

# 13,17-Dimethyltriacontane

<b>Inchi:</b>	InChI=1S/C32H66/c1-5-7-9-11-13-15-17-19-21-23-25-28-32(4)30-26-29-31(3)27-24-22-2
<b>InchiKey:</b>	PBGJCQPDUYBYRS-UHFFFAOYSA-N
<b>Formula:</b>	C32H66
<b>SMILES:</b>	CCCCCCCCCCCC(C)CCCC(C)CCCCCCCCCCCC
<b>Mol. weight [g/mol]:</b>	450.87

## Physical Properties

Property code	Value	Unit	Source
gf	213.68	kJ/mol	Joback Method
hf	-714.37	kJ/mol	Joback Method
hfus	71.59	kJ/mol	Joback Method
hvap	86.05	kJ/mol	Joback Method
log10ws	-12.73		Crippen Method
logp	12.441		Crippen Method
mcvol	461.740	ml/mol	McGowan Method
pc	547.43	kPa	Joback Method
rinpol	3045.00		NIST Webbook
rinpol	3045.00		NIST Webbook
tb	930.68	K	Joback Method
tc	1153.00	K	Joback Method
tf	420.40	K	Joback Method
vc	1.815	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1618.68	J/molxK	930.68	Joback Method
cpg	1647.60	J/molxK	967.73	Joback Method
cpg	1674.72	J/molxK	1004.79	Joback Method
cpg	1700.16	J/molxK	1041.84	Joback Method
cpg	1724.04	J/molxK	1078.89	Joback Method
cpg	1746.45	J/molxK	1115.95	Joback Method
cpg	1767.52	J/molxK	1153.00	Joback Method
dvisc	0.0012543	Paxs	420.40	Joback Method

dvisc	0.0003218	Paxs	505.45	Joback Method
dvisc	0.0001222	Paxs	590.49	Joback Method
dvisc	0.0000592	Paxs	675.54	Joback Method
dvisc	0.0000337	Paxs	760.59	Joback Method
dvisc	0.0000215	Paxs	845.63	Joback Method
dvisc	0.0000149	Paxs	930.68	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=R530542&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R530542&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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