

## **Product** Data Sheet

# **Kobophenol A**

Cat. No.: HY-126419

CAS No.: 124027-58-3

Molecular Formula:  $C_{56}H_{44}O_{13}$ Molecular Weight: 924.94

Target: SARS-CoV; PKC

Pathway: Anti-infection; Epigenetics; TGF-beta/Smad

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

DMSO: 100 mg/mL (108.12 mM; Need ultrasonic)

Preparing Stock Solutions	Solvent Mass Concentration	1 mg	5 mg	10 mg
	1 mM	1.0812 mL	5.4058 mL	10.8115 mL
	5 mM	0.2162 mL	1.0812 mL	2.1623 mL
	10 mM	0.1081 mL	0.5406 mL	1.0812 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 (2.70 mM); Clear solution
- 2. Add each solvent one by one: 10% DMSO >> 90% (20% SBE-β-CD in saline) Solubility: ≥ 2.5 mg/mL (2.70 mM); Clear solution
- 3. Add each solvent one by one: 10% DMSO >> 90% corn oil Solubility: ≥ 2.5 mg/mL (2.70 mM); Clear solution

#### **BIOLOGICAL ACTIVITY**

Description

Kobophenol A, an oligomeric stilbene, blocks the interaction between the ACE2 receptor and S1-RBD with an IC $_{50}$  of 1.81  $\mu$ M and inhibits SARS-CoV-2 viral infection in cells with an EC $_{50}$  of 71.6  $\mu$ M. Kobophenol A inhibits the activity of partially purified rat brain protein kinase C (PKC) with an IC $_{50}$  of 52  $\mu$ M $^{[1][2]}$ .

#### **REFERENCES**

[1]. Gangadevi S, et al. Kobophenol A Inhibits Binding of Host ACE2 Receptor with Spike RBD Domain of SARS-CoV-2, a Lead Compound for Blocking COVID-19. J Phys

Chem Lett. 2021;12(7):1793-1802. [2]. Xu G, Zhang LP, Chen LF, Hu CQ. Yao Xue Xue Bao. 1994;29(11):818-822. Caution: Product has not been fully validated for medical applications. For research use only. Tel: 609-228-6898 Fax: 609-228-5909 E-mail: tech@MedChemExpress.com Address: 1 Deer Park Dr, Suite Q, Monmouth Junction, NJ 08852, USA

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