

# 3,7,11-trimethyl-6E,10-dodecadienal (dihydrofarnesal)

Other names:	Dihydrofarnesal
Inchi:	InChI=1S/C15H26O/c1-13(2)7-5-8-14(3)9-6-10-15(4)11-12-16/h9,11-13H,5-8,10H2,1-4H
InchiKey:	WBNJCBPWAMSDMR-YFVJMOTDSA-N
Formula:	C15H26O
SMILES:	CC(=CC=O)CCC=C(C)CCCC(C)C
Mol. weight [g/mol]:	222.37

## Physical Properties

Property code	Value	Unit	Source
gf	116.80	kJ/mol	Joback Method
hf	-228.93	kJ/mol	Joback Method
hfus	31.16	kJ/mol	Joback Method
hvap	55.39	kJ/mol	Joback Method
log10ws	-4.85		Crippen Method
logp	4.684		Crippen Method
mcvol	215.180	ml/mol	McGowan Method
pc	1649.77	kPa	Joback Method
rinpol	1620.00		NIST Webbook
rinpol	1594.00		NIST Webbook
ripol	1961.00		NIST Webbook
tb	598.90	K	Joback Method
tc	783.10	K	Joback Method
tf	247.73	K	Joback Method
vc	0.849	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	551.05	J/molxK	598.90	Joback Method
cpg	568.45	J/molxK	629.60	Joback Method
cpg	584.98	J/molxK	660.30	Joback Method
cpg	600.68	J/molxK	691.00	Joback Method
cpg	615.60	J/molxK	721.70	Joback Method
cpg	629.78	J/molxK	752.40	Joback Method

## Sources

<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
<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=R276312&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R276312&amp;Units=SI</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<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>mccvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>rinpol:</b>	Non-polar retention indices
<b>ripol:</b>	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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