

L-Proline, N-(2-chlorobenzoyl)-, isoheptyl ester

Inchi: InChI=1S/C18H24ClNO3/c1-13(2)7-6-12-23-18(22)16-10-5-11-20(16)17(21)14-8-3-4-9-1
InchiKey: XENZNEXYHOZGHW-UHFFFAOYSA-N
Formula: C18H24ClNO3
SMILES: CC(C)CCCOC(=O)C1CCCN1C(=O)c1ccccc1Cl
Mol. weight [g/mol]: 337.84

Physical Properties

Property code	Value	Unit	Source
log10ws	-4.70		Crippen Method
logp	3.924		Crippen Method
mcvol	261.090	ml/mol	McGowan Method
rinpol	2519.00		NIST Webbook
rinpol	2519.00		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U346053&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
Crippen Method: https://www.cheméo.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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