Benzoic acid, 4-trimethylsilyloxy-, trifluoroacetyl anhydride

Inchi: InChI=1S/C12H13F3O4Si/c1-20(2,3)19-9-6-4-8(5-7-9)10(16)18-11(17)12(13,14)15/h4-7H

InchiKey: ZHPVZBLCRICRQV-UHFFFAOYSA-N

Formula: C12H13F3O4Si

SMILES: C[Si](C)(C)Oc1ccc(C(=O)OC(=O)C(F)(F)F)cc1

Mol. weight [g/mol]: 306.31

Physical Properties

Property code	Value	Unit	Source
log10ws	-1.58		Crippen Method
logp	3.146		Crippen Method
rinpol	1417.00		NIST Webbook

Sources

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=U375026&Units=SI

Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l

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