Proteins

Screening Libraries

Product Data Sheet

Mal-NH-ethyl-SS-propionic acid

Cat. No.: HY-140120 CAS No.: 2128735-24-8 Molecular Formula: $C_{12}H_{16}N_2O_5S_2$

Molecular Weight: 332.4 **ADC Linker**

Pathway: Antibody-drug Conjugate/ADC Related

Storage: -20°C, protect from light, stored under nitrogen

* In solvent: -80°C, 6 months; -20°C, 1 month (protect from light, stored under

nitrogen)

SOLVENT & SOLUBILITY

| Vitro | |
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Target:

DMSO: 100 mg/mL (300.84 mM; Need ultrasonic)

| Preparing Stock Solutions | Solvent Mass Concentration | 1 mg | 5 mg | 10 mg |
|------------------------------|-------------------------------|-----------|------------|------------|
| | 1 mM | 3.0084 mL | 15.0421 mL | 30.0842 mL |
| | 5 mM | 0.6017 mL | 3.0084 mL | 6.0168 mL |
| | 10 mM | 0.3008 mL | 1.5042 mL | 3.0084 mL |

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

In Vivo

- 1. Add each solvent one by one: 10% DMSO >> 40% PEG300 >> 5% Tween-80 >> 45% saline Solubility: ≥ 2.5 mg/mL (7.52 mM); Clear solution
- 2. Add each solvent one by one: 10% DMSO >> 90% (20% SBE-β-CD in saline) Solubility: ≥ 2.5 mg/mL (7.52 mM); Clear solution
- 3. Add each solvent one by one: 10% DMSO >> 90% corn oil Solubility: ≥ 2.5 mg/mL (7.52 mM); Clear solution

BIOLOGICAL ACTIVITY

| Description | ${\it Mal-NH-ethyl-SS-propionic\ acid\ is\ a\ cleavable\ ADC\ linker\ used\ in\ the\ synthesis\ of\ antibody-drug\ conjugates\ (ADCs)^{[1]}.}$ | | |
|---------------------------|---|-----------|--|
| IC ₅₀ & Target | Disulfide Cleavable | Cleavable | |
| In Vitro | ADCs are comprised of an antibody to which is attached an ADC cytotoxin through an ADC linker ^[1] . MCE has not independently confirmed the accuracy of these methods. They are for reference only. | | |

| REFERENCES | | | |
|--|------------------------------------|--|----|
| [1]. Beck A, et al. Strategies and challenges for the next | generation of antibody-drug conjug | ates. Nat Rev Drug Discov. 2017 May;16(5):315-33 | 7. |
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| Caution: Product ha | as not been fully validated for m | edical applications. For research use only. | |
| Tel: 609-228-6898 | Fax: 609-228-5909 | E-mail: tech@MedChemExpress.com | |
| Addres | ss: 1 Deer Park Dr, Suite Q, Monm | outh Junction, NJ 08852, USA | |
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