

Silane, dimethyl(dimethyl(3-methylphenoxy)silyloxy)(3-m

Inchi: InChI=1S/C18H26O3Si2/c1-15-9-7-11-17(13-15)19-22(3,4)21-23(5,6)20-18-12-8-10-16(2)
InchiKey: PRTFCHBVTUTLGL-UHFFFAOYSA-N
Formula: C18H26O3Si2
SMILES: Cc1cccc(O[Si](C)(C)O[Si](C)(C)Oc2cccc(C)c2)c1
Mol. weight [g/mol]: 346.57

Physical Properties

Property code	Value	Unit	Source
log10ws	-1.50		Crippen Method
logp	5.181		Crippen Method
rinpol	1945.00		NIST Webbook
rinpol	1945.00		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=U347342&Units=SI>

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/124-665-6/Silane-dimethyl-dimethyl-3-methylphenoxy-silyloxy-3-methylphenoxy.pdf>

Generated by Cheméo on 2024-05-01 06:18:58.799142615 +0000 UTC m=+16833587.719719937.

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