

Silane, diphenylheptyloxy(3-oxobut-2-yloxy)-

Inchi: InChI=1S/C23H32O3Si/c1-4-5-6-7-14-19-25-27(26-21(3)20(2)24,22-15-10-8-11-16-22)23
InchiKey: DOCYUJBGMHUPHZ-UHFFFAOYSA-N
Formula: C23H32O3Si
SMILES: CCCCCCO[Si](OC(C)C(C)=O)(c1ccccc1)c1ccccc1
Mol. weight [g/mol]: 384.58

Physical Properties

Property code	Value	Unit	Source
log10ws	-11.81		Crippen Method
logp	4.224		Crippen Method
rinpol	2412.00		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U367739&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

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/15-448-8/Silane-diphenylheptyloxy-3-oxobut-2-yloxy.pdf>

Generated by Cheméo on 2024-05-11 04:44:42.695498864 +0000 UTC m=+17691931.616076177.

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