

o-Synephrine, N-DTFMB-TMS

Inchi: InChI=1S/C24H31F6NO3Si2/c1-31(22(32)16-12-17(23(25,26)27)14-18(13-16)24(28,29)30)
InchiKey: QDKWRYQRJXZNDI-UHFFFAOYSA-N
Formula: C24H31F6NO3Si2
SMILES: CN(CC(O[Si](C)(C)C)c1ccccc1O[Si](C)(C)C)C(=O)c1cc(C(F)(F)F)cc(C(F)(F)F)c1
Mol. weight [g/mol]: 551.67

Physical Properties

Property code	Value	Unit	Source
log10ws	-3.72		Crippen Method
logp	7.603		Crippen Method
rinpol	2152.00		NIST Webbook
rinpol	2152.00		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R165014&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.cheméo.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.cheméo.com/cid/122-972-7/o-Synephrine-N-DTFMB-TMS.pdf>

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