

(6Z)-Nonen-1-ol

Other names:	cis-6-Nonen-1-ol 6-Nonen-1-ol, (Z)- cis-6-Nonenol (6Z)-6-Nonen-1-ol (Z)-6-Nonen-1-ol (Z)-6-Nonenol (Z)-Non-6-en-1-ol 6-Nonen-1-ol, (6Z)-
Inchi:	InChI=1S/C9H18O/c1-2-3-4-5-6-7-8-9-10/h3-4,10H,2,5-9H2,1H3/b4-3-
InchiKey:	XJHRZBIBSSVCEL-ARJAWSKDSA-N
Formula:	C9H18O
SMILES:	CCC=CCCCCO
Mol. weight [g/mol]:	142.24
CAS:	35854-86-5

Physical Properties

Property code	Value	Unit	Source
gf	-31.70	kJ/mol	Joback Method
hf	-264.10	kJ/mol	Joback Method
hfus	23.36	kJ/mol	Joback Method
hvap	52.27	kJ/mol	Joback Method
log10ws	-2.71		Crippen Method
logp	2.505		Crippen Method
mcvol	139.240	ml/mol	McGowan Method
pc	2662.52	kPa	Joback Method
ripol	1172.00		NIST Webbook
ripol	1171.00		NIST Webbook
ripol	1171.00		NIST Webbook
ripol	1171.00		NIST Webbook
ripol	1171.00		NIST Webbook
ripol	1171.00		NIST Webbook
ripol	1172.00		NIST Webbook
ripol	1690.00		NIST Webbook
ripol	1688.00		NIST Webbook
ripol	1711.00		NIST Webbook
ripol	1720.00		NIST Webbook
ripol	1696.00		NIST Webbook

ripol	1676.00		NIST Webbook
ripol	1714.00		NIST Webbook
ripol	1688.00		NIST Webbook
ripol	1714.00		NIST Webbook
ripol	1711.00		NIST Webbook
ripol	1710.00		NIST Webbook
tb	501.66	K	Joback Method
tc	667.34	K	Joback Method
tf	246.93	K	Joback Method
vc	0.538	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	317.11	J/molxK	501.66	Joback Method
cpg	329.17	J/molxK	529.27	Joback Method
cpg	340.72	J/molxK	556.89	Joback Method
cpg	351.78	J/molxK	584.50	Joback Method
cpg	362.36	J/molxK	612.11	Joback Method
cpg	372.49	J/molxK	639.73	Joback Method
cpg	382.18	J/molxK	667.34	Joback Method
dvisc	0.0432087	Paxs	246.93	Joback Method
dvisc	0.0079294	Paxs	289.38	Joback Method
dvisc	0.0022455	Paxs	331.84	Joback Method
dvisc	0.0008466	Paxs	374.30	Joback Method
dvisc	0.0003894	Paxs	416.75	Joback Method
dvisc	0.0002067	Paxs	459.21	Joback Method
dvisc	0.0001222	Paxs	501.66	Joback Method

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=C35854865&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
h vap:	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
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/63-975-0/6Z-Nonen-1-ol.pdf>

Generated by Cheméo on 2023-12-08 13:09:23.703273532 +0000 UTC m=+4330212.623850845.

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