

7-(2-Methylbutyryl)-retronecine

Inchi: InChI=1S/C13H21NO3/c1-3-9(2)13(16)17-11-5-7-14-6-4-10(8-15)12(11)14/h4,9,11-12,15
InchiKey: NSBSOFPYQQNEGC-ZYXZCXLHSA-N
Formula: C13H21NO3
SMILES: CCC(C)C(=O)OC1CCN2CC=C(CO)C12
Mol. weight [g/mol]: 239.31

Physical Properties

Property code	Value	Unit	Source
log10ws	-1.59		Crippen Method
logp	0.951		Crippen Method
mcvol	191.300	ml/mol	McGowan Method
rinpol	1738.00		NIST Webbook

Sources

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=R299588&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>

Legend

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

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<https://www.chemeo.com/cid/41-659-5/7-2-Methylbutyryl-retronecine.pdf>

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