

Ipanguline D13

Inchi: InChI=1S/C18H29NO6/c1-5-11(2)16(21)24-12(3)18(4,23)17(22)25-14-7-9-19-8-6-13(10-7)
InchiKey: XHRLKHKWQGYBDW-LXQJWIEQSA-N
Formula: C18H29NO6
SMILES: CC=C(C)C(=O)OC(C)C(C)(O)C(=O)OC1CCN2CCC(CO)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
rinpol	2358.00		NIST Webbook
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Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
Crippen Method: 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=R394975&Units=SI>

Legend

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

Latest version available from:

<https://www.chemeo.com/cid/70-117-4/lpanguline-D13.pdf>

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