L-Proline, N-(2,4,5-trifluoro-3-methoxybenzoyl)-, undecyl ester

Inchi: InChI=1S/C24H34F3NO4/c1-3-4-5-6-7-8-9-10-11-15-32-24(30)19-13-12-14-28(19)23(29)
InchiKey: OOIBLGIMBPSRMG-UHFFFAOYSA-N
Formula: C24H34F3NO4
SMILES: CCCCCCCCCCCOC(=O)C1CCCN1C(=O)c1cc(F)c(F)c(OC)c1F
Mol. weight [g/mol]: 457.53

Physical Properties

Property code	Value	Unit	Source
log10ws	-7.46		Crippen Method
logp	5.791		Crippen Method
mcvol	344.570	ml/mol	McGowan Method
rinpol	3015.00		NIST Webbook
rinpol	3015.00		NIST Webbook

Sources

McGowan Method: http://link.springer.com/article/10.1007/BF02311772
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=U346028&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

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.