

3-Methyl-3-butenyl (E)-caffeate, bis-TMS

Other names:	3-Methyl-3-butyl (E)-caffeate, TMS
Inchi:	InChI=1S/C20H32O4Si2/c1-16(2)13-14-22-20(21)12-10-17-9-11-18(23-25(3,4)5)19(15-1
InchiKey:	JRYPMVVFJLTEOJ-ZRDIBKRKSA-N
Formula:	C20H32O4Si2
SMILES:	C=C(C)CCOC(=O)C=Cc1ccc(O[Si](C)(C)C)c(O[Si](C)(C)C)c1
Mol. weight [g/mol]:	392.64

Physical Properties

Property code	Value	Unit	Source
log10ws	-1.54		Crippen Method
logp	5.637		Crippen Method
rinpola	2347.00		NIST Webbook
rinpola	2348.00		NIST Webbook
rinpola	2347.00		NIST Webbook

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R55640&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
rinpola:	Non-polar retention indices

Latest version available from:

<https://www.chemeo.com/cid/70-911-2/3-Methyl-3-butenyl-E-caffeate-bis-TMS.pdf>

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