

4-Nonanol

Other names:	nonan-4-ol
Inchi:	InChI=1S/C9H20O/c1-3-5-6-8-9(10)7-4-2/h9-10H,3-8H2,1-2H3
InchiKey:	IXUOEGRSQCCHEB-UHFFFAOYSA-N
Formula:	C9H20O
SMILES:	CCCCC(O)CCC
Mol. weight [g/mol]:	144.25
CAS:	5932-79-6

Physical Properties

Property code	Value	Unit	Source
cpl	367.85	J/mol×K	Vapour pressures and heat capacity measurements on the C7 C9 secondary aliphatic alcohols
gf	-114.36	kJ/mol	Joback Method
hf	-386.60	kJ/mol	Joback Method
hfus	19.63	kJ/mol	Joback Method
hvap	71.50 ± 0.30	kJ/mol	NIST Webbook
log10ws	-2.97		Crippen Method
logp	2.728		Crippen Method
mcvol	143.540	ml/mol	McGowan Method
pc	2545.61	kPa	Joback Method
rhoc	251.00	kg/m3	NIST Webbook
rhoc	251.00 ± 5.77	kg/m3	NIST Webbook
ripol	1078.00		NIST Webbook
ripol	1467.00		NIST Webbook
ripol	1479.00		NIST Webbook
ripol	1474.00		NIST Webbook
ripol	1474.00		NIST Webbook
ripol	1479.00		NIST Webbook
tb	465.65 ± 3.00	K	NIST Webbook
tc	645.10 ± 0.30	K	NIST Webbook
tc	645.10	K	NIST Webbook
tc	645.00 ± 1.00	K	NIST Webbook
tf	237.01	K	Joback Method
vc	0.575	m3/kmol	NIST Webbook

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	334.58	J/molxK	497.06	Joback Method
cpg	347.30	J/molxK	524.07	Joback Method
cpg	359.54	J/molxK	551.09	Joback Method
cpg	371.30	J/molxK	578.10	Joback Method
cpg	382.60	J/molxK	605.11	Joback Method
cpg	393.45	J/molxK	632.13	Joback Method
cpg	403.87	J/molxK	659.14	Joback Method
dvisc	0.0878509	Paxs	237.01	Joback Method
dvisc	0.0131188	Paxs	280.35	Joback Method
dvisc	0.0032599	Paxs	323.69	Joback Method
dvisc	0.0011254	Paxs	367.03	Joback Method
dvisc	0.0004864	Paxs	410.38	Joback Method
dvisc	0.0002468	Paxs	453.72	Joback Method
dvisc	0.0001409	Paxs	497.06	Joback Method

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.64041e+01
Coeff. B	-4.69496e+03
Coeff. C	-6.72920e+01
Temperature range (K), min.	358.61
Temperature range (K), max.	490.54

Sources

NIST Webbook:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C5932796&Units=SI>

The Yaws Handbook of Vapor

<https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>

Pressure:
Crippen Method:

<http://pubs.acs.org/doi/abs/10.1021/ci9903071>

Crippen Method:

https://www.chemeo.com/doc/models/crippen_log10ws

Vapour pressures and heat capacity measurements on the C7 C9 secondary aliphatic alcohols:

<https://www.doi.org/10.1016/j.jct.2006.10.007>

Joback Method:

https://en.wikipedia.org/wiki/Joback_method

McGowan Method:

<http://link.springer.com/article/10.1007/BF02311772>

Legend

cpg:	Ideal gas heat capacity
cpl:	Liquid phase 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
pvap:	Vapor pressure
rhoc:	Critical density
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.cheméo.com/cid/19-838-1/4-Nonanol.pdf>

Generated by Cheméo on 2024-04-27 07:23:46.649746038 +0000 UTC m=+16491875.570323405.

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