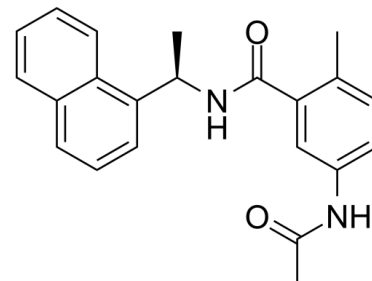


## Data Sheet

<b>Product Name:</b>	PLpro inhibitor
<b>Cat. No.:</b>	CS-2273
<b>CAS No.:</b>	1093070-14-4
<b>Molecular Formula:</b>	C <sub>22</sub> H <sub>22</sub> N <sub>2</sub> O <sub>2</sub>
<b>Molecular Weight:</b>	346.42
<b>Target:</b>	SARS-CoV
<b>Pathway:</b>	Anti-infection
<b>Solubility:</b>	DMSO : ≥ 42 mg/mL (121.24 mM)



### BIOLOGICAL ACTIVITY:

PLpro inhibitor is a potent inhibitor of papain-like protease (PLpro) with IC<sub>50</sub> of 2.6 uM. IC<sub>50</sub> Value: 2.6 uM [1] Target: PLpro; SARS-CoV in vitro: PLpro inhibitor is a potent inhibitor against the papain-like protease (PLpro) from the coronavirus that causes severe acute respiratory syndrome (SARS-CoV). PLpro inhibitor was found to have IC<sub>50</sub> value of 2.6 ± 0.1 μM. PLpro inhibitor display significant antiviral activity with EC<sub>50</sub> values of 13.1±0.7 uM, without toxicity up to the highest concentration tested. Notably, the increasing antiviral potency correlates with the in vitro inhibition of PLpro, suggesting that the compounds work directly on the enzyme in cells [1,2]. in vivo:

### References:

[1]. Ratia, K., et al., A noncovalent class of papain-like protease/deubiquitinase inhibitors blocks SARS virus replication. Proc Natl Acad Sci U S A, 2008. 105(42): p. 16119-24.

[2]. <http://www.google.com/patents/WO2010022355A1cl=en>

### CAIndexNames:

Benzamide, 5-(acetylamino)-2-methyl-N-[(1R)-1-(1-naphthalenyl)ethyl]-

### SMILES:

O=C(C1=CC(NC(C)=O)=CC=C1C)N[C@H](C)C2=C(C=CC=C3)C3=CC=C2

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

Tel: 732-484-9848 Fax: 888-484-5008 E-mail: [sales@ChemScene.com](mailto:sales@ChemScene.com)

Address: 1 Deer Park Dr, Suite Q, Monmouth Junction, NJ 08852, USA