

Data Sheet

Product Name: Topiramate
Cat. No.: CS-1885
CAS No.: 97240-79-4
Molecular Formula: C12H21NO8S

Molecular Weight: 339.36

Target: Calcium Channel; Carbonic Anhydrase; GABA Receptor; iGluR;

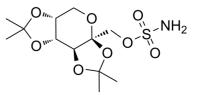
Potassium Channel; Sodium Channel

Pathway: Membrane Transporter/Ion Channel; Metabolic

Enzyme/Protease; Neuronal Signaling

Solubility: H2O: 4 mg/mL (11.79 mM; Need ultrasonic); DMSO: \geq 100

mg/mL (294.67 mM)



BIOLOGICAL ACTIVITY:

Topiramate (McN 4853) is a broad-spectrum antiepileptic agent. Topiramate is a **GluR5 receptor** antagonist. Topiramate produces its antiepileptic effects through enhancement of **GABAergic** activity, inhibition of **kainate/AMPA** receptors, inhibition of voltage-sensitive **sodium and calcium channels**, increases in **potassium** conductance, and inhibition of **carbonic anhydrase**^{[1][2][3]}. IC50 & Target: GluR5 receptor^[1]; GABAergic^[2]; Kainate/AMPA^[2]; Sodium channel^[2]; Calcium channel^[2]; Potassium channel^[2]; Carbonic anhydrase^[2] **In Vitro:** Topiramate has been believed to be a type of antiepileptic drug that blocks spread of seizures. Thus far, the mechanisms of its actions have been proven to include use-dependent inhibition of voltage-dependent Na+ channels in neurons, potentiation of GABA (γ-amino-butyric acid)-induced CI- influx, and inhibitory effects on inward currents by antagonizing kainate/alpha-amino-3-hydroxy-5-methylisoxazole-4-propionic acid (AMPA) receptors^[2].

References:

- [1]. Lyseng-Williamson KA, et al. Topiramate: a review of its use in the treatment of epilepsy. Drugs. 2007;67(15):2231-56.
- [2]. Nakamura J, et al. Target pharmacology of topiramate, a new antiepileptic drug. Nihon Yakurigaku Zasshi. 2000 Jan;115(1):53-7.
- [3]. Kaminski RM, et al. Topiramate selectively protects against seizures induced by ATPA, a GluR5 kainate receptor agonist. Neuropharmacology. 2004 Jun;46(8):1097-104.

CAIndexNames:

β-D-Fructopyranose, 2,3:4,5-bis-O-(1-methylethylidene)-, 1-sulfamate

SMILES:

NS(OC[C@]1(OC(C)(C)O2)[C@@H]2[C@H](OC(C)(C)O3)[C@H]3CO1)(=O)=O

Caution: Product has not been fully validated for medical applications. For research use only.

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