l-Proline, n-heptafluorobutyryl-, decyl ester

Inchi:	InChI=1S/C19H28F7NO3/c1-2-3-4-5-6-7-8-9-13-30-15(28)14-11-10-12-27(14)16(29)17(2
InchiKey:	RBWZUOAVMQQOAD-UHFFFAOYSA-N
Formula:	C19H28F7NO3
SMILES:	CCCCCCCCCCOC(=O)C1CCCN1C(=O)C(F)(F)C(F)(F)C(F)(F)F
Mol. weight [g/mol]:	451.42

Physical Properties

Property code	Value	Unit	Source
log10ws	-6.28		Crippen Method
logp	5.494		Crippen Method
mcvol	299.090	ml/mol	McGowan Method
rinpol	2073.00		NIST Webbook

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U321104&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

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