

# tert-Butyldimethylsilyl pentan-2-yl phthalate

<b>Other names:</b>	1,2-Benzenedicarboxylic acid, pentan-2-yl tert-butyldimethylsilyl ester
<b>Inchi:</b>	InChI=1S/C19H30O4Si/c1-8-11-14(2)22-17(20)15-12-9-10-13-16(15)18(21)23-24(6,7)19
<b>InchiKey:</b>	ULWQXKUUWRTVLH-UHFFFAOYSA-N
<b>Formula:</b>	C19H30O4Si
<b>SMILES:</b>	CCCC(C)OC(=O)c1ccccc1C(=O)O[Si](C)(C)C(C)C(C)C
<b>Mol. weight [g/mol]:</b>	350.52

## Physical Properties

Property code	Value	Unit	Source
log10ws	-3.89		Crippen Method
logp	5.194		Crippen Method
rinpol	2079.00		NIST Webbook

## Sources

<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U373616&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U373616&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>
<b>Crippen Method:</b>	<a href="https://www.cheméo.com/doc/models/crippen_log10ws">https://www.cheméo.com/doc/models/crippen_log10ws</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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