

# Silane, dimethyl(4-phenylphenoxy)octadecyloxy-

**Inchi:** InChI=1S/C32H52O2Si/c1-4-5-6-7-8-9-10-11-12-13-14-15-16-17-18-22-29-33-35(2,3)34-  
**InchiKey:** JQAYNEFWSPCMPP-UHFFFAOYSA-N  
**Formula:** C32H52O2Si  
**SMILES:** CCCCCCCCCCCCCCCCCO[Si](C)(C)Oc1ccc(-c2ccccc2)cc1  
**Mol. weight [g/mol]:** 496.84

## Physical Properties

Property code	Value	Unit	Source
log10ws	-9.87		Crippen Method
logp	10.712		Crippen Method
rinpol	3669.00		NIST Webbook
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## Sources

**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=U347499&Units=SI>  
**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci990307l>

## Legend

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

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