

# L-Proline, N-(3-bromobenzoyl)-, pentyl ester

**Inchi:** InChI=1S/C17H22BrNO3/c1-2-3-4-11-22-17(21)15-9-6-10-19(15)16(20)13-7-5-8-14(18)1  
**InchiKey:** KMIYYWAFOLFZSG-UHFFFAOYSA-N  
**Formula:** C17H22BrNO3  
**SMILES:** CCCCCOC(=O)C1CCCN1C(=O)c1cccc(Br)c1  
**Mol. weight [g/mol]:** 368.26

## Physical Properties

Property code	Value	Unit	Source
log10ws	-5.00		Crippen Method
logp	3.787		Crippen Method
mcvol	252.260	ml/mol	McGowan Method
rinpol	2609.00		NIST Webbook
rinpol	2609.00		NIST Webbook

## Sources

**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=U346140&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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