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# **DATASHEET**

uPSEM792 hydrochloride

#### **Product overview**

Name uPSEM792 hydrochloride

Cat No HB8542
Biological action Agonist
Purity >99%

**Description** Ultrapotent PSEM agonist for PSAM<sup>4</sup>-GlyR and PSAM<sup>4</sup>-5HT3. Brain penetrant.

### **Biological Data**

#### Biological description Overview

Ultrapotent PSEM agonist for PSAM $^4$ -GlyR and PSAM $^4$ -5HT3 ( $K_i$  values are 0.7 nM and 10,000-fold agonist selectivity for PSAM $^4$ -GlyR over  $\alpha$ -7-GlyR,  $\alpha$ 7-5HT3R and 5-HT3R.

uPSEM792 is a very weak agonist at  $\alpha 4\beta 2$  nAChR and shows 230-fold selectivity for PSAM<sup>4</sup>-GlyR over  $\alpha 4\beta 2$  nAChR.

It retains the potency of varenicline for PSAM<sup>4</sup>-GlyR with enhanced chemogenetic selectivity.

It does not act as a P-glycoprotein pump (PgP) substrate.

#### **Uses and applications**

It strongly suppresses layer 2/3 cortical neurons expressing PSAM<sup>4</sup>-GlyR in brain slices at low concentrations (ranging from 1-15 nM).

## **Solubility & Handling**

Solubility overview Storage instructions Storage of solutions

**Shipping Conditions** 

**Important** 

Soluble in water (100 mM)

-20°C

Prepare and use solutions on the same day if possible. Store solutions at -20°C for up to one month if storage is required. Equilibrate to RT and ensure the solution is precipitate free before use.

Stable for ambient temperature shipping. Follow storage instructions on receipt.

This product is for RESEARCH USE ONLY and is not intended for therapeutic or diagnostic use. Not

for human or veterinary use

#### **Chemical Data**

Chemical name Molecular Weight Chemical structure 1-Methyl-7,8,9,10-tetrahydro-1H-6,10-methanoazepino[4,5-g]quinoxalin-2(6H)-one hydrochloride 277.75

Molecular Formula CAS Number PubChem identifier

C<sub>14</sub>H<sub>15</sub>N<sub>3</sub>O · HCl 2341841-08-3 138991792 **SMILES** CN1C2=C(C=C3C4CC(C3=C2)CNC4)N=CC1=O.Cl

InChi InChi=1S/C14H15N3O.CIH/c1-17-13-4-11-9-2-8(5-15-6-9)10(11)3-12(13)16-7-14(17)18;/h3-4,7-9,15

H,2,5-6H2,1H3;1H

InChiKey CDHPEJUYEXNGCV-UHFFFAOYSA-N

Appearance Yellow solid

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scientific research use only. This product may not be used to research, develop, make, use, offer to

sell, sell, or import any products for human therapeutic uses.

#### References

Ultrapotent chemogenetics for research and potential clinical applications.

Magnus CJ *et al* (2019) Science 364(6436) **PubMedID**30872534