

# 3-(2-Methylbutyryloxy)-tropane

**Inchi:** InChI=1S/C13H23NO2/c1-4-9(2)13(15)16-12-7-10-5-6-11(8-12)14(10)3/h9-12H,4-8H2,1-3H3  
**InchiKey:** OGQXAZJUVVPCRL-WSVSKBAQSA-N  
**Formula:** C13H23NO2  
**SMILES:** CCC(C)C(=O)OC1CC2CCC(C1)N2C  
**Mol. weight [g/mol]:** 225.33

## Physical Properties

Property code	Value	Unit	Source
log10ws	-2.58		Crippen Method
logp	2.201		Crippen Method
mcvol	189.730	ml/mol	McGowan Method
rinpol	1493.00		NIST Webbook
rinpol	1493.00		NIST Webbook

## Sources

**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=R421339&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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