

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

Inchi: InChI=1S/C31H57NO4Si2/c1-10-12-18-24-31(3,36-38(7,8)9)25-23-27-21-22-29(32-34-26)
InchiKey: BOMRYIBKVKYZFO-QQTMMTGTSA-N
Formula: C31H57NO4Si2
SMILES: CCCCCC(C)(C=CC1=C(CC=CCCC(=O)O[Si](C)(C)C)C(=NOCCCC)CC1)O[Si](C)(C)C
Mol. weight [g/mol]: 563.96

Physical Properties

Property code	Value	Unit	Source
log10ws	-5.66		Crippen Method
logp	9.491		Crippen Method
rinpol	2964.00		NIST Webbook

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
Crippen Method: https://www.cheméo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R581327&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.cheméo.com/cid/45-692-4/15-S-15-Methyl-PGB2-BO-TMS.pdf>

Generated by Cheméo on 2024-04-29 10:14:10.643465323 +0000 UTC m=+16674899.564042638.

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