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

Inchi: InChI=1S/C23H34FNO3/c1-2-3-4-5-6-7-8-9-12-18-28-23(27)21-16-13-17-25(21)22(26)19
InchiKey: MIYQGQWXCOHLSV-UHFFFAOYSA-N
Formula: C23H34FNO3
SMILES: CCCCCCCCCCCOC(=O)C1CCCN1C(=O)c1ccccc1F
Mol. weight [g/mol]: 391.52

Physical Properties

Property code	Value	Unit	Source
log10ws	-6.68		Crippen Method
logp	5.504		Crippen Method
mcvol	321.070	ml/mol	McGowan Method
rinpol	2927.00		NIST Webbook
rinpol	2927.00		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=U346101&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.