

# Lanost-8-en-3-ol, acetate, (3«beta»,13«alpha»,14«beta»,17«alpha»)-

Other names:

Euphenyl acetate

Lanost-8-en-3-yl acetate, (3«beta»,13«alpha»,14«beta»,17«alpha»)-

24-Dihydroeuphol acetate

**Inchi:** InChI=1S/C32H54O2/c1-21(2)11-10-12-22(3)24-15-19-32(9)26-13-14-27-29(5,6)28(34-2

**InchiKey:** VARRUGKCHMYWET-RBNNMGMASA-N

**Formula:** C32H54O2

**SMILES:** CC(=O)OC1CCC2(C)C3=C(CCC2C1(C)C)C1(C)CCC(C(C)CCCC(C)C)C1(C)CC3

**Mol. weight [g/mol]:** 470.77

**CAS:** 38602-31-2

## Physical Properties

Property code	Value	Unit	Source
gf	127.87	kJ/mol	Joback Method
hf	-663.99	kJ/mol	Joback Method
hfus	34.88	kJ/mol	Joback Method
hvap	91.80	kJ/mol	Joback Method
log10ws	-9.69		Crippen Method
logp	9.130		Crippen Method
mcvol	421.440	ml/mol	McGowan Method
pc	818.20	kPa	Joback Method
rinsol	3250.00		NIST Webbook
tb	1051.35	K	Joback Method
tc	1290.99	K	Joback Method
tf	655.40	K	Joback Method
vc	1.603	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1680.25	J/molxK	1051.35	Joback Method
cpg	1731.90	J/molxK	1091.29	Joback Method
cpg	1787.46	J/molxK	1131.23	Joback Method
cpg	1847.54	J/molxK	1171.17	Joback Method
cpg	1912.77	J/molxK	1211.11	Joback Method

cpg	1983.75	J/mol×K	1251.05	Joback Method
cpg	2061.12	J/mol×K	1290.99	Joback Method

## Sources

<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=C38602312&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C38602312&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>

## 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>hvac:</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>mccol:</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

Latest version available from:

<https://www.chemeo.com/cid/45-734-7/Lanost-8-en-3-ol-acetate-3-beta-13-alpha-14-beta-17-alpha.pdf>

Generated by Cheméo on 2024-04-19 18:56:57.839201898 +0000 UTC m=+15842266.759779213.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.