

1-Pentacosene

Inchi:	InChI=1S/C25H50/c1-3-5-7-9-11-13-15-17-19-21-23-25-24-22-20-18-16-14-12-10-8-6-4-2
InchiKey:	BDWBGSCCECOPTTH-UHFFFAOYSA-N
Formula:	C25H50
SMILES:	C=CCCCCCCCCCCCCCCCCCCCCCC
Mol. weight [g/mol]:	350.66
CAS:	16980-85-1

Physical Properties

Property code	Value	Unit	Source
gf	247.46	kJ/mol	Joback Method
hf	-433.90	kJ/mol	Joback Method
hfus	59.23	kJ/mol	Joback Method
hvap	70.57	kJ/mol	Joback Method
log10ws	-10.14		Crippen Method
logp	9.775		Crippen Method
mcvol	358.810	ml/mol	McGowan Method
pc	782.00	kPa	Joback Method
rinpol	2492.00		NIST Webbook
rinpol	2496.00		NIST Webbook
rinpol	2488.00		NIST Webbook
rinpol	2486.00		NIST Webbook
rinpol	2496.00		NIST Webbook
rinpol	2483.70		NIST Webbook
rinpol	2492.00		NIST Webbook
rinpol	2492.00		NIST Webbook
ripol	2488.00		NIST Webbook
ripol	2488.00		NIST Webbook
tb	768.08	K	Joback Method
tc	941.27	K	Joback Method
tf	369.75	K	Joback Method
vc	1.417	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1246.59	J/mol×K	941.27	Joback Method
cpg	1228.63	J/mol×K	912.41	Joback Method
cpg	1209.78	J/mol×K	883.54	Joback Method
cpg	1190.02	J/mol×K	854.68	Joback Method
cpg	1169.29	J/mol×K	825.81	Joback Method
cpg	1147.55	J/mol×K	796.95	Joback Method
cpg	1124.75	J/mol×K	768.08	Joback Method
dvisc	0.0020492	Paxs	369.75	Joback Method
dvisc	0.0000535	Paxs	768.08	Joback Method
dvisc	0.0000737	Paxs	701.69	Joback Method
dvisc	0.0001085	Paxs	635.30	Joback Method
dvisc	0.0001749	Paxs	568.91	Joback Method
dvisc	0.0003198	Paxs	502.53	Joback Method
dvisc	0.0007028	Paxs	436.14	Joback Method
hvapt	103.70	kJ/mol	550.00	NIST Webbook

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.33819e+01
Coeff. B	-4.09824e+03
Coeff. C	-2.00116e+02
Temperature range (K), min.	513.09
Temperature range (K), max.	707.92

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=C16980851&Units=SI
The Yaws Handbook of Vapor Pressure:	https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure

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
hvac:	Enthalpy of vaporization at standard conditions
hvapt:	Enthalpy of vaporization at a given temperature
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mccvol:	McGowan's characteristic volume
pc:	Critical Pressure
pvap:	Vapor pressure
rinpol:	Non-polar retention indices
ripol:	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/72-256-8/1-Pentacosene.pdf>

Generated by Cheméo on 2024-04-26 03:36:22.22615973 +0000 UTC m=+16391831.146737043.

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