

Tetracosane, 2,6,10,14,18,22-hexamethyl

Inchi:

InchiKey:

Formula:

SMILES:

Mol. weight [g/mol]:

InChI=1S/C30H62/c1-9-26(4)16-11-18-28(6)20-13-22-30(8)24-14-23-29(7)21-12-19-27(5)

PEDCUYIFPGMASW-UHFFFAOYSA-N

C30H62

CCC(C)CCCC(C)CCCC(C)CCCC(C)CCCC(C)CCCC(C)C

422.81

Physical Properties

Property code	Value	Unit	Source
gf	187.08	kJ/mol	Joback Method
hf	-694.21	kJ/mol	Joback Method
hfus	52.32	kJ/mol	Joback Method
hvap	80.05	kJ/mol	Joback Method
log10ws	-10.93		Crippen Method
logp	11.084		Crippen Method
mcvol	433.560	ml/mol	McGowan Method
pc	609.66	kPa	Joback Method
rinpol	2627.00		NIST Webbook
rinpol	2627.00		NIST Webbook
rinpol	2627.00		NIST Webbook
rinpol	2627.00		NIST Webbook
rinpol	2627.00		NIST Webbook
tb	883.16	K	Joback Method
tc	1082.52	K	Joback Method
tf	337.86	K	Joback Method
vc	1.679	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1484.10	J/molxK	883.16	Joback Method
cpg	1510.47	J/molxK	916.39	Joback Method
cpg	1535.32	J/molxK	949.61	Joback Method
cpg	1558.73	J/molxK	982.84	Joback Method
cpg	1580.78	J/molxK	1016.06	Joback Method

cpg	1601.55	J/molxK	1049.29	Joback Method
cpg	1621.11	J/molxK	1082.52	Joback Method
dvisc	0.0066197	Paxs	337.86	Joback Method
dvisc	0.0008030	Paxs	428.74	Joback Method
dvisc	0.0002037	Paxs	519.63	Joback Method
dvisc	0.0000778	Paxs	610.51	Joback Method
dvisc	0.0000381	Paxs	701.39	Joback Method
dvisc	0.0000220	Paxs	792.28	Joback Method
dvisc	0.0000142	Paxs	883.16	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
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=R213846&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/48-225-9/Tetracosane-2-6-10-14-18-22-hexamethyl.pdf>

Generated by Cheméo on 2024-04-23 12:06:31.543139692 +0000 UTC m=+16163240.463717008.

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