

# L-Proline, N-(pentafluorobenzoyl)-, tetradecyl ester

Inchi:	InChI=1S/C26H36F5NO3/c1-2-3-4-5-6-7-8-9-10-11-12-13-17-35-26(34)18-15-14-16-32(1
InchiKey:	GUZISAIESKIJLG-UHFFFAOYSA-N
Formula:	C26H36F5NO3
SMILES:	CCCCCCCCCCCCCOC(=O)C1CCCN1C(=O)c1c(F)c(F)c(F)c(F)c1F
Mol. weight [g/mol]:	505.56

## Physical Properties

Property code	Value	Unit	Source
log10ws	-9.26		Crippen Method
logp	7.231		Crippen Method
mcvol	370.420	ml/mol	McGowan Method
rinpol	3014.00		NIST Webbook

## Sources

Crippen Method:	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
Crippen Method:	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
McGowan Method:	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
NIST Webbook:	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U346303&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U346303&amp;Units=SI</a>

## 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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