

Propane-1,2,3-tricarboxylic acid, tris(trimethylsilyl) ester

Other names: Tricarballic acid, TMS
Inchi: InChI=1S/C15H32O6Si3/c1-22(2,3)19-13(16)10-12(15(18)21-24(7,8)9)11-14(17)20-23(4,5)
InchiKey: VEXAYJWEIRRBOE-UHFFFAOYSA-N
Formula: C15H32O6Si3
SMILES: C[Si](C)(C)OC(=O)CC(CC(=O)O[Si](C)(C)C)C(=O)O[Si](C)(C)C
Mol. weight [g/mol]: 392.67

Physical Properties

Property code	Value	Unit	Source
log10ws	3.40		Crippen Method
logp	3.517		Crippen Method
rinpol	1736.00		NIST Webbook
rinpol	1733.00		NIST Webbook
rinpol	1736.00		NIST Webbook
rinpol	1733.00		NIST Webbook

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.cheméo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U153601&Units=SI>

Legend

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

Latest version available from:

<https://www.cheméo.com/cid/121-925-0/Propane-1-2-3-tricarboxylic-acid-tris-trimethylsilyl-ester.pdf>

Generated by Cheméo on 2024-05-16 07:47:15.526383731 +0000 UTC m=+18134884.446961047.

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