

Pinobanksin-3-propanoate, bis-TMS

Inchi: InChI=1S/C24H32O6Si2/c1-8-20(25)28-24-22(26)21-18(27-23(24)16-12-10-9-11-13-16)1
InchiKey: SVXAIBOEPUZKDQ-BJKOFHAPSA-N
Formula: C24H32O6Si2
SMILES: CCC(=O)OC1C(=O)c2c(cc(O[Si](C)(C)C)cc2O[Si](C)(C)C)OC1c1ccccc1
Mol. weight [g/mol]: 472.68

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.47		Crippen Method
logp	5.752		Crippen Method
rinpol	2713.00		NIST Webbook

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=R55985&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/31-382-3/Pinobanksin-3-propanoate-bis-TMS.pdf>

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