

# Benzamide, 2-fluoro-N-tetradecyl-

**Inchi:** InChI=1S/C21H34FNO/c1-2-3-4-5-6-7-8-9-10-11-12-15-18-23-21(24)19-16-13-14-17-20(20)  
**InchiKey:** VYZGAKXLSSMCLT-UHFFFAOYSA-N  
**Formula:** C21H34FNO  
**SMILES:** CCCCCCCCCCCCCCN=C(O)c1ccccc1F  
**Mol. weight [g/mol]:** 335.50

## Physical Properties

Property code	Value	Unit	Source
hf	-527.62	kJ/mol	Joback Method
hvap	84.53	kJ/mol	Joback Method
log10ws	-7.13		Crippen Method
logp	6.831		Crippen Method
mcvol	296.310	ml/mol	McGowan Method
pc	1132.91	kPa	Joback Method
rinpol	2647.00		NIST Webbook
rinpol	2647.00		NIST Webbook
tb	879.55	K	Joback Method
tc	1078.81	K	Joback Method

## 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)  
**Joback Method:** [https://en.wikipedia.org/wiki/Joback\\_method](https://en.wikipedia.org/wiki/Joback_method)  
**McGowan Method:** <http://link.springer.com/article/10.1007/BF02311772>  
**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=U407143&Units=SI>

## Legend

**hf:** Enthalpy of formation at standard conditions  
**hvap:** Enthalpy of vaporization at standard conditions

<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>rinpol:</b>	Non-polar retention indices
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature

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