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

Inchi: InChI=1S/C16H17F4NO3/c1-2-9-24-15(23)12-7-4-8-21(12)14(22)10-5-3-6-11(13(10)17)1

InchiKey: JCMLXVHCASNOGD-UHFFFAOYSA-N

Formula: C16H17F4NO3

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

Mol. weight [g/mol]: 347.30

Physical Properties

Property code	Value	Unit	Source
log10ws	-4.43		Crippen Method
logp	3.402		Crippen Method
mcvol	227.750	ml/mol	McGowan Method
rinpol	2025.00		NIST Webbook
rinpol	2025.00		NIST Webbook

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=U345991&Units=SI

Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l

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.