

# Furan-2-carboxamide, N-ethyl-N-(3-methylphenyl)-

**Inchi:** InChI=1S/C14H15NO2/c1-3-15(12-7-4-6-11(2)10-12)14(16)13-8-5-9-17-13/h4-10H,3H2,1  
**InchiKey:** NPWWNLYUVPYRIS-UHFFFAOYSA-N  
**Formula:** C14H15NO2  
**SMILES:** CCN(C(=O)c1ccco1)c1cccc(C)c1  
**Mol. weight [g/mol]:** 229.27

## Physical Properties

Property code	Value	Unit	Source
log10ws	-7.97		Crippen Method
logp	3.255		Crippen Method
mcvol	182.320	ml/mol	McGowan Method
rinpol	1738.00		NIST Webbook
rinpol	1738.00		NIST Webbook

## Sources

**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci9903071>  
**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=U308203&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/117-785-1/Furan-2-carboxamide-N-ethyl-N-3-methylphenyl.pdf>

Generated by Cheméo on 2024-04-26 21:56:17.183030305 +0000 UTC m=+16457826.103607627.

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