

# Pyruvic acid ditbdms

<b>Other names:</b>	2-Ketopropionic acid ditbdms Pyruvic acid, DMTBS Pyruvic acid, TBDMS
<b>Inchi:</b>	InChI=1S/C15H32O3Si2/c1-12(17-19(8,9)14(2,3)4)13(16)18-20(10,11)15(5,6)7/h1H2,2-1
<b>InchiKey:</b>	IWHIWOKJVFOXFC-UHFFFAOYSA-N
<b>Formula:</b>	C15H32O3Si2
<b>SMILES:</b>	C=C(O[Si](C)(C)C(C)(C)C)C(=O)O[Si](C)(C)C(C)(C)C
<b>Mol. weight [g/mol]:</b>	316.58

## Physical Properties

Property code	Value	Unit	Source
log10ws	-0.50		Crippen Method
logp	5.070		Crippen Method
rinpol	1590.00		NIST Webbook
rinpol	1535.00		NIST Webbook

## Sources

<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U332347&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U332347&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>

## Legend

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

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<https://www.chemeo.com/cid/55-294-5/Pyruvic-acid-ditbdms.pdf>

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