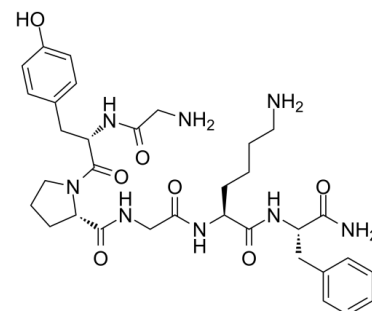


Data Sheet

Product Name:	Protease-Activated Receptor-4
Cat. No.:	CS-7046
CAS No.:	245443-52-1
Molecular Formula:	C33H46N8O7
Molecular Weight:	666.77
Target:	Protease-Activated Receptor (PAR)
Pathway:	GPCR/G Protein
Solubility:	H2O : 50 mg/mL (74.99 mM; Need ultrasonic)



BIOLOGICAL ACTIVITY:

Protease-Activated Receptor-4 is the agonist of **proteinase-activated receptor-4 (PAR4)**. Sequence: Gly-Tyr-Pro-Gly-Lys-Phe-NH₂. **In Vitro:** GYPGKF-NH₂ significantly reduces the agonistic potency of AYPGKF-NH₂ by 25-fold^[1]. GYPGKF-NH₂ (500 μM) does not cause contraction or relaxation of the guinea pig IAS strips^[2].

References:

- [1]. Moschonas IC, et al. Molecular requirements involving the human platelet protease-activated receptor-4 mechanism of activation by peptide analogues of its tethered-ligand. Platelets. 2017 Mar 7:1-10. doi: 10.1080/09537104.2017.1282607. [Epub ahead of print]
- [2]. Huang SC, et al. Proteinase-activated receptor-1 (PAR1) and PAR2 mediate relaxation of guinea pig internal anal sphincter. Regul Pept. 2014 Feb 10;189:46-50.

CAIndexNames:

L-Phenylalaninamide, glycyl-L-tyrosyl-L-prolylglycyl-L-lysyl-

SMILES:

O=C(N)[C@H](CC1=CC=CC=C1)NC([C@H](CCCCN)NC(CNC([C@H]2N(C([C@H](CC3=CC=C(C=C3)O)NC(CN)=O)=O)CCC2)=O)=O)=O

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

Tel: 732-484-9848 Fax: 888-484-5008 E-mail: sales@ChemScene.com

Address: 1 Deer Park Dr, Suite Q, Monmouth Junction, NJ 08852, USA