

# Metandienone, per-TMS

**Other names:** Methandienone, bis-TMS  
Methylboldenone, TMS  
Methandienone, di-TMS

**Inchi:** InChI=1S/C26H44O2Si2/c1-24-15-12-20(27-29(4,5)6)18-19(24)10-11-21-22(24)13-16-25

**InchiKey:** BSSYZSWTKYGGGLG-IEHATGJWSA-N

**Formula:** C26H44O2Si2

**SMILES:** CC12C=CC(O[Si](C)(C)C)=CC1=CCC1C2CCC2(C)C1CCC2(C)O[Si](C)(C)C

**Mol. weight [g/mol]:** 444.80

## Physical Properties

Property code	Value	Unit	Source
log10ws	-3.52		Crippen Method
logp	7.681		Crippen Method
rinpol	2733.00		NIST Webbook
rinpol	2780.00		NIST Webbook
rinpol	2733.00		NIST Webbook

## Sources

**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=R61661&Units=SI>

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

**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](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

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<https://www.chemeo.com/cid/48-570-6/Metandienone-per-TMS.pdf>

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