

PIMARIC ACID TMS

Other names:	Pimaric acid TMS ester Pimaric acid, tms derivative Podocarp-8(14)-en-15-oic acid, 13-methyl-13-vinyl-, trimethylsilyl ester
Inchi:	InChI=1S/C23H38O2Si/c1-8-21(2)15-12-18-17(16-21)10-11-19-22(18,3)13-9-14-23(19,4)
InchiKey:	FYQMSVLENZHTHD-UHFFFAOYSA-N
Formula:	C23H38O2Si
SMILES:	C=CC1(C)C=C2CCC3C(C)(C(=O)O[Si](C)(C)C)CCCC3(C)C2CC1
Mol. weight [g/mol]:	374.64

Physical Properties

Property code

logp

Sources

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):	
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C21414460&Units=SI

Legend

logp: Octanol/Water partition coefficient

Latest version available from:

<https://www.chemeo.com/cid/46-370-0/PIMARIC-ACID-TMS.pdf>

Generated by Cheméo on 2026-07-19 18:59:37.201699308 +0000 UTC m=+36848.414996796.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.