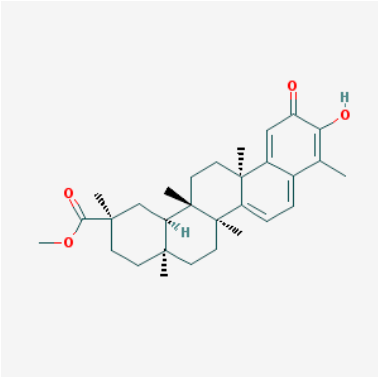
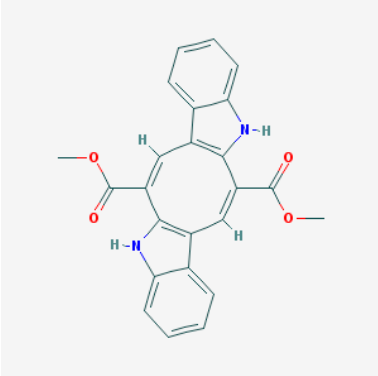
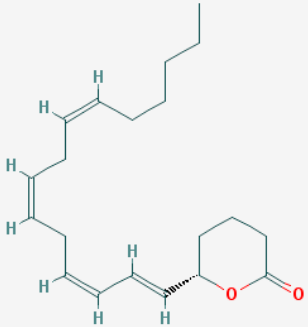
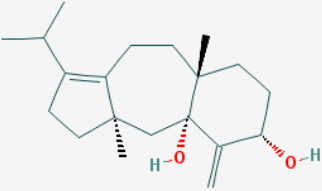
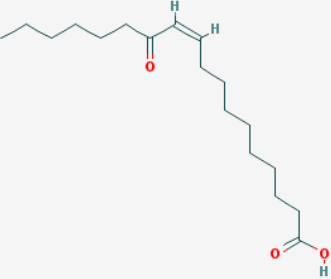
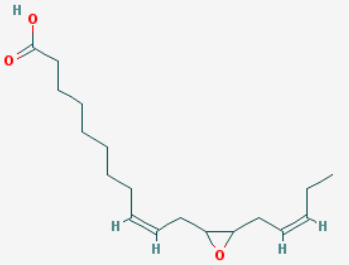
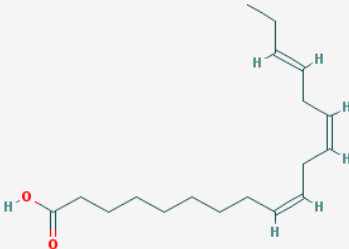
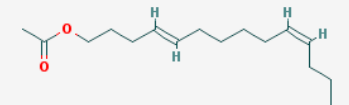
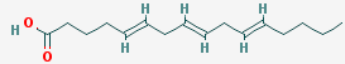
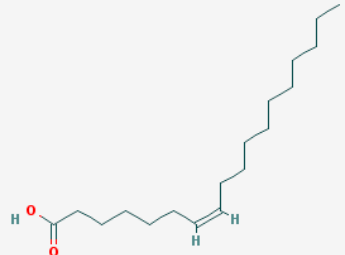
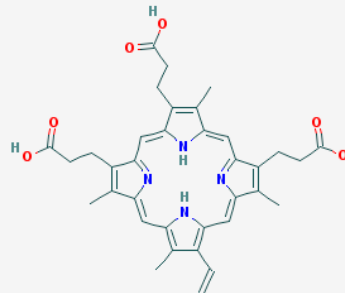


Supplementary Material Table S1: Proposed compounds present in *C. racemosa* chloroform extract identified via LCMS analysis.

No.	Compound name	Formula	Structure	Reported activity
1	Pristimerin; Celastrol methyl ester	C ₃₀ H ₄₀ O ₄		<p>Pristimerin is a triterpenoid commonly found in various members from the Celastraceae family, including <i>Maytenus ilicifolia</i>. Anticancer activities of Pristimerin have been reported in uveal melanoma [1], glioma [2], colorectal cancer [3], osteosarcoma [4], cervical cancer [5], ovarian cancer [62], pancreatic cancer [6, 7], prostate cancer [8, 9], and breast cancer [10, 11].</p> <p>Pristimerin has also been shown to have anti-inflammation effects [12-14] and antioxidant activities [62]. Very little is known about its antibacterial effects but this triterpenoid has been shown to inhibit SARS-CoV and cytomegalovirus <i>in vitro</i> [15, 16].</p>
2	Caulerpin	C ₄₀ H ₁₈ N ₂ O ₄		<p>Caulerpin is a red pigmented bis-indole alkaloid isolated from marine <i>Caulerpa</i> algae spp. This molecule has been demonstrated to have anticancer effects [17-19]. It is also known to have antinociceptive effects [20, 21]. This alkaloid is known to possess antibacterial activity [44] and anti-tuberculosis activity [22].</p>

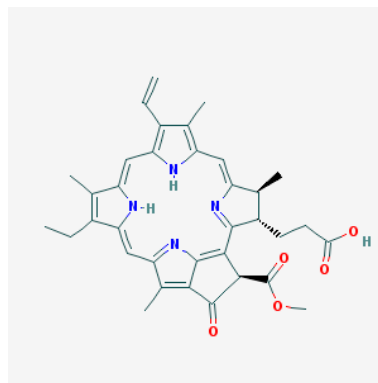
3	5(S)-HETE lactone	$C_{20}H_{30}O_2$	 <p>The structure shows a long hydrocarbon chain with four double bonds. The first double bond is at the end of a pentyl group. The second double bond is at the end of a propyl group. The third double bond is at the end of a propyl group. The fourth double bond is at the end of a propyl group. The chain ends in a lactone ring (a six-membered ring with an oxygen atom and a carbonyl group).</p>	Bioactivity has not been reported.
4	Isoamijiol	$C_{20}H_{30}O_2$	 <p>The structure is a complex diterpenoid with a central ring system consisting of two fused seven-membered rings and two fused six-membered rings. It has several methyl groups, a double bond, and two hydroxyl groups (one on a double bond, one on a saturated carbon).</p>	Isoamijiol is a diterpenoid. This molecule is commonly found in brown algae of genus <i>Dictyota</i> [23]. Its bioactivity is not known.
5	12-oxo-10Z-octadecenoic acid	$C_{18}H_{32}O_3$	 <p>The structure is a long-chain fatty acid with 18 carbons. It has a ketone group at C-12 and a Z double bond at C-10. The chain ends in a carboxylic acid group.</p>	Bioactivity has not been reported.

6	12(13)-epoxy-9Z,15Z-octadecadienoic acid or α -9(10)-EpODE	$C_{18}H_{30}O_3$		Bioactivity has not been reported.
7	9Z,12Z,15E-octadecatrienoic acid	$C_{18}H_{30}O_2$		Bioactivity has not been reported.
8	4E,10Z-Tetradecadienyl acetate	$C_{16}H_{28}O_2$		A sex pheromone found in Apple Leafminer Moth, <i>Phyllonorycter ringoniella</i> [24].

9	trans-5, trans-8, trans-11- hexadecatrienoic acid	$C_{16}H_{26}O_2$		Bioactivity has not been reported.
10	7-Octadecenoic acid	$C_{18}H_{32}O_2$		This monounsaturated fatty acid is found in pistachio nuts [25] but its bioactivity is not known.
11	Harderoporphyrin	$C_{35}H_{36}N_4O_6$		Harderoporphyrin was first identified and isolated from rat Harderian glands. Its presence and amount in the rodent harderian glands has been studied to be used as a biomarker for arsenic exposure [26].

12

Pheophorbide a

 $C_{35}H_{36}N_4O_5$ 

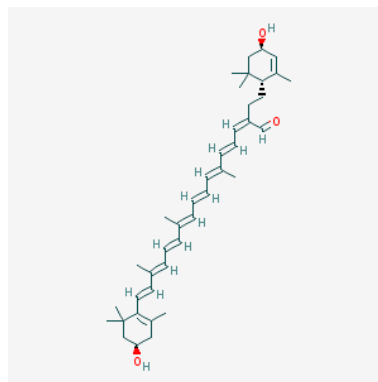
This bioactive molecule has been isolated from marine algae species and terrestrial plants. It is known to have antioxidant and anti-inflammation effects [63].

This molecule also induces photodynamic inactivation of *Trypanosoma cruzi* [27]. It is also known to have anti-*S. aureus* [64] and anti-hepatitis C virus [28].

On its own, pheophorbide-a has been shown to induce apoptosis in human hepatocellular carcinoma cells [29].

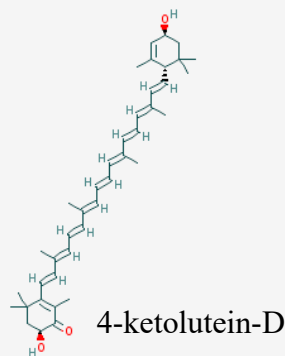
13

Micromonal

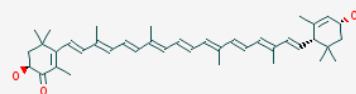
 $C_{40}H_{56}O_3$ 

This carotenoid is found in green algae Prasinophyceae [30].

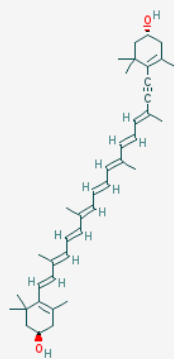
14 4-Ketolutein D and
4-ketolutein F $C_{40}H_{54}O_3$



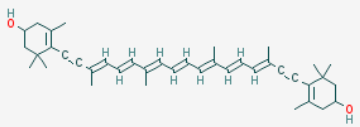
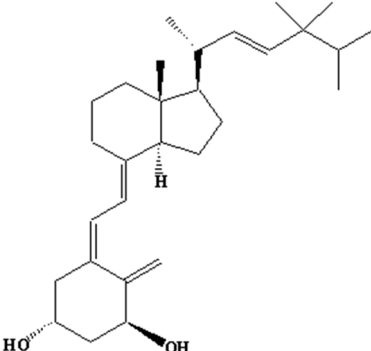
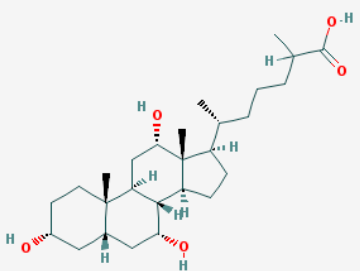
4-ketolutein is a ketolated carotenoid found in algae and red tilefish (*Branchiostegus japonicas*) [31]. This carotenoid is now patented to as a food additive in egg yolk [32].

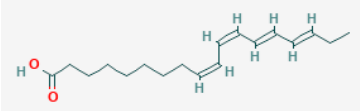
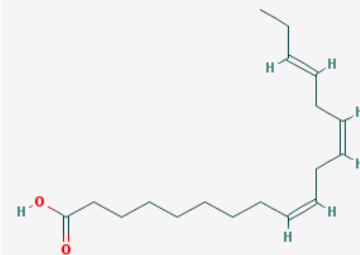
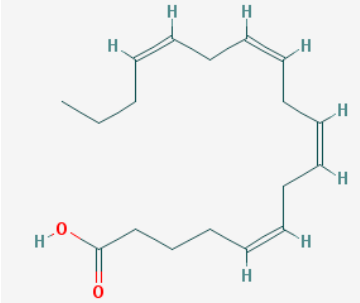


15 Diatoxanthin/ 7,8-
Didehydrozeaxanthin $C_{40}H_{54}O_2$

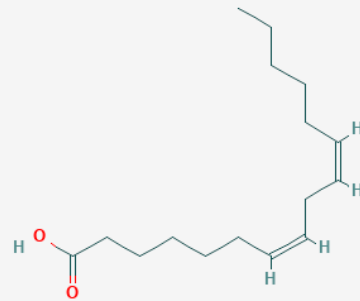
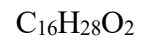


Diatoxanthin is a carotenoid in algae that plays a role in dissipating excess energy in algae species [33]. This carotenoid has been shown to have anti-inflammation in the LPS-induced pro-inflammatory in RAW264.7 cells [34].

16	Pectenoxanthin; Alloxanthin; Manixanthin;	$C_{40}H_{52}O_2$		<p>Alloxanthin is a carotenoid that has been isolated from various marine species, including sea pineapple (<i>Halocynthia roretzi</i>), channel catfish, giant scallop (<i>Pecten maximus</i>) and edible mussel (<i>Mytilus edulis</i>) [35]. Similarly, to Diatoxanthin, this carotenoid was shown to be anti-inflammation <i>in vitro</i> [34].</p>
17	1 α -hydroxy-24-methylvitamin D2 / 1 α -hydroxy-24-methylergocalciferol	$C_{29}H_{46}O_2$		Vitamin D.
18	Coprocholic acid; 3 α ,12 α -Dihydroxy-5 β -chol-8(14)-en-24-oic Acid	$C_{24}H_{38}O_4$		<p>Coprocholic acid, also called 3α, 7α, 12α-Trihydroxy-5β-cholestan-26-oic acid, is a bile acid.</p>

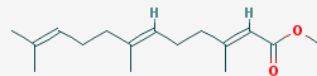
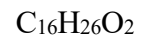
19	5Z,8Z,11Z,14Z- Octadecatetraenoic acid	$C_{18}H_{30}O_2$	 <p>The structure shows an 18-carbon chain with a carboxylic acid group at C1. Double bonds are located at C5, C8, C11, and C14, all in the Z configuration. The chain is drawn in a zig-zag pattern to represent the cis configuration of the double bonds.</p>	Bioactivity has not been reported.
20	9Z,12Z,15E- octadecatrienoic acid	$C_{18}H_{30}O_2$	 <p>The structure shows an 18-carbon chain with a carboxylic acid group at C1. Double bonds are located at C9, C12, and C15. C9 and C12 are in the Z configuration, while C15 is in the E configuration. The chain is drawn in a zig-zag pattern.</p>	Bioactivity has not been reported.
21	5Z,8Z,11Z,14Z- octadecatetraenoic acid	$C_{18}H_{28}O_2$	 <p>The structure shows an 18-carbon chain with a carboxylic acid group at C1. Double bonds are located at C5, C8, C11, and C14, all in the Z configuration. The chain is drawn in a zig-zag pattern.</p>	Bioactivity has not been reported.

22 10E, 12Z-
Tetradecadienyl
acetate



A sex pheromone found in the avocado pest, *Amorbia cuneana* (Walsingham) [36].

23 Methyl farnesoate



A sesquiterpenoid hormone that is found in the *Drosophila* metamorphosis and stingless bee [37, 38].

Reference List for Supplementary Material Table S1: Proposed compounds present in *C. racemosa* chloroform extract identified via LCMS analysis.

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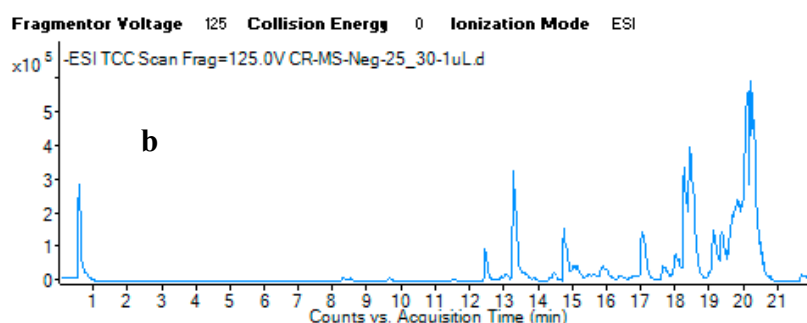
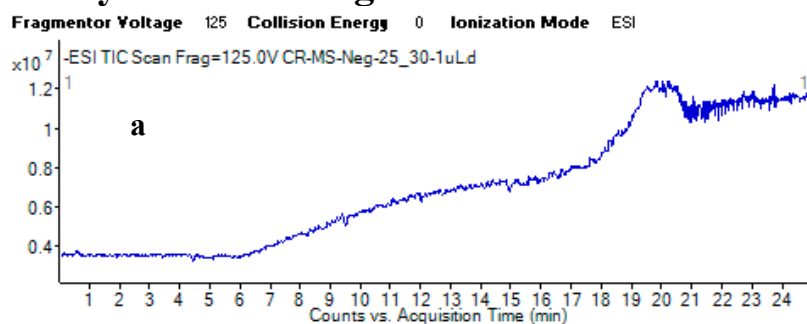
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*references are cited in main text.

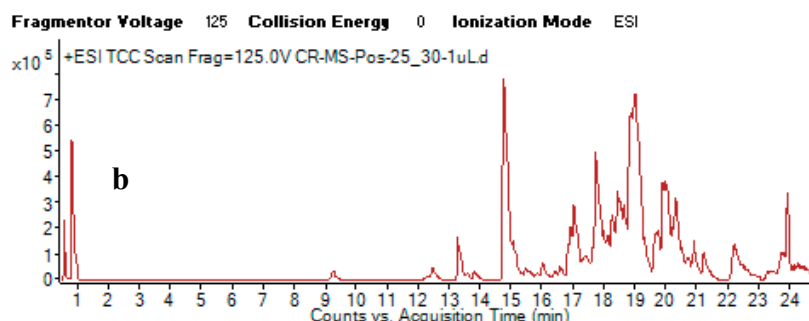
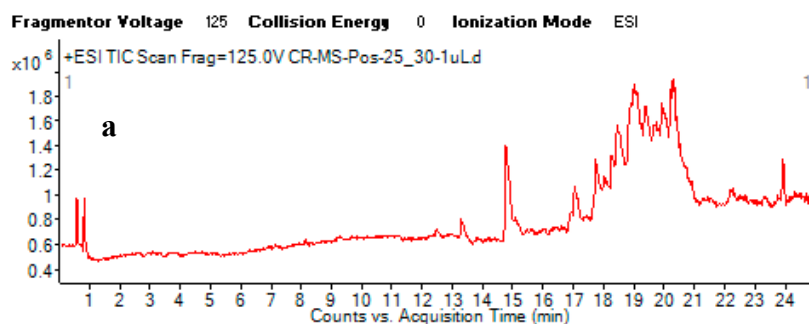
Supplementary Material Table S2: Molecular Formula of 25 Unknown Compounds from *C. racemosa* Chloroform Extract.

No.	Formula	Mass	m/z
1	C ₄ H ₁₂ O ₂ S ₂	156.02781	157.03509
2	C ₂₁ H ₄₀ O S	340.28029	341.28732
3	C ₂₃ H ₁₆ N ₂ O ₄	384.11126	385.11829
4	C ₂₂ H ₂₅ N ₅ O ₂	391.20081	392.20754
5	C ₁₄ H ₃₃ N ₉ O S	375.25339	393.28682
6	C ₂₈ H ₄₉ N O ₆	495.35626	496.36363
7	C ₂₈ H ₅₁ N O ₆	497.37183	498.3793
8	C ₂₈ H ₅₃ N O ₆	499.38752	500.39495
9	C ₂₄ H ₄₉ N ₁₁ O S	539.38349	540.39043
10	C ₃₃ H ₅₅ N O ₆	561.40325	562.41052
11	C ₃₈ H ₅₃ N ₃ O	567.4197	568.42687
12	C ₂₉ H ₅₃ N ₅ O ₇	583.39411	584.40116
13	C ₃₁ H ₅₃ N ₉ S	583.41502	584.42149
14	C ₃₆ H ₅₃ N ₅ O ₂	587.4194	588.4262
15	C ₃₅ H ₅₉ N O ₆	589.43467	590.44186
16	C ₃₈ H ₅₃ N ₃ O ₃	599.40895	600.41655
17	C ₃₉ H ₅₁ N ₇	617.4208	618.42806
18	C ₁₄ H ₃ N O ₄ S ₂	312.95072	311.94352
19	C ₂₅ H ₁₄ N ₆ S	430.10007	429.09259
20	C ₂₄ H ₂₇ N O ₃ S ₂	441.14341	440.13623
21	C ₂₆ H ₂₇ N O ₃ S ₂	465.14329	464.1363
22	C ₂₅ H ₁₆ N ₆ O ₂ S ₂	496.07662	495.06912
23	C ₁₅ H ₈ N ₄ O ₈ S ₄	499.92316	498.9159
24	C ₂₀ H ₄₆ N ₁₀ O ₅ S	538.33685	537.32925
25	C ₃₆ H ₆ O S ₅	613.90261	612.89583

Supplementary Materials Figures S1 & S2: Chromatograms



Supplementary Material Figure S1: *C. racemosa* chloroform extract was subjected to LCMS qualitative analysis using negative ion mode. The compounds were separated based on m/z ratio and retention time in the column, a) showing total ion current (TIC) chromatogram and b) showing total compound chromatogram (TCC). 48 peaks are observed in the extract.

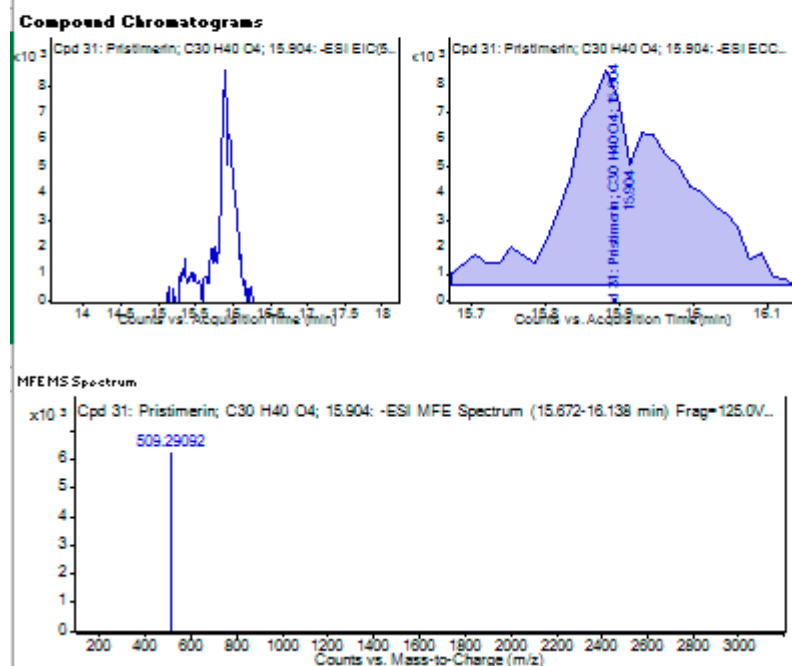


Supplementary Material Figure S2: *C. racemosa* chloroform extract was subjected to LCMS qualitative analysis using positive ion mode. The compounds were separated based on m/z ratio and retention time in the column, a) showing total ion current (TIC) chromatogram and b) showing total compound chromatogram (TCC). 74 peaks are observed in the extract.

Supplementary Material Figure S3: Mass spectra of individual compounds listed in Supplementary Material Table S1.

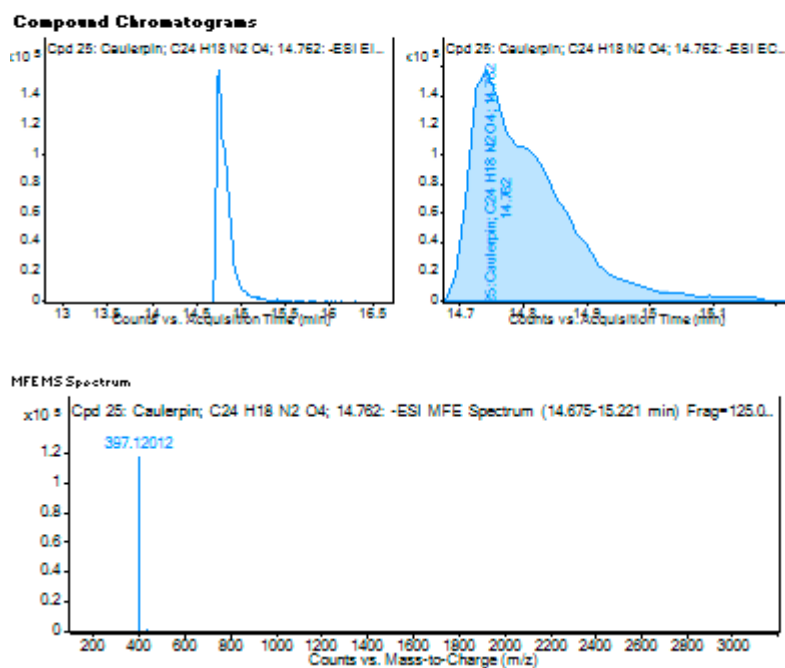
Compound 1: Pristimerin

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 31: Pristimerin; C30 H40 O4; 15.904	Pristimerin	509.29092	15.9	Find by Molecular Feature	464.29287



Compound 2: Caulerpin

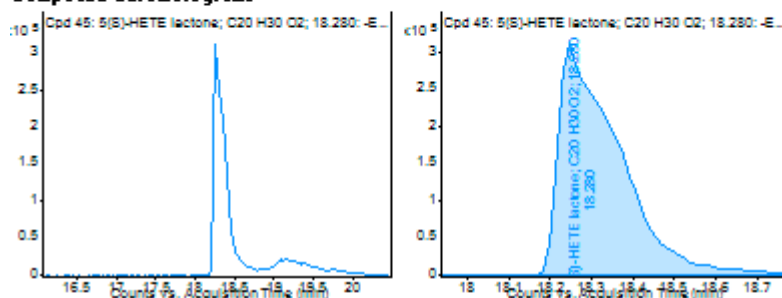
Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 25: Caulerpin; C24 H18 N2 O4; 14.762	Caulerpin	397.12012	14.76	Find by Molecular Feature	398.12732



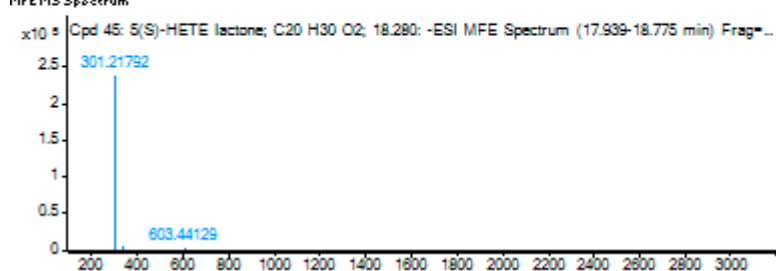
Compound 3: 5(S)-HETE lactone

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 45: 5(S)-HETE lactone; C20 H30 O2; 18.280	5(S)-HETE lactone	301.21792	18.28	Find by Molecular Feature	302.22512

Compound Chromatograms



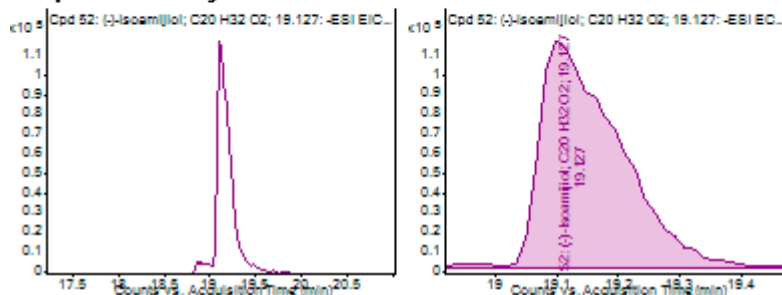
MFEMS Spectrum



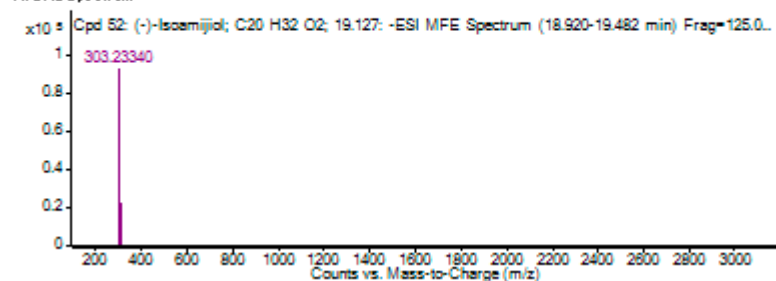
Compound 4: Isoamijiol

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 52: (-)-Isoamijiol; C20 H32 O2; 19.127	(-)-Isoamijiol	303.2334	19.13	Find by Molecular Feature	304.24063

Compound Chromatograms



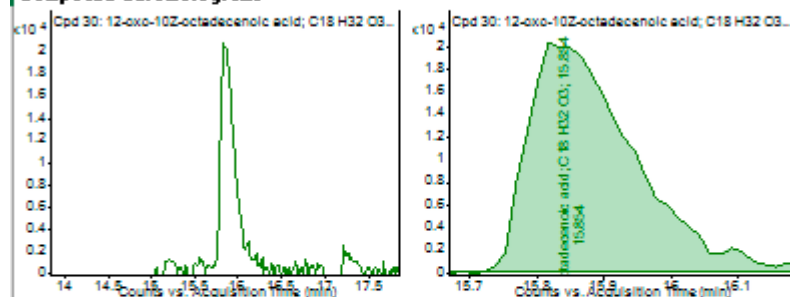
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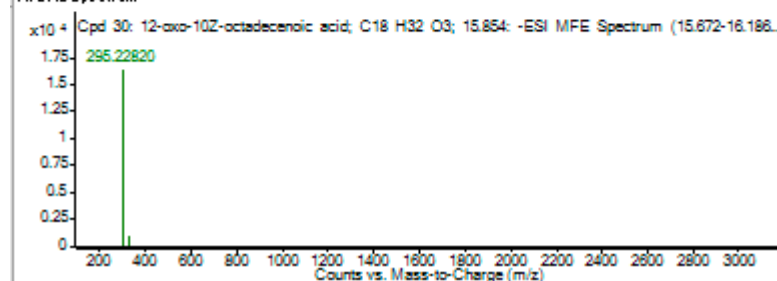
Compound 5: 12-oxo-10Z-octadecenoic acid

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 30: 12-oxo-10Z-octadecenoic acid; C18 H32 O3; 15.854	12-oxo-10Z-octadecenoic acid	295.2282	15.85	Find by Molecular Feature	296.2353

Compound Chromatograms



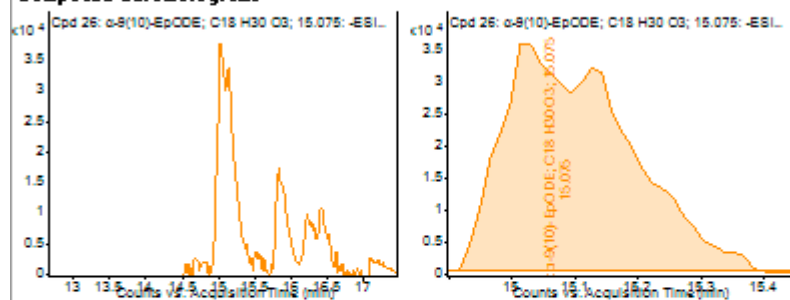
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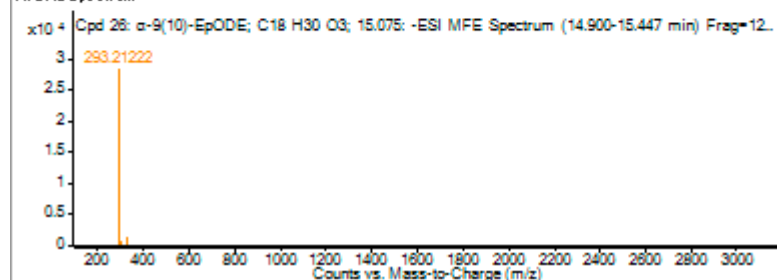
Compound 6: 12(13)epoxy-9Z,15Z-octadecadienoic acid

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 26: α-9(10)-EpODE; C18 H30 O3; 15.075	α-9(10)-EpODE	293.21222	15.08	Find by Molecular Feature	294.21965

Compound Chromatograms



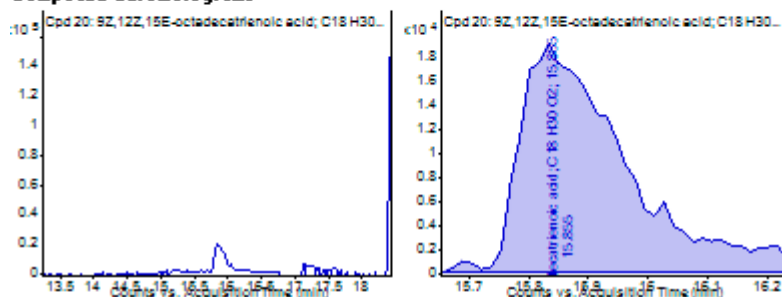
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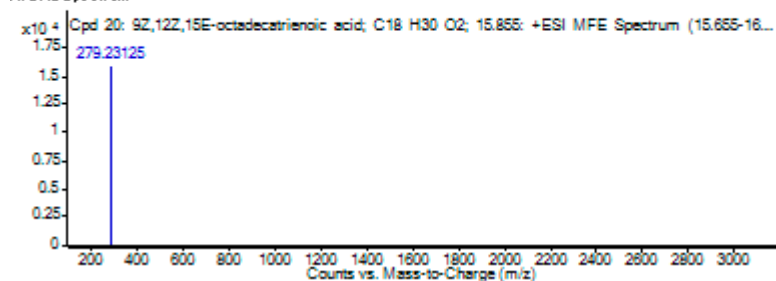
Compound 7: 9Z, 12Z, 15E-octadecatrienoic acid

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 20: 9Z,12Z,15E-octadecatrienoic acid; C18 H30 O2; 15.855	9Z,12Z,15E-octadecatrienoic acid	279.23125	15.86	Find by Molecular Feature	278.22405

Compound Chromatograms



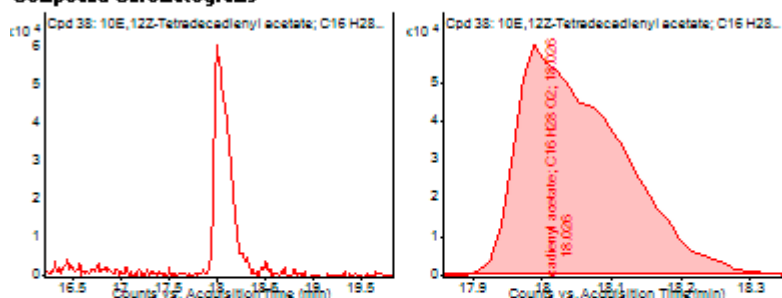
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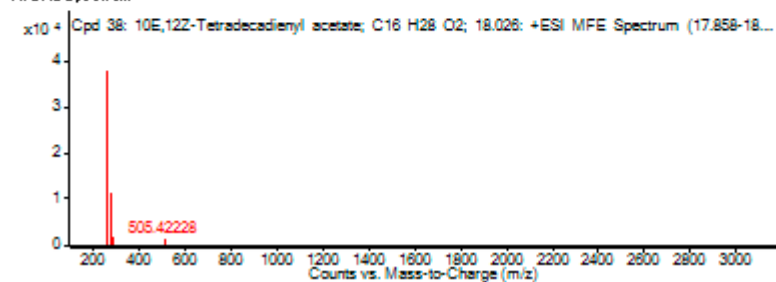
Compound 8: 4E, 10Z-tetradecadienyl acetate

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 38: 10E,12Z-Tetradecadienyl acetate; C16 H28 O2; 18.026	10E,12Z-Tetradecadienyl acetate	253.21626	18.03	Find by Molecular Feature	252.20921

Compound Chromatograms



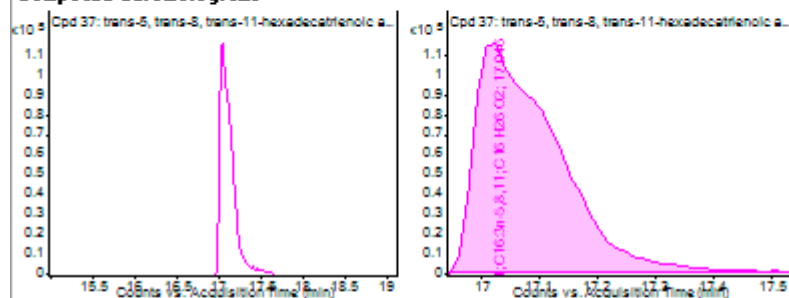
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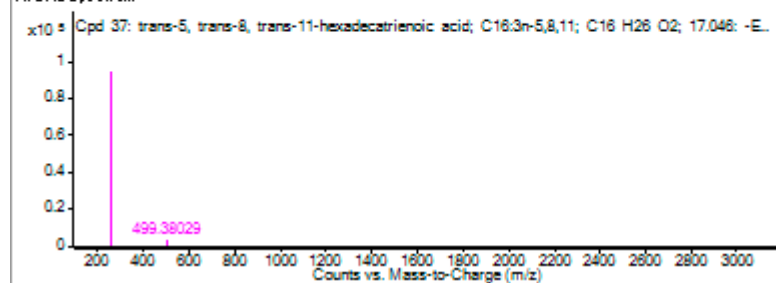
Compound 9: trans-5, trans-8, trans-11, hexadecatrienoic acid

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 37: trans-5, trans-8, trans-11-hexadecatrienoic acid; C16:3n-5,8,11; C16 H26 O2; 17.046	trans-5, trans-8, trans-11-hexadecatrienoic acid; C16:3n-5,8,11	249.18627	17.05	Find by Molecular Feature	250.19348

Compound Chromatograms



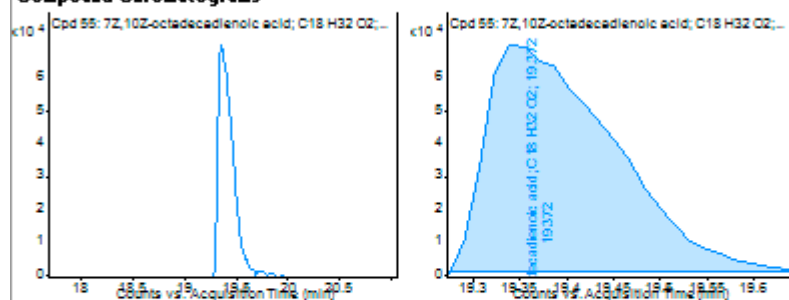
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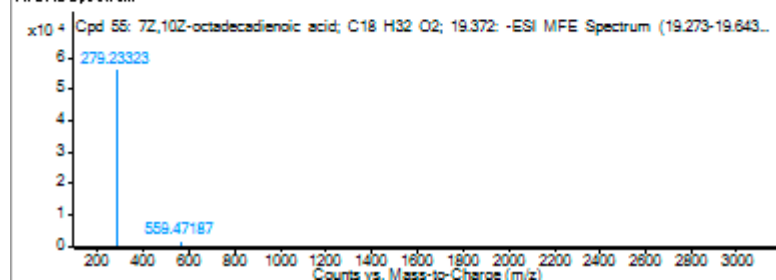
Compound 10: 7-octadecenoic acid

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 55: 7Z,10Z-octadecadienoic acid; C18 H32 O2; 19.372	7Z,10Z-octadecadienoic acid	279.23323	19.37	Find by Molecular Feature	280.24054

Compound Chromatograms



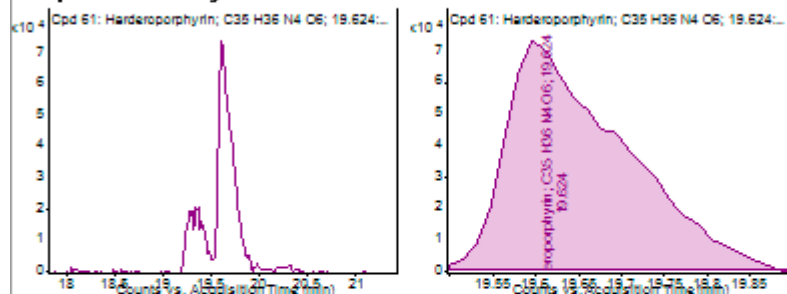
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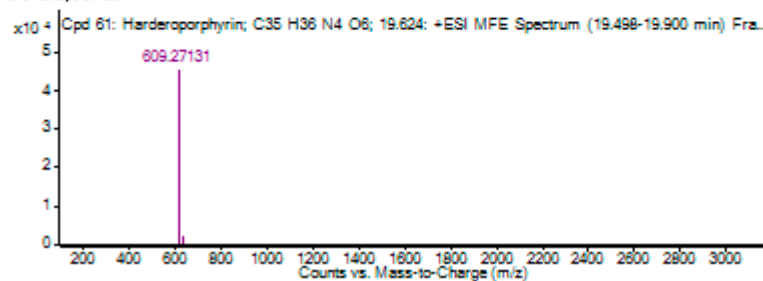
Compound 11: Harderoporphyrin

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 61: Harderoporphyrin; C35 H36 N4 O6; 19.624	Harderoporphyrin	609.27131	19.62	Find by Molecular Feature	608.26415

Compound Chromatograms



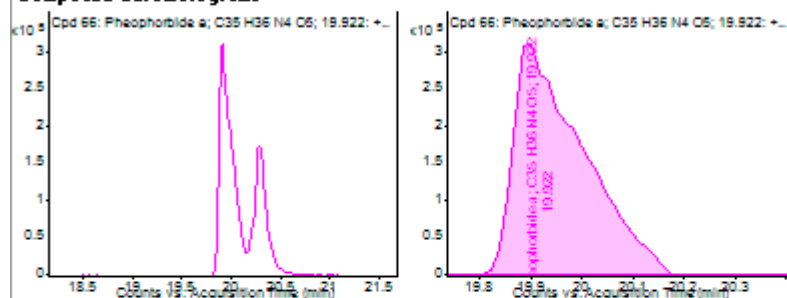
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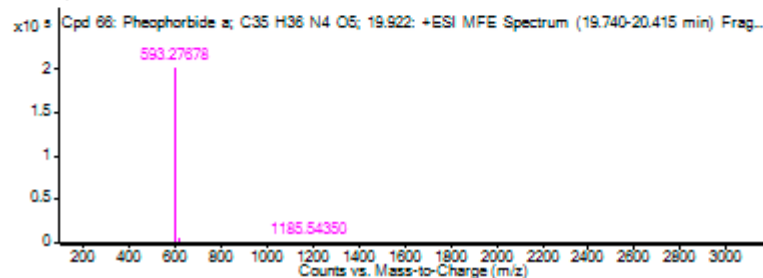
Compound 12: Pheophorbide a

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 66: Pheophorbide a; C35 H36 N4 O5; 19.922	Pheophorbide a	593.27678	19.92	Find by Molecular Feature	592.26945

Compound Chromatograms



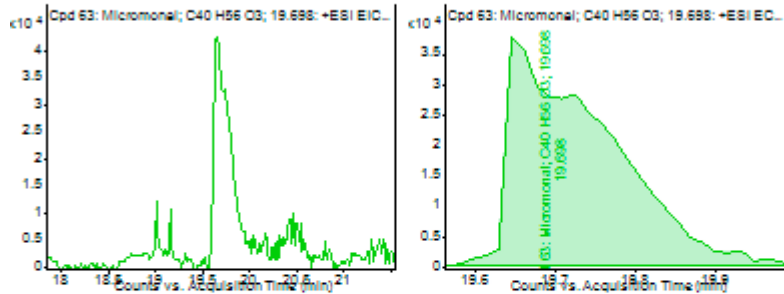
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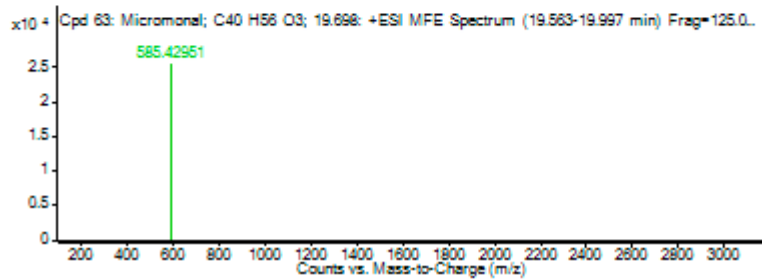
Compound 13: Micromonal

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 63: Micromonal; C40 H56 O3; 19.698	Micromonal	585.42951	19.7	Find by Molecular Feature	584.42184

Compound Chromatograms



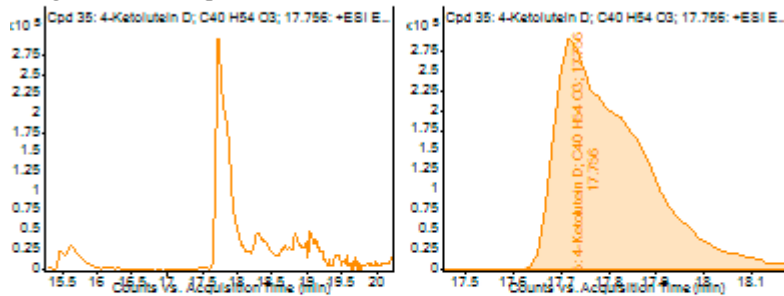
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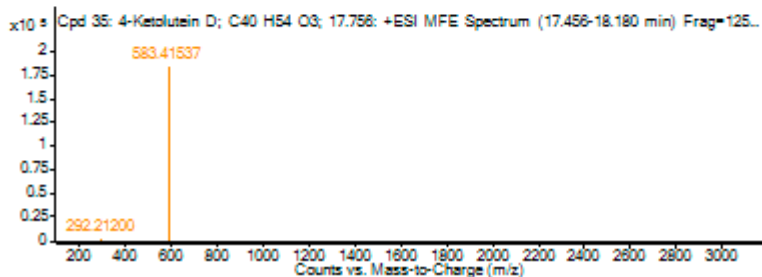
Compound 14: 4-ketolutein D and 4-ketolutein F

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 35: 4-Ketolutein D; C40 H54 O3; 17.756	4-Ketolutein D	583.41537	17.76	Find by Molecular Feature	582.40796

Compound Chromatograms

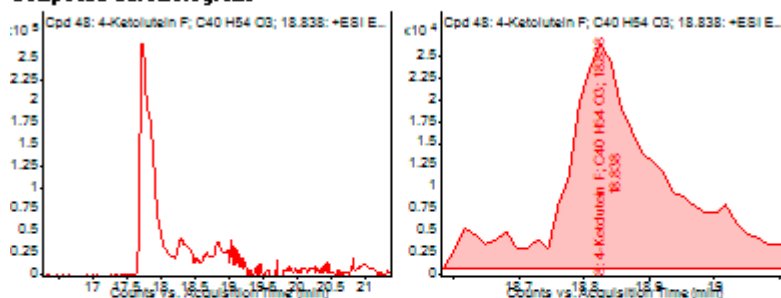


MFEMS Spectrum

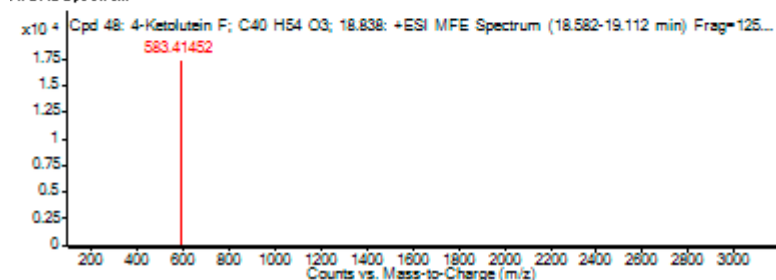


Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 48: 4-Ketolutein F; C40 H54 O3; 18.838	4-Ketolutein F	583.41452	18.84	Find by Molecular Feature	582.40748

Compound Chromatograms



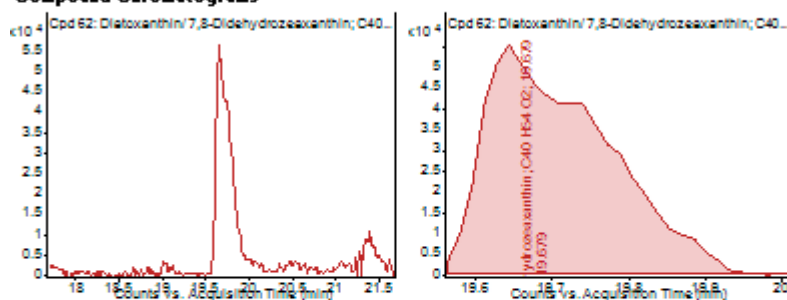
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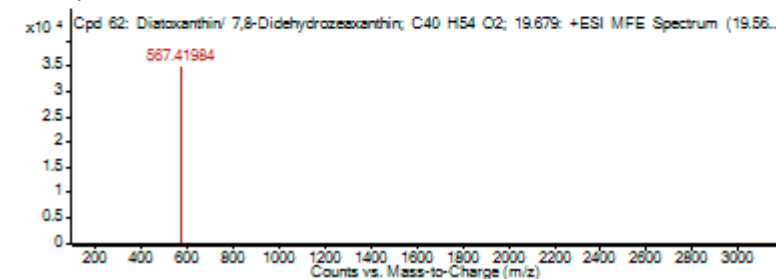
Compound 15: Diatoxanthin/7,8-Didehydrozeaxanthin

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 62: Diatoxanthin/ 7,8-Didehydrozeaxanthin; C40 H54 O2; 19.679	Diatoxanthin/ 7,8-Didehydrozeaxanthin	567.41984	19.68	Find by Molecular Feature	566.4117

Compound Chromatograms



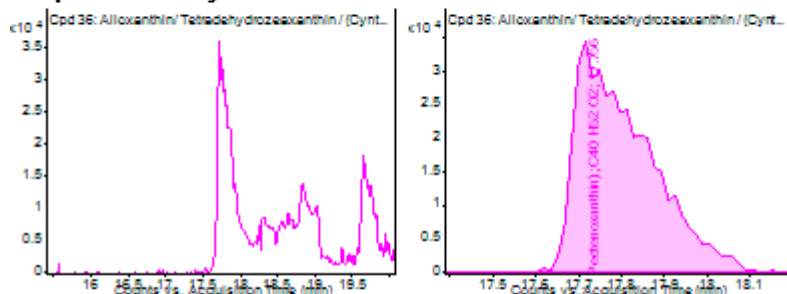
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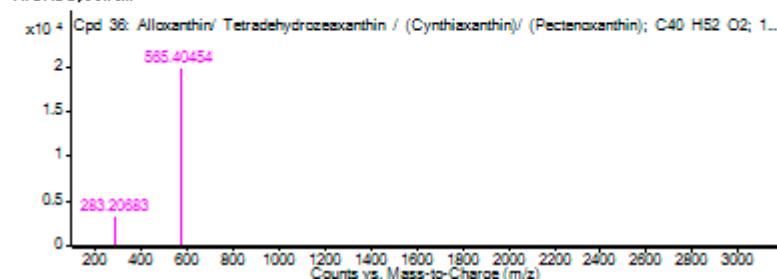
Compound 16: Pectenoxanthin; Alloxanthin; Manixanthin

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 36: Alloxanthin/ Tetradehydrozeaxanthin / (Cynthiixanthin)/ (Pectenoxanthin); C40 H52 O2; 17.756	Alloxanthin/ Tetradehydrozeaxanthin / (Cynthiixanthin)/ (Pectenoxanthin)	565.40454	17.76	Find by Molecular Feature	564.33704

Compound Chromatograms



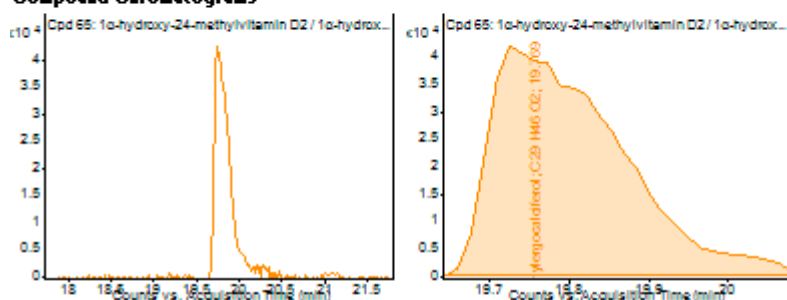
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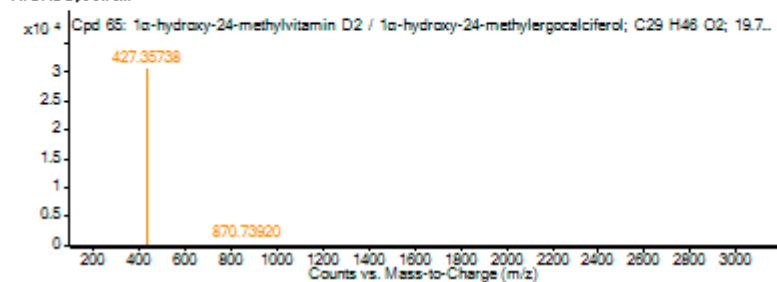
Compound 17: 1α-hydroxy-24-methylvitamin D2/1α-hydroxy-24-methylergocalciferol

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 65: 1α-hydroxy-24-methylvitamin D2 / 1α-hydroxy-24-methylergocalciferol; C29 H46 O2; 19.763	1α-hydroxy-24-methylvitamin D2 / 1α-hydroxy-24-methylergocalciferol	427.35738	19.77	Find by Molecular Feature	426.34966

Compound Chromatograms



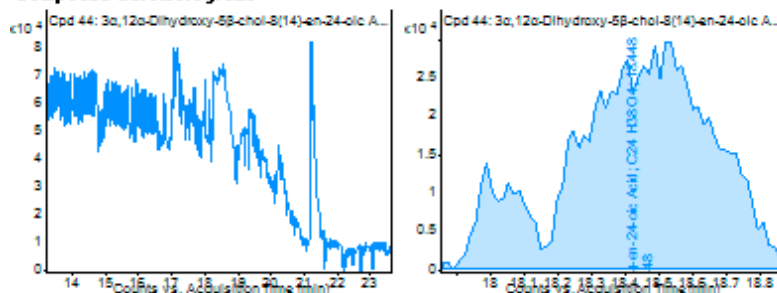
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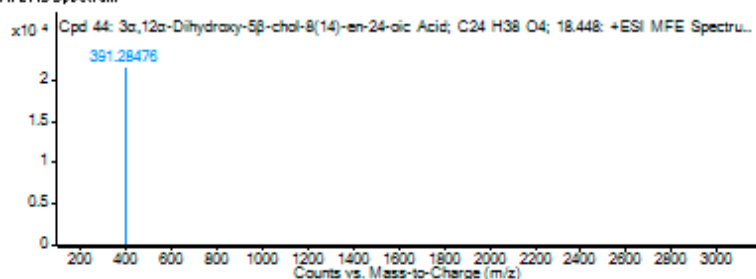
Compound 18: Coprocholic acid; 3 α ,12 α -Dihydroxy-5 β -chol-8(14)-en-24-oic Acid

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 44: 3 α ,12 α -Dihydroxy-5 β -chol-8(14)-en-24-oic Acid; C24 H38 O4; 18.448	3α,12α-Dihydroxy-5β-chol-8(14)-en-24-oic Acid	391.28476	18.45	Find by Molecular Feature	390.27804

Compound Chromatograms



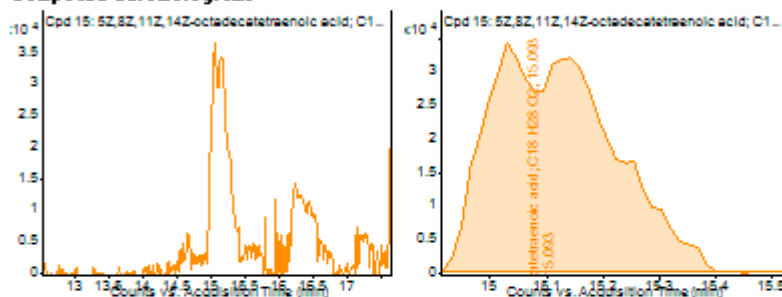
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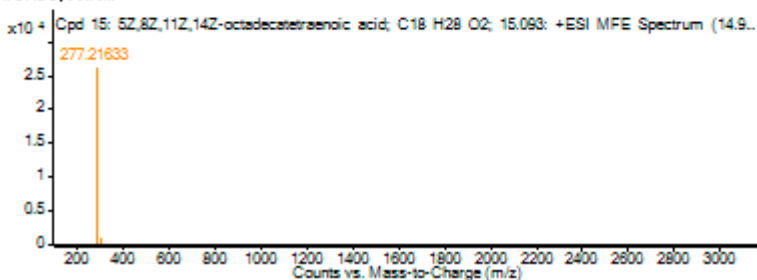
Compound 19: 9Z, 11Z, 13E, 15E-Octadecatetraenoic acid

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 15: 5Z,8Z,11Z,14Z-octadecatetraenoic acid; C18 H28 O2; 15.093	5Z,8Z,11Z,14Z-octadecatetraenoic acid	277.21633	15.09	Find by Molecular Feature	276.20918

Compound Chromatograms



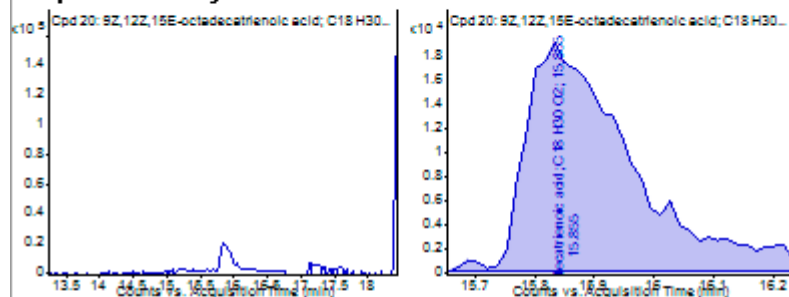
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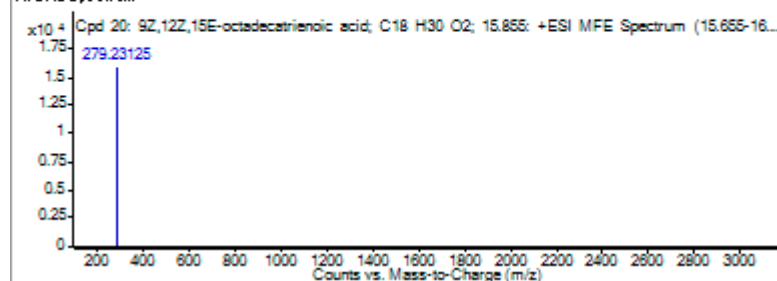
Compound 20: 9Z, 12Z, 15E-octadecatrienoic acid

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 20: 9Z,12Z,15E-octadecatrienoic acid; C18 H30 O2; 15.855	9Z,12Z,15E-octadecatrienoic acid	279.23125	15.86	Find by Molecular Feature	278.22405

Compound Chromatograms



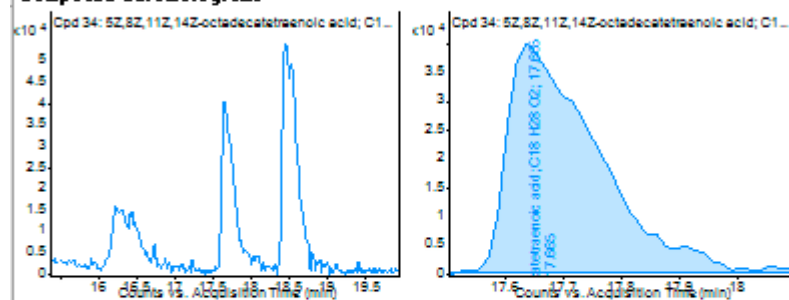
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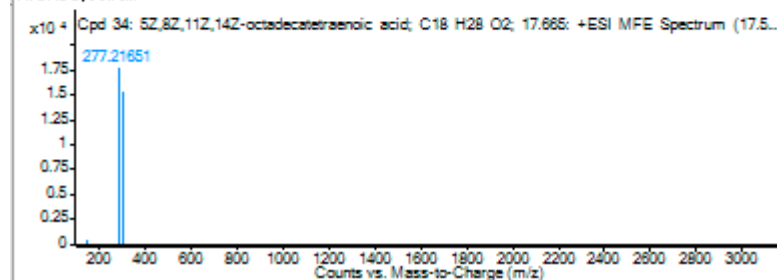
Compound 21: 5Z, 8Z, 11Z, 14Z-octadecatetraenoic acid

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 34: 5Z,8Z,11Z,14Z-octadecatetraenoic acid; C18 H28 O2; 17.665	5Z,8Z,11Z,14Z-octadecatetraenoic acid	277.21651	17.67	Find by Molecular Feature	276.2094

Compound Chromatograms



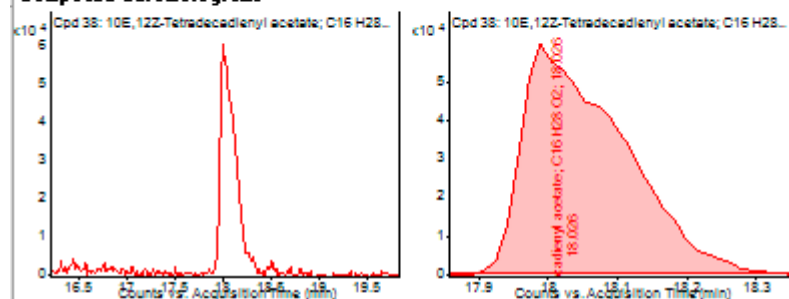
MFEMS Spectrum



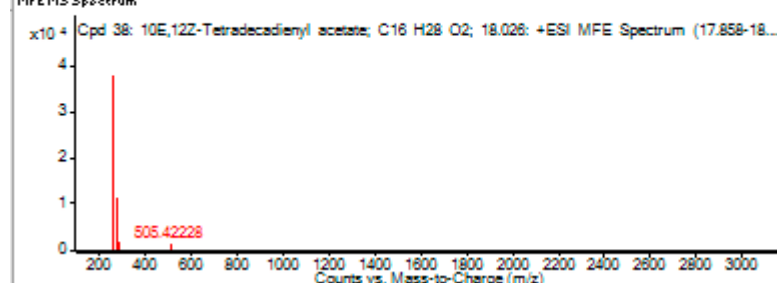
Compound 22: 10E, 12Z-tetradecadienyl acetate

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 38: 10E,12Z-Tetradecadienyl acetate; C16 H28 O2; 18.026	10E,12Z-Tetradecadienyl acetate	253.21626	18.03	Find by Molecular Feature	252.20921

Compound Chromatograms



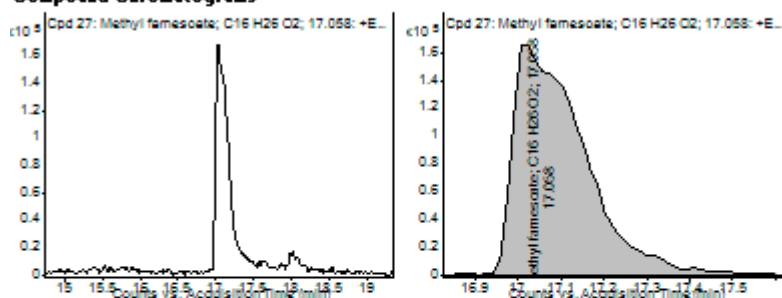
MFEMS Spectrum



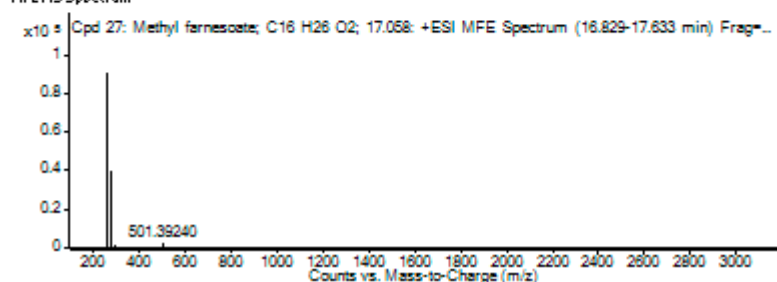
Compound 23: Methyl farnesoate

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 27: Methyl farnesoate; C16 H26 O2; 17.058	Methyl farnesoate	251.2006	17.06	Find by Molecular Feature	250.19341

Compound Chromatograms



MFEMS Spectrum



Supplementary Material Figure S3: Mass spectra of individual compounds listed in Supplementary Material Table S1.