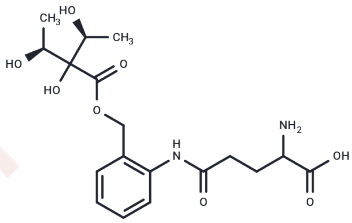
# Data Sheet (Cat.No.T130291)

TargetMol

2-amino-4-{[2-({[(3S)-dihydroxy-2-[hydroxyethyl]butanoyl]oxy}methyl)phenyl]carbamoyl} butanoic acid

## Chemical Properties



CAS No. :

Formula: C18H26N2O8

Molecular Weight: 398.41

Storage: Powder: -20°C for 3 years | In solvent: -80°C for 1 year
Actual storage temperature shall be subject to the COA.

## Biological Description

| | |
|---|---|
| Description | 2-amino-4-{[2-({[(3S)-2,3-dihydroxy-2-[(1S)-1-hydroxyethyl]butanoyl]oxy}methyl) phenyl]carbamoyl}butanoic acid is a high-purity biochemical reagent suitable for biochemical experiments and drug synthesis research. |

Preparing Stock Solutions

| | 1mg | 5mg | 10mg |
|---|---|---|---|
| 1 mM | 2.510 mL | 12.5499 mL | 25.0998 mL |
| 5 mM | 0.502 mL | 2.510 mL | 5.020 mL |
| 10 mM | 0.251 mL | 1.255 mL | 2.510 mL |
| 50 mM | 0.0502 mL | 0.251 mL | 0.502 mL |

Please select the appropriate solvent to prepare the stock solution, according to the solubility of the product in different solvents. Please use it as soon as possible.

Note: The dilution table applies only to solid products. For liquid products, please calculate the stock solution based on the stated concentration and/or density.



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