

Octacosane, 11-methyl

Inchi:	InChI=1S/C29H60/c1-4-6-8-10-12-14-15-16-17-18-19-20-22-24-26-28-29(3)27-25-23-21-
InchiKey:	QCOWNEUAMDSCFW-UHFFFAOYSA-N
Formula:	C29H60
SMILES:	CCCCCCCCCCCCCCCC(C)CCCCCCCC
Mol. weight [g/mol]:	408.79

Physical Properties

Property code	Value	Unit	Source
gf	190.86	kJ/mol	Joback Method
hf	-647.17	kJ/mol	Joback Method
hfus	67.34	kJ/mol	Joback Method
hvap	79.76	kJ/mol	Joback Method
log10ws	-11.72		Crippen Method
logp	11.415		Crippen Method
mvol	419.470	ml/mol	McGowan Method
pc	626.88	kPa	Joback Method
rinpol	2833.00		NIST Webbook
rinpol	2835.90		NIST Webbook
tb	862.48	K	Joback Method
tc	1058.99	K	Joback Method
tf	401.59	K	Joback Method
vc	1.653	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1415.09	J/molxK	862.48	Joback Method
cpg	1532.71	J/molxK	1026.24	Joback Method
cpg	1511.74	J/molxK	993.49	Joback Method
cpg	1489.57	J/molxK	960.74	Joback Method
cpg	1466.11	J/molxK	927.98	Joback Method
cpg	1441.31	J/molxK	895.23	Joback Method
cpg	1552.54	J/molxK	1058.99	Joback Method
dvisc	0.0000258	Paxs	862.48	Joback Method

dvisc	0.0000366	Paxs	785.66	Joback Method
dvisc	0.0000560	Paxs	708.85	Joback Method
dvisc	0.0000949	Paxs	632.04	Joback Method
dvisc	0.0001863	Paxs	555.22	Joback Method
dvisc	0.0004537	Paxs	478.40	Joback Method
dvisc	0.0015539	Paxs	401.59	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
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=R558946&Units=SI

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/34-736-7/Octacosane-11-methyl.pdf>

Generated by Cheméo on 2024-05-02 22:23:22.636190746 +0000 UTC m=+16977851.556768058.

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