

Silane, dimethyl(1-phenylpropoxy)undecyloxy-

Inchi: InChI=1S/C22H40O2Si/c1-5-7-8-9-10-11-12-13-17-20-23-25(3,4)24-22(6-2)21-18-15-14-
InchiKey: LJQANQMLZNAHDN-UHFFFAOYSA-N
Formula: C22H40O2Si
SMILES: CCCCCCCCCCO[Si](C)(C)OC(CC)c1ccccc1
Mol. weight [g/mol]: 364.64

Physical Properties

Property code	Value	Unit	Source
log10ws	-5.40		Crippen Method
logp	7.403		Crippen Method
rinpol	2174.00		NIST Webbook
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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=U347280&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/59-471-4/Silane-dimethyl-1-phenylpropoxy-undecyloxy.pdf>

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