

L-Proline, N-(2-fluoro-3-trifluoromethylbenzoyl)- undecyl ester

InChI: InChI=1S/C24H33F4NO3/c1-2-3-4-5-6-7-8-9-10-17-32-23(31)20-15-12-16-29(20)22(30)1
InChIKey: FGSIXHPINYQYNW-UHFFFAOYSA-N

Formula: C₂₄H₃₃F₄NO₃

SMILES: CCCCCCCCCCOC(=O)C1CCCN1C(=O)c1cccc(C(F)(F)F)c1F

Mol. weight [g/mol]: 459.52

Physical Properties

Property code	Value	Unit	Source
log10ws	-7.78		Crippen Method
logp	6.523		Crippen Method
mcvol	340.470	ml/mol	McGowan Method
rinpol	2818.00		NIST Webbook
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Sources

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=U345999&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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