

# 2,6,10,14-Hexadecatetraen-1-ol, 3,7,11,15-tetramethyl-, acetate, (E,E,E)-

Other names:	(2E,6E,10E)-3,7,11,15-Tetramethylhexadeca-2,6,10,14-tetraen-1-yl acetate
Inchi:	InChI=1S/C22H36O2/c1-18(2)10-7-11-19(3)12-8-13-20(4)14-9-15-21(5)16-17-24-22(6)23
InchiKey:	NHYSRKKHKHCRGR-RFRQLJORSA-N
Formula:	C22H36O2
SMILES:	CC(=O)OCC=C(C)CCC=C(C)CCC=C(C)CCC=C(C)C
Mol. weight [g/mol]:	332.52
CAS:	61691-98-3

## Physical Properties

Property code	Value	Unit	Source
gf	187.12	kJ/mol	Joback Method
hf	-312.49	kJ/mol	Joback Method
hfus	51.09	kJ/mol	Joback Method
hvap	73.87	kJ/mol	Joback Method
log10ws	-7.31		Crippen Method
logp	6.695		Crippen Method
mcvol	311.080	ml/mol	McGowan Method
pc	1070.76	kPa	Joback Method
rinpol	2301.50		NIST Webbook
rinpol	2301.50		NIST Webbook
tb	795.21	K	Joback Method
tc	988.86	K	Joback Method
tf	333.70	K	Joback Method
vc	1.216	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	931.49	J/molxK	795.21	Joback Method
cpg	950.47	J/molxK	827.48	Joback Method
cpg	968.54	J/molxK	859.76	Joback Method
cpg	985.79	J/molxK	892.03	Joback Method
cpg	1002.28	J/molxK	924.31	Joback Method
cpg	1018.09	J/molxK	956.58	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=C61691983&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C61691983&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>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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