

# 1-Stearoyl-2-Arachidonoyl-sn-Glycero-3-Phosphatidylcholine

**Catalog Number:** LP-2155

**CAS Number:** 35418-59-8

**Price:** [Online Inquiry](#)

| Catalog Number | Price                          |
|----------------|--------------------------------|
| LP-2155        | <a href="#">Online Inquiry</a> |

| Product Information |                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                   |
|---------------------|---------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| Molecular Formula   | C46H84NO8P                                                                                                                                                                                                                                                                                                                                                                                                                                                                                        |
| Molecular Weight    | 810.14                                                                                                                                                                                                                                                                                                                                                                                                                                                                                            |
| Purity              | >98%                                                                                                                                                                                                                                                                                                                                                                                                                                                                                              |
| SMILES              | <chem>O=C(OCC(OC(=O)CCCC=CCC=CCC=CCC=CCCCC)COP(=O)([O-])OCC[N+](C)(C)C)CCCCC</chem>                                                                                                                                                                                                                                                                                                                                                                                                               |
| InChi               | InChI=1S/C46H84NO8P/c1-6-8-10-12-14-16-18-20-22-23-25-27-29-31-33-35-37-39-46(49)55-44(43-54-56(50,51)53-41-40-47(3,4)5)42-52-45(48)38-36-34-32-30-28-26-24-21-19-17-15-13-11-9-7-2/h14,16,20,22,25,27,31,33,44H,6-13,15,17-19,21,23-24,26,28-30,32,34-43H2,1-5H3/b16-14-,22-20-,27-25-,33-31-/t44-/m1/s1                                                                                                                                                                                         |
| InChi Key           | InChIKey=PSVRFUPOQYJOOZ-QNPWAGBNSA-N                                                                                                                                                                                                                                                                                                                                                                                                                                                              |
| Alternate Name      | 3,5,8-Trioxa-4-phosphaoctacosa-13,16,19,22-tetraen-1-aminium, 4-hydroxy-N,N [...],N-trimethyl-9-oxo-7-[[[(1-oxooctadecyl)oxy]methyl]-, inner salt, 4-oxide, [R-(all-Z)], 18:0/20:4 Phosphatidylcholine, 1-Stearoyl-2-arachidonoyl-sn-glycero-3-phosphorylcholine, 1-Stearoyl-2-arachidonoyl-sn-glycero-3-phosphatidylcholine, 1-Octadecanoyl-2-arachidonoyl-sn-glycero-3-phosphocholine, 1-Stearoyl-2-arachidonoyl-sn-glycero-3-phosphocholine, 1-Stearoyl-2-archidonoyl-3-sn-phosphatidylcholine |
| Appearance          | Solid                                                                                                                                                                                                                                                                                                                                                                                                                                                                                             |
| IUPAC Name          | 3,5,8-Trioxa-4-phosphaoctacosa-13,16,19,22-tetraen-1-aminium, 4-hydroxy-N,N,N-trimethyl-9-oxo-7-[[[(1-oxooctadecyl)oxy]methyl]-, inner salt, 4-oxide, (7R,13Z,16Z,19Z,22Z)-                                                                                                                                                                                                                                                                                                                       |

| Properties         |         |
|--------------------|---------|
| Storage Conditions | Freezer |