

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

Other names:	trans-3-Decen-1-ol (3E)-3-Decen-1-ol (E)-3-Decen-1-ol (E)-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-BQYQJAHWSA-N
Formula:	C10H20O
SMILES:	CCCCCCC=CCCO
Mol. weight [g/mol]:	156.27
CAS:	10339-60-3

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	78.80	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	1231.00		NIST Webbook
rinpol	1230.00		NIST Webbook
rinpol	1233.00		NIST Webbook
rinpol	1231.00		NIST Webbook
rinpol	1233.00		NIST Webbook
rinpol	1233.00		NIST Webbook
rinpol	1232.00		NIST Webbook
ripol	1765.00		NIST Webbook
ripol	1759.00		NIST Webbook
ripol	1765.00		NIST Webbook
ripol	1765.00		NIST Webbook
tb	524.54	K	Joback Method
tc	689.26	K	Joback Method
tf	258.20	K	Joback Method
vc	0.595	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	362.81	J/mol×K	524.54	Joback Method
cpg	421.86	J/mol×K	661.81	Joback Method
cpg	411.06	J/mol×K	634.36	Joback Method
cpg	399.77	J/mol×K	606.90	Joback Method
cpg	387.98	J/mol×K	579.45	Joback Method
cpg	375.67	J/mol×K	551.99	Joback Method
cpg	432.20	J/mol×K	689.26	Joback Method
dvisc	0.0001027	Pa×s	524.54	Joback Method
dvisc	0.0001727	Pa×s	480.15	Joback Method
dvisc	0.0003228	Pa×s	435.76	Joback Method
dvisc	0.0006952	Pa×s	391.37	Joback Method
dvisc	0.0018223	Pa×s	346.98	Joback Method
dvisc	0.0063373	Pa×s	302.59	Joback Method
dvisc	0.0338296	Pa×s	258.20	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=C10339603&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

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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