

3-Decen-1-ol, (Z)-

Other names:	cis-3-Decen-1-Ol 3-Dedecen-1-ol, (Z)- (3Z)-3-Decen-1-ol (Z)-3-Decen-1-ol (Z)-3-Decenol
Inchi:	InChI=1S/C10H20O/c1-2-3-4-5-6-7-8-9-10-11/h7-8,11H,2-6,9-10H2,1H3/b8-7-
InchiKey:	MTIJDFJGPCJFKE-FPLPWBNLSA-N
Formula:	C10H20O
SMILES:	CCCCCCC=CCCO
Mol. weight [g/mol]:	156.27
CAS:	10340-22-4

Physical Properties

Property code	Value	Unit	Source
gf	-23.28	kJ/mol	Joback Method
hf	-284.74	kJ/mol	Joback Method
hfus	25.95	kJ/mol	Joback Method
hvap	78.70	kJ/mol	NIST Webbook
log10ws	-3.13		Crippen Method
logp	2.895		Crippen Method
mcvol	153.330	ml/mol	McGowan Method
pc	2417.12	kPa	Joback Method
rinpol	1237.00		NIST Webbook
rinpol	1237.00		NIST Webbook
rinpol	1225.00		NIST Webbook
rinpol	1239.00		NIST Webbook
rinpol	1245.00		NIST Webbook
rinpol	1237.00		NIST Webbook
ripol	1789.00		NIST Webbook
ripol	1790.00		NIST Webbook
ripol	1783.00		NIST Webbook
ripol	1783.00		NIST Webbook
ripol	1765.00		NIST Webbook
ripol	1790.00		NIST Webbook
tb	524.54	K	Joback Method
tc	689.26	K	Joback Method
tf	258.20	K	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	362.81	J/molxK	524.54	Joback Method
cpg	375.67	J/molxK	551.99	Joback Method
cpg	387.98	J/molxK	579.45	Joback Method
cpg	399.77	J/molxK	606.90	Joback Method
cpg	411.06	J/molxK	634.36	Joback Method
cpg	421.86	J/molxK	661.81	Joback Method
cpg	432.20	J/molxK	689.26	Joback Method
dvisc	0.0338296	Paxs	258.20	Joback Method
dvisc	0.0063373	Paxs	302.59	Joback Method
dvisc	0.0018223	Paxs	346.98	Joback Method
dvisc	0.0006952	Paxs	391.37	Joback Method
dvisc	0.0003228	Paxs	435.76	Joback Method
dvisc	0.0001727	Paxs	480.15	Joback Method
dvisc	0.0001027	Paxs	524.54	Joback Method

Sources

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=C10340224&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

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
log₁₀w_s:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
m_{cvol}:	McGowan's characteristic volume
p_c:	Critical Pressure
r_{inpol}:	Non-polar retention indices
r_{ipol}:	Polar retention indices
t_b:	Normal Boiling Point Temperature
t_c:	Critical Temperature
t_f:	Normal melting (fusion) point
v_c:	Critical Volume

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