

Silane, dimethyl(4-(2-phenylprop-2-yl)phenoxy)pentadecyl

Inchi: InChI=1S/C32H52O2Si/c1-6-7-8-9-10-11-12-13-14-15-16-17-21-28-33-35(4,5)34-31-26-27
InchiKey: TVGZXTFICFCHSV-UHFFFAOYSA-N
Formula: C32H52O2Si
SMILES: CCCCCCCCCCCCCCO[Si](C)(C)Oc1ccc(C(C)(C)c2ccccc2)cc1
Mol. weight [g/mol]: 496.84

Physical Properties

Property code	Value	Unit	Source
log10ws	-8.62		Crippen Method
logp	10.201		Crippen Method
rinpol	3308.00		NIST Webbook

Sources

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U347197&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>

Legend

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

Latest version available from:

<https://www.chemeo.com/cid/56-155-8/Silane-dimethyl-4-2-phenylprop-2-yl-phenoxy-pentadecyloxy.pdf>

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