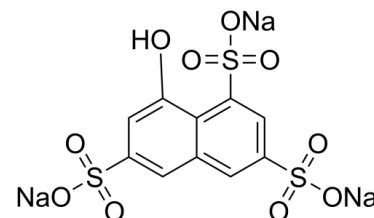


## Data Sheet

<b>Product Name:</b>	ζ-Stat (trisodium)
<b>Cat. No.:</b>	CS-0134051
<b>CAS No.:</b>	31894-34-5
<b>Molecular Formula:</b>	C <sub>10</sub> H <sub>5</sub> Na <sub>3</sub> O <sub>10</sub> S <sub>3</sub>
<b>Molecular Weight:</b>	450.31
<b>Target:</b>	Apoptosis; PKC
<b>Pathway:</b>	Apoptosis; Epigenetics; TGF-beta/Smad
<b>Solubility:</b>	H <sub>2</sub> O : 12.5 mg/mL (ultrasonic;warming;heat to 60°C)



### BIOLOGICAL ACTIVITY:

ζ-Stat trisodium (NSC37044 trisodium) is a specific and atypical **PKC-ζ** inhibitor, with an **IC<sub>50</sub>** of 5 μM. ζ-Stat trisodium can reduce melanoma cell lines proliferation and induce apoptosis, and has antitumor activity in vitro<sup>[1][2]</sup>. IC<sub>50</sub> & Target: IC<sub>50</sub>: 5 μM (PKC-ζ)<sup>[1]</sup> *In Vitro*: ζ-Stat (0.1-20 μM) shows only 13% inhibition on PKC-ι at 20 μM, but shows a significant inhibition on PKC-ζ as 51% at 5 μM level<sup>[1]</sup>.

ζ-Stat (0.1-10 μM; 3 d) significantly decreases cell proliferation of SK-MEL-2 and MeWo upon increasing the concentrations<sup>[1]</sup>.

ζ-Stat (7 or 10 μM; 24-72 h) and 5-FU in combination is able to decrease the viability of LoVo CRC cells by more than 75%<sup>[2]</sup>.

ζ-Stat (5 μM; 3 d) shows a significant diminution of phosphorylated, total PKC-ζ, Bcl-2 and PARP levels, and increases Caspase-3 and cleaved-PARP levels in SK-MEL-2 and MeWo cells<sup>[1]</sup>.

ζ-Stat (5 μM; 1-10 h) does not show significant cytotoxicity on MEL-F-NEO, SK-MEL-2 and MeWo cells<sup>[1]</sup>.

### References:

[1]. Ratnayake WS, et, al. Oncogenic PKC-ι activates Vimentin during epithelial-mesenchymal transition in melanoma; a study based on PKC-ι and PKC-ζ specific inhibitors. Cell Adh Migr. 2018; 12(5):447-463.

[2]. Islam SMA, et, al. Atypical Protein Kinase-C inhibitors exhibit a synergistic effect in facilitating DNA damaging effect of 5-fluorouracil in colorectal cancer cells. Biomed Pharmacother. 2020 Jan; 121:109665.

### CAIndexNames:

1,3,6-Naphthalenetrisulfonic acid, 8-hydroxy-, sodium salt (1:3)

### SMILES:

O=S(C1=C2C(O)=CC(S(=O)(O[Na])=O)=CC2=CC(S(=O)(O[Na])=O)=C1)(O[Na])=O

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

Tel: 610-426-3128

Fax: 888-484-5008

E-mail: sales@ChemScene.com

Address: 1 Deer Park Dr, Suite Q, Monmouth Junction, NJ 08852, USA