

5,17-Dimethylpentacosane

Inchi:	InChI=1S/C27H56/c1-5-7-9-10-16-19-24-27(4)25-21-18-15-13-11-12-14-17-20-23-26(3)2
InchiKey:	CMYHRIYCOQTQRW-UHFFFAOYSA-N
Formula:	C27H56
SMILES:	CCCCCCCC(C)CCCCCCCCCCCC(C)CCCC
Mol. weight [g/mol]:	380.73

Physical Properties

Property code	Value	Unit	Source
gf	171.58	kJ/mol	Joback Method
hf	-611.17	kJ/mol	Joback Method
hfus	58.64	kJ/mol	Joback Method
hvap	74.92	kJ/mol	Joback Method
log10ws	-10.64		Crippen Method
logp	10.491		Crippen Method
mvol	391.290	ml/mol	McGowan Method
pc	694.71	kPa	Joback Method
rinpol	2585.00		NIST Webbook
rinpol	2585.00		NIST Webbook
tb	816.28	K	Joback Method
tc	999.37	K	Joback Method
tf	364.05	K	Joback Method
vc	1.536	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1282.57	J/molxK	816.28	Joback Method
cpg	1307.23	J/molxK	846.80	Joback Method
cpg	1330.65	J/molxK	877.31	Joback Method
cpg	1352.88	J/molxK	907.83	Joback Method
cpg	1373.98	J/molxK	938.34	Joback Method
cpg	1394.00	J/molxK	968.86	Joback Method
cpg	1412.98	J/molxK	999.37	Joback Method
dvisc	0.0027479	Paxs	364.05	Joback Method

dvisc	0.0006915	Paxs	439.42	Joback Method
dvisc	0.0002606	Paxs	514.79	Joback Method
dvisc	0.0001260	Paxs	590.16	Joback Method
dvisc	0.0000719	Paxs	665.54	Joback Method
dvisc	0.0000459	Paxs	740.91	Joback Method
dvisc	0.0000319	Paxs	816.28	Joback Method

Sources

Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R505610&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

Legend

cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/24-417-2/5-17-Dimethylpentacosane.pdf>

Generated by Cheméo on 2024-04-26 08:47:40.336076313 +0000 UTC m=+16410509.256653635.

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