

# 8,14,18-trimethyloctatriacontane

**Inchi:** InChI=1S/C41H84/c1-6-8-10-12-13-14-15-16-17-18-19-20-21-22-23-24-26-29-34-40(4)3  
**InchiKey:** QEKGEXLUFKHLEN-UHFFFAOYSA-N  
**Formula:** C41H84  
**SMILES:** CCCCCCCCCCCCCCCCCCCCCC(C)CCCC(C)CCCCC(C)CCCCCCC  
**Mol. weight [g/mol]:** 577.11

## Physical Properties

Property code	Value	Unit	Source
gf	287.02	kJ/mol	Joback Method
hf	-905.41	kJ/mol	Joback Method
hfus	91.38	kJ/mol	Joback Method
hvap	105.70	kJ/mol	Joback Method
log10ws	-16.26		Crippen Method
logp	15.808		Crippen Method
mcvol	588.550	ml/mol	McGowan Method
pc	379.98	kPa	Joback Method
rinpol	3874.00		NIST Webbook
tb	1136.16	K	Joback Method
tc	1517.43	K	Joback Method
tf	506.83	K	Joback Method
vc	2.313	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	2246.59	J/molxK	1136.16	Joback Method
cpg	2290.12	J/molxK	1199.71	Joback Method
cpg	2329.56	J/molxK	1263.25	Joback Method
cpg	2365.64	J/molxK	1326.80	Joback Method
cpg	2399.11	J/molxK	1390.34	Joback Method
cpg	2430.68	J/molxK	1453.89	Joback Method
cpg	2461.09	J/molxK	1517.43	Joback Method
dvisc	0.0003383	Paxs	506.83	Joback Method
dvisc	0.0000794	Paxs	611.72	Joback Method

dvisc	0.0000285	Paxs	716.61	Joback Method
dvisc	0.0000133	Paxs	821.49	Joback Method
dvisc	0.0000074	Paxs	926.38	Joback Method
dvisc	0.0000046	Paxs	1031.27	Joback Method
dvisc	0.0000031	Paxs	1136.16	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=R280683&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R280683&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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