

Isonipectic acid, N-(2-fluoro-5-trifluoromethylbenzoyl)-, pentyl ester

InChI: InChI=1S/C19H23F4NO3/c1-2-3-4-11-27-18(26)13-7-9-24(10-8-13)17(25)15-12-14(19(20)16)3
InChIKey: JAEIVFPMHPXGPX-UHFFFAOYSA-N
Formula: C19H23F4NO3
SMILES: CCCCCOC(=O)C1CCN(C(=O)c2cc(C(F)(F)F)ccc2F)CC1
Mol. weight [g/mol]: 389.38

Physical Properties

Property code	Value	Unit	Source
log10ws	-5.33		Crippen Method
logp	4.430		Crippen Method
mcvol	270.020	ml/mol	McGowan Method
rinpole	2383.00		NIST Webbook

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.cheméo.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=U361389&Units=SI>

Legend

log10ws: Log10 of Water solubility in mol/l
logp: Octanol/Water partition coefficient
mcvol: McGowan's characteristic volume
rinpole: Non-polar retention indices

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