

# Nefazodone-M (N-desalkyl-HO-) isomer-1 2AC

**Inchi:** InChI=1S/C14H17CIN2O3/c1-10(18)16-5-7-17(8-6-16)12-3-4-14(13(15)9-12)20-11(2)19/  
**InchiKey:** VFJMBFXTQVURJB-UHFFFAOYSA-N  
**Formula:** C14H17CIN2O3  
**SMILES:** CC(=O)Oc1ccc(N2CCN(C(C)=O)CC2)cc1Cl  
**Mol. weight [g/mol]:** 296.75

## Physical Properties

Property code	Value	Unit	Source
log10ws	-2.29		Crippen Method
logp	1.934		Crippen Method
mcvol	214.710	ml/mol	McGowan Method
rinpol	2335.00		NIST Webbook
rinpol	2335.00		NIST Webbook

## Sources

**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci9903071>  
**Crippen Method:** [https://www.cheméo.com/doc/models/crippen\\_log10ws](https://www.cheméo.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=R331132&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

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