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

Inchi: InChI=1S/C27H40F3NO4/c1-3-4-5-6-7-8-9-10-11-12-13-14-18-35-27(33)22-16-15-17-31(
InchiKey: SCXMJLIAWFUVGA-UHFFFAOYSA-N
Formula: C27H40F3NO4
SMILES: CCCCCCCCCCCCCCOC(=O)C1CCCN1C(=O)c1cc(F)c(F)c(OC)c1F
Mol. weight [g/mol]: 499.61

Physical Properties

Property code	Value	Unit	Source
log10ws	-8.72		Crippen Method
logp	6.961		Crippen Method
mcvol	386.840	ml/mol	McGowan Method
rinpol	3325.00		NIST Webbook
rinpol	3325.00		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=U346030&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.