

# Cyclopentane, 1-nonyl, 3-(3-methylpentadecyl)

Inchi:	InChI=1S/C30H60/c1-4-6-8-10-12-13-14-16-17-19-21-28(3)23-24-30-26-25-29(27-30)22
InchiKey:	QIOAXSCBIPZQIO-UHFFFAOYSA-N
Formula:	C30H60
SMILES:	CCCCCCCCCCCC(C)CCC1CCC(CCCCCCCC)C1
Mol. weight [g/mol]:	420.80

## Physical Properties

Property code	Value	Unit	Source
gf	228.12	kJ/mol	Joback Method
hf	-627.67	kJ/mol	Joback Method
hfus	64.94	kJ/mol	Joback Method
hvap	81.93	kJ/mol	Joback Method
log10ws	-11.55		Crippen Method
logp	11.271		Crippen Method
mvol	422.700	ml/mol	McGowan Method
pc	643.20	kPa	Joback Method
rinpol	2945.00		NIST Webbook
rinpol	2945.00		NIST Webbook
tb	895.97	K	Joback Method
tc	1097.94	K	Joback Method
tf	419.52	K	Joback Method
vc	1.649	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1481.02	J/molxK	895.97	Joback Method
cpg	1507.38	J/molxK	929.63	Joback Method
cpg	1532.19	J/molxK	963.29	Joback Method
cpg	1555.51	J/molxK	996.95	Joback Method
cpg	1577.43	J/molxK	1030.61	Joback Method
cpg	1598.04	J/molxK	1064.27	Joback Method
cpg	1617.42	J/molxK	1097.94	Joback Method
dvisc	0.0015963	Paxs	419.52	Joback Method

dvisc	0.0005443	Paxs	498.93	Joback Method
dvisc	0.0002494	Paxs	578.34	Joback Method
dvisc	0.0001379	Paxs	657.74	Joback Method
dvisc	0.0000867	Paxs	737.15	Joback Method
dvisc	0.0000596	Paxs	816.56	Joback Method
dvisc	0.0000438	Paxs	895.97	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=R501207&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R501207&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

Latest version available from:

<https://www.chemeo.com/cid/26-082-2/Cyclopentane-1-nonyl-3-3-methylpentadecyl.pdf>

Generated by Cheméo on 2024-04-27 06:47:56.980386236 +0000 UTC m=+16489725.900963551.

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