

L-Proline, N-(2-fluorobenzoyl)-, ethyl ester

Inchi: InChI=1S/C14H16FNO3/c1-2-19-14(18)12-8-5-9-16(12)13(17)10-6-3-4-7-11(10)15/h3-4,11-13,15-18
InchiKey: UNYWSABQKCNHFR-UHFFFAOYSA-N
Formula: C14H16FNO3
SMILES: CCOC(=O)C1CCCN1C(=O)c1ccccc1F
Mol. weight [g/mol]: 265.28

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|--------|----------------|
| log10ws | -2.91 | | Crippen Method |
| logp | 1.993 | | Crippen Method |
| mcvol | 194.260 | ml/mol | McGowan Method |
| rinpol | 2011.00 | | NIST Webbook |
| rinpol | 2011.00 | | NIST Webbook |

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U346091&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.cheméo.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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