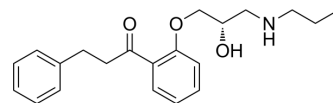


## (S)-Propafenone

Cat. No.:	HY-B0432B		
CAS No.:	107381-32-8		
Molecular Formula:	C <sub>21</sub> H <sub>27</sub> NO <sub>3</sub>		
Molecular Weight:	341.44		
Target:	Sodium Channel		
Pathway:	Membrane Transporter/Ion Channel		
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 (97.62 mM; Need ultrasonic)

Concentration	Solvent	Mass		
		1 mg	5 mg	10 mg
Preparing Stock Solutions	1 mM	2.9288 mL	14.6439 mL	29.2877 mL
	5 mM	0.5858 mL	2.9288 mL	5.8575 mL
	10 mM	0.2929 mL	1.4644 mL	2.9288 mL

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

### BIOLOGICAL ACTIVITY

#### Description

(S)-Propafenone ((S)-SA-79) is the S-enantiomer of Propafenone. (S)-Propafenone ((S)-SA-79) exerts beta-blocking action and the sodium channel-dependent antiarrhythmic class 1 activity<sup>[1]</sup>.

### CUSTOMER VALIDATION

- Clin Chem. 2019 Dec;65(12):1522-1531.

See more customer validations on [www.MedChemExpress.com](http://www.MedChemExpress.com)

### REFERENCES

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[1]. Stoschitzky K, et al. Different stereoselective effects of (R)- and (S)-propafenone: clinical pharmacologic, electrophysiologic, and radioligand binding studies. Clin Pharmacol Ther. 1990 Jun;47(6):740-6.

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

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