LEVORPHANOL TMS

Other names:	Levorphanol O-TMS
	Levorphanol TMS
	Levorphanol, tms derivative
Inchi:	InChI=1S/C20H31NOSi/c1-21-12-11-20-10-6-5-7-17(20)19(21)13-15-8-9-16(14-18(15)20
InchiKey:	SMNMCZXTILNDLC-UHFFFAOYSA-N
Formula:	C20H31NOSi
SMILES:	CN1CCC23CCCCC2C1Cc1ccc(O[Si](C)(C)C)cc13
Mol. weight [g/mol]:	329.56

Physical Properties

Property code	Value	Unit	Source
logp	4.588		Crippen Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U99023&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):	
Cheméo Relay (1.0, beta):	

Legend

logp:	Octanol/Water partition coefficient