

# 4-Octene, 2,3,6-trimethyl-

<b>Inchi:</b>	InChI=1S/C11H22/c1-6-10(4)7-8-11(5)9(2)3/h7-11H,6H2,1-5H3/b8-7+
<b>InchiKey:</b>	BPHPRVXZLWQTOR-BQYQJAHWSA-N
<b>Formula:</b>	C11H22
<b>SMILES:</b>	CCC(C)C=CC(C)C(C)C
<b>Mol. weight [g/mol]:</b>	154.29
<b>CAS:</b>	63830-65-9

## Physical Properties

Property code	Value	Unit	Source
gf	114.64	kJ/mol	Joback Method
hf	-168.99	kJ/mol	Joback Method
hfus	13.88	kJ/mol	Joback Method
hvap	38.87	kJ/mol	Joback Method
log10ws	-3.56		Crippen Method
logp	3.881		Crippen Method
mvol	161.550	ml/mol	McGowan Method
pc	2069.88	kPa	Joback Method
tb	453.92	K	Joback Method
tc	634.30	K	Joback Method
tf	163.65	K	Joback Method
vc	0.614	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	344.74	J/molxK	453.92	Joback Method
cpg	361.86	J/molxK	483.98	Joback Method
cpg	378.20	J/molxK	514.05	Joback Method
cpg	393.80	J/molxK	544.11	Joback Method
cpg	408.68	J/molxK	574.18	Joback Method
cpg	422.87	J/molxK	604.24	Joback Method
cpg	436.39	J/molxK	634.30	Joback Method
dvisc	0.0447839	Paxs	163.65	Joback Method
dvisc	0.0060487	Paxs	212.03	Joback Method

dvisc	0.0017189	Paxs	260.41	Joback Method
dvisc	0.0007246	Paxs	308.78	Joback Method
dvisc	0.0003859	Paxs	357.16	Joback Method
dvisc	0.0002389	Paxs	405.54	Joback Method
dvisc	0.0001638	Paxs	453.92	Joback Method

## Sources

<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C63830659&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C63830659&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>dvisc:</b>	Dynamic viscosity
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>hvap:</b>	Enthalpy of vaporization at standard conditions
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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