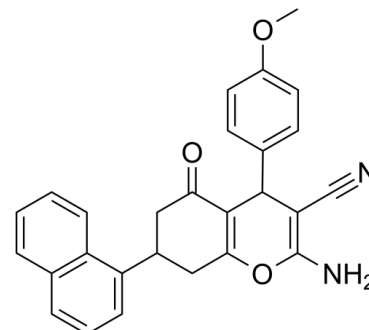


## Data Sheet

Product Name:	UCPH-101
Cat. No.:	CS-7960
CAS No.:	1118460-77-7
Molecular Formula:	C <sub>27</sub> H <sub>22</sub> N <sub>2</sub> O <sub>3</sub>
Molecular Weight:	422.48
Target:	Others
Pathway:	Others
Solubility:	DMSO : 100 mg/mL (236.70 mM; Need ultrasonic)



### BIOLOGICAL ACTIVITY:

UCPH-101 is an **excitatory amino acid transporter subtype 1 (EAAT1)** inhibitor with an  $IC_{50}$  of 0.66  $\mu$ M.  $IC_{50}$  & Target:  $IC_{50}$ : 0.66  $\mu$ M (EAAT1)<sup>[1]</sup> **In Vitro:** UCPH-101 and UCPH-102 inhibit EAAT1 anion currents in a concentration-dependent manner, with  $K_D$  values of  $0.34 \pm 0.03$   $\mu$ M (Hill=1.3 $\pm$ 0.13,  $n \geq 9$ ) for UCPH-101 and  $0.17 \pm 0.02$   $\mu$ M (Hill=0.97 $\pm$ 0.11,  $n \geq 7$ ) for UCPH-102. A small but significant decrease in the total expression levels of both HA-EAAT1 and HA-GLAST is observed in cells preincubated with 100  $\mu$ M UCPH-101 ( $p=0.048$ )<sup>[1]</sup>.

### PROTOCOL (Extracted from published papers and Only for reference)

**Cell Assay:** <sup>[1]</sup> Cells are split into poly-D-lysine-coated black 96-well plates with clear bottom. At 16 to 24 h later, the medium is aspirated, and the cells are washed with 100  $\mu$ L Krebs buffer [140 mM NaCl/4.7 mM KCl/2.5 mM CaCl<sub>2</sub>/1.2 mM MgCl<sub>2</sub>/11 mM HEPES/10 mM D-glucose, pH 7.4]; 50  $\mu$ L Krebs buffer supplemented with various concentrations of UCPH-101 or TBOA is added to the wells, after which an additional 50  $\mu$ L Krebs buffer supplemented with the FMP assay dye (1 mg/mL) is added to each well. The plate is incubated at 37°C in a humidified 5% CO<sub>2</sub> incubator for 30 min and assayed in a reader measuring emission at 560 nm caused by excitation at 530 nm before and up to 1 min after addition of 33  $\mu$ L Glu solution<sup>[1]</sup>.

### References:

[1]. Abrahamsen B, et al. Allosteric modulation of an excitatory amino acid transporter: the subtype-selective inhibitor UCPH-101 exerts sustained inhibition of EAAT1 through an intramonomeric site in the trimerization domain. J Neurosci. 2013 Jan 16;33(3):1068-87.

### CAIndexNames:

4H-1-Benzopyran-3-carbonitrile, 2-amino-5,6,7,8-tetrahydro-4-(4-methoxyphenyl)-7-(1-naphthalenyl)-5-oxo-

### SMILES:

O=C(C1)C2=C(OC(N)=C(C#N)C2C3=CC=C(OC)C=C3)CC1C4=CC=CC5=C4C=CC=C5

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

Tel: 732-484-9848 Fax: 888-484-5008 E-mail: sales@ChemScene.com

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