

4-Androsten-17«beta»-ol-3-one (testosterone), MO-TMS

Inchi: InChI=1S/C23H39NO2Si/c1-22-13-11-17(24-25-3)15-16(22)7-8-18-19-9-10-21(26-27(4,5
InchiKey: SFXUCUOVEDBETC-FMDUFRAASA-N
Formula: C23H39NO2Si
SMILES: CON=C1C=C2CCC3C(CCC4(C)C(O[Si](C)(C)C)CCC34)C2(C)CC1
Mol. weight [g/mol]: 389.65

Physical Properties

Property code	Value	Unit	Source
log10ws	-5.26		Cheméo Relay (1.0)
logp	6.172		Crippen Method
rinpol	2645.00		NIST Webbook
tf	379.32	K	Cheméo Relay (1.0)

Sources

Cheméo Relay (1.0):

Cheméo Relay (1.0, beta):

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R523158&Units=SI>

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

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

Legend

log10ws: Log10 of Water solubility in mol/l
logp: Octanol/Water partition coefficient
rinpol: Non-polar retention indices
tf: Normal melting (fusion) point

Latest version available from:

<https://www.chemeo.com/cid/67-060-1/4-Androsten-17-beta-ol-3-one-testosterone-MO-TMS.pdf>

Generated by Cheméo on 2026-07-21 18:15:44.643149138 +0000 UTC m=+168840.485034517.

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