

Androstan-17-one, 3-hydroxy-, (3«beta»,5«beta»)-

Other names:	5«beta»-Androstan-17-one, 3«beta»-hydroxy- epi-5«beta»-Androsterone Epietiocholanolone Etiocholan-3«beta»-ol-17-one 3«beta»-Etiocholanolone 3«beta»-Hydroxy-5«beta»-androstane-17-one 3«beta»-Hydroxyetiocholan-17-one 5«beta»-Epiandrosterone Androstan-3«beta»-ol-17-one, (5«beta»)- 3-Hydroxyandrostan-17-one, (3«beta»,5«beta»)- 5-beta-Androstan-17-one, 3beta-hydroxy-
Inchi:	InChI=1S/C19H30O2/c1-18-9-7-13(20)11-12(18)3-4-14-15-5-6-17(21)19(15,2)10-8-16(14)
InchiKey:	QGXBDMJGAMFCBF-LISVHTKMSA-N
Formula:	C19H30O2
SMILES:	CC12CCC3C(CCC4CC(O)CCC43C)C1CCC2=O
Mol. weight [g/mol]:	290.44
CAS:	571-31-3

Physical Properties

Property code	Value	Unit	Source
gf	-1.92	kJ/mol	Joback Method
hf	-495.56	kJ/mol	Joback Method
hfus	21.22	kJ/mol	Joback Method
hvap	76.10	kJ/mol	Joback Method
log10ws	-4.56		Crippen Method
logp	3.959		Crippen Method
mcvol	242.570	ml/mol	McGowan Method
pc	1925.36	kPa	Joback Method
tb	828.90	K	Joback Method
tc	1062.98	K	Joback Method
tf	522.17	K	Joback Method
vc	0.906	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	870.61	J/mol×K	828.90	Joback Method
cpg	895.49	J/mol×K	867.91	Joback Method
cpg	920.14	J/mol×K	906.93	Joback Method
cpg	944.87	J/mol×K	945.94	Joback Method
cpg	969.98	J/mol×K	984.95	Joback Method
cpg	995.80	J/mol×K	1023.97	Joback Method
cpg	1022.62	J/mol×K	1062.98	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
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=C571313&Units=SI

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
hvap:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.cheméo.com/cid/57-648-0/Androstan-17-one-3-hydroxy-3-beta-5-beta.pdf>

Generated by Cheméo on 2025-12-21 23:21:39.761043591 +0000 UTC m=+6107497.291084286.

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