

Hexadecane, 6-nonyl

Inchi:	InChI=1S/C25H52/c1-4-7-10-12-14-16-18-21-24-25(22-19-9-6-3)23-20-17-15-13-11-8-5-
InchiKey:	ZJFLWBKTNCVPRR-UHFFFAOYSA-N
Formula:	C25H52
SMILES:	CCCCCCCCCCC(CCCCC)CCCCCCCCCC
Mol. weight [g/mol]:	352.68

Physical Properties

Property code	Value	Unit	Source
gf	157.18	kJ/mol	Joback Method
hf	-564.61	kJ/mol	Joback Method
hfus	56.98	kJ/mol	Joback Method
hvap	70.86	kJ/mol	Joback Method
log10ws	-10.05		Crippen Method
logp	9.854		Crippen Method
mcvol	363.110	ml/mol	McGowan Method
pc	767.34	kPa	Joback Method
rinpol	2357.00		NIST Webbook
rinpol	2357.00		NIST Webbook
tb	770.96	K	Joback Method
tc	944.71	K	Joback Method
tf	356.51	K	Joback Method
vc	1.429	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1151.89	J/molxK	770.96	Joback Method
cpg	1175.34	J/molxK	799.92	Joback Method
cpg	1197.69	J/molxK	828.88	Joback Method
cpg	1218.99	J/molxK	857.83	Joback Method
cpg	1239.27	J/molxK	886.79	Joback Method
cpg	1258.59	J/molxK	915.75	Joback Method
cpg	1276.97	J/molxK	944.71	Joback Method
dvisc	0.0027240	Paxs	356.51	Joback Method

dvisc	0.0007960	Paxs	425.58	Joback Method
dvisc	0.0003280	Paxs	494.66	Joback Method
dvisc	0.0001679	Paxs	563.73	Joback Method
dvisc	0.0000995	Paxs	632.81	Joback Method
dvisc	0.0000654	Paxs	701.88	Joback Method
dvisc	0.0000463	Paxs	770.96	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=R47906&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

Legend

cp_g:	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:	Log ₁₀ of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
m_{cvol}:	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/72-247-8/Hexadecane-6-nonyl.pdf>

Generated by Cheméo on 2024-04-20 12:33:23.361677309 +0000 UTC m=+15905652.282254630.

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