

Inchi:	InChI=1S/C32H48F4O2Si/c1-2-3-4-5-6-7-8-9-10-11-12-13-14-15-22-27-37-39(29-23-18-
InchiKey:	GWRCPWKGHISZHN-UHFFFAOYSA-N
Formula:	C32H48F4O2Si
SMILES:	CCCCCCCCCCCCCCCCCCO[Si](OCC(F)(F)C(F)F)(c1ccccc1)c1ccccc1
Mol. weight [g/mol]:	568.80

Property code	Value	Unit	Source
log10ws	-16.82		Crippen Method
logp	9.048		Crippen Method
rinpol	3139.00		NIST Webbook
rinpol	3139.00		NIST Webbook

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=U368113&Units=SI>

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

<https://www.chemeo.com/cid/57-552-6/Silane-diphenylheptadecyloxy-2-2-3-3-tetrafluoropropoxy.pdf>

Generated by Cheméo on 2025-12-05 11:32:43.047931309 +0000 UTC m=+4682560.577971964.

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