

# 7-tigloylheliotridine

**Inchi:** InChI=1S/C13H19NO3/c1-3-9(2)13(16)17-11-5-7-14-6-4-10(8-15)12(11)14/h3-4,11-12,15  
**InchiKey:** TYGYPIIOOQNWBU-NEIHAMAVSA-N  
**Formula:** C13H19NO3  
**SMILES:** CC=C(C)C(=O)OC1CCN2CC=C(CO)C12  
**Mol. weight [g/mol]:** 237.29

## Physical Properties

Property code	Value	Unit	Source
log10ws	-1.68		Crippen Method
logp	0.871		Crippen Method
mcvol	187.000	ml/mol	McGowan Method
rinpol	1873.00		NIST Webbook
rinpol	1872.00		NIST Webbook
rinpol	1917.00		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=R178015&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/23-589-3/7-tigloylheliotridine.pdf>

Generated by Cheméo on 2024-12-10 21:57:59.23633898 +0000 UTC m=+8448741.873308228.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.