

# 14,18,22-Trimethyltetracontane

**Inchi:** InChI=1S/C43H88/c1-6-8-10-12-14-16-18-19-20-21-22-24-26-28-30-32-36-42(4)38-34-40  
**InchiKey:** KERBDCMSTFMOCU-UHFFFAOYSA-N  
**Formula:** C43H88  
**SMILES:** CCCCCCCCCCCCCCCCCC(C)CCCC(C)CCCC(C)CCCCCCCCCCCCC  
**Mol. weight [g/mol]:** 605.16

## Physical Properties

Property code	Value	Unit	Source
gf	303.86	kJ/mol	Joback Method
hf	-946.69	kJ/mol	Joback Method
hfus	96.56	kJ/mol	Joback Method
hvap	110.15	kJ/mol	Joback Method
log10ws	-17.10		Crippen Method
logp	16.588		Crippen Method
mcvol	616.730	ml/mol	McGowan Method
pc	353.06	kPa	Joback Method
rinpol	4071.00		NIST Webbook
rinpol	4071.00		NIST Webbook
tb	1181.92	K	Joback Method
tc	1624.44	K	Joback Method
tf	529.37	K	Joback Method
vc	2.425	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	2388.44	J/molxK	1181.92	Joback Method
cpg	2437.38	J/molxK	1255.67	Joback Method
cpg	2481.48	J/molxK	1329.43	Joback Method
cpg	2521.94	J/molxK	1403.18	Joback Method
cpg	2559.95	J/molxK	1476.94	Joback Method
cpg	2596.73	J/molxK	1550.69	Joback Method
cpg	2633.46	J/molxK	1624.44	Joback Method
dvisc	0.0002373	Paxs	529.37	Joback Method

dvisc	0.0000562	Paxs	638.13	Joback Method
dvisc	0.0000203	Paxs	746.89	Joback Method
dvisc	0.0000095	Paxs	855.64	Joback Method
dvisc	0.0000052	Paxs	964.40	Joback Method
dvisc	0.0000033	Paxs	1073.16	Joback Method
dvisc	0.0000022	Paxs	1181.92	Joback Method

## Sources

<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>
<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=R505150&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R505150&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>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

Latest version available from:

<https://www.chemeo.com/cid/65-782-2/14-18-22-Trimethyltetracontane.pdf>

Generated by Cheméo on 2024-04-20 11:14:25.765900741 +0000 UTC m=+15900914.686478069.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.