback Method
dvisc	0.0000054	Paxs	1010.33	Joback Method
dvisc	0.0000037	Paxs	1113.28	Joback Method

## Sources

<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>
<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=R271877&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R271877&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>mvol:</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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