1-Tetrahydrocannabinol, 6.alpha.-hydroxy, TBDMS

Inchi: InChI=1S/C33H58O3Si2/c1-15-16-17-18-25-22-27-29(28(23-25)35-37(11,12)30(3,4)5)26
InchiKey: DXPGKUPLISGZBQ-OOOSNSGVSA-N
Formula: C33H58O3Si2
SMILES: CCCCCc1cc2c(c(O[Si](C)(C)C(C)(C)C)c1)[C@@H]1C=C(C)CC[C@]1(O[Si](C)(C)C(C)(C
Mol. weight [g/mol]: 559.00

Physical Properties

Property code	Value	Unit	Source
logp	10.559		Crippen Method

Sources

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):
Cheméo Relay (1.0, beta):
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?Str2File=R525956
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

logp: Octanol/Water partition coefficient