

allo-Cholanic acid, 7«beta»-hydroxy, Me-DMES

Inchi: InChI=1S/C29H52O3Si/c1-8-33(6,7)32-25-19-21-11-9-10-17-28(21,3)24-16-18-29(4)22(1)
InchiKey: RKNVOTKYHVSSQK-MRYKGAJVSA-N
Formula: C29H52O3Si
SMILES: CC[Si](C)(C)OC1CC2CCCCC2(C)C2CCC3(C)C(C(C)CCC(=O)OC)CCC3C12
Mol. weight [g/mol]: 476.81

Physical Properties

Property code	Value	Unit	Source
log10ws	-5.72		Crippen Method
logp	7.845		Crippen Method
rinpol	3229.00		NIST Webbook
rinpol	3229.00		NIST Webbook
ripol	3752.00		NIST Webbook

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R533803&Units=SI>

Legend

log10ws: Log10 of Water solubility in mol/l
logp: Octanol/Water partition coefficient
rinpol: Non-polar retention indices
ripol: Polar retention indices

Latest version available from:

<https://www.chemeo.com/cid/47-614-8/allo-Cholanic-acid-7-beta-hydroxy-Me-DMES.pdf>

Generated by Cheméo on 2024-05-13 18:20:25.490536335 +0000 UTC m=+17913674.411113652.

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