

Norethandrolone

Other names:	19-Norpregn-4-en-3-one, 17-hydroxy-, (17«alpha»)- 19-Nor-17«alpha»-pregn-4-en-3-one, 17-hydroxy- Nileva Nilevar Pronabol SC 5914 17«alpha»-Ethyl-19-nortestosterone 17«alpha»-Ethylnortestosterone 17-Ethyl-19-nortestosterone 17-ENT 19-Nortestosterone, 17-Ethyl- 8022 C. B. Estr-4-en-3-one, 17«alpha»-ethyl-17-hydroxy- 17-«alpha»-Ethyl-17-hydroxynorandrostenone 17-«alpha»-Ethyl-17-hydroxy-4-norandrosten-3-one 17-«alpha»-Ethyl-17-hydroxy-19-norandrost-4-en-3-one 17-Hydroxy-19-norpregn-4-en-3-one 17-Hydroxy-19-nor-17-«alpha»-pregn-4-en-3-one 17-«beta»-Hydroxy-19-nor-17-«alpha»-pregn-4-en-3-one 19-Norethyltestosterone 19-Nor-17-«alpha»-ethyltestosterone 17-«alpha»-Nortestosterone Solevar 17«alpha»-Ethyl-17«beta»-hydroxyestr-4-en-3-one CB 8022 NSC 70581 U 6817 17-Ethyl-17-hydroxyestr-4-en-3-one
Inchi:	InChI=1S/C20H30O2/c1-3-20(22)11-9-18-17-6-4-13-12-14(21)5-7-15(13)16(17)8-10-19(
InchiKey:	ZDHCJEIGTNNEMY-BROHZWGRSA-N
Formula:	C20H30O2
SMILES:	CCC1(O)CCC2C3CCC4=CC(=O)CCC4C3CCC21C
Mol. weight [g/mol]:	302.45
CAS:	52-78-8

Physical Properties

Property code	Value	Unit	Source
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gf	34.54		kJ/mol	Joback Method
hf	-449.55		kJ/mol	Joback Method
hfus	23.57		kJ/mol	Joback Method
hvap	79.59		kJ/mol	Joback Method
log10ws	-5.07			Crippen Method
logp	4.269			Crippen Method
mcvol	252.360		ml/mol	McGowan Method
pc	1856.31		kPa	Joback Method
rinpol	2822.00			NIST Webbook
rinpol	2822.00			NIST Webbook
tb	860.59		K	Joback Method
tc	1093.77		K	Joback Method
tf	412.00 ± 1.00		K	NIST Webbook
vc	0.950		m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	901.36	J/mol×K	860.59	Joback Method
cpg	925.87	J/mol×K	899.45	Joback Method
cpg	950.46	J/mol×K	938.32	Joback Method
cpg	975.45	J/mol×K	977.18	Joback Method
cpg	1001.15	J/mol×K	1016.05	Joback Method
cpg	1027.88	J/mol×K	1054.91	Joback Method
cpg	1055.94	J/mol×K	1093.77	Joback Method

Sources

McGowan Method:

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

NIST Webbook:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C52788&Units=SI>

Crippen Method:

<http://pubs.acs.org/doi/abs/10.1021/ci9903071>

Crippen Method:

https://www.chemeo.com/doc/models/crippen_log10ws

Joback Method:

https://en.wikipedia.org/wiki/Joback_method

Legend

cp_g:	Ideal gas heat capacity
g_f:	Standard Gibbs free energy of formation
h_f:	Enthalpy of formation at standard conditions
h_{fus}:	Enthalpy of fusion at standard conditions
h_{vap}:	Enthalpy of vaporization at standard conditions
log₁₀w_s:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mc_{vol}:	McGowan's characteristic volume
pc:	Critical Pressure
rin_{pol}:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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