

# 2-Nonene, 4,6-dimethyl

<b>Inchi:</b>	InChI=1S/C11H22/c1-5-7-10(3)9-11(4)8-6-2/h5,7,10-11H,6,8-9H2,1-4H3/b7-5+
<b>InchiKey:</b>	ONGPAIHWJMXYLDFNORWQNLSA-N
<b>Formula:</b>	C11H22
<b>SMILES:</b>	CC=CC(C)CC(C)CCC
<b>Mol. weight [g/mol]:</b>	154.29

## Physical Properties

Property code	Value	Unit	Source
gf	117.08	kJ/mol	Joback Method
hf	-163.71	kJ/mol	Joback Method
hfus	17.40	kJ/mol	Joback Method
hvap	39.26	kJ/mol	Joback Method
log10ws	-3.80		Crippen Method
logp	4.025		Crippen Method
mcvol	161.550	ml/mol	McGowan Method
pc	2054.89	kPa	Joback Method
rinpol	994.00		NIST Webbook
tb	454.36	K	Joback Method
tc	630.87	K	Joback Method
tf	178.65	K	Joback Method
vc	0.620	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	344.70	J/molxK	454.36	Joback Method
cpg	420.97	J/molxK	601.45	Joback Method
cpg	407.08	J/molxK	572.03	Joback Method
cpg	392.54	J/molxK	542.62	Joback Method
cpg	377.31	J/molxK	513.20	Joback Method
cpg	361.37	J/molxK	483.78	Joback Method
cpg	434.23	J/molxK	630.87	Joback Method
dvisc	0.0001718	Paxs	454.36	Joback Method
dvisc	0.0002429	Paxs	408.41	Joback Method

dvisc	0.0003749	Paxs	362.46	Joback Method
dvisc	0.0006563	Paxs	316.50	Joback Method
dvisc	0.0013897	Paxs	270.55	Joback Method
dvisc	0.0040002	Paxs	224.60	Joback Method
dvisc	0.0198348	Paxs	178.65	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=R568304&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R568304&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>rinpol:</b>	Non-polar retention indices
<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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