

MATREYA



Lipids, Biochemicals,
and Standards for
Life Science Research

2013 - 2014

About Matreya LLC

- Matreya strives to develop, manufacture and deliver products of the highest value to our customers.
- Quality will always be the best achievable by state-of-the-art techniques, typically greater than 98%
- We strive for rapid delivery. 95+% of our products are shipped within 24 hours of receipt of an order.
- Within the area of sphingolipids and glycolipids, we have earned a reputation as the preferred problem solver and technology leader.
- When you demand quality and consistency, you may rely on Matreya lipids.

Matreya Products for Biochemistry Research.

We offer one of the widest selection of ceramides for intracellular signaling research available. We stock antibodies to glycosphingolipids as well as inhibitors of enzymes involved in glycosphingolipid metabolism.

Our products provide the valuable tools for the study of cell membrane and its structure, growth regulators in the cellular metabolism, and intracellular mediators.

We are able to make our products better and better with the latest technology in Chromatography, Mass Spectrometry, and NMR techniques.

We are proud to offer our products as a valuable tool for your life science research needs.

Matreya Products for Microbiology Research.

Matreya stocks many unusual fatty acid standards produced by bacteria that are useful for culture characterization.

Matreya Products for the Food and Agriculture Industries.

Many of Matreya's fatty acid products have been industry standards for many years. The acids and their methyl esters are used as standards in analysis and quality control.

Custom Preparations.

Our experience in chemical synthesis and the extraction and purification of natural products allows us to produce custom preparations with the same high quality and purity as the products listed in the catalog. Depending on the complexity of the molecule, delivery will be 4 to 12 weeks after receipt of an order, usually less than 6 weeks.

If you can't find a product in the catalog, please check the INDEX, where we also try to list common synonyms for our products.

TABLE OF CONTENTS

TABLE OF CONTENTS.....	i
Technical Service.....	.iii
Natural Products.....	.iii
Storage.....	.iii
Sphingolipid Structures and Pathways.....	.iii
Package Weightiii
Matreya's Mission.....	.iii
Sphingoid Bases, Sphingolipids and Glycosphingolipids	1
Sphingosines	2
Synthetic Sphingosines with C18 Sphingoid Base	2
Synthetic Sphingosines with Sphingoid Bases other than C18.....	2
Synthetic Dihydrosphingosines.....	3
3-Keto-Dihydrosphingosines.....	4
Phytosphingosines	5
Other Sphingosine Derivatives and Precursors.....	5
Ceramides	6
Synthetic Ceramides Derived from C18-Sphingosine	7
2-Hydroxy Ceramides	9
Ceramide Made from Sphingosines with Sphingoid Bases Other Than C18.....	10
Dihydroceramides	10
2-Hydroxy Dihydroceramides	11
Ceramides From Natural Sources	11
Phytoceramides	12
Fluorescent Ceramides.....	13
Phosphosphingolipids	15
Sphingomyelins.....	15
Sphingosylphosphorylcholines (SPC).....	17
Sphingosine Phosphates	17
Fluorescent Sphingomyelins.....	18
Glycosphingolipids	19
Galactosylceramides and Glucosylceramides	19
Lactosylceramides	24
Ceramide Trihexosides	25
Globosides	26
Stable Isotopes Labeled Glycolipids.....	27
Fluorescent Compounds.....	27
Gangliosides	29
Glycosphingolipid Reference Mixtures for TLC	32
Antibodies Directed Against Glycolipids	33
Enzyme Inhibitors	35
Glycerolipids	40
Glycerophospholipids	40
Natural Phospholipids	40
Synthetic Phospholipids	42
Phosphatidylcholines	42
Phosphatidylglycerols	44
Phosphatidylethanolamines	45
Phosphatidylinositols	46
Bacterial Tetraethers.....	47
Glycosyl Glycerides.....	48
Fatty Acids	48

Simple Fatty Acids.....	48
Saturated Fatty Acids and Methyl Esters.....	48
Unsaturated Fatty Acids and Methyl Esters.....	53
Trans Fatty Acids and Methyl Esters	59
Conjugated Linoleic Acid Isomers (CLA)	61
Other CLA Products and Derivatives.....	63
Hydroxy Fatty Acids	63
2-Hydroxy Fatty Acids and Methyl Esters.....	63
3-Hydroxy Fatty Acids and Methyl Esters.....	66
Omega Hydroxy Fatty Acids	68
Other Hydroxy Fatty Acids	70
Branched and Cyclic Fatty Acids	70
Iso-Fatty Acids and Esters	70
Anteiso-Fatty Acids and Esters.....	72
Methylated Fatty Acids.....	72
Cyclopropyl Fatty Acids and Esters	73
Unusual Fatty Acids and Derivatives	73
Other Lipids.....	74
Tocopherols	74
Cholestane Derivatives	76
Plant Sterols and Steryl Glucosides	76
Standards and Reference Compounds	78
Food Industry Mixtures	78
Each methyl ester mixture is carefully prepared by weight.....	78
Polyunsaturated Fatty Acid Methyl Esters Mixtures	79
Carbohydrate Mixtures.....	79
Other Fatty Acid Methyl Ester Mixtures	80
AOCS Animal and Vegetable Oil Reference Mixtures (RM Mixtures).....	81
Custom Mixtures	83
GLC Standard Mixtures	84
Water Soluble Fatty Acid Mixtures.....	85
Microbiology Standard Mixtures	86
Biochemical Research Standard Mixtures.....	87
Glycosphingolipid Reference Mixtures for TLC	87
Biochemicals and Reagents.....	88
Stable Isotope Labeled Compounds.....	88
Fluorescent Compounds.....	89
Sphingolipid Structures and Pathways Wall Chart.....	98
Literature References.....	99
Cross Reference for Product Numbers and Catalog Pages	100
Product Name Index.....	103
Tables	
Table I. AOCS Oil Reference Mixes.....	81
Table II. Standards for GC Analysis	84
Table III. Typical Fatty Acid Composition of Natural Lipids Made by Matreya LLC.	93

All chemicals listed in this catalog are for investigational use only. Chemicals are not for human consumption or to be used in food or food additives. Matreya assumes no liability for any use of these chemicals by the end user. We believe the information in this catalog, offered in good faith, is accurate.

Limited Warranty: All Matreya Products, except those specifically exempted, are warranted (for 30 days) to be free of defects in materials and workmanship, if properly stored. Any replacements required as a result of such defects will be made without charge provided that such defective products are returned with a written explanation. Please request a Returned Goods Authorization before returning products under this warranty.

Technical Service

Our technical service department may be contacted by telephone at 800.342.3595, or by e-mail at techservice@matreya.com.

Natural Products

Some of our glycolipids are extracted from natural sources. These products have a normal heterogeneity in their lipid components, particularly in the fatty acids. Variations include carbon chain length as well as the presence or absence of 2-hydroxy fatty acids. Products based on sphingosine may contain longer chain sphingoid bases as well as chains with multiple double bonds. This heterogeneity may result in additional spots showing on TLC plates or multiple peaks in LC analyses. We have listed the typical fatty acid compositions of our natural products in the appendix.

Storage

Catalog items in unopened containers are stable for at least one year when stored under the conditions indicated in the catalog listing. Items containing unsaturated fatty acids are subject to oxidation and should be stored in a solution of organic solvents or under argon. Glycolipids and phospholipids should not be stored in aqueous solutions due to potential hydrolysis.

Sphingolipid Structures and Pathways

In a clear and straightforward manner, this wall chart indicates the structures and relationships between most commonly discussed sphingolipids. A one-page thumbnail version of the chart is shown on page 98. Full size copies (approximately 35 x 26 inches) are available on request to customer service.

Package Weight

Unless otherwise specified, the package will contain at least the indicated amount and usually slightly more. The user is cautioned to always measure the required amount from the container.

Matreya's Mission

Matreya is committed to manufacturing high purity lipids to be used as research standards in the biotechnology and pharmaceutical areas. These lipids will be offered world-wide at a fair market price, and at a profit sufficient to assure company growth, for the benefit of its customers, employees, share holders, and community. Matreya will also be committed to fast delivery, excellent technical backup, new product development, safety, and environmentally friendly.

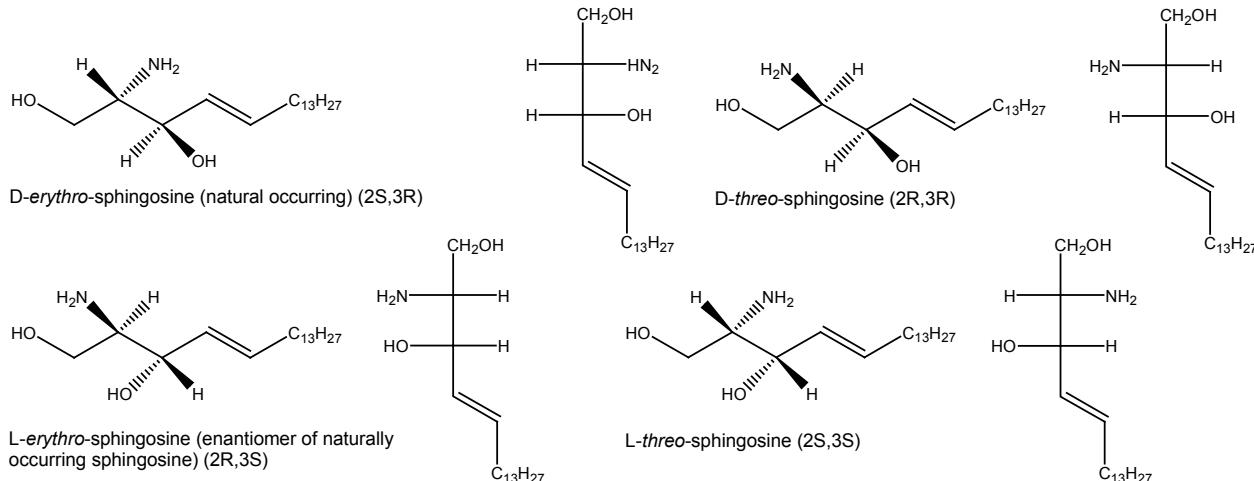
Sphingoid Bases, Sphingolipids and Glycosphingolipids

Sphingoid bases such as sphingosine are the characteristic structural unit of the sphingolipids. The bases are long chain aliphatic amines, containing two or three hydroxyl groups, and typically a *trans*-double bond at C4. In animal tissues, the most abundant base is sphingosine with a C18 aliphatic chain containing a double bond in position 4. The saturated analogue is dihydrosphingosine or sphinganine. In plants, the common long chain base is the 4 hydroxy saturated base phytosphingosine.

Sphingolipids are widely distributed in animal tissues, particularly cell membranes. Sphingoid bases linked to fatty acids via an amide bond at C2 are ceramides and are present in trace amounts in most tissues. Glycosphingolipids (ceramides having various mono- and oligosaccharides on the OH group at C1) are neutral glycosphingolipids (i.e., cerebrosides and globosides). Those with sialic acid derivatized sugars are acidic glycolipids (i.e., gangliosides). They are amphiphilic and can be solubilized in buffers via sonication and micelle formation.

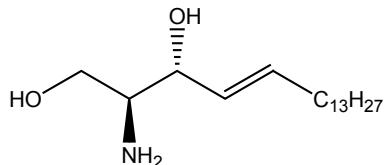
Gangliosides are present in substantial amounts in nerve cell membranes, and together with globosides are found in the membranes of white and red blood cells. These plus the glycosphingolipids of the lacto- and neolacto-series are involved in cell recognition (e.g. blood group determinants). Glycolipid expression on the surface of cells determines their antigenicity as well as their status, i.e. differentiated vs. undifferentiated (embryonic), normal vs. malignant, etc. (1). The ganglioside GM1 stimulates nerve growth (2,3) and has been reported to have a curative effect on experimental Parkinsonism (4). For an overview, see (5). Gangliosides are also being investigated as potential anti-tumor vaccines (6). Glycosphingolipids are also essential for the correct functioning of cell surface receptors (7). Matreya is your best source for many sphingolipids. Most of Matreya's sphingosines and ceramides are fully synthetic and as such 98%+ pure. Others, particularly the glycosphingolipids are highly purified natural products (98%+), and can be used either as standards or biochemical reagents without further purification.

Through total synthesis, all four isomers of sphingosine are available as well as a number of sphingosines with other than 18 carbons and a number of ceramides (for details in using ceramides in cell culture see Hauser et al. [9]). Fluorescent labeled ceramides, glycosphingolipids and sphingomyelins are also available for study. D. N. Brindley and his group have been exploring the interaction of ceramides, sphingosine and sphingosine 1-phosphate in regulating DNA synthesis and phospholipase D activity. **See Literature References on page 99.**



Sphingosines

Synthetic Sphingosines with C18 Sphingoid Base



Catalog number 1802

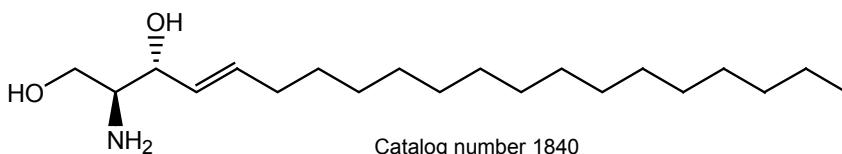
			25 mg	77.00
1802	D-erythro-Sphingosine Sphingosine with C18 chain C ₁₈ H ₃₇ NO ₂ CAS#: 123-78-4		25 mg	77.00
	Source: synthetic Mol. Wt.: 299 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol, DMSO Storage: -20°C			
	Selective inhibitor of phosphokinase C			
1806	L-threo-Sphingosine L-threo-Sphingosine, C18 chain C ₁₈ H ₃₇ NO ₂		10 mg	232.00
	Source: synthetic Mol. Wt.: 299 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol, DMSO Storage: -20°C			
1826	L-erythro-Sphingosine L-erythro-Sphingosine, C18 chain C ₁₈ H ₃₇ NO ₂ CAS#: 6036-75-5		5 mg	262.00
	Source: synthetic Mol. Wt.: 299 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol, DMSO Storage: -20°C			
1827	D-threo-Sphingosine D-threo-Sphingosine, C18 chain C ₁₈ H ₃₇ NO ₂ CAS#: 6036-85-7		5 mg	262.00
	Source: synthetic Mol. Wt.: 299 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol, DMSO Storage: -20°C			

Synthetic Sphingosines with Sphingoid Bases other than C18

Varying chain lengths allow the study of translocation effects of sphingosines and ceramides into cells.

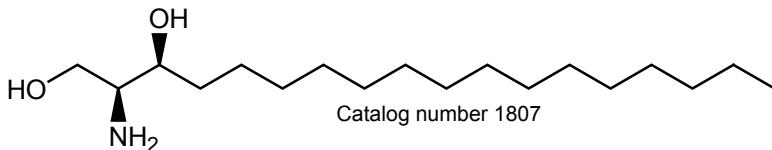
1833	D-erythro-C14-Sphingosine Sphingosine with C14 chain C ₁₄ H ₂₉ NO ₂		5 mg	232.00
	Source: synthetic Mol. Wt.: 243 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol, DMSO Storage: -20°C			
1634	omega-N-NBD-D-erythro-C14-Sphingosine <i>omega</i> -N-(7-nitrobenzo-2-oxa-1,3-diazol-4-yl)-(2S)-amino-tetradec-(4E)-ene-(1,3R)-diol C ₂₀ H ₃₁ N ₅ O ₅		1 mg	522.00
	Source: synthetic Mol. Wt.: 422 Purity: 98+% by TLC Appearance: solid Solubility: methanol, ethanol, chloroform/methanol, 9:1 Storage: -20°C			

1835	D-erythro-C16-Sphingosine Sphingosine with C16 chain C ₁₆ H ₃₃ NO ₂	5 mg	232.00
	Source: synthetic Mol. Wt.: 271 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: -20°C		
1838	D-erythro-C12-Sphingosine Sphingosine with C12 chain C ₁₂ H ₂₅ NO ₂ CAS#: 6918-49-6	5 mg	232.00
	Source: synthetic Mol. Wt.: 215 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: -20°C		
1840	D-erythro-C20-Sphingosine Sphingosine with C20 chain C ₂₀ H ₄₁ NO ₂	5 mg	227.00
	Source: synthetic Mol. Wt.: 328 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol, DMSO Storage: -20°C		



Synthetic Dihydrosphingosines

D,L-*threo*- Dihydrosphingosine has also been found to be a significant inhibitor of sphingosine kinase (8). The D,L-*erythro*-isomer has been used as an inactive control. We offer all four isomers in pure form making detailed studies possible. Safingol, the L-*threo*-isomer is a potent inhibitor of PKC and as such is capable of reversing multi-drug resistance in cancer cells (9). See Literature References on page 99.



1807	L-<i>threo</i>-Dihydrosphingosine (Safingol)	5 mg	174.00
1807-025	L- <i>threo</i> -Sphinganine, C18 chain C ₁₈ H ₃₉ NO ₂ CAS#: 15639-50-6	25 mg	556.00

Source: synthetic Mol. Wt.: 301 Melting Point (°C): 103-114 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol, DMSO Storage: -20°C

Inhibitor of PKC

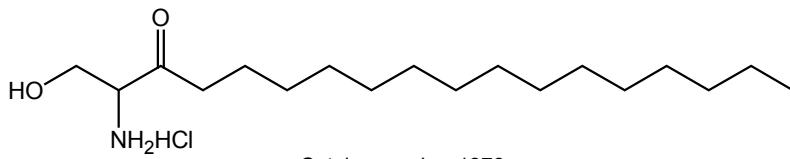
1831	D-<i>erythro</i>-Dihydrosphingosine	25 mg	144.00
1831-1	D- <i>erythro</i> -Sphinganine, C18 chain C ₁₈ H ₃₉ NO ₂ CAS#: 764-22-7	1 g	2,494.00

Source: synthetic Mol. Wt.: 301 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol, DMSO Storage: -20°C

Inhibitor of PLA₂ and PLD

1846	L-erythro-Dihydrosphingosine L- <i>erythro</i> -Sphinganine, C18 chain C ₁₈ H ₃₉ NO ₂	1 mg	111.00
Source: synthetic Mol. Wt.: 301 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol, DMSO Storage: -20°C			
1851	D-threo-Dihydrosphingosine D- <i>threo</i> -Sphinganine, C18 chain C ₁₈ H ₃₉ NO ₂ CAS#: 6036-86-8	1 mg	169.00
Source: synthetic Mol. Wt.: 301 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol, DMSO Storage: -20°C			
1324	D,L-erythro-Dihydrosphingosine D,L- <i>erythro</i> -Sphinganine, C18 chain C ₁₈ H ₃₉ NO ₂ CAS#: 3102-56-5	25 mg	103.00
Source: synthetic Mol. Wt.: 301 Purity: erythro 77%; threo 23% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol, DMSO Storage: -20°C			
Inhibitor of sphingosine kinase			
1326	D,L-C16-Dihydrosphingosine (mixed isomers) D,L-Sphinganine with C16 chain C ₁₆ H ₃₅ NO ₂	10 mg	111.00
Source: synthetic Mol. Wt.: 273 Purity: <i>erythro</i> 90%, <i>threo</i> 10% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol, DMSO Storage: -20°C			
1845	D-erythro-C20-Dihydrosphingosine D- <i>erythro</i> -Sphinganine, C20 chain C ₂₀ H ₄₃ NO ₂ CAS#: 24006-62-0	5 mg	144.00
Source: synthetic Mol. Wt.: 330 Purity: 98+% by TLC, GC Appearance: solid Solubility: warm ethanol, chloroform/methanol, 5:1 Storage: -20°C			
1839	D,L-erythro-C20-Dihydrosphingosine D,L- <i>erythro</i> -Sphinganine, C20 chain C ₂₀ H ₄₃ NO ₂	10 mg	103.00
Source: synthetic Mol. Wt.: 330 Purity: 98+% by TLC, GC Appearance: solid Solubility: warm ethanol, chloroform/methanol, 5:1 Storage: -20°C			

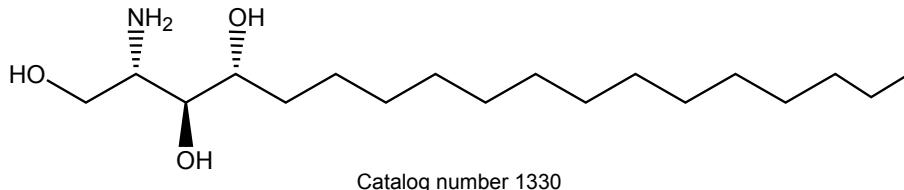
3-Keto-Dihydrosphingosines



1876	3-keto-Dihydrosphingosine•HCl 3-keto-Sphinganine hydrochloride C ₁₈ H ₃₇ NO ₂ •HCl CAS#: 18944-28-0	10 mg	405.00
Source: synthetic Mol. Wt.: 299 + HCl Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: -20°C			

1891	3-keto-C6-Dihydrosphingosine•HCl 1-Hydroxy-2-amino-3-keto-hexane • HCl C ₆ H ₁₃ NO ₂ •HCl	10 mg	405.00
	Source: synthetic Mol. Wt.: 168 Purity: 98+% by TLC Appearance: solid Solubility: ethanol, methanol, DI water Storage: -20°C		
1892	3-keto-C8-Dihydrosphingosine•HCl 1-Hydroxy-2-amino-3-keto-octane • HCl C ₈ H ₁₇ NO ₂ •HCl	10 mg	405.00
	Source: synthetic Mol. Wt.: 196 Purity: 98+% by TLC Appearance: solid Solubility: chloroform, ethanol, methanol, DI water Storage: -20°C		
1893	3-keto-C12-Dihydrosphingosine•HCl 1-Hydroxy-2-amino-3-keto-dodecane • HCl C ₁₂ H ₂₅ NO ₂ •HCl	10 mg	405.00
	Source: synthetic Mol. Wt.: 252 Purity: 98+% by TLC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: -20°C		

Phytosphingosines

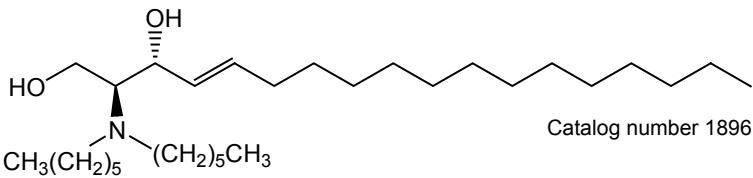


1330	Phytosphingosine	50 mg	111.00
1330-1	4-Hydroxysphinganine C ₁₈ H ₃₉ NO ₃ CAS# 554-62-1	1 g	752.00
	Source: natural, yeast (Pichia ciferri) Mol. Wt.: 318 Purity: 98+% by TLC, GC Appearance: solid Solubility: ethanol, methanol, chloroform/methanol, 2:1 (warm) Storage: -20°C		

Other Sphingosine Derivatives and Precursors

1320	N,N-Dimethyl-D-erythro-sphingosine C ₂₀ H ₄₁ NO ₂ CAS#: 119567-63-4	5 mg/ml, 1 ml	103.00
	Source: synthetic Mol. Wt.: 328 Purity: 98+% by TLC Appearance: liquid Solvent: isopropanol Solubility: chloroform, ethanol, isopropanol, methanol Storage: -20°C		

Inhibitor of phosphokinase C



1896	N,N-Dihexyl-D-erythro-sphingosine Sphingosine with tertiary amine group	5 mg/ml, 1 ml	160.00
		C ₃₀ H ₆₁ NO ₂	

Source: synthetic **Mol. Wt.:** 468 **Purity:** 95% by TLC **Appearance:** liquid
Solvent: ethanol **Solubility:** chloroform, ethanol, methanol **Storage:** -20°C

1805	N-Palmitoyl serinol C ₁₉ H ₃₉ NO ₃ CAS#: 126127-31-9	10 mg	126.00
-------------	---	--------------	---------------

Source: synthetic **Mol. Wt.:** 329 **Purity:** 98+% by TLC, GC **Appearance:** solid
Solubility: chloroform, methanol, ethanol **Storage:** -20°C

Sphingosine precursor

Ceramides

Ceramide is a fatty acid amide of sphingosine. It may be formed by dehydrogenation of dihydroceramide; by hydrolysis of sphingomyelin or glycosphingolipids; or by acylation of free sphingosine. Ceramide functions as a precursor in the synthesis of sphingomyelin (by an exchange reaction with phosphatidylcholine and phosphatidylethanolamine); of glycosphingolipids (by glycosylation with UDP-hexose); and of free sphingosine and fatty acid by hydrolysis. The sphingosine can be phosphorylated by a kinase to form sphingosine-1-phosphate, which may undergo further hydrolysis or cleavage.

Control of sphingolipid metabolism maintains vital balance points in cell physiology. Two of ceramide's metabolites, sphingosine-1-phosphate and glucosylceramide, produce cell proliferation. Sphingosine-1-phosphate is also a highly active regulator of angiogenesis, vascular maturation, cardiac development, immunity, and directed cell movement. Sphingosine, the free base, is a potent inhibitor of protein kinase C and is involved in intracellular calcium regulation.

Sphingolipid enzymes seem to be particularly active in cancers, so modifying their activities by exogenous action may provide alternatives to chemical therapies. These enzymes are controlled by many known agents, such as 1,25-dihydroxy-vitamin D₃, tumor necrosis factor- α , nerve growth factor, interleukin 1, endothelial growth factor, glutathione, arachidonic acid, dexamethasone, many anticancer drugs, therapeutic radiation, and activators of the FAS receptor.

Ceramide exerts numerous biological effects, including induction of cell maturation, cell cycle arrest, terminal cell differentiation, cell senescence, and cell death. Other effects include producing reactive oxygen in mitochondria (followed by apoptosis) and stimulating phosphorylation of certain proteins (especially mitogen activated protein). It also stimulates some protein phosphatases (especially protein phosphatase 2A). Thus ceramide is an important controller of protein activity.

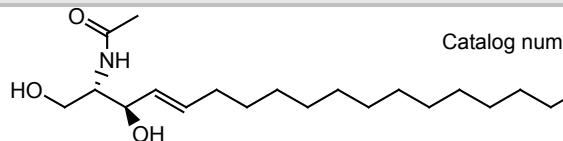
It is apparent from these relationships that ceramide exists at the crux of several enzyme reaction cycles and that experiments involving sphingolipid function call for control of all of the cycles and their branch-off points. Matreya is the major supplier of these lipids, which can be used as standards for analysis of tissues (a much needed part of modern research) and identification of major sphingolipids.

Ceramides with short side chains have been shown to enter easily into cells where they are biologically active. Ceramides with longer side chains, however, also enter cells if dissolved in dodecane-isopropanol first. Fluorescent labeled ceramides and sphingomyelin made from fluorescent labeled acids instead of plain fatty acids are also available for the study of the localization and metabolism of sphingolipids in the cell. Matreya now offers all four isomers of C2, C4, C6 and C18 ceramides. The corresponding dihydroceramides are being used as inactive controls

In three major reviews, Radin (10-12) has discussed the biochemistry and chemistry of ceramide and outlined many potential approaches to cancer therapy using ceramides and related compounds as generators of apoptosis.

See Literature References on page 99.

Synthetic Ceramides Derived from C18-Sphingosine



Catalog number 1901

1901	N-Acetyl-D-erythro-sphingosine		10 mg	103.00
1901-100	N-C2:0-D- <i>erythro</i> -Ceramide	C ₂₀ H ₃₉ NO ₃	100 mg	667.00

Source: synthetic **Mol. Wt.:** 342 **Purity:** 98+ by TLC, GC **Appearance:** solid
Solubility: chloroform, ethanol, methanol, DMSO (up to 5 mg/ml) **Storage:** -20°C

1829	N-Acetyl-L-threo-sphingosine		1 mg	118.00
	N-C2:0-L- <i>threo</i> -Ceramide	C ₂₀ H ₃₉ NO ₃		

Source: synthetic **Mol. Wt.:** 342 **Purity:** 98+% by TLC, GC **Appearance:** solid
Solubility: chloroform, ethanol, methanol, DMSO, DMF (up to 5 mg/ml)
Storage: -20°C

1847	N-Acetyl-L-erythro-sphingosine		1 mg	103.00
	N-C2:0-L- <i>erythro</i> -Ceramide	C ₂₀ H ₃₉ NO ₃		

Source: synthetic **Mol. Wt.:** 342 **Purity:** 98+% by TLC, GC **Appearance:** solid
Solubility: chloroform, ethanol, methanol, DMSO, DMF (up to 5 mg/ml)
Storage: -20°C

1900	N-Hexanoyl-D-erythro-sphingosine		10 mg	103.00
1900-100	N-C6:0-D- <i>erythro</i> -Ceramide	C ₂₄ H ₄₇ NO ₃	100 mg	667.00

Source: synthetic **Mol. Wt.:** 398 **Purity:** 98+% by TLC, GC **Appearance:** solid
Solubility: chloroform, ethanol, DMSO (up to 5 mg/ml) **Storage:** -20°C

1828	N-Hexanoyl-L-threo-sphingosine		1 mg	135.00
	N-C6:0-L- <i>threo</i> -Ceramide	C ₂₄ H ₄₇ NO ₃		

Source: synthetic **Mol. Wt.:** 398 **Purity:** 98+% by TLC, GC **Appearance:** solid
Solubility: chloroform, ethanol, DMSO, DMF (up to 5mg/ml) **Storage:** -20°C

1848	N-Hexanoyl-L-erythro-sphingosine		1 mg	118.00
	N-C6:0-L- <i>erythro</i> -Ceramide	C ₂₄ H ₄₇ NO ₃		

Source: synthetic **Mol. Wt.:** 398 **Purity:** 98+% by TLC, GC **Appearance:** solid
Solubility: chloroform, ethanol, DMSO, DMF (up to 5mg/ml) **Storage:** -20°C

1809	N-Hexanoyl-D-threo-sphingosine N-C6:0-D- <i>threo</i> -Ceramide C ₂₄ H ₄₇ NO ₃	1 mg	135.00
Source: synthetic Mol. Wt.: 398 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, methanol, DMSO (up to 5 mg/ml) Storage: -20°C			
1903	N-Octanoyl-D-erythro-sphingosine	10 mg	103.00
1903-100	N-C8:0-D- <i>erythro</i> -Ceramide C ₂₆ H ₅₁ NO ₃ CAS#: 74713-59-0	100 mg	667.00
Source: synthetic Mol. Wt.: 426 Purity: 98+ by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol, DMSO (up to 5 mg/ml) Storage: -20°C			
1830	N-Octanoyl-L-threo-sphingosine N-C8:0-L- <i>threo</i> -Ceramide C ₂₆ H ₅₁ NO ₃	1 mg	135.00
Source: synthetic Mol. Wt.: 426 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, DMSO, DMF (up to 5 mg/ml) Storage: -20°C			
1810	N-Octanoyl-D-threo-sphingosine N-C8:0-D- <i>threo</i> -Ceramide C ₂₆ H ₅₁ NO ₃	1 mg	135.00
Source: synthetic Mol. Wt.: 426 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, DMSO, DMF (up to 5 mg/ml) Storage: -20°C			
1333	N-Decanoyl-D-erythro-sphingosine	10 mg	89.00
1333-100	N-C10:0-D- <i>erythro</i> -Ceramide C ₂₈ H ₅₅ NO ₃	100 mg	667.00
Source: synthetic Mol. Wt.: 454 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol, DMSO (up to 5 mg/ml) Storage: -20°C			
2037	N-Pentadecanoyl-D-erythro-sphingosine	10 mg	99.00
2037-100	N-C15:0-D- <i>erythro</i> -Ceramide C ₃₃ H ₆₅ NO ₃	100 mg	637.00
Source: synthetic Mol. Wt.: 524 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, warm ethanol, warm methanol Storage: -20°C			
1915	N-Hexadecanoyl-D-erythro-sphingosine	10 mg	103.00
1915-100	N-C16:0-D- <i>erythro</i> -Ceramide; N-Palmitoyl-D- <i>erythro</i> -sphingosine C ₃₄ H ₆₇ NO ₃ CAS#: 24696-26-2	100 mg	637.00
Source: synthetic Mol. Wt.: 538 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, warm ethanol, warm methanol Storage: -20°C			
2038	N-Heptadecanoyl-D-erythro-sphingosine	10 mg	99.00
2038-100	N-C17:0-D- <i>erythro</i> -Ceramide C ₃₅ H ₆₉ NO ₃ CAS#: 67492-16-4	100 mg	637.00
Source: synthetic Mol. Wt.: 552 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, warm ethanol, warm methanol Storage: -20°C			

1832	N-Octadecanoyl-D-<i>erythro</i>-sphingosine		10 mg	93.00
1832-100	N-C18:0-D- <i>erythro</i> -Ceramide; N-Stearoyl-D- <i>erythro</i> -sphingosine C ₃₆ H ₇₁ NO ₃ CAS#: 2304-81-6		100 mg	637.00
	Source: synthetic Mol. Wt.: 566 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, hot ethanol, chloroform/methanol, 2:1 (up to 5mg/ml) Storage: -20°C			
2039	N-Nonadecanoyl-D-<i>erythro</i>-sphingosine		10 mg	111.00
2039-100	N-C19:0-D- <i>erythro</i> -Ceramide C ₃₇ H ₇₃ NO ₃		100 mg	695.00
	Source: synthetic Mol. Wt.: 580 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, warm ethanol, warm methanol Storage: -20°C			
1843	N-Octadecanoyl-L-<i>threo</i>-sphingosine		1 mg	118.00
	N-C18:0-L- <i>threo</i> -Ceramide; N-Stearoyl-L- <i>threo</i> -sphingosine C ₃₆ H ₇₁ NO ₃			
	Source: synthetic Mol. Wt.: 566 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, DMSO, DMF (up to 5mg/ml) Storage: -20°C			
1850	N-Octadecanoyl-L-<i>erythro</i>-sphingosine		1 mg	118.00
	N-C18:0-L- <i>erythro</i> -Ceramide; N-Stearoyl-L- <i>erythro</i> -sphingosine C ₃₆ H ₇₁ NO ₃			
	Source: synthetic Mol. Wt.: 566 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, DMSO, DMF (up tp 5mg/ml) Storage: -20°C			
1855	N-Octadecanoyl-D-<i>threo</i>-sphingosine		1 mg	135.00
	N-C18:0-D- <i>threo</i> -Ceramide; N-Stearoyl-D- <i>threo</i> -sphingosine C ₃₆ H ₇₁ NO ₃			
	Source: synthetic Mol. Wt.: 566 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, DMSO, DMF (up to 5mg/ml) Storage: -20°C			
1916	N-Tetracosanoyl-D-<i>erythro</i>-sphingosine		5 mg	111.00
1916-25	N-C24:0-D- <i>erythro</i> -Ceramide; N-Lignoceroyl-D- <i>erythro</i> -sphingosine C ₄₂ H ₈₃ NO ₃ CAS#: 34435-05-7		25 mg	475.00
	Source: synthetic Mol. Wt.: 650 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform Storage: -20°C			
1930	N-Tetracosenoyl-D-<i>erythro</i>-sphingosine		5 mg	111.00
1930-25	N-cis-15-C24:1-D- <i>erythro</i> -Ceramide; N-Nervonoyl-D- <i>erythro</i> -sphingosine C ₄₂ H ₈₁ NO ₃ CAS#: 54164-50-0		25 mg	510.00
	Source: synthetic Mol. Wt.: 648 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, DMSO, warm methanol Storage: -20°C			

2-Hydroxy Ceramides

2042	N-(R,S)-alpha-Hydroxydodecanoyl-D-<i>erythro</i>-sphingosine		5 mg	174.00
	N-(R,S)-alpha-Hydroxy-C12:0-D- <i>erythro</i> -ceramide C ₃₀ H ₅₉ NO ₄			
	Source: synthetic Mol. Wt.: 498 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, methanol, ethanol, DMSO Storage: -20°C			

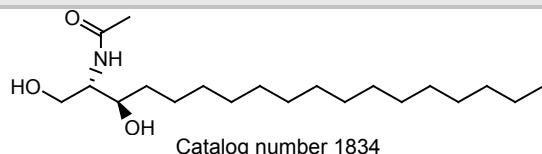
2044	N-(R,S)-alpha-Hydroxyoctadecanoyl-D-erythro-sphingosine N-(R,S)-alpha-Hydroxy-C18:0-D- <i>erythro</i> -ceramide; N-(R,S)-alpha-Hydroxystearoyl-D- <i>erythro</i> -sphingosine C ₃₆ H ₇₁ NO ₄	5 mg	174.00
Source: synthetic Mol. Wt.: 582 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.5 Storage: -20°C			

Ceramide Made from Sphingosines with Sphingoid Bases Other Than C18

1842	N-Acetyl-D-erythro-sphingosine (C14 sphingoid base) N-C2:0 Ceramide of D- <i>erythro</i> -C14-sphingosine C ₁₆ H ₃₁ NO ₃	5 mg	204.00
Source: synthetic Mol. Wt.: 285 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, DMSO, DMF (up to 5 mg/ml) Storage: -20°C			
2077	N-Hexadecanoyl-D-erythro-C16-sphingosine (C16 sphingoidbase) N-Palmitoyl-D- <i>erythro</i> -C16-sphingosine; N-C16:0 Ceramide of D- <i>erythro</i> -C16-sphingosine C ₃₂ H ₆₃ NO ₃	1 mg	264.00

Source: synthetic **Mol. Wt.:** 510 **Purity:** 98% by TLC, MS **Appearance:** solid
Solubility: chloroform, warm ethanol, warm methanol **Storage:** -20°C

Dihydroceramides



1834	N-Acetyl-D-erythro-dihydrosphingosine N-C2:0-D- <i>erythro</i> -Dihydroceramide; N-Acetyl-D- <i>erythro</i> -sphinganine C ₂₀ H ₄₁ NO ₃	5 mg	76.00
Source: synthetic Mol. Wt.: 344 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: -20°C			
1910	N-Hexanoyl-D-erythro-dihydrosphingosine N-C6:0-D- <i>erythro</i> -Dihydroceramide; N-Hexanoyl-D- <i>erythro</i> -sphinganine C ₂₄ H ₄₉ NO ₃	5 mg	76.00

Source: synthetic **Mol. Wt.:** 400 **Purity:** 98+% by TLC, GC **Appearance:** solid
Solubility: chloroform, ethanol, methanol, DMSO **Storage:** -20°C

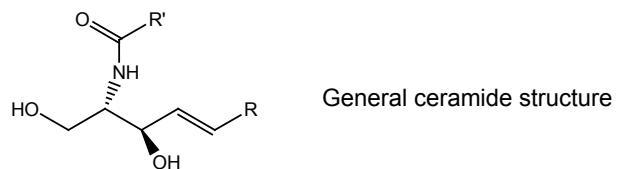
1854	N-Octanoyl-D-erythro-dihydrosphingosine N-C8:0-D- <i>erythro</i> -Dihydroceramide; N-Octanoyl-D- <i>erythro</i> -sphinganine C ₂₆ H ₅₃ NO ₃	5 mg	76.00
Source: synthetic Mol. Wt.: 428 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, DMSO Storage: -20°C			

2041	N-Octadecanoyl-D-erythro-dihydrosphingosine N-C18:0-D- <i>erythro</i> -Dihydroceramide; N-Octadecanoyl-D- <i>erythro</i> -sphinganine; N-Stearoyl-D- <i>erythro</i> -dihydrosphingosine C ₃₆ H ₇₃ NO ₃	10 mg	118.00
Source: synthetic Mol. Wt.: 568 Purity: 98% by TLC Appearance: solid Solubility: hot ethanol, DMSO, warm chloroform/methanol, 5:1 Storage: -20°C			

2-Hydroxy Dihydroceramides

2043	N-(R,S)-alpha-Hydroxydodecanoyl-D-erythro-dihydrosphingosine N-(R,S)- <i>alpha</i> -Hydroxy-C12:0-D- <i>erythro</i> -dihydroceramide C ₃₀ H ₆₁ NO ₄	5 mg	186.00
Source: synthetic Mol. Wt.: 500 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.5 Storage: -20°C			
2045	N-(R,S)-alpha-Hydroxyoctadecanoyl-D-erythro-dihydrosphingosine N-(R,S)- <i>alpha</i> -Hydroxy-C18:0-D- <i>erythro</i> -dihydroceramide; N-(R,S)- <i>alpha</i> -Hydroxystearoyl-D- <i>erythro</i> -dihydrosphingosine C ₃₆ H ₇₃ NO ₄	5 mg	186.00
Source: synthetic Mol. Wt.: 584 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.5 Storage: -20°C			
2047	N-(R,S)-alpha-Hydroxyhexadecanoyl-D-erythro-dihydrosphingosine N-(R,S)- <i>alpha</i> -Hydroxy-C16:0-D- <i>erythro</i> -dihydroceramide; N-(R,S)- <i>alpha</i> -Hydroxypalmitoyl-D- <i>erythro</i> -dihydrosphingosine C ₃₄ H ₆₉ NO ₄	5 mg	115.00
Source: synthetic Mol. Wt.: 556 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.5 Storage: -20°C			

Ceramides From Natural Sources



1056	Ceramides Ceramides with hydroxy and non-hydroxy acyl groups C ₄₂ H ₈₃ NO ₄ CAS#: 104404-17-13	25 mg	103.00
Source: natural, bovine Mol. Wt.: 666(2-hydroxylignoceroyl) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C			
1322	Ceramides	10 mg	77.00
1322-05	Ceramides with mostly non-hydroxy acyl groups C ₃₆ H ₇₁ NO ₃	50 mg	259.00

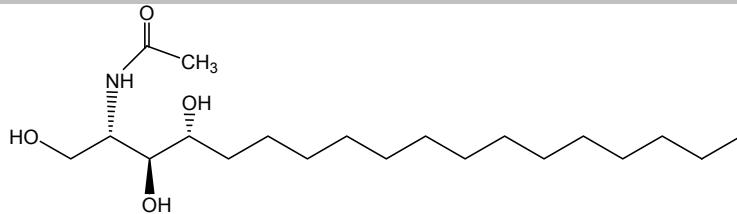
Source: natural, bovine **Mol. Wt.:** 566 (stearoyl) **Purity:** 98+% by TLC
Appearance: solid **Solubility:** ethanol, chloroform/methanol, 2:1 **Storage:** -20°C

1323	Ceramides		10 mg	77.00
1323-05	Ceramides with mostly hydroxy acyl groups	$C_{36}H_{71}NO_4$	50 mg	259.00

Source: natural, bovine **Mol. Wt.:** 582 (2-hydroxystearoyl) **Purity:** 98+% by TLC
Appearance: solid **Solubility:** methanol, chloroform/methanol, 2:1 **Storage:** -20°C

See Table III (pg. 93-97) for typical fatty acid content of products prepared from natural sources.

Phytoceramides



1897	N-Acetyl-phytosphingosine		5 mg	103.00
	N-C2:0-Phytoceramide	$C_{20}H_{41}NO_4$		

Source: semisynthetic, yeast (*Pichia ciferri*) **Mol. Wt.:** 360 **Purity:** 98+% by TLC,
GC Appearance: solid **Solubility:** ethanol, methanol, warm DMSO,
chloroform/methanol, 1:1 (warm) **Storage:** -20°C

1895	N-Hexanoyl-phytosphingosine		5 mg	103.00
	N-C6:0-Phytoceramide	$C_{24}H_{49}NO_4$		

Source: semisynthetic, yeast (*Pichia ciferri*) **Mol. Wt.:** 416 **Purity:** 98+% by TLC,
GC Appearance: solid **Solubility:** ethanol, methanol, chloroform/methanol, 1:1
(warm) **Storage:** -20°C

1894	N-Octanoyl-phytosphingosine		5 mg	103.00
	N-C8:0-Phytoceramide	$C_{26}H_{53}NO_4$		

Source: semisynthetic, yeast (*Pichia ciferri*) **Mol. Wt.:** 444 **Purity:** 98+% by TLC,
GC Appearance: solid **Solubility:** ethanol, methanol, chloroform/methanol, 1:1
(warm) **Storage:** -20°C

2035	N-Hexadecanoyl-phytosphingosine		5 mg	99.00
	N-C16:0-Phytoceramide; N-Palmitoyl-phytosphingosine	$C_{34}H_{69}NO_4$		

Source: semisynthetic, yeast (*Pichia ciferri*) **Mol. Wt.:** 556 **Purity:** 98+% by TLC,
GC Appearance: solid **Solubility:** chloroform/methanol, 5:1 **Storage:** -20°C

2034	N-Octadecanoyl-phytosphingosine		5 mg	99.00
	N-C18:0-Phytoceramide; N-Stearoyl-phytosphingosine	$C_{36}H_{73}NO_4$		

Source: semisynthetic, yeast (*Pichia ciferri*) **Mol. Wt.:** 584 **Purity:** 98+% by TLC,
MS Appearance: solid **Solubility:** chloroform/methanol, 1:1 (warm)
Storage: -20°C

2036	N-Tetracosanoyl-phytosphingosine N-C24:0-Phytoceramide; N-Lignoceroyl-phytosphingosine	5 mg	126.00
	C ₄₂ H ₈₅ NO ₄		

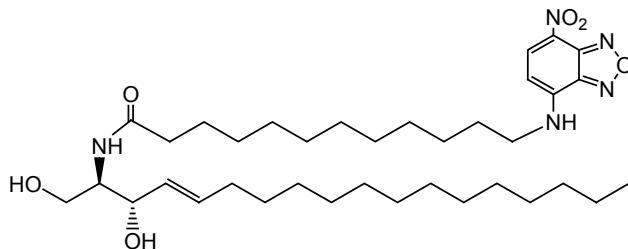
Source: semisynthetic, yeast (*Pichia ciferrri*) **Mol. Wt.:** 668 **Purity:** 98+% by TLC,
MS Appearance: solid **Solubility:** chloroform/methanol, 5:1 **Storage:** -20°C

Fluorescent Ceramides

1841	N-Hexanoyl-NBD-D-<i>erythro</i>-sphingosine	100 µg	204.00
1841-001	N-C6:0-NBD-Ceramide; N-C6:0-NBD-D- <i>erythro</i> -Sphingosine	1 mg	302.00

CAS#: 86701-10-2

Source: synthetic **Mol. Wt.:** 576 **Melting Point (°C):** 85.7-87.9 **Purity:** 98+% by TLC
Appearance: solid **Solubility:** chloroform, ethanol, methanol
Storage: -20°C



Catalog number 1618

Excitation: 460 nm
Emission: 535 nm

1618	N-Dodecanoyl-NBD-D-<i>erythro</i>-sphingosine	100 µg	147.00
1618-001	N-C12:0-NBD-Ceramide; N-C12:0-NBD-D- <i>erythro</i> -Sphingosine	1 mg	526.00

C₃₆H₆₁N₅O₆

Source: synthetic **Mol. Wt.:** 660 **Purity:** 98+% by TLC **Appearance:** solid
Solubility: methanol, chloroform/methanol, 2:1 **Storage:** -20°C

1857	N-Hexanoyl-NBD-L-<i>threo</i>-sphingosine	100 µg	162.00
1857-001	N-C6:0-NBD-Ceramide; N-C6:0-NBD-L- <i>threo</i> -Sphingosine	1 mg	526.00

Source: synthetic **Mol. Wt.:** 576 **Purity:** 98+% by TLC **Appearance:** solid
Solubility: chloroform, ethanol, methanol **Storage:** -20°C

1620	N-Dodecanoyl-NBD-L-<i>threo</i>-sphingosine	100 µg	162.00
1620-001	N-C12:0-NBD-Ceramide; N-C12:0-NBD-L- <i>threo</i> -Sphingosine	1 mg	526.00

Source: synthetic **Mol. Wt.:** 660 **Purity:** 98+% by TLC **Appearance:** solid
Solubility: methanol, chloroform/methanol, 2:1 **Storage:** -20°C

1624	N-Hexanoyl-NBD-L-<i>threo</i>-dihydrosphingosine	100 µg	147.00
1624-001	N-C6:0-NBD-Dihydroceramide; N-C6:0-NBD-L- <i>threo</i> -Dihydrosphingosine	1 mg	526.00

C₃₀H₅₁N₅O₆

Source: synthetic **Mol. Wt.:** 578 **Purity:** 98+% by TLC **Appearance:** solid
Solubility: methanol, chloroform/methanol, 2:1 **Storage:** -20°C

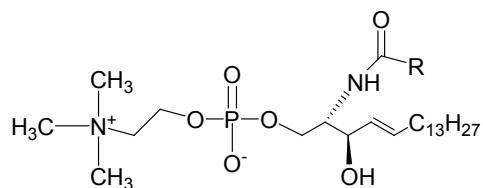
1623	N-Dodecanoyl-NBD-L-<i>threo</i>-dihydrosphingosine		100 µg	147.00
1623-001	N-C12:0-NBD-Dihydroceramide; N-C12:0-NBD-L- <i>threo</i> -Dihydrosphingosine C ₃₆ H ₆₃ N ₅ O ₆		1 mg	526.00
	Source: synthetic Mol. Wt.: 662 Purity: 98+% by TLC Appearance: solid Solubility: methanol, chloroform/methanol, 2:1 Storage: -20°C			
1626	N-Hexanoyl-NBD-D-<i>erythro</i>-dihydrosphingosine		100 µg	126.00
1626-001	N-C6:0-NBD-Dihydroceramide; N-C6:0-NBD-D- <i>erythro</i> -Dihydrosphingosine C ₃₀ H ₅₁ N ₅ O ₆		1 mg	526.00
	Source: synthetic Mol. Wt.: 578 Purity: 98+% by TLC Appearance: solid Solubility: methanol, chloroform/methanol, 2:1 Storage: -20°C			
1625	N-Dodecanoyl-NBD-D-<i>erythro</i>-dihydrosphingosine		100 µg	147.00
1625-001	N-C12:0-NBD-Dihydroceramide; N-C12:0-NBD-D- <i>erythro</i> -Dihydrosphingosine C ₃₆ H ₆₃ N ₅ O ₆		1 mg	526.00
	Source: synthetic Mol. Wt.: 662 Purity: 98+% by TLC Appearance: solid Solubility: methanol, chloroform/methanol, 2:1 Storage: -20°C			
1628	N-Hexanoyl-NBD-phytosphingosine		100 µg	126.00
1628-001	N-C6:0-NBD-Phytoceramide; N-C6:0-NBD-Phytosphingosine C ₃₀ H ₅₁ N ₅ O ₇		1 mg	526.00
	Source: semisynthetic, bacteria Mol. Wt.: 594 Purity: 98+% by TLC Appearance: solid Solubility: methanol, chloroform/methanol, 2:1 Storage: -20°C			
1627	N-Dodecanoyl-NBD-phytosphingosine		100 µg	133.00
1627-001	N-C12:0-NBD-Phytoceramide; N-C12:0-NBD-Phytosphingosine C ₃₆ H ₆₃ N ₅ O ₇		1 mg	526.00
	Source: semisynthetic, bacteria Mol. Wt.: 678 Purity: 98+% by TLC Appearance: solid Solubility: methanol, chloroform/methanol, 2:1 Storage: -20°C			

See Biochemicals and Reagents section (page 88) for additional fluorescent labeled products.

Compounds with fluorescent labels other than NBD are available on custom basis. Contact Technical Service for more information.

Phosphosphingolipids

Sphingomyelins



1051	Sphingomyelin		25 mg	47.00
1051-1	SPM; Ceramide-1-phosphorylcholine	C ₄₁ H ₈₃ N ₂ O ₆ P CAS#: 85187-10-6	1 g	522.00

Source: natural, bovine **Mol. Wt.:** 731 (stearoyl) **Purity:** 98+% by TLC
Appearance: solid **Solubility:** chloroform, ethanol **Storage:** -20°C

See Table III (pg. 93-97) for fatty acid content

1328	Sphingomyelin		25 mg	47.00
	SPM; Ceramide-1-phosphorylcholine	C ₄₇ H ₉₅ N ₂ O ₆ P CAS#: 85187-10-6		

Source: natural, porcine RBC **Mol. Wt.:** 815 (lignoceroyl) **Purity:** 98+% by TLC
Appearance: solid **Solubility:** chloroform, ethanol **Storage:** -20°C

See Table III (pg. 93-97) for fatty acid content

1329	Sphingomyelin		25 mg	47.00
1329-1	SPM; Ceramide-1-phosphorylcholine	C ₄₆ H ₉₃ N ₂ O ₆ P CAS#: 85187-10-6	1 gram	474.00

Source: natural, bovine buttermilk **Mol. Wt.:** 801 (tricosanoyl) **Purity:** 98+% by TLC
Appearance: solid **Solubility:** chloroform, ethanol **Storage:** -20°C

See Table III (pg. 93-97) for fatty acid content

1332	Sphingomyelin		25 mg	47.00
1332-1	SPM; Ceramide-1-phosphorylcholine	C ₃₉ H ₇₉ N ₂ O ₆ P	1 gram	522.00

Source: natural, chicken, egg **Mol. Wt.:** 703 (palmitoyl) **Purity:** 98+% by TLC
Appearance: solid **Solubility:** chloroform, methanol, warm ethanol **Storage:** -20°C

See Table III (pg. 93-97) for fatty acid content

1907	N-Acetyl-sphingosylphosphorylcholine		5 mg	184.00
	N-C2:0-Sphingomyelin	C ₂₅ H ₅₁ N ₂ O ₆ P		

Source: semisynthetic, bovine buttermilk **Mol. Wt.:** 506 **Purity:** 98+% by TLC
Appearance: solid **Solubility:** ethanol, chloroform/methanol, 2:1 **Storage:** -20°C

Mixture of D-erythro and L-threo isomers

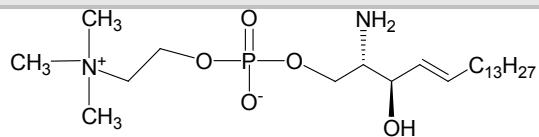
1909	N-Hexanoyl-sphingosylphosphorylcholine		5 mg	184.00
	N-C6:0-Sphingomyelin	C ₂₉ H ₅₉ N ₂ O ₆ P		

Source: semisynthetic, bovine buttermilk **Mol. Wt.:** 563 **Purity:** 98+% by TLC
Appearance: solid **Solubility:** ethanol, chloroform/methanol, 2:1 **Storage:** -20°C

Mixture of D-erythro and L-threo isomers

1911	N-Octadecanoyl-sphingosylphosphorylcholine N-C18:0-Sphingomyelin; N-Stearoyl-sphingosylphosphorylcholine $C_{41}H_{83}N_2O_6P$	5 mg	184.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 731 Purity: 98+% by TLC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: -20°C		
	Mixture of D- <i>erythro</i> and L- <i>threo</i> isomers		
1890	N-Heptadecanoyl-sphingosylphosphorylcholine N-C17:0-Sphingomyelin $C_{40}H_{81}N_2O_6P$	5 mg	184.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 717 Purity: 98+% by TLC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: -20°C		
	Mixture of D- <i>erythro</i> and L- <i>threo</i> isomers		
1917	N-Eicosanoyl-D-<i>erythro</i>-sphingosylphosphorylcholine N-C20:0-Sphingomyelin $C_{43}H_{87}N_2O_6P$	0.5 mg	184.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 759 Purity: 98+% by TLC Appearance: solid Solubility: ethanol, methanol, chloroform/methanol, 14:1 Storage: -20°C		
<hr/>			
<p>Catalog number 1918</p>			
1918	N-Docosanoyl-D-<i>erythro</i>-sphingosylphosphorylcholine N-C22:0-Sphingomyelin $C_{45}H_{91}N_2O_6P$	0.5 mg	184.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 787 Purity: 98+% by TLC Appearance: solid Solubility: ethanol, methanol, chloroform/methanol, 14:1 Storage: -20°C		
2200	N-1-¹³C-Hexadecanoyl-sphingosylphosphorylcholine D- <i>erythro</i> -Sphingomyelin with 1- ¹³ C-palmitic acid; N-1- ¹³ C-Palmitoyl-sphingosylphosphorylcholine $^{12}C_{38}^{13}CH_{79}N_2O_6P$	1mg	184.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 703 Purity: 98+% by TLC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: -20°C		
1327	N-Acyl-sphingosylphosphorylethanolamine Ceramide phosphorylethanolamine $C_{43}H_{87}N_2O_6P$	5 mg	154.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 773 (tricosanoyl) Purity: 98+%		
	by TLC Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C		

Sphingosylphosphorylcholines (SPC)



Catalog number 1318

1318	D-erythro-Sphingosylphosphorylcholine D- <i>erythro</i> -SPC $C_{23}H_{49}N_2O_5P$	5 mg	254.00
------	--	------	--------

Source: semisynthetic, bovine buttermilk **Mol. Wt.:** 465 **Purity:** 98+% by TLC
Appearance: solid **Solubility:** chloroform/methanol, 2:1 **Storage:** -20°C

1319	L-threo-Sphingosylphosphorylcholine L- <i>threo</i> -SPC $C_{23}H_{49}N_2O_5P$	5 mg	254.00
------	--	------	--------

Source: semisynthetic, bovine buttermilk **Mol. Wt.:** 465 **Purity:** 98+% by TLC
Appearance: solid **Solubility:** chloroform/methanol, 2:1 **Storage:** -20°C

1321	Sphingosylphosphorylcholine	10 mg	234.00
1321-05	<i>lyso</i> -Sphingomyelin; SPC (mixture of D- <i>erythro</i> and L- <i>threo</i> isomers) $C_{23}H_{49}N_2O_5P$ CAS#: 82970-80-7	50 mg	926.00

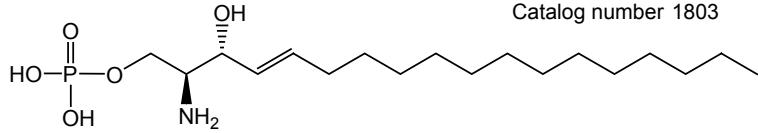
Source: semisynthetic, bovine buttermilk **Mol. Wt.:** 465 **Purity:** 98+% by TLC
Appearance: solid **Solubility:** chloroform/methanol, 2:1 **Storage:** -20°C

1913	<i>lyso</i>-Dihydrophingomyelin <i>Dihydro</i> sphingosylphosphorylcholine (mixture of D- <i>erythro</i> and L- <i>threo</i> isomers) $C_{23}H_{51}N_2O_5P$	1 mg	133.00
------	---	------	--------

Source: semisynthetic, bovine buttermilk **Mol. Wt.:** 467 **Purity:** 98+% by TLC
Appearance: solid **Solubility:** chloroform/methanol, 2:1 **Storage:** -20°C

Sphingosine Phosphates

Catalog number 1803

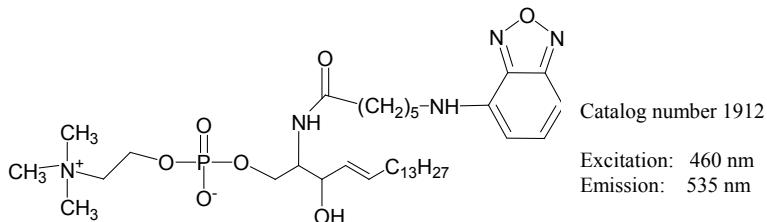


1803	D-erythro-Sphingosine-1-phosphate S-1-P $C_{18}H_{38}NO_5P$ CAS#: 26993-30-6	5 mg	278.00
------	--	------	--------

Source: synthetic **Mol. Wt.:** 380 **Purity:** 98+% by TLC **Appearance:** solid
Solubility: chloroform plus a few drops of TFA, chloroform/methanol/40% dimethylamine, 5:15:3, 1mg/ml **Storage:** -20°C

1852	D-erythro-Dihydrosphingosine-1-phosphate C ₁₈ H ₄₀ NO ₅ P CAS#: 19794-97-9	5 mg	278.00
	Source: synthetic Mol. Wt.: 382 Purity: 98+% by TLC Appearance: solid Solubility: chloroform plus a few drops of TFA, chloroform/methanol/40% dimethylamine, 5:15:3, 1mg/ml Storage: -20°C		
2046	N-Hexadecanoyl-D-erythro-sphingosine-1-phosphate, NH₄⁺ salt N-C16:0-Ceramide-1-phosphate C ₃₄ H ₆₈ NO ₆ P•NH ₃	5 mg	377.00
	Source: synthetic Mol. Wt.: 618+NH ₃ Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/acetic acid, 60:15:25 Storage: -20°C		

Fluorescent Sphingomyelins



1912	N-Hexanoyl-NBD-sphingosylphosphorylcholine	100 µg	111.00
1912-001	N-C6:0-NBD-Sphingomyelin; N-C6:0-NBD-Sphingosylphosphorylcholine C ₃₅ H ₆₁ N ₆ O ₉ P CAS#: 94885-04-8	1 mg	290.00

Source: semisynthetic, bovine buttermilk **Mol. Wt.:** 740 **Purity:** 98+% by TLC
Appearance: solid **Solubility:** chloroform, ethanol, methanol **Storage:** -20°C

Mixture of D-*erythro* and L-*threo* isomers

1619	N-Dodecanoyl-NBD-sphingosylphosphorylcholine	100 µg	147.00
1619-001	N-C12:0-NBD-Sphingomyelin; N-C12:0-NBD-Sphingosylphosphorylcholine C ₄₁ H ₇₃ N ₆ O ₉ P	1 mg	319.00

Source: semisynthetic, bovine buttermilk **Mol. Wt.:** 825 **Purity:** 98+% by TLC
Appearance: solid **Solubility:** methanol, chloroform/methanol, 2:1 **Storage:** -20°C

Mixture of D-*erythro* and L-*threo* isomers

See Biochemicals and Reagents section (page 88) for additional fluorescent labeled products.

Compounds with fluorescent labels other than NBD are available on custom basis. Contact Technical Service for more information.

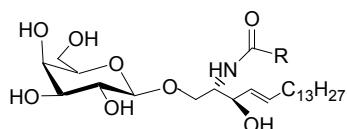
Glycosphingolipids

Glycosphingolipids are widely distributed in animal and plant tissues. They consist of a ceramide (Cer) bound in glycosidic linkage through the primary hydroxyl to a mono- or oligosaccharide which may contain substituents such as a sulfate, acetate, or phosphate group. They are amphiphilic and the less glycosylated compounds can be dispersed in buffers by dissolving them in a detergent or organic solvent (EtOH, DMSO, isoPrOH) and mixing by sonication.

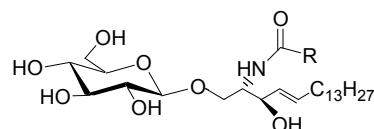
Galactosphingolipids, mainly GalCer (cerebrosides) and its sulfate ester, occur in large amounts in the nervous system. Glucosphingolipids, the simplest of which is GlcCer (glucocerebrosides), are very widely distributed, particularly in nerve cell membranes. GlcCer is isolated from a variety of natural sources including human, bovine, and plant. Each of these sources has a heterogeneity in the fatty acid content of the ceramide as well as an occasional variation in the sphingoid chain. Globosides (containing both glucose and galactose) are a prominent group of glycosphingolipids, they contain an α -linked galactose moiety and are typically located in blood cell membranes. Gangliosides are another prominent group of glycosphingolipids; they are acidic because of substitution with sialic (neuraminic) acid. The glycosphingolipids function in a wide range of enzyme and structural interactions, such as immunological or membrane recognition phenomena, binding of microbial pathogens, hormone and growth factor actions, cancer cell growth and malignancy, atherosclerosis, genetic disease errors, blood group determinants, etc. Tissues change in glycosphingolipid composition during embryogenesis, maturation, aging, and other vital physiological processes. Some glycosphingolipids stimulate cell proliferation, others induce apoptosis, effects of great significance to cancer therapy and maturational development. Marked differences in glycosphingolipid composition are seen in normal and cancerous cells. See references (13-25).

See Literature References on page 99.

Galactosylceramides and Glucosylceramides



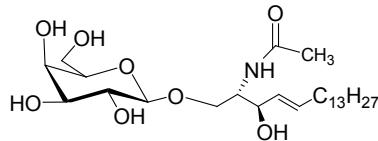
Galactosylceramide



Glucosylceramide

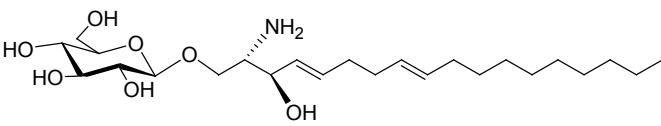
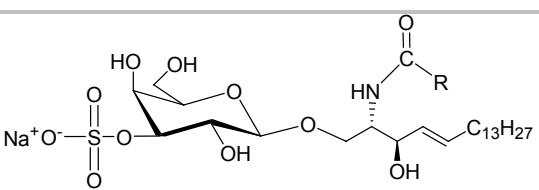
1050	Cerebrosides Galactosylceramide; Ceramide <i>beta</i> -D-galactoside C ₄₈ H ₉₃ NO ₉ CAS#: 85305-88-0	50 mg	89.00
	Source: natural, bovine Mol. Wt.: 828 (2-hydroxytetraicosanoyl) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C		
	See Table III (pg. 93-97) for fatty acid content		
1066	Cerebroside; Kerasin (top spot) Galactosylceramide with mostly non-hydroxy fatty acid side chain C ₄₂ H ₈₁ NO ₈ CAS#: 536-13-0	10 mg	111.00
	Source: natural, bovine Mol. Wt.: 810 (nervonyl, [24:1]) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.5 Storage: -20°C		
	See Table III (pg. 93-97) for fatty acid content		

1138	Cerebroside; Phrenosin (bottom spot) Galactosylceramide with mostly 2-hydroxy fatty acid side chains C ₄₂ H ₈₁ NO ₉ CAS#: 37211-11-3	10 mg	111.00
	Source: natural, bovine Mol. Wt.: 744 (2-hydroxystearoyl) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.5 Storage: -20°C		
See Table III (pg. 93-97) for fatty acid content			
1305	Psychosine (free amine form) lyso-Cerebroside; 1- <i>beta</i> -D-Galactosylsphingosine C ₂₄ H ₄₇ NO ₇ CAS#: 2238-90-6	10 mg	184.00
	Source: semisynthetic, bovine Mol. Wt.: 461 Purity: 98+% by TLC Appearance: solid Solubility: ethanol, chloroform/methanol/DI water, 5:1:0.1 Storage: -20°C		
1914	N-Octadecanoyl-D₃₅-psychosine (perdeuterated C18:0 fatty acid) N-C18:0-D ₃₅ -Cerebroside, perdeuterated; N-Stearoyl-D ₃₅ -Psychosine, perdeuterated C ₄₂ H ₄₆ D ₃₅ NO ₈	5 mg	306.00
	Source: semisynthetic, bovine Mol. Wt.: 762 Purity: 98+% by TLC Appearance: solid Solubility: chloroform, hot ethanol, chloroform/methanol, 2:1 Storage: -20°C		
	Deuterium labeled stearoyl sidechain		
1325	N-Acetyl-psychosine N-C2:0-Cerebroside C ₂₆ H ₄₉ NO ₈	10 mg	154.00
	Source: semisynthetic, bovine Mol. Wt.: 503 Purity: 98+% by TLC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: -20°C		
1335	N-Pentadecanoyl-psychosine N-C15:0-Cerebroside C ₃₉ H ₇₅ NO ₈	5 mg	111.00
	Source: semisynthetic, bovine Mol. Wt.: 686 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/ methanol, 2:1 Storage: -20°C		
1334 1334-50	N-Octanoyl-beta-D-galactosylceramide N-C8:0-Galactosylceramide C ₃₂ H ₆₁ NO ₈	10 mg 50 mg	271.00 942.00
	Source: semisynthetic, bovine Mol. Wt.: 588 Purity: 98+% by TLC Appearance: solid Solubility: ethanol, methanol, chloroform/methanol, 9:1 Storage: -20°C		



Catalog number 1325

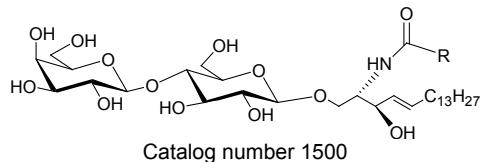
1621	N-Hexanoyl-NBD-galactosylceramide		100 µg	147.00
1621-001	N-C6:0-NBD- <i>beta</i> -D-Galactosylsphingosine; N-C6:0-NBD-Cerebroside C ₃₆ H ₅₉ N ₅ O ₁₁		1 mg	526.00
Source: semisynthetic, bovine Mol. Wt.: 738 Purity: 98+% by TLC Appearance: solid Solubility: methanol, chloroform/methanol, 5:1 Storage: -20°C				
1633	N-Dodecanoyl-NBD-galactosylceramide		100 µg	118.00
1633-001	N-C12:0-NBD- <i>beta</i> -D-Galactosylsphingosine; N-C12:0-NBD-Cerebroside C ₄₂ H ₇₁ N ₅ O ₁₁		1 mg	526.00
Source: semisynthetic, bovine Mol. Wt.: 822 Purity: 98+% by TLC Appearance: solid Solubility: chloroform, DMSO, chloroform/methanol, 2:1 Storage: -20°C				
1057	Glucocerebrosides, Gaucher's spleen		5 mg	204.00
1057-25	Glucosylceramide; Ceramide <i>beta</i> -D-glucoside C ₄₈ H ₉₃ NO ₈ CAS# 85305-87-9		25 mg	753.00
Source: natural, human Mol. Wt.: 812 (lignoceryl) Purity: 98+% by TLC, GC, MS Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C				
See Table III (pg. 93-97) for fatty acid content				
1521	Glucocerebrosides		5 mg	133.00
1521-50	Glucosylceramide; Ceramide <i>beta</i> -D-glucoside C ₄₆ H ₈₉ NO ₈		50 mg	1,013.00
Source: natural, bovine buttermilk Mol. Wt.: 784 (docosanoyl) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C				
See Table III (pg. 93-97) for side chain variants				
<p style="text-align: center;">Catalog number 1522</p>				
1522	Glucocerebrosides, plant		5 mg	60.00
1522-100	Glucosylceramide; Ceramide <i>beta</i> -D-glucoside C ₄₀ H ₇₅ NO ₉		100 mg	826.00
Source: natural, plant Mol. Wt.: 714 (2-hydroxyhexadecanoyl) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C				
Sphingoid backbone is >95% 4,8-sphingadiene (C18:2 t,t-4,8) and most of the fatty acids are of the 2-hydroxy type. See Table III (pg. 93-97) for fatty acid content				
1622	N-Hexanoyl-NBD-glucosylceramide		100 µg	147.00
1622-001	N-C6:0-NBD- <i>beta</i> -D-Glucosylsphingosine; N-C6:0-NBD-Glucosylceramide C ₃₆ H ₅₉ N ₅ O ₁₁		1 mg	526.00
Source: semisynthetic, bovine Mol. Wt.: 738 Purity: 98+% by TLC Appearance: solid Solubility: methanol, chloroform/methanol, 5:1 Storage: -20°C				

1306	Glucospsychosine Glucosylsphingosine; <i>lyso</i> -Glucocerebroside; 1- <i>beta</i> -D-Glucosylsphingosine C ₂₄ H ₄₇ NO ₇ CAS#: 52050-17-6	5 mg	254.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 461 Purity: 98+% by TLC Appearance: solid Solubility: ethanol, methanol, chloroform/methanol, 2:1 Storage: -20°C		
1310		Catalog number 1310	
1310	Glucospsychosine Glucosylsphingosine; <i>lyso</i> -Glucocerebroside; 1- <i>beta</i> -D-Glucosylsphingadienine C ₂₄ H ₄₅ NO ₇ CAS#: 52050-17-6	5 mg	184.00
	Source: natural, plant Mol. Wt.: 460 (based on 1- <i>beta</i> -D-glucosylsphinga-4,8-dienine) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 4:1 Storage: -20°C		
	Sphingoid backbone is >95% 4,8-sphingadiene (C18:2 t,t-4, 8)		
1531	N-Docosanoyl-glucospsychosine N-C22:0-Glucocerebroside; N-Docosanoyl- <i>beta</i> -glucosylsphingosine C ₄₆ H ₈₉ NO ₈	1 mg	306.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 784 Purity: 98+% by TLC Appearance: solid Solubility: chloroform Storage: -20°C		
1533	N-Hexadecanoyl-D₃-glucospsychosine, deuterated N-C16:0-D ₃ -Glucospsychosine, deuterated; N-C16:0-D ₃ -Glucocerebroside, deuterated; N-Palmitoyl-D ₃ -glucospsychosine, deuterated C ₄₀ H ₇₄ D ₃ NO ₈	1 mg	290.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 703 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C		
1049		Catalog number 1049	
1049	Sulfatides Ceramide-galactoside-3-sulfate; Cerebroside sulfate C ₄₂ H ₈₀ NNaO ₁₁ S CAS#: 85496-63-5	50 mg	377.00
	Source: natural, bovine Mol. Wt.: 830 (stearoyl) Na ⁺ Salt Purity: 98+% by TLC Appearance: solid Solubility: DMSO, chloroform/methanol/DI water, 2:1:0.1 (if needed, a few drops of acetic acid) Storage: -20°C		
	See Table III (pg. 93-97) for fatty acid content		

1904	Iyso-Sulfatide (NH_4^+ salt) Sphingosine-1-galactoside-3-sulfate $\text{C}_{24}\text{H}_{47}\text{NO}_{10}\text{S} \cdot \text{NH}_3$ CAS#: 38621-58-8	1 mg	429.00
	Source: semisynthetic, bovine Mol. Wt.: 542 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C		
2076	N-Acetyl-sulfatide N-C2:0-Sulfatide; N-Acetyl-sphingosyl- <i>beta</i> -D-galactoside-3-sulfatide $\text{C}_{26}\text{H}_{49}\text{NO}_{11}\text{S}$	1 mg	290.00
	Source: semisynthetic, bovine Mol. Wt.: 584 Purity: 98+% by TLC Appearance: solid Solubility: ethanol, methanol, chloroform/methanol, 1:1 Storage: -20°C		
1875	N-Hexadecanoyl-sulfatide N-C16:0-Sulfatide; N-Palmitoyl-sulfatide; N-Palmitoyl-sphingosyl- <i>beta</i> -D-galactoside-3-sulfate $\text{C}_{40}\text{H}_{77}\text{NO}_{11}\text{S}$	1 mg	290.00
	Source: semisynthetic, bovine Mol. Wt.: 780 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C		
1932	N-Octadecanoyl-sulfatide N-C18:0-Sulfatide; N-Octadecanoyl-sphingosyl- <i>beta</i> -D-galactoside-3-sulfate $\text{C}_{42}\text{H}_{81}\text{NO}_{11}\text{S}$	1 mg	319.00
	Source: semisynthetic, bovine Mol. Wt.: 808 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 5:1 Storage: -20°C		
1933	N-Octadecenoyl-sulfatide N-C18:1-Sulfatide; N-Octadecenoyl-sphingosyl- <i>beta</i> -D-galactoside-3-sulfate $\text{C}_{42}\text{H}_{79}\text{NO}_{11}\text{S}$	1 mg	319.00
	Source: semisynthetic, bovine Mol. Wt.: 806 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 5:1 Storage: -20°C		
1888	N-Tetracosanoyl sulfatide N-C24:0-Sulfatide; N-Tetracosanoyl-sphingosyl- <i>beta</i> -D-galactoside-3-sulfate; N-Lignoceroyl-sulfatide $\text{C}_{48}\text{H}_{93}\text{NO}_{11}\text{S}$	1 mg	319.00
	Source: semisynthetic, bovine Mol. Wt.: 892 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 5:1 Storage: -20°C		
1931	N-Tetracosenoyl-sulfatide N-Nervonyl-sulfatide; N-C24:1-Sulfatide; N-Tetracosenoyl-sphingosyl- <i>beta</i> -D-galactoside-3-sulfate $\text{C}_{48}\text{H}_{91}\text{NO}_{11}\text{S}$	1 mg	319.00
	Source: semisynthetic, bovine Mol. Wt.: 890 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 5:1 Storage: -20°C		
1536	N-Octadecanoyl-D₃-sulfatide, deuterated N-C18:0-D ₃ -Sulfatide, deuterated; N-Stearoyl-D ₃ -sulfatide, deuterated $\text{C}_{42}\text{H}_{78}\text{D}_3\text{NO}_{11}\text{S}$	1 mg	414.00
	Source: semisynthetic, bovine Mol. Wt.: 811 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1 Storage: -20°C		

1632	N-Dodecanoyl-NBD-sulfatide		100 µg	119.00
1632-001	N-C12:0-NBD-Sulfatide; N-Dodecanoyl-NBD- <i>lys</i> -sulfatide; N-Dodecanoyl-NBD-sphingosyl- <i>beta</i> -D-galactoside-3-sulfate C ₄₂ H ₇₁ N ₅ O ₁₄ S		1 mg	526.00
	Source: semisynthetic, bovine Mol. Wt.: 901 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C			

Lactosylceramides



1500	Lactosylceramides		1 mg	147.00
	LC; Lactocerebrosides; CDH; Ceramide <i>beta</i> -lactoside C ₄₈ H ₉₁ NO ₁₃ CAS#: 4682-48-8			
	Source: natural, porcine RBC Mol. Wt.: 890 (stearoyl) Purity: 98+% by TLC Appearance: solid Solubility: DMSO, chloroform/methanol/DI water, 5:1:0.1 Storage: -20°C			

See Table III (pg. 93-97) for fatty acid content

1507	Lactosylceramides		5 mg	204.00
1507-50	LC; Lactocerebrosides; CDH; Ceramide <i>beta</i> -lactoside C ₅₃ H ₁₀₁ NO ₁₃ CAS#: 4682-48-8		50 mg	1,390.00
	Source: natural, bovine buttermilk Mol. Wt.: 960 (tricosanoyl) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 5:1:0.1 Storage: -20°C			

See Table III (pg. 93-97) for fatty acid content

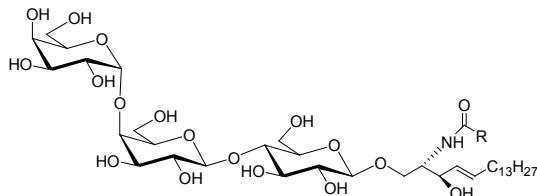
1517	<i>lys</i>-Lactosylceramide		1 mg	186.00
	Lactosylsphingosine; <i>lys</i> -LC C ₃₀ H ₅₇ NO ₁₂			
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 623 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1 Storage: -20°C			

1532	N-Hexadecanoyl-lactosylceramide		1 mg	134.00
	N-C16:0-Lactosylceramide; N-Palmitoyl-lactosylceramide C ₄₆ H ₈₇ NO ₁₃			
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 862 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1 Storage: -20°C			

1534	N-Hexadecanoyl-D₃-lactosylceramide, deuterated		1 mg	376.00
	N-C16:0-D ₃ -Lactosylceramide, deuterated; N-Palmitoyl-D3-lactosylceramide, deuterated C ₄₆ H ₈₄ D ₃ NO ₁₃			
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 865 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 5:1:0.1 Storage: -20°C			

1538	N-Heptadecanoyl-lactosylceramide N-C17:0-Lactosylceramide; Lactosylceramide with C17:0 fatty acid side chain $C_{47}H_{89}NO_{13}$	1 mg	167.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 876 Purity: 98+% by TLC, MS Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1 Storage: -20°C		
1629	N-Hexanoyl-NBD-lactosylceramide	50 ug	184.00
1629-001	N-Hexanoyl-NBD-beta-D-lactosylphingosine; N-C6:0-NBD-beta-D-Lactosylphingosine; N-C6:0-NBD-Lactosylceramide $C_{42}H_{69}N_5O_{16}$	1 mg	824.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 900 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C		
1630	N-Dodecanoyl-NBD-lactosylceramide	50 ug	204.00
1630-001	N-Dodecanoyl-NBD-beta-D-lactosylphingosine; N-C12:0-NBD-beta-D-Lactosylphingosine; N-C12:0-NBD-Lactosylceramide $C_{48}H_{81}N_5O_{16}$	1 mg	793.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 984 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C		

Ceramide Trihexosides



Catalog number 1067

1067	Ceramide trihexosides	1 mg	285.00
1067-10	CTH; Gb3; Globotriaosylceramide $C_{60}H_{113}NO_{18}$ CAS#: 71965-57-6	10 mg	2,242.00

Source: natural, porcine RBC **Mol. Wt.:** 1137 (tetracosanoyl) **Purity:** 98+% by TLC
Appearance: solid **Solubility:** DMSO, hot methanol, chloroform/methanol, 2:1
Storage: -20°C

See Table III (pg. 93-97) for fatty acid content

1513	Ceramide trihexosides (top spot)	0.5 mg	204.00
	CTH with non-hydroxy fatty acid side chain $C_{54}H_{101}NO_{18}$		

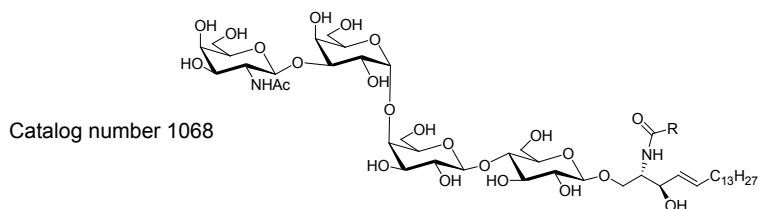
Source: natural, porcine RBC **Mol. Wt.:** 1052 (stearoyl) **Purity:** 98+% by TLC
Appearance: solid **Solubility:** chloroform/methanol, 2:1 **Storage:** -20°C

1514	Ceramide trihexosides (bottom spot)	0.5 mg	232.00
	CTH with hydroxy fatty acid side chain $C_{54}H_{101}NO_{19}$		

Source: natural, porcine RBC **Mol. Wt.:** 1068 (2-hydroxystearoyl) **Purity:** 98+% by TLC **Appearance:** solid **Solubility:** DMSO, hot methanol, chloroform/methanol, 1:1
Storage: -20°C

1520	<i>lyso-Ceramide trihexoside</i> <i>lyso-CTH; lyso-Globotriaosylsphingosine</i> C ₃₆ H ₆₇ NO ₁₇ CAS# 126550-86-5	1 mg	429.00
	Source: semisynthetic, porcine RBC Mol. Wt.: 786 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1 Storage: -20°C		
1523	N-Heptadecanoyl-ceramide trihexoside N-C17:0-Ceramide trihexoside; N-Heptadecanoyl globotriaosylceramide C ₅₃ H ₉₉ NO ₁₈	0.5 mg	319.00
	Source: semisynthetic, porcine RBC Mol. Wt.: 1038 Purity: 98+% by TLC Appearance: solid Solubility: DMSO, hot methanol, chloroform/methanol, 2:1 Storage: -20°C		
1524	N-Tricosanoyl-ceramide trihexoside N-C23:0-Ceramide trihexoside; N-Tricosanoyl globotriaosylceramide C ₅₉ H ₁₁₁ NO ₁₈	0.5 mg	319.00
	Source: semisynthetic, porcine RBC Mol. Wt.: 1122 Purity: 98+% by TLC Appearance: solid Solubility: DMSO, hot methanol, chloroform/methanol, 2:1 Storage: -20°C		
1631 1631-001	N-Dodecanoyl-NBD-ceramide trihexoside N-C12:0-NBD-CTH; N-C12:0-NBD-Globotriaosylceramide C ₅₄ H ₉₁ N ₅ O ₂₁	100 µg 1 mg	184.00 793.00
	Source: semisynthetic, porcine RBC Mol. Wt.: 1145 Purity: 98+% by TLC Appearance: solid Solubility: DMSO, hot methanol, chloroform/methanol, 2:1 Storage: -20°C		
1537	N-Octadecanoyl-D₃-ceramide trihexoside, deuterated N-C18:0-D ₃ -CTH, deuterated; N-C18:0-D ₃ -Gb3, deuterated; N-Octadecanoyl-D ₃ -globotriaosylceramide,deuterated; N-Stearoyl-D ₃ -ceramide trihexoside, deuterated C ₅₄ H ₉₈ D ₃ NO ₁₈	0.5 mg	348.00
	Source: semisynthetic, porcine RBC Mol. Wt.: 1055 Purity: 98+% by TLC Appearance: solid Solubility: DMSO, chloroform/methanol, 2:1 Storage: -20°C		

Globosides



1068	Globosides Gb4; Globotetrahexosylceramide C ₆₈ H ₁₂₆ N ₂ O ₂₃ CAS#: 11034-93-8	5 mg	365.00
	Source: natural, porcine RBC Mol. Wt.: 1340 (tetracosanoyl) Purity: 98+% by TLC Appearance: solid Solubility: DMSO, hot methanol, chloroform/methanol, 2:1 Storage: -20°C		

See Table III (pg. 93-97) for fatty acid content

Stable Isotopes Labeled Glycolipids

1914	N-Octadecanoyl-D₃₅-psychosine, perdeuterated N-C18:0-D ₃₅ -Cerebrosides, perdeuterated; N-Stearoyl-D ₃₅ -psychosine, perdeuterated C ₄₂ H ₄₆ D ₃₅ NO ₈	5 mg	306.00
	Source: semisynthetic, bovine Mol. Wt.: 762 Purity: 98+% by TLC Appearance: solid Solubility: chloroform, hot ethanol, chloroform/methanol, 2:1 Storage: -20°C		
	Deuterium labeled stearoyl sidechain		
1533	N-Hexadecanoyl-D₃-glucopsychosine, deuterated N-C16:0-D ₃ -Glucopsychosine, deuterated; N-C16:0-D ₃ -Glucocerebroside, deuterated; N-Palmitoyl-D ₃ -glucopsychosine, deuterated C ₄₀ H ₇₄ D ₃ NO ₈	1 mg	290.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 703 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C		
1534	N-Hexadecanoyl-D₃-lactosylceramide, deuterated N-C16:0-D ₃ -Lactosylceramide, deuterated; N-Palmitoyl-D ₃ -lactosylceramide, deuterated C ₄₆ H ₈₄ D ₃ NO ₁₃	1 mg	376.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 865 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 5:1:0.1 Storage: -20°C		
1536	N-Octadecanoyl-D₃-sulfatide, deuterated N-C18:0-D ₃ -Sulfatide, deuterated; N-Stearoyl-D ₃ -sulfatide, deuterated C ₄₂ H ₇₈ D ₃ NO ₁₁ S	1 mg	414.00
	Source: semisynthetic, bovine Mol. Wt.: 811 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1 Storage: -20°C		
1537	N-Octadecanoyl-D₃-ceramide trihexoside, deuterated N-C18:0-D ₃ -CTH, deuterated; N-C18:0-D ₃ -Gb3, deuterated; N-Octadecanoyl-D ₃ -globotriaosylceramide, deuterated; N-Stearoyl-D ₃ -ceramide trihexoside, deuterated C ₅₄ H ₉₈ D ₃ NO ₁₈	0.5 mg	348.00
	Source: semisynthetic, porcine Mol. Wt.: 1055 Purity: 98+% by TLC Appearance: solid Solubility: DMSO, chloroform/methanol, 2:1 Storage: -20°C		

Fluorescent Compounds

1621	N-Hexanoyl-NBD-galactosylceramide	100 µg	147.00
1621-001	N-C6:0-NBD-beta-D-Galactosylsphingosine; N-C6:0-NBD-Cerebroside C ₃₆ H ₅₉ N ₅ O ₁₁	1 mg	526.00
	Source: semisynthetic, bovine Mol. Wt.: 738 Purity: 98+% by TLC Appearance: solid Solubility: methanol, chloroform/methanol, 5:1 Storage: -20°C		

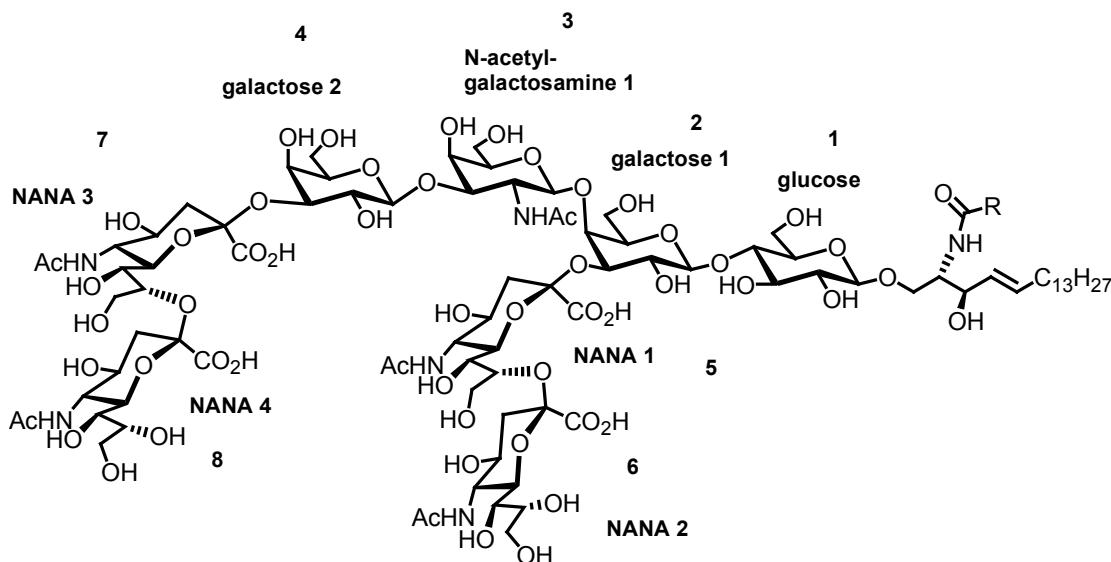
1633	N-Dodecanoyl-NBD-galactosylceramide		100 µg	118.00
1633-001	N-C12:0-NBD- <i>beta</i> -D-Galactosylsphingosine; N-C12:0-NBD-Cerebroside C ₄₂ H ₇₁ N ₅ O ₁₁		1 mg	526.00
	Source: semisynthetic, bovine Mol. Wt.: 822 Purity: 98+% by TLC Appearance: solid Solubility: chloroform, DMSO, chloroform/methanol, 2:1 Storage: -20°C			
1622	N-Hexanoyl-NBD-glucosylceramide		100 µg	147.00
1622-001	N-C6:0-NBD- <i>beta</i> -D-Glucosylsphingosine; N-C6:0-NBD-Glucosylceramide C ₃₆ H ₅₉ N ₅ O ₁₁		1 mg	526.00
	Source: semisynthetic, bovine Mol. Wt.: 738 Purity: 98+% by TLC Appearance: solid Solubility: methanol, chloroform/methanol, 5:1 Storage: -20°C			
1629	N-Hexanoyl-NBD-lactosylceramide		50 ug	184.00
1629-001	N-Hexanoyl-NBD- <i>beta</i> -D-lactosylsphingosine; N-C6:0-NBD- <i>beta</i> -D-Lactosylsphingosine; N-C6:0-NBD-Lactosylceramide C ₄₂ H ₆₉ N ₅ O ₁₆		1 mg	824.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 900 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C			
1630	N-Dodecanoyl-NBD-lactosylceramide		50 µg	204.00
1630-001	N-Dodecanoyl-NBD- <i>beta</i> -D-lactosylsphingosine; N-C12:0-NBD- <i>beta</i> -D-Lactosylsphingosine; N-C12:0-NBD-Lactosylceramide C ₄₈ H ₈₁ N ₅ O ₁₆		1 mg	793.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 984 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C			
1631	N-Dodecanoyl-NBD-ceramide trihexoside		100 µg	184.00
1631-001	N-C12:0-NBD-CTH; N-C12:0-NBD-Globotriaosylceramide C ₅₄ H ₉₁ N ₅ O ₂₁		1 mg	793.00
	Source: semisynthetic, porcine RBC Mol. Wt.: 1145 Purity: 98+% by TLC Appearance: solid Solubility: DMSO, hot methanol, chloroform/methanol, 2:1 Storage: -20°C			
1632	N-Dodecanoyl-NBD-sulfatide		100 µg	119.00
1632-001	N-C12:0-NBD-Sulfatide; N-Dodecanoyl-NBD- <i>lyso</i> -sulfatide; N-Dodecanoyl-NBD-sphingosyl- <i>beta</i> -D-galactoside-3-sulfate C ₄₂ H ₇₁ N ₅ O ₁₄ S		1 mg	526.00
	Source: semisynthetic, bovine Mol. Wt.: 901 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C			

See Biochemicals and Reagents section (page 88) for additional fluorescent labeled products.

Compounds with fluorescent labels other than NBD are available on custom basis. Contact Technical service for more information.

Gangliosides

The diagram below can be used with the general formulas given in the ganglioside descriptions to construct the individual structures.



1064	Gangliotetraosylceramide Asialo GM ₁ ; Gg4 C ₆₂ H ₁₁₄ N ₂ O ₂₃ CAS#: 71012-19-6	1 mg	335.00
	Source: semisynthetic, bovine Mol. Wt.: 1256 (stearoyl) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1; forms micellar solution in water Storage: -20°C		
	General formula: 1,2,3,4 See Table III (pg. 93-97) for fatty acid content		
1512	Gangliotriosylceramide Asialo GM ₂ ; Gg3 C ₅₆ H ₁₀₄ N ₂ O ₁₈	100 µg	447.00
	Source: semisynthetic, human Mol. Wt.: 1093 (stearoyl) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1; forms micellar solution in water Storage: -20°C		
	General formula: 1,2,3		
1061 1061-50	Monosialoganglioside GM₁ (NH₄⁺ salt) GM ₁ C ₇₃ H ₁₃₁ N ₃ O ₃₁ •NH ₃ CAS#: 37758-47-7	5 mg 50 mg	249.00 1,995.00
	Source: natural, bovine Mol. Wt.: 1547 + NH ₃ (stearoyl) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1; forms micellar solution in water Storage: -20°C		
	General formula: 1,2,3,4,5 See Table III (pg. 93-97) for fatty acid content		

1518	<i>lyso</i>-Monosialoganglioside GM₁ (NH₄⁺ salt) <i>lyso</i> -GM ₁ C ₅₅ H ₉₇ N ₃ O ₃₀ •NH ₃ CAS#: 171483-40-2	500 µg	271.00
	Source: semisynthetic, bovine Mol. Wt.: 1280 + NH ₃ Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.2 Storage: -20°C		
1526	Fucosylated monosialoganglioside GM₁ (NH₄⁺ salt) Fucosyl-GM ₁ C ₇₉ H ₁₄₁ N ₃ O ₃₅ •NH ₃ CAS#: 71812-11-8	500 µg	494.00
	Source: natural, porcine Mol. Wt.: 1693 + NH ₃ (stearoyl) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1; forms micellar solution in water Storage: -20°C		
2050	N-omega-CD₃-Octadecanoyl monosialoganglioside GM₁ (NH₄⁺ salt) N-CD ₃ -Stearoyl GM ₁ , C ₇₃ H ₁₂₈ N ₃ O ₃₁ D ₃ •NH ₃	0.5 mg	435.00
	Source: semisynthetic, bovine Mol. Wt.: 1550 + NH ₃ Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1; forms micellar solution in water Storage: -20°C		
1502	Monosialoganglioside GM₂ (NH₄⁺ salt) GM ₂ C ₆₇ H ₁₂₁ N ₃ O ₂₆ •NH ₃ CAS#: 19600-01-2	500 µg	231.00
	Source: natural, human Tay-Sachs Mol. Wt.: 1385+ NH ₃ (stearoyl) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1; forms micellar solution in water Storage: -20°C		
	General formula: 1,2,3,5 See Table III (pg. 93-97) for fatty acid content		
2051	N-omega-CD₃-Octadecanoyl monosialoganglioside GM₂ (NH₄⁺ salt) N-CD ₃ -Stearoyl GM ₂ C ₆₇ H ₁₁₈ D ₃ N ₃ O ₂₆ •NH ₃	250 µg	319.00
	Source: semisynthetic, human Tay-Sachs Mol. Wt.: 1388 + NH ₃ Purity: 98+% by TLC, MS Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1; forms micellar solution in water Storage: -20°C		
1503	Monosialoganglioside GM₃ (NH₄⁺ salt) GM ₃ C ₆₄ H ₁₁₈ N ₂ O ₂₁ •NH ₃ CAS#: 54827-14-4	1 mg	127.00
	Source: natural, bovine buttermilk Mol. Wt.: 1252+ NH ₃ (tricosanoyl) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1; forms micellar solution in water Storage: -20°C		
	General formula: 1,2,5 See Table III (pg. 93-97) for fatty acid content		
2052	N-omega-CD₃-Octadecanoyl monosialoganglioside GM₃ (NH₄⁺ salt) N-CD ₃ -Stearoyl GM ₃ C ₅₉ H ₁₀₅ D ₃ N ₂ O ₂₁ •NH ₃	250 µg	290.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 1185 + NH ₃ Purity: 98+% by TLC, MS Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.2; forms micellar solution in water Storage: -20°C		

1535	Monosialoganglioside GM4, egg (NH_4^+ salt) GM ₄ C ₅₇ H ₁₀₆ N ₂ O ₁₇ •NH ₃ CAS#: 66456-69-7	0.5 mg	319.00
Source: natural, egg, chicken Mol. Wt.: 1091+NH ₃ (2-hydroxydocosanoyl) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1; forms micellar solution in water Storage: -20°C			
General formula: 2,5 See Table III (pg. 93-97) for fatty acid content			
1062	Disialoganglioside GD_{1a} (NH_4^+ salt) GD _{1a} C ₈₄ H ₁₄₈ N ₄ O ₃₉ •2NH ₃ CAS#: 12707-58-3	5 mg	262.00
Source: natural, bovine Mol. Wt.: 1838 + 2NH ₃ (stearoyl) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1; forms micellar solution in water Storage: -20°C			
General formula: 1,2,3,4,5,7 See Table III (pg. 93-97) for fatty acid content			
1501	Disialoganglioside GD_{1b} (NH_4^+ salt) GD _{1b} C ₈₄ H ₁₄₈ N ₄ O ₃₉ •2NH ₃ CAS#: 19553-76-5	1 mg	199.00
Source: natural, bovine Mol. Wt.: 1838+ 2NH ₃ (stearoyl) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1; forms micellar solution in water Storage: -20°C			
General formula: 1,2,3,4,5,6 See Table III (pg. 93-97) for fatty acid content			
1527	Disialoganglioside GD₂ (NH_4^+ salt) GD ₂ C ₇₈ H ₁₃₈ N ₄ O ₃₄ •2NH ₃ CAS#: 65988-71-8	0.5 mg	579.00
Source: semisynthetic, rabbit Mol. Wt.: 1676 + 2NH ₃ (stearoyl) Purity: 98+% by TLC, MS Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1; forms micellar solution in water Storage: -20°C			
General formula: 1,2,3,5,6			
1504	Disialoganglioside GD₃ (NH_4^+ salt) GD ₃ C ₇₅ H ₁₃₅ N ₃ O ₂₉ •2NH ₃ CAS#: 62010-37-1	5 mg	248.00
Source: natural, bovine buttermilk Mol. Wt.: 1543+2NH ₃ (tricosanoyl) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1; forms micellar solution in water Storage: -20°C			
General formula: 1,2,5,6 See Table III (pg. 93-97) for fatty acid content			
1063	Trisialoganglioside GT_{1b} (NH_4^+ salt) GT _{1b} C ₉₅ H ₁₆₅ N ₅ O ₄₇ •3NH ₃ CAS#: 59247-13-1	5 mg	319.00
Source: natural, bovine Mol. Wt.: 2129 + 3NH ₃ (stearoyl) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1; forms micellar solution in water Storage: -20°C			
General Formula: 1,2,3,4,5,6,7 See Table III (pg. 93-97) for fatty acid content			

1516	Tetrasialoganglioside GQ_{1b} (NH₄⁺ salt) GQ _{1b} C ₁₀₆ H ₁₈₂ N ₆ O ₅₅ •4NH ₃ CAS#: 68652-37-9	100 µg	169.00
Source: natural, bovine Mol. Wt.: 2421+4NH ₃ (stearoyl) Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1; forms micellar solution in water Storage: -20°C			
General formula: 1,2,3,4,5,6,7,8 See Table III (pg. 93-97) for fatty acid content			
1065	Mixed Gangliosides, purified, bovine (NH₄⁺ salt) Mixed Gangliosides	25 mg	193.00
Source: natural, bovine Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1; forms micellar solution in water Storage: -20°C			
Approximately 98% GM ₁ , GD _{1a} , GD _{1b} and GT _{1b} , remaining 2% other gangliosides See Table III (pg. 93-97) for fatty acid content			

1525	Mixed Gangliosides, purified, porcine, (NH₄⁺ salt)	25 mg	226.00
Source: natural, porcine Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1; forms micellar solution in water Storage: -20°C			
Approximately 98% GM ₁ , GD _{1a} , GD _{1b} and GT _{1b} , remaining 2% other gangliosides			

Glycosphingolipid Reference Mixtures for TLC

These mixtures are qualitative standards prepared from our purified glycosphingolipids.

1505	Neutral Glycosphingolipid Mixture	1 mg/ml, 1 ml	136.00
Source: natural, bovine and porcine Appearance: liquid Solvent: chloroform/methanol, 2:1 Solubility: chloroform/methanol, 2:1 Storage: -20°C			
Contains: cerebrosides, lactosylceramides, ceramide trihexosides, globosides			
1508	Monosialoganglioside Mixture	0.5 mg/ml, 1 ml	136.00
Source: natural, bovine, human Appearance: liquid Solvent: chloroform/methanol/DI water, 2:1:0.1 Solubility: chloroform/methanol/DI water, 2:1:0.1 Storage: -20°C			
Contains: GM ₃ , GM ₂ , GM ₁			
1509	Disialoganglioside Mixture	0.5 mg/ml, 1 ml	147.00
Source: natural, bovine Appearance: liquid Solvent: chloroform/methanol/DI water, 2:1:0.1 Solubility: chloroform/methanol/DI water, 2:1:0.1 Storage: -20°C			
Contains: GD ₃ , GD _{1a} , GD _{1b}			

1510	Lactosylceramide and Sialosyl Derivatives Mixture	0.5 mg/ml, 1 ml	169.00
Source: natural, bovine buttermilk Appearance: liquid Solvent: chloroform/methanol/DI water, 2:1:0.1 Solubility: chloroform/methanol/DI water, 2:1:0.1 Storage: -20°C			
	Contains: LC, GM ₃ , GD ₃		
<hr/>			
1511	Gangliotetraosylceramide and Sialosyl Derivatives Mixture	0.5 mg/ml, 1 ml	126.00
Source: natural, bovine Appearance: liquid Solvent: chloroform/methanol/DI water, 2:1:0.1 Solubility: chloroform/methanol/DI water, 2:1:0.1 Storage: -20°C			
	Contains: asialo GM ₁ , GM ₁ , GD _{1a} , GD _{1b} , GT _{1b}		

Antibodies Directed Against Glycolipids

These monoclonal and polyclonal antibodies are directed against the carbohydrate chains of Matreya's glycolipids. The same carbohydrate moieties are found on many glycoproteins. The antibodies are for use in ELISA or TLC immunoblotting applications (9). All antibodies are quality tested by actual performance in ELISA and TLC immunoblotting. The antibodies contain no preservatives and are shipped on dry ice.

See Literature References on page 99.

1977	Anti-ganglioside GD₃	50 µl	271.00
Monoclonal antibody to GD ₃ , isotype IgG/IgM			
Source: natural, mouse hybridoma R-24 cell line Appearance: liquid Solubility: DI water Storage: -20°C Dry Ice Charge Applies			
Suitable for TLC immunoblotting, ELISA			
1950	Anti-ganglioside asialo GM₁	100 µl	342.00
Polyclonal antibody to asialo GM ₁ , isotype IgG/IgM			
Source: natural, rabbit Appearance: liquid Solubility: DI water Storage: -20°C Dry Ice Charge Applies			
Suitable for ELISA, TLC-immunoblotting. Slight cross reaction to GM ₁			
1951	Anti-ganglioside asialo GM₂	50 µl	271.00
Polyclonal antibody to asialo GM ₂ , isotype IgG/IgM			
Source: natural, rabbit Appearance: liquid Solubility: DI water Storage: -20°C Dry Ice Charge Applies			
Suitable for ELISA, TLC-immunoblotting			
1954	Anti-ganglioside GM₁	100 µl	271.00
Polyclonal antibody to GM ₁ , isotype IgG/IgM			
Source: natural, rabbit Appearance: liquid Solubility: DI water Storage: -20°C Dry Ice Charge Applies			
Suitable for ELISA, TLC-immunoblotting. Slight cross reaction to asialo-GM ₁			

1961	Anti-ganglioside GM₂ (NANA) Polyclonal antibody to GM ₂ (NANA), isotype IgG/IgM	50 µl	271.00
Source: natural, rabbit Appearance: liquid Solubility: DI water Storage: -20°C Dry Ice Charge Applies			
Suitable for ELISA, TLC-immunoblotting			
1957	Anti-ganglioside GM₄ Polyclonal antibody to GM ₄ , isotype IgG/IgM	50 µl	271.00
Source: natural, rabbit Appearance: liquid Solubility: DI water Storage: -20°C Dry Ice Charge Applies			
Suitable for ELISA, TLC-immunoblotting			
1963	Anti-ganglioside GD₂ Polyclonal antibody to GD ₂ , isotype IgG/IgM	50 µl	285.00
Source: natural, rabbit Appearance: liquid Solubility: DI water Storage: -20°C Dry Ice Charge Applies			
Suitable for ELISA, TLC-immunoblotting			
1964	Anti-ganglioside GD_{1b} Polyclonal antibody to GD _{1b} , isotype IgG/IgM	50 µl	285.00
Source: natural, rabbit Appearance: liquid Solubility: DI water Storage: -20°C Dry Ice Charge Applies			
Suitable for ELISA, TLC-immunoblotting			
1960	Anti-globoside GL-4 Polyclonal antibody to GL-4, isotype IgG/IgM	50 µl	271.00
Source: natural, rabbit Appearance: liquid Solubility: DI water Storage: -20°C Dry Ice Charge Applies			
Suitable for ELISA, TLC-immunoblotting			

Enzyme Inhibitors

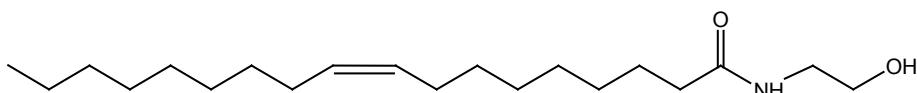
Ceramide: UDPglucose Transferase. PDMP (D,L-*threo*-1-phenyl-2-decanoylamino-3-morpholino-1-propanol-HCl) closely resembles the natural sphingolipid substrate of brain glucosyl transferase and is a very potent and competitive inhibitor of the enzyme (26). It has been shown to block outgrowth of neurites in cultured retina and to block glucolipid synthesis in cultured 3T3 cells (27). N.S. Radin and co-workers have shown (28) that PPMP has activity equivalent to that of PDMP when cell homogenates and brain and liver microsomes are used, but it is about 20 times more potent when used with intact cells. In another paper (29), Radin's group has shown that PDMP has substantial activity against Ehrlich ascites tumors in mice. Recent publications from the laboratory of Myles Cabot (30, 31) show that PPMP can reverse multi-drug resistance in cancer cells by causing a build-up of ceramide and preventing the synthesis of glycosylated ceramides. **See Literature References on page 99.**

Matreya also offers the resolved D- and L-*threo*-isomers of PDMP and PPMP.

Protein Kinase C Inhibitor. Sphingosine is a potent and reversible inhibitor of protein kinase C (32); it also has been shown at low concentrations to stimulate DNA synthesis and act synergistically with known growth factors (33). Note that Safingol (our L-*threo*-dihydrosphingosine) has also been shown to partially reverse multi-drug resistance in cancer cells (31) via inhibition of protein kinase C.

Dihydroceramide desaturase Inhibitor. Cyclopropenylceramide is the first known inhibitor of this enzyme and may allow significant studies on the role of ceramide in apoptosis. Matreya is the only source for this inhibitor. (34)

Ceramidase Inhibitors. N-Oleoyl ethanolamine has been shown to be an efficacious inhibitor of the ceramidase found in human kidney and cerebellum (35). It is specifically an inhibitor of acid ceramidase (36) with an IC₅₀ of ca. 500 µM. N-Hexadecanoyl ethanolamine can be used as an inactive control. D-MAPP is a potent (IC₅₀ approximately 5 µM) inhibitor of alkaline ceramidase. Its enantiomer L-MAPP is inactive as an inhibitor and acts as a substrate for this enzyme (36, 37). **See Literature References on page 99.**



Catalog number 1751

1751	N-Oleoyl ethanolamine NOE C ₂₀ H ₃₉ NO ₂ CAS#: 111-58-0	100 mg	163.00
------	--	--------	--------

Source: synthetic **Mol. Wt.:** 326 **Melting Point (°C):** 63-66 **Purity:** 98+% by TLC,
GC Appearance: solid **Solubility:** chloroform, ethanol, methanol, ethyl ether, DMSO
Storage: -20°C

Activity: acid ceramidase inhibitor

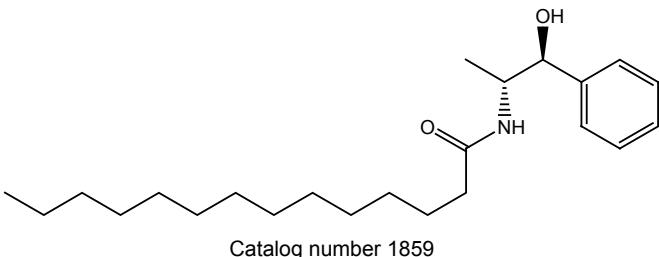
1786	N-Hexadecanoyl ethanolamine C ₁₈ H ₃₇ NO ₂ CAS# 544-31-0	100 mg	114.00
------	---	--------	--------

Source: synthetic **Mol. Wt.:** 299 **Melting Point (°C):** 99-102 **Purity:** 98+% by TLC
Appearance: solid **Solubility:** chloroform, ethanol, methanol, DMSO
Storage: -20°C

Activity: inactive as acid ceramidase inhibitor

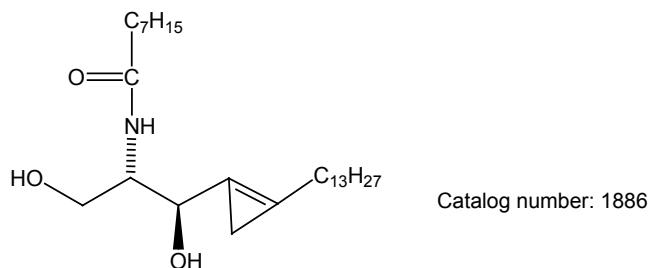
1807	L-<i>threo</i>-Dihydrosphingosine (Safingol)	5 mg	174.00
1807-025	L- <i>threo</i> -Sphinganine, C18 chain C ₁₈ H ₃₉ NO ₂ CAS#: 15639-50-6	25 mg	556.00

Source: synthetic **Mol. Wt.:** 301 **Melting Point (°C):** 103-114 **Purity:** 98+% by TLC, GC
Appearance: solid **Solubility:** chloroform, ethanol, methanol
Storage: -20°C



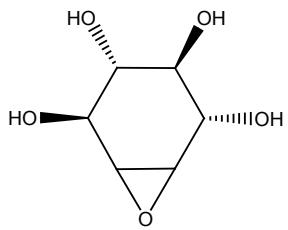
1859	D-MAPP D- <i>erythro</i> -2-Tetradecanoylamino-1-phenyl-1-propanol C ₂₃ H ₃₉ NO ₂ CAS#: 143492-39-1	100 mg	154.00
Source: synthetic Mol. Wt.: 361 Purity: 98+% by TLC Appearance: solid Solubility: ethanol Storage: -20°C			
Activity: alkaline ceramidase inhibitor			

1860	L-MAPP L- <i>erythro</i> -2-Tetradecanoylamino-1-phenyl-1-propanol C ₂₃ H ₃₉ NO ₂ CAS#: 143492-38-0	100 mg	154.00
Source: synthetic Mol. Wt.: 361 Purity: 98+% by TLC Appearance: solid Solubility: ethanol Storage: -20°C			
Activity: inactive as alkaline ceramidase inhibitor			



1886	N-C8:0-Cyclopropenylceramide	1 mg	194.00
1886-005	N-C8:0-CPPC; N-[(1R, 2S)-2-Hydroxy-1-hydroxymethyl-2-(2-tridecyl-1-cyclopropenyl) ethyl] octanamide; GT ₁₁ C ₂₇ H ₅₁ NO ₃	5 mg	724.00
Source: synthetic Mol. Wt.: 438 Melting Point (°C): 69-70 Purity: 98+% by ¹ H NMR; HPLC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: -20°C			
Activity: Dihydroceramide desaturase inhibitor			

1887	N-C16:0-Cyclopropenylceramide	1 mg	197.00
1887-005	N-C16:0-CPPC; N-[(1R, 2S)-2-Hydroxy-1-hydroxymethyl-2-(2-tridecyl-1-cyclopropenyl) ethyl] hexadecamide C ₃₅ H ₆₇ NO ₃	5 mg	724.00
Source: synthetic Mol. Wt.: 550 Melting Point (°C): 156-157 Purity: 98+% by ¹ H NMR; HPLC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: -20°C			
Activity: Dihydroceramide desaturase inhibitor			



Catalog number 1889

1889	Conduritol B Epoxide C ₆ H ₁₀ O ₅ CAS#: 6090-95-5	25 mg	195.00
-------------	--	--------------	---------------

Source: synthetic **Mol. Wt.:** 162 **Melting Point (°C):** 164-166 **Purity:** 98+% by TLC, NMR **Appearance:** solid **Solubility:** DI water, DMSO, methanol (slightly) **Storage:** -20°C

Inhibits α -and β -glucosidase activity; specific inhibitor of glucocerebrosidase in cultured cells.

1719	D,L-threo-PDMP D,L- <i>threo</i> -1-Phenyl-2-decanoylamino-3-morpholino-1-propanol•HCl C ₂₃ H ₃₈ N ₂ O ₃ •HCl CAS#: 80938-69-8	100 mg	184.00
-------------	---	---------------	---------------

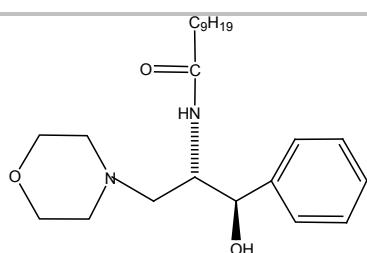
Source: synthetic **Mol. Wt.:** 427 **Purity:** 98+% by TLC **Appearance:** solid **Solubility:** ethanol, methanol, chloroform, DMSO **Storage:** -20°C

Activity: glucosylceramide synthase inhibitor

1720	D,L-threo-PPMP D,L- <i>threo</i> -1-Phenyl-2-hexadecanoylamino-3-morpholino-1-propanol•HCl C ₂₉ H ₅₀ N ₂ O ₃ •HCl CAS#: 149022-18-4	100 mg	184.00
-------------	--	---------------	---------------

Source: synthetic **Mol. Wt.:** 511 **Purity:** 98+% by TLC **Appearance:** solid **Solubility:** alcohols, chloroform **Storage:** -20°C

Activity: glucosylceramide synthase inhibitor

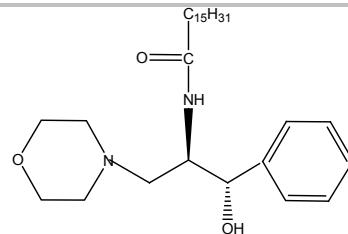


Catalog number 1749

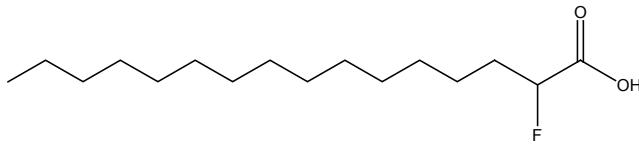
1749	L-threo-PDMP L- <i>threo</i> -1-Phenyl-2-decanoylamino-3-morpholino-1-propanol•HCl C ₂₃ H ₃₈ N ₂ O ₃ •HCl CAS#: 109836-81-9	10 mg	250.00
-------------	--	--------------	---------------

Source: synthetic **Mol. Wt.:** 427 **Purity:** 98+% by TLC **Appearance:** solid **Solubility:** ethanol, methanol **Storage:** -20°C

1753	D,L-erythro-PPMP D,L- <i>erythro</i> -1-Phenyl-2-hexadecanoylamino-3-morpholino-1-propanol•HCl C ₂₉ H ₅₀ N ₂ O ₃ •HCl	100 mg	169.00
	Source: synthetic Mol. Wt.: 511 Purity: 98+% by TLC Appearance: solid Solubility: chloroform, ethanol, methanol, DSMO Storage: -20°C		
1755	D,L-erythro-PDMP D,L- <i>erythro</i> -1-Phenyl-2-decanoylamino-3-morpholino-1-propanol•HCl C ₂₃ H ₃₈ N ₂ O ₃ •HCl CAS#: 109760-77-2	100 mg	169.00
	Source: synthetic Mol. Wt.: 427 Purity: 98+% by TLC Appearance: solid Solubility: chloroform, ethanol, methanol, DSMO Storage: -20°C		
1756	D-threo-PDMP D- <i>threo</i> -1-Phenyl-2-decanoylamino-3-morpholino-1-propanol•HCl C ₂₃ H ₃₈ N ₂ O ₃ •HCl CAS#: 109836-82-0	10 mg	250.00
	Source: synthetic Mol. Wt.: 427 Purity: 98+% by TLC Appearance: solid Solubility: ethanol, methanol Storage: -20°C		
	Activity: glucosylceramide synthase inhibitor		
1865	D-threo-PPMP D- <i>threo</i> -1-Phenyl-2-hexadecanoylamino-3-morpholino-1-propanol•HCl C ₂₉ H ₅₀ N ₂ O ₃ •HCl	10 mg	250.00
	Source: synthetic Mol. Wt.: 511 Purity: 98+% by TLC Appearance: solid Solubility: ethanol, methanol Storage: -20°C		
	Activity: glucosyl ceramide synthase inhibitor		
1868	L-threo-PPMP L- <i>threo</i> -1-Phenyl-2-hexadecanoylamino-3-morpholino-1-propanol•HCl C ₂₉ H ₅₀ N ₂ O ₃ •HCl	10 mg	250.00
	Source: synthetic Mol. Wt.: 511 Purity: 98+% by TLC Appearance: solid Solubility: ethanol, methanol Storage: -20°C		
1800	Castanospermine 1,6,7,8-Tetrahydroxyoctahydroindolizine C ₈ H ₁₅ NO ₄ CAS#: 79831-76-8	25 mg	65.00
	Source: natural, plant Mol. Wt.: 189 Melting Point (°C): 210-215 Purity: 98+%		
	by TLC, NMR Appearance: solid Solubility: DI water, methanol/DI water, 90:10		
	Storage: -20°C		



Catalog number 1865



Catalog number 1717

1717 2-Fluoropalmitic acid **25 mg** **162.00**
C₁₆H₃₁FO₂ CAS#: 89270-22-4

Source: synthetic **Mol. Wt.:** 274 **Melting Point (°C):** 83-85 **Purity:** 98+% by TLC
Appearance: solid **Solubility:** chloroform, ethanol, methanol **Storage:** -20°C

Activity: Acyl-CoA synthase inhibitor

1718 Methyl 2-fluoropalmitate **10 mg** **162.00**
C₁₇H₃₃FO₂

Source: synthetic **Mol. Wt.:** 288 **Melting Point (°C):** 36-38 **Purity:** 98+% by TLC
Appearance: solid **Solubility:** chloroform, ethanol, methanol **Storage:** -20°C

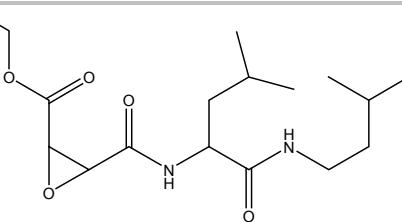
Activity: inactive ester of 2-fluoropalmitic acid

1750 2,2-Difluoropalmitic acid **25 mg** **147.00**
C₁₆H₃₀F₂O₂

Source: synthetic **Mol. Wt.:** 292 **Melting Point (°C):** 50.8-53 **Purity:** 98+% by TLC, GC
Appearance: solid **Solubility:** chloroform, ethanol, methanol
Storage: -20°C

1858 2-Acetyl-4-(1R, 2S, 3R, 4-tetrahydroxybutyl)-imidazole **1 mg** **176.00**
THI C₉H₁₄N₂O₅ CAS#: 94944-70-4

Source: synthetic **Mol. Wt.:** 230 **Melting Point (°C):** n/a **Purity:** 99% by HPLC, MS, NMR
Appearance: solid **Solubility:** DI water **Storage:** -20°C



Catalog number 1752

1752 EST **5 mg** **319.00**
E-64-d; Loxastatin C₁₇H₃₀N₂O₅ CAS#: 88321-09-9

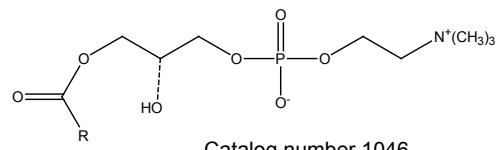
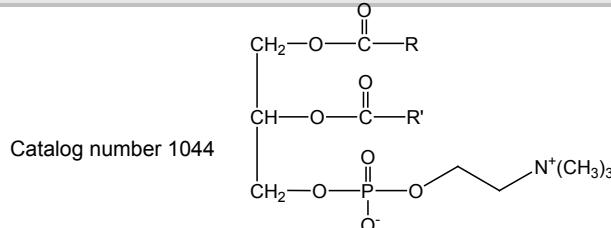
Source: synthetic **Mol. Wt.:** 342 **Melting Point (°C):** 125-127 **Purity:** 98+% by TLC, GC
Appearance: solid **Solubility:** chloroform, ethanol, methanol
Storage: -20°C

Activity: cysteine protease inhibitor

Glycerolipids

Glycerophospholipids

Natural Phospholipids



Catalog number 1046

1044	Lecithin Phosphatidylcholine; PC C ₄₄ H ₈₄ NO ₈ P CAS#: 8002-43-5	50 mg/ml, 1 ml	50.00
------	--	----------------	-------

Source: natural, chicken, egg **Mol. Wt.:** 787 (oleoyl) **Purity:** 98+% by TLC
Appearance: liquid **Solvent:** chloroform **Solubility:** chloroform, ethyl ether, ethanol
Storage: -20°C

See Table III (pg. 93-97) for fatty acid content

1070	Lecithin Phosphatidylcholine; PC C ₄₄ H ₈₄ NO ₈ P CAS#: 8002-43-5	50 mg/ml, 1 ml	174.00
------	--	----------------	--------

Source: natural, bovine **Mol. Wt.:** 787 (oleoyl) **Purity:** 98+% by TLC
Appearance: liquid **Solvent:** chloroform **Solubility:** chloroform, ethyl ether
Storage: -20°C

See Table III (pg. 93-97) for fatty acid content

1302	Lecithin Phosphatidylcholine; PC C ₄₄ H ₈₀ NO ₈ P CAS#: 8002-43-5	50 mg/ml, 1 ml	86.00
------	--	----------------	-------

Source: natural, plant **Mol. Wt.:** 782 (linoleoyl) **Purity:** 98+% by TLC
Appearance: liquid **Solvent:** chloroform **Solubility:** chloroform, ethyl ether
Storage: -20°C

See Table III (pg. 93-97) for fatty acid content

1046	lyso-Lecithin lyso-Phosphatidylcholine C ₂₄ H ₅₀ NO ₇ P CAS#: 9008-30-4	50 mg	85.00
------	--	-------	-------

Source: semisynthetic, chicken, egg **Mol. Wt.:** 496 (palmitoyl) **Purity:** 98+% by TLC
Appearance: solid **Solvent:** chloroform/methanol, 2:1 **Storage:** -20°C

See Table III (pg. 93-97) for fatty acid content

1047	Phosphatidylserine PS C ₄₂ H ₇₈ NO ₁₀ P	50 mg/ml, 1 ml	197.00
------	--	----------------	--------

Source: natural, bovine **Mol. Wt.:** 788 (oleoyl) **Purity:** 98+% by TLC
Appearance: liquid **Solvent:** chloroform **Solubility:** chloroform, toluene
Storage: -20°C

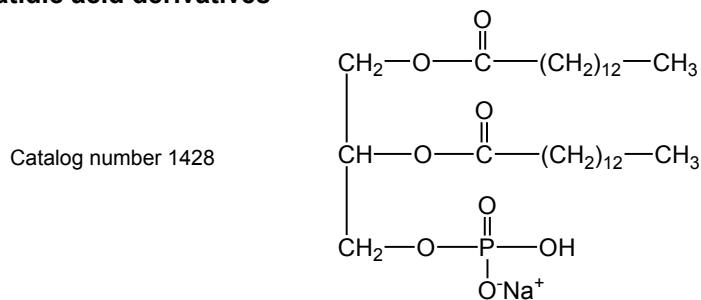
See Table III (pg. 93-97) for fatty acid content

1048	Phosphatidylinositol (Na⁺ salt) PI C ₄₅ H ₇₈ O ₁₃ P•Na CAS# 383907-36-6	10 mg/ml, 1 ml	126.00
Source: natural, plant Mol. Wt.: 858 + Na (linoleoyl) Purity: 98+% by TLC Appearance: liquid Solvent: chloroform Solubility: chloroform, ethyl ether Storage: -20°C			
See Table III (pg. 93-97) for fatty acid content			
<hr/>			
1336	Phosphatidylinositol, plant, soy, (Na salt) C ₄₆ H ₈₀ O ₁₃ P•Na CAS# 383907-36-6	50 mg/ml, 1ml	328.00
Source: natural, plant, soy Mol. Wt.: 872 +Na (linoleoyl) Purity: 98+% by TLC Appearance: liquid Solvent: chloroform Solubility: chloroform, ethyl ether Storage: -20°C			
<hr/>			
1053	Phosphatidic acid (NH₄⁺ salt) PA C ₃₉ H ₇₂ O ₈ P•NH ₄ ⁺	50 mg	111.00
Source: semisynthetic, chicken, egg Mol. Wt.: 718 (oleoyl, NH ₄ ⁺) Purity: 98+% by TLC Appearance: solid Solubility: chloroform, ethyl ether Storage: -20°C			
See Table III (pg. 93-97) for fatty acid content			
<hr/>			
1045	Phosphatidylethanolamine PE C ₄₁ H ₇₈ NO ₈ P CAS#: 39382-08-6	50 mg/ml, 1 ml	151.00
Source: natural, chicken, egg Mol. Wt.: 744 (oleoyl) Purity: 98+% by TLC Appearance: liquid Solvent: chloroform Solubility: chloroform Storage: -20°C			
See Table III (pg. 93-97) for fatty acid content			
<hr/>			
1069	Phosphatidylethanolamine PE C ₄₁ H ₇₈ NO ₈ P CAS#: 90989-93-8	50 mg/ml, 1 ml	219.00
Source: natural, bovine Mol. Wt.: 744 (oleoyl) Purity: 98+% by TLC Appearance: liquid Solvent: chloroform Solubility: chloroform Storage: -20°C			
<hr/>			
1301	Phosphatidylethanolamine PE C ₄₁ H ₇₄ NO ₈ P CAS#: 90989-93-8	50 mg/ml, 1 ml	133.00
Source: natural, plant Mol. Wt.: 740 (linoleoyl) Purity: 98+% by TLC Appearance: liquid Solvent: chloroform Solubility: chloroform Storage: -20°C			
See Table III (pg. 93-97) for fatty acid content			
<hr/>			
1052	Phosphoglycerides Kit	1 each	503.00
Source: natural, chicken egg, bovine, plant Purity: 98+% by TLC Appearance: liquid/solid Solvent: various Storage: -20°C			
Individually packed in ampules and vials (Purity 98+%) 10mg; Phosphatidylethanolamine, egg (in 1 ml CHCl ₃) 10mg; Sphingomyelin, bovine 10mg; Phosphatidylserine, bovine (in 1 ml CHCl ₃) 10 mg; Lecithin, egg (in 1 ml CHCl ₃) 10 mg; lyso-Lecithin, egg 10 mg; Cerebroside, bovine 10mg; Sulfatides, bovine 10mg; Phosphatidylinositol, Na ⁺ salt, plant (in 1 ml CHCl ₃) 3mg			

Synthetic Phospholipids

These phospholipids have 98+% chemical purity except where stated and 99% fatty acid chain purity. Store at -20°C.
Solubility: see individual entries

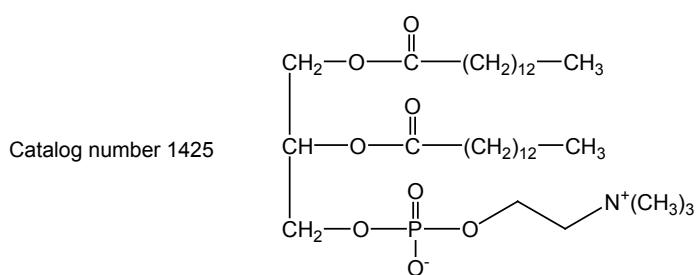
Phosphatidic acid derivatives



1428	1,2-Dimyristoyl-sn-glycero-3-phosphatidic acid (Na⁺ salt) DMPA C ₃₁ H ₆₀ O ₈ P•Na CAS#: 80724-31-8	100 mg	64.00
Source: synthetic Mol. Wt.: 615 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/acetic acid, 4:1:0.1 Storage: -20°C			
1429	1,2-Dipalmitoyl-sn-glycero-3-phosphatidic acid (Na⁺ salt) DPPA C ₃₅ H ₆₈ O ₈ P•Na CAS#: 71065-87-7	100 mg	64.00
Source: synthetic Mol. Wt.: 671 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/acetic acid, 4:1:0.1 Storage: -20°C			
1430	1,2-Distearoyl-sn-glycero-3-phosphatidic acid (Na⁺ salt) DSPA C ₃₉ H ₇₆ O ₈ P•Na CAS#: 108321-18-2	100 mg	64.00
Source: synthetic Mol. Wt.: 727 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/acetic acid, 4:1:0.1 Storage: -20°C			

Phosphatidylcholines

1442	1,2-Dilauroyl-sn-glycero-3-phosphorylcholine DLPC C ₃₂ H ₆₄ NO ₈ P CAS#: 18194-25-7	100 mg	64.00
Source: synthetic Mol. Wt.: 622 Purity: 98+% by TLC Appearance: solid Solubility: methylene chloride, methanol Storage: -20°C			

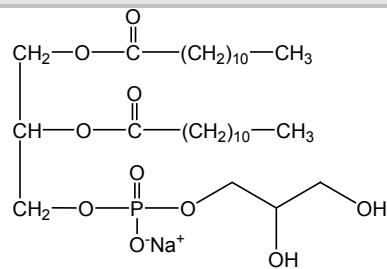


1425	1,2-Dimyristoyl-sn-glycero-3-phosphorylcholine DMPC C ₃₆ H ₇₂ NO ₈ P CAS#: 18194-24-6	100 mg	64.00
Source: synthetic Mol. Wt.: 678 Purity: 98+% by TLC Appearance: solid Melting Point: 130-139°C Solubility: methylene chloride, methanol Storage: -20°C			
1426	1,2-Dipalmitoyl-sn-glycero-3-phosphorylcholine DPPC C ₄₀ H ₈₀ NO ₈ P CAS#: 63-89-8	100 mg	64.00
Source: synthetic Mol. Wt.: 734 Purity: 98+% by TLC Appearance: solid Solubility: methylene chloride, methanol Storage: -20°C			
1400	1,2-Diheptadecanoyl-sn-glycero-3-phosphorylcholine DHDPG C ₄₂ H ₈₄ NO ₈ P CAS#: 70897-27-7	50 mg	63.00
Source: synthetic Mol. Wt.: 762 Purity: 98+% by TLC Appearance: solid Solubility: methylene chloride, methanol Storage: -20°C			
1427	1,2-Distearoyl-sn-glycero-3-phosphorylcholine DSPC C ₄₄ H ₈₈ NO ₈ P CAS#: 816-94-4	100 mg	64.00
Source: synthetic Mol. Wt.: 790 Purity: 98+% by TLC Appearance: solid Solubility: methylene chloride, methanol Storage: -20°C			
1437	1-Palmitoyl-2-oleoyl-sn-glycero-3-phosphorylcholine POPC C ₄₂ H ₈₂ NO ₈ P CAS#: 26853-31-6	100 mg	83.00
Source: synthetic Mol. Wt.: 760 Purity: 98+% by TLC Appearance: solid Solubility: methylene chloride, methanol Storage: -20°C			
1445	1-Palmitoyl-sn-glycero-3-phosphorylcholine lyso-PPG C ₂₄ H ₅₀ NO ₇ P CAS#: 17364-16-8	100 mg	83.00
Source: synthetic Mol. Wt.: 496 Purity: 98+% by TLC Appearance: solid Solubility: methylene chloride, methanol Storage: -20°C			
1409	1-Stearoyl-2-linoleoyl-sn-glycero-3-phosphorylcholine C ₄₄ H ₈₄ NO ₈ P	25 mg/ml, 1ml	204.00
Source: synthetic Mol. Wt.: 786 Purity: 98+% by TLC Appearance: liquid Solvent: chloroform Solubility: chloroform, ethanol Storage: -20°C			

1410	1-Stearoyl-2-[9(Z),11(E)-octadecadienoyl]-sn-glycero-3-phosphorylcholine C ₄₄ H ₈₄ NO ₈ P	25 mg/ml, 1ml	204.00
	Source: synthetic Mol. Wt.: 786 Purity: 98+% by TLC Appearance: liquid Solvent: chloroform Solubility: chloroform, ethanol Storage: -20°C		
1411	1-Stearoyl-2-[10(E),12(Z)-octadecadienoyl]-sn-glycero-3-phosphorylcholine C ₄₄ H ₈₄ NO ₈ P	25 mg/ml, 1ml	204.00
	Source: synthetic Mol. Wt.: 786 Purity: 98+% by TLC Appearance: liquid Solvent: chloroform Solubility: chloroform, ethanol Storage: -20°C		

Phosphatidylglycerols

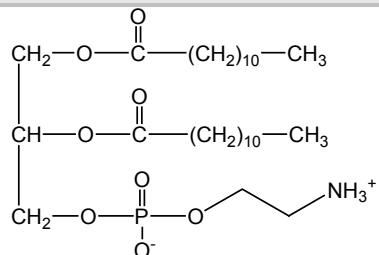
Catalog number 1443



1443	1,2-Dilauroyl-sn-glycero-3-phosphorylglycerol (Na⁺ salt) DLPG C ₃₀ H ₅₈ O ₁₀ P•Na CAS#: 73548-69-3	100 mg	64.00
	Source: synthetic Mol. Wt.: 632 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 5:1 Storage: -20°C		
1431	1,2-Dimyristoyl-sn-glycero-3-phosphorylglycerol (Na⁺ salt) DMPG C ₃₄ H ₆₆ O ₁₀ P•Na CAS#: 200880-40-6	100 mg	64.00
	Source: synthetic Mol. Wt.: 689 Purity: 98+% by TLC Appearance: solid Melting Point: 120-129°C Solubility: chloroform/methanol, 5:1 Storage: -20°C		
1432	1,2-Dipalmitoyl-sn-glycero-3-phosphorylglycerol (Na⁺ salt) DPPG C ₃₈ H ₇₄ O ₁₀ P•Na CAS#: 200880-41-7	100 mg	64.00
	Source: synthetic Mol. Wt.: 745 Purity: 98+% by TLC Appearance: solid Melting Point: 122-127°C Solubility: chloroform/methanol, 5:1 Storage: -20°C		
1433	1,2-Distearoyl-sn-glycero-3-phosphorylglycerol (Na⁺ salt) DSPG C ₄₂ H ₈₂ O ₁₀ P•Na CAS#: 4537-78-4	100 mg	64.00
	Source: synthetic Mol. Wt.: 801 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 5:1 Storage: -20°C		
1438	1-Palmitoyl-2-oleoyl-sn-glycero-3-phosphorylglycerol (Na⁺ salt) POPG C ₄₀ H ₇₆ O ₁₀ P•Na CAS#: 202070-86-8	100 mg	138.00
	Source: synthetic Mol. Wt.: 771 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 5:1 Storage: -20°C		

Phosphatidylethanolamines

Catalog number 1444

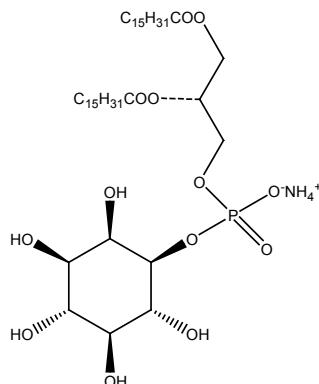


1444	1,2-Dilauroyl-sn-glycero-3-phosphorylethanolamine DLPE C ₂₉ H ₅₈ NO ₈ P CAS#: 42436-56-6	100 mg	70.00
Source: synthetic Mol. Wt.: 580 Purity: 98+% by TLC Appearance: solid Solubility: chloroform + methanol mixture Storage: -20°C			
1434	1,2-Dimyristoyl-sn-glycero-3-phosphorylethanolamine DMPE C ₃₃ H ₆₆ NO ₈ P CAS# 998-07-2	100 mg	70.00
Source: synthetic Mol. Wt.: 636 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/acetic acid, 95:5; chloroform/methanol/DI water/acetic acid, 100:30:10:2.5 Storage: -20°C			
1435	1,2-Dipalmitoyl-sn-glycero-3-phosphorylethanolamine DPPE C ₃₇ H ₇₄ NO ₈ P CAS#: 923-61-5	100 mg	70.00
Source: synthetic Mol. Wt.: 692 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/acetic acid, 95:5; chloroform/methanol/DI water/acetic acid, 100:30:10:2.5 Storage: -20°C			
1436	1,2-Distearoyl-sn-glycero-3-phosphorylethanolamine DSPE C ₄₁ H ₈₂ NO ₈ P CAS#: 1069-79-0	100 mg	70.00
Source: synthetic Mol. Wt.: 748 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/acetic acid, 95:5; chloroform/methanol/DI water/acetic acid, 100:30:10:2.5 Storage: -20°C			
1439	1,2-Distearoyl-phosphatidylethanolamine-methyl-polyethyleneglycol conjugate-2000 (Na⁺ salt) DSPE-MPEG-2000 CAS#: 147867-65-0	100 mg	108.00
Source: synthetic Purity: 98+% by TLC Appearance: solid Solubility: chloroform Storage: -20°C			

Phosphatidylinositols

The metabolism of inositol lipids is involved in the signal transduction of many hormones, neurotransmitters and growth factors (1, 2). In the classical pathway, phosphatidylinositol-specific phospholipase C (PI-PLC) hydrolyzes phosphatidyl 4,5-biphosphate (PIP_2) to yield 1,2-diacylglycerol (DAG) and inositol 1,4,5-triphosphate (IP_3). The role of IP_3 and DAG as second messengers is well recognized.

In a second, more recently discovered pathway, the activation of phosphoinositide (PI) 3-kinase results in the formation of three novel phosphatidyl (PI) lipids phosphorylated at the D3 position of the inositol ring: PI-3-P, PI-3,4-P₂ and PI-3,4,5-P₃ (3). These D3 lipids are not known substrates for any of the phospholipase C enzymes and function as second messengers. PI 3-kinase activity is correlated with many cellular processes, including the regulation of cell growth, oncogenic transformation, chemotaxis and receptor down-regulation among others (4, 5, 7). The recent paper on the effect of PI3,4-P₂ on the *Akt* proto-oncogene product (38) also contains protocols for applying PIP's to cell cultures. Matreya's synthetic phosphatidylinositols and inositol phosphates are excellent tools for investigating these second messengers, understanding the enzyme mechanisms involved in phosphoinositide metabolism (39, 40) and for designing therapeutic pharmacological agents. The compounds are evaluated by ¹H and ³¹P NMR to guarantee enantiomeric purity of >98%. See Literature References on page 99.



Catalog number 1779

1779	Phosphatidylinositol, dipalmitoyl, (NH_4^+ salt)	0.5 mg	290.00
1779-1	PI; DPPI (NH_4^+ salt) $\text{C}_{41}\text{H}_{78}\text{O}_{13}\text{P} \cdot \text{NH}_4^+$	1 mg	516.00

Source: synthetic **Mol. Wt.:** 828 **Purity:** 98+% by ¹H NMR, ³¹P NMR

Appearance: solid **Solubility:** chloroform/methanol/DI water, 5:1:0.1 **Storage:** -20°C

1773	Phosphatidylinositol 3-phosphate, dipalmitoyl, (NH_4^+ salt)	100 µg	57.00
1773-1	DPPI-3-P; PI-3-P dipalmitoyl (NH_4^+ salt) $\text{C}_{41}\text{H}_{77}\text{O}_{16}\text{P}_2 \cdot 3\text{NH}_4^+$	1 mg	388.00
1773-5		5 mg	1,815.00

Source: synthetic **Mol. Wt.:** 942 **Purity:** 98+% by ¹H NMR, ³¹P NMR

Appearance: solid **Solubility:** chloroform/methanol/DI water, 1:1:0.3

Storage: -20°C

1919	Phosphatidylinositol 4-phosphate, dipalmitoyl, (NH_4^+ salt)	100 µg	57.00
1919-1	DPPI-4-P; PI-4-P dipalmitoyl (NH_4^+ salt) $\text{C}_{41}\text{H}_{77}\text{O}_{16}\text{P}_2 \cdot 3\text{NH}_4^+$	1 mg	388.00
1919-5		5 mg	1,815.00

Source: synthetic **Mol. Wt.:** 942 **Purity:** 98+% by ¹H NMR, ³¹P NMR

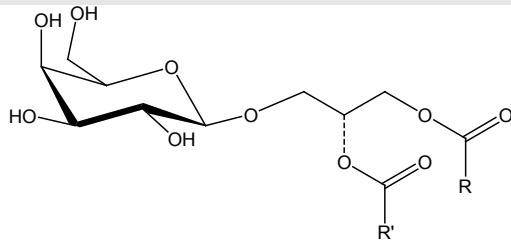
Appearance: solid **Solubility:** methanol, chloroform/methanol/DI water, 1:1:0.3; slightly soluble in DI water **Storage:** -20°C

1784	Phosphatidylinositol bis-4,5-phosphate, dioctanoyl, (NH_4^+ salt)		100 µg	57.00
1784-1	DOPI-4,5-P2; PI-4,5-P2 dioctanoyl (NH_4^+ salt) C ₂₅ H ₄₉ O ₁₉ P ₃ •5NH ₄		1 mg	388.00
1784-5			5 mg	1,815.00
	Source: synthetic Mol. Wt.: 831 Purity: 98+% by ¹ H NMR, ³¹ P NMR			
	Appearance: solid Solubility: chloroform/methanol/DI water, 1:1:0.3			
	Storage: -20°C			
1778	Phosphatidylinositol bis-4,5-phosphate, dioctanoyl, (Na^+ salt)		100 µg	57.00
1778-1	DOPI-4,5-P2; PI-4,5-P2 dioctanoyl (Na^+ salt) C ₂₅ H ₄₄ O ₁₉ P ₃ •5Na		1 mg	388.00
1778-5			5 mg	1,815.00
	Source: synthetic Mol. Wt.: 856 Purity: 98+% by ¹ H NMR, ³¹ P NMR			
	Appearance: solid Solubility: DI water Storage: -20°C			
1783	Phosphatidylinositol tris-3,4,5-phosphate, dipalmitoyl, (NH_4^+ salt)		100 µg	49.00
1783-1	DPPI-3,4,5-P3; PI-3,4,5-P3 dipalmitoyl (NH_4^+ salt) C ₄₁ H ₇₅ O ₂₂ P ₄ •7NH ₄		1 mg	363.00
1783-5			5 mg	1,573.00
	Source: synthetic Mol. Wt.: 1170 Purity: 98+% by ¹ H NMR, ³¹ P NMR			
	Appearance: solid Solubility: chloroform/methanol/DI water, 1:1:0.3			
	Storage: -20°C			
1775	Phosphatidylinositol tris-3,4,5-phosphate, dipalmitoyl, (Na^+ salt)		100 µg	43.00
1775-1	DPPI-3,4,5-P3; PI-3,4,5-P3 dipalmitoyl (Na^+ salt) C ₄₁ H ₇₅ O ₂₂ P ₄ •7Na		1 mg	303.00
1775-5			5 mg	1,392.00
	Source: synthetic Mol. Wt.: 1205 Purity: 98+% by ¹ H NMR, ³¹ P NMR			
	Appearance: solid Solubility: DI water Storage: -20°C			

Bacterial Tetraethers

1303	Main phospholipid (MPL) of <i>Thermoplasma acidophilum</i>, (>95% pure)		5 mg	790.00
	Purified MPL of <i>Thermoplasma acidophilum</i> (>95% pure) C ₉₅ H ₁₈₈ O ₁₆ P			
	Source: natural, Archaebacteria Mol. Wt.: 1618 Purity: >95% by TLC, HPLC			
	Appearance: solid Solubility: chloroform/methanol, 2:1; hexane/2-propanol/DI water, 30:40:5 Storage: 4-8°C			
1303-2	Main phospholipid (MPL) of <i>Thermoplasma acidophilum</i>, (>50% pure)		50 mg	2,024.00
	MPL of <i>Thermoplasma acidophilum</i> (>50% pure) C ₉₅ H ₁₈₈ O ₁₆ P			
	Source: natural, Archaebacteria Mol. Wt.: 1618 Purity: >50% by TLC			
	Appearance: liquid Solubility: chloroform/methanol, 2:1; hexane/2-propanol/DI water, 30:40:5 Storage: 4-8°C highly hygroscopic			

Glycosyl Glycerides



Catalog number 1058

1058	Monogalactosyldiglyceride MGDG (hydrogenated) C ₄₅ H ₈₆ O ₁₀ CAS#: 41670-62-6	10 mg	197.00
------	--	-------	--------

Source: natural, plant Mol. Wt.: 787 (stearoyl) Purity: 98+% by TLC
Appearance: solid Solubility: chloroform/methanol/DI water, 4:1:0.1
Storage: -20°C

1059	Digalactosyldiglyceride DGDG (hydrogenated) C ₅₁ H ₉₆ O ₁₅ CAS#: 92457-02-8	5 mg	234.00
------	--	------	--------

Source: natural, plant Mol. Wt.: 949 (distearoyl) Purity: 98+% by TLC
Appearance: solid Solubility: chloroform/methanol/DI water, 4:1:0.1
Storage: -20°C

Fatty Acids

Simple Fatty Acids

Saturated Fatty Acids and Methyl Esters

These products are 99% pure by GC. They are stable at room temperature and are supplied neat.

1200	Methyl hexanoate Methyl caproate; C6:0 Methyl ester C ₇ H ₁₄ O ₂ CAS#: 106-70-7	1 g	42.00
------	--	-----	-------

Source: natural, plant Mol. Wt.: 130 Purity: 99% by TLC, GC Appearance: liquid
Solubility: chloroform, ethanol, ethyl ether Storage: room temperature

1196	Heptanoic acid C7:0 Fatty acid C ₇ H ₁₄ O ₂ CAS#: 111-14-8	1 g	66.00
------	---	-----	-------

Source: natural, plant Mol. Wt.: 130 Purity: 99% by TLC, GC Appearance: liquid
Solubility: chloroform, ethanol, ethyl ether Storage: room temperature

1197	Methyl heptanoate C7:0 Methyl ester C ₈ H ₁₆ O ₂ CAS#: 106-73-0	1 g	66.00
------	--	-----	-------

Source: natural, plant Mol. Wt.: 144 Purity: 99% by TLC, GC Appearance: liquid
Solubility: chloroform, ethanol, ethyl ether Storage: room temperature

1198	Octanoic acid Caprylic acid; C8:0 Fatty acid C ₈ H ₁₆ O ₂ CAS#: 124-07-2	1 g	42.00
	Source: natural, plant Mol. Wt.: 144 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		
1199	Methyl octanoate Methyl caprylate; C8:0 Methyl ester C ₉ H ₁₈ O ₂ CAS#: 111-11-5	1 g	42.00
	Source: natural, plant Mol. Wt.: 158 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		
1163	Nonanoic acid C9:0 Fatty acid; Pelargonic acid C ₉ H ₁₈ O ₂ CAS#: 112-05-0	100 mg	33.00
	Source: synthetic Mol. Wt.: 158 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		
1164	Methyl nonanoate C9:0 Methyl ester C ₁₀ H ₂₀ O ₂ CAS#: 1731-84-6	100 mg	33.00
	Source: synthetic Mol. Wt.: 172 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		
1261	Methyl decanoate Methyl caprate; C10:0 Methyl ester C ₁₁ H ₂₂ O ₂ CAS#: 110-42-9	500 mg	33.00
	Source: natural, plant Mol. Wt.: 186 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane Storage: room temperature		
1165	Undecanoic acid C11:0 Fatty acid C ₁₁ H ₂₂ O ₂ CAS#: 112-37-8	100 mg	33.00
	Source: synthetic Mol. Wt.: 186 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		
1166	Methyl undecanoate C11:0 Methyl ester C ₁₂ H ₂₄ O ₂ CAS#: 1731-86-8	100 mg	33.00
	Source: synthetic Mol. Wt.: 200 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		
1008	Dodecanoic acid Lauric acid; C12:0 Fatty acid C ₁₂ H ₂₄ O ₂ CAS#: 143-07-7	1 g	42.00
	Source: natural, plant Mol. Wt.: 200 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		
1009	Methyl dodecanoate Methyl laurate; C12:0 Methyl ester C ₁₃ H ₂₆ O ₂ CAS#: 111-82-0	1 g	42.00
	Source: natural, plant Mol. Wt.: 214 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		

1161	Tridecanoic acid C13:0 Fatty acid C ₁₃ H ₂₆ O ₂ CAS#: 638-53-9	100 mg	60.00
	Source: synthetic Mol. Wt.: 214 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		
1162	Methyl tridecanoate C13:0 Methyl ester C ₁₄ H ₂₈ O ₂ CAS#: 1731-88-0	100 mg	60.00
	Source: synthetic Mol. Wt.: 228 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		
1010	Tetradecanoic acid Myristic acid; C14:0 Fatty acid C ₁₄ H ₂₈ O ₂ CAS#: 544-63-8	1 g	39.00
	Source: natural, plant Mol. Wt.: 228 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		
1011	Methyl tetradecanoate Methyl myristate; C14:0 Methyl ester C ₁₅ H ₃₀ O ₂ CAS#: 124-10-7	1 g	39.00
	Source: natural, plant Mol. Wt.: 242 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		
1012	Pentadecanoic acid C15:0 Fatty acid C ₁₅ H ₃₀ O ₂ CAS#: 1002-84-2	1 g	97.00
	Source: synthetic Mol. Wt.: 242 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		
1013	Methyl pentadecanoate C15:0 Methyl ester C ₁₆ H ₃₂ O ₂ CAS#: 7132-64-1	1 g	97.00
	Source: synthetic Mol. Wt.: 256 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		
1014	Hexadecanoic acid Palmitic acid; C16:0 Fatty acid C ₁₆ H ₃₂ O ₂ CAS#: 57-10-3	1 g	33.00
	Source: natural, plant Mol. Wt.: 256 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		
1015	Methyl hexadecanoate Methyl palmitate; C16:0 Methyl ester C ₁₇ H ₃₄ O ₂ CAS#: 112-39-0	1 g	33.00
	Source: natural, plant Mol. Wt.: 270 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		
1018	Heptadecanoic acid Margaric acid; C17:0 Fatty acid C ₁₇ H ₃₄ O ₂ CAS#: 506-12-7	1 g	146.00
	Source: synthetic Mol. Wt.: 270 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		

1019	Methyl heptadecanoate Methyl margarate; C17:0 Methyl ester C ₁₈ H ₃₆ O ₂ CAS#: 1731-92-6	1 g	146.00
	Source: synthetic Mol. Wt.: 284 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		
1020	Octadecanoic acid Stearic acid; C18:0 Fatty acid C ₁₈ H ₃₆ O ₂ CAS#: 57-11-4	1 g	33.00
	Source: natural, plant Mol. Wt.: 284 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		
1021	Methyl octadecanoate Methyl stearate; C18:0 Methyl ester C ₁₉ H ₃₈ O ₂ CAS#: 112-61-8	1 g	33.00
	Source: natural, plant Mol. Wt.: 298 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, ethyl ether Storage: room temperature		
1028	Nonadecanoic acid C19:0 Fatty acid C ₁₉ H ₃₈ O ₂ CAS#: 646-30-0	100 mg	71.00
	Source: synthetic Mol. Wt.: 298 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether Storage: room temperature		
1029	Methyl nonadecanoate C19:0 Methyl ester C ₂₀ H ₄₀ O ₂ CAS#: 1731-94-8	100 mg	71.00
	Source: synthetic Mol. Wt.: 312 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether Storage: room temperature		
1030	Eicosanoic acid Arachidic acid; C20:0 Fatty acid C ₂₀ H ₄₀ O ₂ CAS#: 506-30-9	500 mg	71.00
	Source: natural, plant Mol. Wt.: 312 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether Storage: room temperature		
1031	Methyl eicosanoate Methyl arachidate; C20:0 Methyl ester C ₂₁ H ₄₂ O ₂ CAS#: 1120-28-1	500 mg	71.00
	Source: natural, plant Mol. Wt.: 327 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether Storage: room temperature		
1241	Heneicosanoic acid C21:0 Fatty acid C ₂₁ H ₄₂ O ₂ CAS#: 2363-71-5	100 mg	66.00
	Source: synthetic Mol. Wt.: 326 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether Storage: room temperature		
1242	Methyl heneicosanoate C21:0 Methyl ester C ₂₂ H ₄₄ O ₂ CAS#: 6064-90-0	100 mg	66.00
	Source: synthetic Mol. Wt.: 341 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether Storage: room temperature		

1035	Docosanoic acid Behenic acid; C22:0 Fatty acid C ₂₂ H ₄₄ O ₂ CAS#: 112-85-6	500 mg	66.00
	Source: natural, plant Mol. Wt.: 341 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether Storage: room temperature		
1036	Methyl docosanoate Methyl behenate; C22:0 Methyl ester C ₂₃ H ₄₆ O ₂ CAS#: 929-77-1	500 mg	66.00
	Source: natural, plant Mol. Wt.: 354 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether Storage: room temperature		
1186	Tricosanoic acid C23:0 Fatty acid C ₂₃ H ₄₆ O ₂ CAS#: 2433-96-7	100 mg	77.00
	Source: synthetic Mol. Wt.: 355 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether Storage: room temperature		
1187	Methyl tricosanoate C23:0 Methyl ester C ₂₄ H ₄₈ O ₂ CAS#: 2433-97-8	100 mg	77.00
	Source: synthetic Mol. Wt.: 368 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether Storage: room temperature		
1037	Tetracosanoic acid Lignoceric acid; C24:0 Fatty acid C ₂₄ H ₄₈ O ₂ CAS#: 557-59-5	100 mg	77.00
	Source: synthetic Mol. Wt.: 369 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether Storage: room temperature		
1038	Methyl tetracosanoate Methyl lignocerate; C24:0 Methyl ester C ₂₅ H ₅₀ O ₂ CAS#: 2442-49-1	100 mg	77.00
	Source: synthetic Mol. Wt.: 382 Purity: 99% by GC Appearance: solid Solubility: chloroform, ethyl ether Storage: room temperature		
1251	Hexacosanoic acid Cerotic acid; C26:0 Fatty acid C ₂₆ H ₅₂ O ₂ CAS#: 506-46-7	25 mg	89.00
	Source: synthetic Mol. Wt.: 397 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether Storage: room temperature		
1252	Methyl hexacosanoate Methyl cerotate; C26:0 Methyl ester C ₂₇ H ₅₄ O ₂ CAS#: 5802-82-4	25 mg	89.00
	Source: synthetic Mol. Wt.: 411 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether Storage: room temperature		
1271	Methyl octacosanoate Methyl montanate; C28:0 Methyl ester C ₂₉ H ₅₈ O ₂ CAS#: 55682-92-3	50 mg	88.00
	Source: synthetic Mol. Wt.: 439 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, methylene chloride Storage: room temperature		

1273	Methyl triacontanoate Methyl melissate; C30:0 Methyl ester C ₃₁ H ₆₂ O ₂ CAS#: 629-83-4	50 mg	99.00
	Source: synthetic Mol. Wt.: 467 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, methylene chloride Storage: room temperature		
1275	Methyl dotriacontanoate Methyl lacceroate; C32:0 Methyl ester C ₃₃ H ₆₆ O ₂ CAS#: 41755-79-7	50 mg	111.00
	Source: synthetic Mol. Wt.: 495 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, methylene chloride Storage: room temperature		
Unsaturated Fatty Acids and Methyl Esters			
Unsaturated fatty acids are easily oxidized. Flush open containers with argon or nitrogen and store at -20°C, in dark.			
1157	Tetradecenoic acid (<i>cis</i>-9) Myristoleic acid; C14:1 (<i>cis</i> -9) fatty acid C ₁₄ H ₂₆ O ₂ CAS#: 544-64-9	100 mg	92.00
	Source: natural, plant Mol. Wt.: 226 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1040	Methyl tetradecenoate (<i>cis</i>-9) Methyl myristoleate; C14:1 (<i>cis</i> -9) Methyl ester C ₁₅ H ₂₈ O ₂ CAS#: 56219-06-8	100 mg	92.00
	Source: natural, plant Mol. Wt.: 240 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1243	Hexadecenoic acid (<i>cis</i>-6) Sapienic acid C ₁₆ H ₃₀ O ₂	25 mg	174.00
	Source: synthetic Mol. Wt.: 254 Purity: 99% by TLC, GC Appearance: liquid Solubility: ethanol, methanol, chloroform, ethyl ether Storage: -20°C		
1016	Hexadecenoic acid (<i>cis</i>-9) Palmitoleic acid; C16:1 (<i>cis</i> -9) Fatty acid C ₁₆ H ₃₀ O ₂ CAS#: 373-49-9	100 mg	57.00
	Source: natural, plant Mol. Wt.: 254 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1017	Methyl hexadecenoate (<i>cis</i>-9) Methyl palmitoleate; C16:1 (<i>cis</i> -9) Methyl ester C ₁₇ H ₃₂ O ₂ CAS#: 1120-25-8	100 mg	57.00
	Source: natural, plant Mol. Wt.: 268 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1147	Hexadecenoic acid (<i>trans</i>-9) Palmitelaidic acid; C16:1 (<i>trans</i> -9) Fatty acid C ₁₆ H ₃₀ O ₂ CAS#: 10030-73-6	100 mg	77.00
	Source: synthetic Mol. Wt.: 254 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		

1148	Methyl hexadecenoate (<i>trans</i>-9) Methyl palmitelaidate; C16:1 (<i>trans</i> -9) Methyl ester C ₁₇ H ₃₂ O ₂ CAS#: 10030-74-7	100 mg	77.00
	Source: synthetic Mol. Wt.: 268 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1208	11-Hexadecenoic acid (92% <i>cis</i>, 8% <i>trans</i>) C16:1 (<i>cis</i> -11) Fatty acid C ₁₆ H ₃₀ O ₂	50 mg	131.00
	Source: synthetic Mol. Wt.: 254 Purity: >98%, by TLC Appearance: liquid Solubility: chloroform, ethanol, hexane, methanol Storage: -20°C 92% <i>cis</i> , 8% <i>trans</i> by GC		
1204	Heptadecenoic acid (<i>cis</i>-10) C17:1 (<i>cis</i> -10) Fatty acid C ₁₇ H ₃₂ O ₂ CAS#: 29743-97-3	100 mg	73.00
	Source: synthetic Mol. Wt.: 268 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1203	Methyl heptadecenoate (<i>cis</i>-10) Methyl heptadecenoate; C17:1 (<i>cis</i> -10) Methyl ester C ₁₈ H ₃₄ O ₂ CAS#: 75190-82-8	100 mg	73.00
	Source: synthetic Mol. Wt.: 282 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1022	Octadecenoic acid (<i>cis</i>-9) Oleic acid; C18:1 (<i>cis</i> -9) Fatty acid C ₁₈ H ₃₄ O ₂ CAS#: 112-80-1	1 g	46.00
	Source: natural, plant Mol. Wt.: 282 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1023	Methyl octadecenoate (<i>cis</i>-9) Methyl oleate; C18:1 (<i>cis</i> -9) Methyl ester C ₁₉ H ₃₆ O ₂ CAS#: 112-62-9	1 g	46.00
	Source: natural, plant Mol. Wt.: 296 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1149	Octadecenoic acid (<i>trans</i>-9) Elaidic acid; C18:1 (<i>trans</i> -9) Fatty acid C ₁₈ H ₃₄ O ₂ CAS#: 112-79-8	1 g	105.00
	Source: synthetic Mol. Wt.: 282 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1150	Methyl octadecenoate (<i>trans</i>-9) Methyl elaidate; C18:1 (<i>trans</i> -9) Methyl ester C ₁₉ H ₃₆ O ₂ CAS#: 1937-62-8	1 g	105.00
	Source: natural, plant Mol. Wt.: 296 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		

1266	Octadecenoic acid (<i>cis</i>-11) <i>cis</i> -Vaccenic acid; C18:1(<i>cis</i> -11) Fatty acid C ₁₈ H ₃₄ O ₂ CAS#: 506-17-2	100 mg	92.00
	Source: natural, plant Mol. Wt.: 282 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1267	Methyl octadecenoate (<i>cis</i>-11) Methyl <i>cis</i> -vaccenate; C18:1(<i>cis</i> -11) Methyl ester C ₁₉ H ₃₆ O ₂ CAS#: 1937-63-9	100 mg	92.00
	Source: semisynthetic, plant Mol. Wt.: 296 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1262	Octadecenoic acid (<i>trans</i>-11) <i>trans</i> Vaccenic acid; C18:1 (<i>trans</i> -11) Fatty acid C ₁₈ H ₃₄ O ₂ CAS#: 693-72-1	100 mg	92.00
	Source: synthetic Mol. Wt.: 282 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1263	Methyl octadecenoate (<i>trans</i>-11) Methyl <i>trans</i> vaccenate; C18:1 (<i>trans</i> -11) Methyl ester C ₁₉ H ₃₆ O ₂ CAS#: 6198-58-9	100 mg	92.00
	Source: synthetic Mol. Wt.: 296 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1024	Octadecadienoic acid (all <i>cis</i>-9,12) Linoleic acid; C18:2 (all <i>cis</i> -9,12) Fatty acid C ₁₈ H ₃₂ O ₂ CAS#: 60-33-3	1 g	46.00
	Source: natural, plant Mol. Wt.: 280 Purity: 99% by TLC, GC Appearance: liquid Solubility: ethyl ether, ethanol, hexane Storage: -20°C		
1025	Methyl octadecadienoate (all <i>cis</i>-9,12) Methyl linoleate; C18:2 (all <i>cis</i> -9,12) Methyl ester C ₁₉ H ₃₄ O ₂ CAS#: 112-63-0	1 g	46.00
	Source: natural, plant Mol. Wt.: 294 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1151	Linoelaidic acid (all <i>trans</i>-9,12) C18:2 (all <i>trans</i> -9, 12) Fatty acid C ₁₈ H ₃₂ O ₂ CAS#: 506-21-8	100 mg	57.00
	Source: natural, plant Mol. Wt.: 280 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1152	Methyl octadecadienoate (all <i>trans</i>-9,12) Methyl linoelaidate; C18:2 (all <i>trans</i> -9,12) Methyl ester C ₁₉ H ₃₄ O ₂ CAS#: 2566-97-4	100 mg	57.00
	Source: natural, plant Mol. Wt.: 294 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		

1026	Octadecatrienoic acid (all <i>cis</i>-9,12,15) Linolenic acid; C18:3 (all <i>cis</i> -9,12,15) Fatty acid C ₁₈ H ₃₀ O ₂ CAS#: 463-40-1	500 mg	105.00
	Source: natural, plant Mol. Wt.: 278 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1027	Methyl octadecatrienoate (all <i>cis</i>-9,12,15) Methyl linolenate; C18:3 (all <i>cis</i> -9,12,15) Methyl ester C ₁₉ H ₃₂ O ₂ CAS#: 301-00-8	500 mg	105.00
	Source: natural, plant Mol. Wt.: 292 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1153	Octadecatrienoic acid (all <i>cis</i>-6,9,12) <i>gamma</i> -Linolenic acid; C18:3 (all <i>cis</i> -6,9,12) Fatty acid C ₁₈ H ₃₀ O ₂ CAS#: 506-26-3	100 mg	64.00
	Source: natural, plant Mol. Wt.: 278 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1154	Methyl octadecatrienoate (all <i>cis</i>-6,9,12) Methyl <i>gamma</i> -linolenate; C18:3 (all <i>cis</i> -6,9,12) Methyl ester C ₁₉ H ₃₂ O ₂ CAS#: 16326-32-2	100 mg	64.00
	Source: natural, plant Mol. Wt.: 292 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1276	Stearidonic acid (all <i>cis</i>-6,9,12,15) Morocitic acid; C18:4 (all <i>cis</i> -6,9,12,15) Fatty acid C ₁₈ H ₂₈ O ₂ CAS#: 20290-75-9	25 mg	264.00
	Source: natural, plant Mol. Wt.: 276 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C Dry Ice Charge Applies		
1277	Methyl stearidonate (all <i>cis</i>-6,9,12,15) Morocitic acid methyl ester; C18:4 (all <i>cis</i> -6,9,12,15) Methyl ester C ₁₉ H ₃₀ O ₂ CAS#: 73097-00-4	25 mg	264.00
	Source: natural, plant Mol. Wt.: 290 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C Dry Ice Charge Applies		
1205	Nonadecenoic acid (<i>cis</i>-10) C19:1 (<i>cis</i> -10) Fatty acid C ₁₉ H ₃₆ O ₂ CAS#: 73033-09-7	100 mg	83.00
	Source: synthetic Mol. Wt.: 296 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1206	Methyl nonadecenoate (<i>cis</i>-10) C19:1 (<i>cis</i> -10) Methyl ester C ₂₀ H ₃₈ O ₂ CAS#: 19788-74-0	100 mg	83.00
	Source: synthetic Mol. Wt.: 310 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		

1032	Eicosenoic acid (<i>cis</i>-11) C20:1 (<i>cis</i> -11) Fatty acid C ₂₀ H ₃₈ O ₂ CAS#: 5561-99-9	100 mg	57.00
	Source: natural, plant Mol. Wt.: 310 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1033	Methyl eicosenoate (<i>cis</i>-11) Methyl eicosenoate; C20:1 (<i>cis</i> -11) Methyl ester C ₂₁ H ₄₀ O ₂ CAS#: 2390-09-2	100 mg	57.00
	Source: natural, plant Mol. Wt.: 324 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1192	Eicosadienoic acid (all <i>cis</i>-11,14) C20:2 (all <i>cis</i> -11,14) Fatty acid C ₂₀ H ₃₆ O ₂ CAS#: 2091-39-6	100 mg	64.00
	Source: synthetic Mol. Wt.: 309 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1193	Methyl eicosadienoate (all <i>cis</i>-11,14) Methyl eicosadienoate; C20:2 (all <i>cis</i> -11,14) Methyl ester C ₂₁ H ₃₈ O ₂ CAS#: 2463-02-7	100 mg	64.00
	Source: synthetic Mol. Wt.: 322 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1179	Methyl eicosatrienoate (all <i>cis</i>-5,8,11) Mead acid methyl ester; C20:3 (all <i>cis</i> -5,8,11) Methyl ester C ₂₁ H ₃₆ O ₂ CAS#: 14602-39-2	1 mg/ml, 1 ml	180.00
	Source: natural, plant Mol. Wt.: 320 Purity: 90% by TLC, GC Appearance: liquid Solvent: hexane Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1269	Methyl eicosatrienoate (all <i>cis</i>-8,11,14) Methyl homogamma linolenate; C20:3 (all <i>cis</i> -8,11,14) Methyl ester C ₂₁ H ₃₆ O ₂	50 mg	99.00
	Source: semi-synthetic, plant Mol. Wt.: 321 Purity: 99% by TLC, GC Appearance: liquid Solubility: hexane, ethyl ether, chloroform Storage: -20°C		
1042	Arachidonic acid (all <i>cis</i>-5,8,11,14) C20:4 (all <i>cis</i> -5,8,11,14) Fatty acid C ₂₀ H ₃₂ O ₂ CAS#: 506-32-1	100 mg	88.00
	Source: natural, fungal Mol. Wt.: 304 Purity: 99% by TLC, GC Appearance: liquid Solubility: ethyl ether, hexane, methylene chloride Storage: -20°C Dry Ice Charge Applies		
1034	Methyl eicosatetraenoate (all <i>cis</i>-5,8,11,14) Methyl arachidonate; C20:4 (all <i>cis</i> -5,8,11,14) Methyl ester C ₂₁ H ₃₄ O ₂ CAS#: 2566-89-4	100 mg	88.00
	Source: natural, fungal Mol. Wt.: 318 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C Dry Ice Charge Applies		

1167	Eicosapentaenoic acid (all <i>cis</i>-5,8,11,14,17) EPA ; <i>omega</i> -3 Fatty acid; C20:5 (all <i>cis</i> -5,8,11,14,17) Fatty acid C ₂₀ H ₃₀ O ₂ CAS#: 10417-94-4	25 mg	97.00
	Source: natural, fish oil Mol. Wt.: 302 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether, ethanol, DMSO, DMF Storage: -20°C Dry Ice Charge Applies		
	Anti-hyperlipoproteinemic agent; 5-LOX inhibitor		
1194	Methyl eicosapentaenoate (all <i>cis</i>-5,8,11,14,17) Methyl ester of <i>omega</i> -3 Fatty acid; C20:5 (all <i>cis</i> -5,8,11,14,17) Methyl ester C ₂₁ H ₃₂ O ₂ CAS#: 2734-47-6	25 mg	97.00
	Source: natural, fish oil Mol. Wt.: 316 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, ethyl ether, hexane Storage: -20°C Dry Ice Charge Applies		
1264	Docosenoic acid (<i>cis</i>-13) Erucic acid; C22:1 (<i>cis</i> -13) Fatty acid C ₂₂ H ₄₂ O ₂ CAS#: 112-86-7	100 mg	49.00
	Source: natural, plant Mol. Wt.: 339 Purity: >99% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether, hexane Storage: -20°C		
1265	Methyl docosenoate (<i>cis</i>-13) Methyl erucate; C22:1 (<i>cis</i> -13) Methyl ester C ₂₃ H ₄₄ O ₂ CAS#: 1120-34-9	100 mg	49.00
	Source: natural, plant Mol. Wt.: 352 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, ethyl ether, hexane Storage: -20°C		
1175	Docosapentaenoic acid (all <i>cis</i>-7,10,13,16,19) C22:5 (all <i>cis</i> -7,10,13,16,19) Fatty acid C ₂₂ H ₃₄ O ₂ CAS#: 24880-45-3	25 mg	97.00
	Source: semi-synthetic Mol. Wt.: 330 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, ethyl ether, hexane Storage: -20°C Dry Ice Charge Applies		
1244	Methyl docosapentaenoate (all <i>cis</i>-7,10,13,16,19) C22:5 (all <i>cis</i> -7,10,13,16,19) Methyl ester C ₂₃ H ₃₆ O ₂ CAS#: 108698-02-8	25 mg	97.00
	Source: semi-synthetic Mol. Wt.: 344 Purity: 99% by TLC, GC Appearance: liquid Solubility: ethyl ether, ethanol, hexane, Storage: -20°C Dry Ice Charge Applies		
1136	Docosahexaenoic acid (all <i>cis</i>-4,7,10,13,16,19) DHA; C22:6 (all <i>cis</i> -4,7,10,13,16,19) <i>omega</i> -3 Fatty acid C ₂₂ H ₃₂ O ₂ CAS#: 6217-54-5	100 mg	88.00
	Source: natural, algae Mol. Wt.: 328 Purity: 99% by TLC, GC Appearance: liquid Solubility: ethyl ether, hexane, methylene chloride, ethanol, DMSO, DMF Storage: -20°C Dry Ice Charge Applies		

1041	Methyl docosahexaenoate (all cis-4,7,10,13,16,19) Methyl ester of omega-3 fatty acid; C22:6 (all <i>cis</i> -4,7,10,13,16,19) Methyl ester C ₂₃ H ₃₄ O ₂ CAS#: 2566-90-7	100 mg	88.00
	Source: natural, algae Mol. Wt.: 342 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C Dry Ice Charge Applies		
1155	Nervonic acid (cis-15) Tetracosenoic acid (<i>cis</i> -15); C24:1 (<i>cis</i> -15) Fatty acid C ₂₄ H ₄₆ O ₂ CAS#: 506-37-6	100 mg	88.00
	Source: synthetic Mol. Wt.: 367 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1156	Methyl tetracosenoate (cis-15) Methyl nervonate; C24:1 (<i>cis</i> -15) Methyl ester C ₂₅ H ₄₈ O ₂ CAS#: 2733-88-2	100 mg	88.00
	Source: synthetic Mol. Wt.: 381 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		

Trans Fatty Acids and Methyl Esters

1147	Hexadecenoic acid (trans-9) Palmitelaidic acid; C16:1 (<i>trans</i> -9) Fatty acid C ₁₆ H ₃₀ O ₂ CAS#: 10030-73-6	100 mg	77.00
	Source: synthetic Mol. Wt.: 254 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1148	Methyl hexadecenoate (trans-9) Methyl palmitelaidate; C16:1 (<i>trans</i> -9) Methyl ester C ₁₇ H ₃₂ O ₂ CAS#: 10030-74-7	100 mg	77.00
	Source: synthetic Mol. Wt.: 268 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1149	Octadecenoic acid (trans-9) Elaidic acid; C18:1 (<i>trans</i> -9) Fatty acid C ₁₈ H ₃₄ O ₂ CAS#: 112-79-8	1 g	105.00
	Source: synthetic Mol. Wt.: 282 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1150	Methyl octadecenoate (trans-9) Methyl elaidate; C18:1 (<i>trans</i> -9) Methyl ester C ₁₉ H ₃₆ O ₂ CAS#: 1937-62-8	1 g	105.00
	Source: natural, plant Mol. Wt.: 296 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1262	Octadecenoic acid (trans-11) <i>trans</i> Vaccenic acid; C18:1 (<i>trans</i> -11) Fatty acid C ₁₈ H ₃₄ O ₂ CAS#: 693-72-1	100 mg	92.00
	Source: synthetic Mol. Wt.: 282 Purity: 99% by TLC, GC Appearance: solid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		

1263	Methyl octadecenoate (<i>trans</i>-11) Methyl <i>trans</i> vaccenate; C18:1 (<i>trans</i> -11) Methyl ester C ₁₉ H ₃₆ O ₂ CAS#: 6198-58-9	100 mg	92.00
	Source: synthetic Mol. Wt.: 296 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1151	Linoelaidic acid (all <i>trans</i>-9,12) C18:2 (all <i>trans</i> -9,12) Fatty acid C ₁₈ H ₃₂ O ₂ CAS#: 506-21-8	100 mg	57.00
	Source: natural, plant Mol. Wt.: 280 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1152	Methyl octadecadienoate (all <i>trans</i>-9,12) Methyl linoelaidate; C18:2 (all <i>trans</i> -9,12) Methyl ester C ₁₉ H ₃₄ O ₂ CAS#: 2566-97-4	100 mg	57.00
	Source: natural, plant Mol. Wt.: 294 Purity: 99% by TLC, GC Appearance: liquid Solubility: chloroform, hexane, ethyl ether Storage: -20°C		
1131	Cis-Trans Isomer Standard Mixture	5 mg/ml, 5 ml	135.00
	Source: margarine Appearance: liquid Solvent: 5ml methylene chloride Solubility: methylene chloride, chloroform Storage: -20°C		
	Analysis of positional cis-trans fatty acid isomers is ever more important in light of the new food industry rules. These isomers can be resolved on Supelco SP-2560 or an equivalent capillary GC column. Use this specially formulated mix to ensure proper operation of your column for this tricky separation. Mix consists of cis-trans fatty acid isomers as methyl esters in methylene chloride.		
	This is a qualitative standard containing in order of elution: C16:0, C18:0, C18:1 trans isomers (4 peaks), C18:1 cis & trans isomers (2 peaks), C18:1 cis isomers (4 peaks), C18:2, C20:0, C20:1 and C18:3 (same peak), C22:0		
1181	9(E),11(E)-Octadecadienoic acid 9- <i>trans</i> , 11- <i>trans</i> CLA C ₁₈ H ₃₂ O ₂ CAS#: 544-71-8	25 mg	111.00
	Source: synthetic Mol. Wt.: 280 Melting Point (°C): 55-57 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, hexane, methanol Storage: -20°C		

Conjugated Linoleic Acid Isomers (CLA)

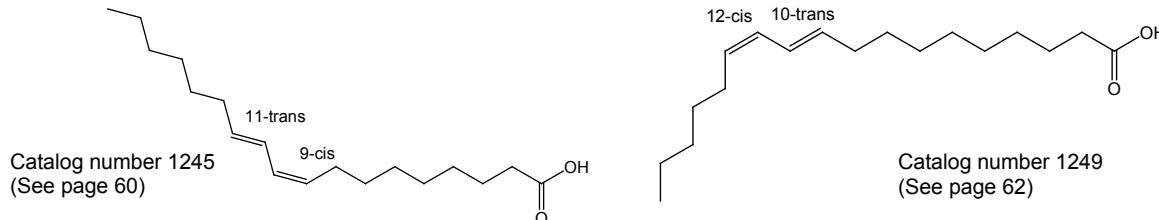
Linoleic acid is an essential fatty acid (18:2 ω6) of which several naturally occurring conjugated derivatives have been identified. These derivatives, called "conjugated linoleic acid" or CLA can have the two double bonds mainly in the 9 and 11 or in the 10 and 12 positions, resulting in eight possible geometric isomers. CLA occurs in meat (41) and dairy products (42, 43). In both cases, the 9(Z),11(E)-isomer is predominant and is thought to be the biologically active form. CLA assimilated through the diet of animals is found in the intestinal musosa, liver and adipose tissue (44). See also review article by Parodi (43). CLA has several biological properties. Its anti-carcinogenic activity has been demonstrated by its ability to inhibit chemically induced tumor formation in animal models of carcinogenesis (41,45-47). The addition of CLA to culture medium suppresses the *in vitro* growth of human melanoma, colorectal and breast cancer cells (48). CLA also exhibits anti-atherogenic activity. Addition of CLA to a controlled atherogenic diet significantly reduced the development of atherosclerosis in hamsters and rabbits (49,50). Animals fed a diet containing CLA also had lower levels of low-density-lipoprotein (LDL) cholesterol. CLA may be involved in regulating fat and protein metabolism (51,52). Several species of animals fed CLA-supplemented diets showed improved feed efficiency. Lean body mass increased while body fat was reduced. This seems to be due, mainly or exclusively, to the 10(E),12(Z)-isomer (catalog # 1249, see below). CLA competes with linoleate for Δ6 desaturase (53). Dietary CLA normalizes impaired glucose tolerance in the Zucker diabetic fatty *fa/fa* rat (54) via activation of PPAR γ, a result which bears on the possible amelioration or prevention of NIDDM. The 11(Z),13(E)-isomer (catalog # 1259) has been shown to be concentrated in the heart and in mitochondria.

See Literature References on page 99.

CLA Research is Being Redone With Our Highly Pure Isomers

Most studies to date have utilized a mixture of CLA isomers containing less than 30% of the presumed active 9(Z),11(E)-isomer (55,56). In addition to the 9,11- and 10,12-isomers, 8,10- and 11,13-isomers have recently been identified in the widely used mixture (56,57). Matreya offers a highly pure CLA which is 98+% the active 9,11-“cis, trans” isomer. The corresponding “trans,trans” and “cis,cis” isomers are also available. In addition, we now offer the pure 10(E),12(Z)-isomer, which has been widely sought for comparison studies.

See Literature References on page 99.



1245	9(Z),11(E)-Octadecadienoic acid	25 mg	111.00
1245-1	9-cis, 11-trans CLA C ₁₈ H ₃₂ O ₂ CAS#: 2540-56-9	1 g	353.00
1245-10		10 g	1,707.00

Source: synthetic Mol. Wt.: 280 Purity: 98% by TLC, GC Appearance: liquid
Solubility: chloroform, ethanol, hexane, methanol Storage: -20°C

1255	Methyl 9(Z), 11(E)-octadecadienoate	25 mg	111.00
	Methyl ester of CLA (9-cis, 11-trans) C ₁₉ H ₃₄ O ₂		

Source: synthetic Mol. Wt.: 294 Purity: 98% by TLC, GC Appearance: liquid
Solubility: chloroform, ethanol, hexane, methanol Storage: -20°C

1181	9(E),11(E)-Octadecadienoic acid	25 mg	111.00
	9-trans, 11-trans CLA C ₁₈ H ₃₂ O ₂ CAS#: 544-71-8		

Source: synthetic Mol. Wt.: 280 Melting Point (°C): 55-57 Purity: 98% by TLC,
GC Appearance: solid Solubility: chloroform, ethanol, hexane, methanol
Storage: -20°C

1257	Methyl 9(E),11(E)-octadecadienoate Methyl ester of CLA (9- <i>trans</i> , 11- <i>trans</i>) C ₁₉ H ₃₄ O ₂	25 mg	111.00
	Source: synthetic Mol. Wt.: 294 Purity: 98+% by TLC, GC Appearance: liquid Solubility: chloroform, ethanol, hexane, methanol Storage: -20°C		
1248	9(Z),11(Z)-Octadecadienoic acid	25 mg	111.00
1248-1	9-cis, 11-cis CLA C₁₈H₃₂O₂ CAS#: 544-40-7	1 g	2,024.00
	Source: synthetic Mol. Wt.: 280 Melting Point (°C): 40-42 Purity: 96+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol, ethyl ether Storage: -20°C		
1256	Methyl 9(Z), 11(Z)-octadecadienoate Methyl ester of CLA (9- <i>cis</i> , 11- <i>cis</i>) C ₁₉ H ₃₄ O ₂	25 mg	111.00
	Source: synthetic Mol. Wt.: 294 Purity: 96+% by TLC, GC Appearance: liquid Solubility: chloroform, ethanol, hexane, methanol Storage: -20°C		
1249	10(E),12(Z)-Octadecadienoic acid	25 mg	111.00
1249-1	10-trans, 12-cis CLA C₁₈H₃₂O₂ CAS#: 2420-44-2	1 g	319.00
1249-10	Source: synthetic Mol. Wt.: 280 Purity: 98+% by TLC, GC Appearance: liquid Solubility: chloroform, ethanol, hexane, methanol Storage: -20°C	10 g	1,591.00
1254	Methyl 10(E), 12(Z)-octadecadienoate Methyl ester of CLA (10- <i>trans</i> , 12- <i>cis</i>) C ₁₉ H ₃₄ O ₂	25 mg	111.00
	Source: synthetic Mol. Wt.: 294 Purity: 98+% by TLC, GC Appearance: liquid Solubility: chloroform, ethanol, hexane, methanol Storage: -20°C		
1259	11(Z), 13(E)-Octadecadienoic acid 11- <i>cis</i> , 13- <i>trans</i> CLA C ₁₈ H ₃₂ O ₂	25 mg	111.00
	Source: synthetic Mol. Wt.: 280 Purity: 77% cis, trans; 2 % cis, cis; 6% trans, trans by TLC, GC Appearance: liquid Solubility: chloroform, ethanol, hexane, methanol Storage: -20°C		
1247-1	9(Z),11(E)-Octadecadienoic acid	1 g	174.00
1247-10	9-cis, 11-trans CLA C₁₈H₃₂O₂ CAS#: 2540-56-9	10 g	1,100.00
	Source: synthetic Mol. Wt.: 280 Purity: 74% 9(Z),11(E); 17%(Z),(Z); 1% (E),(E) by TLC, GC Appearance: liquid Solubility: ethanol, ethyl ether, hexane Storage: -20°C		
1258	Methyl 9(Z),11(E)-octadecadienoate Methyl ester of CLA (9- <i>cis</i> , 11- <i>trans</i>) C ₁₉ H ₃₄ O ₂	25 mg	85.00
	Source: synthetic Mol. Wt.: 294 Purity: 74% 9(Z),11(E); 17%(Z),(Z); 1% (E),(E) by TLC, GC Appearance: liquid Solubility: chloroform, ethyl ether, hexane Storage: -20°C		

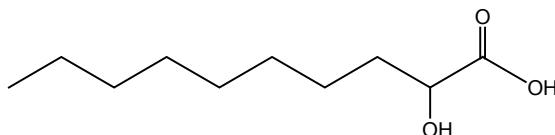
Other CLA Products and Derivatives

1409	1-Stearoyl-2-linoleoyl-sn-glycero-3-phosphorylcholine C ₄₄ H ₈₄ NO ₈ P	25 mg/ml, 1ml	204.00
Source: synthetic Mol. Wt.: 786 Purity: 98+% by TLC Appearance: liquid Solvent: chloroform Solubility: chloroform, ethanol Storage: -20°C			
1410	1-Stearoyl-2-[9(Z),11(E)-octadecadienoyl]-sn-glycero-3-phosphorylcholine C ₄₄ H ₈₄ NO ₈ P	25 mg/ml, 1ml	204.00
Source: synthetic Mol. Wt.: 786 Purity: 98+% by TLC Appearance: liquid Solvent: chloroform Solubility: chloroform, ethanol Storage: -20°C			
1411	1-Stearoyl-2-[10(E),12(Z)-octadecadienoyl]-sn-glycero-3-phosphorylcholine C ₄₄ H ₈₄ NO ₈ P	25 mg/ml, 1ml	204.00
Source: synthetic Mol. Wt.: 786 Purity: 98+% by TLC Appearance: liquid Solvent: chloroform Solubility: chloroform, ethanol Storage: -20°C			

Hydroxy Fatty Acids

2-Hydroxy Fatty Acids and Methyl Esters

These products are racemic and 98+% pure by GC and TLC. The 2-hydroxy fatty acids are components of glycosphingolipids and are involved in fatty acid degradation. They are stable and are supplied neat in vials.



Catalog number 1758

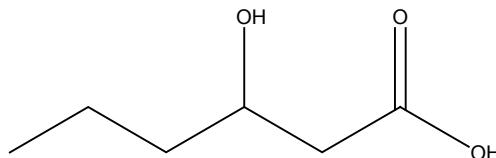
1758	2-Hydroxydecanoic acid	50 mg	114.00
1758-1	2-Hydroxy C10:0 fatty acid C ₁₀ H ₂₀ O ₃ CAS#: 5393-81-7	1 g	655.00
Source: synthetic Mol. Wt.: 188 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, methanol Storage: -20°C			
1759	Methyl 2-hydroxydecanoate	50 mg	114.00
1759-1	2-Hydroxy C10:0 methyl ester C ₁₁ H ₂₂ O ₃ CAS#: 71271-24-4	1 g	655.00
Source: synthetic Mol. Wt.: 202 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, methanol Storage: -20°C			

1701	2-Hydroxydodecanoic acid		50 mg	114.00
1701-1	2-Hydroxy C12:0 fatty acid C ₁₂ H ₂₄ O ₃ CAS#: 2984-55-6		1 g	655.00
Source: synthetic Mol. Wt.: 216 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, methanol Storage: -20°C				
1702	Methyl 2-hydroxydodecanoate		50 mg	114.00
1702-1	2-Hydroxy C12:0 methyl ester C ₁₃ H ₂₆ O ₃ CAS#: 51067-85-7		1 g	655.00
Source: synthetic Mol. Wt.: 230 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether, methanol Storage: -20°C				
1703	2-Hydroxytetradecanoic acid		50 mg	114.00
1703-1	2-Hydroxy C14:0 fatty acid C ₁₄ H ₂₈ O ₃ CAS#: 2507-55-3		1 g	655.00
Source: synthetic Mol. Wt.: 244 Melting Point (°C): 81-82 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, methanol Storage: -20°C				
1704	Methyl 2-hydroxytetradecanoate		50 mg	114.00
1704-1	2-Hydroxy C14:0 methyl ester C ₁₅ H ₃₀ O ₃ CAS#: 56009-40-6		1 g	655.00
Source: synthetic Mol. Wt.: 258 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether, methanol Storage: -20°C				
1705	2-Hydroxyhexadecanoic acid		50 mg	114.00
1705-1	2-Hydroxy C16:0 fatty acid C ₁₆ H ₃₂ O ₃ CAS#: 764-67-0		1 g	655.00
Source: synthetic Mol. Wt.: 272 Melting Point (°C): 86-87 Purity: 98+% by TLC, GC Appearance: solid Solubility: methanol, chloroform/methanol, 2:1 Storage: -20°C				
1706	Methyl 2-hydroxyhexadecanoate		50 mg	114.00
1706-1	2-Hydroxy C16:0 methyl ester C ₁₇ H ₃₄ O ₃ CAS#: 16742-51-1		1 g	655.00
Source: synthetic Mol. Wt.: 286 Melting Point (°C): 59-60 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether, methanol Storage: -20°C				
1707	2-Hydroxyoctadecanoic acid		50 mg	114.00
1707-1	2-Hydroxy C18:0 fatty acid C ₁₈ H ₃₆ O ₃ CAS#: 629-22-1		1 g	655.00
Source: synthetic Mol. Wt.: 300 Melting Point (°C): 92-93 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform/methanol, 5:1 Storage: -20°C				
1708	Methyl 2-hydroxyoctadecanoate		50 mg	114.00
1708-1	2-Hydroxy C18:0 methyl ester C ₁₉ H ₃₈ O ₃ CAS#: 2420-35-1		1 g	655.00
Source: synthetic Mol. Wt.: 315 Melting Point (°C): 64-66 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether, methanol Storage: -20°C				
1709	2-Hydroxyeicosanoic acid		25 mg	114.00
1709-0.5	2-Hydroxy C20:0 fatty acid C ₂₀ H ₄₀ O ₃ CAS#: 16742-48-6		0.5 g	655.00
Source: synthetic Mol. Wt.: 329 Melting Point (°C): 91-92 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform/methanol, 5:1 Storage: -20°C				

1710	Methyl 2-hydroxyeicosanoate	25 mg	114.00
1710-0.5	2-Hydroxy C20:0 methyl ester C ₂₁ H ₄₂ O ₃ CAS#: 16742-49-7	0.5 g	655.00
Source: synthetic Mol. Wt.: 343 Melting Point (°C): 62-64 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether Storage: -20°C			
1711	2-Hydroxydocosanoic acid	25 mg	114.00
1711-0.5	2-Hydroxy C22:0 fatty acid C ₂₂ H ₄₄ O ₃ CAS#: 13980-14-8	0.5 g	655.00
Source: synthetic Mol. Wt.: 366 Melting Point (°C): 96-97 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform/methanol, 5:1 Storage: -20°C			
1712	Methyl 2-hydroxydocosanoate	25 mg	114.00
1712-0.5	2-Hydroxy C22:0 methyl ester C ₂₃ H ₄₆ O ₃ CAS#: 13980-17-1	0.5 g	655.00
Source: synthetic Mol. Wt.: 371 Melting Point (°C): 72-73 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether Storage: -20°C			
1713	2-Hydroxytricosanoic acid	10 mg	162.00
	2-Hydroxy C23:0 fatty acid C ₂₃ H ₄₆ O ₃ CAS#: 2718-37-8		
Source: synthetic Mol. Wt.: 371 Melting Point (°C): 98-99 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform/methanol, 5:1 Storage: -20°C			
1714	Methyl 2-hydroxytricosanoate	10 mg	162.00
	2-Hydroxy C23:0 methyl ester C ₂₄ H ₄₈ O ₃ CAS#: 118745-41-8		
Source: synthetic Mol. Wt.: 385 Melting Point (°C): 68-70 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether Storage: -20°C			
1715	2-Hydroxytetacosanoic acid	5 mg	162.00
	2-Hydroxy C24:0 fatty acid; Cerebronic acid C ₂₄ H ₄₈ O ₃ CAS#: 544-57-0		
Source: synthetic Mol. Wt.: 385 Melting Point (°C): 101-104 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform/methanol, 5:1 Storage: -20°C			
1716	Methyl 2-hydroxytetacosanoate	5 mg	162.00
	2-Hydroxy C24:0 methyl ester C ₂₅ H ₅₀ O ₃ CAS#: 2433-95-6		
Source: synthetic Mol. Wt.: 399 Melting Point (°C): 77-80 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether Storage: -20°C			
1722	2-Hydroxy Methyl Ester Mixture	10 mg/ml, 1 ml	114.00
	Quantitative mixture		
Source: synthetic Appearance: liquid Solvent: chloroform Solubility: chloroform Storage: -20°C			
Contains: 2-OH C14:0, 20.0%; 2-OH C16:0, 20.0%; 2-OH C18:0, 15.0%; 2-OH C20:0, 15.0%; 2-OH C22:0, 10.0%; 2-OH C23:0, 10.0%; 2-OH C24:0, 10.0%			

3-Hydroxy Fatty Acids and Methyl Esters

These products are racemic and 98+% pure by GC and TLC. 3-Hydroxy fatty acids occur in the lipid fraction of many microorganisms and are useful in the typing of microbial isolates. They are stable and are supplied neat in vials.



Catalog number 1747

1747	3-Hydroxyhexanoic acid	25 mg	138.00
1747-0.5	3-Hydroxy C6:0 fatty acid C ₆ H ₁₂ O ₃ CAS#: 10191-24-9	0.5 g	689.00

Source: synthetic **Mol. Wt.:** 132 **Purity:** 98+% by TLC, GC **Appearance:** liquid
Solubility: chloroform, ethanol, methanol **Storage:** -20°C

1748	Methyl 3-hydroxyhexanoate	25 mg	138.00
1748-0.5	3-Hydroxy C6:0 methyl ester C ₇ H ₁₄ O ₃ CAS#: 21188-58-9	0.5 g	689.00

Source: synthetic **Mol. Wt.:** 146 **Purity:** 98+% by TLC, GC **Appearance:** liquid
Solubility: chloroform, ethanol, methanol **Storage:** -20°C

1745	3-Hydroxyoctanoic acid	25 mg	126.00
1745-0.5	3-Hydroxy C8:0 fatty acid C ₈ H ₁₆ O ₃ CAS#: 88930-08-9	0.5 g	689.00

Source: synthetic **Mol. Wt.:** 160 **Purity:** 98+% by TLC, GC **Appearance:** liquid
Solubility: chloroform, ethanol, methanol **Storage:** -20°C

1746	Methyl 3-hydroxyoctanoate	25 mg	126.00
1746-0.5	3-Hydroxy C8:0 methyl ester C ₉ H ₁₈ O ₃ CAS#: 85549-54-8	0.5 g	689.00

Source: synthetic **Mol. Wt.:** 174 **Purity:** 98+% by TLC, GC **Appearance:** liquid
Solubility: chloroform, ethanol, ethyl ether **Storage:** -20°C

1725	3-Hydroxynonanoic acid	25 mg	138.00
1725-0.5	3-Hydroxy C9:0 fatty acid C ₉ H ₁₈ O ₃ CAS#: 88930-09-0	0.5 g	689.00

Source: synthetic **Mol. Wt.:** 174 **Melting Point (°C):** 60-62 **Purity:** 98+% by TLC,
GC Appearance: solid **Solubility:** chloroform, ethanol, methanol
Storage: -20°C

1726	Methyl 3-hydroxynonanoate	25 mg	138.00
1726-0.5	3-Hydroxy C9:0 methyl ester C ₁₀ H ₂₀ O ₃ CAS#: 83968-06-3	0.5 g	689.00

Source: synthetic **Mol. Wt.:** 188 **Purity:** 98+% by TLC, GC **Appearance:** liquid
Solubility: chloroform, ethanol, ethyl ether **Storage:** -20°C

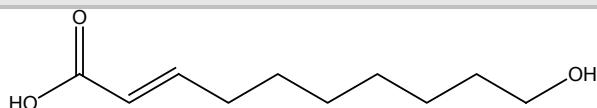
1727	3-Hydroxydecanoic acid	25 mg	162.00
1727-0.5	3-Hydroxy C10:0 fatty acid C ₁₀ H ₂₀ O ₃ CAS#: 5561-87-5	0.5 g	689.00

Source: synthetic **Mol. Wt.:** 188 **Melting Point (°C):** 57-60 **Purity:** 98+% by TLC,
GC Appearance: solid **Solubility:** chloroform, ethanol, methanol **Storage:** -20°C

1728	Methyl 3-hydroxydecanoate		25 mg	162.00
1728-0.5	3-Hydroxy C10:0 methyl ester C ₁₁ H ₂₂ O ₃ CAS#: 62675-82-5		0.5 g	689.00
	Source: synthetic Mol. Wt.: 202 Purity: 98+% by TLC, GC Appearance: liquid Solubility: chloroform, ethanol, methanol Storage: -20°C			
1729	3-Hydroxyundecanoic acid		25 mg	147.00
1729-0.5	3-Hydroxy C11:0 fatty acid C ₁₁ H ₂₂ O ₃ CAS#: 40165-88-6		0.5 g	689.00
	Source: synthetic Mol. Wt.: 202 Melting Point (°C): 74-76 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: -20°C			
1730	Methyl 3-hydroxyundecanoate		25 mg	147.00
1730-0.5	3-Hydroxy C11:0 methyl ester C ₁₂ H ₂₄ O ₃ CAS#: 127593-21-9		0.5 g	689.00
	Source: synthetic Mol. Wt.: 216 Purity: 98+% by TLC, GC Appearance: liquid Solubility: chloroform, ethanol, methanol Storage: -20°C			
1731	3-Hydroxydodecanoic acid		25 mg	138.00
1731-0.5	3-Hydroxy C12:0 fatty acid C ₁₂ H ₂₄ O ₃ CAS#: 1883-13-2		0.5 g	689.00
	Source: synthetic Mol. Wt.: 216 Melting Point (°C): 71-72 Purity: 98+% by TLC, GC Appearance: solid Solubility: ethanol, methanol Storage: -20°C			
1732	Methyl 3-hydroxydodecanoate		25 mg	147.00
1732-0.5	3-Hydroxy C12:0 methyl ester C ₁₃ H ₂₆ O ₃ CAS#: 85464-97-7		0.5 g	689.00
	Source: synthetic Mol. Wt.: 230 Purity: 98+% by TLC, GC Appearance: liquid Solubility: chloroform, ethanol, ethyl ether Storage: -20°C			
1733	3-Hydroxytridecanoic acid		25 mg	147.00
1733-0.5	3-Hydroxy C13:0 fatty acid C ₁₃ H ₂₆ O ₃ CAS#: 32602-69-0		0.5 g	689.00
	Source: synthetic Mol. Wt.: 230 Melting Point (°C): 80-83 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: -20°C			
1734	Methyl 3-hydroxytridecanoate		25 mg	147.00
1734-0.5	3-Hydroxy C13:0 methyl ester C ₁₄ H ₂₈ O ₃		0.5 g	689.00
	Source: synthetic Mol. Wt.: 244 Purity: 98+% by TLC, GC Appearance: liquid Solubility: chloroform, ethyl ether Storage: -20°C			
1735	3-Hydroxytetradecanoic acid		25 mg	138.00
1735-0.5	3-Hydroxy C14:0 fatty acid C ₁₄ H ₂₈ O ₃ CAS#: 3422-31-9		0.5 g	689.00
	Source: synthetic Mol. Wt.: 244 Melting Point (°C): 80-81 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: -20°C			
1736	Methyl 3-hydroxytetradecanoate		25 mg	138.00
1736-0.5	3-Hydroxy C14:0 methyl ester C ₁₅ H ₃₀ O ₃ CAS#: 55682-83-2		0.5 g	689.00
	Source: synthetic Mol. Wt.: 258 Melting Point (°C): 36-37 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethyl ether, methanol Storage: -20°C			

1739	3-Hydroxyhexadecanoic acid		25 mg	138.00
1739-0.5	3-Hydroxy C16:0 fatty acid C ₁₆ H ₃₂ O ₃ . CAS#: 928-17-6		0.5 g	689.00
Source: synthetic Mol. Wt.: 272 Melting Point (°C): 85-86 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: -20°C				
1740	Methyl 3-hydroxyhexadecanoate		25 mg	138.00
1740-0.5	3-Hydroxy C16:0 methyl ester C ₁₇ H ₃₄ O ₃ CAS#: 51883-36-4		0.5 g	689.00
Source: synthetic Mol. Wt.: 286 Melting Point (°C): 43-45 Purity: 98+% by TLC, GC Appearance: solid Solubility: ethanol, methanol Storage: -20°C				
1741	3-Hydroxyheptadecanoic acid		25 mg	126.00
1741-0.5	3-Hydroxy C17:0 fatty acid C ₁₇ H ₃₄ O ₃ CAS#: 40165-89-7		0.5 g	689.00
Source: synthetic Mol. Wt.: 286 Melting Point (°C): 93-95 Purity: 98+% by TLC, GC Appearance: solid Solubility: ethanol, methanol Storage: -20°C				
1742	Methyl 3-hydroxyheptadecanoate		25 mg	126.00
1742-0.5	3-Hydroxy C17:0 methyl ester C ₁₈ H ₃₆ O ₃		0.5 g	689.00
Source: synthetic Mol. Wt.: 300 Melting Point (°C): 53-55 Purity: 98+% by TLC, GC Appearance: solid Solubility: ethanol, methanol Storage: -20°C				
1743	3-Hydroxyoctadecanoic acid		25 mg	126.00
1743-0.5	3-Hydroxy C18:0 fatty acid C ₁₈ H ₃₆ O ₃ CAS#: 45261-96-9		0.5 g	689.00
Source: synthetic Mol. Wt.: 300 Melting Point (°C): 52-54 Purity: 98+% by TLC, GC Appearance: solid Solubility: ethanol, methanol Storage: -20°C				
1744	Methyl 3-hydroyoctadecanoate		25 mg	126.00
1744-0.5	3-Hydroxy C18:0 methyl ester C ₁₉ H ₃₈ O ₃ CAS#: 14531-40-9		0.5 g	689.00
Source: synthetic Mol. Wt.: 314 Melting Point (°C): 52-54 Purity: 98+% by TLC, GC Appearance: solid Solubility: ethanol, methanol Storage: -20°C				

Omega Hydroxy Fatty Acids



Catalog number 1754

1754	Royal Jelly acid		50 mg	184.00
	10-Hydroxy-2-(E)-decenoic acid; <i>omega</i> -Hydroxy C10:1 (2- <i>trans</i>) fatty acid C ₁₀ H ₁₈ O ₃ CAS#: 14113-05-4			
Source: synthetic Mol. Wt.: 186 Melting Point (°C): 63-65 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: -20°C				

1881	15-Hydroxypentadecanoic acid <i>omega</i> -Hydroxy C15:0 fatty acid C ₁₅ H ₃₀ O ₃ CAS#: 4617-33-8	25 mg	75.00
	Source: synthetic Mol. Wt.: 258 Melting Point (°C): 84-86 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: room temperature		
1882	Methyl 15-hydroxypentadecanoate <i>omega</i> -Hydroxy C15:0 methyl ester C ₁₆ H ₃₂ O ₃ CAS#: 76529-42-5	25 mg	75.00
	Source: synthetic Mol. Wt.: 272 Melting Point (°C): 50-52 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, warm ethanol, ethyl ether Storage: room temperature		
1760	17-Hydroxyheptadecanoic acid <i>omega</i> -Hydroxy C17:0 fatty acid C ₁₇ H ₃₄ O ₃ CAS#: 13099-34-8	25 mg	184.00
	Source: synthetic Mol. Wt.: 286 Melting Point (°C): 93-95 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, warm ethanol, ethyl ether Storage: room temperature		
1761	Methyl 17-hydroxyheptadecanoate <i>omega</i> -Hydroxy C17:0 methyl ester C ₁₈ H ₃₆ O ₃ CAS#: 94036-00-7	25 mg	184.00
	Source: synthetic Mol. Wt.: 300 Melting Point (°C): 59-63 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, warm ethanol, ethyl ether Storage: room temperature		
1877	20-Hydroxyeicosanoic acid <i>omega</i> -Hydroxy C20:0 fatty acid C ₂₀ H ₄₀ O ₃	25 mg	162.00
	Source: synthetic Mol. Wt.: 328 Melting Point (°C): 96-98 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, warm ethanol Storage: room temperature		
1878	Methyl 20-hydroxyeicosanoate <i>omega</i> -Hydroxy C20:0 methyl ester C ₂₁ H ₄₂ O ₃	25 mg	162.00
	Source: synthetic Mol. Wt.: 342 Melting Point (°C): 69-71 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, warm ethanol, ethyl ether Storage: room temperature		
1880	Methyl 21-hydroxyheneicosanoate <i>omega</i> -Hydroxy C21:0 methyl ester C ₂₂ H ₄₄ O ₃	25 mg	184.00
	Source: synthetic Mol. Wt.: 356 Melting Point (°C): 73-76 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, warm ethanol, ethyl ether Storage: room temperature		
1818	22-Hydroxydocosanoic acid <i>omega</i> -Hydroxy C22:0 fatty acid C ₂₂ H ₄₄ O ₃	25 mg	194.00
	Source: synthetic Mol. Wt.: 356 Melting Point (°C): 100-102 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, warm ethanol Storage: room temperature		

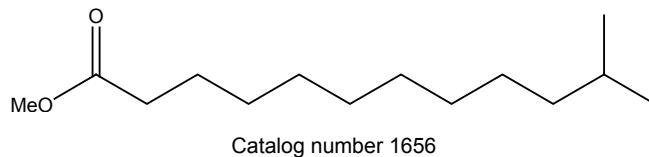
1819	Methyl 22-hydroxydocosanoate <i>omega</i> -Hydroxy C22:0 methyl ester C ₂₃ H ₄₆ O ₃	25 mg	194.00
Source: synthetic Mol. Wt.: 370 Melting Point (°C): 73-75 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, warm ethanol, ethyl ether Storage: room temperature			
1883	Methyl 27-hydroxyheptacosanoate <i>omega</i> -Hydroxy C27:0 methyl ester C ₂₈ H ₅₆ O ₃	25 mg	194.00
Source: synthetic Mol. Wt.: 440 Melting Point (°C): 85-89 Purity: 97+% by TLC, GC Appearance: solid Solubility: chloroform Storage: room temperature			
1884	Methyl 30-hydroxytriacontanoate <i>omega</i> -Hydroxy C30:0 methyl ester C ₃₁ H ₆₂ O ₃	25 mg	194.00
Source: synthetic Mol. Wt.: 482 Melting Point (°C): 88-91 Purity: 97+% by TLC, GC Appearance: solid Solubility: chloroform Storage: room temperature			

Other Hydroxy Fatty Acids

1182	Ricinelaidic acid 12-Hydroxy C18:1 (9- <i>trans</i>) fatty acid C ₁₈ H ₃₄ O ₃ CAS#: 82188-83-8	100 mg	60.00
Source: synthetic Mol. Wt.: 298 Melting Point (°C): 50-53 Purity: 98+% by TLC, GC Appearance: solid Solubility: ethanol, methanol Storage: -20°C			
1183	Methyl ricinelaideate 12-Hydroxy C18:1 (9- <i>trans</i>) methyl ester C ₁₉ H ₃₆ O ₃ CAS#: 7706-01-6	100 mg	60.00
Source: synthetic Mol. Wt.: 312 Purity: 98+% by TLC, GC Appearance: solid Solubility: ethanol, methanol Storage: -20°C			
1766	6-Hydroxyoctadecanoic acid 6-Hydroxy C18:0 fatty acid C ₁₈ H ₃₆ O ₃	10 mg	167.00
Source: synthetic Mol. Wt.: 300 Melting Point (°C): 80-82 Purity: 98+% by TLC, GC Appearance: solid Solubility: ethanol, methanol Storage: room temperature			

Branched and Cyclic Fatty Acids

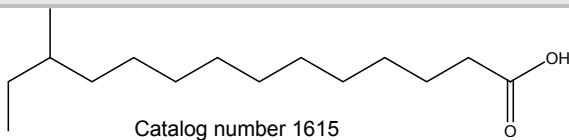
Iso-Fatty Acids and Esters



1656	Methyl 11-methyldodecanoate iso-Tridecanoic methyl ester; iso C13 Methyl ester C ₁₄ H ₂₈ O ₂ CAS#: 5129-57-7	20 mg	290.00
Source: synthetic Mol. Wt.: 228 Purity: 98+% by GC Appearance: liquid Solubility: hexane, ethyl ether, methylene chloride Storage: -20°C			

1657	Methyl 12-methyltridecanoate iso-Tetradecanoic methyl ester; iso C14 Methyl ester C ₁₅ H ₃₀ O ₂ CAS#: 5129-58-8	20 mg	290.00
	Source: synthetic Mol. Wt.: 242 Purity: 98+% by GC Appearance: liquid Solubility: chloroform, ethyl ether, ethanol Storage: -20°C		
1605	13-Methyltetradecanoic acid iso-Pentadecanoic acid; iso C15 Fatty acid C ₁₅ H ₃₀ O ₂ CAS#: 27836-87-9	20 mg	290.00
	Source: synthetic Mol. Wt.: 242 Purity: 98+% by GC Appearance: solid Solubility: chloroform, ethyl ether, ethanol Storage: -20°C		
1600	Methyl 13-methyltetradecanoate iso-Pentadecanoic methyl ester; iso C15 Methyl ester C ₁₆ H ₃₂ O ₂ CAS#: 5129-59-9	20 mg	290.00
	Source: synthetic Mol. Wt.: 256 Purity: 98+% by GC Appearance: liquid Solubility: chloroform, ethyl ether, ethanol Storage: -20°C		
1601	Methyl 14-methylpentadecanoate iso-Palmitic methyl ester; iso C16 Methyl ester C ₁₇ H ₃₄ O ₂ CAS#: 5129-60-2	20 mg	290.00
	Source: synthetic Mol. Wt.: 270 Purity: 98+% by GC Appearance: liqiuud Solubility: chloroform, ethyl ether, ethanol Storage: -20°C		
1606	15-Methylhexadecanoic acid iso-Heptadecanoic acid; iso C17 Fatty acid C ₁₇ H ₃₄ O ₂ CAS#: 1603-03-8	20 mg	290.00
	Source: synthetic Mol. Wt.: 270 Purity: 98+% by GC Appearance: solid Solubility: chloroform, ethyl ether, ethanol Storage: -20°C		
1602	Methyl 15-methylhexadecanoate iso-Heptadecanoic methyl ester; iso C17 Methyl ester C ₁₈ H ₃₆ O ₂ CAS#: 6929-04-0	20 mg	290.00
	Source: synthetic Mol. Wt.: 284 Purity: 98+% by GC Appearance: liquid Solubility: chloroform, ethyl ether, ethanol Storage: -20°C		
1603	Methyl 17-methyloctadecanoate iso-Nonadecanoic methyl ester; iso C19 Methyl ester C ₂₀ H ₄₀ O ₂ CAS#: 55124-97-5	20 mg	290.00
	Source: synthetic Mol. Wt.: 313 Purity: 98+% by GC Appearance: solid Solubility: chloroform, ethyl ether, ethanol Storage: -20°C		

Anteiso-Fatty Acids and Esters



1615	12-Methyltetradecanoic acid anteiso-Pentadecanoic acid; anteiso C15 Fatty acid C ₁₅ H ₃₀ O ₂ CAS#: 5502-94-3	20 mg	290.00
	Source: synthetic Mol. Wt.: 242 Purity: 98+% by GC Appearance: solid Solubility: chloroform, ethyl ether, ethanol Storage: -20°C		
1612	Methyl 12-methyltetradecanoate anteiso-Pentadecanoic methyl ester; anteiso C15 Methyl ester C ₁₆ H ₃₂ O ₂ CAS#: 5129-66-8	20 mg	290.00
	Source: synthetic Mol. Wt.: 256 Purity: 98+% by GC Appearance: liquid Solubility: chloroform, ethyl ether, ethanol Storage: -20°C		
1613	Methyl 13-methylpentadecanoate anteiso-Palmitic methyl ester; anteiso C16 Methyl ester C ₁₇ H ₃₄ O ₂ CAS#: 5487-50-3	20 mg	290.00
	Source: synthetic Mol. Wt.: 270 Purity: 98+% by GC Appearance: liquid Solubility: chloroform, ethyl ether, ethanol Storage: -20°C		
1616	14-Methylhexadecanoic acid anteiso-Heptadecanoic acid; anteiso C17 Fatty acid C ₁₇ H ₃₄ O ₂ CAS#: 5918-29-6	20 mg	290.00
	Source: synthetic Mol. Wt.: 270 Purity: 98+% by GC Appearance: solid Solubility: chloroform, ethyl ether, ethanol Storage: -20°C		
1614	Methyl 14-methylhexadecanoate anteiso-Heptadecanoic methyl ester; anteiso C17 Methyl ester C ₁₈ H ₃₆ O ₂ CAS#: 2490-49-5	20 mg	290.00
	Source: synthetic Mol. Wt.: 284 Purity: 98+% by GC Appearance: liquid Solubility: chloroform, ethyl ether, ethanol Storage: -20°C		

Methylated Fatty Acids

1207	D,L-2,6-Dimethylheptanoic acid 2,6-Dimethyl C7:0 fatty acid C ₉ H ₁₈ O ₂	50 mg	120.00
	Source: synthetic Mol. Wt.: 158 Purity: 98+% by TLC, GC Appearance: liquid Solubility: chloroform Storage: room temperature		

1791	10-Methylhexadecanoic acid 10-Methyl C16:0 fatty acid C ₁₇ H ₃₄ O ₂	25 mg	176.00
	Source: synthetic Mol. Wt.: 270 Purity: 98+% by TLC, GC Appearance: liquid Solubility: chloroform Storage: room temperature		
1792	Methyl 10-methylhexadecanoate 10-Methyl C16:0 methyl ester C ₁₈ H ₃₆ O ₂	25 mg	176.00
	Source: synthetic Mol. Wt.: 284 Purity: 98+% by TLC, GC Appearance: liquid Solubility: chloroform Storage: room temperature		
1195	Phytanic acid 3,7,11,15-Tetramethylhexadecanoic acid C ₂₀ H ₄₀ O ₂ CAS#: 14721-66-5	25 mg	251.00
	Source: semi-synthetic Mol. Wt.: 312 Purity: 97+% by GC Appearance: solid Solubility: chloroform, methanol Storage: -20°C		

Cyclopropyl Fatty Acids and Esters

1822	Methyleneoctadecanoic acid (all cis-9,10) Dihydrosterculic acid C ₁₉ H ₃₆ O ₂ CAS#: 4675-61-0	25 mg	135.00
	Source: synthetic Mol. Wt.: 296 Melting Point (°C): 38-42 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol, hexane Storage: -20°C		
1823	Methyl cis-9,10-methyleneoctadecanoate, C19:0 delta (all cis-9,10) Methyl dihydrosterculate C ₂₀ H ₃₈ O ₂ CAS#: 3971-54-8	25 mg	135.00
	Source: synthetic Mol. Wt.: 310 Purity: 98+% by TLC, GC Appearance: liquid Solubility: chloroform, ethanol, methanol, hexane Storage: -20°C		

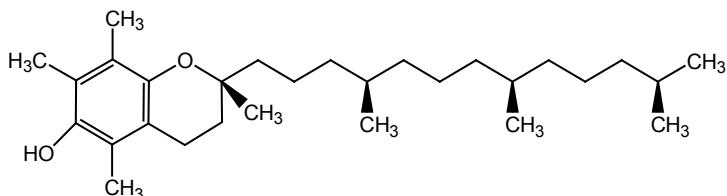
Unusual Fatty Acids and Derivatives

1751	N-Oleoylethanolamine NOE C ₂₀ H ₃₉ NO ₂ CAS#: 111-58-0	100 mg	163.00
	Source: synthetic Mol. Wt.: 326 Melting Point (°C): 63-66 Purity: 98+% by TLC, GC Appearance: solid Solubility: chloroform, ethanol, methanol, ethyl ether, DMSO Storage: -20°C		
	Activity: acid ceramidase inhibitor		
1786	N-Hexadecanoylethanolamine C ₁₈ H ₃₇ NO ₂ CAS# 544-31-0	100 mg	114.00
	Source: synthetic Mol. Wt.: 299 Melting Point (°C): 99-102 Purity: 98+% by TLC Appearance: solid Solubility: chloroform, ethanol, methanol, Storage: -20°C		
	Activity: inactive as acid ceramidase inhibitor		

Other Lipids

Tocopherols

Catalog number 1072



1072

rac-alpha-Tocopherol

5,7,8-Trimethyltocol C₂₉H₅₀O₂ CAS#: 59-02-9

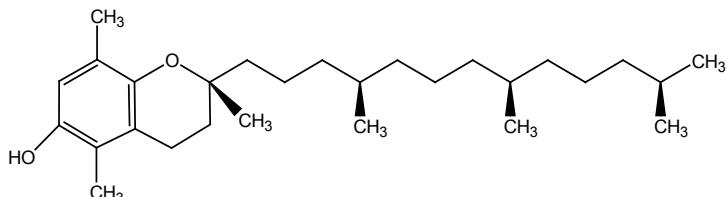
50 mg/ml, 1 ml

103.00

Source: synthetic Mol. Wt.: 431 Purity: 95% by TLC, 98% by GC

Appearance: liquid Solvent: hexane Solubility: chloroform, ethanol, hexane, methanol Storage: -20°C

Catalog number 1071



1071

rac-beta-Tocopherol

5,8-Dimethyltocol C₂₈H₄₈O₂ CAS#: 148-03-8

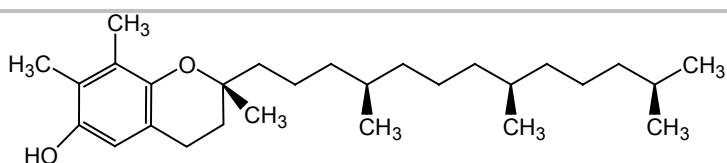
50 mg/ml, 1 ml

154.00

Source: synthetic Mol. Wt.: 417 Purity: 95% by TLC, 98% by GC

Appearance: liquid Solvent: hexane Solubility: chloroform, ethanol, hexane, methanol Storage: -20°C

Catalog number 1073



1073

rac-gamma-Tocopherol

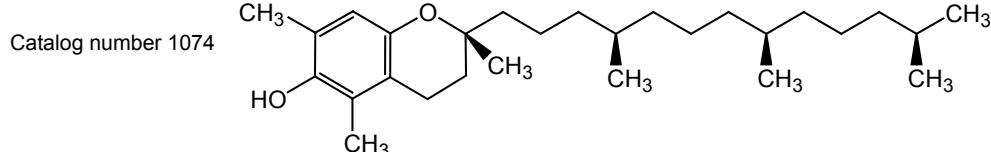
7,8-Dimethyltocol C₂₈H₄₈O₂ CAS#: 73980-80-0

50 mg/ml, 1 ml

154.00

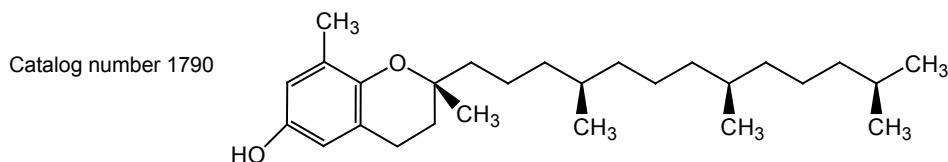
Source: synthetic Mol. Wt.: 417 Purity: 95% by TLC, 97% by GC

Appearance: liquid Solvent: hexane Solubility: chloroform, ethanol, hexane, methanol Storage: -20°C



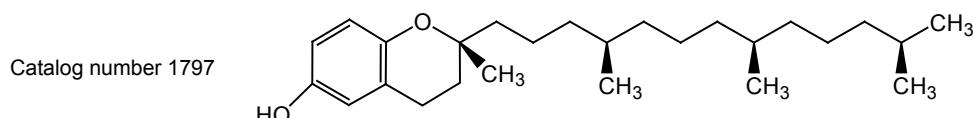
1074 *rac*-5,7-Dimethyltocol **50 mg/ml, 1 ml** **154.00**
 $C_{28}H_{48}O_2$ CAS#: 493-35-6

Source: synthetic **Mol. Wt.:** 417 **Purity:** 95% by TLC, 98% by GC
Appearance: liquid **Solvent:** hexane **Solubility:** hexane, ethyl ether, chloroform, alcohols **Storage:** -20°C



1790 (+)-*delta*-Tocopherol **50 mg/ml, 1 ml** **138.00**
 8-Methyltocol $C_{27}H_{46}O_2$ CAS#: 119-13-1

Source: natural, plant **Mol. Wt.:** 403 **Purity:** 95% by TLC, 98% by GC
Appearance: liquid **Solvent:** hexane **Solubility:** chloroform, ethanol, hexane, methanol **Storage:** -20°C

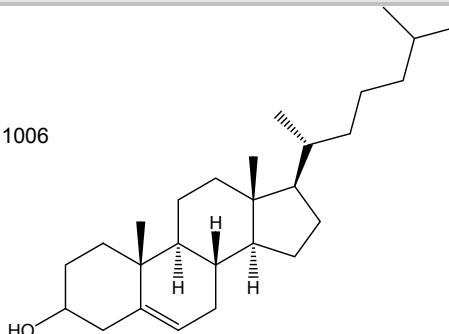


1797 Tocol **50 mg/ml, 1 ml** **145.00**
 $rac\text{-Tocol}$ $C_{26}H_{44}O_2$

Source: synthetic **Mol. Wt.:** 389 **Purity:** 95% by TLC, 98% by GC
Appearance: liquid **Solvent:** hexane **Solubility:** hexane, methanol, ethanol **Storage:** -20°C

Cholestane Derivatives

Catalog number 1006



1006	Cholesterol C ₂₇ H ₄₆ O CAS#: 57-88-5	500 mg	53.00
-------------	---	---------------	--------------

Source: natural, ovine **Mol. Wt.:** 387 **Melting Point (°C):** 147-148 **Purity:** 98+% by TLC, GC **Appearance:** solid **Solubility:** chloroform, ethanol **Storage:** -20°C

1115	5-alpha-Cholestane C ₂₇ H ₄₈ CAS#: 481-21-0	100 mg	85.00
-------------	---	---------------	--------------

Source: synthetic **Mol. Wt.:** 373 **Purity:** 98+% by GC **Appearance:** solid
Solubility: chloroform, ethyl ether, hexane **Storage:** -20°C

1116	Coprostanol 5-beta-Cholestan-3-beta-ol C ₂₇ H ₄₈ O CAS#: 360-68-9	25 mg	154.00
-------------	---	--------------	---------------

Source: semisynthetic **Mol. Wt.:** 389 **Melting Point (°C):** 101-103 **Purity:** 98+% by GC **Appearance:** solid **Solubility:** chloroform, ethyl ether, warm methanol **Storage:** -20°C

Plant Sterols and Steryl Glucosides

1119	Plant Sterol Mixture Sterol mixture, qualitative	25 mg/ml, 1 ml	93.00
-------------	--	-----------------------	--------------

Source: natural, plant **Appearance:** liquid **Solvent:** chloroform
Solubility: chloroform **Storage:** -20°C

Contains: Brassicasterol, Campesterol, Stigmasterol, *beta*-Sitosterol, in order of elution

1123	Plant Sterols Kit	1 kit	387.00
-------------	--------------------------	--------------	---------------

Source: synthetic or plant **Appearance:** liquid/solid **Solvent:** chloroform
Solubility: chloroform **Storage:** -20°C

Contains in individual packages: sterol glucosides 25 mg, esterified sterol glucosides 10 mg, plant sterol mixture 25 mg, β -sitosterol (55%) 100 mg, desmosterol (85%) 2 mg, lanosterol (55%) 100 mg, stigmasterol 25 mg, ergosterol 25 mg, coprostanol 5 mg, cholestanol 100 mg

1113	beta-Sitostanol Stigmastanol C ₂₉ H ₅₂ O CAS#: 19466-47-8	50 mg	73.00
Source: synthetic Mol. Wt.: 417 Melting Point (°C): 127-132 Purity: 98+% by TLC, 97+% by GC Appearance: solid Solubility: chloroform Storage: -20°C			
1120	Lanosterol C ₃₀ H ₅₀ O CAS#: 79-63-0	500 mg	73.00
Source: synthetic or plant Mol. Wt.: 427 Purity: 55% by TLC, GC Appearance: solid Solubility: chloroform Storage: -20°C			
1121	Stigmasterol 5,22-Cholestadien-24-beta-ethyl-3-beta-ol C ₂₉ H ₄₈ O CAS#: 83-48-7	100 mg	73.00
Source: synthetic Mol. Wt.: 413 Melting Point (°C): 165-167 Purity: 95% by TLC, GC Appearance: solid Solubility: chloroform Storage: -20°C			
1122	Ergosterol C ₂₈ H ₄₄ O CAS#: 57-87-4	100 mg	54.00
Source: synthetic or plant Mol. Wt.: 397 Melting Point (°C): 156-158 Purity: 95% by TLC, GC Appearance: solid Solubility: chloroform Storage: -20°C			
1117	Steryl glucosides C ₃₅ H ₆₀ O ₆	25 mg	88.00
Source: natural, plant Mol. Wt.: 577 (based on β-sitosteryl glucoside) Melting Point (°C): 283-287 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1 (warm) Storage: -20°C			
1118	Esterified Steryl Glucosides 1:1:1, sterol:glucose:fatty acid C ₅₁ H ₉₀ O ₇	10 mg	99.00
Source: natural, plant Mol. Wt.: 815 (based on β-sitosteryl glucoside palmitate) Purity: 98+% by TLC Appearance: solid Solubility: chloroform, ethyl ether, pyridine Storage: -20°C			
Sterol, glucose and fatty acid in a molar ratio 1:1:1.			

Standards and Reference Compounds

Food Industry Mixtures

Each methyl ester mixture is carefully prepared by weight.

4210	KEL-FIM-FAME-5 Mixture Methyl ester mixture	15.5 mg/ml, 1 ml	85.00
Source: synthetic or plant Appearance: liquid Solvent: heptane Solubility: heptane Storage: -20°C			
Contains the methyl esters of the following fatty acids (mg/ml in brackets): C8:0 [0.3], C10:0 [0.5], C12:0 [1.0], C13:0 [0.5], C14:0 [0.5], C14:1 [0.3], C15:0 [0.3], C16:0 [2.0], C16:1 [1.0], C17:0 [0.5], C18:0 [1.0], C18:1tr [0.4], C18:1c [3.0], C18:2 [2.0], C20:0 [0.3], C18:3 [1.0], C20:1 [0.3], C22:0 [0.3], C22:1 [0.3], listed in order of their elution.			
2009	FIM-FAME-6 Mixture Methyl ester mixture	33 mg/ml, 1 ml	114.00
Source: synthetic or plant Appearance: liquid Solvent: heptane Storage: -20°C			
Contains the methyl esters of these fatty acids. Each methyl ester is 3.03% of the mixture except C16:0 which is 6.06%. C4:0 , C6:0 , C8:0, C10:0 , C11:0, C12:0, C13:0 , C14:0, C14:1(<i>cis</i> -9), C15:0, C15:1(<i>cis</i> -10), C16:0, C16:1(<i>cis</i> -9), C17:0, C17:1(<i>cis</i> -10), C18:0, C18:1(<i>trans</i> -9), C18:1(<i>cis</i> -9), C18:2(all <i>cis</i> -9,12), C20:0, C18:3(all <i>cis</i> -6,9,12), C20:1(<i>cis</i> -11), C18:3(all <i>cis</i> -9,12,15), C20:2(all <i>cis</i> -11,14), C22:0, C20:3(all <i>cis</i> -8,11,14), C22:1(<i>cis</i> -13), C20:3(all <i>cis</i> -11,14,17), C20:4(all <i>cis</i> -5,8,11,14), C22:2(all <i>cis</i> -13,16), C24:1(<i>cis</i> -15), C22:6(all <i>cis</i> -4,7,10,13,16,19), listed in order of their elution.			
2010	FIM-FAME-7 Mixture Methyl ester mixture	30 mg/ml, 1 ml	114.00
Source: synthetic or plant Appearance: liquid Solvent: methylene chloride Solubility: methylene chloride Storage: -20°C			
Contains the methyl esters of these fatty acids (weight percent in [brackets]): C4:0 [4.0], C6:0 [4.0], C8:0 [4.0], C10:0 [4.0], C11:0 [2.0], C12:0 [4.0], C13:0 [2.0], C14:0 [4.0], C14:1(<i>cis</i> -9) [2.0], C15:0 [2.0], C15:1(<i>cis</i> -10) [2.0], C16:0 [6.0], C16:1(<i>cis</i> -9) [2.0], C17:0 [2.0], C17:1(<i>cis</i> -10) [2.0], C18:0 [4.0], C18:1(<i>trans</i> -9) [2.0], C18:1(<i>cis</i> -9) [4.0], C18:2(all <i>trans</i> -9,12) [2.0], C18:2(all <i>cis</i> -9,12) [2.0], C20:0 [4.0], C18:3(all <i>cis</i> -6,9,12) [2.0], C20:1(<i>cis</i> -11) [2.0], C18:3(all <i>cis</i> -9,12,15) [2.0], C21:0 [2.0], C20:2(all <i>cis</i> -11,14) [2.0], C22:0 [4.0], C20:3 (all <i>cis</i> -8,11,14) [2.0], C22:1(<i>cis</i> -13) [2.0], C20:3(all <i>cis</i> -11,14,17) [2.0], C20:4(all <i>cis</i> -5,8,11,14) [2.0], C23:0 [2.0], C22:2(all <i>cis</i> -13,16) [2.0], C24:0 [4.0], C20:5(all <i>cis</i> -5,8,11,14,17) [2.0], C24:1(<i>cis</i> -15) [2.0], C22:6(all <i>cis</i> -4,7,10,13,16,19) [2.0], listed in order of their elution.			

Polyunsaturated Fatty Acid Methyl Esters Mixtures

These are complex qualitative standard mixtures of polyunsaturated fatty acid methyl esters. Because they are extracted from natural materials, relative peak sizes may vary from lot to lot.

1093	PUFA-1 Qualitative mixture	100 mg	111.00
	Source: natural, fish oil Appearance: liquid Solubility: chloroform, ethanol, hexane, methanol Storage: -20°C		
	Contains: C14:0, C16:0, C16:1ω7, C18:1ω9, C18:1ω7, C18:2ω6, C20:1ω9, C18:4ω3, C22:1ω11, C22:1ω9, C20:5ω3, C22:5ω3, C22:6ω3		
1081	PUFA-2 Qualitative mixture	100 mg	122.00
	Source: natural, porcine Appearance: liquid Solubility: alcohols, hexane, chloroform Storage: -20°C		
	Contains: C14:0, C16:0, C16:1ω7, C18:0, C18:1ω9, C18:1ω7, C18:2ω6, C18:3ω6, C18:3ω3, C20:1ω9, C20:2ω6, C20:3ω6, C20:4ω6, C20:5ω3, C22:4ω6, C22:5ω3, C22:6ω3		
1177	PUFA-3 Qualitative mixture	100 mg	122.00
	Source: natural, menhaden oil Appearance: liquid Solubility: alcohols, hexane, chloroform Storage: -20°C		
	Contains: C14:0, C16:0, C16:1ω7, C16:2ω4, C16:3ω4, C16:4ω1, C18:0, C18:1ω9, C18:1ω7, C18:2ω6, C18:2ω4, C18:3ω4, C18:3ω3, C18:4ω3, C20:1ω9, C20:4ω6, C20:4ω3, C20:5ω3, C21:5ω3, C22:5ω3, C22:6ω3		

Carbohydrate Mixtures

1124	Alditol Acetate Mixture-1 Quantitative carbohydrate mixture	50 mg/ml, 1 ml	139.00
	Source: synthetic Appearance: liquid Solvent: chloroform Solubility: chloroform Storage: -20°C		
	Contains: rhamnitol, fucitol, ribitol and arabinitol pentaacetates, 12.5 mg/ml each		
1125	Alditol Acetate Mixture-2 Quantitative carbohydrate mixture	50 mg/ml, 1 ml	139.00
	Source: synthetic Appearance: liquid Solvent: chloroform Solubility: chloroform Storage: -20°C		
	Contains: mannitol, galactitol, glucitol and inositol hexaacetates, 12.5 mg/ml each		

Other Fatty Acid Methyl Ester Mixtures

1722	2-Hydroxy Methyl Ester Mixture Source: synthetic Appearance: liquid Solvent: chloroform Solubility: chloroform Storage: -20°C Quantitative mix contains: C14:0, 20.0%; C16:0, 20.0%; C18:0, 15.0%; C20:0, 15.0%; C22:0, 10.0%; C23:0, 10.0%; C24:0, 10.0%	10 mg/ml, 1 ml	114.00
1131	Cis-Trans Isomer Standard Mixture Source: margarine Appearance: liquid Solvent: 5ml methylene chloride Solubility: methylene chloride, chloroform Storage: -20°C Analysis of positional <i>cis-trans</i> fatty acid isomers is ever more important in light of the new food industry rules. These isomers can be resolved on Supelco SP-2560 or an equivalent capillary GC column. Use this specially formulated mixture to ensure proper operation of your column for this tricky separation. Mixture consists of <i>cis-trans</i> fatty acid isomers as methyl esters in methylene chloride. This is a qualitative standard containing in order of elution: C16:0, C18:0, C18:1 <i>trans</i> isomers (4 peaks), C18:1 <i>cis</i> & <i>trans</i> isomers (2 peaks), C18:1 <i>cis</i> isomers (4 peaks), C18:2, C20:0, C20:1 and C18:3 (same peak), C22:0	5 mg/ml, 5 ml	135.00
2011	Long Chain Fatty Acid Methyl Ester Mixture C24:0, C26:0, C28:0, C30:0, C32:0 Fatty acid methyl ester mixture Source: synthetic Appearance: liquid Solvent: methylene chloride Solubility: methylene chloride Storage: -20°C Quantitative mixture contains: C24:0, 20.0%; C26:0, 20.0%; C28:0, 20.0%; C30:0, 20.0%; C32:0, 20.0%	25 mg/ml, 1 ml	99.00

AOCS Animal and Vegetable Oil Reference Mixtures (RM Mixtures)

By studying problems with the quantitative analysis of animal and vegetable oils and fats, the American Oil Chemists' Society has found certain mixtures to be useful as reference standards. The composition of each mixture (see Table I below) is similar to the fatty acid distribution of certain oils. All mixtures are in methyl ester form and ready for GC analysis

Table I. AOCS Oil Reference Mixtures

Each methyl ester mixture is carefully prepared by weight and the composition verified by gas chromatography. The weight percentage of each component is indicated in the Table.

Mix No. Catalog No.	RM-1 1084	RM-2 1085	RM-3 1086	Rapeseed 1083	RM-4 1087	RM-5 1088	RM-6 1089
C8:0 Caprylate						7.0	
C10:0 Caprate						5.0	
C12:0 Laurate						48.0	
C14:0 Myristate			1.0	1.0		15.0	2.0
C16:0 Palmitate	6.0	7.0	4.0	4.0	11.0	7.0	30.0
C16:1 Palmitoleate (<i>cis</i> -9)							3.0
C18:0 Stearate	3.0	5.0	3.0	3.0	3.0	3.0	14.0
C18:1 Oleate (<i>cis</i> -9)	35.0	18.0	45.0	60.0	80.0	12.0	41.0
C18:2 Linoleate (all <i>cis</i> -9,12)	50.0	36.0	15.0	12.0	6.0	3.0	7.0
C18:3 Linolenate (all <i>cis</i> -9,12,15)	3.0	34.0	3.0	5.0			3.0
C20:0 Arachidate	3.0		3.0	3.0			
C20:1 Eicosenoate (<i>cis</i> -11)				1.0			
C22:0 Behenate			3.0	3.0			
C22:1 Erucate (<i>cis</i> -13)			20.0	5.0			
C24:0 Lignocerate			3.0	3.0			

1083 Rapeseed Oil Reference Mixture 25 mg/ml, 1 ml 54.00

Source: synthetic or plant **Appearance:** liquid **Solvent:** methylene chloride
Solubility: ethyl ether, methylene chloride **Storage:** -20°C

Suitable standard for low erucic acid oil

1084 RM-1 Mixture 50 mg 54.00

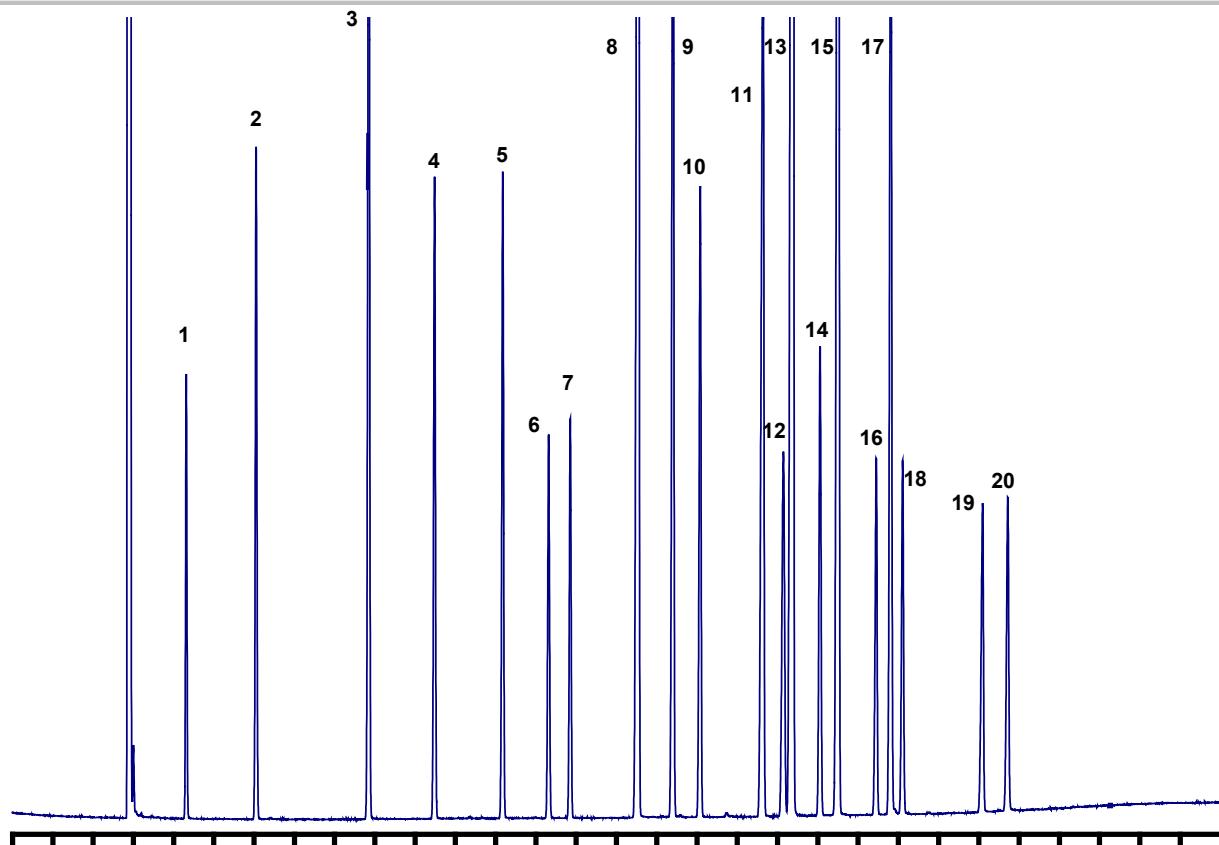
Source: synthetic or plant **Appearance:** liquid **Solubility:** chloroform, ethyl ether
Storage: -20°C

Suitable standard for corn, cottonseed, soybean, safflower, sesame, poppy seed, walnut kapok, and rice oils

1085	RM-2 Mixture	50 mg	54.00
Source: synthetic or plant Appearance: liquid Solubility: chloroform, ethanol, ethyl ether Storage: -20°C			
Suitable standard for linseed, perilla, hempseed, and rubberseed oils			
1086	RM-3 Mixture	50 mg/ml, 1 ml	54.00
Source: synthetic or plant Appearance: liquid Solvent: methylene chloride Solubility: ethyl ether, methylene chloride Storage: -20°C			
Suitable standards for peanut, rapeseed, and mustard seed oils			
1087	RM-4 Mixture	50 mg	54.00
Source: synthetic or plant Appearance: liquid Solubility: chloroform, ethyl ether Storage: -20°C			
Suitable standard for olive, teaseed, and neatsfoot oils			
1088	RM-5 Mixture	50 mg	54.00
Source: synthetic or plant Appearance: liquid Solubility: chloroform Storage: -20°C			
Suitable standard for coconut, palm kernel, babassu and ouri-ouri oils			
1089	RM-6 Mixture	50 mg	54.00
Source: synthetic or plant Appearance: liquid Solubility: ethyl ether, methylene chloride Storage: -20°C			
Suitable standard for lard, beef tallow, mutton tallow, and palm oil			

Custom Mixtures

Custom fatty acid methyl ester mixtures can be prepared to your specification. Minimum quantity requirements apply to these orders.



Cat# 4210 spiked with 0.4 mg/ml C18:2t ester (methyl linoelaidate) and chromatographed on a Supelco SP 2330 fused silica column.

Peak number	FAME
1	C8:0
2	C10:0
3	C12:0
4	C13:0
5	C14:0
6	C14:1
7	C15:0
8	C16:0
9	C16:1
10	C17:0
11	C18:0
12	C18:1t-9
13	C18:1c-9
14	C18:2t,t-9,12
15	C18:2c,c-9,12
16	C20:0
17	C18:3
18	C20:1
19	C22:0
20	C22:1

Table II. Standards for GC Analysis**GLC Standard Mixtures**

GLC-10 through GLC-100 standards are equal weight measures of fatty acid methyl esters. They are quantitative standards, useful for determining relative retention times and response factors.

Each methyl ester mixture is carefully prepared by weight and the composition verified by gas chromatography. The weight percentage of each component is indicated below. All double bonds are *cis*.

Mixture Number Catalog Number	GLC-10 1095	GLC-30 1097	GLC-40 1098	GLC-50 1099	GLC-60 1100	GLC-70 1101	GLC-80 1102	GLC-90 1103	GLC-100 1104
C8:0 Caprylate		20.0				20.0			
C9:0 Nonanoate						20.0			
C10:0 Caprate		20.0				20.0			
C11:0 Undecanoate						20.0			
C12:0 Laurate		20.0				20.0			
C13:0 Tridecanoate							20.0	20.0	
C14:0 Myristate		20.0					20.0		
C15:0 Pentadecanoate							20.0	20.0	
C16:0 Palmitate	20.0	20.0	25.0				20.0		
C16:1 Palmitoleate (<i>cis</i> -9)				20.0					
C17:0 Heptadecanoate							20.0	20.0	
C18:0 Stearate	20.0		25.0						20.0
C18:1 Oleate (<i>cis</i> -9)	20.0			20.0					
C18:2 Linoleate (all <i>cis</i> -9,12)	20.0								
C18:3 Linolenate (all <i>cis</i> -9,12,15)	20.0								
C19:0 Nonadecanoate								20.0	20.0
C20:0 Arachidate			25.0		25.0				20.0
C20:1 Eicosenoate (<i>cis</i> -11)				20.0	25.0				
C20:2 Eicosadienoate (all <i>cis</i> -11,14)					25.0				
C20:3 Eicosatrienoate (all <i>cis</i> -11,14,17)					25.0				
C21:0 Heneicosanoate								20.0	20.0
C22:0 Behenate			25.0						20.0
C22:1 Erucate (<i>cis</i> -13)				20.0					
C24:1 Nervonate (<i>cis</i> -15)				20.0					

1095 GLC-10 Mixture 50 mg 54.00

Source: synthetic or plant **Appearance:** liquid **Solubility:** methylene chloride
Storage: -20°C

1097 GLC-30 Mixture 50 mg 54.00

Source: synthetic or plant **Appearance:** liquid **Solubility:** methylene chloride
Storage: -20°C

1098 GLC-40 Mixture 50 mg/ml, 1 ml 54.00

Source: synthetic or plant **Appearance:** liquid **Solvent:** methylene chloride
Solubility: methylene chloride **Storage:** -20°C

1099	GLC-50 Mixture	50 mg/ml, 1 ml	54.00
Source: synthetic or plant Appearance: liquid Solvent: methylene chloride Solubility: methylene chloride Storage: -20°C			
1100	GLC-60 Mixture	50 mg/ml, 1 ml	54.00
Source: synthetic or plant Appearance: liquid Solvent: methylene chloride Solubility: methylene chloride Storage: -20°C			
1101	GLC-70 Mixture	50 mg	54.00
Source: synthetic or plant Appearance: liquid Solubility: methylene chloride Storage: -20°C			
1102	GLC-80 Mixture	50 mg/ml, 1 ml	54.00
Source: synthetic or plant Appearance: liquid Solvent: methylene chloride Solubility: methylene chloride Storage: -20°C			
1103	GLC-90 Mixture	50 mg/ml, 1 ml	54.00
Source: synthetic or plant Appearance: liquid Solvent: methylene chloride Solubility: methylene chloride Storage: -20°C			
1104	GLC-100 Mixture	50 mg/ml, 1 ml	128.00
Source: synthetic or plant Appearance: liquid Solvent: methylene chloride Solubility: methylene chloride Storage: -20°C			

Water Soluble Fatty Acid Mixtures

1106	WSFA-2 Mixture	5 ml	107.00
Water soluble fatty acid qualitative mixture			
Appearance: liquid Solvent: DI water Solubility: DI water Storage: Room Temp			
Contains: acetic, propionic, isobutyric, n-butyrlic, isovaleric and n-valeric acids			
1108	WSFA-4 Mixture	5 ml	107.00
Water soluble fatty acid qualitative mixture			
Appearance: liquid Solvent: DI water Solubility: DI water Storage: Room Temp			
Contains: acetic, propionic, isobutyric, n-butyrlic, 2-methylbutyric, isovaleric and n-valeric acids			

Microbiology Standard Mixtures

1105	GLC-110 Mixture Bacterial lipid standard, qualitative mixture	10 mg/ml, 1 ml	128.00
Source: various Appearance: liquid Solvent: chloroform Solubility: methylene chloride, chloroform Storage: -20°C			
Contains:			
Methyl 12-methyltridecanoate	(iso-C14:0)	Methyl 14-methylpentadecanoate	(iso-C16:0)
Methyl tetradecanoate (myristate)	(C14:0)	Methyl hexadecanoate (palmitate)	(C16:0)
Methyl 12-methyltetradecanoate	(anteiso-C15:0)	Methyl 14-methylhexadecanoate	(anteiso-C17:0)
Methyl pentadecanoate	(C15:0)		
 1114 Bacterial Acid Methyl Esters CP Mixture			
Qualitative mixture			
Source: various Appearance: liquid Solvent: methyl caproate Solubility: hexane, ethanol, methanol Storage: -20°C			
A qualitative standard. Mixture consists of equal amounts of the compounds listed.			
Methyl undecanoate	C11:0	Methyl cis-9-hexadecenoate (palmitoleate)	C16:1(cis-9)
Methyl 2-hydroxydecanoate	2-OH C10:0	Methyl hexadecanoate (palmitate)	C16:0
Methyl dodecanoate (laurate)	C12:0	Methyl 15-methylhexadecanoate	iso-C17:0
Methyl tridecanoate	C13:0	Methyl cis-9,10-methylenehexadecanoate	C17:0Δ (all cis-9,10)
Methyl 2-hydroxydodecanoate	2-OH C12:0	Methyl heptadecanoate (margarate)	C17:0
Methyl 3-hydroxydodecanoate	3-OH C12:0	Methyl 2-hydroxyhexadecanoate	2-OH C16:0
Methyl tetradecanoate (myristate)	C14:0	Methyl cis-9,12-octadecadienoate (linoleate)	C18:2 (all cis-9,12)
Methyl 13-methyltetradecanoate	iso-C15:0	Methyl cis-9-octadecenoate (oleate)	C18:1(cis-9)
Methyl 12-methyltetradecanoate	anteiso-C15:0	Methyl trans-9-octadecenoate (elaidate)	C18:1 (trans-9)
Methyl pentadecanoate	C15:0	Methyl octadecanoate (stearate)	C18:0
Methyl 2-hydroxytetradecanoate	2-OH C14:0	Methyl cis-9,10-methyleneoctadecanoate	C19:0Δ (all cis-9,10)
Methyl 3-hydroxytetradecanoate	3-OH C14:0	Methyl nonadecanoate	C19:0
Methyl 14-methylpentadecanoate	iso-C16:0	Methyl eicosanoate (arachidate)	C20:0
1075	Volatile Acid Mixture Qualitative mixture	100 ml	133.00
Appearance: liquid Solvent: DI water Solubility: DI water Storage: 4-8°C			
Contains: formic, acetic, propionic, isobutyric, n-butyric, isovaleric, n-valeric, isocaproic, n-caproic, and heptanoic acids			
1077	Non-Volatile Acid Mixture Qualitative mixture	100 ml	133.00
Appearance: liquid Solvent: DI water Solubility: DI water Storage: 4-8°C			
Contains: pyruvic, lactic, oxalacetic, oxalic, methyl malonic, malonic, fumaric and succinic acids.			

Biochemical Research Standard Mixtures

These mixtures are prepared by precise gravimetric technique. All mixtures contain equal amounts of listed components. A data sheet is supplied with each mixture.

1127	Polar Lipid Mixture TLC standards mixture	25 mg/ml, 1 ml	103.00
	Source: natural, egg, ovine Appearance: liquid Solvent: chloroform/methanol, 2:1 Solubility: chloroform/methanol, 2:1 Storage: -20°C		
	Contains: cholesterol, phosphatidylethanolamine, lecithin, and <i>lyso</i> -lecithin		
1128	Sphingolipid Mixture TLC standards mixture	25 mg/ml, 1 ml	103.00
	Source: natural, bovine Appearance: liquid Solvent: chloroform/methanol, 2:1 Solubility: chloroform/methanol, 2:1 Storage: -20°C		
	Contains: cerebrosides, sulfatides, and sphingomyelin		
1129	Non-Polar Lipid Mixture A TLC standards mixture	25 mg/ml, 1 ml	93.00
	Source: natural, plant, ovine Appearance: liquid Solvent: chloroform Solubility: chloroform Storage: -20°C		
	Contains: cholesteryl palmitate, tripalmitin, palmitic acid, and cholesterol		
1130	Non-Polar Lipid Mixture B TLC standards mixture	25 mg/ml, 1 ml	93.00
	Source: natural, plant, ovine Appearance: liquid Solvent: chloroform Solubility: chloroform Storage: -20°C		
	Contains: cholesteryl oleate, methyl oleate, triolein, oleic acid, and cholesterol		

Glycosphingolipid Reference Mixtures for TLC

These mixtures are qualitative standards prepared from our purified glycosphingolipids.

1505	Neutral Glycosphingolipid Mixture Glycosylceramides, qualitative mixture	1 mg/ml, 1 ml	136.00
	Source: natural, bovine and porcine Appearance: liquid Solvent: chloroform/methanol, 2:1 Solubility: chloroform/methanol, 2:1 Storage: -20°C		
	Contains: cerebrosides, lactosylceramides, ceramide trihexosides, globosides		

1508	Monosialoganglioside Mixture	0.5 mg/ml, 1 ml	136.00
Source: natural, bovine, human Appearance: liquid Solvent: chloroform/methanol/DI water, 2:1:0.1 Solubility: chloroform/methanol/DI water, 2:1:0.1 Storage: -20°C			
	Contains: GM ₃ , GM ₂ , GM ₁		
<hr/>			
1509	Disialoganglioside Mixture	0.5 mg/ml, 1 ml	147.00
Source: natural, bovine Appearance: liquid Solvent: chloroform/methanol/DI water, 2:1:0.1 Solubility: chloroform/methanol/DI water, 2:1:0.1 Storage: -20°C			
	Contains: GD ₃ , GD _{1a} , GD _{1b}		
<hr/>			
1510	Lactosylceramides and Sialosyl Derivatives Mixture	0.5 mg/ml, 1 ml	169.00
Source: natural, bovine Appearance: liquid Solvent: chloroform/methanol/DI water, 2:1:0.1 Solubility: chloroform/methanol/DI water, 2:1:0.1 Storage: -20°C			
	Contains: LC, GM ₃ , GD ₃		
<hr/>			
1511	Gangliotetraosylceramide and Sialosyl Derivatives Mixture	0.5 mg/ml, 1 ml	126.00
Source: natural, bovine Appearance: liquid Solvent: chloroform/methanol/DI water, 2:1:0.1 Solubility: chloroform/methanol/DI water, 2:1:0.1 Storage: -20°C			
	Contains: asialo-GM ₁ , GM ₁ , GD _{1a} , GD _{1b} , GT _{1b}		

Biochemicals and Reagents

Stable Isotope Labeled Compounds

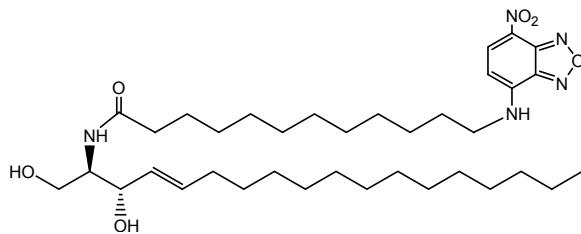
1914	N-Octadecanoyl-D₃₅-psychosine, (perdeuterated, C18:0 fatty acid) N-C18:0-D ₃₅ -Cerebrosides, perdeuterated; N-Stearoyl-D ₃₅ -psychosine, perdeuterated C ₄₂ H ₄₆ D ₃₅ NO ₈	5 mg	306.00
Source: semisynthetic, bovine Mol. Wt.: 762 Purity: 98+% by TLC Appearance: solid Solubility: chloroform, hot ethanol, chloroform/methanol, 2:1 Storage: -20°C			
1533	N-Hexadecanoyl-D₃-glucopsychosine, deuterated N-C16:0-D ₃ -Glucopsychosine, deuterated; N-C16:0-D ₃ -Glucocerebroside, deuterated; N-Palmitoyl-D ₃ -glucopsychosine, deuterated C ₄₀ H ₇₄ D ₃ NO ₈	1 mg	290.00
Source: semisynthetic, bovine buttermilk Mol. Wt.: 703 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C			
1534	N-Hexadecanoyl-D₃-lactosylceramide, deuterated N-C16:0-D ₃ -Lactosylceramide, deuterated; N-Palmitoyl-D ₃ -lactosylceramide, deuterated C ₄₆ H ₈₄ D ₃ NO ₁₃	1 mg	376.00
Source: semisynthetic, bovine buttermilk Mol. Wt.: 865 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 5:1:0.1 Storage: -20°C			

2200	N-1-¹³C-Hexadecanoyl-sphingosylphosphorylcholine D- <i>erythro</i> -Sphingomyelin with 1- ¹³ C-palmitic acid; N-1- ¹³ C-Palmitoyl-sphingosylphosphorylcholine ¹² C ₃₈ ¹³ CH ₇₉ N ₂ O ₆ P	1 mg	184.00
	Source: semisynthetic, bovine Mol. Wt.: 703 Purity: 98+% by TLC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: -20°C		
1536	N-Octadecanoyl-D₃-sulfatide, deuterated N-C18:0-D ₃ -Sulfatide, deuterated; N-Stearoyl-D ₃ -sulfatide, deuterated C ₄₂ H ₇₈ D ₃ NO ₁₁ S	1 mg	414.00
	Source: semisynthetic, bovine Mol. Wt.: 833 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1 Storage: -20°C		
1537	N-Octadecanoyl-D₃-ceramide trihexoside, deuterated C18:0-D ₃ -CTH, deuterated; N-C18:0-D ₃ -Gb3, deuterated; N-Octadecanoyl-D ₃ -globotriaosylceramide, deuterated; N-Stearoyl-D ₃ -ceramide trihexoside, deuterated C ₅₄ H ₉₈ D ₃ NO ₁₈	0.5 mg	348.00
	Source: semisynthetic, porcine Mol. Wt.: 1055 Purity: 98+% by TLC Appearance: solid Solubility: DMSO, chloroform/methanol, 2:1 Storage: -20°C		
2050	N-omega-CD₃-Octadecanoyl monosialoganglioside GM₁ (NH₄⁺ salt) N-CD ₃ -Stearoyl GM ₁ , C ₇₃ H ₁₂₈ N ₃ O ₃₁ D ₃ •NH ₃	0.5 mg	435.00
	Source: semisynthetic, bovine Mol. Wt.: 1550 + NH ₃ Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1; forms micellar solution in water Storage: -20°C		
2051	N-omega-CD₃-Octadecanoyl monosialoganglioside GM₂ (NH₄⁺ salt) N-CD ₃ -Stearoyl GM ₂ C ₆₇ H ₁₁₈ D ₃ N ₃ O ₂₆ •NH ₃	250 µg	319.00
	Source: semisynthetic, human Tay-Sachs Mol. Wt.: 1388 + NH ₃ Purity: 98+% by TLC, MS Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.1; forms micellar solution in water Storage: -20°C		
2052	N-omega-CD₃-Octadecanoyl monosialoganglioside GM₃ (NH₄⁺ salt) N-CD ₃ -Stearoyl GM ₃ C ₅₉ H ₁₀₅ D ₃ N ₂ O ₂₁ •NH ₃	250 µg	290.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 1185 + NH ₃ Purity: 98+% by TLC, MS Appearance: solid Solubility: chloroform/methanol/DI water, 2:1:0.2; forms micellar solution in water Storage: -20°C		

Fluorescent Compounds

1841	N-Hexanoyl-NBD-D-<i>erythro</i>-sphingosine	100 µg	204.00
1841-001	N-C6:0-NBD-Ceramide; N-C6:0-NBD-D- <i>erythro</i> -Sphingosine C ₃₀ H ₄₉ N ₅ O ₆ CAS#: 86701-10-2	1 mg	302.00
	Source: synthetic Mol. Wt.: 576 Melting Point (°C): 85-88 Purity: 98+% by TLC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: -20°C		

Catalog number 1618

Excitation: 460 nm
Emission: 535 nm

1618	N-Dodecanoyl-NBD-D-erythro-sphingosine	100 µg	147.00
1618-001	N-C12:0-NBD-Ceramide; N-C12:0-NBD-D- <i>erythro</i> -Sphingosine C ₃₆ H ₆₁ N ₅ O ₆	1 mg	526.00
Source: synthetic Mol. Wt.: 660 Purity: 98+% by TLC Appearance: solid			
Solubility: methanol, chloroform/methanol, 2:1 Storage: -20°C			
1857	N-Hexanoyl-NBD-L-<i>threo</i>-sphingosine	100 µg	162.00
1857-001	N-C6:0-NBD-Ceramide; N-C6:0-NBD-L- <i>threo</i> -Sphingosine C ₃₀ H ₄₉ N ₅ O ₆	1 mg	526.00
Source: synthetic Mol. Wt.: 575 Purity: 98+% by TLC Appearance: solid			
Solubility: chloroform, ethanol, methanol Storage: -20°C			
1620	N-Dodecanoyl-NBD-L-<i>threo</i>-sphingosine	100 µg	162.00
1620-001	N-C12:0-NBD-Ceramide; N-C12:0-NBD-L- <i>threo</i> -Sphingosine C ₃₆ H ₆₁ N ₅ O ₆	1 mg	526.00
Source: synthetic Mol. Wt.: 660 Purity: 98+% by TLC Appearance: solid			
Solubility: methanol, chloroform/methanol, 2:1 Storage: -20°C			
1624	N-Hexanoyl-NBD-L-<i>threo</i>-dihydrosphingosine	100 µg	147.00
1624-001	N-C6:0-NBD-Dihydroceramide; N-C6:0-NBD-L- <i>threo</i> -Dihydrosphingosine C ₃₀ H ₅₁ N ₅ O ₆	1 mg	526.00
Source: synthetic Mol. Wt.: 578 Purity: 98+% by TLC Appearance: solid			
Solubility: methanol, chloroform/methanol, 2:1 Storage: -20°C			
1623	N-Dodecanoyl-NBD-L-<i>threo</i>-dihydrosphingosine	100 µg	147.00
1623-001	N-C12:0-NBD-Dihydroceramide; N-C12:0-NBD-L- <i>threo</i> -Dihydrosphingosine C ₃₆ H ₆₃ N ₅ O ₆	1 mg	526.00
Source: synthetic Mol. Wt.: 662 Purity: 98+% by TLC Appearance: solid			
Solubility: methanol, chloroform/methanol, 2:1 Storage: -20°C			
1626	N-Hexanoyl-NBD-D-<i>erythro</i>-dihydrosphingosine	100 µg	126.00
1626-001	N-C6:0-NBD-Dihydroceramide; N-C6:0-NBD-D- <i>erythro</i> -Dihydrosphingosine C ₃₀ H ₅₁ N ₅ O ₆	1 mg	526.00
Source: synthetic Mol. Wt.: 578 Purity: 98+% by TLC Appearance: solid			
Solubility: methanol, chloroform/methanol, 2:1 Storage: -20°C			
1625	N-Dodecanoyl-NBD-D-<i>erythro</i>-dihydrosphingosine	100 µg	147.00
1625-001	N-C12:0-NBD-Dihydroceramide; N-C12:0-NBD-D- <i>erythro</i> -Dihydrosphingosine C ₃₆ H ₆₃ N ₅ O ₆	1 mg	526.00
Source: synthetic Mol. Wt.: 662 Purity: 98+% by TLC Appearance: solid			
Solubility: methanol, chloroform/methanol, 2:1 Storage: -20°C			

1628	N-Hexanoyl-NBD-phytosphingosine		100 µg	126.00
1628-001	N-C6:0-NBD-Phytoceramide; N-C6:0-NBD-Phytosphingosine C ₃₀ H ₅₁ N ₅ O ₇		1 mg	526.00
	Source: semisynthetic, bacteria Mol. Wt.: 594 Purity: 98+% by TLC Appearance: solid Solubility: methanol, chloroform/methanol, 2:1 Storage: -20°C			
1627	N-Dodecanoyl-NBD-phytosphingosine		100 µg	133.00
1627-001	N-C12:0-NBD-Phytoceramide; N-C12:0-NBD-Phytosphingosine C ₃₆ H ₆₃ N ₅ O ₇		1 mg	526.00
	Source: semisynthetic, bacteria Mol. Wt.: 678 Purity: 98+% by TLC Appearance: solid Solubility: methanol, chloroform/methanol, 2:1 Storage: -20°C			
1912	N-Hexanoyl-NBD-sphingosylphosphorylcholine		100 µg	111.00
1912-001	N-C6:0-NBD-Sphingomyelin; N-C6:0-NBD-Sphingosylphosphorylcholine C ₃₅ H ₆₁ N ₆ O ₉ P CAS#: 94885-04-8		1 mg	290.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 740 Purity: 98+% by TLC Appearance: solid Solubility: chloroform, ethanol, methanol Storage: -20°C Mixture of D-erythro and L-threo isomers			
1619	N-Dodecanoyl-NBD-sphingosylphosphorylcholine		100 µg	147.00
1619-001	N-C12:0-NBD-Sphingomyelin; N-C12:0-NBD-Sphingosylphosphorylcholine C ₄₁ H ₇₃ N ₆ O ₉ P		1 mg	319.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 825 Purity: 98+% by TLC Appearance: solid Solubility: methanol, chloroform/methanol, 2:1 Storage: -20°C Mixture of D-erythro and L-threo isomers			
1621	N-Hexanoyl-NBD-galactosylceramide		100 µg	147.00
1621-001	N-C6:0-NBD- <i>beta</i> -D-Galactosylsphingosine; N-C6:0-NBD-Cerebrosides C ₃₆ H ₅₉ N ₅ O ₁₁		1 mg	526.00
	Source: semisynthetic, bovine Mol. Wt.: 738 Purity: 98+% by TLC Appearance: solid Solubility: methanol, chloroform/methanol, 5:1 Storage: -20°C			
1633	N-Dodecanoyl-NBD-galactosylceramide		100 µg	118.00
1633-001	N-C12:0-NBD- <i>beta</i> -D-Galactosylsphingosine; N-C12:0-NBD-Cerebroside C ₄₂ H ₇₁ N ₅ O ₁₁		1 mg	526.00
	Source: semisynthetic, bovine spinal cord Mol. Wt.: 822 Purity: 98+% by TLC Appearance: solid Solubility: chloroform, DMSO, chloroform/methanol, 2:1 Storage: -20°C			
1622	N-Hexanoyl-NBD-glucosylceramide		100 µg	147.00
1622-001	N-C6:0-NBD- <i>beta</i> -D-Glucosylsphingosine; N-C6:0-NBD-Glucosylceramide, fluorescent C ₃₆ H ₅₉ N ₅ O ₁₁		1 mg	526.00
	Source: semisynthetic, bovine Mol. Wt.: 738 Purity: 98+% by TLC Appearance: solid Solubility: methanol, chloroform/methanol, 5:1 Storage: -20°C			
1629	N-Hexanoyl-NBD-lactosylceramide		50 ug	184.00
1629-001	N-Hexanoyl-NBD- <i>beta</i> -D-lactosylsphingosine; N-C6:0-NBD- <i>beta</i> -D-Lactosylsphingosine; N-C6:0-NBD-Lactosylceramide C ₄₂ H ₆₉ N ₅ O ₁₆		1 mg	824.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 900 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C			

1630	N-Dodecanoyl-NBD-lactosylceramide		50 µg	204.00
1630-001	N-Dodecanoyl-NBD- <i>beta</i> -D-lactosylsphingosine; N-C12:0-NBD- <i>beta</i> -D-Lactosylsphingosine; N-C12:0-NBD-Lactosylceramide C ₄₈ H ₈₁ N ₅ O ₁₆		1 mg	793.00
	Source: semisynthetic, bovine buttermilk Mol. Wt.: 984 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C			
1631	N-Dodecanoyl-NBD-ceramide trihexoside		100 µg	184.00
1631-001	N-C12:0-NBD-CTH; N-C12:0-NBD-Globotriaosylceramide C ₅₄ H ₉₁ N ₅ O ₂₁		1 mg	793.00
	Source: semisynthetic, porcine Mol. Wt.: 1145 Purity: 98+% by TLC Appearance: solid Solubility: DMSO; hot methanol, chloroform/methanol, 2:1 Storage: -20°C			
1632	N-Dodecanoyl-NBD-sulfatide		100 µg	119.00
1632-001	N-C12:0-NBD-Sulfatide; N-Dodecanoyl-NBD- <i>lyso</i> -sulfatide; N-Dodecanoyl-NBD-sphingosyl- <i>beta</i> -D-galactoside-3-sulfate C ₄₂ H ₇₁ N ₅ O ₁₄ S		1 mg	526.00
	Source: semisynthetic, bovine Mol. Wt.: 901 Purity: 98+% by TLC Appearance: solid Solubility: chloroform/methanol, 2:1 Storage: -20°C			
1634	omega-N-NBD-D-erythro-C14-Sphingosine		1 mg	522.00
	omega-N-(7-nitrobenzo-2-oxa-1,3-diazol-4-yl)-(2S)-amino-tetradec-(4E)-ene-(1,3R)-diol C ₂₀ H ₃₁ N ₅ O ₅			
	Source: synthetic Mol. Wt.: 422 Purity: 98+% by TLC Appearance: solid Solubility: methanol, ethanol, chloroform/methanol, 9:1 Storage: -20°C			

Table III. Typical Fatty Acid Composition of Natural Lipids Made by Matreya LLC
(actual composition may vary according to dietary history and growth condition of the source)

	Lecithin (egg)	Phosphatidyl-ethanolamine (egg)	lyso-Lecithin (egg)	Phosphatidylserine (bovine)	Phosphatidylinositol (plant)	Sulfatides (bovine)	Cerebrosides (bovine)	Sphingomyelin (bovine)	Phosphatidic acid (semi-synthetic)	Ceramides (bovine)
Catalog Number	#1044	#1045	#1046	#1047	#1048	#1049	#1050	#1051	#1053	#1056
Fatty Acids										
C14:0										
C16:0	31	17	72	1	42	trace	trace	4	39	trace
C16:1										
C18:0	16	29	24	42		5	4	40	12	4
C18:1	31	17	3	27	6	trace			34	
C18:2	16	11			47				15	
C18:3					5					
C20:0				1		1	1	3		1
C20:1				4						
C20:4		12		4						
C21:0										
C22:0				1		7	4	13		4
C22:1				1		trace				
C22:6				7						
C23:0							2	2		2
C24:0						18	10	9		10
C24:1						29	15	22		15
C25:0						2	3			9
C25:1						2	1			1
C26:0						1	2			2
C26:1						3	1			1
C27:0						1	2			2
C27:1							2			2
C14:0 2-OH										
C16:0 2-OH										
C18:0 2-OH						5	15			15
C20:0 2-OH						trace	1			1
C22:0 2-OH						3	6			6
C23:0 2-OH							5			5
C24:0 2-OH						10	17			17
C24:1 2-OH						6	6			
C25:0 2-OH						2	3			3
C25:1 2-OH										
C26:0 2-OH										
C26:1 2-OH										
C16 cis 9,10 methylene										
C18 cis 9,10 methylene										
Others	6	14	1	12	0	5	0	7	0	0
Total	100	100	100	100	100	100	100	100	100	100

Typical Fatty Acid Composition of Natural Lipids Made by Matreya LLC (continued)
 (actual composition may vary according to dietary history and growth condition of the source)

	Glucocerebrosides (Gaucher's spleen)	Monogalactosyl- diglycerides (plant)	Digalactosyl- diglyceride (plant)	Monosialo- ganglioside GM ₁	Disialoganglioside GD _{1a}	Trisialoganglio- side GT _{1b}	Gangliotetraosy- ceramide	Purified mixed gangliosides (bovine)	Cerebrosides Kerasin (bovine)	Ceramide trihexoside (porcine)
Catalog Number	#1057	#1058	#1059	#1061	#1062	#1063	#1064	#1065	#1066	#1067
Fatty Acids										
C14:0					1		trace	trace		
C16:0	26	23	9	2	1	1	1	1	trace	3
C16:1										
C18:0	9	77	91	90	89	87	86	86	5	2
C18:1						1	3	3		2
C18:2										
C18:3										
C20:0	5				3	2	4	4	1	2
C20:1										
C20:4										
C21:0										
C22:0	26				1	1	1	2	9	17
C22:1									trace	
C22:6										
C23:0	5					1	1	1	5	1
C24:0	22					1	1	1	25	29
C24:1	6			1		1	2	2	43	5
C25:0									3	
C25:1									3	
C26:0									2	
C26:1									4	
C27:0										
C27:1										
C14:0 2-OH										
C16:0 2-OH										
C18:0 2-OH										
C20:0 2-OH										
C22:0 2-OH									3	
C23:0 2-OH									1	
C24:0 2-OH									19	
C24:1 2-OH									10	
C25:0 2-OH										
C25:1 2-OH										
C26:0 2-OH										
C26:1 2-OH										
C16 cis 9,10 methylene										
C18 cis 9,10 methylene										
Others	1	0	0	3	6	3	0	0	0	6
Total	100	100	100	100	100	100	100	100	100	100

Typical Fatty Acid Composition of Natural Lipids Made by Matreya LLC (continued)
 (actual composition may vary according to dietary history and growth condition of the source)

	Globosides (porcine)	Lecithin (bovine)	Esterified steryl glucoside	Cerebrosides Phrenosin (bovine)	Phosphatidyl- ethanolamine (plant)	Lecithin (plant)	Ceramides (non- hydroxy)	Ceramides (hydroxy)	Sphingomyelin (porcine RBC)	Sphingomyelin (buttermilk)
Catalog Number	#1068	#1070	#1118	#1138	#1301	#1302	#1322	#1323	#1328	#1329
Fatty Acids										
C14:0		trace								1
C16:0	2	35	34		22	14			25	14
C16:1		1								
C18:0	2	14	8		3	4	11		7	3
C18:1		33	8		7	11				
C18:2			36		60	65				
C18:3			4		8	6				
C20:0	2		1				2		3	1
C20:1										
C20:4										
C21:0										
C22:0	20		4				10		9	26
C22:1										
C22:6										
C23:0	2		2				6		1	30
C24:0	33		2				24		22	21
C24:1	5						31		22	3
C25:0							3			
C25:1							3			
C26:0	2						2			
C26:1							3			
C27:0										
C27:1										
C14:0 2-OH										
C16:0 2-OH										
C18:0 2-OH				36			24			
C20:0 2-OH				1				1		
C22:0 2-OH	4			8				8		
C23:0 2-OH				6				6		
C24:0 2-OH	19			25				35		
C24:1 2-OH	9			9				17		
C25:0 2-OH				4				4		
C25:1 2-OH				2						
C26:0 2-OH				2						
C26:1 2-OH				2				2		
C16 cis 9,10 methylene										
C18 cis 9,10 methylene										
Others	0	17	1	5	0	0	5	3	11	1
Total	100	100	100	100	100	100	100	100	100	100

Typical Fatty Acid Composition of Natural Lipids Made by Matreya LLC (continued)
 (actual composition may vary according to dietary history and growth condition of the source)

	Sphingomyelin, (egg, chicken)	Phosphatidylinositol	Lactosyl ceramide (porcine)	Disialoganglioside GD1b	Monosialoganglioside GM2	Monosialoganglioside GM3 (buttermilk)	Disialoganglioside GD3 (buttermilk)	Lactosyl ceramide (buttermilk)	Ceramide trihexosides (top spot)
Catalog Number	#1332	#1336	#1500	#1501	#1502	#1503	#1504	#1507	#1513
Fatty Acids									
C14:0	trace			trace					
C16:0	72	32	14	1	2	6	8	12	1
C16:1									
C18:0	8	7	6	86	82	1	1	1	1
C18:1	3	7	4	3					
C18:2		47							
C18:3		6							
C20:0	2		1	4	7	1	1	1	2
C20:1									
C20:4									
C21:0						1	2		
C22:0	5		9	2	4	23	24	25	22
C22:1									
C22:6									
C23:0	1		1	1	trace	36	35	36	2
C24:0	2		15	1	1	22	21	21	58
C24:1	4		5	2	2	3	3		7
C25:0								1	1
C25:1									
C26:0									5
C26:1									
C27:0									
C27:1									
C14:0 2-OH									
C16:0 2-OH									
C18:0 2-OH			trace						
C20:0 2-OH									
C22:0 2-OH			8						
C23:0 2-OH									
C24:0 2-OH			24						
C24:1 2-OH			13						
C25:0 2-OH									
C25:1 2-OH									
C26:0 2-OH									
C26:1 2-OH									
C16 cis 9,10 methylene									
C18 cis 9,10 methylene									
Others	3	1	0	0	2	7	5	3	1
Total	100	100	100	100	100	100	100	100	100

Typical Fatty Acid Composition of Natural Lipids Made by Matreya LLC (continued)
 (actual composition may vary according to dietary history and growth condition of the source)

	Ceramide trihexosides (bottom spot)	Tetrasialoganglioside GC1 _b	Glucocerebrosides (buttermilk)	Glucocerebrosides (plant)	Mixed Gangliosides, purified (porcine)	Fucosylated monosialoganglioside GM ₁	Disialogangliosides GD ₂	Monosialoganglioside GM ₄
Catalog Number	#1514	#1516	#1521	#1522	#1525	#1526	#1527	#1535
Fatty Acids								
C14:0								
C16:0	3	5	7		1	8	1	4
C16:1		1						
C18:0		80	2		87	2	89	2
C18:1		2						
C18:2		3						
C18:3								
C20:0		4	1		4	13	7	trace
C20:1								trace
C20:4								
C21:0			1					
C22:0	2	2	27		1	43	1	3
C22:1								4
C22:6								
C23:0			36		1	3	1	4
C24:0	3		23		1	26		6
C24:1					2	5	1	4
C25:0			1					
C25:1								
C26:0								
C26:1								
C27:0								
C27:1								
C14:0 2-OH				trace				
C16:0 2-OH				79				
C18:0 2-OH	1			trace				1
C20:0 2-OH	1							3
C22:0 2-OH	11			8				25
C23:0 2-OH	1			1				17
C24:0 2-OH	52			9				18
C24:1 2-OH	25							7
C25:0 2-OH								
C25:1 2-OH								
C26:0 2-OH								
C26:1 2-OH								
C16 cis 9,10 methylene								
C18 cis 9,10 methylene								
Others	1	3	2	3	3			2
Total	100	100	100	100	100	100	100	100

Matreya

Lipids for Research

168 Tressler Street

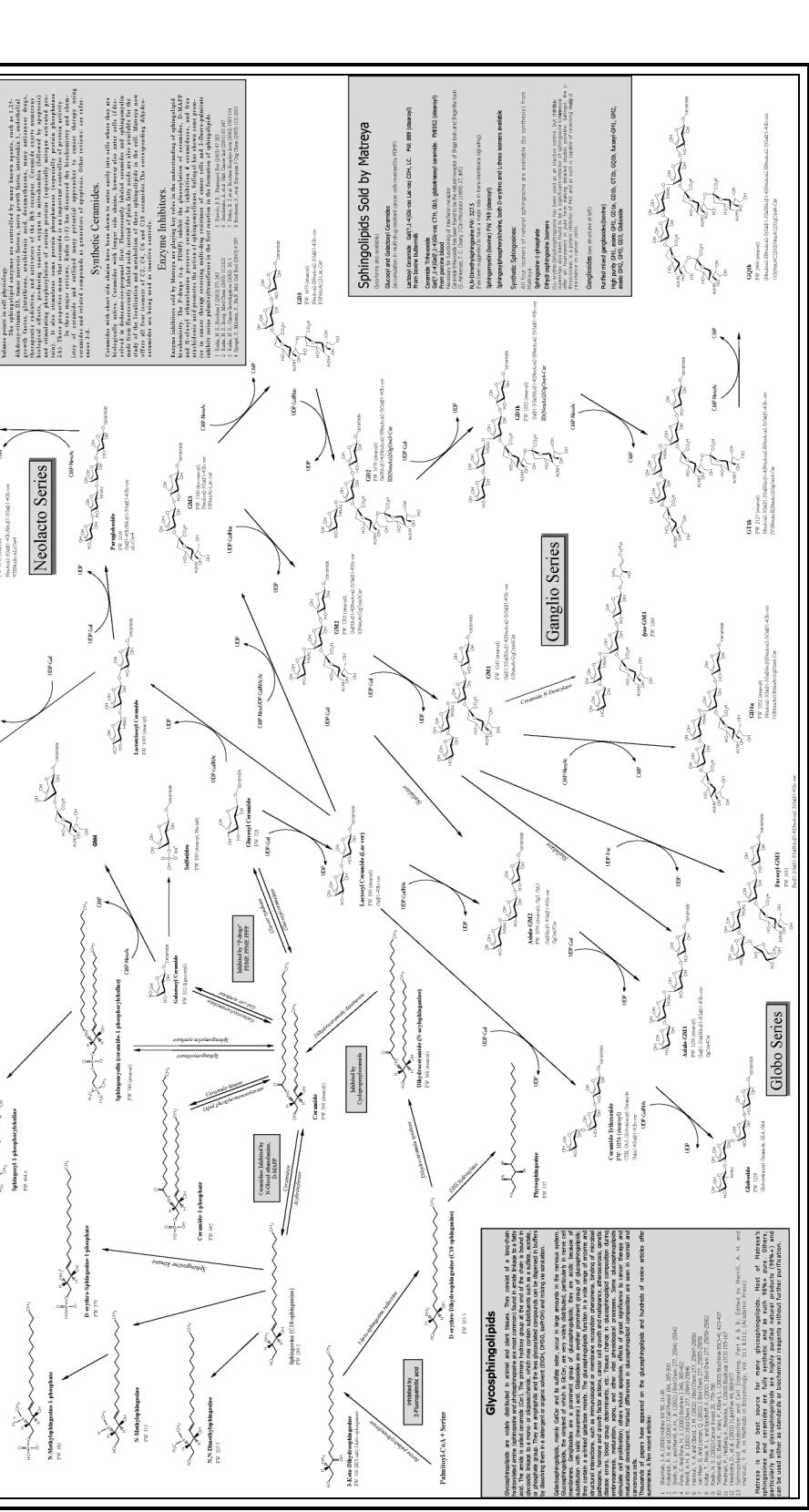
Pleasant Gap, PA 16823 USA

Tel: 814.359.5060 Fax: 814.359.5062

e-mail: customer service@matreya.com

www.matreya.com

Corporate 2000, Matreya LLC



Literature References

1. B.A. Fenderson, E.M. Eddy, S.Hakomori, BioEssays 12, 173, 1990
2. A. Gorio et al., Exp. Brain Res., Suppl. 13, 283, 1984
3. F. Di Gregorio et al., Neuropediatrics, Suppl. 15, 93, 1984
4. J.S. Schneider et al., Science, 256, 843, 1992
5. R.W. Leeden, R.K. Yu, Methods Enzymol, 83, 139, 1982
6. Borman, S., Chemical and Engineering News 82:32, 31-35, 2004
7. M. Faucher et al., J. Biol. Chem. 263, 5319, 1988
8. B. M. Buehrer, R.M. Bell, Adv. in Lipid Res., 26, 59, 1993
9. C. W. Sachs et al., J. Biol. Chem., 270, 26639, 1995
10. Radin, N. S., Biochem. J. 371:243, 2003
11. Radin, N. S., Biorg. Med. Chem. 11:2123, 2003
12. Radin, N. S., Cancer Investigation 20:1, 2002
13. Sietsma H., Veldman R.J., Vander Kolk D., Ausema B., Nijhof W., Kamps W., Vellenga E., Kok J.W. Clin. Cancer Res. 6, 942-948, 2000
14. Shayman, J. A., Kidney Int. 58, 11-26, 2000
15. Kolesnick, R. N. et al. J. Cell Physiol. 184, 285-300, 2000
16. Smith, W. L. and Merrill, A. H., Jr. J. Biol. Chem. 277, 25841-25842, 2002
17. Pyne, S. And Pyne, N. J. Biochem. J. 349, 385-402, 2000
18. Merrill, A. H. Jr. J. Biol. Chem. 277, 25843-25846, 2002
19. Hannun, Y. A. and Obeid, L., J. Biol. Chem. 277, 25847-25850, 2002
20. van Meer, G. And Lisman, Q. J. Biol. Chem. 277, 25855-25858, 2002
21. Kolter, T., Proia, R. L. and Sandhoff, K. J. Biol. Chem. 277, 25859-25862, 2002
22. Radin, N. S., Cancer Invest. 20, 779-786, 2002
23. Tettamanti, G., Bassi R., Viani, P., Riboni, L., Biochimie 85(3-4), 423-437, 2003
24. Fredman, P., Hedberg, K., Brezicka, T., Biodrugs 17(3) 155-167, 2003
25. Hoeckstra, D., et al. J. Lipid Res. 44, 869-877, 2003
26. R.R. Vunnam, N.S. Radin, Chem. Phys. Lipids, 265, 1980
27. J. Inokuchi and N.S. Radin, J. Lipid Res., 28, 565, 1987
28. N.S. Radin et al. J. Biochem, 111, 191, 1992
29. J. Inokuchi et al., Cancer letters, 38, 23, 1987
30. Y. Lavie et al., J. Biol. Chem., 271, 19530, 1996
31. Y. Lavie et al., J. Biol. Chem., 271, in press, 1996
32. Y. Hannun et al., Science, 235, 670, 1987
33. S. Spiegel et al., Proc. Intern. Conf. Biol. Function Glycosphingolipids, Santa Barbara, CA, 1990
34. Jacqueline M. Kravcka, Li Li, Zdzislaw M. Szulc, Jacek Bielawska, Besum Ogretmen, Yusuf A. Hannun, Lina M. Obeid, and Alicja Bielawska. J. Biol. Chem., 10, 1074/jbc. M700647200, February 5, 2007
35. M. Sugita et al. Biochim. Biophys. Acta, 398, 125, 1975
36. A. Bielawska et al. J. Biol. Chem., 271, 12646, 1996
37. A. Bielawska et al. ibid. 267, 18493, 1992
38. J. M. L. Hauser et al., J. Biol. Chem. 269, 6803, 1994
39. Y. Hannun, ibid, 269, 3125, 1994
40. R. Kolesnick, D.W. Golde, Cell, 77, 325, 1984
41. M.W. Pariza et al., Cancer Res. 43, 2444s, 1983
42. Y. L Ha et al., J. Agr. Food Chem., 37, 75, 1989
43. P.W. Parodi, Austral. J. Dairy Tech., 49, 93, 1994
44. S. Banni et al. Abstr. 87th AOCS Mtg., 28, 1996
45. C. Ip et al., Cancer Res. 51, 6118, 1991
46. M. A. Belury, Nutr. Rev. 53 (4 Pt. 1), 83, 1995
47. C. Liew et al. Carcinogenesis, 16, 3037, 1995
48. T. D. Shultz et al. Cancer Let, 63, 125, 1992
49. R.J. Nicolosi, Abstr. 87th AOCS Mtg., 1996
50. K.N. Lee et al., Atheroscl., 108(1), 19, 1994
51. B.F. Haumann, Inform, 7(2), 152, 1996
52. M.W. Pariza et al., Abstr., 87th AOCS Mtg., 1996
53. M.A. Belury et al., Lipids, 32, 199, 1997
54. Houseknecht, K.L. et al., Biochem. Biophys. Res. Comm., 224, 678, 1998
55. W.W. Christie et al. JAOCS, 74, 1231, 1997
56. Sehat, N. et al., Lipids, 33, 217, 1998
57. Sébédio, J. L. et al., Biochim. Biophys. Acta, 1345,5, 1997

Note: We are pleased and honored to have the above list of scientists in the field of Lipid Research. In case your name and contribution is not listed above, we apologize. If you would like your publication listed in our next catalog, please send your name and publication to the attention of Marketing Department.

Cross Reference for Product Numbers and Catalog Pages

1006	76	1077	86	1192	57
1008	49	1081	79	1193	57
1009	49	1083	81	1194	58
1010	50	1084	81	1195	73
1011	50	1085	81, 82	1196	48
1012	50	1086	81, 82	1197	48
1013	50	1087	81, 82	1198	49
1014	50	1088	81, 82	1199	49
1015	50	1089	81, 82	1200	48
1016	53	1093	79	1203	54
1017	53	1095	84	1204	54
1018	50	1097	84	1205	56
1019	51	1098	84	1206	56
1020	51	1099	84, 85	1207	72
1021	51	1100	84, 85	1208	54
1022	54	1101	84, 85	1241	51
1023	54	1102	84, 85	1242	51
1024	55	1103	84, 85	1243	53
1025	55	1104	84, 85	1244	58
1026	56	1105	86	1245	61
1027	56	1106	85	1245-1	61
1028	51	1108	85	1245-10	61
1029	51	1113	77	1247-1	62
1030	51	1114	86	1247-10	62
1031	51	1115	76	1248	62
1032	57	1116	76	1248-1	62
1033	57	1117	77	1249	62
1034	57	1118	77, 97	1249-1	62
1035	52	1119	76	1249-10	62
1036	52	1120	77	1251	52
1037	52	1121	77	1252	52
1038	52	1122	77	1254	62
1040	53	1123	76	1255	61
1041	59	1124	79	1256	62
1042	57	1125	79	1257	62
1044	40, 94	1127	87	1258	62
1045	41, 94	1128	87	1259	62
1046	40, 94	1129	87	1261	49
1047	40, 94	1130	87	1262	55, 59
1048	41, 94	1131	60, 80	1263	55, 59
1049	22, 96	1136	58	1264	58
1050	19, 96	1138	20, 96	1265	58
1051	15, 95	1147	53, 59	1266	55
1051-1	15	1148	54, 59	1267	55
1052	41	1149	54, 59	1269	57
1053	41, 95	1150	54, 59	1271	52
1056	11, 97	1151	55, 60	1273	53
1057	21, 98	1152	55, 60	1275	53
1057-25	21	1153	56	1276	56
1058	48, 95	1154	56	1277	56
1059	48, 95	1155	59	1301	41, 94
1061	29, 95	1156	59	1302	40, 94
1061-50	29	1157	53	1303	47
1062	31, 95	1161	50	1303-2	47
1063	31, 96	1162	50	1305	20
1064	29, 96	1163	49	1306	22
1065	32, 96	1164	49	1310	22
1066	19, 96	1165	49	1318	17
1067	25, 97	1166	49	1319	17
1067-10	25	1167	58	1320	5
1068	26, 97	1175	58	1321	17
1069	41	1177	79	1321-05	17
1070	40, 94	1179	57	1322	11, 97
1071	74	1181	60, 61	1322-05	11
1072	74	1182	70	1323	12, 97
1073	74	1183	70	1323-05	12
1074	75	1186	52	1324	4
1075	86	1187	52	1325	20

1326	4	1537	26, 27, 89	1713	65
1327	16	1538	25	1714	65
1328	15, 95	1600	71	1715	65
1329	15, 98	1601	71	1716	65
1329-1	15	1602	71	1717	39
1330	5	1603	71	1718	39
1330-1	5	1605	71	1719	37
1332	15, 97	1606	71	1720	37
1332-1	15	1612	72	1722	65, 80
1333	8	1613	72	1725	66
1333-100	8	1614	72	1725-0.5	66
1334	20	1615	72	1726	66
1334-50	20	1616	72	1726-0.5	66
1335	20	1618	13, 90	1727	66
1336	41	1618-001	13, 90	1727-0.5	66
1400	43	1619	18, 91	1728	67
1409	43, 63	1619-001	18, 91	1728-0.5	67
1410	43, 63	1620	13, 90	1729	67
1411	44, 63	1620-001	13, 90	1729-0.5	67
1425	43	1621	21, 27, 91	1730	67
1426	43	1621-001	21, 27, 91	1730-0.5	67
1427	43	1622	21, 28, 91	1731	67
1428	42	1622-001	21, 28, 91	1731-0.5	67
1429	42	1623	14, 90	1732	67
1430	42	1623-001	14, 90	1732-0.5	67
1431	44	1624	13, 90	1733	67
1432	44	1624-001	13, 90	1733-0.5	67
1433	44	1625	14, 90	1734	67
1434	45	1625-001	14, 90	1734-0.5	67
1435	45	1626	14, 90	1735	67
1436	45	1626-001	14, 90	1735-0.5	67
1437	43	1627	14, 91	1736	67
1438	44	1627-001	14, 91	1736-0.5	67
1439	45	1628	14, 91	1739	68
1442	42	1628-001	14, 91	1739-0.5	68
1443	44	1629	25, 28, 91	1740	68
1444	45	1629-001	25, 28, 91	1740-0.5	68
1445	43	1630	25, 28, 92	1741	68
1500	24, 98	1630-001	25, 28, 92	1741-0.5	68
1501	31, 96	1631	26, 28, 92	1742	68
1502	30, 98	1631-001	26, 28, 92	1742-0.5	68
1503	30, 98	1632	24, 28, 92	1743	68
1504	31, 98	1632-01	24, 28, 92	1743-0.5	68
1505	32, 87	1633	21, 28, 91	1744	68
1507	24, 98	1633-001	21, 28, 91	1744-0.5	68
1507-50	24	1634	2, 92	1745	66
1508	32, 88	1656	70	1745-0.5	66
1509	32, 88	1657	71	1746	66
1510	33, 88	1701	64	1746-0.5	66
1511	33, 88	1701-1	64	1747	66
1512	29	1702	64	1747-0.5	66
1513	25	1702-1	64	1748	66
1514	25	1703	64	1748-0.5	66
1516	32, 97	1703-1	64	1749	37
1517	24	1704	64	1750	39
1518	30	1704-1	64	1751	35, 73
1520	26	1705	64	1752	39
1521	21, 98	1705-1	64	1753	38
1521-50	21	1706	64	1754	68
1522	21, 98	1706-1	64	1755	38
1522-100	21	1707	64	1756	38
1523	26	1707-1	64	1758	63
1524	26	1708	64	1758-1	63
1525	32	1708-1	64	1759	64
1526	30	1709	65	1759-1	64
1527	31	1709-0.5	65	1760	69
1531	22	1710	65	1761	69
1532	24	1710-0.5	65	1766	70
1533	22, 27, 88	1711	65	1773	46
1534	24, 27, 88	1711-0.5	65	1773-1	46
1535	31, 97	1712	65	1773-5	46
1536	23, 27, 89	1712-0.5	65	1775	47

1775-1	47	1882	69	2046	18
1775-5	47	1883	70	2047	11
1778	47	1884	70	2050	30,89
1778-1	47	1886	36	2051	30,89
1778-5	47	1886-005	36	2052	30,89
1779	46	1887	36	2076	23
1779-1	46	1887-005	36	2077	10
1783	47	1888	23	2200	16, 89
1783-1	47	1889	37	4210	78
1783-5	47	1890	16		
1784	46	1891	5		
1784-1	46	1892	5		
1784-5	46	1893	5		
1786	35, 73	1894	12		
1790	75	1895	12		
1791	73	1896	6		
1792	73	1897	12		
1797	75	1900	7		
1800	38	1900-100	7		
1802	2	1901	7		
1803	17	1901-100	7		
1805	6	1903	8		
1806	2	1903-100	8		
1807	3, 35	1904	23		
1807-025	3, 35	1907	15		
1809	8	1909	15		
1810	8	1910	10		
1818	69	1911	16		
1819	70	1912	18, 91		
1822	73	1912-001	18, 91		
1823	73	1913	17		
1826	2	1914	20, 27, 88		
1827	2	1915	8		
1828	7	1915-100	8		
1829	7	1916	9		
1830	8	1916-25	9		
1831	3	1917	16		
1831-1	3	1918	16		
1832	9	1919	46		
1832-100	9	1919-1	46		
1833	2	1919-5	46		
1834	10	1931	23		
1835	3	1930	9		
1838	3	1930-25	9		
1839	4	1932	23		
1840	3	1933	23		
1841	13, 89	1950	33		
1841-001	13, 89	1951	33		
1842	10	1954	33		
1843	9	1957	34		
1845	4	1960	34		
1846	4	1961	34		
1847	7	1963	34		
1848	7	1964	34		
1850	9	1977	33		
1851	4	2009	78		
1852	18	2010	78		
1854	10	2011	80		
1855	9	2034	12		
1857	13, 90	2035	12		
1857-001	13, 90	2036	13		
1858	39	2037	8		
1859	36	2037-100	8		
1860	36	2038	8		
1865	38	2038-100	8		
1868	38	2039	9		
1875	23	2039-100	9		
1876	4	2041	11		
1877	69	2042	9		
1878	69	2043	11		
1880	69	2044	10		
1881	69	2045	11		

Product Name Index

10-Methylhexadecanoic acid	73	Behenic acid	52	Ceramide trihexosides (bottom spot)
12-Methyltetradecanoic acid	72	Behenic acid methyl ester	52	25,98
13-Methyltetradecanoic acid	71			Ceramide trihexosides (top spot)
14-Methylhexadecanoic acid	72			25,97
15-Hydroxypentadecanoic acid	69			Ceramides (hydroxy and non-
15-Hydroxypentadecanoic acid methyl ester	69			hydroxy acyl groups)
15-Methylhexadecanoic acid	71	Capric acid methyl ester	49	11,94
17-Hydroxyheptadecanoic acid	69	Caproic acid methyl ester	48	Ceramides (mostly hydroxy
17-Hydroxyheptadecanoic acid methyl ester	69	Caprylic acid	49	acyl groups)
20-Hydroxyeicosanoic acid	69	Caprylic acid methyl ester	49	12,96
20-Hydroxyeicosanoic acid methyl ester	69	Castanospermine	38	Ceramides (mostly non-hydroxy acyl groups)
21-Hydroxyheneicosanoic acid methyl ester	69	Ceramide-1-phosphate, N-C16:0-D- <i>erythro</i>	18	11,96
22-Hydroxydocosanoic acid methyl ester	69	Ceramide, N-C10:0-D- <i>erythro</i>	8	Cerebroside; Kerasin (top spot)
22-Hydroxydocosanoic acid methyl ester	69	Ceramide, N-C12:0-NBD-D- <i>erythro</i>	13,90	19,95
27-Hydroxyheptacosanoic acid methyl ester	70	Ceramide, N-C12:0-NBD-L- <i>threo</i>	13,90	Cerebroside, N-C12:0-NBD
30-Hydroxytriacontanoic acid methyl ester	70	Ceramide, N-C15:0-D- <i>erythro</i>	8	21,28,91
6-Hydroxyoctadecanoic acid	70	Ceramide, N-C16:0-D- <i>erythro</i>	8	Cerebroside, N-C15:0
		Ceramide, N-C17:0-D- <i>erythro</i>	8	20
		Ceramide, N-C18:0-D- <i>erythro</i>	9	Cerebroside, N-C18:0-D- ₃₅
		Ceramide, N-C18:0-D- <i>threo</i>	9	20,27,88
		Ceramide, N-C18:0-L- <i>erythro</i>	9	Cerebroside, N-C2:0
		Ceramide, N-C18:0-L- <i>threo</i>	9	20
		Ceramide, N-C19:0-D- <i>erythro</i>	9	Cerebroside, N-C6:0-NBD
		Ceramide, N-C2:0-D- <i>erythro</i>	7	21,27,91
		Ceramide, N-C2:0-D- <i>erythro</i>	60,80	Cerebroside, N-C8:0
		(C14 sphingoid base)		Cerebroside, Phrenosin
A				(bottom spot)
Acetyl-4-(1R,2S,3R,4-tetrahydroxybutyl)-imidazole	39	Ceramide, N-C2:0-L- <i>erythro</i>	7	20,96
Alditol Acetate Mixture-1	79	Ceramide, N-C2:0-L- <i>threo</i>	7	Cerebroside sulfate
Alditol Acetate Mixture-2	79	Ceramide, N-C24:0-D- <i>erythro</i>	9	22,94
anteiso-Heptadecanoic acid	72	Ceramide, N-C24:1 (cis-15)-D- <i>erythro</i>	9	Cerebrosides
anteiso-Heptadecanoic acid methyl ester	72	Ceramide, N-C6:0-D- <i>erythro</i>	7	19,94
anteiso-Palmitic acid methyl ester	72	Ceramide, N-C6:0-L- <i>erythro</i>	7	Cerotic acid
anteiso-Pentadecanoic acid	72	Ceramide, N-C6:0-L- <i>threo</i>	7	52
anteiso-Pentadecanoic acid methyl ester	72	Ceramide, N-C6:0-NBD-D- <i>erythro</i>	13,89	Cerotic acid methyl ester
Anti-ganglioside asialo GM ₁	33	Ceramide, N-C6:0-NBD-L- <i>threo</i>	13,90	52
Anti-ganglioside asialo GM ₂	33	Ceramide, N-C8:0-D- <i>erythro</i>	8	Cholestan, 5- <i>alpha</i>
Anti-ganglioside GD _{1b}	34	Ceramide, N-C8:0-D- <i>threo</i>	8	76
Anti-ganglioside GD ₂	34	Ceramide, N-C8:0-L- <i>threo</i>	8	Cholesterol
Anti-ganglioside GD ₃	33	Ceramide, N-(R,S)- <i>alpha</i> -hydroxy-C12:0-D- <i>erythro</i>	9	76
Anti-ganglioside GM ₁	33	Ceramide, N-(R,S)- <i>alpha</i> -hydroxy-C18:0-D- <i>erythro</i>	10	Cis-Trans Isomer Standard Mixture
Anti-ganglioside GM ₂ (NANA),	34	Ceramide, N-hexadecanoyl-D- <i>erythro</i> (C16 sphingoid base)	10	60,80
Anti-ganglioside GM ₄	34	Ceramide, N-hexanoyl-D- <i>threo</i>	8	Conduritol B Epoxide
Anti-globoside GL-4	34	Ceramide trihexoside, lyso	26	37
Antibody asialo GM ₁	33	Ceramide trihexoside, N-C17:0	26	Coprostanol
Antibody asialo GM ₂	33	Ceramide trihexoside, N-C12:0-NBD	26,28,92	Cyclopropenylceramide, C8:0-N
Antibody disialoganglioside GD _{1b}	34	Ceramide trihexoside, N-C18:0-D3	26,27,89	36
Antibody disialoganglioside GD ₂	34	Ceramide trihexoside, N-C23:0	26	Dihydroceramide, N-C2:0-
Antibody ganglioside GD ₃	33	Ceramide trihexoside, N- dodecanoyl-NBD	26,28,92	D- <i>erythro</i>
Antibody globoside	34	Ceramide trihexoside, N-heptadecanoyl	26	10
Antibody monosialoganglioside GM ₁	33	Ceramide trihexoside, N-octadecanoyl-D ₃	26,27,89	Dihydroceramide, N-C6:0-
Antibody monosialoganglioside GM ₂ (NANA)	34	Ceramide trihexoside, N-stearoyl-D ₃	26,27	D- <i>threo</i>
Antibody monosialoganglioside GM ₄	34	Ceramide trihexoside, N-tricosanoyl	26	10
Arachidic acid	51	Ceramide trihexosides	25,95	Dihydroceramide, N-(R,S)- <i>alpha</i> -hydroxy-C16:0-D- <i>erythro</i>
Arachidic acid methyl ester	51			11
Arachidonic acid	57			Dihydroceramide, N-(R,S)- <i>alpha</i> -hydroxy-C18:0-D- <i>erythro</i>
Arachidonic acid methyl ester	57			11
Asialo GM ₁	29			Dihydroceramide, N-(R,S)- <i>alpha</i> -hydroxyoctadecanoyl-D- <i>erythro</i>
Asialo GM ₂	29			11
				Dihydrophosphatidylethanolamine, D- <i>erythro</i>
B				3
Bacterial Acid Methyl Esters CP Mixture	86			Dihydrophosphatidylethanolamine, D- <i>erythro</i> -C ₂₀
				4
				Dihydrophosphatidylethanolamine, D- <i>threo</i>
				4
				Dihydrophosphatidylethanolamine, D,L-C ₁₆
				4
				Dihydrophosphatidylethanolamine, D,L- <i>erythro</i>
				4
				Dihydrophosphatidylethanolamine, L- <i>erythro</i>
				4

Dihydrophingosine, L- <i>threo</i>	3,35	Distearoyl-sn-glycero-3-phosphorylcholine	43	GLC-110 Mixture	86
Dihydrophingosine, N-acetyl-D- <i>erythro</i>	10	Distearoyl-sn-glycero-3-phosphorylethanamine	45	GLC-30 Mixture	84
Dihydrophingosine, N-dodecanoyl-NBD-D- <i>erythro</i>	14,90	Distearoyl-sn-glycero-3-phosphorylglycerol	44	GLC-40 Mixture	84
Dihydrophingosine, N-dodecanoyl-NBD-L- <i>threo</i>	14,90	Docosahexaenoic acid	58	GLC-50 Mixture	84,85
Dihydrophingosine, N-hexanoyl-D- <i>erythro</i>	10	Docosahexaenoic acid methyl ester	58	GLC-60 Mixture	84,85
Dihydrophingosine, N-hexanoyl-NBD-D- <i>erythro</i>	14,90	Docosanoic acid	52	GLC-70 Mixture	84,85
Dihydrophingosine, N-hexanoyl-NBD-L- <i>threo</i>	13,90	Docosanoic acid methyl ester	52	GLC-80 Mixture	84,85
Dihydrophingosine, N-octadecanoyl-D- <i>erythro</i>	11	Docosapentaenoic acid	58	GLC-90 Mixture	84,85
Dihydrophingosine, N-octanoyl-D- <i>erythro</i>	10	Docosapentaenoic acid methyl ester	58	Globosides	26,96
Dihydrophingosine, N-(R,S)-alpha-hydroxydodecanoyl-D- <i>erythro</i>	11	Docosenoic acid	58	Globotriaosylceramide	25,95
Dihydrophingosine, N-(R,S)-alpha-hydroxyhexadecanoyl-D- <i>erythro</i>	11	Docosenoic acid methyl ester	58	Glucocerebroside, N-C16:0-D ₃	22,27,88
Dihydrophingosine, N-(R,S)-alpha-hydroxyoctadecanoyl-D- <i>erythro</i>	11	Dodecanoic acid	49	Glucocerebroside, N-C22:0	22
Dihydrophingosine-1-phosphate, D- <i>erythro</i>	17	Dodecanoic acid methyl ester	49	Glucocerebroside-lyso, bovine buttermilk	22
Dihydrophingosine•HCl (3-keto)	4	Dotriacontanoic acid methyl ester	53	Glucocerebroside-lyso, plant	22
Dihydrophingosine•HCl (3-keto-C ₆)	5	E		Glucocerebrosides, bovine buttermilk	21,98
Dihydrophingosine•HCl (3-keto-C ₈)	5	E-64-d	39	Glucocerebrosides, Gaucher's spleen	21,95
Dihydrophingosine•HCl (3-keto-C ₁₂)	5	Eicosadienoic acid	57	Glucocerebrosides, plant	21,98
Dihydrophingosylphosphoryl-choline	17	Eicosadienoic acid methyl ester	57	Glucopsychosine, bovine buttermilk	22
Dihydrosterculic acid	73	Eicosanoic acid	51	Glucopsychosine, N-docosanoyl	22
Dilauroyl-sn-glycero-3-phosphorylcholine	42	Eicosanoic acid methyl ester	51	Glucopsychosine, N-hexadecanoyl-D ₃	22,27,88
Dilauroyl-sn-glycero-3-phosphorylethanamine	45	Eicosapentaenoic acid	58	Glucopsychosine, plant	22
Dilauroyl-sn-glycero-3-phosphorylglycerol	44	Eicosapentaenoic acid methyl ester	58	Glucosylceramide, bovine buttermilk	21,98
Dimethylheptanoic, D,L-2,6 acid	72	Eicosatetraenoic acid	57	Glucosylceramide, Gaucher's spleen	21,95
Dimethyltocol-5,8	74	Eicosatrienoic acid methyl ester	57	Glucosylceramide, N-hexanoyl-NBD	21,28,91
Dimethyltocol-7,8	74	Eicosatrienoic acid methyl ester (all cis-8,11,14)	57	Glucosylceramide, plant	21,98
Dimethyltocol, rac-5,7	75	Elaidic acid	54,59	Glucosylceramide, N-C6:0-NBD-beta-D	21,28,91
Dimyristoyl-sn-glycero-3-phosphatidic acid	42	Elaidic acid methyl ester	54	Glucosylceramide, plant	22
Dimyristoyl-sn-glycero-3-phosphorylcholine	43	EPA	58	Glycosphingolipid, Mixture, neutral	32,87
Dimyristoyl-sn-glycero-3-phosphorylethanamine	45	Ergosterol	77	GM ₁	29,95
Dimyristoyl-sn-glycero-3-phosphorylglycerol	44	Erucic acid	58	GT ₁₁	36
Dipalmitoyl-sn-glycero-3-phosphatidic acid	42	Erucic acid methyl ester	58		
Dipalmitoyl-sn-glycero-3-phosphorylcholine	43	EST	39		
Dipalmitoyl-sn-glycero-3-phosphorylethanamine	45	Esterified Steryl Glucosides	77,96		
F					
FIM-FAME-6 Mixture		F			
FIM-FAME-7 Mixture		Galactosylceramide, N-dodecanoyl-NBD	21,28,91	H	
Fucosylated monosialoganglioside GM ₁	30,98	Galactosylceramide, N-hexanoyl-NBD	21,27,91	Heneicosanoic acid	51
Fluoropalmitic acid	39	Galactosylceramide, N-octanoyl-beta-D-	20	Heneicosanoic acid methyl ester	51
Fluoropalmitic acid methyl ester	39	Gangliosides, mixed, purified, bovine	32,95	Heptadecanoic acid	50
G		Gangliosides, mixed, purified, porcine	32,98	Heptadecanoic acid methyl ester	50
Galactosylceramide, N-dodecanoyl-NBD	21,28,91	Gangliotetraosylceramide	29,95	Hexadecanoic acid	50
Galactosylceramide, N-hexanoyl-NBD	21,27,91	Gangliotetraosylceramide and Sialosyl Derivatives Mixture	33,88	Hexadecanoic acid methyl ester	50
Galactosylceramide, N-octanoyl-beta-D-	20	Gangliotriosylceramide	29	Hexanoic acid methyl ester	48
Gangliosides, mixed, purified, bovine	32,95	Gb ₃	25,95	Hydroxy Methyl Ester Mixture	80
Gangliosides, mixed, purified, porcine	32,98	Gb ₄	26,96	Hydroxy Acid (α) Methyl Ester Mixture	65
Gangliotetraosylceramide	29,95	GLC-10 Mixture	84	Hydroxydecanoic acid (α)	63
Gangliotetraosylceramide and Sialosyl Derivatives Mixture	33,88	GLC-100 Mixture	84,85	Hydroxydecanoic acid (α) methyl ester	63
Gangliotriosylceramide	29			Hydroxydecanoic acid (β)	66
Gb ₃	25,95				
Gb ₄	26,96				

Hydroxydocosanoic acid (α)	65	K	KEL-FIM-FAME-5 Mixture	78	M	
methyl ester	65				Margaric acid	50
Hydroxydodecanoic acid (α)	64	L			Margaric acid methyl ester	51
methyl ester	64	L-MAPP	36		Mead acid methyl ester	57
Hydroxydodecanoic acid (β)	67	Laccernoic acid methyl ester	53		Melissnoic acid methyl ester	53
methyl ester	67	Lactosylceramide and Sialosyl			Methyl 10(E),12(Z)-octadecadienoate	62
Hydroxyeicosanoic acid (α)	64	Derivatives Mixture	33,88		Methyl 10-methylhexadecanoate	73
methyl ester	65	Lactosylceramide, lyso	24		Methyl 11-methylhexadecanoate	70
Hydroxyeicosanoic acid (α)	64	Lactosylceramide, N-C12:0-NBD	25,28,92		Methyl 12-methyltetradecanoate	72
methyl ester	65	Lactosylceramide, N-C16:0	24		Methyl 12-methyltridecanoate	71
Hydroxyheptadecanoic acid (β)	68	Lactosylceramide, N-C16:0-D ₃	24,27,88		Methyl 13-methylpentadecanoate	72
methyl ester	68	Lactosylceramide, N-C17:0	25		Methyl 13-methyltetradecanoate	71
Hydroxyhexadecanoic acid (α)	64	Lactosylceramide, N-C6:0-NBD	25,28,91		Methyl 14-methylhexadecanoate	72
methyl ester	64	Lactosylceramide, N-dodecanoyl			Methyl 14-methylpentadecanoate	71
Hydroxyhexadecanoic acid (β)	68	Lactosylceramide, N-heptadecanoyl	25		Methyl 15-hydroxypentadecanoate	69
methyl ester	68	Lactosylceramide, N-hexadecanoyl	24		Methyl 15-methylhexadecanoate	71
Hydroxyhexadecanoic acid (β)	68	Lactosylceramide, N-hexadecanoyl-D ₃	24,27,88		Methyl 17-hydroxyheptadecanoate	69
methyl ester	68	Lactosylceramide, N-hexanol-NBD	25,28,91		Methyl 17-methyloctadecanoate	71
Hydroxyoctadecanoic acid (α)	64	Lactosylceramide, N-palmitoyl	24		Methyl 20-hydroxyeicosanoate	69
methyl ester	64	Lactosylceramide, N-palmitoyl-D ₃	24,27,88		Methyl 21-hydroxyheneicosanoate	69
Hydroxyoctadecanoic acid (α)	64	Lactosylceramides, bovine			Methyl 22-hydroxydocosanoate	70
methyl ester	64	buttermilk	24,97		Methyl 27-hydroxyheptacosanoate	70
Hydroxyoctadecanoic acid (β)	68	Lactosylceramides, porcine	24,97		Methyl 2-fluoropalmitate	39
methyl ester	68	Lactocerebrosides, bovine			Methyl 2-hydroxydecanoate	63
Hydroxyoctanoic acid (β)	66	buttermilk	24,97		Methyl 2-hydroxydocosanoate	65
Hydroxyoctanoic acid (β)	66	Lactocerebrosides, porcine	24,97		Methyl 2-hydroxydodecanoate	64
methyl ester	66	Lanosterol	77		Methyl 2-hydroxyeicosanoate	65
Hydroxytetracosanoic acid (α)	65	Lauric acid	49		Methyl 2-hydroxyhexadecanoate	64
Hydroxytetracosanoic acid (α)	65	Lauric acid methyl ester	49		Methyl 2-hydroxyoctadecanoate	64
methyl ester	65	Lecithin, bovine	40,96		Methyl 2-hydroxytetraicosanoate	65
Hydroxytetradecanoic acid (α)	64	Lecithin, egg	40,94		Methyl 2-hydroxytetradecanoate	64
Hydroxytetradecanoic acid (α)	64	Lecithin, lyso, egg	40,94		Methyl 3-hydroxydodecanoate	67
methyl ester	64	Lecithin, plant	40,96		Methyl 3-hydroxyheptadecanoate	68
Hydroxytetradecanoic acid (β)	67	Lignoceric acid	52		Methyl 3-hydroxyhexadecanoate	68
Hydroxytetradecanoic acid (β)	67	Lignoceric acid methyl ester	52		Methyl 3-hydroxyhexanoate	66
methyl ester	67	Linoelaidic acid	55,60		Methyl 3-hydroxynonanoate	66
Hydroxytricosanoic acid (α)	65	Linoelaidic acid methyl ester	55,60		Methyl 3-hydroyoctadecanoate	68
Hydroxytricosanoic acid (α)	65	Linoleic acid	55		Methyl 3-hydroyoctanoate	66
methyl ester	65	Linoleic acid methyl ester	55		Methyl 3-hydroxytetradecanoate	67
Hydroxytridecanoic acid (β)	67	Linolenic acid	56		Methyl 3-hydroxytridecanoate	67
Hydroxytridecanoic acid (β)	67	Linolenic acid-gamma	56		Methyl 3-hydroxyundecanoate	67
methyl ester	67	Linolenic acid-gamma methyl ester	56		Methyl 9(E),11(E)-octadecadienoate	62
Hydroxyundecanoic acid (β)	67	Linolenic acid methyl ester	56		Methyl 9(Z) 11(E)-octadecadienoate	61,62
Hydroxyundecanoic acid (β)	67	Long Chain Fatty Acid Methyl Ester Mixture	80		Methyl 9(Z) 11(Z)-octadecadienoate	62
methyl ester	67	Loxastatin	39		Methyl arachidate	51
I		lyso-Ceramide trihexoside	26		Methyl arachidonate	57
iso-Heptadecanoic acid	71	lyso-Cerebroside	20		Methyl behenate	52
iso-Heptadecanoic acid	71	lyso-Dihydrosphingomyelin	17		Methyl caprate	49
methyl ester	71	lyso-Glucocerebroside, bovine			Methyl caproate	48
iso-Nonadecanoic acid		buttermilk	22		Methyl cerotate	52
methyl ester	71	lyso-Glucocerebroside, plant	22		Methyl cis-9,10-methyleneoctadecanoate, C19:0 delta	73
iso-Palmitic acid methyl ester	71	lyso-Lactosylceramide	24		Methyl decanoate	49
iso-Pentadecanoic acid	71	lyso-Lecithin, egg	40,94		Methyl dihydrosterculate	73
iso-Pentadecanoic acid	71	lyso-Monosialoganglioside GM ₁	30		Methyl docosahexaenoate	59
methyl ester	71	lyso-Phosphatidylcholine, egg	40,94		Methyl docosanoate	52
iso-Tetradecanoic acid	71	lyso-Sphingomyelin	17		Methyl docosapentaenoate	58
methyl ester	71	lyso-Sulfatide	23		Methyl docosenoate	58
iso-Tridecanoic acid methyl ester	70				Methyl dodecanoate	49
					Methyl dotriacontanoate	53
					Methyl eicosadienoate	57
					Methyl eicosanoate	51
					Methyl eicosapentaenoate	58

Phosphatidylinositol, plant, soy	41,97	Sphinganine, D- <i>erythro-C</i> ₂₀	4	Sphingosine, N-heptadecanoyl- <i>D-erythro</i>	8
Phosphatidylinositol tris-3,4,5-phosphate, dipalmitoyl (Na^+ salt)	47	Sphinganine, D- <i>threo</i>	4	Sphingosine, N-hexadecanoyl- <i>D-erythro</i>	8
Phosphatidylinositol tris-3,4,5-phosphate, dipalmitoyl (NH_4^+ salt)	47	Sphinganine, D,L- <i>C</i> ₁₆	4	Sphingosine, N-hexadecanoyl- <i>D-erythro</i> (C16 sphingoid base)	10
Phosphatidylserine, bovine	40,94	Sphinganine, D,L- <i>erythro</i>	4	Sphingosine, N-hexanoyl-D- <i>erythro</i>	7
Phosphoglycerides Kit	41	Sphinganine, D,L- <i>erythro-C</i> ₂₀	4	Sphingosine, N-hexanoyl-D- <i>threo</i>	8
Phospholipid (MPL) of <i>Thermoplasma acidophilum</i> (>50%)	47	Sphinganine, L- <i>erythro</i>	4	Sphingosine, N-hexanoyl- <i>erythro</i>	7
Phospholipid (MPL) of <i>Thermoplasma acidophilum</i> (>95%)	47	Sphinganine, L- <i>threo</i>	3,35	Sphingosine, N-hexanoyl-NBD- <i>erythro</i>	14,90
Phytanic acid	73	Sphinganine, N-acetyl-D- <i>erythro</i>	10	Sphingosine, N-hexanoyl-NBD- <i>threo</i>	14,90
Phytoceramide, N-C12:0-NBD	14,91	Sphinganine, N-dodecanoyl-NBD-D- <i>erythro</i>	14,90	Sphingosine, N-hexanoyl-NBD- <i>threo</i>	13,89
Phytoceramide, N-C16:0	12	Sphinganine, N-dodecanoyl-NBD-L- <i>threo</i>	14,90	Sphingosine, N-hexanoyl-NBD-L- <i>threo</i>	13,90
Phytoceramide, N-C18:0	12	Sphinganine, N-hexanoyl-NBD-L- <i>threo</i>	14,90	Sphingosine, N-N, dihexyl-D- <i>erythro</i>	6
Phytoceramide, N-C2:0	12	Sphinganine, N-hexanoyl-NBD-L- <i>threo</i>	13,90	Sphingosine, N,N-dimethyl-D- <i>erythro</i>	5
Phytoceramide, N-C24:0	13	Sphinganine, N-octadecanoyl-D- <i>erythro</i>	11	Sphingosine, N-nonadecanoyl-D- <i>erythro</i>	9
Phytoceramide, N-C6:0	12	Sphinganine, N-octanoyl-D- <i>erythro</i>	10	Sphingosine, N-octadecanoyl-D- <i>erythro</i>	9
Phytoceramide, N-C6:0-NBD	14,91	Sphinganine, N-(R,S)-alpha-hydroxydodecanoyl-D- <i>erythro</i>	11	Sphingosine, N-octadecanoyl-D- <i>threo</i>	9
Phytoceramide, N-C8:0	12	Sphinganine, N-(R,S)-alpha-hydroxyhexadecanoyl-D- <i>erythro</i>	11	Sphingosine, N-octadecanoyl-L- <i>erythro</i>	9
Phytosphingosine	5	Sphinganine, N-(R,S)-alpha-hydroxyoctadecanoyl-D- <i>erythro</i>	11	Sphingosine, N-octadecanoyl-L- <i>threo</i>	9
Phytosphingosine, N-acetyl	12	Sphingolipid Mixture	87	Sphingosine, N-octanoyl-D- <i>erythro</i>	8
Phytosphingosine, N-dodecanoyl-NBD	14,91	Sphingomyelin, bovine	15,94	Sphingosine, N-octanoyl-D- <i>threo</i>	8
Phytosphingosine, N-hexadecanoyl	12	Sphingomyelin, bovine buttermilk	15,96	Sphingosine, N-octanoyl-L- <i>erythro</i>	8
Phytosphingosine, N-hexanoyl	12	Sphingomyelin, D- <i>erythro</i> with 1- ¹³ C-palmitic acid	16,89	Sphingosine, omega-N-NBD-D- <i>erythro-C</i> ₁₄	2,92
Phytosphingosine, N-hexanoyl-NBD	14,91	Sphingomyelin, egg	15,97	Sphingosine, N-pentadecanoyl-D- <i>erythro</i>	8
Phytosphingosine, N-octadecanoyl	12	Sphingomyelin, N-C12:0-NBD	18,91	Sphingosine, N-(R,S)-alpha-hydroxydodecanoyl-D- <i>erythro</i>	9
Phytosphingosine, N-octanoyl	12	Sphingomyelin, N-C17:0	16	Sphingosine, N-(R,S)-alpha-hydroxyoctadecanoyl-D- <i>erythro</i>	10
Phytosphingosine, N-tetraacosanoyl	13	Sphingomyelin, N-C18:0	16	Sphingosine, N-tetraacosanoyl-D- <i>erythro</i>	9
Plant Sterol Mixture	76	Sphingomyelin, N-C2:0	15	Sphingosine, N-tetracosenoyl-D- <i>erythro</i>	9
Plant Sterols Kit	76	Sphingomyelin, N-C20:0-D- <i>erythro</i>	16	Sphingosylphosphorylcholine	17
Polar Lipid Mixture	87	Sphingomyelin, N-C22:0-D- <i>erythro</i>	16	Sphingosylphosphorylcholine, D- <i>erythro</i>	17
PPMP, D- <i>threo</i>	38	Sphingosine-1-phosphate, N-hexadecanoyl-D- <i>erythro</i>	18	Sphingosylphosphorylcholine, L- <i>threo</i>	17
PPMP, D,L- <i>erythro</i>	38	Sphingosine-1-phosphate, N-C6:0	15	Sphingosylphosphorylcholine, N-1- ¹³ C-hexadecanoyl	16,89
PPMP, D,L- <i>threo</i>	37	Sphingosine-1-phosphate, N-C6:0-NBD	18,91	Sphingosylphosphorylcholine, N-acetyl	15
PPMP, L- <i>threo</i>	38	Sphingomyelin, porcine	15,96	Sphingosylphosphorylcholine, N-eicosanoyl-D- <i>erythro</i>	16
Psychosine	20	Sphingosine-1-phosphate, D- <i>erythro</i>	17	Sphingosylphosphorylcholine, N-docosanoyl-D- <i>erythro</i>	16
Psychosine, N-acetyl	20	Sphingosine, D- <i>erythro</i>	2	Sphingosylphosphorylcholine, N-dodecanoyl-NBD	18,91
Psychosine, N-octadecanoyl-D ₃₅	20,27,88	Sphingosine, D- <i>erythro-C</i> ₁₂	3	Sphingosylphosphorylcholine, N-heptadecanoyl	16
Psychosine, N-pentadecanoyl	20	Sphingosine, D- <i>erythro-C</i> ₁₄	2	Sphingosylphosphorylcholine, N-hexanoyl	15
PUFA-1 Mixture	79	Sphingosine, D- <i>erythro-C</i> ₁₆	3	Sphingosylphosphorylcholine, N-hexanoyl-NBD	18,91
PUFA-2 Mixture	79	Sphingosine, D- <i>erythro-C</i> ₂₀	3	Sphingosylphosphorylcholine, N-octadecanoyl	16
PUFA-3 Mixture	79	Sphingosine, D- <i>threo</i>	2	Sphingosylphosphorylethanolamine, N-acyl	16
R					
Rapeseed Oil Reference Mixture	81	Sphingosine, L- <i>erythro</i>	2	Stearic acid	51
Ricinelaidic acid	70	Sphingosine, L- <i>threo</i>	2	Stearic acid methyl ester	51
RM-1 Mixture	81	Sphingosine, N-acetyl-D- <i>erythro</i>	7	Stearidonic acid	56
RM-2 Mixture	82	Sphingosine, N-acetyl-D- <i>erythro</i> (C14 sphingoid base)	10	Stearidonic acid methyl ester	56
RM-3 Mixture	82	Sphingosine, N-acetyl-L- <i>erythro</i>	7		
RM-4 Mixture	82	Sphingosine, N-acetyl-L- <i>threo</i>	7		
RM-5 Mixture	82	Sphingosine, N-decanoil-D- <i>erythro</i>	8		
RM-6 Mixture	82	Sphingosine, N-dodecanoyl-NBD-D- <i>erythro</i>	13,90		
Royal Jelly acid	68	Sphingosine, N-dodecanoyl-NBD-D- <i>threo</i>	13,90		
S					
Safingol	3,35	Sphingosine, N-dodecanoyl-NBD-L- <i>threo</i>	13,90		
Sapienic acid	53	Sphingosine, N-N, dihexyl-D- <i>erythro</i>	6		
Sitostanol- <i>beta</i>	77	Sphingosine, N-N, dihexyl-D- <i>threo</i>	13,90		
Sphinganine•HCl (3-keto)	4	Sphingosine, N-octadecanoyl-L- <i>erythro</i>	9		
Sphinganine, D- <i>erythro</i>	3	Sphingosine, N-octadecanoyl-L- <i>threo</i>	9		

Steraryl-2-[10(E),12(Z)-octadecadienoyl]-sn-glycero-3-phosphorylcholine	44,63	Sulfatide, N-octadecanoyl-D ₃	23,27,89	Tricosanoic acid	52
Steraryl-2-[9(Z),11(E)-octadecadienoyl]-sn-glycero-3-phosphorylcholine	44,63	Sulfatide, N-octadecenoyl	23	Tricosanoic acid methyl ester	52
Steraryl-2-linoleoyl-sn-glycero-phosphorylcholine	43,63	Sulfatide, N-oleoyl	23	Tridecanoic acid	50
Steryl glucosides	77	Sulfatide, N-palmitoyl	23	Tridecanoic acid methyl ester	50
Stigmastanol	77	Sulfatide, N-stearoyl	23	Trimethyltocol-5,7,8	74
Stigmasterol	77	Sulfatide, N-stearoyl-D ₃	23,27,89	Trisialoganglioside GT _{1b}	31,95
Sulfatide-lyso	23	Sulfatide, N-tetracosanoyl	23		
Sulfatide, N-acetyl	23	Sulfatide, tetracosenoyl	23		
Sulfatide, N-C12:0-NBD	24,28,92	Sulfatides	22,94		
Sulfatide, N-C16:0	23				
Sulfatide, N-C18:0	23				
Sulfatide, N-C18:0-D ₃	23,27,89				
Sulfatide, N-C18:1	23				
Sulfatide, N-C2:0	23				
Sulfatide, N-C24:0	23				
Sulfatide, N-C24:1	23				
Sulfatide, N-dodecanoyl-NBD	24,28,92				
Sulfatide, N-hexadecanoyl	23				
Sulfatide, N-lignoceroyl	23				
Sulfatide, N-nervonyl	23				
Sulfatide, N-octadecanoyl	23				

T

Tetracontanoic acid methyl ester	53
Tetracosanoic acid	52
Tetracosanoic acid methyl ester	52
Tetracosenoic acid	59
Tetracosenoic acid methyl ester	59
Tetradecanoic acid	50
Tetradecanoic acid methyl ester	50
Tetradecenoic acid methyl ester	53
Tetrasialoganglioside GQ _{1b}	32,98
THI	39
Tocol-rac	75
Tocopherol, (+)-delta	75
Tocopherol, rac-alpha	74
Tocopherol, rac-beta	74
Tocopherol, rac-gamma	74

U

Undecanoic acid	49
Undecanoic acid methyl ester	49

V

Vaccenic acid (<i>cis</i> -11)	55
Vaccenic acid methyl ester (<i>cis</i> -11)	55
Vaccenic acid methyl ester (<i>trans</i> -11)	55,59
Vaccenic acid (<i>trans</i> -11)	55,59
Volatile Acid Mixture	86

W

WSFA-2 Mixture	85
WSFA-4 Mixture	85

Matreya LLC Ordering Information

Office Hours: Monday - Friday, 8:00am - 4:00pm, Eastern Time

• Ordering and Customer Service

Telephone (Toll Free)	(800) 342 3595
Telephone (worldwide)	(814) 359 5060
Fax (24/7 worldwide)	(814) 359 5062
e-mail	customerservice@matreya.com
e-commerce website	www.matreya.com

• Technical Service

Telephone (worldwide)	(814) 359 5060
Fax (worldwide)	(814) 359 5062
e-mail	techservice@matreya.com

• Terms

Prices and discounts are subject to change without notice. Freight charges are prepaid and added as a separate charge on the invoice. Orders placed are F.O.B. Pleasant Gap, PA. **Net 30 Days.**

• Shipping

Standard orders:	FedEx 2 nd -Day
Rush Orders:	FedEx Priority or Standard overnight
International Orders:	FedEx International Priority
Items requiring dry ice for shipping will be charged a \$40 fee.	

• Return Policy

All returns must have prior authorization. Items incorrectly ordered by you and returned to Matreya are subject to a 25% (\$25 minimum) restocking charge. The following items may not be returned: frozen products, items which have passed their expiration dates, custom synthesized products and accommodation orders. If a product has been incorrectly sent to you due to an error on our part, a credit will be issued to your account immediately. Please inspect and verify your order upon receipt. No products will be returned after 30 days (assuming proper storage and handling by customer).

• Disclaimer

The information and data included in this catalog are correct and reliable to the best of our knowledge. However, we offer no guarantees and assume no responsibility for this information. No license or immunity under any patent is granted or implied through our sale of any material. Any information sent to us by the customer, relative to his or her interest in Matreya's products, must be submitted voluntarily and without obligation on the part of Matreya.

Matreya LLC
168 Tressler Street
Pleasant Gap, PA 16823 USA
Federal ID No. 20-1237500

International Dealers & Representatives

Australia (South Australia)

Adelab Scientific
36 Holland Street
Thebarton 5031
Tel: 61-88-234-7955
Fax: 61-88-234-7897
Email: info@adelab.com.au
Web: www.adelab.com.au

Canada

MJS Biolynx
PO Box 1150
300 Laurier Blvd
Brockville, Ontario K6V 5W1
Tel: 613-498-2126
Toll free: 888-593-5969
Fax: 613 342 1341
Email: sales@biolynx.ca
Web: www.biolynx.ca

China

Shenzhen Bolin Chemical Co. LTD
Rm 1309, No. 1 News Bldg
2 Shennan Zhong Road
Shenzhen, 518027
Tel: 867-556-165-1418 (1428)
Fax: 867-558-209-6552
Email: lh@bolin-lktlabs.com
Web: www.bolin-lktlabs.com

France

Biovalley
18 Route de Tournan
Marne la Vallée Cedex 3
Conches, 77601
Tel: 33-16-007-2020
Fax: 33-16-007-5051
Email: biovalley@biovalley.fr
Web: www.biovalley.fr

Germany

BIOTREND Chemikalien GmbH
Eupenerstr. 157
Köln, 50933
Tel: 49-221-949-8320
Fax: 49-221-949-8325
Email: info@biotrend.com
Web: www.biotrend.com

Italy

Superchrom S.r.l.
Via C. Menotti, 11
Milano, 20129
Tel: 39-02-738-6315
Fax: 39-027-010-0100
Email: Superchrom@iol.it

Japan

Techno Chemical Corp
27-9, Honkomagome 1 Chome
Bunkyo-ku, Tokyo 113
Tel: 81-33-947-7310
Fax: 81-33-947-7306
Email: info@technochemical.com
Web: www.technochemical.com

Korea

Kim & Friends, Inc.
SK Twintech Tower B-304
345-9 Gasan-dong, Geumcheon-gu
Seoul 153-773
Tel: 82-2-2647-6611
Fax: 82-2-2647-6687
Email: kslee@kimnfriends.co.kr
Web: www.kimnfriends.co.kr

Netherlands

Distrilab BV
Olmenlaan 6-C
Leusden, 3833 AV
Tel: 31-33-494-7834
Fax: 31-33-432-1441
Email: info@distrilab.nl
Web: www.distrilab.nl

New Zealand

Phenomenex NZ Ltd.
PO Box 31-6-1
Milford, Auckland
Tel: 649-478-0951
Fax: 649-478-0952
Email: info@phenomenex.co.za
Web: www.phenomenex.com

Spain

Teknokroma S.C.C.L.
Camí de Can Calders, 14
Apartado de Correos, 147
Sant Cugat del Vallès
Barcelona ES-08173
Tel: 34-93-674-8800
Fax: 34-93-675-2405
Email: commercial@teknokroma.es
Web: www.teknokroma.es

Sweden

Larodan Fine Chemicals AB
Limhamnsgardens Alle'9
Malmö, S-21616
Tel: 464-016-4155
Fax: 464-015-5498
Email: info@larodan.se
Web: www.larodan.se

Switzerland

VWR International AG-Life Sciences
Lerzenstrasse 16/18
Dietikon, 8953
Tel: 41-44-745-1436
Fax: 41-44-745-1410
Email: info@ch.vwr.com
Web: www.vwr.com

Taiwan

Super Chroma Enterprise Ltd
5F.226 Roosevelt Rd, Section 5
Taipei ROC, 116
Tel: 88-622-930-7110
Fax: 88-622-930-7112
Email: sc@superchroma.com.tw
Web: www.superchroma.com.tw

United Kingdom (England)

Universal Biologicals (Cambridge) Ltd.
Passhouse Farmhouse
Papworth St. Agnes
Cambridge CB23 3QU
Tel: 44-148-083-9015
Fax: 44-148-803-1912
Email: info@universalbiologicals.com
Web: www.universalbiologicals.com