

# Cholan-24-oic acid, 3,7-bis(acetyloxy)-, methyl ester, (3«alpha»,5«beta»,7«alpha»)-

Other names:

Methyl diacetylchenodesoxycholate

Chenodesoxycholic acid diacetate methyl ester

Cholane-24-oic acid, 3,7-bis-acetoxy-

12-Desoxycholic acid diacetate methyl ester

Chenodeoxycholic acid, acetate-methyl ester

Inchi: InChI=1S/C29H46O6/c1-17(7-10-26(32)33-6)22-8-9-23-27-24(12-14-29(22,23)5)28(4)13

InchiKey: ZKHVKSAMEUAGEN-JGYHGBFCSA-N

Formula: C29H46O6

SMILES: COC(=O)CCC(C)C1CCC2C3C(OC(C)=O)CC4CC(OC(C)=O)CCC4(C)C3CCC12C

Mol. weight [g/mol]: 490.67

CAS: 2616-71-9

## Physical Properties

Property code	Value	Unit	Source
gf	-377.93	kJ/mol	Joback Method
hf	-1192.39	kJ/mol	Joback Method
hfus	50.50	kJ/mol	Joback Method
hvap	103.89	kJ/mol	Joback Method
log10ws	-6.42		Crippen Method
logp	5.708		Crippen Method
mcvol	398.350	ml/mol	McGowan Method
pc	926.11	kPa	Joback Method
rinpol	3316.00		NIST Webbook
rinpol	3316.00		NIST Webbook
tb	1116.79	K	Joback Method
tc	1367.31	K	Joback Method
tf	698.83	K	Joback Method
vc	1.504	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1610.75	J/molxK	1116.79	Joback Method
cpg	1646.27	J/molxK	1158.54	Joback Method

cpg	1682.95	J/mol×K	1200.30	Joback Method
cpg	1721.15	J/mol×K	1242.05	Joback Method
cpg	1761.24	J/mol×K	1283.81	Joback Method
cpg	1803.59	J/mol×K	1325.56	Joback Method
cpg	1848.57	J/mol×K	1367.31	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=C2616719&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C2616719&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>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>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/31-106-9/Cholan-24-oic-acid-3-7-bis-acetyloxy-methyl-ester-3-alpha-5-beta-7-alpha.pdf>

Generated by Cheméo on 2024-04-23 11:18:17.564701605 +0000 UTC m=+16160346.485278921.

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