

Silane, diethyldecyloxy(1-phenylpropoxy)-

Inchi: InChI=1S/C23H42O2Si/c1-5-9-10-11-12-13-14-18-21-24-26(7-3,8-4)25-23(6-2)22-19-16-
InchiKey: OCRNRLGIWNZTGO-UHFFFAOYSA-N
Formula: C23H42O2Si
SMILES: CCCCCCCCCCO[Si](CC)(CC)OC(CC)c1ccccc1
Mol. weight [g/mol]: 378.66

Physical Properties

Property code	Value	Unit	Source
log10ws	-5.82		Crippen Method
logp	7.794		Crippen Method
rinpol	2253.00		NIST Webbook

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
Crippen Method: https://www.cheméo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U363271&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.cheméo.com/cid/65-663-4/Silane-diethyldecyloxy-1-phenylpropoxy.pdf>

Generated by Cheméo on 2025-12-05 15:45:15.389136827 +0000 UTC m=+4697712.919177492.

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