

4-Nonene, 5-methyl-

Other names:	5-Methyl-4-nonene
Inchi:	InChI=1S/C10H20/c1-4-6-8-10(3)9-7-5-2/h8H,4-7,9H2,1-3H3/b10-8+
InchiKey:	TZHLNYXJNGVBOJ-CSKARUKUSA-N
Formula:	C10H20
SMILES:	CCCC=C(C)CCCC
Mol. weight [g/mol]:	140.27
CAS:	15918-07-7

Physical Properties

Property code	Value	Unit	Source
gf	104.99	kJ/mol	Joback Method
hf	-142.30	kJ/mol	Joback Method
hfus	20.55	kJ/mol	Joback Method
hvap	37.89	kJ/mol	Joback Method
log10ws	-3.86		Crippen Method
logp	3.923		Crippen Method
mcvol	147.460	ml/mol	McGowan Method
pc	2222.89	kPa	Joback Method
rinpol	1002.00		NIST Webbook
rinpol	1015.00		NIST Webbook
tb	432.24	K	Joback Method
tc	605.63	K	Joback Method
tf	183.42	K	Joback Method
vc	0.577	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	300.23	J/molxK	432.24	Joback Method
cpg	315.43	J/molxK	461.14	Joback Method
cpg	329.98	J/molxK	490.04	Joback Method
cpg	343.91	J/molxK	518.94	Joback Method
cpg	357.23	J/molxK	547.83	Joback Method
cpg	369.97	J/molxK	576.73	Joback Method

cpg

382.16

J/mol×K

605.63

Joback Method

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.53844e+01
Coeff. B	-4.05361e+03
Coeff. C	-6.34580e+01
Temperature range (K), min.	331.96
Temperature range (K), max.	465.88

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=C15918077&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
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
mccvol:	McGowan's characteristic volume
pc:	Critical Pressure
pvap:	Vapor pressure
rinpolar:	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.cheméo.com/cid/76-257-3/4-Nonene-5-methyl.pdf>

Generated by Cheméo on 2024-04-18 13:53:54.60445669 +0000 UTC m=+15737683.525034003.

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