cis-7-Dodecen-1-ol

Other names:	7-Dodecen-1-ol, (Z)-
	Looplure inhibitor
	(Z)-7-Dodecen-1-ol
	(Z)-7-Dodecenyl alcohol
	(7Z)-7-Dodecen-1-ol
	7-Dodecenol, Z-
	(Z)-dodec-7-enol
Inchi:	InChI=1S/C12H24O/c1-2-3-4-5-6-7-8-9-10-11-12-13/h5-6,13H,2-4,7-12H2,1H3/b6-5-
InchiKey:	WWDOVTHLTQFGOZ-WAYWQWQTSA-N
Formula:	C12H24O
SMILES:	CCCCC=CCCCCCCO
Mol. weight [g/mol]:	184.32
CAS:	20056-92-2

Physical Properties

Property code	Value	Unit	Source
gf	-6.44	kJ/mol	Joback Method
hf	-326.02	kJ/mol	Joback Method
hfus	31.13	kJ/mol	Joback Method
hvap	90.50	kJ/mol	NIST Webbook
log10ws	-3.96		Crippen Method
logp	3.676		Crippen Method
mcvol	181.510	ml/mol	McGowan Method
pc	2018.13	kPa	Joback Method
rinpol	1463.00		NIST Webbook
rinpol	1463.00		NIST Webbook
ripol	2009.00		NIST Webbook
ripol	2009.00		NIST Webbook
ripol	2040.00		NIST Webbook
ripol	1997.00		NIST Webbook
ripol	2040.00		NIST Webbook
tb	570.30	K	Joback Method
tc	733.56	K	Joback Method
tf	280.74	K	Joback Method
vc	0.707	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	459.51	J/mol×K	570.30	Joback Method
cpg	473.74	J/mol×K	597.51	Joback Method
cpg	487.37	J/mol×K	624.72	Joback Method
cpg	500.43	J/mol×K	651.93	Joback Method
cpg	512.93	J/mol×K	679.14	Joback Method
cpg	524.90	J/mol×K	706.35	Joback Method
cpg	536.36	J/mol×K	733.56	Joback Method
dvisc	0.0209597	Pa×s	280.74	Joback Method
dvisc	0.0040712	Pa×s	329.00	Joback Method
dvisc	0.0012026	Pa×s	377.26	Joback Method
dvisc	0.0004685	Pa×s	425.52	Joback Method
dvisc	0.0002211	Pa×s	473.78	Joback Method
dvisc	0.0001199	Pa×s	522.04	Joback Method
dvisc	0.0000721	Pa×s	570.30	Joback Method

Sources

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=C20056922&Units=SI

Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l

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

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