

# Campesterol

<b>Other names:</b>	Ergost-5-en-3-ol, (3«beta»,24R)- Ergost-5-en-3«beta»-ol, (24R)- (24R)-5-Ergosten-3«beta»-ol Campesterin 24«alpha»-Methylcholesterol 24«alpha»-Methyl-5-cholesten-3«beta»-ol (24R)-Methylcholest-5-en-3«beta»-ol «DELTA»5-24-Isoergosten-3«beta»-ol Ergost-5-en-3-ol-, (24R, 3«beta»)- NSC 224330 (24R)-ergost-5-en-3«beta»-ol
<b>Inchi:</b>	InChI=1S/C28H48O/c1-18(2)19(3)7-8-20(4)24-11-12-25-23-10-9-21-17-22(29)13-15-27(2)
<b>InchiKey:</b>	SGNBVLSWZMBQTH-JJZZAQJISA-N
<b>Formula:</b>	C28H48O
<b>SMILES:</b>	<chem>CC(C)C(C)CCC(C)C1CCC2C3CC=C4CC(O)CCC4(C)C3CCC12C</chem>
<b>Mol. weight [g/mol]:</b>	400.68
<b>CAS:</b>	474-62-4

## Physical Properties

Property code	Value	Unit	Source
gf	209.46	kJ/mol	Joback Method
hf	-513.15	kJ/mol	Joback Method
hfus	35.28	kJ/mol	Joback Method
hvap	91.67	kJ/mol	Joback Method
log10ws	-8.18		Crippen Method
logp	7.635		Crippen Method
mcvol	363.510	ml/mol	McGowan Method
pc	1018.78	kPa	Joback Method
rinpole	3138.00		NIST Webbook
rinpole	3116.00		NIST Webbook
rinpole	3117.00		NIST Webbook
rinpole	3135.00		NIST Webbook
rinpole	3103.00		NIST Webbook
rinpole	3099.00		NIST Webbook
rinpole	3112.00		NIST Webbook
rinpole	3165.00		NIST Webbook
rinpole	3146.00		NIST Webbook

rnpol	3131.00		NIST Webbook
rnpol	3110.00		NIST Webbook
rnpol	3165.00		NIST Webbook
rnpol	3110.00		NIST Webbook
rnpol	3105.00		NIST Webbook
rnpol	3150.00		NIST Webbook
rnpol	3114.00		NIST Webbook
rnpol	3147.00		NIST Webbook
rnpol	3131.00		NIST Webbook
rnpol	3103.00		NIST Webbook
rnpol	3131.00		NIST Webbook
rnpol	3116.00		NIST Webbook
tb	969.82	K	Joback Method
tc	1193.49	K	Joback Method
tf	523.66	K	Joback Method
vc	1.371	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1386.76	J/mol×K	969.82	Joback Method
cpg	1418.35	J/mol×K	1007.10	Joback Method
cpg	1450.56	J/mol×K	1044.38	Joback Method
cpg	1483.71	J/mol×K	1081.65	Joback Method
cpg	1518.15	J/mol×K	1118.93	Joback Method
cpg	1554.19	J/mol×K	1156.21	Joback Method
cpg	1592.18	J/mol×K	1193.49	Joback Method

## Sources

**Joback Method:**

[https://en.wikipedia.org/wiki/Joback\\_method](https://en.wikipedia.org/wiki/Joback_method)

**McGowan Method:**

<http://link.springer.com/article/10.1007/BF02311772>

**NIST Webbook:**

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C474624&Units=SI>

**Crippen Method:**

<http://pubs.acs.org/doi/abs/10.1021/ci990307l>

**Crippen Method:**

[https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)

# 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

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