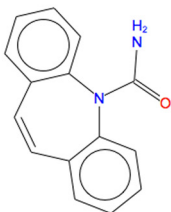


carbamazepine.mol

Decision Support



Properties

Property	Value
Target: a688239	
Description	PubChem BioAssay. qHTS Assay for Lipid Storage Modulators in Drosophil...
Interaction type	Potency
Target type	N/A
Unit	nM
Value	=1458.099975585938
Target: a688289	
Description	PubChem BioAssay. An HTS Cytotoxicity Screen to evaluate New Inhibitors...
Interaction type	CC50
Target type	N/A
Unit	uM
Value	=9.460000038146973
Target: a688658	
Description	PubChem BioAssay. A Cell Based HTS Approach for the Discovery of New I...
Interaction type	IC50
Target type	N/A
Unit	nM
Value	=2020
Target: a688867	
Description	PubChem BioAssay. Luminescence Cell-Based Dose Response HTS to Iden...
Interaction type	EC50
Target type	N/A
Unit	nM
Value	=30034
Target: a752359	
Description	PubChem BioAssay. qHTS for inhibitors of binding or entry into cells for ...
Interaction type	Potency
Target type	N/A
Unit	uM
Value	1.121999979019165
Target: a752475	
Description	PubChem BioAssay. Confirmation screen for delayed death inhibitors of t...
Interaction type	Potency
Target type	N/A

ChempSpider

Nearest neighbours

Nearest neighbours in ChEMBL

- Carbamazepine (tanimoto=1.00)
- N-methyl-5H-dibenzo[b,f]azepin-5-carboxamide (tanimoto=1.00)
- 5H-Dibenzo[b,f]azepin-5-yl(4-methylpiperazin-1-yl)methanone (tanimoto=1.00)
- N,N-diethyl-5H-dibenzo[b,f]azepin-5-carboxamide (tanimoto=0.92)

2D-Structure

Progress

Outline

