

# Geranyl heptanoate

<b>Other names:</b>	Geranyl-n-heptanoate Heptanoic acid, (2E)-3,7-dimethyl-2,6-octadien-1-yl ester [(2E)-3,7-dimethylocta-2,6-dienyl] heptanoate
<b>Inchi:</b>	InChI=1S/C17H30O2/c1-5-6-7-8-12-17(18)19-14-13-16(4)11-9-10-15(2)3/h10,13H,5-9,11H
<b>InchiKey:</b>	NSMHPPLPBQPIQJ-DTQAZKPQSA-N
<b>Formula:</b>	C17H30O2
<b>SMILES:</b>	CCCCCCC(=O)OCC=C(C)CCC=C(C)C
<b>Mol. weight [g/mol]:</b>	266.42
<b>CAS:</b>	73019-15-5

## Physical Properties

Property code	Value	Unit	Source
gf	1.68	kJ/mol	Joback Method
hf	-424.15	kJ/mol	Joback Method
hfus	40.36	kJ/mol	Joback Method
hvap	62.67	kJ/mol	Joback Method
log10ws	-5.51		Crippen Method
logp	5.193		Crippen Method
mcvol	249.230	ml/mol	McGowan Method
pc	1381.96	kPa	Joback Method
rinpol	1828.00		NIST Webbook
rinpol	1820.00		NIST Webbook
rinpol	1820.00		NIST Webbook
rinpol	1826.00		NIST Webbook
rinpol	1827.00		NIST Webbook
rinpol	1831.00		NIST Webbook
rinpol	1824.00		NIST Webbook
rinpol	1857.00		NIST Webbook
rinpol	1855.00		NIST Webbook
rinpol	1826.00		NIST Webbook
rinpol	1810.00		NIST Webbook
rinpol	1835.00		NIST Webbook
ripol	2138.00		NIST Webbook
ripol	2157.00		NIST Webbook
tb	672.73	K	Joback Method
tc	854.15	K	Joback Method
tf	315.43	K	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	688.75	J/mol×K	672.73	Joback Method
cpg	706.58	J/mol×K	702.97	Joback Method
cpg	723.56	J/mol×K	733.20	Joback Method
cpg	739.73	J/mol×K	763.44	Joback Method
cpg	755.11	J/mol×K	793.68	Joback Method
cpg	769.75	J/mol×K	823.92	Joback Method
cpg	783.68	J/mol×K	854.15	Joback Method

## Sources

<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=C73019155&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C73019155&amp;Units=SI</a>
<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>

## 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>rinpola:</b>	Non-polar retention indices
<b>ripola:</b>	Polar retention indices
<b>tb:</b>	Normal Boiling Point Temperature

**tc:** Critical Temperature  
**tf:** Normal melting (fusion) point  
**vc:** Critical Volume

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