# **PRODUCT** INFORMATION



16-phenyl tetranor Prostaglandin F<sub>2a</sub>

Item No. 16770

CAS Registry No.:	38315-48-9	
Formal Name:	(5Z)-7-[(1R,2R,3R,5S)-3,5-dihydroxy-2-	
	[(1E,3S)-3-hydroxy-4-phenyl-1-buten-	OH
	1-yl]cyclopentyl]-5-heptenoic acid	$\wedge \wedge - \wedge \wedge$
Synonym:	16-phenyl tetranor PGF <sub>2α</sub>	СООН
MF:	C <sub>22</sub> H <sub>30</sub> O <sub>5</sub>	
FW:	374.5	но
Purity:	≥99%	ОН
Supplied as:	A crystalline solid	
Storage:	-20°C	
Stability:	≥4 years	
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Information represents the product specifications. Batch specific analytical results are provided on each certificate of analysis.

# Laboratory Procedures

16-phenyl tetranor Prostaglandin  $F_{2\alpha}$  (16-phenyl tetranor  $PGF_{2\alpha}$ ) is supplied as a crystalline solid. A stock solution may be made by dissolving the 16-phenyl tetranor  $PGF_{2\alpha}$  in the solvent of choice, which should be purged with an inert gas. 16-phenyl tetranor  $PGF_{2\alpha}$  is soluble in organic solvents such as ethanol, DMSO, and dimethyl formamide. The solubility of 16-phenyl tetranor  $PGF_{2\alpha}$  in these solvents is approximately 100 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. Organic solvent-free aqueous solutions of 16-phenyl tetranor  $PGF_{2\alpha}$  can be prepared by directly dissolving the crystalline solid in aqueous buffers. The solubility of 16-phenyl tetranor  $PGF_{2\alpha}$  in PBS (pH 7.2) is approximately 0.8 mg/ml. We do not recommend storing the aqueous solution for more than one day.

# Description

16-phenyl tetranor  $PGF_{2\alpha}$  is a metabolically stable analog of  $PGF_{2\alpha}$ . The affinity of 16-phenyl tetranor  $PGF_{2\alpha}$  at the FP receptor of ovine luteal cells is poor (8.7%) compared to  $PGF_{2\alpha}$ .<sup>1</sup>

# 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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