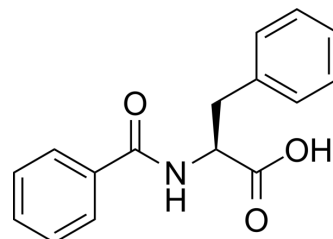


## Benzoyl-L-phenylalanine

Cat. No.:	HY-118391
CAS No.:	2566-22-5
Molecular Formula:	C <sub>16</sub> H <sub>15</sub> NO <sub>3</sub>
Molecular Weight:	269.3
Target:	Others
Pathway:	Others
Storage:	<div> <div>Powder</div> <div>-20°C    3 years</div> <div>4°C    2 years</div> </div> <div> <div>In solvent</div> <div>-80°C    6 months</div> <div>-20°C    1 month</div> </div>



### SOLVENT & SOLUBILITY

In Vitro	DMSO : 100 mg/mL (371.33 mM; Need ultrasonic)					
	Preparing Stock Solutions	<div><div>Solvent</div><div>Concentration</div></div>	Mass	1 mg	5 mg	10 mg
		1 mM		3.7133 mL	18.5667 mL	37.1333 mL
		5 mM		0.7427 mL	3.7133 mL	7.4267 mL
		10 mM		0.3713 mL	1.8567 mL	3.7133 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 (9.28 mM); Clear solution					
	2. Add each solvent one by one: 10% DMSO >> 90% (20% SBE-β-CD in saline) Solubility: ≥ 2.5 mg/mL (9.28 mM); Clear solution					
	3. Add each solvent one by one: 10% DMSO >> 90% corn oil Solubility: ≥ 2.5 mg/mL (9.28 mM); Clear solution					

### BIOLOGICAL ACTIVITY

Description	Benzoyl-L-phenylalanine (N-Benzoyl-L-phenylalanine) is a chorisate mutase-prephenate dehydrogenase inhibitor <sup>[1]</sup> .
IC <sub>50</sub> & Target	Chorisate Mutase-prephenate Dehydrogenase <sup>[1]</sup>

### REFERENCES

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[1]. Smith GD, et al. Affinity chromatography and inhibition of chorismate mutase-prephenate dehydrogenase by derivatives of phenylalanine and tyrosine. Biochem J. 1977 Jul 1;165(1):121-6.

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

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