l-Proline, N-(2-fluorobenzoyl)-, methyl ester

Inchi: InChI=1S/C13H14FNO3/c1-18-13(17)11-7-4-8-15(11)12(16)9-5-2-3-6-10(9)14/h2-3,5-6,1
InchiKey: YMNDBZPDECAGSK-UHFFFAOYSA-N
Formula: C13H14FNO3
SMILES: COC(=O)C1CCCN1C(=O)c1ccccc1F
Mol. weight [g/mol]: 251.25

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.49		Crippen Method
logp	1.603		Crippen Method
mcvol	180.170	ml/mol	McGowan Method
rinpol	1863.00		NIST Webbook
rinpol	1863.00		NIST Webbook

Sources

Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method: 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=U299622&Units=SI

Legend

log10ws: Log10 of Water solubility in mol/l
logp: Octanol/Water partition coefficient
mcvol: McGowan's characteristic volume
rinpol: Non-polar retention indices

Cheméo (https://www.chemeo.com) is the biggest free database of chemical and physical data for the process industry.