

# N-Acetylloine

**Inchi:** InChI=1S/C10H16N2O2/c1-6(13)11(2)9-8-5-12-4-3-7(14-8)10(9)12/h7-10H,3-5H2,1-2H3  
**InchiKey:** YIZSKLHCDNIMHK-WSSOQDAQSA-N  
**Formula:** C10H16N2O2  
**SMILES:** CC(=O)N(C)C1C2CN3CCC(O2)C13  
**Mol. weight [g/mol]:** 196.25

## Physical Properties

| Property code | Value   | Unit   | Source         |
|---------------|---------|--------|----------------|
| log10ws       | -0.14   |        | Crippen Method |
| logp          | -0.311  |        | Crippen Method |
| mcvol         | 146.580 | ml/mol | McGowan Method |
| rinpol        | 1657.00 |        | NIST Webbook   |
| rinpol        | 1657.00 |        | 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=R268341&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/23-280-5/N-Acetylloine.pdf>

Generated by Cheméo on 2025-12-23 09:48:35.846741056 +0000 UTC m=+6231513.376781712.

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