

15(S)-15-Methyl-PGB2, MO-TMS

Inchi: InChI=1S/C28H51NO4Si2/c1-10-11-16-22-28(2,33-35(7,8)9)23-21-24-19-20-26(29-31-3)
InchiKey: SETGIJFTUOUV FV-LLGZSRKNSA-N
Formula: C28H51NO4Si2
SMILES: CCCCCC(C)(C=CC1=C(CC=CCCCC(=O)O[Si](C)(C)C)C(=NOC)CC1)O[Si](C)(C)C
Mol. weight [g/mol]: 521.88

Physical Properties

Property code	Value	Unit	Source
log10ws	-4.40		Crippen Method
logp	8.321		Crippen Method
rinpol	2758.00		NIST Webbook
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Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R581347&Units=SI>

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.chemeo.com/cid/125-305-4/15-S-15-Methyl-PGB2-MO-TMS.pdf>

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