

24-Methyl-31-nor-9(11)-lanostenol acetate

Inchi:	InChI=1S/C32H54O2/c1-10-21(2)11-12-22(3)24-15-19-32(9)26-13-14-27-29(5,6)28(34-2
InchiKey:	WLHVUZHSICZLPG-FOIYUBGNSA-N
Formula:	C32H54O2
SMILES:	CCC(C)CCC(C)C1CCC2(C)C3CCC4C(C)(CCC(OC(C)=O)C4(C)C)C3=CCC12C
Mol. weight [g/mol]:	470.77

Physical Properties

Property code	Value	Unit	Source
gf	129.79	kJ/mol	Joback Method
hf	-672.86	kJ/mol	Joback Method
hfus	36.34	kJ/mol	Joback Method
hvap	90.83	kJ/mol	Joback Method
log10ws	-9.45		Crippen Method
logp	8.986		Crippen Method
mcvol	421.440	ml/mol	McGowan Method
pc	809.83	kPa	Joback Method
rinsol	3369.00		NIST Webbook
tb	1041.70	K	Joback Method
tc	1279.77	K	Joback Method
tf	638.64	K	Joback Method
vc	1.601	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1677.94	J/mol×K	1041.70	Joback Method
cpg	1728.28	J/mol×K	1081.38	Joback Method
cpg	1782.26	J/mol×K	1121.06	Joback Method
cpg	1840.47	J/mol×K	1160.73	Joback Method
cpg	1903.53	J/mol×K	1200.41	Joback Method
cpg	1972.04	J/mol×K	1240.09	Joback Method
cpg	2046.59	J/mol×K	1279.77	Joback Method

Sources

Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
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=R110435&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

cpg:	Ideal gas heat capacity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvac:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mccol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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