

# L-Methionine, N-(trimethylsilyl)-, trimethylsilyl ester

**Other names:**

Methionine, N-(trimethylsilyl)-, trimethylsilyl ester, L-  
N,O-Bis-(trimethylsilyl)methionine  
Trimethylsilyl 4-(methylsulfanyl)-2-[(trimethylsilyl)amino]butanoate, (S)-  
Methionine, di-TMS  
Methionine, (N,O-TMS)  
Methionine, bis-TMS  
Met, (2TMS)  
Met, TMS  
Methionine, TMS  
L-methionine, 2tms derivative

**Inchi:**

InChI=1S/C11H27NO2SSi2/c1-15-9-8-10(12-16(2,3)4)11(13)14-17(5,6)7/h10,12H,8-9H2

**InchiKey:**

JATQZWJTOMUJCG-SNVBAGLBSA-N

**Formula:**

C11H27NO2SSi2

**SMILES:**

CSCCC(N[Si](C)(C)C)C(=O)O[Si](C)(C)C

**Mol. weight [g/mol]:**

293.57

**CAS:**

27844-10-6

## Physical Properties

Property code	Value	Unit	Source
log10ws	1.42		Crippen Method
logp	2.911		Crippen Method
rinpol	1530.00		NIST Webbook
rinpol	1527.51		NIST Webbook
rinpol	1541.58		NIST Webbook
rinpol	1513.50		NIST Webbook
rinpol	1530.00		NIST Webbook
rinpol	1531.10		NIST Webbook
rinpol	1508.20		NIST Webbook
rinpol	1534.00		NIST Webbook
rinpol	1541.58		NIST Webbook
rinpol	1516.78		NIST Webbook
rinpol	1518.48		NIST Webbook
rinpol	1516.78		NIST Webbook
ripol	1741.00		NIST Webbook

# Sources

**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci990307l>  
**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)  
**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=C27844106&Units=SI>

# Legend

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

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