

1-Octene, 3,7-dimethyl-

Other names:	2,6-Dimethyl-7-octene 3,7-Dimethyl-1-octene 3,7-Dimethyloct-1-ene NSC 157589
Inchi:	InChI=1S/C10H20/c1-5-10(4)8-6-7-9(2)3/h5,9-10H,1,6-8H2,2-4H3
InchiKey:	KSXTZYRIJKDCEA-UHFFFAOYSA-N
Formula:	C10H20
SMILES:	<chem>C=CC(C)CCCC(C)C</chem>
Mol. weight [g/mol]:	140.27
CAS:	4984-01-4

Physical Properties

Property code	Value	Unit	Source
gf	116.28	kJ/mol	Joback Method
hf	-134.86	kJ/mol	Joback Method
hfus	13.33	kJ/mol	Joback Method
hvap	36.41	kJ/mol	Joback Method
log10ws	-3.38		Crippen Method
logp	3.635		Crippen Method
mcvol	147.460	ml/mol	McGowan Method
pc	2224.99	kPa	Joback Method
rinpol	963.00		NIST Webbook
rinpol	963.00		NIST Webbook
tb	424.00	K	Joback Method
tc	597.01	K	Joback Method
tf	170.70	K	Joback Method
vc	0.565	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	299.08	J/mol×K	424.00	Joback Method
cpg	314.52	J/mol×K	452.84	Joback Method
cpg	329.33	J/mol×K	481.67	Joback Method

cpg	343.52	J/mol×K	510.51	Joback Method
cpg	357.12	J/mol×K	539.34	Joback Method
cpg	370.13	J/mol×K	568.18	Joback Method
cpg	382.59	J/mol×K	597.01	Joback Method
dvisc	0.0178939	Paxs	170.70	Joback Method
dvisc	0.0041287	Paxs	212.92	Joback Method
dvisc	0.0015477	Paxs	255.13	Joback Method
dvisc	0.0007666	Paxs	297.35	Joback Method
dvisc	0.0004522	Paxs	339.57	Joback Method
dvisc	0.0002998	Paxs	381.78	Joback Method
dvisc	0.0002157	Paxs	424.00	Joback Method

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.39187e+01
Coeff. B	-3.53486e+03
Coeff. C	-5.98980e+01
Temperature range (K), min.	319.22
Temperature range (K), max.	470.58

Sources

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
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C4984014&Units=SI
The Yaws Handbook of Vapor Pressure:	https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure

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
pvap:	Vapor pressure
rinpolar:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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