

# 11-Methylheptacosane

<b>Other names:</b>	Heptacosane, 11-methyl
<b>Inchi:</b>	InChI=1S/C28H58/c1-4-6-8-10-12-14-15-16-17-18-19-21-23-25-27-28(3)26-24-22-20-13
<b>InchiKey:</b>	NBBBXASYEKYDOB-UHFFFAOYSA-N
<b>Formula:</b>	C28H58
<b>SMILES:</b>	CCCCCCCCCCCCCCCC(C)CCCCCCCC
<b>Mol. weight [g/mol]:</b>	394.76
<b>CAS:</b>	15689-68-6

## Physical Properties

Property code	Value	Unit	Source
gf	182.44	kJ/mol	Joback Method
hf	-626.53	kJ/mol	Joback Method
hfus	64.75	kJ/mol	Joback Method
hvap	77.53	kJ/mol	Joback Method
log10ws	-11.30		Crippen Method
logp	11.025		Crippen Method
mcvol	405.380	ml/mol	McGowan Method
pc	658.14	kPa	Joback Method
rinpol	2733.00		NIST Webbook
rinpol	2734.00		NIST Webbook
rinpol	2734.00		NIST Webbook
rinpol	2735.00		NIST Webbook
rinpol	2733.00		NIST Webbook
rinpol	2732.00		NIST Webbook
rinpol	2731.00		NIST Webbook
rinpol	2733.00		NIST Webbook
rinpol	2735.00		NIST Webbook
rinpol	2736.20		NIST Webbook
rinpol	2731.00		NIST Webbook
rinpol	2733.00		NIST Webbook
rinpol	2734.00		NIST Webbook
rinpol	2735.00		NIST Webbook
rinpol	2731.00		NIST Webbook
rinpol	2728.00		NIST Webbook
rinpol	2730.00		NIST Webbook
rinpol	2743.00		NIST Webbook
tb	839.60	K	Joback Method

tc	1029.02	K	Joback Method
tf	390.32	K	Joback Method
vc	1.597	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1348.34	J/mol×K	839.60	Joback Method
cpg	1373.75	J/mol×K	871.17	Joback Method
cpg	1397.85	J/mol×K	902.74	Joback Method
cpg	1420.68	J/mol×K	934.31	Joback Method
cpg	1442.31	J/mol×K	965.88	Joback Method
cpg	1462.81	J/mol×K	997.45	Joback Method
cpg	1482.23	J/mol×K	1029.02	Joback Method
dvisc	0.0017949	Paxs	390.32	Joback Method
dvisc	0.0005242	Paxs	465.20	Joback Method
dvisc	0.0002154	Paxs	540.08	Joback Method
dvisc	0.0001099	Paxs	614.96	Joback Method
dvisc	0.0000649	Paxs	689.84	Joback Method
dvisc	0.0000425	Paxs	764.72	Joback Method
dvisc	0.0000300	Paxs	839.60	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=C15689686&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C15689686&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>hvac:</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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