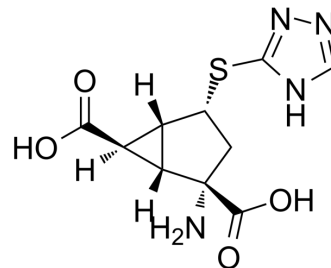


LY2812223

| | |
|---------------------------|--------------------------------------------------------------------------------------------------------|
| Cat. No.: | HY-18760 |
| CAS No.: | 1311385-20-2 |
| Molecular Formula: | C ₁₀ H ₁₂ N ₄ O ₄ S |
| Molecular Weight: | 284.29 |
| Target: | mGluR |
| Pathway: | GPCR/G Protein; Neuronal Signaling |
| Storage: | -20°C, stored under nitrogen * In solvent : -80°C, 6 months; -20°C, 1 month (stored under nitrogen) |



SOLVENT & SOLUBILITY

| | | | | | | |
|-------------------------------------------------------------------------------|-----------------------------------------------------------------------------------------------------------|----------------------|-------------|-------------|-------------|--------------|
| In Vitro | H ₂ O : 50 mg/mL (175.88 mM; Need ultrasonic) | | | | | |
| | DMSO : 28.75 mg/mL (101.13 mM; Need ultrasonic) | | | | | |
| | Preparing Stock Solutions | Solvent | Mass | 1 mg | 5 mg | 10 mg |
| | | Concentration | | | | |
| | | 1 mM | | 3.5175 mL | 17.5877 mL | 35.1753 mL |
| 5 mM | | | 0.7035 mL | 3.5175 mL | 7.0351 mL | |
| | 10 mM | | 0.3518 mL | 1.7588 mL | 3.5175 mL | |
| Please refer to the solubility information to select the appropriate solvent. | | | | | | |
| In Vivo | 1. Add each solvent one by one: PBS Solubility: 140 mg/mL (492.45 mM); Clear solution; Need ultrasonic | | | | | |

BIOLOGICAL ACTIVITY

| | | |
|-------------------------------------|------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|-----------------------|
| Description | LY2812223 is a highly potent, functionally selective mGlu ₂ receptor agonist with mGlu ₂ binding affinity for mGlu ₂ and mGlu ₃ (K _i =144 nM and 156 nM, respectively) ^[1] . | |
| IC₅₀ & Target | mGluR2 144 nM (Ki) | mGluR3 156 nM (Ki) |

REFERENCES

[1]. Monn JA, et al. Synthesis and Pharmacological Characterization of C4-(Thiotriazolyl)-substituted-2-aminobicyclo[3.1.0]hexane-2,6-dicarboxylates. Identification of (1R,2S,4R,5R,6R)-2-Amino-4-(1H-1,2,4-triazol-3-ylsulfanyl)bicyclo[3.1.0]hexane-2,6-dicarboxy

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

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