

# L-Proline, N-(2,6-difluoro-3-methylbenzoyl)-, tetradecyl ester

**Inchi:** InChI=1S/C27H41F2NO3/c1-3-4-5-6-7-8-9-10-11-12-13-14-20-33-27(32)23-16-15-19-30  
**InchiKey:** CWWACCQEOMVGJU-UHFFFAOYSA-N  
**Formula:** C27H41F2NO3  
**SMILES:** CCCCCCCCCCCCCOC(=O)C1CCCN1C(=O)c1c(F)ccc(C)c1F  
**Mol. weight [g/mol]:** 465.62

## Physical Properties

Property code	Value	Unit	Source
log10ws	-8.74		Crippen Method
logp	7.122		Crippen Method
mcvol	379.200	ml/mol	McGowan Method
rinpol	3330.00		NIST Webbook
rinpol	3330.00		NIST Webbook

## Sources

**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=U345897&Units=SI>  
**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>

## 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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