3,4-Dimethoxycinnamic acid, tert-butyldimethylsilyl ester

Other names: 3,4-Dimethoxycinnamic acid, tbdms derivative
Inchi: InChI=1S/C17H26O4Si/c1-17(2,3)22(6,7)21-16(18)11-9-13-8-10-14(19-4)15(12-13)20-5/
InchiKey: PJIIVOSNHDORRP-PKNBQFBNSA-N
Formula: C17H26O4Si
SMILES: COc1ccc(C=CC(=O)O[Si](C)(C)C(C)(C)C)cc1OC
Mol. weight [g/mol]: 322.47

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.38		Crippen Method
logp	4.265		Crippen Method
rinpol	2280.60		NIST Webbook
rinpol	2280.60		NIST Webbook

Sources

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=U352523&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

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