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DATASHEET

(R)-CPP

Product overview

Name	(R)-CPP
Cat No	HB0021
Biological action	Antagonist
Customer comments	<i>Good quality and good price</i> Verified customer, Seoul National University

Description	Potent NMDA receptor antagonist
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Images



Biological Data

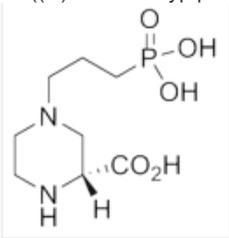
Biological description	Potent NMDA receptor antagonist. More active enantiomer of (RS)-CPP . Selective for GluN2A subtypes over GluN2B, GluN2C and GluN2D subtypes (K_i values are 41 nM, 0.27, 0.63 and 1.99 μ M respectively). Blocks NMDA receptor-mediated EPSCs. Shows antinociceptive effect. Active in vivo.
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Solubility & Handling

Storage instructions	Room temperature (desiccate)
Solubility overview	Soluble in water (100mM)
Important	This product is for RESEARCH USE ONLY and is not intended for therapeutic or diagnostic use. Not for human or veterinary use.

Chemical Data

Chemical name	3-((R)-2-Carboxypiperazin-4-yl)-propyl-1-phosphonic acid
Molecular Weight	252.21
Chemical structure	

Chemical name	3-((R)-2-Carboxypiperazin-4-yl)-propyl-1-phosphonic acid
	
Molecular Formula	C ₈ H ₁₇ N ₂ O ₅ P
CAS Number	126453-07-4
PubChem identifier	6603754
SMILES	C1CN(C[C@@H](N1)C(=O)O)CCCP(=O)(O)O
InChi	InChI=1S/C8H17N2O5P/c11-8(12)7-6-10(4-2-9-7)3-1-5-16(13,14)15/h7,9H,1-6H2,(H,11,12)(H2,13,14,15)/t7-/m1/s1
InChiKey	CUVGUPIVTLGRGI-SSDOTTSWSA-N
Formulation	White solid

References

Structure-activity analysis of a novel NR2C/NR2D-preferring NMDA receptor antagonist: 1-(phenanthrene-2-carbonyl) piperazine-2,3-dicarboxylic acid.

Feng B *et al* (2004) Br J Pharmacol 141(3)

PubMedID [14718249](#)

Long-term potentiation promotes proliferation/survival and neuronal differentiation of neural stem/progenitor cells.

Cho T *et al* (2013) PLoS One 8(10)

PubMedID [24146937](#)

Sertindole restores attentional performance and suppresses glutamate release induced by the NMDA receptor antagonist CPP.

Carli M *et al* (2011) Psychopharmacology (Berl) 214(3)

PubMedID [21049266](#)