

3,13-dimethyl-tritriacontane

Inchi:	InChI=1S/C35H72/c1-5-7-8-9-10-11-12-13-14-15-16-17-18-19-20-22-26-29-32-35(4)33-3
InchiKey:	RVEJECPLPDUOMZ-UHFFFAOYSA-N
Formula:	C35H72
SMILES:	CCCCCCCCCCCCCCCCCCCC(C)CCCCCCCC(C)CC
Mol. weight [g/mol]:	492.95

Physical Properties

Property code	Value	Unit	Source
gf	238.94	kJ/mol	Joback Method
hf	-776.29	kJ/mol	Joback Method
hfus	79.36	kJ/mol	Joback Method
hvap	92.73	kJ/mol	Joback Method
log10ws	-13.99		Crippen Method
logp	13.611		Crippen Method
mcvol	504.010	ml/mol	McGowan Method
pc	480.50	kPa	Joback Method
rinpol	3403.00		NIST Webbook
rinpol	3403.00		NIST Webbook
tb	999.32	K	Joback Method
tc	1260.43	K	Joback Method
tf	454.21	K	Joback Method
vc	1.984	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1825.56	J/mol×K	999.32	Joback Method
cpg	1858.31	J/mol×K	1042.84	Joback Method
cpg	1888.72	J/mol×K	1086.36	Joback Method
cpg	1916.96	J/mol×K	1129.87	Joback Method
cpg	1943.25	J/mol×K	1173.39	Joback Method
cpg	1967.79	J/mol×K	1216.91	Joback Method
cpg	1990.77	J/mol×K	1260.43	Joback Method
dvisc	0.0007702	Paxs	454.21	Joback Method

dvisc	0.0001994	Paxs	545.06	Joback Method
dvisc	0.0000760	Paxs	635.91	Joback Method
dvisc	0.0000368	Paxs	726.76	Joback Method
dvisc	0.0000210	Paxs	817.62	Joback Method
dvisc	0.0000134	Paxs	908.47	Joback Method
dvisc	0.0000093	Paxs	999.32	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R300262&Units=SI
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

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/66-808-2/3-13-dimethyl-tritriacontane.pdf>

Generated by Cheméo on 2024-04-27 19:16:13.899595804 +0000 UTC m=+16534622.820173120.

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