

# MPPA, TBDMS

**Inchi:** InChI=1S/C16H37O4PSi2/c1-15(2,3)22(8,9)19-14(17)12-13-21(7,18)20-23(10,11)16(4,5)  
**InchiKey:** PGMGZVVHAJOMES-UHFFFAOYSA-N  
**Formula:** C16H37O4PSi2  
**SMILES:** CC(C)(C)[Si](C)(C)OC(=O)CCP(C)(=O)O[Si](C)(C)C(C)(C)C  
**Mol. weight [g/mol]:** 380.61

## Physical Properties

Property code	Value	Unit	Source
log10ws	-1.86		Crippen Method
logp	5.855		Crippen Method
rinpol	2031.00		NIST Webbook

## Sources

**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci990307l>  
**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)  
**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=R76064&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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