

# Kaempferol, TMS

**Other names:** KAEMPFEROL 4TMS  
**Inchi:** InChI=1S/C27H44O6Si4/c1-34(2,3)30-20-15-13-19(14-16-20)26-27(33-37(10,11)12)25(2  
**InchiKey:** AGXLLZKIGNXIBX-UHFFFAOYSA-N  
**Formula:** C27H44O6Si4  
**SMILES:** C[Si](C)(C)Oc1ccc(C2Oc3cc(O[Si](C)(C)C)cc(O[Si](C)(C)C)c3C(=O)C2O[Si](C)(C)C)cc1  
**Mol. weight [g/mol]:** 576.98

## Physical Properties

Property code	Value	Unit	Source
log10ws	0.25		Crippen Method
logp	7.864		Crippen Method
rinpol	3101.40		NIST Webbook
rinpol	3065.00		NIST Webbook
rinpol	3101.40		NIST Webbook
rinpol	3065.00		NIST Webbook

## Sources

**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=R835&Units=SI>  
**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci9903071>

## Legend

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

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<https://www.chemeo.com/cid/86-349-0/Kaempferol-TMS.pdf>

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