

acid, EO-TMS #2
InchiKey: CGOPOQNBPGACLI-XDNICMEOSA-N

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Formula: C₃₂H₆₅NO₅Si₃

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

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/ci9903071>

Crippen Method: <https://www.chemeo.com/doc/models/crippen> log10ws

Legend

logp: Octanol/Water partition coefficient

Latest version available from:

<https://www.chemeo.com/cid/121-277-0/11-alpha-15-alpha-dihydroxy-9-ketoprost-13-enoic-acid-EO-TMS-2.pdf>

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