PRODUCT INFORMATION



ENPP1 Inhibitor C

Item No. 29809

CAS Registry No.: 2378640-92-5

Formal Name: 6-[(3-aminophenyl)methyl]-N,N,5-

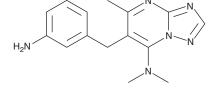
trimethyl-[1,2,4]triazolo[1,5-a]

pyrimidin-7-amine

MF: $C_{15}H_{18}N_{6}$ 282.3 FW: ≥98% **Purity:** UV/Vis.: λ_{max} : 231 nm Supplied as: A crystalline solid

Storage: -20°C Stability: ≥2 years

Information represents the product specifications. Batch specific analytical results are provided on each certificate of analysis.



Laboratory Procedures

ENPP1 inhibitor C is supplied as a crystalline solid. A stock solution may be made by dissolving the ENPP1 inhibitor C in the solvent of choice, which should be purged with an inert gas. ENPP1 inhibitor C is soluble in organic solvents such as ethanol, DMSO, and dimethyl formamide. The solubility of ENPP1 inhibitor C in these solvents is approximately 30 mg/ml.

Description

ENPP1 inhibitor C is an inhibitor of ectonucleotide pyrophosphatase/phosphodiesterase 1 (ENPP1; $IC_{50} = 0.26 \mu M$ in a cell-free assay). It is selective for ENPP1 over ENPP2-7 at 10 μM . ENPP1 inhibitor C decreases ENPP1 activity in MDA-MB-231 human breast and C6 rat glioma cancer cells when used at a concentration of 10 μM.

Reference

1. Kawaguchi, M., Han, X., Hisada, T., et al. Development of an ENPP1 fluorescence probe for inhibitor screening, cellular imaging, and prognostic assessment of malignant breast cancer. J. Med. Chem. 62(20), 9254-9269 (2019).

WARNING
THIS PRODUCT IS FOR RESEARCH ONLY - NOT FOR HUMAN OR VETERINARY DIAGNOSTIC OR THERAPEUTIC USE.

This material should be considered hazardous until further information becomes available. Do not ingest, inhale, get in eyes, on skin, or on clothing. Wash thoroughly after handling. Before use, the user must review the complete Safety Data Sheet, which has been sent via email to your institution.

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CAYMAN CHEMICAL

1180 EAST ELLSWORTH RD ANN ARBOR, MI 48108 · USA PHONE: [800] 364-9897

[734] 971-3335

FAX: [734] 971-3640 CUSTSERV@CAYMANCHEM.COM WWW.**CAYMANCHEM**.COM