

13,14-Dihydro-15-keto-PGF2A, MO-TMS, isomer # 2

Inchi: InChI=1S/C30H61NO5Si3/c1-12-13-16-19-25(31-33-2)22-23-27-26(20-17-14-15-18-21-3
InchiKey: CGDBJQGJWDFROJ-YQVMCYEWSA-N
Formula: C30H61NO5Si3
SMILES: CCCCCC(CCC1C(O[Si](C)(C)C)CC(O[Si](C)(C)C)C1CC=CCCCC(=O)O[Si](C)(C)C)=NO
Mol. weight [g/mol]: 600.07

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.30		Crippen Method
logp	8.920		Crippen Method
rinpol	2712.00		NIST Webbook
rinpol	2712.00		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R580760&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/122-731-4/13-14-Dihydro-15-keto-PGF2A-MO-TMS-isomer-2.pdf>

Generated by Cheméo on 2024-05-01 08:38:30.268194951 +0000 UTC m=+16841959.188772262.

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