

2-Norbornanol, 5-methyl-6-methylene, acetate

Inchi:	InChI=1S/C11H16O2/c1-6-7(2)10-4-9(6)5-11(10)13-8(3)12/h6,9-11H,2,4-5H2,1,3H3
InchiKey:	OWLBHTFZSISYRA-UHFFFAOYSA-N
Formula:	C11H16O2
SMILES:	C=C1C(C)C2CC(OC(C)=O)C1C2
Mol. weight [g/mol]:	180.24

Physical Properties

Property code	Value	Unit	Source
gf	-45.12	kJ/mol	Joback Method
hf	-332.17	kJ/mol	Joback Method
hfus	22.19	kJ/mol	Joback Method
hvap	48.78	kJ/mol	Joback Method
log10ws	-2.32		Crippen Method
logp	2.150		Crippen Method
mcvol	147.270	ml/mol	McGowan Method
pc	2525.19	kPa	Joback Method
rinpol	1227.00		NIST Webbook
rinpol	1227.00		NIST Webbook
tb	534.94	K	Joback Method
tc	739.15	K	Joback Method
tf	323.45	K	Joback Method
vc	0.564	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	371.07	J/molxK	534.94	Joback Method
cpg	448.06	J/molxK	705.12	Joback Method
cpg	434.37	J/molxK	671.08	Joback Method
cpg	419.87	J/molxK	637.05	Joback Method
cpg	404.50	J/molxK	603.01	Joback Method
cpg	388.25	J/molxK	568.98	Joback Method
cpg	460.95	J/molxK	739.15	Joback Method
dvisc	0.0009718	Paxs	534.94	Joback Method

dvisc	0.0010041	Paxs	499.69	Joback Method
dvisc	0.0010427	Paxs	464.44	Joback Method
dvisc	0.0010896	Paxs	429.20	Joback Method
dvisc	0.0011475	Paxs	393.95	Joback Method
dvisc	0.0012208	Paxs	358.70	Joback Method
dvisc	0.0013165	Paxs	323.45	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=R129805&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
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
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/86-369-8/2-Norbornanol-5-methyl-6-methylene-acetate.pdf>

Generated by Cheméo on 2024-04-20 15:59:56.663608519 +0000 UTC m=+15918045.584185830.

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