

# Heneicosane, 7,11-dimethyl-

<b>Inchi:</b>	InChI=1S/C23H48/c1-5-7-9-11-12-13-14-16-19-23(4)21-17-20-22(3)18-15-10-8-6-2/h22-2
<b>InchiKey:</b>	CKJPPWBOKYPKSZ-UHFFFAOYSA-N
<b>Formula:</b>	C23H48
<b>SMILES:</b>	CCCCCCCCCCC(C)CCCC(C)CCCCC
<b>Mol. weight [g/mol]:</b>	324.63

## Physical Properties

Property code	Value	Unit	Source
gf	137.90	kJ/mol	Joback Method
hf	-528.61	kJ/mol	Joback Method
hfus	48.28	kJ/mol	Joback Method
hvap	66.02	kJ/mol	Joback Method
log10ws	-8.97		Crippen Method
logp	8.930		Crippen Method
mcvol	334.930	ml/mol	McGowan Method
pc	859.99	kPa	Joback Method
rinpol	2172.00		NIST Webbook
tb	724.76	K	Joback Method
tc	892.98	K	Joback Method
tf	318.97	K	Joback Method
vc	1.312	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1025.05	J/molxK	724.76	Joback Method
cpg	1047.64	J/molxK	752.80	Joback Method
cpg	1069.23	J/molxK	780.83	Joback Method
cpg	1089.83	J/molxK	808.87	Joback Method
cpg	1109.50	J/molxK	836.90	Joback Method
cpg	1128.25	J/molxK	864.94	Joback Method
cpg	1146.14	J/molxK	892.98	Joback Method
dvisc	0.0049702	Paxs	318.97	Joback Method
dvisc	0.0012262	Paxs	386.60	Joback Method

dvisc	0.0004590	Paxs	454.23	Joback Method
dvisc	0.0002216	Paxs	521.87	Joback Method
dvisc	0.0001265	Paxs	589.50	Joback Method
dvisc	0.0000810	Paxs	657.13	Joback Method
dvisc	0.0000564	Paxs	724.76	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=R195154&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R195154&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/53-379-3/Heneicosane-7-11-dimethyl.pdf>

Generated by Cheméo on 2024-04-23 14:23:23.126404638 +0000 UTC m=+16171452.046981968.

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