

# L-Proline, N-(2-bromobenzoyl)-, undecyl ester

**Inchi:** InChI=1S/C23H34BrNO3/c1-2-3-4-5-6-7-8-9-12-18-28-23(27)21-16-13-17-25(21)22(26)1  
**InchiKey:** WAMDMYBYGIGPFF-UHFFFAOYSA-N  
**Formula:** C23H34BrNO3  
**SMILES:** CCCCCCCCCCOC(=O)C1CCCN1C(=O)c1ccccc1Br  
**Mol. weight [g/mol]:** 452.43

## Physical Properties

Property code	Value	Unit	Source
log10ws	-7.51		Crippen Method
logp	6.128		Crippen Method
mcvol	336.800	ml/mol	McGowan Method
rinpol	3216.00		NIST Webbook
rinpol	3216.00		NIST Webbook

## Sources

**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=U346228&Units=SI>  
**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci9903071>

## 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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