PRODUCT INFORMATION



U-44069 serinol amide

Item No. 40872

Formal Name: (Z)-N-(2,3-dihydroxypropyl)-7-

> ((1S,4R,5S,6R)-5-((S,E)-3-hydroxyoct-1en-1-yl)-2-oxabicyclo[2.2.1]heptan-6-yl)

hept-5-enamide

Synonyms: 9,11-dideoxy-9α,11α-epoxymethano

PGF_{2a} serinol amide,

9,11-dideoxy- $9\alpha,11\alpha$ -epoxymethano Prostaglandin $\boldsymbol{F}_{2\alpha}$ serinol amide, 9,11-epoxymethano Prostaglandin H₂

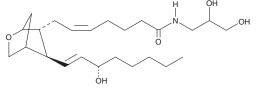
serinol amide

MF: $C_{24}H_{41}NO_5$ 423.6 FW: **Purity:** ≥95%

Supplied as: A solution in methyl acetate

Storage: -20°C Stability: ≥1 year

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



Laboratory Procedures

U-44069 serinol amide is supplied as a solution in methyl acetate. To change the solvent, simply evaporate the methyl acetate under a gentle stream of nitrogen and immediately add the solvent of choice. U-44069 serinol amide is slightly soluble (0.1-1 mg/ml) in chloroform.

Description

U-44069 serinol amide is a derivative of the TP receptor agonist U-44069 (Item No. 16440).

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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