

# 11,17-Dimethyl-heptacosane

<b>Inchi:</b>	InChI=1S/C29H60/c1-5-7-9-11-13-15-17-20-24-28(3)26-22-19-23-27-29(4)25-21-18-16-1
<b>InchiKey:</b>	YYKHRDACVGIVKU-UHFFFAOYSA-N
<b>Formula:</b>	C29H60
<b>SMILES:</b>	CCCCCCCCCCC(C)CCCCC(C)CCCCCCCCC
<b>Mol. weight [g/mol]:</b>	408.79

## Physical Properties

Property code	Value	Unit	Source
gf	188.42	kJ/mol	Joback Method
hf	-652.45	kJ/mol	Joback Method
hfus	63.82	kJ/mol	Joback Method
hvap	79.37	kJ/mol	Joback Method
log10ws	-11.48		Crippen Method
logp	11.271		Crippen Method
mvol	419.470	ml/mol	McGowan Method
pc	629.40	kPa	Joback Method
rinpol	2770.00		NIST Webbook
rinpol	2774.00		NIST Webbook
tb	862.04	K	Joback Method
tc	1057.60	K	Joback Method
tf	386.59	K	Joback Method
vc	1.647	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1415.45	J/molxK	862.04	Joback Method
cpg	1441.54	J/molxK	894.63	Joback Method
cpg	1466.21	J/molxK	927.23	Joback Method
cpg	1489.54	J/molxK	959.82	Joback Method
cpg	1511.59	J/molxK	992.41	Joback Method
cpg	1532.42	J/molxK	1025.01	Joback Method
cpg	1552.12	J/molxK	1057.60	Joback Method
dvisc	0.0020179	Paxs	386.59	Joback Method

dvisc	0.0005120	Paxs	465.83	Joback Method
dvisc	0.0001936	Paxs	545.07	Joback Method
dvisc	0.0000937	Paxs	624.31	Joback Method
dvisc	0.0000534	Paxs	703.56	Joback Method
dvisc	0.0000341	Paxs	782.80	Joback Method
dvisc	0.0000236	Paxs	862.04	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=R248114&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R248114&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>cp<sub>g</sub>:</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>log<sub>10</sub>ws:</b>	Log <sub>10</sub> of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>m<sub>cvol</sub>:</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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