

Data Sheet

Product Name:	DR2313	\circ
Cat. No.:	CS-0026455	Ŭ,
CAS No.:	284028-90-6	
Molecular Formula:	C ₈ H ₁₀ N ₂ OS	c~~~
Molecular Weight:	182.24	S II
Target:	PARP	l ll
Pathway:	Cell Cycle/DNA Damage; Epigenetics	
Solubility:	DMSO : 12.5 mg/mL (ultrasonic);H ₂ O : ≥ 100 mg/mL	

BIOLOGICAL ACTIVITY:

DR2313 is a potent, selective, competitive and brain-penetrant inhibitor of **poly(ADP-ribose) polymerase (PARP)**, with **IC**₅₀s of 0.20 μ M and 0.24 μ M for **PARP-1** and **PARP-2**, respectively. DR2313 exhibits neuroprotective effects on ischemic injuries in vitro and in vivo^{[1][2]}. *In Vitro:* DR2313 (0.016-16.4 μ M; 30 min) inhibits poly(ADP-ribosyl)ation reaction in nuclear extracts of rat brain, with a K_i of 0.23 μ M^[1].

DR2313 shows more powerful inhibition of the poly(ADP-ribosyl)ation in the nuclear extracts of the rat brain (IC₅₀=0.20 μ M) than 3AB (35.4 μ M), PND (0.56 μ M), DIQ (2.96 μ M), and DPQ (0.96 μ M)^[1].

DR2313 (1-100 μ M; 10 min) shows a weak inhibition of the mono(ADP-ribosyl)ation in a concentration-dependent manner (IC₅₀=59 μ M)^[1].

DR2313 (0.1-30 μ M; pretreated for 30 min) reduces hydrogen peroxide (500 μ M; 4 h) or glutamate (1 mM; 30 min) induced excessive formation of poly(ADP-ribose) and cell death^[1]. *In Vivo:* DR2313 (3-10 mg/kg i.v. bolus or infusion for 6 h) significantly reduces the cortical infarct volume in both permanent and transient focal ischemia models in rats^[1].

References:

[1]. Nakajima H, et, al. A newly synthesized poly(ADP-ribose) polymerase inhibitor, DR2313 [2-methyl-3,5,7,8-tetrahydrothiopyrano[4,3-d]-pyrimidine-4one]: pharmacological profiles, neuroprotective effects, and therapeutic time window in cerebral ischemia in rats. J Pharmacol Exp Ther. 2005 Feb;312(2):472-81.

[2]. Xu Z, et, al. Endonuclease G does not play an obligatory role in poly(ADP-ribose) polymerase-dependent cell death after transient focal cerebral ischemia. Am J Physiol Regul Integr Comp Physiol. 2010 Jul;299(1):R215-21.

CAIndexNames:

4H-Thiopyrano[4,3-d]pyrimidin-4-one, 3,5,7,8-tetrahydro-2-methyl-

SMILES:

O=C1C(CSCC2)=C2NC(C)=N1

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

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