L-Proline, N-(2,4-difluorobenzoyl)-, hexadecyl ester

Inchi: InChI=1S/C28H43F2NO3/c1-2-3-4-5-6-7-8-9-10-11-12-13-14-15-21-34-28(33)26-17-16-2

InchiKey: UWQPXMWDFITLIE-UHFFFAOYSA-N

Formula: C28H43F2NO3

SMILES: CCCCCCCCCCCCCCCCOC(=O)C1CCCN1C(=O)c1ccc(F)cc1F

Mol. weight [g/mol]: 479.64

Physical Properties

Property code	Value	Unit	Source
log10ws	-9.10		Crippen Method
logp	7.594		Crippen Method
mcvol	393.290	ml/mol	McGowan Method
rinpol	3381.00		NIST Webbook
rinpol	3381.00		NIST Webbook

Sources

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=U346046&Units=SI

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

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.