

(+)-«alpha»-Tocopherol, O-tert.-butyldimethylsilyl-

Inchi: InChI=1S/C35H64O2Si/c1-25(2)17-14-18-26(3)19-15-20-27(4)21-16-23-35(11)24-22-31-
InchiKey: LCSOAOILNQNJJN-UHFFFAOYSA-N
Formula: C35H64O2Si
SMILES: Cc1c(C)c2c(c(C)c1O[Si](C)(C)C(C)(C)C)CCC(C)(CCCC(C)CCCC(C)CCCC(C)C)O2
Mol. weight [g/mol]: 544.97

Physical Properties

Property code	Value	Unit	Source
log10ws	-10.48		Crippen Method
logp	11.519		Crippen Method
rinpol	3419.00		NIST Webbook

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U374757&Units=SI>

Legend

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

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<https://www.chemeo.com/cid/67-564-2/alpha-Tocopherol-O-tert-butyldimethylsilyl.pdf>

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