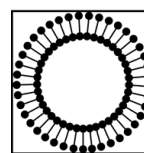
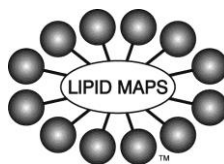


1-heptadecanoyl-2-(9Z-tetradecenoyl)-sn-glycero-3-phosphoethanolamine

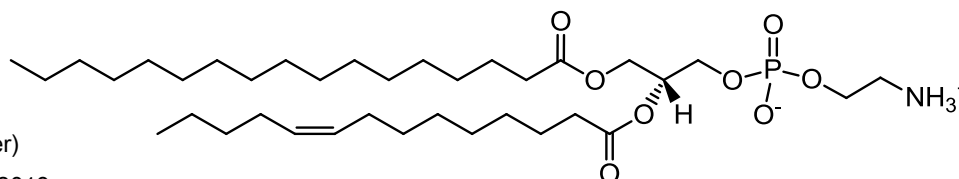
17:0-14:1 PE

LIPIDMAPS™ Compound ID: LMGP02010005

**Avanti®**
POLAR LIPIDS, INC.

Molecular Weight: 675.917
 Exact Mass: 675.484
 Molecular Formula: C₃₆H₇₀NO₈P
 Storage: -20° C (Freezer)

Manufacture Date: September 4, 2013
 Expiration Date: One year from date of receipt
 Physical Form: Methanol Solution (1ml per vial)
 Concentration: 9.64 µg/ml (14.26 µM)
 Lot Number: LM1104-LM27-001C
 MLOT: 6411WAC001

**Product Specifications**

Test	Limits	Result
Mass Spec	Consistent with Structure, [M+H] ⁺ = 676.5 ±1amu	Pass
Quantitative LC/MS	Run and Report	9.64 µg/ml 14.26 µM
HPLC/ELSD	NLT 99% Purity (AUC)	100% AUC
GC-FAME	NLT 99% Purity (AUC) (14:1+17:0)	100% AUC
Phosphorus NMR	Consistent with Structure	Pass
Proton NMR	Consistent with Structure	Pass
TLC Analysis (Mobile Phase 65:25:4 Chloroform:Methanol:Water)	R _f consistent with structure, NLT 99% Purity	Pass
Ninhydrin	Positive	Positive
Iodine	One Major Spot	One Spot
Phosphorus	Positive	Positive
Char	Positive	Positive
Wet Dip	One Major Spot	One Spot

Approved by:

Avanti Polar Lipids, Inc. 700 Industrial Park Drive, Alabaster, AL 35007-9105, USA

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