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Product Information



Fluprostenol-d₄

Item No. 316767

Formal Name: 9 α ,11 α ,15R-trihydroxy-16-(3-(trifluoromethyl)phenoxy)-17,18,19,20-tetranor-prosta-5Z,13E-dien-1-oic-3,3,4,4-d₄ acid

Synonym: 16-*m*-trifluoromethylphenoxy tetranor Prostaglandin F_{2 α} -d

MF: C₂₃H₂₅D₄F₃O₆

FW: 462.5

Chemical Purity: \geq 98%

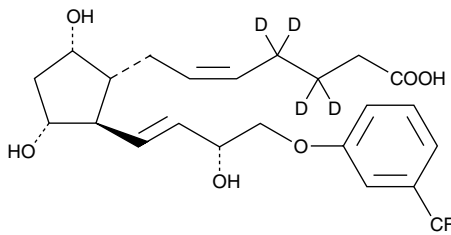
Deuterium

Incorporation: \geq 99% deuterated forms (d₁-d₄); \leq 1% d₀

Stability: \geq 2 years at -20°C

Supplied as: A solution in methyl acetate

UV/Vis.: λ_{\max} : 222, 280 nm



Laboratory Procedures

Fluprostenol-d₄ contains four deuterium atoms at the 3, 3', 4, and 4' positions. It is intended for use as an internal standard for the quantification of Fluprostenol by GC- or LC-mass spectrometry (MS). For long term storage, we suggest that Fluprostenol-d₄ be stored as supplied at -20°C. It should be stable for at least two years.

Fluprostenol-d₄ 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. Solvents such as ethanol, DMSO, and dimethyl formamide purged with an inert gas can be used. The solubility of fluprostenol-d₄ in these solvents is approximately 100 mg/ml.

Fluprostenol-d₄ is used as an internal standard for the quantification of fluprostenol by stable isotope dilution MS. The accuracy of the sample weight in this vial is between 5% over and 2% under the amount shown on the vial. If better precision is required, the deuterated standard should be quantitated against a more precisely weighed unlabeled standard by constructing a standard curve of peak intensity ratios (deuterated *versus* unlabeled).

Fluprostenol is a metabolically stable analog of prostaglandin F_{2 α} (PGF_{2 α}) with potent FP receptor agonist activity.^{1,2} It inhibits PGF_{2 α} binding to human and rat FP receptors with IC₅₀ values of 3.5 and 7.5 nM, respectively.^{1,2} Fluprostenol is a much more potent luteolytic agent than PGF_{2 α} in rats with a minimum fully effective dose of 270 μ g/kg to terminate pregnancy.³ It is also an effective inhibitor of rat adipose precursor differentiation in primary cultures with an IC₅₀ value of 3-10 x 10⁻¹¹ M.⁴

References

1. Abramovitz, M., Boie, Y., Nguyen, T., *et al.* Cloning and expression of a cDNA for the human prostanoid FP receptor. *J. Biol. Chem.* **269**, 2632-2636 (1994).
2. Lake, S., Gullberg, H., Wahlqvist, J., *et al.* Cloning of the rat and human prostaglandin F_{2 α} receptors and the expression of the rat prostaglandin F_{2 α} receptor. *FEBS Lett.* **355**, 317-325 (1994).
3. Dukes, M., Russell, W., and Walpole, A.L. Potent luteolytic agents related to prostaglandin F_{2 α} . *Nature* **250**, 330-331 (1974).
4. Serrero, G. and Lepak, N.M. Prostaglandin F_{2 α} receptor (FP receptor) agonists are potent adipose differentiation inhibitors for primary culture of adipocyte precursors in defined medium. *Biochem. Biophys. Res. Commun.* **233**, 200-202 (1997).

Related Products

For a list of related products please visit: www.caymanchem.com/catalog/316767

WARNING: THIS PRODUCT IS FOR LABORATORY RESEARCH ONLY: NOT FOR ADMINISTRATION TO HUMANS. NOT FOR HUMAN OR VETERINARY DIAGNOSTIC OR THERAPEUTIC USE.

MATERIAL SAFETY DATA

This material should be considered hazardous until information to the contrary becomes available. Do not ingest, swallow, or inhale. Do not get in eyes, on skin, or on clothing. Wash thoroughly after handling. This information contains some, but not all, of the information required for the safe and proper use of this material. Before use, the user must review the complete Material Safety Data Sheet, which has been sent *via* email to your institution.

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