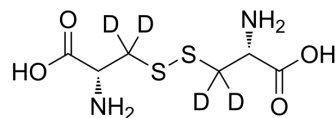


L-Cystine-d₄

| | |
|---------------------------|--|
| Cat. No.: | HY-N0394S1 |
| CAS No.: | 1192736-38-1 |
| Molecular Formula: | C ₆ H ₈ D ₄ N ₂ O ₄ S ₂ |
| Molecular Weight: | 244.33 |
| Target: | Ferroptosis; Endogenous Metabolite; Isotope-Labeled Compounds |
| Pathway: | Apoptosis; Metabolic Enzyme/Protease; Others |
| Storage: | 4°C, sealed storage, away from moisture and light * In solvent : -80°C, 6 months; -20°C, 1 month (sealed storage, away from moisture and light) |



SOLVENT & SOLUBILITY

In Vitro

H₂O : 4.55 mg/mL (18.62 mM); ultrasonic and warming and adjust pH to 1 with HCl and heat to 60°C)

| Concentration | Mass | | |
|---------------|-----------|------------|------------|
| | 1 mg | 5 mg | 10 mg |
| 1 mM | 4.0928 mL | 20.4641 mL | 40.9283 mL |
| 5 mM | 0.8186 mL | 4.0928 mL | 8.1857 mL |
| 10 mM | 0.4093 mL | 2.0464 mL | 4.0928 mL |

Please refer to the solubility information to select the appropriate solvent.

BIOLOGICAL ACTIVITY

Description

L-Cystine-d₄ is the deuterium labeled L-Cystine. L-Cystine is an amino acid and intracellular thiol, which plays a critical role in the regulation of cellular processes.

In Vitro

Stable heavy isotopes of hydrogen, carbon, and other elements have been incorporated into drug molecules, largely as tracers for quantitation during the drug development process. Deuteration has gained attention because of its potential to affect the pharmacokinetic and metabolic profiles of drugs^[1].
MCE has not independently confirmed the accuracy of these methods. They are for reference only.

REFERENCES

[1]. Russak EM, et al. Impact of Deuterium Substitution on the Pharmacokinetics of Pharmaceuticals. Ann Pharmacother. 2019;53(2):211-216.

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

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