

# 14-Heptatricosene, 25-methyl

<b>Inchi:</b>	InChI=1S/C38H76/c1-4-6-8-10-12-14-16-17-18-19-20-21-22-23-24-25-26-27-29-31-33-35
<b>InchiKey:</b>	BLUMBGGQOKMUQGN-GHVJWSGMSA-N
<b>Formula:</b>	C38H76
<b>SMILES:</b>	CCCCCCCCCCCC=CCCCCCCCCCC(C)CCCCCCCCCCCC
<b>Mol. weight [g/mol]:</b>	533.01

## Physical Properties

Property code	Value	Unit	Source
gf	346.86	kJ/mol	Joback Method
hf	-715.71	kJ/mol	Joback Method
hfus	90.86	kJ/mol	Joback Method
hvap	99.75	kJ/mol	Joback Method
log10ws	-15.34		Crippen Method
logp	14.702		Crippen Method
mcvol	541.980	ml/mol	McGowan Method
pc	432.94	kPa	Joback Method
rinsol	3696.00		NIST Webbook
tb	1072.56	K	Joback Method
tc	1385.66	K	Joback Method
tf	497.94	K	Joback Method
vc	2.138	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	2007.56	J/molxK	1072.56	Joback Method
cpg	2045.43	J/molxK	1124.74	Joback Method
cpg	2080.66	J/molxK	1176.93	Joback Method
cpg	2113.68	J/molxK	1229.11	Joback Method
cpg	2144.88	J/molxK	1281.30	Joback Method
cpg	2174.70	J/molxK	1333.48	Joback Method
cpg	2203.55	J/molxK	1385.66	Joback Method
dvisc	0.0003678	Paxs	497.94	Joback Method
dvisc	0.0001039	Paxs	593.71	Joback Method

dvisc	0.0000417	Paxs	689.48	Joback Method
dvisc	0.0000209	Paxs	785.25	Joback Method
dvisc	0.0000122	Paxs	881.02	Joback Method
dvisc	0.0000079	Paxs	976.79	Joback Method
dvisc	0.0000055	Paxs	1072.56	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=R608492&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R608492&amp;Units=SI</a>
<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>

## 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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