11.alpha.,15.alpha.-dihydroxy-9-ketoprost-13-enoic acid, EO-TMS # 2

Inchi: InChI=1S/C32H65NO5Si3/c1-13-15-20-24-32(3,38-41(10,11)12)25-23-28-27(21-18-16-17

InchiKey: CGOPOQNBPGACLI-XDNICMEOSA-N

Formula: C32H65NO5Si3

SMILES: CCCCC[C@@](C)(/C=C/[C@H]1[C@H](O[Si](C)(C)C)CC(=NOCC)[C@@H]1CCCCCC

Mol. weight [g/mol]: 628.13

Physical Properties

Property code	Value	Unit	Source
logp	9.700		Crippen Method

Sources

Cheméo Relay (1.0):

Cheméo Relay (1.0, beta):

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?Str2File=R385510

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

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

Legend

logp: Octanol/Water partition coefficient