

9-Dodecen-1-ol, acetate, (Z)-

Other names:	(Z)-9-Dodecen-1-ol acetate Z-9-Dodecen-1-yl acetate Z-9-Dodecenyl acetate acetic acid dodec-9-enyl ester, cis (Z)-dodec-9-enyl acetate
Inchi:	InChI=1S/C14H26O2/c1-3-4-5-6-7-8-9-10-11-12-13-16-14(2)15/h4-5H,3,6-13H2,1-2H3/b
InchiKey:	MFFQOUCMBNXXSBK-PLNGDYQASA-N
Formula:	C14H26O2
SMILES:	CCC=CCCCCCCCOC(C)=O
Mol. weight [g/mol]:	226.35
CAS:	16974-11-1

Physical Properties

Property code	Value	Unit	Source
gf	-86.70	kJ/mol	Joback Method
hf	-459.87	kJ/mol	Joback Method
hfus	35.00	kJ/mol	Joback Method
hvap	80.70	kJ/mol	NIST Webbook
log10ws	-4.40		Crippen Method
logp	4.246		Crippen Method
mcvol	211.260	ml/mol	McGowan Method
pc	1647.09	kPa	Joback Method
rinpol	1611.00		NIST Webbook
rinpol	1611.00		NIST Webbook
ripol	1986.00		NIST Webbook
ripol	1986.00		NIST Webbook
tb	600.17	K	Joback Method
tc	774.06	K	Joback Method
tf	314.62	K	Joback Method
vc	0.824	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
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cpg	547.45	J/molxK	600.17	Joback Method
cpg	563.96	J/molxK	629.15	Joback Method
cpg	579.75	J/molxK	658.13	Joback Method
cpg	594.83	J/molxK	687.12	Joback Method
cpg	609.24	J/molxK	716.10	Joback Method
cpg	622.98	J/molxK	745.08	Joback Method
cpg	636.08	J/molxK	774.06	Joback Method
dvisc	0.0025872	Paxs	314.62	Joback Method
dvisc	0.0011349	Paxs	362.21	Joback Method
dvisc	0.0006028	Paxs	409.80	Joback Method
dvisc	0.0003653	Paxs	457.40	Joback Method
dvisc	0.0002432	Paxs	504.99	Joback Method
dvisc	0.0001737	Paxs	552.58	Joback Method
dvisc	0.0001309	Paxs	600.17	Joback Method

Sources

McGowan Method:

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

NIST Webbook:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C16974111&Units=SI>

Crippen Method:

<http://pubs.acs.org/doi/abs/10.1021/ci990307I>

Crippen Method:

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

Joback Method:

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

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
ripol:	Polar retention indices
tb:	Normal Boiling Point Temperature

tc: Critical Temperature
tf: Normal melting (fusion) point
vc: Critical Volume

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