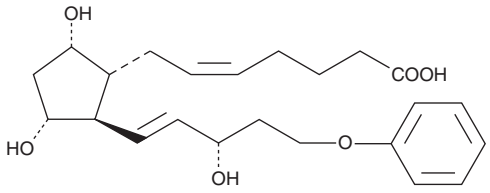
PRODUCT INFORMATION

17-phenoxy trinor Prostaglandin $F_{2\alpha}$

Item No. 10010839

CAS Registry No.: 2162157-41-5
Formal Name: (5Z)-7-[(1R,2R,3R,5S)-3,5-dihydroxy-2-[(1E,3S)-3-hydroxy-5-phenoxy-1-penten-1-yl]cyclopentyl]-5-heptenoic acid
Synonym: 17-phenoxy trinor $PGF_{2\alpha}$
MF: $C_{23}H_{32}O_6$
FW: 404.5
Purity: ≥98%
Supplied as: A solution in methyl acetate
Storage: -20°C
Stability: ≥2 years



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

Laboratory Procedures

17-phenoxy trinor prostaglandin $F_{2\alpha}$ (17-phenoxy trinor $PGF_{2\alpha}$) 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 17-phenoxy trinor $PGF_{2\alpha}$ in these solvents is approximately 50 mg/ml.

Further dilutions of the stock solution into aqueous buffers or isotonic saline should be made prior to performing biological experiments. Ensure that the residual amount of organic solvent is insignificant, since organic solvents may have physiological effects at low concentrations. If an organic solvent-free solution of 17-phenoxy trinor $PGF_{2\alpha}$ is needed, it can be prepared by evaporating the methyl acetate and directly dissolving the neat oil in aqueous buffers. The solubility of 17-phenoxy trinor $PGF_{2\alpha}$ in PBS (pH 7.2) is approximately 10 mg/ml. We do not recommend storing the aqueous solution for more than one day.

Description

17-phenoxy trinor $PGF_{2\alpha}$ is a novel analog of $PGF_{2\alpha}$. A similar analog, 16-phenoxy tetranor $PGF_{2\alpha}$, binds to the FP receptor on luteal cells with much greater affinity (440%) than $PGF_{2\alpha}$.[1] There are no published reports on the biological activity of 17-phenoxy trinor $PGF_{2\alpha}$.

Reference

1. Balapure, A.K., Rexroad, C.E., Jr., Kawada, K., *et al.* Structural requirements for prostaglandin analog interaction with the ovine corpus luteum prostaglandin $F_{2\alpha}$ receptor. *Biochem. Pharmacol.* **38(14)**, 2375-2381 (1989).

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

SAFETY DATA
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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