

# Ipanguline D18

**Inchi:** InChI=1S/C18H29NO6/c1-5-11(2)16(21)25-12(3)18(4,23)17(22)24-10-13-6-8-19-9-7-14(1)  
**InchiKey:** QZDRHOHMWBXGER-LXQJWIEQSA-N  
**Formula:** C18H29NO6  
**SMILES:** CC=C(C)C(=O)OC(C)C(C)(O)C(=O)OCC1CCN2CCC(O)C12  
**Mol. weight [g/mol]:** 355.43

## Physical Properties

Property code	Value	Unit	Source
log10ws	-2.03		Crippen Method
logp	0.634		Crippen Method
mcvol	275.060	ml/mol	McGowan Method
rinpola	2480.00		NIST Webbook

## Sources

**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci990307l>  
**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)  
**McGowan Method:** <http://link.springer.com/article/10.1007/BF02311772>  
**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=R414232&Units=SI>

## Legend

**log10ws:** Log10 of Water solubility in mol/l  
**logp:** Octanol/Water partition coefficient  
**mcvol:** McGowan's characteristic volume  
**rinpola:** Non-polar retention indices

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