

17-«beta»-Hydroxy-7-«alpha»,17-«alpha»-dimethyl

17-TMS

InChI: InChI=1S/C24H42O2Si/c1-16-14-17-15-18(25)8-11-22(17,2)19-9-12-23(3)20(21(16)19)1
InchiKey: WOJBEUHZDDQLML-NZMYFIAWSA-N
Formula: C24H42O2Si
SMILES: CC1CC2CC(=O)CCC2(C)C2CCC3(C)C(CCC3(C)O[Si](C)(C)C)C12
Mol. weight [g/mol]: 390.67

Physical Properties

Property code	Value	Unit	Source
log10ws	-4.29		Crippen Method
logp	6.454		Crippen Method
rinpol	2757.00		NIST Webbook
rinpol	2757.00		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R257420&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.cheméo.com/doc/models/crippen_log10ws

Legend

log10ws: Log10 of Water solubility in mol/l
logp: Octanol/Water partition coefficient
rinpol: Non-polar retention indices

Latest version available from:

<https://www.cheméo.com/cid/60-425-3/17-beta-Hydroxy-7-alpha-17-alpha-dimethyl-5-beta-androstan-3-one-17-TMS>

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