



CHEMFACES



Chem Faces 🔅

By axxo chemicals and services co., ltd

Chamberry (?)



@axxochem

Chem Frees 🗭

Separation and Purification

axe

ChemFaces is a professional high-purity natural products manufacturer, Our products are sold to research institutes, university laboratories, and research and development departments of enterprises and other scientific research units all of the world. Natural compounds have good biological activity, the scientists and scientific researchers use ChemFaces products found that most of the products have a very good research value and obtained very good scientific research results. ChemFaces updates its product inventory every month, mainly with new additions and new batches. Because our products are unique for a long time. So we hope that our newly developed products can help the scientists and scientific researchers around the world. For more information please refer to ChemFaces' official website (www.chemfaces.com).





Our Services

ax

- Construct high-quality natural product library for drug research and development. More than 4000 compounds are available, many of them are our unique products.
- Provide reference substances and active components of Traditional Chinese Medicine.
- Provide more products of inhibitors in addition to a variety of screening libraries.
- Provide isolation and structure determination of natural products.
- Contract research, contract manufacture and process development services from lab scale, pilot scale to commercial scale.

Quality & Validation

ChemFaces Natural Products pay careful attention to the purity and stability our products. We can provide the chemical test data by NMR, LC-MS, and HPLC. If you are in any way unsatisfied with the quality of our products, we will unconditionally refund or replace the products in question.

Tech Support

ChemFaces helps our customers use the products to their full potential. Our team is dedicated to customer satisfaction and you can expect helpful advice on all aspects of our products when you Contact Us. (service@chemfaces.com)

ChemFaces is a professional high-purity natural products manufacturer.

Product Intended Use

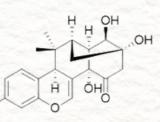
- 1. Reference standards
- 2. Pharmacological research
- 3. Inhibitors



qr-code for ChemFaces website



Pueraria mirifica/ White kwao krua



Miroestrol



-10

CAS No.	2618-41-9	Purity :	>=98%
Catalog No.	CFN91612	Type of Compound:	Diterpenoids
Molecular Formula:	C ₂₀ H ₂₂ O ₆	Physical Desc.:	Powder
Molecular Weight:	358.39 g/mol	Source:	The herbs of Pueraria mirifica
Solvent:	Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.	Price:	Inquiry

Biological Activity

Description: Miroestrol is a highly active phytoestrogen. Miroestrol can produce mammogenic effect. Miroestrol exhibits bone loss prevention and neuroprotective in ovariectomized mice. Miroestrol also can reduce cancer risk. Miroestrol Description

Source: The herbs of Pueraria mirifica

Solvent: Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.

Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months (2-8°C).

Wherever possible, you should prepare and use solutions on the same day. However, if you need to make up stock solutions in advance, we recommend that you store the solution as aliquots in tightly sealed vials at -20°C. Generally, these will be useable for up to two weeks. Before use, and prior to opening the vial we recommend that you allow your product to equilibrate to room temperature for at least 1 hour.

After receiving: The packaging of the product may have turned upside down during transportation, resulting in the natural compounds adhering to the neck or cap of the vial. take the vial out of its packaging and gently shake to let the compounds fall to the bottom of the vial. for liquid products, centrifuge at 200-500 RPM to gather the liquid at the bottom of the vial. try to avoid loss or contamination during handling.

CERTIFIC	ATE OF ANALYSIS
	Miroestrol
[Catalog No.] CFN91612	
[CAS.No.] 2018-41-9	
[Purity] >=98%	
[M.E.] C.,H.,O,	
[Physical Description] Powder	
[Solvent] Chioroform, Dichloromethan	e, Ethyl Acetate, DMSO, Acetone, etc.
[Wisight] 10mg	
[Lot.No.] CFS202101	
[Storage] Protected from air and light,	refrigerate or freeze(2-8 °C)
[Intended Use] For laboratory use onl	y
[Shelf Life] 2 years [Chemical Structure]	HO OHO HO
[Characterization Data Summary]	
Analytical Test	Results
identification by NMR (Purity tested	Consistent with the above structure
frond more	>=05%
Authorized Signature:	
Date:	
Carlo.	
	chemfaces, com i Technological Development Zone, Wahan, Habel 4300
	: service@chemfaces.com Tel: 0086-27-84)



Kaempferia parviflora; KP



5,7-dimethoxyflavone/ Chrysin dimethylether

CAS No.	21392-57-4	Purity	>=98%
Catalog No.	CFN90896	Type of Compound	Flavonoids
Molecular Formula	$C_{17}H_{14}O_4$	Physical Desc.	Powder
Molecular Weight	282.3 g/mol	Source	The herbs of Pueraria mirifica
Solvent	Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.	Price	Inquiry

Biological Activity

Description: Chrysin dimethylether shows moderate activity against Trypanosoma brucei, and the IC50s less than 5.0 microg/ml.

Targets: Antifection Chrysin dimethylether Description

Source: The herbs of Piper methysticum Forst.

Solvent: Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.

Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months(2-8°C).

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	dimethylether
[Catalog No.] CFN90896	
[CAS.No.] 21392-57-4	
[Parity] >+98%	
[M.E] C.H.O.	
[Physical Description] Powder	
[Solvent] Chloroform, Dichloromethane, E	Ethyl Acetate, DMSO, Acetone, etc.
[Weight] 10mg	
[Lot.No.] CFS202201	
[Storage] Protected from air and light, refi	rigerate or freeze(2-8 °C)
[Intended Use] For laboratory use only	
[Shelf Life] 2 years [Ghemical Structure]	
[Characterization Data Summary]	
Analytical Test	Results
Identification by "H-NMR, HPLC Purity tested	Consistent with the above structure >=98%
	219979



Kaempferia parviflora; KP



5,7,4'-trimethoxyflavone/ Trimethylapigenin

CAS No.	5631-70-9	Purity	>=98%
Catalog No.	CFN91890	Type of Compound	Flavonoids
Molecular Formula	$C_{18}H_{16}O_5$	Physical Desc.	Powder
Molecular Weight	312.32 g/mol	Source	The roots of Scutellaria indica L.
Solvent	Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.	Price	Inquiry

Biological Activity

Description: Trimethyapigenin has anti-inflammatory activity, it can moderately inhibit production of TNF- α ., iNOS mRNA and iNOS protein in a dose-dependent manner. It may be a potential candidate for anti-atrial fibrillation, it can significantly inhibit the atrial potassium currents hKv1.5/I(Kur) and I(KACh).

Targets: Potassium Channel | AChR | TNF-α | NO | NOS | ERK | JNK | IkB | MAPK | IKK Trimethylapigenin Description

Source: The roots of Scutellaria indica L.

Solvent: Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.

Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months(2-8°C).

Wherever possible, you should prepare and use solutions on the same day. However, if you need to make up stock solutions in advance, we recommend that you store the solution as aliquots in tightly sealed vials at -20°C. Generally, these will be useable for up to two weeks. Before use, and prior to opening the vial we recommend that you allow your product to equilibrate to room temperature for at least 1 hour.

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	E OF ANALYSIS hylapigenin
[Catalog.No.] CFN91890	
[CAS.No.] 5631-70-9	
[Purity]>=98%	
[M.E] C.H.O.	
[Physical Description] Powder	
[Solvent] Chloroform, Dichloromethane, Ett	tyl Acetate, DMSO, Acetone, etc.
[Weight] 10mg	
[Lot.No.] CFS202201	
[Storage] Protected from air and light, refrig	perate or freeze(2-8 °C)
[Intended Use] For laboratory use only	
[Shelf Life] 2 years [Chemical Structure]	
[Characterization Data Summary]	
Analytical Test	Results
Identification by "H-NMR, HPLC	Consistent with the above structure
Purity tested	>+98%
Authorized Signature:	



Kaempferia parviflora; KP



5,7,3',4'-Tetramethoxyflavone

CAS No.	855-97-0	Purity	>=98%
Catalog No.	CFN91116	Type of Compound	Flavonoids
Molecular Formula	C ₁₉ H ₁₈ O6	Physical Desc.	Powder
Molecular Weight	342.3 g/mol	Source	The herbs of Camellia sinensis
Solvent	Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.	Price	Inquiry

Biological Activity

Description: 5,7,3',4'-Tetramethoxyflavone (TMF) possesses various bioactivities, including antifungal, antimalarial, antimycobacterial, and anti-inflammatory activities; it also exhibits chondroprotective activity by targeting β -catenin signaling in vivo and in vitro. TMF protects chondrocytes from ER stress-induced apoptosis through regulation of the IRE1 α pathway.TMF inhibits the expression of tyrosinase, tyrosine-related protein (TRP)-1, and TRP-2 mRNA, which could be the mechanism of its melanogenesis inhibitory activity.

Targets: β-catenin | IRE1α | TRP-2 | TRP-1 | Antifection | TNF-α | PGE | PKA | cAMP | JNK | Caspase | Bcl/Bax | Antifection

Source: The herbs of Camellia sinensis

Solvent: Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.

Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months(2-8°C).

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CERTIFICAT	E OF ANALYSIS
5,7,3',4'-Tetr	amethoxyflavone
[Catalog No.] CFN01116	
[CAS.No.] 855-97-0	
[Parity] >=96%	
[ME]CJH.O.	
[Physical Description] Powder	
[Solvent] Chloroform, Dichloromethane, Et	hyl Acetale, DMISO, Acetone, etc.
[Weight] 50mg	
[Lot.No.] CF5202201	
[Storage] Protected from air and light, refri	perate or freeze(2-8 °C)
[Intended Use] For laboratory use only	
[ShelfLife] 2 years	~~
	and the
	- 44
[Chemical Structure]	0.0
[Characterization Data Summary]	
Analytical Test	Results
Identification by 'H-NMR, HPLC	Consistent with the above structure
Purity lesied	>+90%
Authorized Classifications	
Authorized Signature:	
Date:	



Curcuma longa L.

Curcumin

			Ref: M
CAS No.	458-37-7	Purity	>=98%
Catalog No.	CFN98686	Type of Compound	Phenols
Molecular Formula	$C_{21}H_{20}O_6$	Physical Desc.	Orange powder
Molecular Weight	368.4 g/mol	Source	The rhizomatous of Curcuma longa L.
Solvent	Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.	Price	Inquiry

Biological Activity

Description: Curcumin is a natural phenolic compound with diverse pharmacologic effects including antitumour, anti-bacteria, anti-fungicidal ,anti-edemic, hepatoprotective, anti-inflammatory, antioxidant, antiproliferative and antiangiogenic activities. Curcumin is an inhibitor of p300 histone acetylatransferase ((HATs)) and also shows inhibitory effects on NF- κ B and MAPKs.Curcumin application causes markedly fast wound closure with well-formed granulation tissue dominated by fibroblast proliferation, collagen deposition, and complete early regenerated epithelial layer.

Targets: DNA Methyltransferase | TNF- α | IFN- γ | IL Receptor | Nrf2 | HO-1 | GLUT | Caspase | STAT | NF-kB | HDAC | VEGFR | TGF- β /Smad

Source: The rhizomatous of Curcuma longa L.

Solvent: Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.

Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months(2-8°C).

Wherever possible, you should prepare and use solutions on the same day. However, if you need to make up stock solutions in advance, we recommend that you store the solution as aliquots in tightly sealed vials at -20°C. Generally, these will be useable for up to two weeks. Before use, and prior to opening the vial we recommend that you allow your product to equilibrate to room temperature for at least 1 hour.

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OLIVITION	TE OF ANALYSIS
0	Curcumin
[Catalog No.] CFN96686	
[CAS.No.] 458-37-7	
(Purity) >=98%	
[M.E.] C., H.O.	
(Physical Description) Orange powder	
[Solvent] Chloroform, Dichloromethane,	Ethyl Acetate, DMSO, Acetone, etc.
[Wisight] 20mg	
[Lot.No.] CFS202202	
(Storage) Protected from air and light, ref	rigerate or freeze(2-8 °C)
[Intended Use] For laboratory use only	
[Shelf Life] 2 years	princ:
[Chemical Structure]	
[Characterization Data Summary]	
(Characterization Data Summary)	Results Consistent with the above structure



Curcuma longa L.

Demethoxycurcumin

CAS No.	458-37-7	Purity	>=98%
Catalog No.	CFN98686	Type of Compound	Phenols
Molecular Formula	C ₂₁ H ₂₀ O ₆	Physical Desc.	Orange powder
Molecular Weight	368.4 g/mol	Source	The rhizomatous of Curcuma longa L.
Solvent	Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.	Price	Inquiry

Biological Activity

Description: Demethoxycurcumin is a potential additive natural product in combination with chemotherapeutic agents in drug-resistant cancers, which has anti-acanthamoebic, anti-proliferative, antimetastatic, anti-inflammatory, antioxidant activities. It inhibited P-glycoprotein-mediated ATP hydrolysis under concentrations of <1 μ M and efficiently inhibited 200 μ M verapamil-stimulated ATPase activity.

Targets: MMP(e.g.TIMP) | COX | TNF-α | NF-kB | EGFR | HSP (e.g. HSP90) | AMPK | ATPase | IL Receptor | P-gp

Source: The herbs of Curcuma longa L.

Solvent: Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.

Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months(2-8°C).

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CERTIFICATE OF ANALYSIS
Demethoxycurcumin
[Catalog No,] CFN00185
[CAS.No.] 22608-11-3
[Parity] >=90%
[ME]C_H_O,
[Physical Description] Yellow powder
[Solvent] Chioroform, Dichloromethane, Ethyl Acetale, DMSO, Acetone, etc.
[Waight] 100mg
[Lot.No.] CF5202202
[Storage] Protected from air and light, refrigerate or freeze(2-8 °C)
[Intended.Use] For laboratory use only
[ShellLife]2 years
но
[Chemical Structure]
[Characterization Data Summary]
Analytical Test Results
Identification by TrAMIR IPEC Consistent with the above structure
Purity tested >=(0)%
Authorized Signature:
Date:



Curcuma longa L.

Bisdemethoxycurcumin HO

CAS No.	33171-05-0	Purity	>=98%
Catalog No.	CFN99186	Type of Compound	Phenols
Molecular Formula	C ₁₉ H ₁₆ O ₄	Physical Desc.	Yellow powder
Molecular Weight	308.33 g/mol	Source	The rhizomatous of Curcuma longa L.
Solvent	Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.	Price	Inquiry

Biological Activity

Description: Bisdemethoxycurcumin is a natural derivative of curcumin with antiulcer, antioxidant, anti-inflammatory and anti-cancer activities, it suppresses MCF-7 cells proliferation by inducing ROS accumulation and modulating senescence-related pathways. Bisdemethoxycurcumin induces apoptosis in activated HSCs, but not in hepatocytes, by impairing cellular energetics and causing a downregulation of cytoprotective proteins, likely through a mechanism that involves CBR2.

Targets: NOS | COX | NF-kB | ROS | MMP(e.g.TIMP) | VEGFR | TNF-α | Akt | ERK | Bcl-2/Bax | p53 | p21

Source: The rhizomes of Curcuma longa L.

Solvent: Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.

Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months(2-8°C).

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	E OF ANALYSIS
[Catalog.No,] CFN99186	
[CAS.No.] 33171-05-0	
[Purity] >=06%	
[ME]CJHO.	
[Physical Description] Yellow powder	
[Solvent] Chloroform, Dichloromethane, Eth	yl Acetale, DMSO, Acetone, etc.
[Weight] 20mg	
[Lot No.] CFS202201	
[Storage] Protected from air and light, refrig	erate or freeze(2-8 °C)
[Intended Use] For laboratory use only	
[Shelf Life] 2 years	orita
[Chemical Structure]	
[Characterization Data Summary]	
Analytical Test	Results
Identification by "H-NMR, HPLC	Consistent with the above structure
Purity tested	>×98%
Authorized Signature:	



10/13 Krungthepkritha 7, Huamak, Bangkapi,Bangkok 10240
 02–3469239
 Sales_chem@axxo.co.th
 www.axxo.co.th

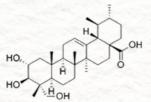
Ref: M-Group: Article



Plant library of Curcuma longa L.

Catalog	Product Name	CAS Number	Manual
CFN98686	Curcumin	458-37-7	PDF
CFN99186	Bisdemethoxycurcumin	33171-05-0	PDF
CFN99540	Dihydrocurcumin	76474-56-1	PDF
CFN99187	Curcumol	4871-97-0	PDF
CFN98133	Germacrone	6902-91-6	PDF
CFN99582	Linalool	78-70-6	PDF
CFN97326	Stigmasterol	83-48-7	PDF
CFN99916	Beta-Sitosterol	83-46-5	PDF
CFN90583	Tetrahydrocurcumin	36062-04-1	PDF
CFN97749	Hexahydrocurcumin	36062-05-2	PDF
CFN90584	Octahydrocurcumin	36062-07-4	PDF
CFN96093	1,5-Bis(4-hydroxy-3-methoxyphenyl)penta-1,4-diene	63644-68-8	PDF
CFN92778	Coronadiene	1145689-64-0	PDF
CFN92779	Zerumin A	176050-48-9	PDF
CFN99185	Demethoxycurcumin	22608-11-3	PDF
CFN97433	5-Hydroxy-1,7-diphenyl-6-hepten-3-one	87095-74-7	PDF
CFN89015	Bisacurone C	127214-86-2	n/a
CFN89231	ar-Turmerone	532-65-0	n/a
CFN96432	Bisacurone	120681-81-4	n/a
CFN96210	1,7-Bis(4-hydroxyphenyl)-3-hydroxy-1,3-heptadien-5-one	207792-17-4	n/a
CFN96744	5"-Methoxyhexahydrocurcumin	138870-96-9	n/a
CFN89244	4-(6-Methyl-4-oxohept-5-en-2-yl)cyclohex-2-en-1-one	170380-68-4	n/a
CFN96841	1-(3,4-Dihydroxyphenyl)-7-(4-hydroxyphenyl)hept-6-en-3-ol	1206788-61-5	n/a









CAS No.	464-92-6	Purity	>==98%
Catalog No.	CFN98688	Type of Compound	Triterpenoids
Molecular Formula	C ₃₀ H ₄₈ O ₅	Physical Desc.	Powder
Molecular Weight	488.7 g/mol	Source	The herbs of Centella asiatica.
Solvent	Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.	Price	Inquiry

Biological Activity

Description: Asiatic acid shows antihyperlipidemic, anti-inflammatory, antioxidant, and antitumorigenesis effects, it inhibits NLRP3 inflammasome activation, NO and COX-2 signals. Asiatic acid inhibits the expression NDR1/2 kinase and promotes the stability of p21WAF1/CIP1 protein through attenuating NDR1/2 dependent phosphorylation of p21WAF1/CIP1 in HepG2 cells.

Targets: TNF-α | IFN-γ | IL Receptor | Caspase | p21 | NOS | NADPH-oxidase | NO | HMG-CoA reductase | COX

Source: The herbs of Centella asiatica.

Solvent: Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.

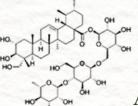
Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months(2-8°C).

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	E OF ANALYSIS atic acid
[Catalog No.] CFN08688	
[CAS.No.] 464-92-6	
[Parity] 99.8%	
[M.E] C.H.O.	
[Physical Description] Powder	
[Solvent] Chloroform, Dichloromethane, Eth	tyl Acetate, DMSO, Acetone, etc.
[Weight] 20mg	
[Lot No.] CFS202201	
[Storage] Protected from air and light, refrig	perate or freeze(2-8 °C)
[Intended.Use] For laboratory use only	1
[Shelf Life] 2 years	
(Chamical Structure)	NO HALL OH
Analytical Test	Results
Identification by "H-NMR, HPLC	Consistent with the above structure
Purity tested	99.8%
Authorized Signature: Date:	





Asiaticoside



CAS No.	16830-15-2	Purity	>=98%
Catalog No.	CFN99912	Type of Compound	Triterpenoids
Molecular Formula	C ₄₈ H ₇₈ O ₁₉	Physical Desc.	White powder
Molecular Weight	959.12 g/mol	Source	The herbs of Centella asiatica (L.) Urban
Solvent	Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.	Price	Inquiry

Biological Activity

Description: Asiaticoside, a biochemical modulator, which has antioxidant, anti-inflammatory, antipyretic, anxiolytic-like, anti-gastric ulcers, hepatoprotective, and antidepressant-like effects, it also exhibits significant wound healing activity in normal as well as delayed healing models. Asiaticoside suppressed collagen expression and TGF- β /Smad signaling through inducing Smad7 and inhibiting TGF- β RI and TGF- β RII in keloid fibroblasts. It and its derivatives can be regarded as reasonable candidates for a therapeutic Alzheimer's disease drug that protects neurons from Abeta toxicity.

Targets: NF-kB | TNF-α | IL Receptor | p65 | IkB | TGF-β/Smad | p38MAPK | JNK | ERK | Caspase | COX | PGE | PPAR | NOS | Bcl-2/Bax | Beta Amyloid | IKK

Source: The herbs of Centella asiatica (L.) Urban

Solvent: DMSO, Pyridine, Methanol, Ethanol, etc.

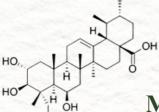
Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months(2-8°C).

Wherever possible, you should prepare and use solutions on the same day. However, if you need to make up stock solutions in advance, we recommend that you store the solution as aliquots in tightly sealed vials at -20°C. Generally, these will be useable for up to two weeks. Before use, and prior to opening the vial we recommend that you allow your product to equilibrate to room temperature for at least 1 hour.

After receiving: The packaging of the product may have turned upside down during transportation, resulting in the natural compounds adhering to the neck or cap of the vial. take the vial out of its packaging and gently shake to let the compounds fall to the bottom of the vial. for liquid products, centrifuge at 200-500 RPM to gather the liquid at the bottom of the vial. try to avoid loss or contamination during handling.

CERTIFICATE	OF ANALYSIS
Asiati	coside
[Catalog No.] OFN88812	
[CAS.No.] 16830-15-2	
(Purity) 99.0%	
IMEICHO.	
[Physical Description] White powder	
[Solvent] DMSO, Pyridine, Methanol, Ethanol	, etc.
[Weight] 20mg	
[Lot.No.] CF5202202	
[Storage] Protected from air and light, refriger	ale or freeze(2-8 °C)
[Intended Use] For laboratory use only	
[ShelfLife]2years	and a star
×0.,	444-7 T
20	No march 10
(Chemical Structure)	Japan Mar
	der Der
[Characterization Data Summary]	
Analytical Test	Results
	Consistent with the above structure
Purity tested	99.0%
Authorized Signature:	
Date:	







Madecassic acid

CAS No.	18449-41-7	Purity	>=98%
Catalog No.	CFN99914	Type of Compound	Triterpenoids
Molecular Formula	C ₃₀ H ₄₈ O ₆	Physical Desc.	Powder
Molecular Weight	504.70 g/mol	Source	The herbs of Centella asiatica (L.) Urban
Solvent	Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.	Price	Inquiry

Biological Activity

Description: Madecassic acid has anti-diabetic, anti- tumor, wound-healing, and anti-inflammatory properties, it can improve glycemic control and hemostatic imbalance, lower lipid accumulation, and attenuate oxidative and inflammatory stress in diabetic mice. It can protect against hypoxia-induced oxidative stress in retinal microvascular endothelial cells via ROS-mediated endoplasmic reticulum stress. It inhibited the esspession of NOS, COX-2, TNF-alpha, IL-1beta, IL-6, and the downregulation of NF-kappaB activation.

Targets: ROS | NOS | COX | TNF-α | IL Receptor | NF-kB | Caspase | Bcl-2/Bax | p65 | NO | PGE | gp120/CD4 | IkB | IKK

Source: The herbs of Centella asiatica (L.) Urban

Solvent: Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.

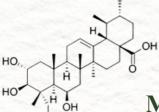
Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months(2-8°C).

Wherever possible, you should prepare and use solutions on the same day. However, if you need to make up stock solutions in advance, we recommend that you store the solution as aliquots in tightly sealed vials at -20°C. Generally, these will be useable for up to two weeks. Before use, and prior to opening the vial we recommend that you allow your product to equilibrate to room temperature for at least 1 hour.

After receiving: The packaging of the product may have turned upside down during transportation, resulting in the natural compounds adhering to the neck or cap of the vial. take the vial out of its packaging and gently shake to let the compounds fall to the bottom of the vial. for liquid products, centrifuge at 200-500 RPM to gather the liquid at the bottom of the vial. try to avoid loss or contamination during handling.

	E OF ANALYSIS assic acid
[Catalog.No.] CFN99914	
[CAS.No.] 18449-41-7	
[Parity] 96.5%	
[MLE]CUHUO,	
[Physical Description] Powder	
[Solvent] Chioroform, Dichioromethane, Eth	yl Acetate, DMSO, Acetone, etc.
[Waight] 20mg	
[Lot.No.] CF5202202	
[Storage] Protected from air and light, refrig	erate-or heeze(2-8 °C)
[Intended Use] For laboratory use only	
(Sheff Life) 2 years (Chemical Bruchum)	A Carlor
[Characterization Data Summary]	
Analytical Test Identification by Tin ABBL, HPLC Purity Instee	Results Consistent with the above structure 98.5%
Authorized Signature: Date:	







Madecassic acid

CAS No.	18449-41-7	Purity	>=98%
Catalog No.	CFN99914	Type of Compound	Triterpenoids
Molecular Formula	C ₃₀ H ₄₈ O ₆	Physical Desc.	Powder
Molecular Weight	504.70 g/mol	Source	The herbs of Centella asiatica (L.) Urban
Solvent	Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.	Price	Inquiry

Biological Activity

Description: Madecassic acid has anti-diabetic, anti- tumor, wound-healing, and anti-inflammatory properties, it can improve glycemic control and hemostatic imbalance, lower lipid accumulation, and attenuate oxidative and inflammatory stress in diabetic mice. It can protect against hypoxia-induced oxidative stress in retinal microvascular endothelial cells via ROS-mediated endoplasmic reticulum stress. It inhibited the esspession of NOS, COX-2, TNF-alpha, IL-1beta, IL-6, and the downregulation of NF-kappaB activation.

Targets: ROS | NOS | COX | TNF-α | IL Receptor | NF-kB | Caspase | Bcl-2/Bax | p65 | NO | PGE | gp120/CD4 | IkB | IKK

Source: The herbs of Centella asiatica (L.) Urban

Solvent: Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.

Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months(2-8°C).

Wherever possible, you should prepare and use solutions on the same day. However, if you need to make up stock solutions in advance, we recommend that you store the solution as aliquots in tightly sealed vials at -20°C. Generally, these will be useable for up to two weeks. Before use, and prior to opening the vial we recommend that you allow your product to equilibrate to room temperature for at least 1 hour.

After receiving: The packaging of the product may have turned upside down during transportation, resulting in the natural compounds adhering to the neck or cap of the vial. take the vial out of its packaging and gently shake to let the compounds fall to the bottom of the vial. for liquid products, centrifuge at 200-500 RPM to gather the liquid at the bottom of the vial. try to avoid loss or contamination during handling.

	E OF ANALYSIS assic acid
[Catalog.No.] CFN99914	
[CAS.No.] 18449-41-7	
[Parity] 96.5%	
[MLE]CUHUO,	
[Physical Description] Powder	
[Solvent] Chioroform, Dichioromethane, Eth	yl Acetate, DMSO, Acetone, etc.
[Waight] 20mg	
[Lot.No.] CF5202202	
[Storage] Protected from air and light, refrig	erate-or heeze(2-8 °C)
[Intended Use] For laboratory use only	
(Sheff Life) 2 years (Chemical Bruchum)	A Carlor
[Characterization Data Summary]	
Analytical Test Identification by Tin ABBL, HPLC Purity Instee	Results Consistent with the above structure 98.5%
Authorized Signature: Date:	





Plant library of Centella asiatica L.

Cat. No.	Information		
CFN97743	41682-30-8	8-Acetoxypentadeca-1,9Z-diene-4,6-diyn-3-ol	
CFN98688	464-92-6	Asiatic acid	
CFN99912	16830-15-2	Asiaticoside	
CFN95124	125265-68-1	Asiaticoside B	
CFN99914	18449-41-7	Madecassic acid	
CFN99913	34540-22-2	Madecassoside	



Phyllanthus emblica L.



2-2'-(Hydroxytetracosanoylamino)-octadecane-1,3,4-triol tetraacetate

CAS No.	340702-68-3	Purity	>=98%
Catalog No.	CFN98439	Type of Compound	Cerebrosides
Molecular Formula	C ₅₀ H ₉₃ NO ₉	Physical Desc.	Powder
Molecular Weight	852.3 g/mol	Source	The herbs of Phyllanthus emblica
Solvent	Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.	Price	Inquiry

Biological Activity

Description: (2S,3S,4R,2'R))-2-(2'-Hydroxytetracosanoylamino) octadecane-1,3,4-triol at concentration of 100 ug/mL shows selectively inhibitory activity against phospholipase A 2 (PLA 2) secreted from Crotalus adamenteus venom, but inactive against PLA 2 of bee venom (Apis mellifera).

Source: The herbs of Phyllanthus emblica

Solvent: DMSO, Pyridine, Methanol, Ethanol, etc.

Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months(2-8°C).

Wherever possible, you should prepare and use solutions on the same day. However, if you need to make up stock solutions in advance, we recommend that you store the solution as aliquots in tightly sealed vials at -20°C. Generally, these will be useable for up to two weeks. Before use, and prior to opening the vial we recommend that you allow your product to equilibrate to room temperature for at least 1 hour.

After receiving: The packaging of the product may have turned upside down during transportation, resulting in the natural compounds adhering to the neck or cap of the vial. take the vial out of its packaging and gently shake to let the compounds fall to the bottom of the vial. for liquid products, centrifuge at 200-500 RPM to gather the liquid at the bottom of the vial. try to avoid loss or contamination during handling.

Chem Faces 😤	ISO 9001:200
CERTIFIC	ATE OF ANALYSIS
2-2'-(Hydroxytetracosanoylar	nino)-octadecane-1,3,4-triol tetraacet
[Catalog No.] CFN98439	
[CAS.No.] 340702-68-3	
[Parity] >=98%	
[ME]C_H_NO,	
(Physical Description)	
[Solvent] Pyridine, DMSO.	
[Weight] Smg	
[Lot.No.] CFS201701	
[Storage] Protected from air and light,	refrigerate or freeze(2-8 °C)
[Intended Use] For laboratory use only	
[ShelfLife] 2 years	A.C.
[Chemical Structure]	÷
[Characterization Data Summary]	
Analytical Test	Results
Identification by NMR	Consistent with the above structure
Purity tested	>+98%



Phyllanthus emblica L.



OH

3,4-Dimethoxybenzoic acid

CAS No.	93-07-2	Purity	>=98%
Catalog No.	CFN97499	Type of Compound	Phenols
Molecular Formula	C ₉ H ₁₀ O ₄	Physical Desc.	Powder
Molecular Weight	182.2 g/mol	Source	The fruits of Phyllanthus emblica L.
Solvent	Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.	Price	Inquiry

Biological Activity

Description: 3,4-Dimethoxybenzoic acid(Veratric acid), a simple benzoic acid derived from plants and fruits, has anti-oxidant, anti-inflammation, and blood pressure-lowering effects.

Source: The fruits of Phyllanthus emblica L.

Solvent: Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.

Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months(2-8°C).

Wherever possible, you should prepare and use solutions on the same day. However, if you need to make up stock solutions in advance, we recommend that you store the solution as aliquots in tightly sealed vials at -20°C. Generally, these will be useable for up to two weeks. Before use, and prior to opening the vial we recommend that you allow your product to equilibrate to room temperature for at least 1 hour.

After receiving: The packaging of the product may have turned upside down during transportation, resulting in the natural compounds adhering to the neck or cap of the vial. take the vial out of its packaging and gently shake to let the compounds fall to the bottom of the vial. for liquid products, centrifuge at 200-500 RPM to gather the liquid at the bottom of the vial. try to avoid loss or contamination during handling.

		E OF ANALYSIS
	Contraction No. 1 OFMOTION	
	[Catalog No.] CFN97499	
	[CAS.No.] 93-07-2	
	[Purity] >=98%	
	[M.E] C,H,,O,	
	[Physical Description] Powder	
	[Solvent] Chloroform, Dichloromethane, Eth	yl Acetate, DMSO, Acetone, etc.
	[Weight] 20mg	
	[Lot.No.] CFS202101	
	[Storage] Protected from air and light, refrige	erate or freeze(2-8 °C)
	[Intended Use] For laboratory use only	
	[Shelf Life] 2 years	OOH
	[Chemical Structure]	0
	[Characterization Data Summary]	
	Analytical Test	Results
	Identification by 'H-NMR, HPLC	Consistent with the above structure
100	Purity tested	>=98%

Authorized Signature: Date:



Phyllanthus emblica L.



Glochicoccin D

CAS No.	927812-23-5	Purity	>=98%
Catalog No.	CFN97495	Type of Compound	Miscellaneous
Molecular Formula	C ₂₁ H ₂₄ O ₁₀	Physical Desc.	Powder
Molecular Weight	436.4 g/mol	Source	The fruits of Phyllanthus emblica L.
Solvent	Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.	Price	Inquiry

Biological Activity

Description: Standard reference

Source: The fruits of Phyllanthus emblica L.

Solvent: Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.

Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months (2-8°C).

Wherever possible, you should prepare and use solutions on the same day. However, if you need to make up stock solutions in advance, we recommend that you store the solution as aliquots in tightly sealed vials at -20°C. Generally, these will be useable for up to two weeks. Before use, and prior to opening the vial we recommend that you allow your product to equilibrate to room temperature for at least 1 hour.

After receiving: The packaging of the product may have turned upside down during transportation, resulting in the natural compounds adhering to the neck or cap of the vial. take the vial out of its packaging and gently shake to let the compounds fall to the bottom of the vial. for liquid products, centrifuge at 200-500 RPM to gather the liquid at the bottom of the vial. try to avoid loss or contamination during handling.

CERTIFICA	TE OF ANALYSIS
Glo	ochicoccin D
[Catalog No.] CFN97495	
CAS No.] 927812-23-5	
[Purity]>=98%	
[M.E] C, H, O.	
[Physical Description]	
[Solvent] Chloroform, Dichloromethane,	Ethyl Acetate, DMSO, Acetone, etc.
[Weight] 5mg	
[Lot No.] CFS201701	
[Storage] Protected from air and light, re	frigerate or freeze(2-8 °C)
[Intended Use] For laboratory use only	OH
[Shelf Life] 2 years	2000 - C
[Chemical Structure]	HOLOO
[Characterization Data Summary]	
Analytical Test	Results
Identification by NMR	Consistent with the above structure



Plant library of *Phyllanthus emblica* L.



Cat. No.	Information	
CFN98439	340702-68-3	2-2'-(Hydroxytetracosanoylamino)-octadecane-1,3,4-triol tetraacetate
CFN97499	93-07-2	3,4-Dimethoxybenzoic acid
CFN97405	84812-00-0	3,4-Dimethoxyphenyl beta-D-glucoside
	927812-23-5	Glochicoccin D
CFN95201	1236310-34-1	Methyl neochebulinate





(2R)-5,7-Dimethoxyflavanone

CAS No.	1277188-85-8	Purity	>=98%
Catalog No.	CFN90899	Type of Compound	Flavonoids
Molecular Formula	C ₁₇ H ₁₆ O ₄	Physical Desc.	Powder
Molecular Weight	284.31 g/mol	Source	The black rhizomes of Boesenbergia panduta.
Solvent	Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.	Price	Inquiry

Biological Activity

Description: 5,7-Dimethoxyflavanone has anti-inflammatory activity.

Targets: TNF-a | IL Receptor | NO | IFN-y

Source: The black rhizomes of Boesenbergia panduta.

Solvent: Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.

Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months(2-8C).

Wherever possible, you should prepare and use solutions on the same day. However, if you need to make up stock solutions in advance, we recommend that you store the solution as aliquots in tightly sealed vials at -20C. Generally, these will be useable for up to two weeks. Before use, and prior to opening the vial we recommend that you allow your product to equilibrate to room temperature for at least 1 hour.

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CERTIFIC	ATE OF ANALYSIS
(2R)-5,7-0	Dimethoxyflavanone
[Catalog No.] CFN90899	
[CAS.No.] 1277188-85-8	
[Purity] >=98%	
[M.E.] C.,H.,O,	
[Physical Description]	
[Solvent] Chioroform, Dichioromethane	, Ethyl Acetate, DMSO, Acetone, etc.
[Weight] 5mg	
[Lot.No.] CFS201701	
[Storage.] Protected from air and light, r	efrigerate or freeze(2-8 °C)
[Intended Use] For laboratory use only	
[Shelf Life] 2 years	$\sim 10^{\circ}$
[Chemical Structure]	11
[Characterization Data Summary]	
Analytical Test	Results
Identification by NMR	Consistent with the above structure
Purity tested	>=98%
Authorized Signature:	



Boesenbergia pandura



Longiferone B

CAS No.	1639810-67-5	Purity	>=98%
Catalog No.	CFN89084	Type of Compound	Sesquiterpenoids
Molecular Formula	C15H22O	Physical Desc.	Oil
Molecular Weight	218.34 g/mol	Source	The rhizomes of Boesenbergia longiflora (Wall.) Kuntze.
Solvent	Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.	Price	Inquiry

Biological Activity

Description: Longiferone B shows anti-inflammatory activity against NO release with IC50 values of 21.0 uM, it also suppresses the iNOS and COX-2 mRNA expression.

Targets: NO | NOS | COX

Source: The rhizomes of Boesenbergia longiflora (Wall.) Kuntze.

Solvent: Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.

Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months (2-8°C).

Wherever possible, you should prepare and use solutions on the same day. However, if you need to make up stock solutions in advance, we recommend that you store the solution as aliquots in tightly sealed vials at -20°C. Generally, these will be useable for up to two weeks. Before use, and prior to opening the vial we recommend that you allow your product to equilibrate to room temperature for at least 1 hour.

After receiving: The packaging of the product may have turned upside down during transportation, resulting in the natural compounds adhering to the neck or cap of the vial. take the vial out of its packaging and gently shake to let the compounds fall to the bottom of the vial. for liquid products, centrifuge at 200-500 RPM to gather the liquid at the bottom of the vial. try to avoid loss or contamination during handling.

Chem Faces 🛠	ISO 9001:2015
	ATE OF ANALYSIS
[Catalog No.] CFN89084	
[CAS.No.] 1639810-67-5	
[Purity]>=90%	
[ME]C.H.O	
[Physical Description] OI	
[Solvent] Chioroform, Dichloromethane	Ethyl Acetate DAISO Acetana atc
[Weight] 10mg	, con normale, preson, normality, esc.
[Lot.No.] CFS202002	
	tioned as for some \$ \$ 100
[Storage] Protected from air and light,	
[Intended Use] For laboratory use only [Shelf Life] 2 years	$d \rightarrow$
[Chemical Structure]	H O
[Characterization Data Summary]	
Analytical Test	Results
Identification by "H-NMR	Consistent with the above structure
Purity tested	>+98%

Date



Boesenbergia pandura



Pinocembrin chalcone

CAS No.	4197-97-1	Purity	>=98%
Catalog No.	CFN96919	Type of Compound	Chalcones
Molecular Formula	C ₁₅ H ₁₂ O ₄	Physical Desc.	Powder
Molecular Weight	256.25 g/mol	Source	The rhizomes of Boesenbergia pandurata.
Solvent	Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.	Price	Inquiry

Biological Activity

Description: Pinocembrin chalcone has tyrosinase inhibitory activity. It also shows antimutagenic effect, which is mainly due to the inhibition of the first step of enzymatic activation of heterocyclic amines. Pinocembrin chalcone shows antimicrobial activity against the antibiotic susceptible NG strain WHO V ; it also displays activity against Candida albicans with a minimal inhibitory concentration value of 100 microg/mL.

Targets: Tyrosinase | Calcium Channel | Antifection

Source: The rhizomes of Boesenbergia pandurata.

Solvent: Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.

Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months(2-8°C).

Wherever possible, you should prepare and use solutions on the same day. However, if you need to make up stock solutions in advance, we recommend that you store the solution as aliquots in tightly sealed vials at -20°C. Generally, these will be useable for up to two weeks. Before use, and prior to opening the vial we recommend that you allow your product to equilibrate to room temperature for at least 1 hour.

After receiving: The packaging of the product may have turned upside down during transportation, resulting in the natural compounds adhering to the neck or cap of the vial. take the vial out of its packaging and gently shake to let the compounds fall to the bottom of the vial. for liquid products, centrifuge at 200-500 RPM to gather the liquid at the bottom of the vial. try to avoid loss or contamination during handling.

	E OF ANALYSIS
[Catalog.No.] CFN96919	
[CAS.No.] 4197-97-1	
[Parity]>=98%	
[M.E] C.,H.,O.	
[Physical Description] Powder	
[Solvent] Chloroform, Dichloromethane, Et	hyl Acetate, DMSO, Acetone, etc.
[Weight] Smg	
[Lot.No.] CF\$202002	
(Storage) Protected from air and light, refri	gerate or freeze(2-8 °C)
[Intended Use] For laboratory use only	
[Shelf Life] 2 years	но он
[Chemical Structure]	OH O
[Characterization Data Summary]	
Analytical Test	Results
Identification by "H-MMR, HPLC	Consistent with the above structure
Purity tested	>=98%
Authorized Signature:	



Andrographis paniculata



Andrographolide

CAS No.	5508-58-7	Purity	>=98%
Catalog No.	CFN98923	Type of Compound	Diterpenoids
Molecular Formula	C ₂₀ H ₃₀ O ₅	Physical Desc.	Powder
Molecular Weight	350.5 g/mol	Source	The herbs of Andrographis paniculata (Burm. f.) Nees
Solvent	Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.	Price	Inquiry

Biological Activity

Description: Andrographolide is an antiinflammatory, antiviral, anti-cancer, hepatoprotective, antithrombotic, hypotensive and antiatherosclerotic drug, it can cure hyperpigmentation disorders. Andrographolide protects against chemical-induced oxidative damage by up-regulating the gene transcription and activity of antioxidant enzymes in various tissues. Andrographolide has potential as a leading compound in the prevention or treatment of obesity and insulin resistance, can ameliorate lipid metabolism and improve glucose use in mice with HFD-induced obesity.

Targets: Wnt/β-catenin | PKC | GSK-3 | Akt | PI3K | TGF-β/Smad | HO-1 | Nrf2 | ERK | JNK | gp120/CD4 | HIV | MMP(e.g.TIMP) | PAFR | AP-1 | ROS | Calcium Channel

Source: The herbs of Andrographis paniculata (Burm. f.) Nees

Solvent: Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.

Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months (2-8°C).

Wherever possible, you should prepare and use solutions on the same day. However, if you need to make up stock solutions in advance, we recommend that you store the solution as aliquots in tightly sealed vials at -20°C. Generally, these will be useable for up to two weeks. Before use, and prior to opening the vial we recommend that you allow your product to equilibrate to room temperature for at least 1 hour.

After receiving: The packaging of the product may have turned upside down during transportation, resulting in the natural compounds adhering to the neck or cap of the vial. take the vial out of its packaging and gently shake to let the compounds fall to the bottom of the vial. for liquid products, centrifuge at 200-500 RPM to gather the liquid at the bottom of the vial. try to avoid loss or contamination during handling.

CERTIFICATI	E OF ANALYSIS
Androg	grapholide
[Catalog No.] CFN98923	
[CAS.No.] 5508-58-7	
[Parity] >=98%	
[ME]C,H,O,	
[Physical Description] Powder	
[Solvent] Chioroform, Dichloromethane, Eth	yl Acetate, DMSO, Acetone, etc.
[Weight] 20mg	
[Lot No.] CFS202201	
[Storage] Protected from air and light, refrig	erate or freeze(2-8 °C)
[Intended Use] For laboratory use only	o
[Shelf Life] 2 years	HO
[Chemical Structure]	HO MOH
Analytical Test Identification by "H-NMR, HPLC	Results Consistent with the above structure
Purity tested	>+98%
Authorized Signature: Date:	



Andrographis paniculata



Dehydroandrographolide

CAS No.	134418-28-3	Purity	>=98%
Catalog No.	CFN98923	Type of Compound	Diterpenoids
Molecular Formula	$C_{20}H_{28}O_4$	Physical Desc.	White powder
Molecular Weight	332.43 g/mol	Source	The herbs of Andrographis paniculata (Burm. f.) Nees
Solvent	Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.	Price	Inquiry

Biological Activity

Description: Dehydroandrographolide is a novel TMEM16A inhibitor and possesses multiple pharmacological activities, including anti-inflammation, anti-cancer, anti-bacterial, anti-virus and anti-hepatitis activity. It possesses activity against HBV DNA replication with IC50 values of 22.58 uM and low SI values of 8.7 ; it can alleviate oxidative stress in LPS-induced acute lung injury possibly by inactivating iNOS.

Targets: IL Receptor | NOS | TNF-a | p38MAPK | HBV

Source: The herbs of Andrographis paniculata (Burm. f.) Nees

Solvent: Chloroform, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.

Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months(2-8°C).

Wherever possible, you should prepare and use solutions on the same day. However, if you need to make up stock solutions in advance, we recommend that you store the solution as aliquots in tightly sealed vials at -20°C. Generally, these will be useable for up to two weeks. Before use, and prior to opening the vial we recommend that you allow your product to equilibrate to room temperature for at least 1 hour.

After receiving: The packaging of the product may have turned upside down during transportation, resulting in the natural compounds adhering to the neck or cap of the vial. take the vial out of its packaging and gently shake to let the compounds fall to the bottom of the vial. for liquid products, centrifuge at 200-500 RPM to gather the liquid at the bottom of the vial. try to avoid loss or contamination during handling.

CERTIFICATE OF ANALYSIS
Dehydroandrographolide
[Catalog No.] C/N00770
[CAS.No.] 134418-28-3
[Parity] >=90%
(ME)CHO.
[Physical Description] White powder
[Solvest] Chicrotom, Dichloromethane, Ethyl Acetate, DMSO, Acetone, etc.
[Waight] 20mg
[Lot.No.] CF8202102
[Storage] Protected from air and light, refrigerate or freeze(2-8 °C)
[Intended Use] For laboratory use only
[ShellLife]2 years
(Chemical Bootlans) HO HO HO
Analytical Test Results
Identification by "H-MBR, HPLC Consistent with the above structure Purey tested >+0()%
Authorized Signature:
Date:



Stevia rebaudiana

Stevioside

CAS No.	57817-89-7	Purity	>=98%
Catalog No.	CFN99548	Type of Compound	Diterpenoids
Molecular Formula	C ₃₈ H ₆₀ O ₁₈	Physical Desc.	White powder
Molecular Weight	804.88 g/mol	Source	The leaves of Stevia rebaudiana
Solvent	DMSO, Pyridine, Methanol, Ethanol, etc.	Price	Inquiry

Biological Activity

Description: Stevioside is a safe natural sweetener, has no allergic reactions, suited for both diabetics, and PKU patients, as well as for obese persons intending to lose weight by avoiding sugar supplements in the diet. Stevioside enjoys a dual positive effect by acting as an antihyperglycemic and a blood pressure-lowering substance, it may have therapeutic potential in the treatment of type 2 diabetes and the metabolic syndrome.Stevioside exerts anti-inflammatory and anti-apoptotic properties by inhibiting the release of cytokines and the activation of TLR2 and proteins of the NF-KB and MAPK signaling pathways, as well as caspase-3 and Bax.

Targets: TLR | NF-kB | TNF-α | IL Receptor | IkB | p38MAPK | ERK | JNK | p65 | Caspase | PDE | IKK

Source: The leaves of Stevia rebaudiana

Solvent: DMSO, Pyridine, Methanol, Ethanol, etc.

Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months(2-8°C).

Wherever possible, you should prepare and use solutions on the same day. However, if you need to make up stock solutions in advance, we recommend that you store the solution as aliquots in tightly sealed vials at -20°C. Generally, these will be useable for up to two weeks. Before use, and prior to opening the vial we recommend that you allow your product to equilibrate to room temperature for at least 1 hour.

After receiving: The packaging of the product may have turned upside down during transportation, resulting in the natural compounds adhering to the neck or cap of the vial. take the vial out of its packaging and gently shake to let the compounds fall to the bottom of the vial. for liquid products, centrifuge at 200-500 RPM to gather the liquid at the bottom of the vial. try to avoid loss or contamination during handling.

CERTIFICATE C Stevio	
[Catalog No.] CFN99548	
[CAS.No.] 57817-89-7	
[Pully]>=90%	
[MLE]CJUO,	
[Physical Description] While powder	
[Solvent] DMSO, Pyridine, Methanol, Ethanol, et	к.
[Weight] 20mg*3	
[Lot.No.] CF5202201	
[Storage] Protected from air and light, refrigerate	or theeze(2-8 °C)
[Inlanded Use] For laboratory use only	-D
[Shelf Life] 2 years	-03-0-
(Chemical Structure)	Sight a
[Characterization Data Summary]	
Analytical Test Identification by "In-AMIT, HPLC C) Purity Instand	Results relatert with the above structure >+06%
Authorized Signature: Date:	



Stevia rebaudiana



Rebaudioside A

CAS No.	58543-16-1	Purity	>=98%
Catalog No.	CFN99110	Type of Compound	Diterpenoids
Molecular Formula	C44H70O23	Physical Desc.	Powder
Molecular Weight	967.01 g/mol	Source	The leaves of Stevia rebaudiana Bertoni.
Solvent	DMSO, Pyridine, Methanol, Ethanol, etc.	Price	Inquiry

Biological Activity

Description: Rebaudioside A, an natural sweetening ingredient, is approximately 250 to 300 times sweeter than sucrose, do not contribute calories or carbohydrates to the diet and do not affect blood glucose or insulin response. It is a α -glucosidase inhibitor with IC50 of 35.01 µg/ml. **Targets:** α -glucosidase

Source: The leaves of Stevia rebaudiana Bertoni.

Solvent: DMSO, Pyridine, Methanol, Ethanol, etc.

Storage: Providing storage is as stated on the product vial and the vial is kept tightly sealed, the product can be stored for up to 24 months(2-8°C).

Wherever possible, you should prepare and use solutions on the same day. However, if you need to make up stock solutions in advance, we recommend that you store the solution as aliquots in tightly sealed vials at -20°C. Generally, these will be useable for up to two weeks. Before use, and prior to opening the vial we recommend that you allow your product to equilibrate to room temperature for at least 1 hour.

After receiving: The packaging of the product may have turned upside down during transportation, resulting in the natural compounds adhering to the neck or cap of the vial. take the vial out of its packaging and gently shake to let the compounds fall to the bottom of the vial. for liquid products, centrifuge at 200-500 RPM to gather the liquid at the bottom of the vial. try to avoid loss or contamination during handling.

CERTIFICA	TE OF ANALYSIS
Reb	audioside A
[Catalog No.] CFN90110	
[CAS No.] 58543-16-1	
[Parity]>+98%	
ME C.H.O.	
[Physical Description] Powder	
[Solvent] DMSO, Pyridine, Methanol, Eth	hanol, etc.
[Weight] 20mg	
[Lot.No.] CF5202201	
(Storage) Protected from air and light, re	frigerate or freeze(2-8 °C)
[Intended Use] For laboratory use only	94
[Shelf Life] 2 years	100° 20°
[Chemical Structure]	entre anti-
[Characterization Data Summary]	_
Analytical Test	Results
Identification by "H-NMR, HPLC Purity lasted	Consistent with the above structur >=08%
	1000
Authorized Signature:	
Date:	



Catal

Ref: cleanjang-healthy.

Plant library of Stevia rebaudiana

Catalog	Product Name	CAS Number	Manual
FN99548	Stevioside	57817-89-7	PDF
CFN99110	Rebaudioside A	58543-16-1	PDF
CFN97326	Stigmasterol	83-48-7	PDF
CFN99916	Beta-Sitosterol	83-46-5	PDF
CFN98843	Apigenin	520-36-5	PDF
CFN98981	Cosmosiin	578-74-5	PDF
CFN98601	Vitexin	3681-93-4	PDF
CFN98784	Luteolin	491-70-3	PDF
CFN98565	Luteolin-7-O-glucoside	5373-11-5	PDF
CFN98756	Kaempferitrin	482-38-2	PDF
CFN98850	Quercitrin	522-12-3	PDF
CFN90544	Isosteviol	27975-19-5	n/a
CFN90470	Rebaudioside B	58543-17-2	PDF
CFN97018	Paniculoside I	60129-63-7	n/a
CFN97019	Paniculoside II	60129-64-8	n/a
CFN90257	Rebaudioside D	63279-13-0	n/a
CFN90258	Rebaudioside C	63550-99-2	n/a
CFN92266	Dulcoside A	64432-06-0	PDF
CFN90984	Rebaudioside G	127345-21-5	n/a



Natural Products

Alkaloids	
23512-53-0	4.5-Dihydropiperlonguminine
2584997-91-9	N-trans-p-CoumaroyI-N-trans-feruloyI-3-hydroxy-cadaverine
1210820-67-9	1-Methyl-6 8-dimethoxyguinoline-2.1H-one
2854-40-2	Cyclo(L-Pro-L-Val)
57089-60-8	Cyclo(L-Pro-L-IIe)
2873-36-1	Cyclo(L-Leu-L-Pro)

Lignans	

126253-42-7	Icariside E4
285136-03-0	alpha: [4: [(1E): 2-Carboxyethenyi]: 2-hydroxyphenoxy]-beta: 3.4-trihydroxybenzene, alpha: [4: [(1E): 2-Carboxyethenyi]: 2-hydroxybenzene, alpha: [4: [4: [4: [4: [4: [4: [4: [4: [4: [4
42123-27-3	Tetradehydropodophyllotoxin
144608-09-3	Schizotenuin A
848844-79-1	Secoisolariciresinol 9.9'-diacetate
159736-38-6	Clinopodic acid E

Phenols

6-O-GalloyIglucose
4-(2-Hydroxyethyl)benzoic acid
New compound 26
(E)-[6]-Dehydroparadol
Dihydroparado)
Ellagic acid-4-O-beta-D-xylopyranoside
(

Chalcones

189299-05-6	Xanthohumol C
152340-67-5	Calythropsin
1776-30-3	2'.4'-Dihydroxychalcone
197227-39-7	Tetrahydroxymethoxychalcone
111316-17-7	2'-O-Methylphloretin.(4.2'.4'-Trihydroxy-6'-methoxydihydrochalcone)
94103-36-3	2:4:6:4-Tetramethoxychalcone

Quinones

4707-32-8	Beta-Lapachone
125906-49-2	$\label{eq:constraint} 3-(beta-D-Glucopyranosyloxy)-1.6-dihydroxy-2-methyl-9,10-anthracenedione$
34298-85-6	Emodin 6-O-beta-D-glucoside
83-72-7	2-Hydroxy-1.4-naphoquinone
2348-82-5	2-Methoxy-1.4-naphthoquinone
635-12-1	1.4-Anthraquinone

 Coumarins

 57601-61-3
 Dipetalolactone

 1857-05-2
 6.7-Dimethoxy-4-phenylcoumarin

 73292-92-9
 2'E-3'-formaldehydylosthole

 135064-04-9
 Umbetilferone 7-O-rutinoside

 477-32-7
 Cardine

 104196-69-2
 Neoartanin

Steroids	
54082-42-7	25.28-Di-epi-cyasterone
562-34-5	Chlorogenin
141360-88-5	Polyporusterone A
38395-01-6	Kidjolanin
20817-72-5	Stigmastadienone
14214-69-8	Cholest-4.6-diene-3-ol

Iridoids	
79172-04-6	8-Epiloganin
93930-20-2	Gardoside methyl ester
182172-02-7	10-Carboxyloganin
158599-51-0	Phloyoside II
1008532-71-5	8-Dehydroxy.shanzhiside
276691-32-8	Phlorigidoside C

Flavonoids

19046-26-5	Wistin
14292-40-1	Taxifolin 7-O-glucoside
18467-06-6	Isokaempferide 7-rutinoside
76060-29-2	Quercetagetin 3-methyl ether 7-glucoside
1627598-00-8	Nepetin 7-O-beta-D-glucopyranoside
69256-15-1	Leucocyanidin
Xanthones	
164365-71-3	11-hydroxy-1-isomangostin
2522597-99-3	Garcixanthones B
2182593-73-1	Mangostanaxanthone IV
138529-04-1	Diphlorethohydroxycarmalol
84272-84-4	3'-O-Angeloyihamaudol
117221-65-5	Neolancerin
Anthraquinones	
473335 33.9	Packalain soluted compared P

473225-22-8	Barbaloin-related compound B
134863-92-6	10-Hydroxyaloin B
473225-21-7	Barbaloin-related compound A
569-05-1	Fallacinol
132210-48-1	Elgonica dimer A
70588-06-6	Chrysoobtusin

Cerebrosides	
88642-46-0	Cerebroside B
606125-07-9	Momor-cerebroside I
340702-68-3	2-2-(Hydroxytetracosanoylamino)-octadecane-1.3.4-triol tetraacetate
295803-03-1	Gynuramide II
154801-30-6	2-(2'-Hydroxytetracosanoylamino)-octadecane-1.3.4-triol
115074-93-6	Soyacerebroside II



Natural Products

Miscellaneous	
50257-98-2	(<u>E)-Tonghaosu</u>
4575-53-5	(Z) <u>-Tonghaosu</u>
6338-41-6	5-Hydroxymethyl-2-furoic acid
77-93-0	Triethyl citrate
129499-78-1	Ascorbyl glucoside
1109-28-0	Maltotriose

Phenylpropanoid	
537-33-7	Sinapyl alcohol
31564-49-5	2-Glucosyloxy-4-methoxycinnamic acid.(Z-GMCA)
150892-86-7	cis-2-Hydroxy_4-methoxycinnamic_acid_2-glucoside
116872-05-0	Magnoloside B
122-97-4	HydrocinnamyLalcohol
478294-16-5	Cimiracemate A

Triterpenoids

Sesquiterpenoids 97643-91-9

63568-73-0

5945-42-6

3668-14-2

20267-92-9

226546-99-2

8-epi-Helenalin

Carpesiolin

Aromaticin

Bigelovin

Oxyphyllol B

Hydroxylinderstrenolide

124961-62-2	Raddeanoside R9
781676-86-6	Raddeanoside R18
228251-25-0	24 - epi-24 - O-acetyl-7.8 - dide hydrohydroshengmanol-3 - O-beta-D-xylopyranoside
150972-77-3	24-epi-7.8-Didehydrocimigenol 3-xyloside
2101977-10-8	Cimiricaside E
197013-75-5	Patrinia saponin H3

Diterpenoid	s
60129-60-4	Steviolmonoside
93380-12-2	Epirosmanol
205440-23-9	Taxezopidine H
76429-85-1	10-Deacetylcephalomannine
243448-72-8	Necorthosiphol A
142741-25-1	Orthosiphol A
142741-25-1 Monoterpenoid	
Monoterpenoid	s .
Monoterpenoid 184103-78-4	s Benzoyla/biflorin
Monoterpenoid 184103-78-4 4221-98-1	s Benzoylalbiflorin (B)::(:):alpha-Phellandrene
Monoterpenold 184103-78-4 4221-98-1 3391-90-0	s Benzoylalbiflorin (B):(:):-alpha-Phellandrene (S):(:):-Pulegone



Bioactive Products

Hot Research

Antioxidants	Anti-Inflammatory
Anticancer	Hepatoprotective
Anti-diabetic	Analgesia
Antimalaric	Anti-proliferation
Anti-Atheroscierotic	Cardioprotective
Antibacterial	Antifungal
Anti-allergy	Inhibitors
Antiviral	Immunomodulators
Anti-hyperlipidemic	Antihypertensive
Antileukaemic	Anti-platelet aggregation
Anti-photosging	Antihyperglycemic
Antiangiogenic	Pro-apoptotic
Neuroprotection	Antidepressant
Anti-osteoporosis	Anti-obesity
Cancer chemoprevention	Cytotoxic
Vasorelaxation	Insecticides
Food additives	Dietary supplement
Toxicity	

Agricultural Research

Antifeedant	Herbicides	
Anthelmintic	Algicides	
Phytoalexin	Insecticides	
Post ingestive	Phytotoxic	
Growth-promoting effects	Anti-plant allelopathy	
Anti-powdery mildew	Phagostimulants	
Acaricidal	Nematicidal	
Nodulation genes inducers	Anti-juvenile hormone	
Moulting hormone activity	Ecdysteroid antagonists	
Anti-plant-germination and growth		



Bioactive Products

Food and Spice Research

Food additives	Fragrance ingredients
Preservatives	Sweeteners
Pigments	Nutraceutical
Dietary supplement	Anti-fatigue
Antinutritional	Hygroscopic

Agricultural Research

Anti-photoaging	Cosmetic additives	
Skin-lightening agents	Skin-whitening agents	
Anti-melanogenic	Preservatives	
Pigments	Anti-wrinkle	
Maudring bermang articles		

Plant Catalog

Screening Library of Plants

Morus alba L	Ginkgo biloba L.	Cinnamomum cassia Presi	Ganoderma lucidum (Curtis) P. Karst
A unique collection of 55 Morus alba L for high throughput screening (HTS) and New drug research	A unique collection of 41 Ginkgo biloba L. for high throughput screening (HTS) and New drug research	A unique collection of 47 Cinnamomum cassia PresI for high throughput screening (HTS) and New drug research	A unique collection of 66 Ganoderma lucidum (Curtis) P. Karst for high throughput screening (HTS) and New drug research
Salvia miltiorrhiza Bge.	Morus alba L.	Panax ginseng C. A. Mey.	Glycyrrhiza uralensis Fisch.
A unique collection of 81 Salvia miltiorrhiza Bge. for high throughput screening (HTS) and New drug research	A unique collection of 69 Morus alba L. for high throughput screening (HTS) and New drug research	A unique collection of 52 Panax ginseng C. A. Mey. for high throughput screening (HTS) and New drug research	A unique collection of 50 Glycyrrhiza uralensis Fisch. for high throughput screening (HTS) and New drug research
Plant Catalog			
Panax ginseng C. A. Mey.	Illicium verum	Lilium pumilum D. C.	Lithosperraum erythrorhizon Sieb. et Zucc.
Fraxinus chinensis Roxb	Foeniculum vulgare	Aplum graveolens	Polygonum cuspidatum Sieb. et Zucc.
/itis vinifera L.	Veratrum grandiflorum Loes.	Veronica linariifolia	Physalis alkekengi
Semiaquilegia adoxoides (DC.) makino.	Rubus corchorifolius	Pinus cerebral	Cherry tree
Artemisia absinthium L.	Camellia sinensis (L.) O. Kuntze.	Tripterygium wilfordii Hook. f.	Rubia cordifolia L.
Polygala japonica Houtt.	Ginkgo biloba L.	Garcinia cambogia	Angelica acutilobac (Sieb. et Zucc.) Ki
Centella asiatica (L.) Urban	Asarum sieboldii Miq.	Scrophularia ningpoensis Hemsl.	Cornus officinalis Sieb. et Zucc.



CONTACT US











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