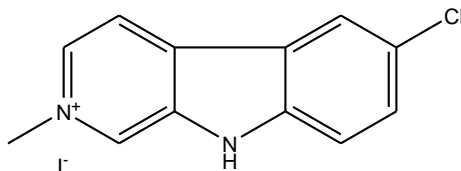


## Nostocarboline iodide

Cat # NB-48-0690



### Product Information

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<b>Batch No.:</b>	0656BN/01
<b>Chemical Name:</b>	6-Chloro-2-methyl-9 <i>H</i> -pyrido[3,4- <i>b</i> ]indol-2-ium iodide; 6-Chloro-2-methylnorharmaline iodide
<b>Batch Molecular Formula:</b>	C <sub>12</sub> H <sub>10</sub> ClIN <sub>2</sub>
<b>Batch Molecular Weight:</b>	344.58
<b>CAS No.:</b>	
<b>Physical Appearance:</b>	Yellow powder
<b>Storage:</b>	Desiccate at +4° C

### Solvent and solubility

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Soluble in water

### Biological activity

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Potent Butyrylcholinesterase (BChE) inhibitor (IC<sub>50</sub> = 13.2 μM), originally isolated from the freshwater cyanobacterium *Nostoc* 78-12A. The inhibitory activity is of the same order of magnitude as that of galanthamine. Very recently, it shows a 50% reduction in parasitaemia at 4 x 50 mg/kg (i.p.) in an in vivo *Plasmodium berghei* mouse model. See also the Des-chloro homologue 2-Methylnorharmaline (Cat. No. NB-48-0063-1MG).

(Sold under license from ETH Zürich; Patent applied)

### References

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1. Becher et al. (2005) *J Nat Prod* 68:1793
2. Blom et al. (2006) *Org Lett* 8:737
3. Bonazzi et al. (2010) *Bioorg Med Chem* 18:1464

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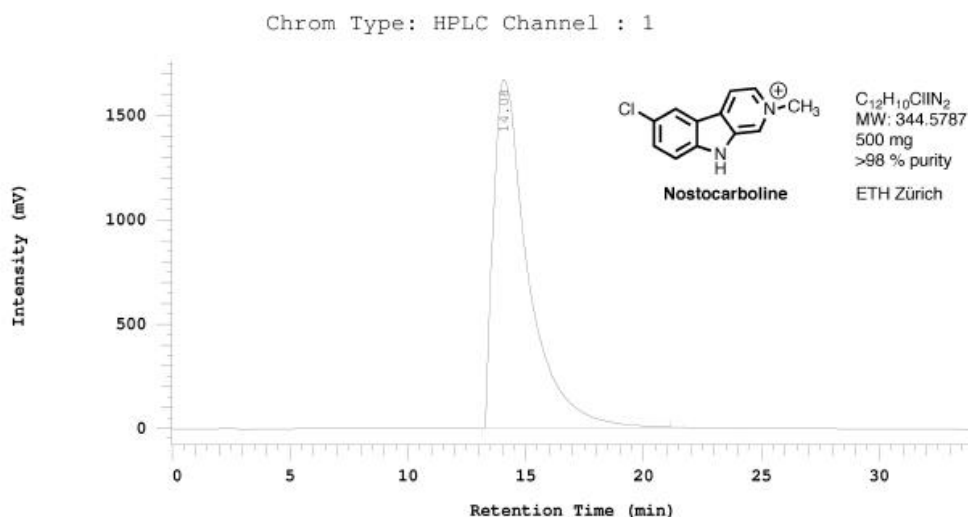
NB-48-0690 Nostocarboline iodide

## Analytical data

HPLC: corresponds to the reference

MS: corresponds to the reference

Tests: HPLC Assay: 98% (complies).



Acquisition Method: method1  
 Column Type: Siehe Vorschrift  
 Pump A Type: L-7100  
 Solvent A: CH<sub>3</sub>CN  
 Solvent C:  
 Method Description:

Developed by: Damien Barbaras  
 Solvent B:  
 Solvent D: H<sub>2</sub>O (TFA 0.1%)

Chrom Type: HPLC Channel : 1

Peak Quantitation: AREA  
 Calculation Method: AREA%

No.	RT	Area	Conc 1	BC
1	14.08	1.698E+08	100.000	BB
		1.698E+08	100.000	

Peak rejection level: 0

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