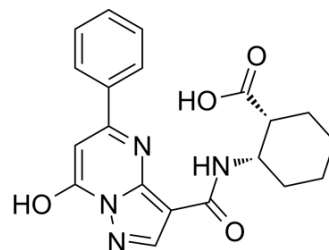


## PF-06928215

Cat. No.:	HY-114182		
Molecular Formula:	C <sub>20</sub> H <sub>20</sub> N <sub>4</sub> O <sub>4</sub>		
Molecular Weight:	380.4		
Target:	Others		
Pathway:	Others		
Storage:	Powder	-20°C	3 years
		4°C	2 years
	In solvent	-80°C	6 months
		-20°C	1 month



### SOLVENT & SOLUBILITY

In Vitro	DMSO : 33.33 mg/mL (87.62 mM; Need ultrasonic)					
	Preparing Stock Solutions	Solvent Concentration	Mass	1 mg	5 mg	10 mg
			1 mM	2.6288 mL	13.1441 mL	26.2881 mL
			5 mM	0.5258 mL	2.6288 mL	5.2576 mL
			10 mM	0.2629 mL	1.3144 mL	2.6288 mL
Please refer to the solubility information to select the appropriate solvent.						
In Vivo	1. Add each solvent one by one: <b>10% DMSO &gt;&gt; 90% (20% SBE-β-CD in saline)</b> Solubility: ≥ 1.43 mg/mL (3.76 mM); Clear solution					
	2. Add each solvent one by one: <b>10% DMSO &gt;&gt; 90% corn oil</b> Solubility: ≥ 1.43 mg/mL (3.76 mM); Clear solution					
	3. Add each solvent one by one: <b>10% DMSO &gt;&gt; 40% PEG300 &gt;&gt; 5% Tween-80 &gt;&gt; 45% saline</b> Solubility: ≥ 1.43 mg/mL (3.76 mM); Clear solution					

### BIOLOGICAL ACTIVITY

Description	PF-06928215 is a cGAS (cyclic GMP-AMP Synthase) inhibitor with an IC <sub>50</sub> of 4.9 μM. PF-06928215 has a high binding affinity of 0.2 μM (K <sub>d</sub> ) <sup>[1]</sup> .
IC <sub>50</sub> & Target	IC <sub>50</sub> : 4.9 μM (cGAS) <sup>[1]</sup> K <sub>d</sub> : 0.2 μM <sup>[1]</sup>

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## REFERENCES

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[1]. Hall J, et al. Discovery of PF-06928215 as a high affinity inhibitor of cGAS enabled by a novel fluorescence polarization assay. PLoS One. 2017 Sep 21;12(9):e0184843.

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

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