

Treating hyperglycemia from *Eryngium caeruleum* M. Bieb: *In-vitro* α -glucosidase, antioxidant, *in-vivo* antidiabetic and molecular docking-based approaches

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Supporting Information-II

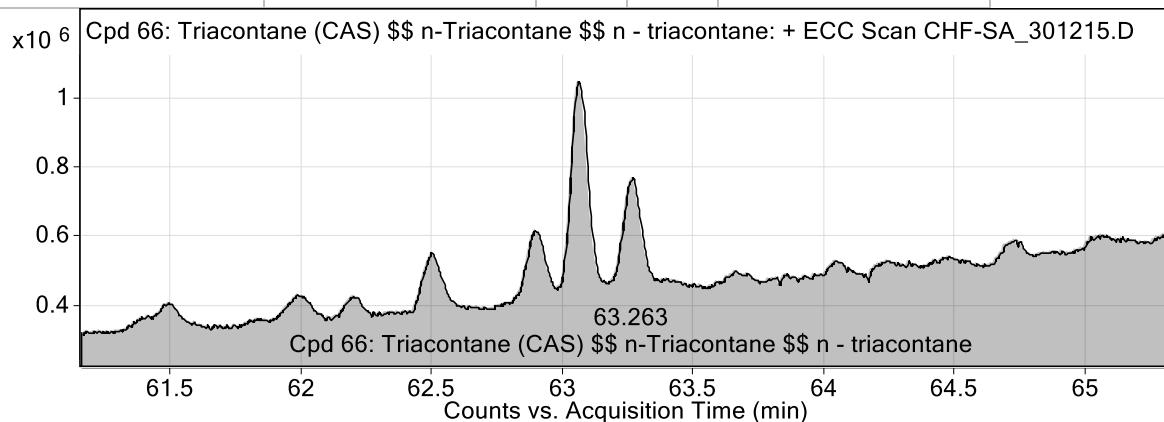
(Bioactive compounds)

GC-MS spectra/details of the compounds

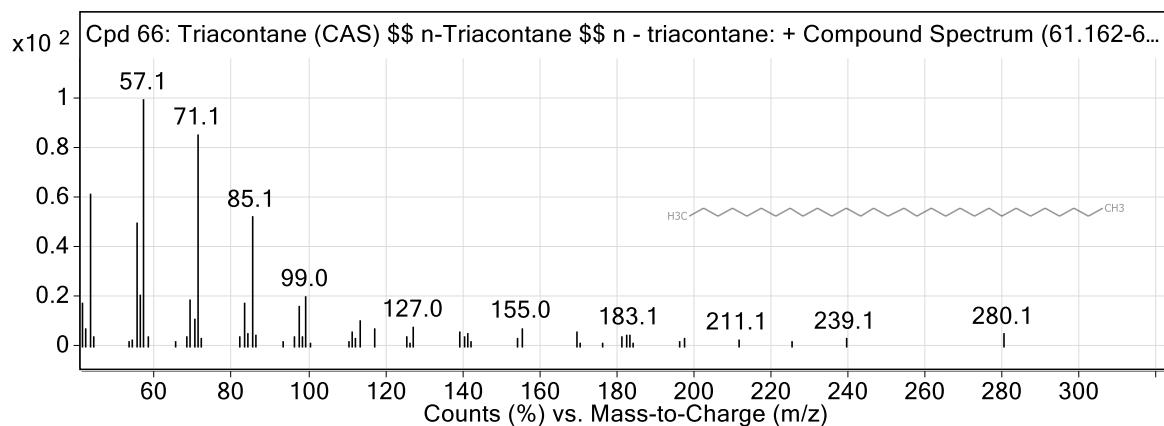


Triacontane

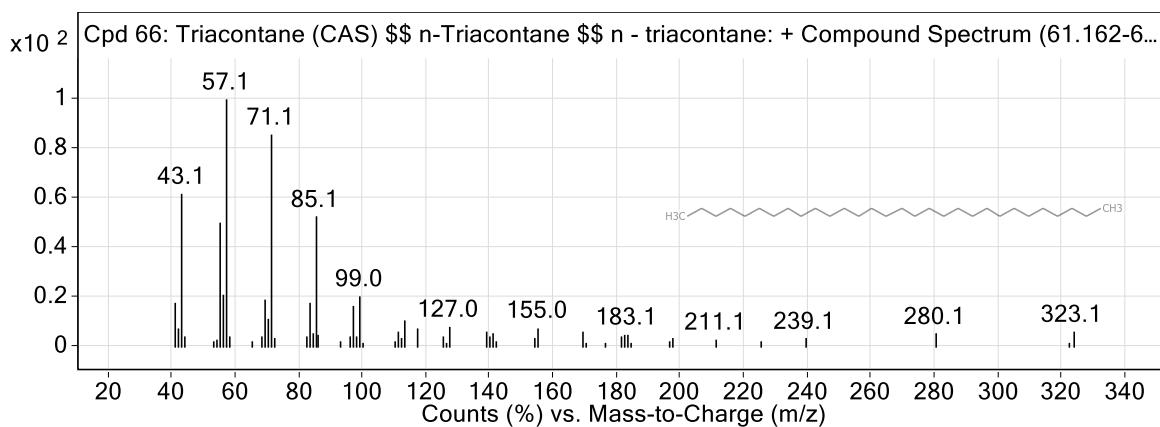
Compound Label	Name	m/z	RT	Algorithm
Cpd 66: Triacontane (CAS) \$\$ n-Triacontane \$\$ n - triacontane	Triacontane (CAS) \$\$ n-Triacontane \$\$ n - triacontane	57,1	63,26	Find by Chromatogram Deconvolution



MS Spectrum



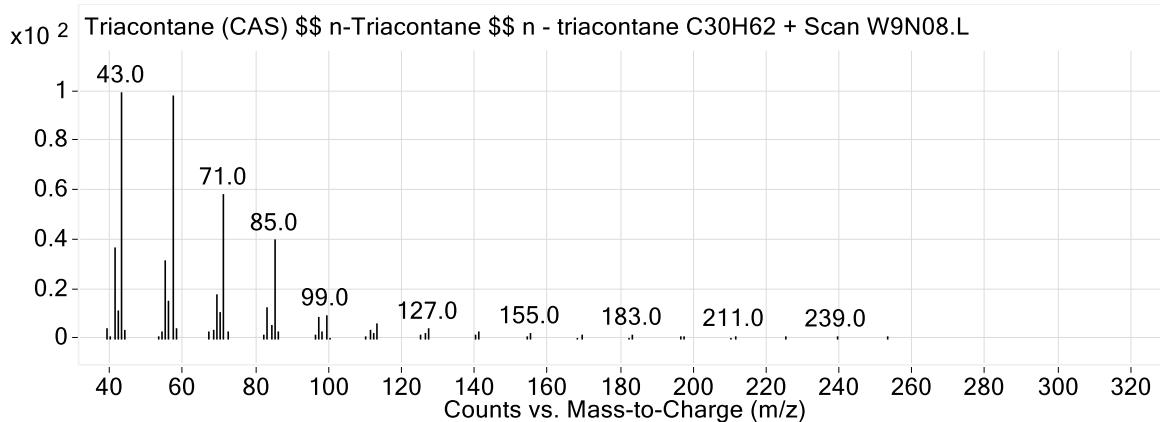
MS Zoomed Spectrum



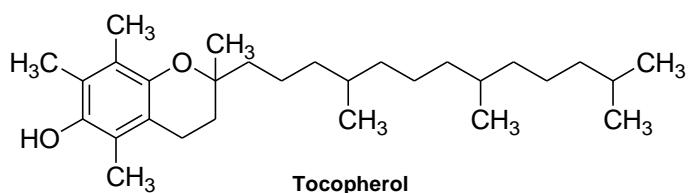
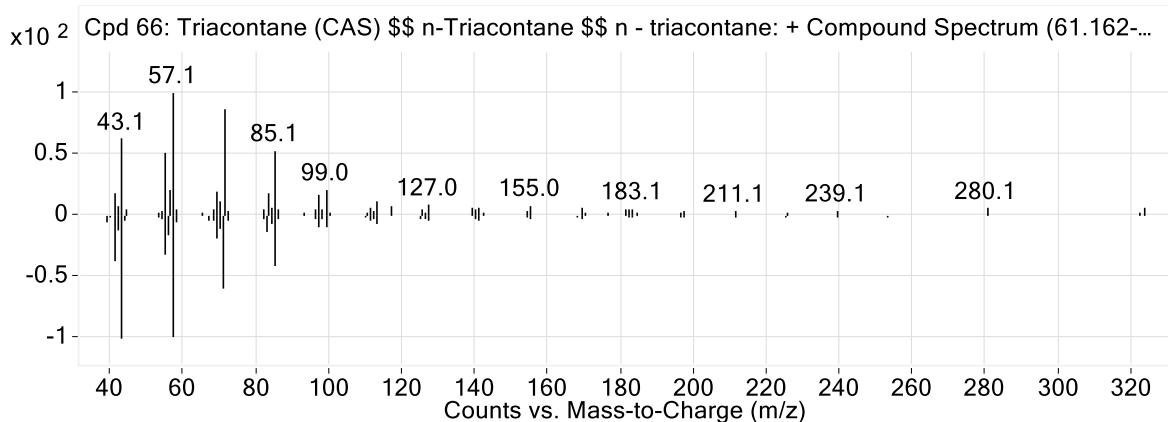
MS Spectrum Peak List

m/z	Abund
41,1	10892
43,1	37206
55,1	30142
56,1	12737
57,1	59929
69,1	11574
71,1	51468
83,1	10932
85,1	31629
99	12534

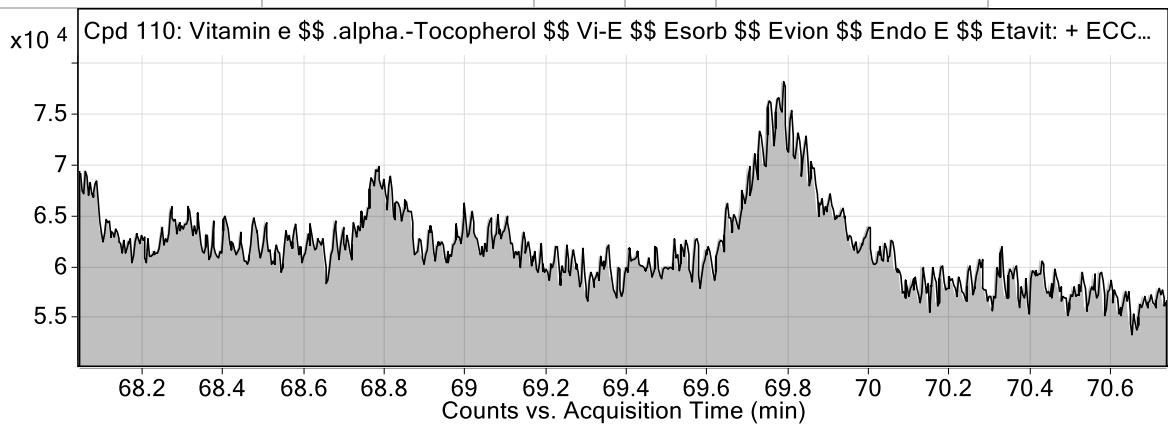
Library Spectrum



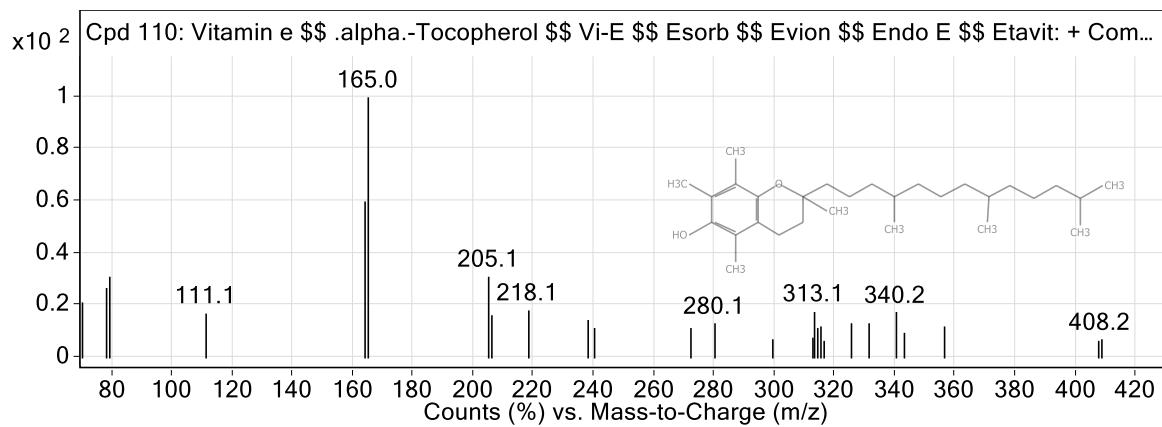
Difference Spectrum



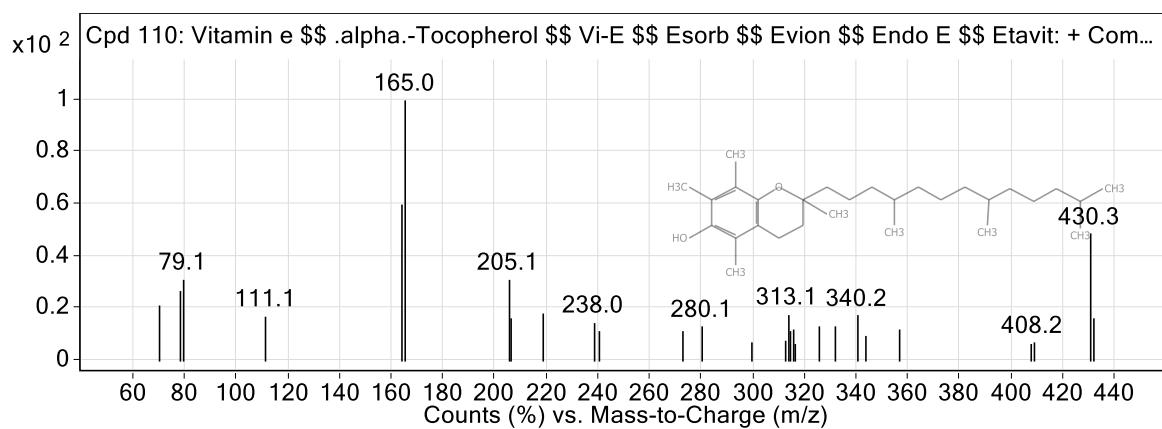
Compound Label	Name	m/z	RT	Algorithm
Cpd 110: Vitamin e \$\$.alpha.-Tocopherol \$\$ Vi-E \$\$ Esorb \$\$ Evion \$\$ Endo E \$\$ Etavit	Vitamin e \$\$.alpha.-Tocopherol \$\$ Vi-E \$\$ Esorb \$\$ Evion \$\$ Endo E \$\$ Etavit	165	69,78	Find by Chromatogram Deconvolution



MS Spectrum



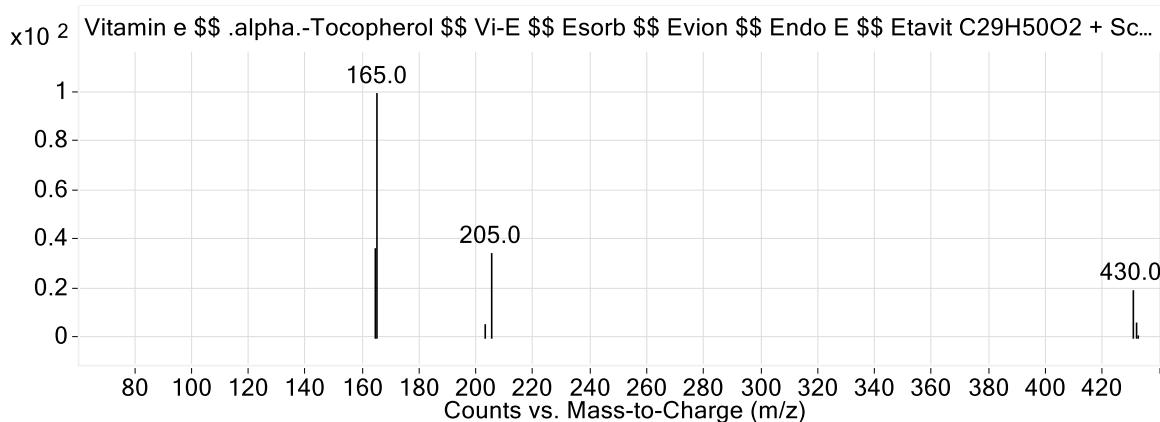
MS Zoomed Spectrum



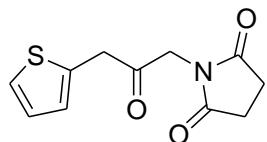
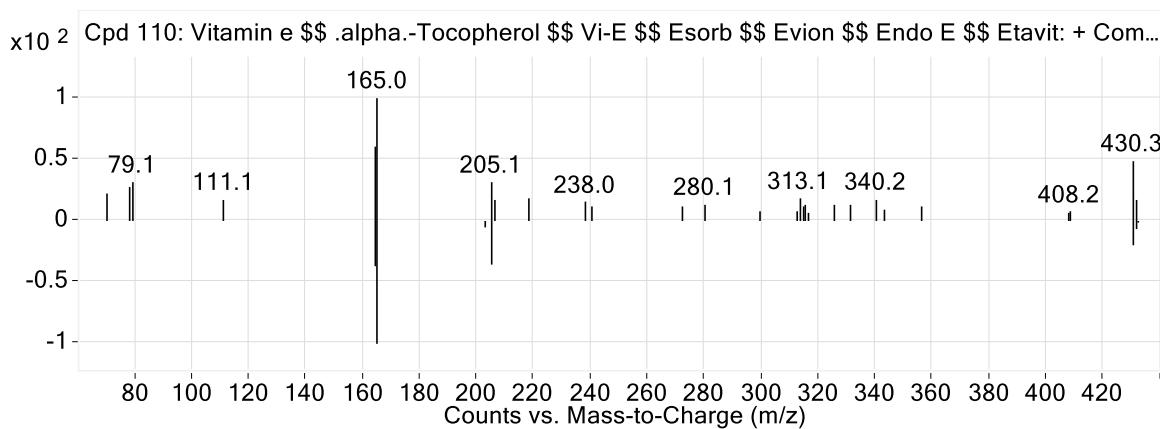
MS Spectrum Peak List

m/z	Abund
70,1	1737
78	2165,5
79,1	2521,9
164	4798,3
165	8035,8
205,1	2516,6
218,1	1478,3
313,1	1448
340,2	1401,5
430,3	3914,6

Library Spectrum

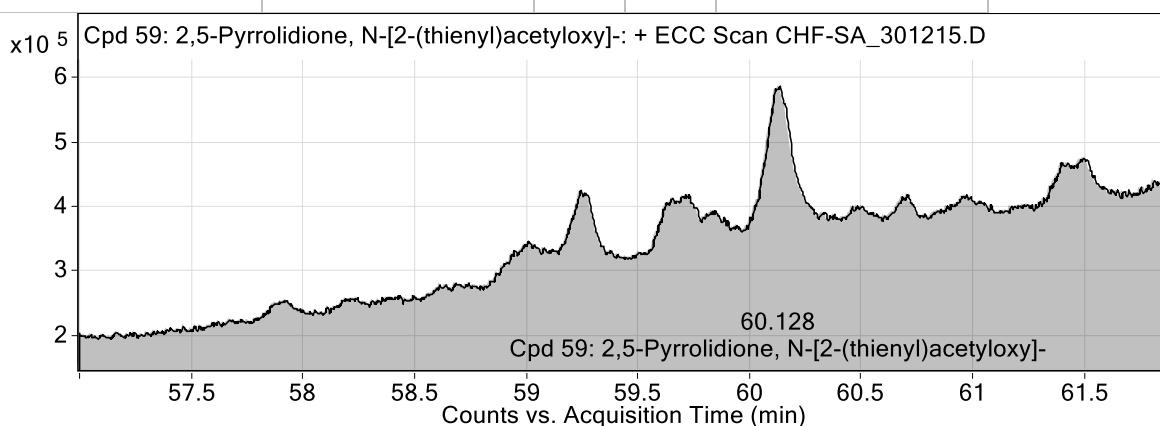


Difference Spectrum

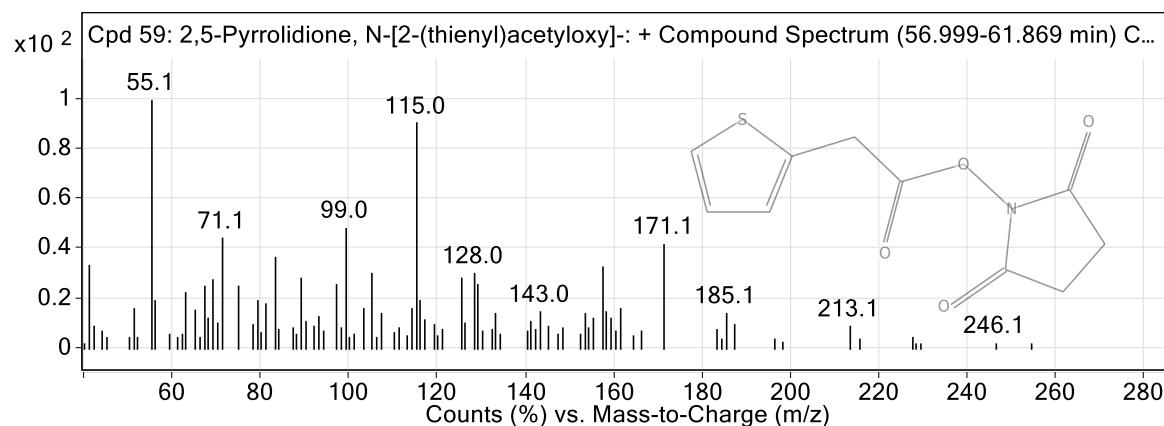


2,5-Pyrrolidione, N-[2-(thienyl)acetoxy]

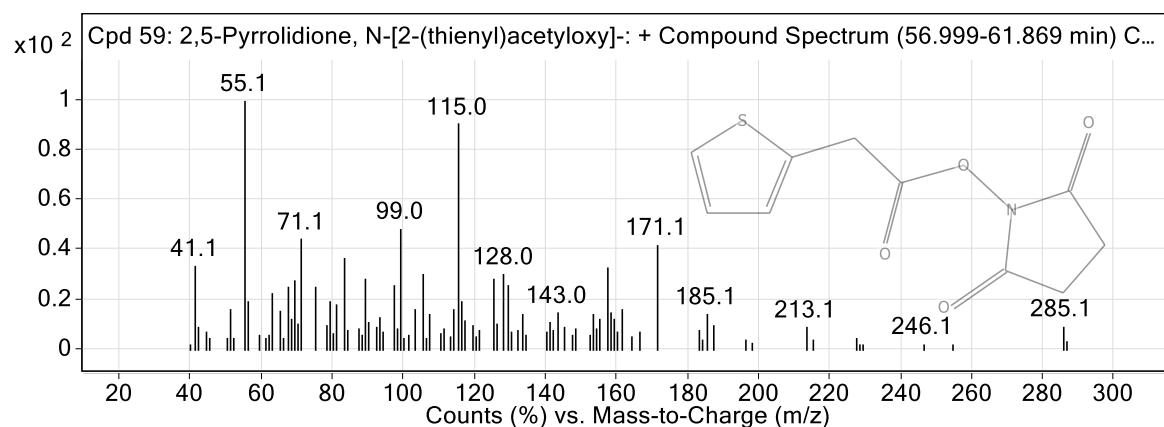
Compound Label	Name	m/z	RT	Algorithm
Cpd 59: 2,5-Pyrrolidione, N-[2-(thienyl)acetoxy]-	2,5-Pyrrolidione, N-[2-(thienyl)acetoxy]-	55,1	60,13	Find by Chromatogram Deconvolution



MS Spectrum



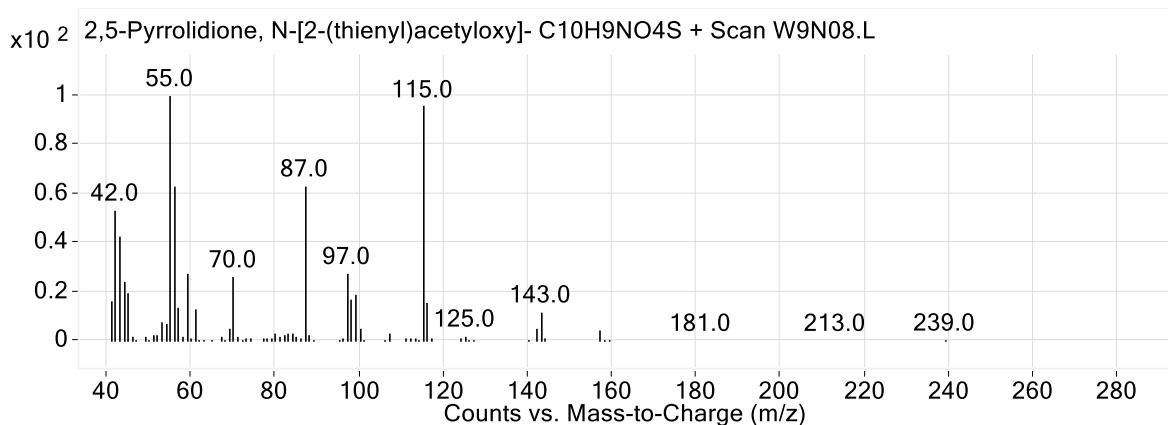
MS Zoomed Spectrum



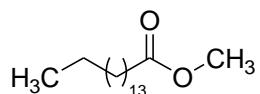
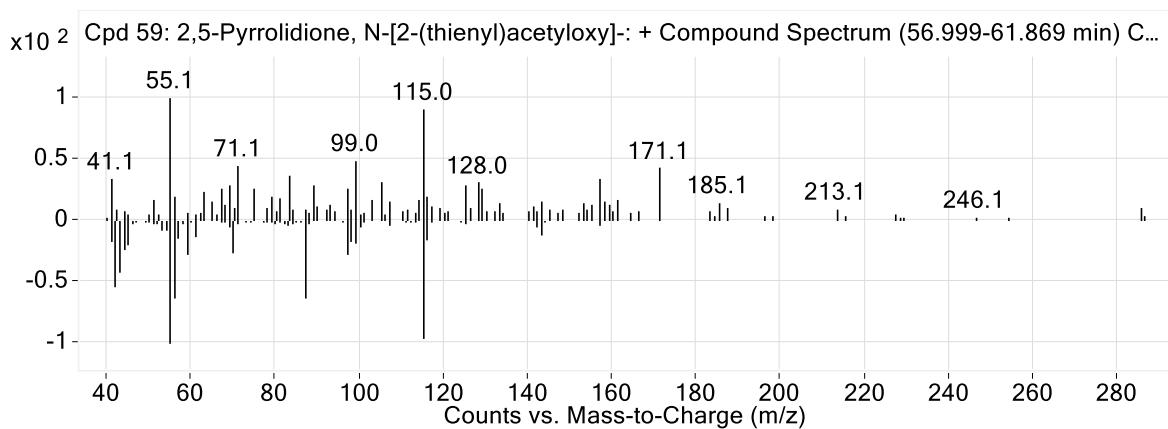
MS Spectrum Peak List

m/z	Abund
41,1	7298,5
55,1	21500
71,1	9583,6
83,1	7919,3
99	10504
105	6645,8
115	19472
128	6584
157	7194,6
171,1	9118,9

Library Spectrum

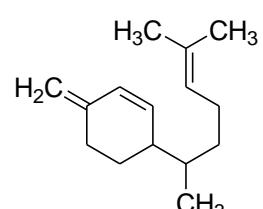


Difference Spectrum



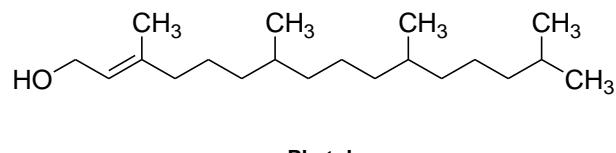
Methyl palmitate

This is a repeated compound. For details please see the supporting information of Compound (K) in Figure 2.

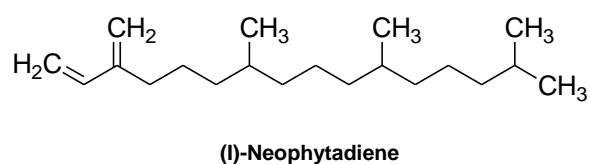


Cyclohexene, 3-(1,5-dimethyl-4-hexenyl)-6-methylene

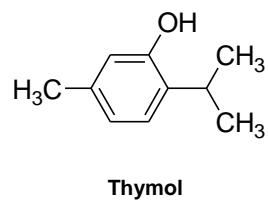
This is a repeated compound. For details please see the supporting information of Compound (C) in Figure 2.



This is a repeated compound. For details please see the supporting information of Compound (J) in Figure 2.

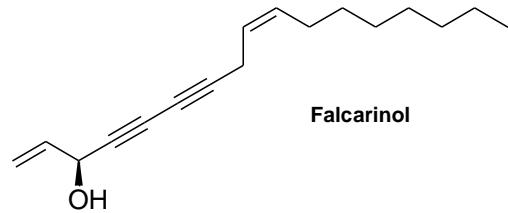


This is a repeated compound. For details please see the supporting information of Compound (I) in Figure 2.

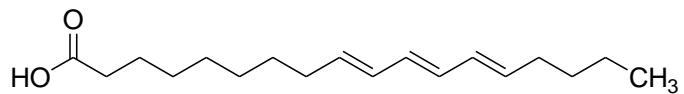


Thymol

This is a repeated compound. For details please see the supporting information of Compound (B) in Figure 2.

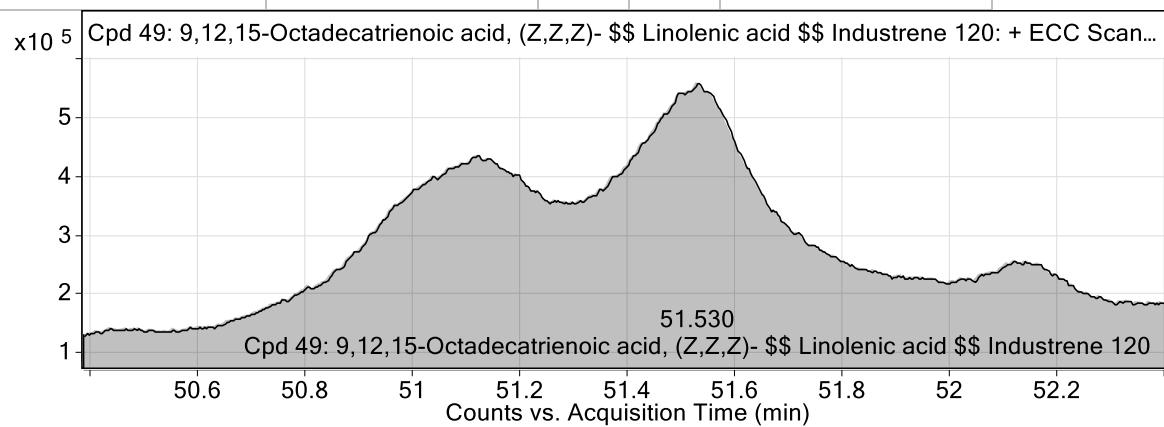


This is a repeated compound. For details please see the supporting information of Compound (M) in Figure 2.

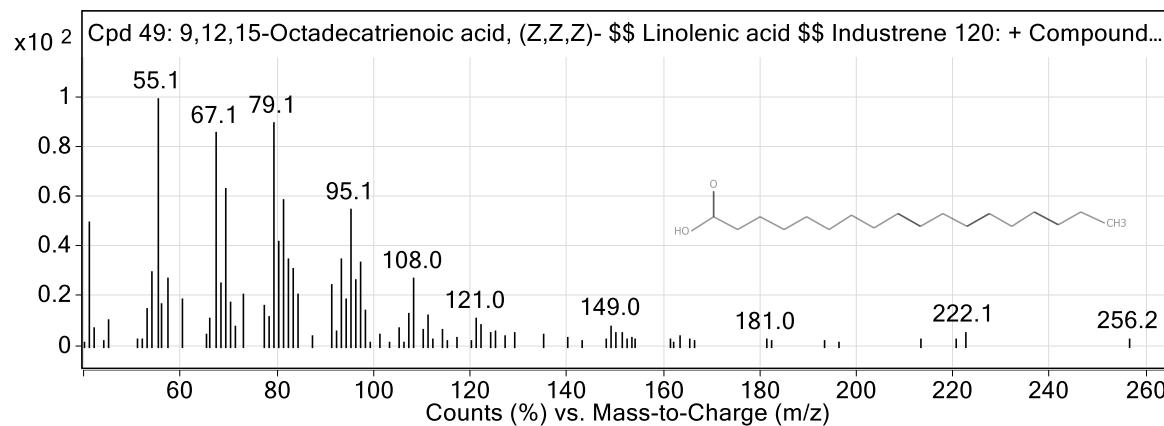


Linolenic acid

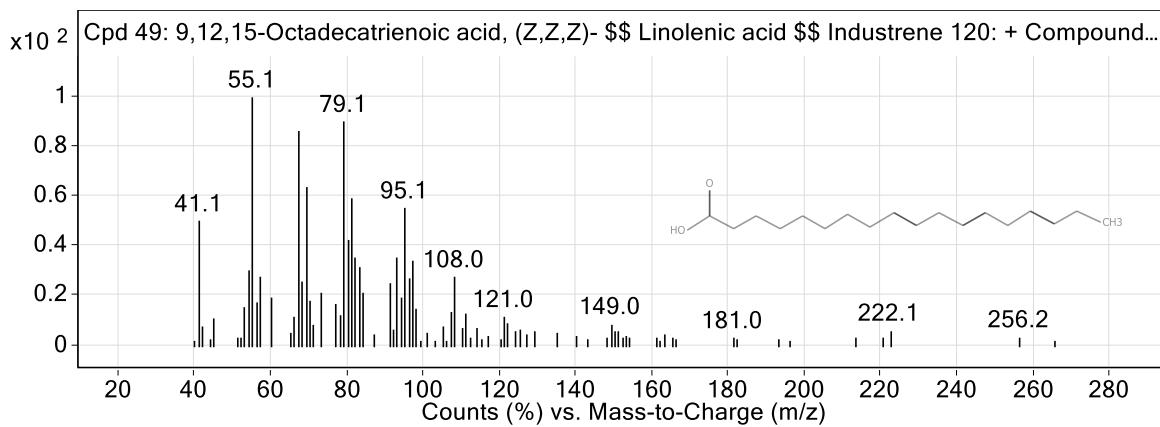
Compound Label	Name	m/z	RT	Algorithm
Cpd 49: 9,12,15-Octadecatrienoic acid, (Z,Z,Z)- \$\$ Linolenic acid \$\$ Industrene 120	9,12,15-Octadecatrienoic acid, (Z,Z,Z)- \$\$ Linolenic acid \$\$ Industrene 120	55,1	51,53	Find by Chromatogram Deconvolution



MS Spectrum



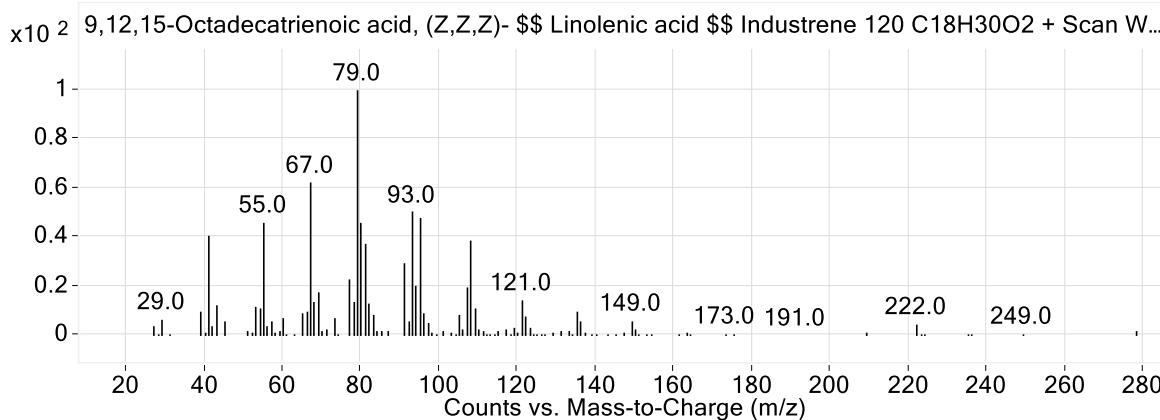
MS Zoomed Spectrum



MS Spectrum Peak List

m/z	Abund
41,1	15854
55,1	31558
67,1	27172
69,1	20050
79,1	28464
80	13457
81,1	18698
82,1	11214
93	11237
95,1	17501

Library Spectrum



Difference Spectrum

