(E/Z)-3.beta.,29-bis-(Trimethylsilyloxy)lup-20(29)-e

#1

InChI: InChI=1S/C36H66O2Si2/c1-25(24-37-39(8,9)10)26-16-19-33(4)22-23-35(6)27(31(26)33)

InChIKey: ZDQRMDWEWXCLJX-JDPBXAALSA-N

Formula: C36H66O2Si2

SMILES: CC(=CO[Si](C)(C)C)[C@@H]1CC[C@]2(C)CC[C@]3(C)C(CCC4[C@@]5(C)CC[C@H](C

Mol. weight [g/mol]: 587.09

Physical Properties

Property code	Value	Unit	Source
logp	11.063		Crippen Method

Sources

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=R583472&Units=SI

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

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

Cheméo Relay (1.0):

Cheméo Relay (1.0, beta):

Legend

logp: Octanol/Water partition coefficient