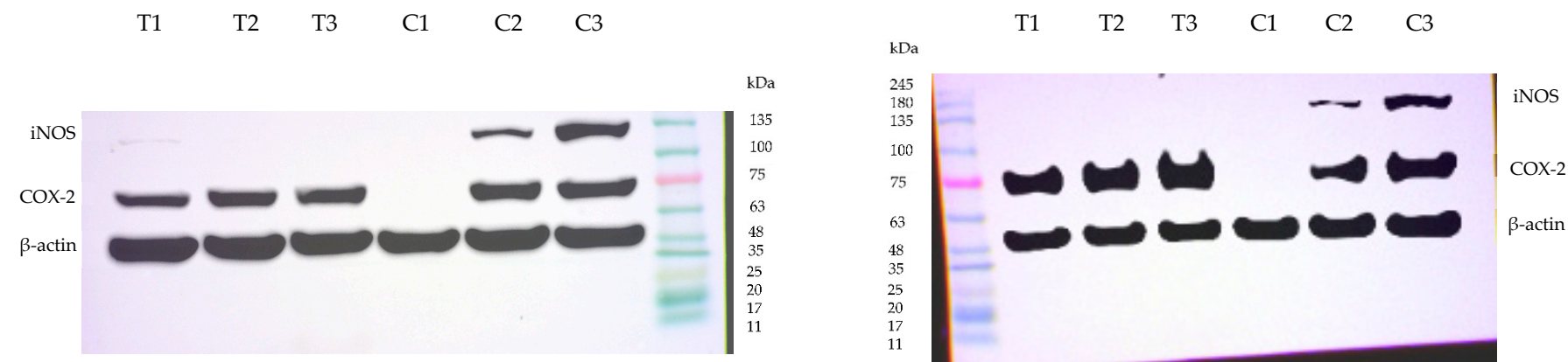
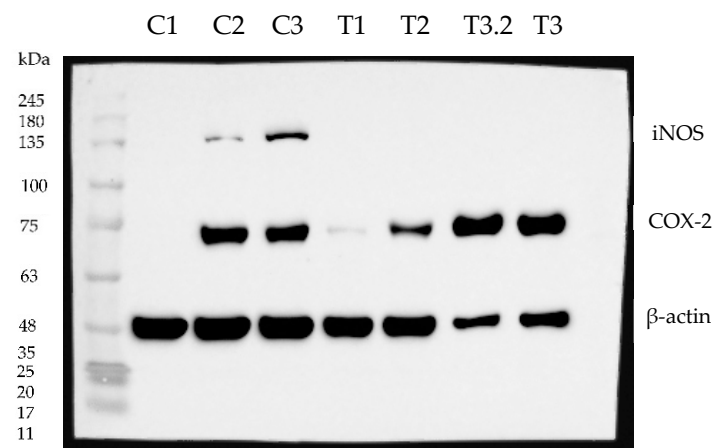


Figure S1: Original Western blot images for three repeats of iNOS, COX-2 and  $\beta$ -actin



Representative blot shown in main text (Repeat 1)



Repeat 3

Repeat 2

T1: *Spirulina platensis* ethanol extract 0.5 mg/mL  
T2: *Spirulina platensis* ethanol extract 1 mg/mL  
T3: *Spirulina platensis* ethanol extract 2 mg/mL  
C1: Negative control (Untreated)  
C2: Lipopolysaccharide (LPS) control (Stimulated with 1  $\mu$ g/mL LPS)

**Excluded in the main text:**

C3: Positive control (Stimulated with 1  $\mu$ g/mL LPS + treated with N( $\gamma$ )-nitro-L-arginine methyl ester (L-NAME))  
T3.2: Replicate of *Spirulina platensis* ethanol extract 2 mg/mL

**Molecular weight of the target proteins:**

iNOS – 130 kDa  
COX-2 – 74 kDa  
 $\beta$ -actin – 45 kDa

Figure S2: Chromatogram of *S. platensis* ethanol extract

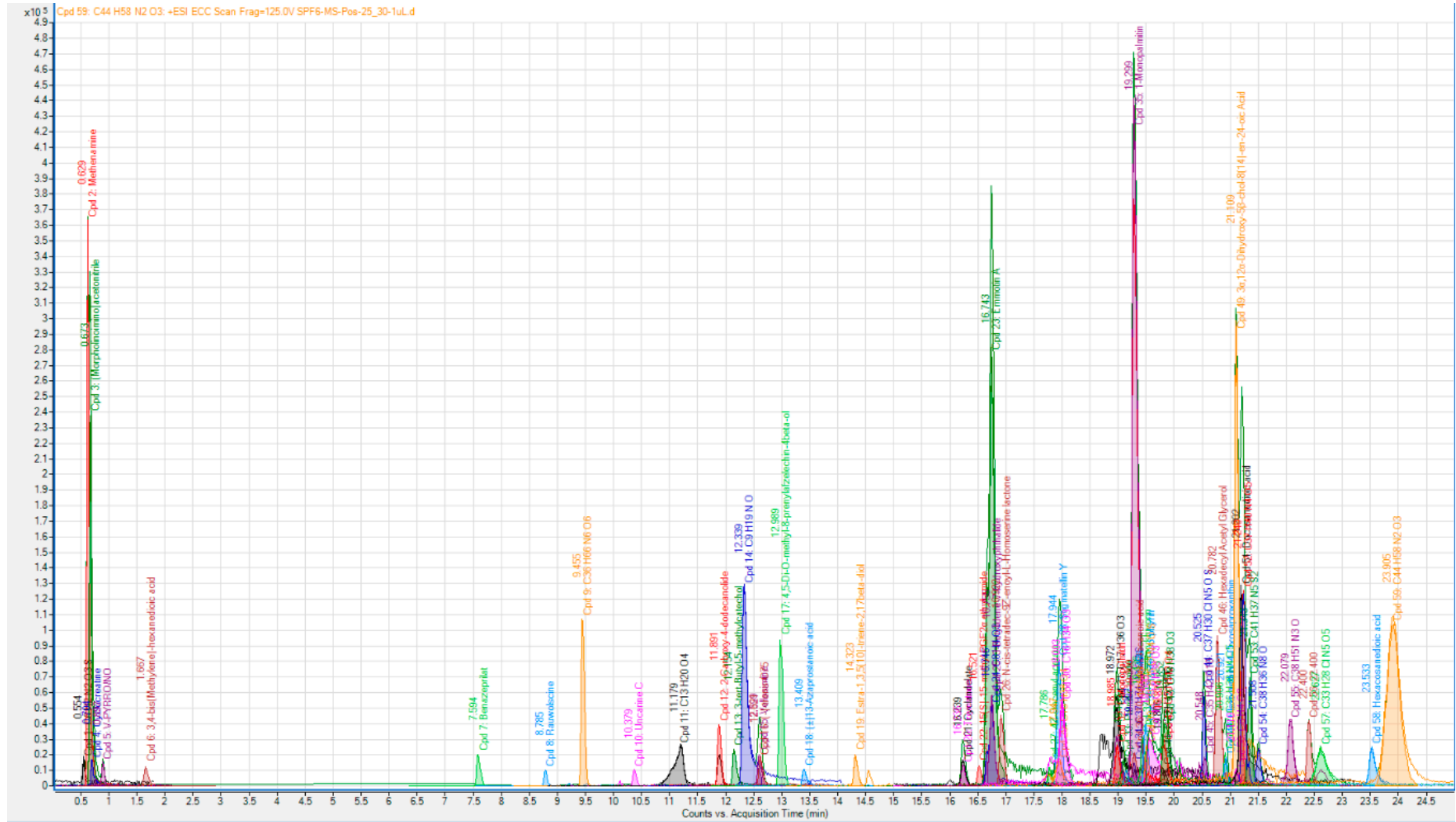
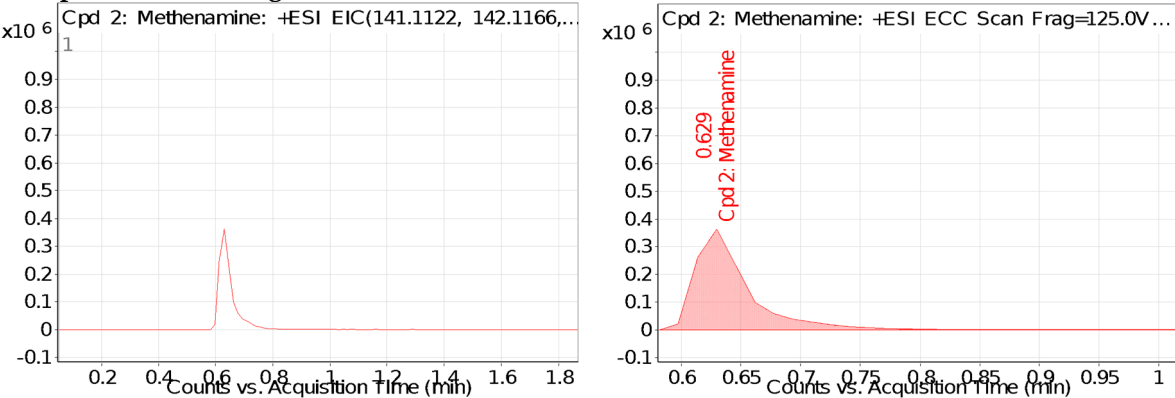


Figure S3: Mass spectra of individual compounds

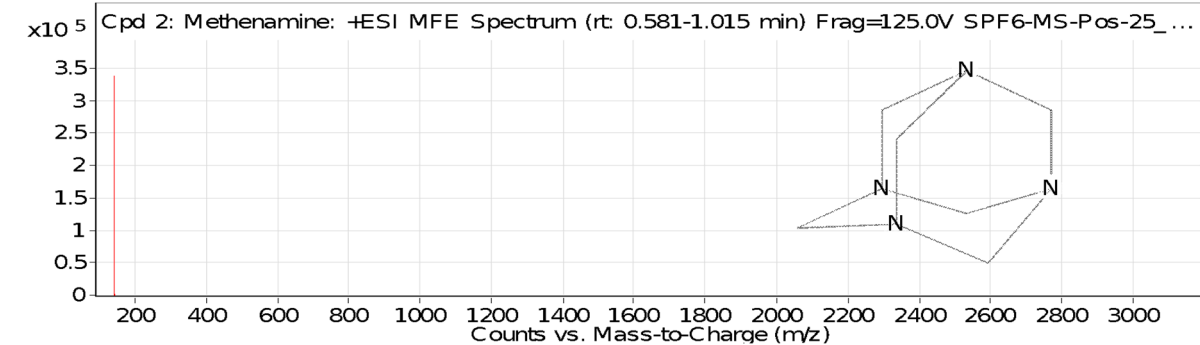
Compound i: Methenamine

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 2: Methenamine	Methenamine	141.1132	0.629	Find by Molecular Feature	140.106

Compound Chromatograms



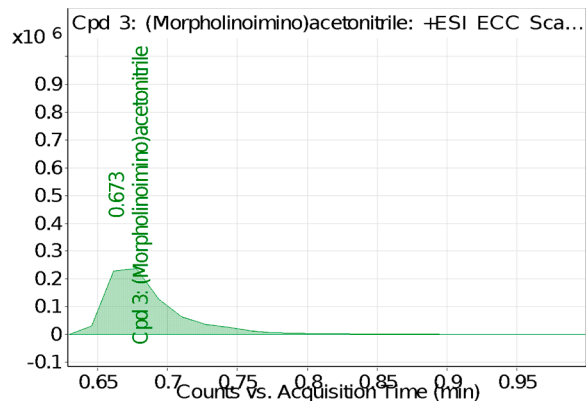
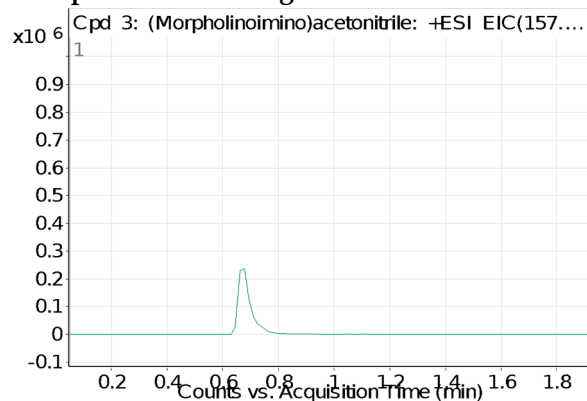
MFE MS Spectrum



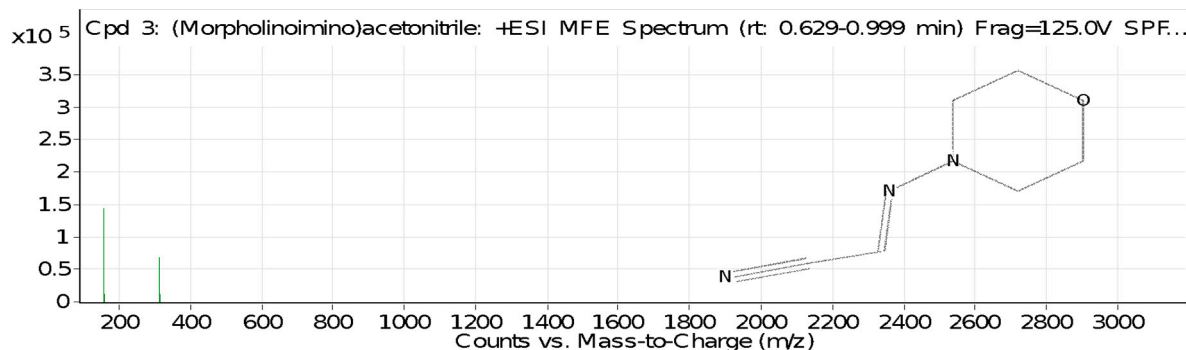
### Compound ii: Morpholinoiminoacetonitrile

Compound Label	Name	<i>m/z</i>	RT	Algorithm	Mass
Cpd 3: (Morpholinoimino)acetonitrile	(Morpholinoimino)acetonitrile	157.1081	0.673	Find by Molecular Feature	139.0742

### Compound Chromatograms



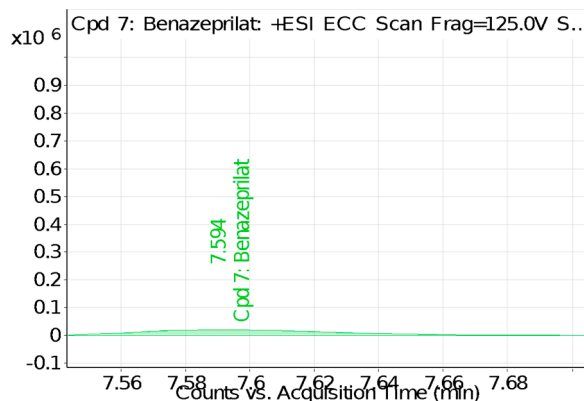
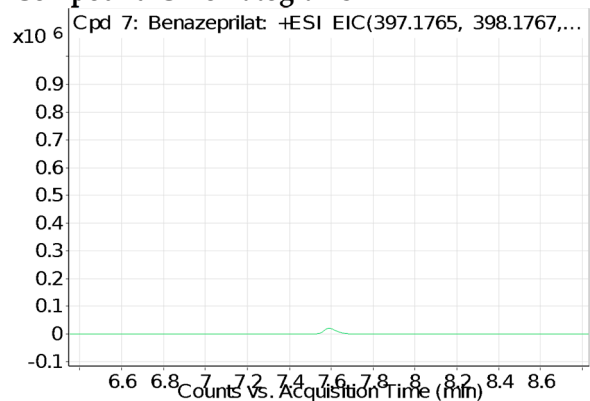
### MFE MS Spectrum



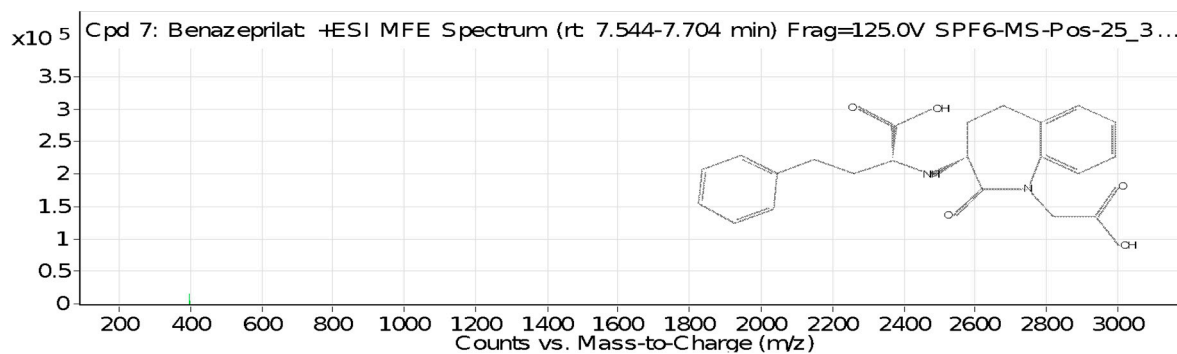
### Compound iii: Benazeprilat

Compound Label	Name	<i>m/z</i>	RT	Algorithm	Mass
Cpd 7: Benazeprilat	Benazeprilat	397.1761	7.594	Find by Molecular Feature	396.1687

### Compound Chromatograms



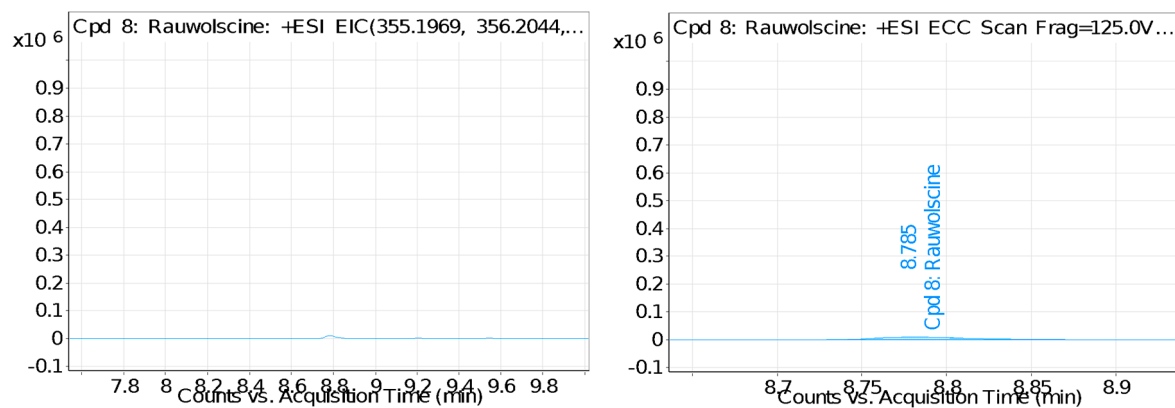
### MFE MS Spectrum



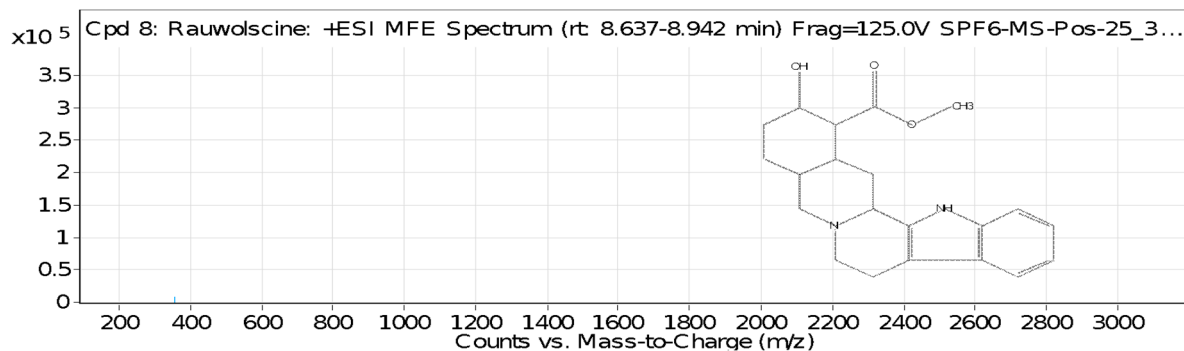
### Compound iv: Rauwolscline

Compound Label	Name	<i>m/z</i>	RT	Algorithm	Mass
Cpd 8: Rauwolscline	Rauwolscline	355.2013	8.785	Find by Molecular Feature	354.1938

### Compound Chromatograms



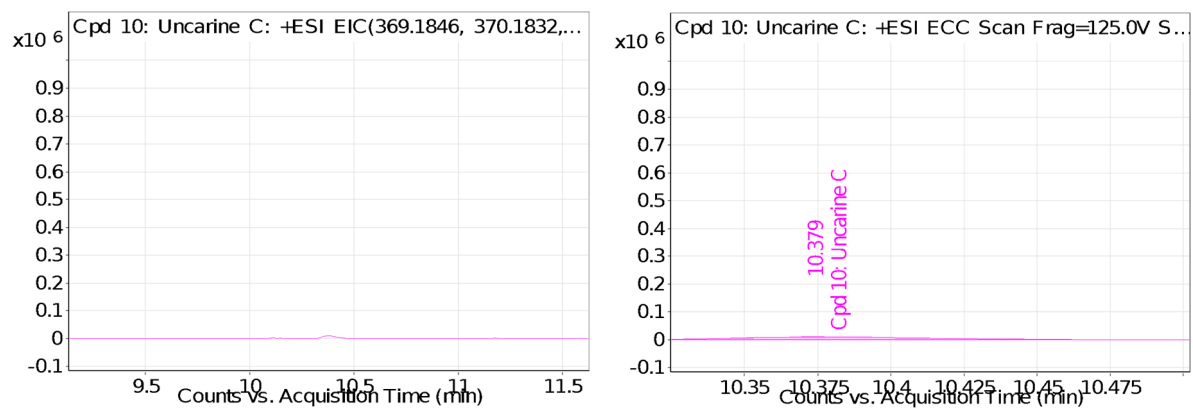
### MFE MS Spectrum



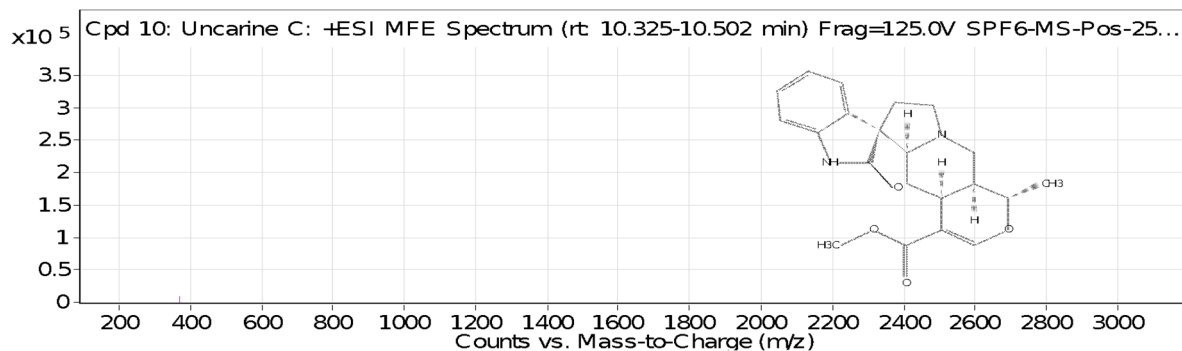
### Compound v: Uncarine C

Compound Label	Name	<i>m/z</i>	RT	Algorithm	Mass
Cpd 10: Uncarine C	Uncarine C	369.1806	10.379	Find by Molecular Feature	368.1734

### Compound Chromatograms



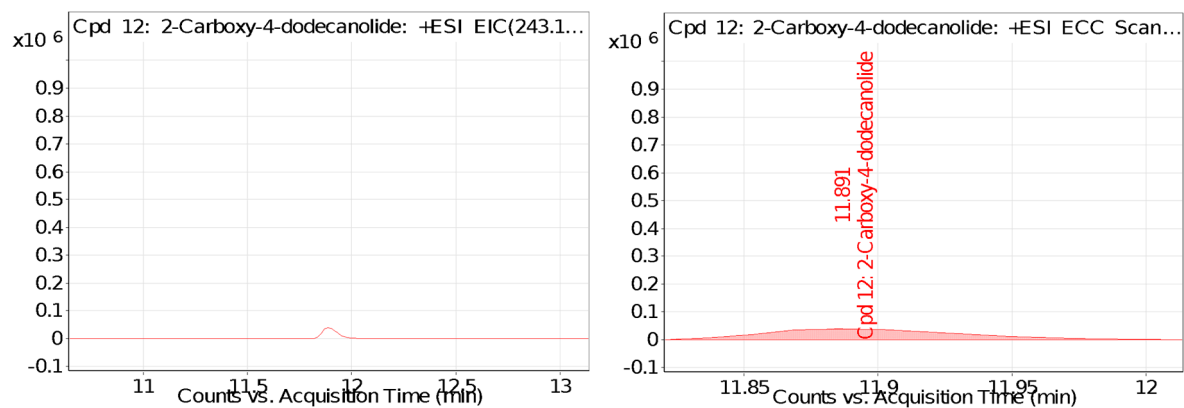
### MFE MS Spectrum



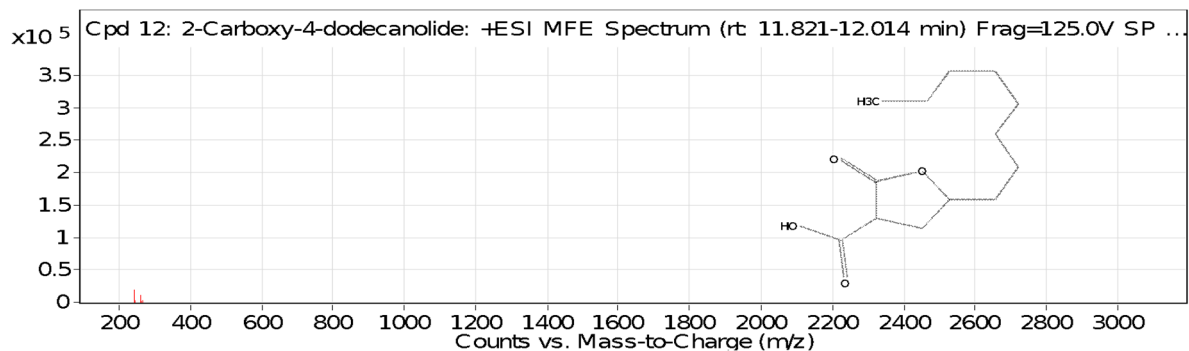
### Compound vi: 2-Carboxy-4-dodecanolide

Compound Label	Name	<i>m/z</i>	RT	Algorithm	Mass
Cpd 12: 2-Carboxy-4-dodecanolide	2-Carboxy-4-dodecanolide	243.1593	11.891	Find by Molecular Feature	242.152

### Compound Chromatograms



### MFE MS Spectrum

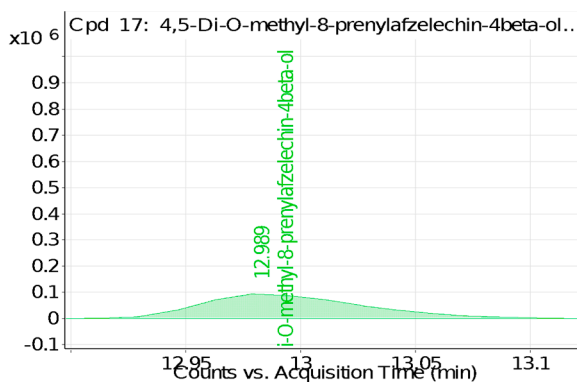
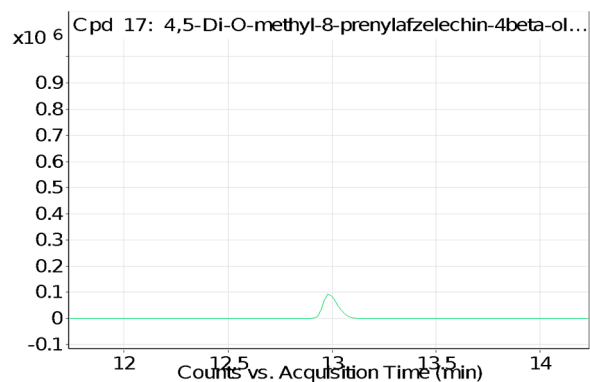




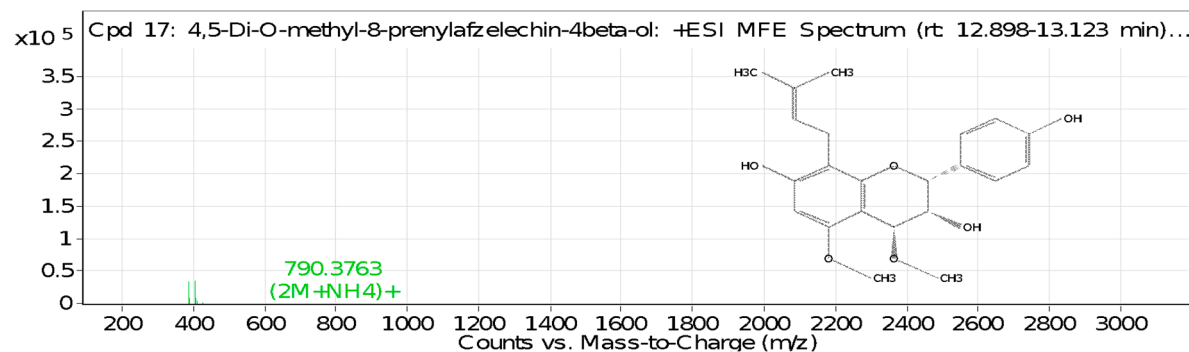
**Compound vii: 4,5-Di-O-methyl-8-prenylafzelechin-4beta-ol**

Compound Label	Name	<i>m/z</i>	RT	Algorithm	Mass
Cpd 17: 4,5-Di-O-methyl-8-prenylafzelechin-4beta-ol	<b>4,5-Di-O-methyl-8-prenylafzelechin-4beta-ol</b>	404.2063	12.989	Find by Molecular Feature	386.1734

**Compound Chromatograms**



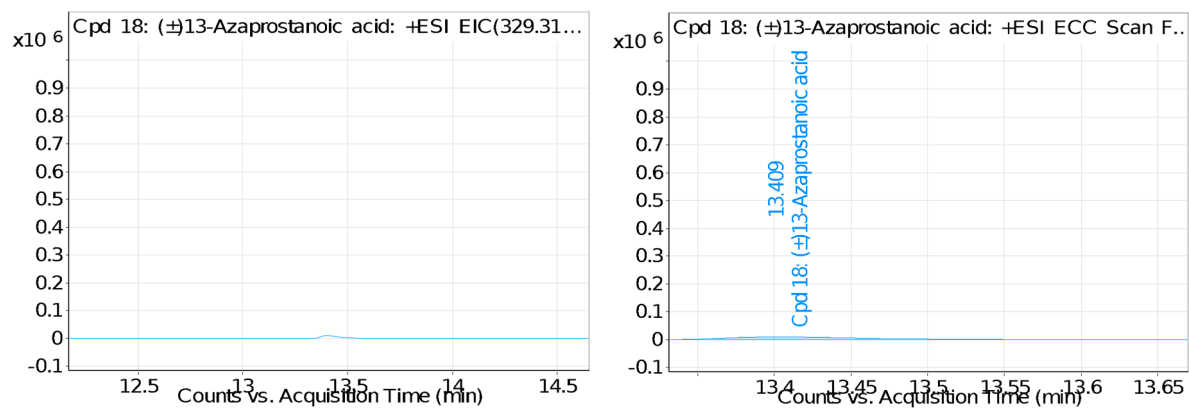
MFE MS Spectrum



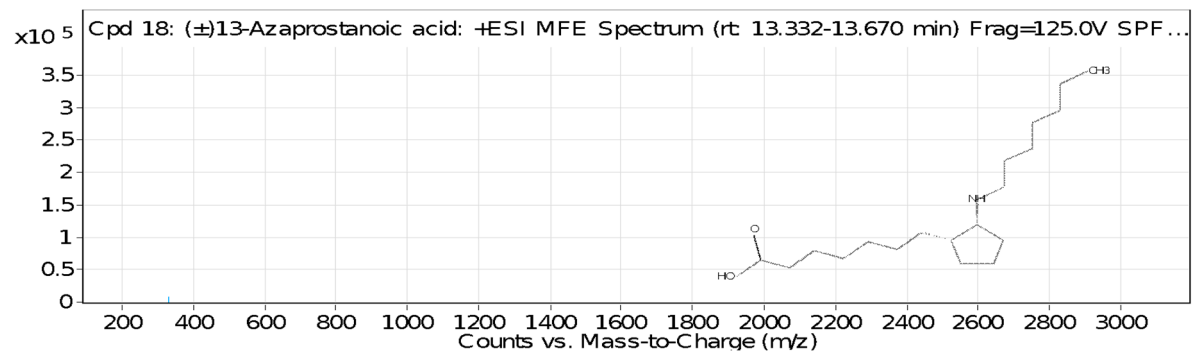
### Compound viii: (±)13-Azaprostanoic acid

Compound Label	Name	<i>m/z</i>	RT	Algorithm	Mass
Cpd 18: (±)13-Azaprostanoic acid	(±)13-Azaprostanoic acid	329.3163	13.409	Find by Molecular Feature	311.2822

### Compound Chromatograms



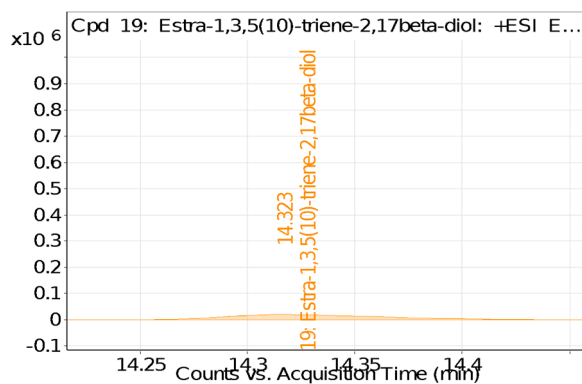
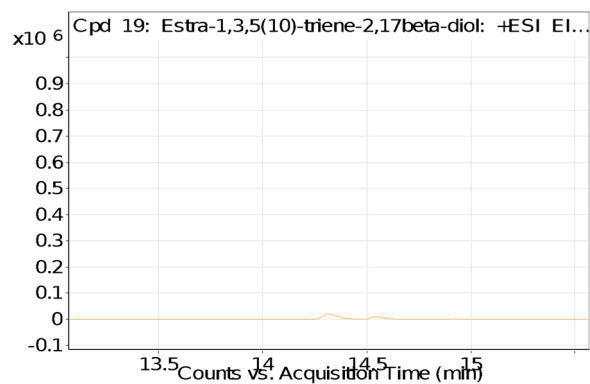
### MFE MS Spectrum



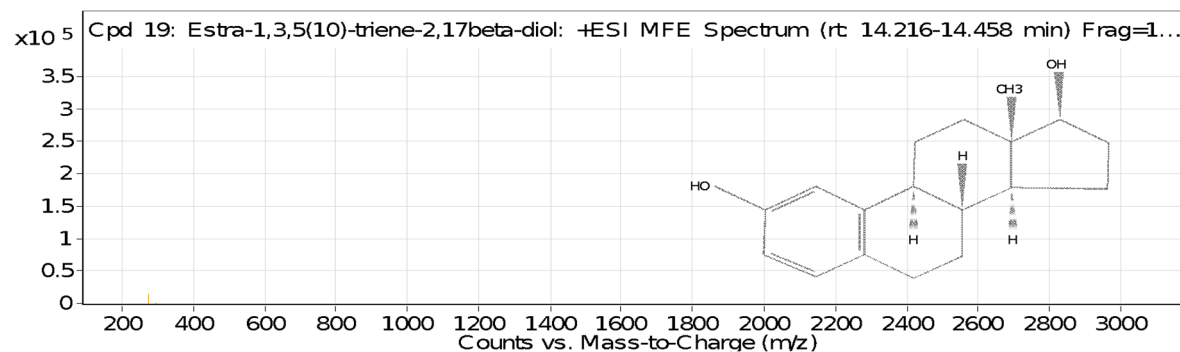
### Compound ix: Estra-1,3,5(10)-triene-2,17beta-diol

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 19: Estra-1,3,5(10)-triene-2,17beta-diol	<b>Estra-1,3,5(10)-triene-2,17beta-diol</b>	273.1848	14.323	Find by Molecular Feature	272.1774

### Compound Chromatograms



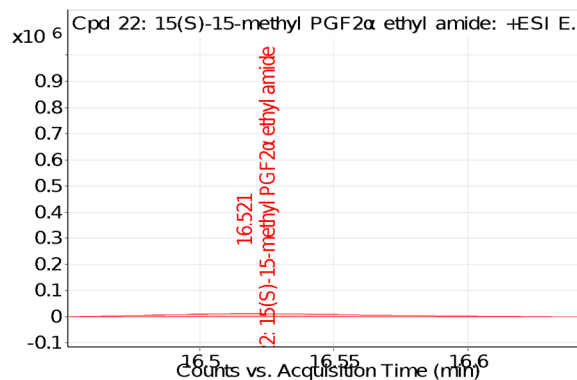
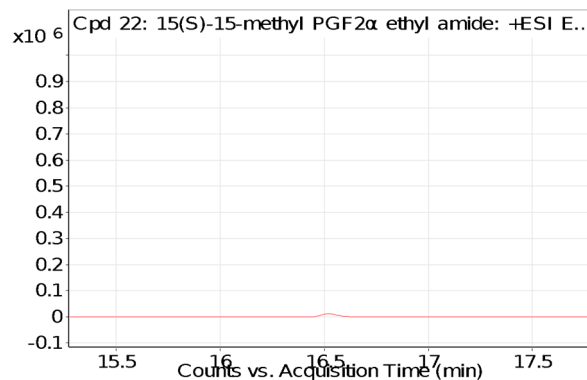
### MFE MS Spectrum



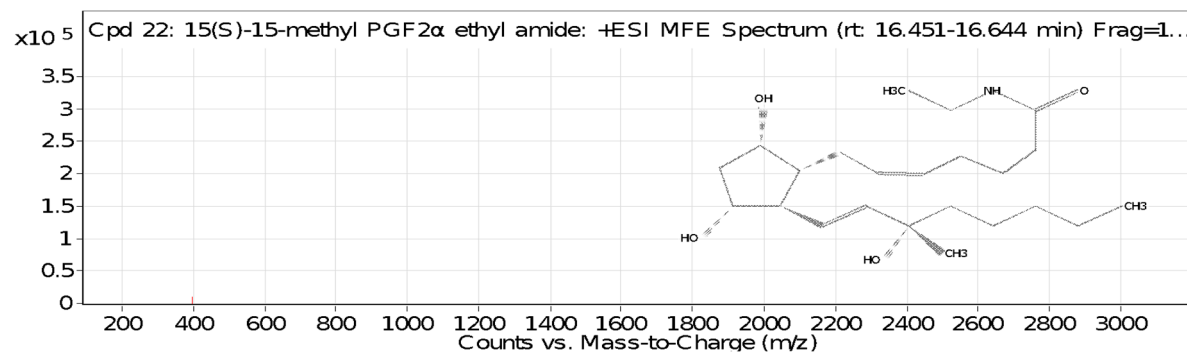
**Compound x: 15(S)-15-methyl PGF2 $\alpha$  ethyl amide**

Compound Label	Name	<i>m/z</i>	RT	Algorithm	Mass
Cpd 22: 15(S)-15-methyl PGF2 $\alpha$ ethyl amide	15(S)-15-methyl PGF2 $\alpha$ ethyl amide	396.3106	16.521	Find by Molecular Feature	395.3029

**Compound Chromatograms**



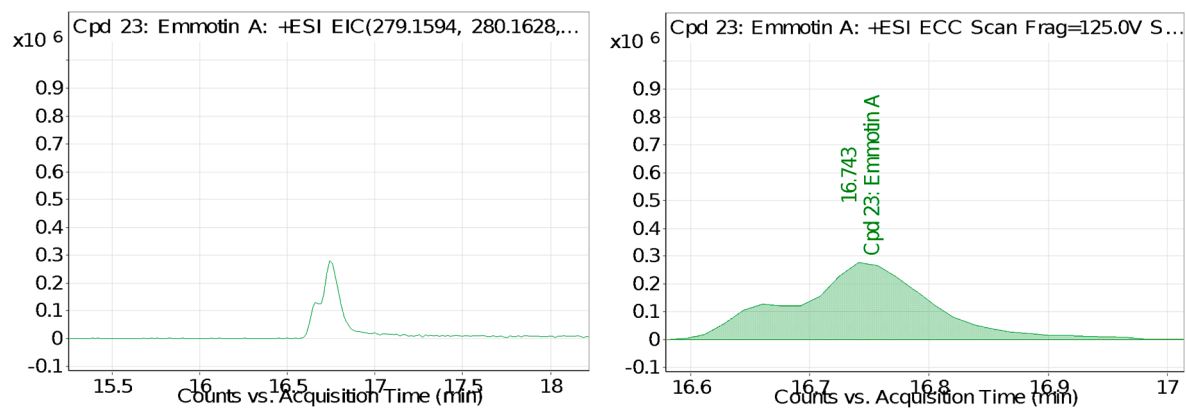
**MFE MS Spectrum**



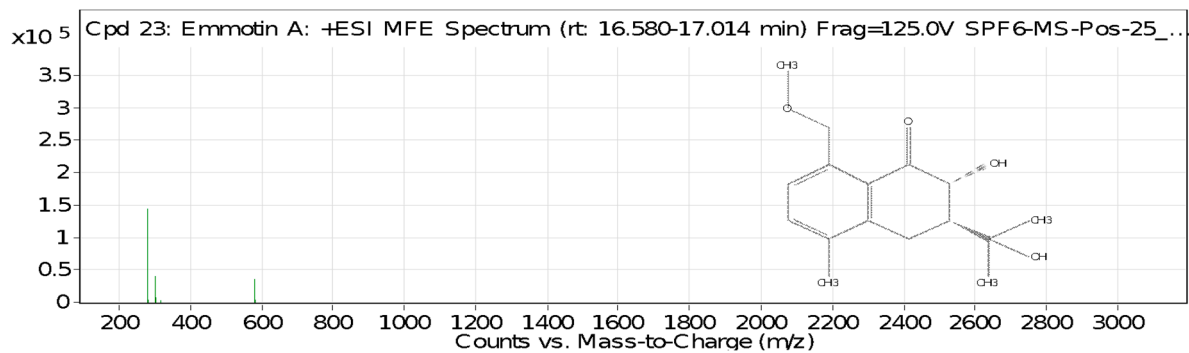
### Compound xi: Emmotin A

Compound Label	Name	<i>m/z</i>	RT	Algorithm	Mass
Cpd 23: Emmotin A	<b>Emmotin A</b>	279.1594	16.743	Find by Molecular Feature	278.1521

### Compound Chromatograms



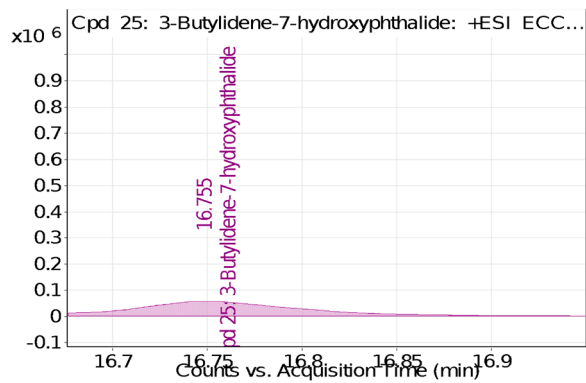
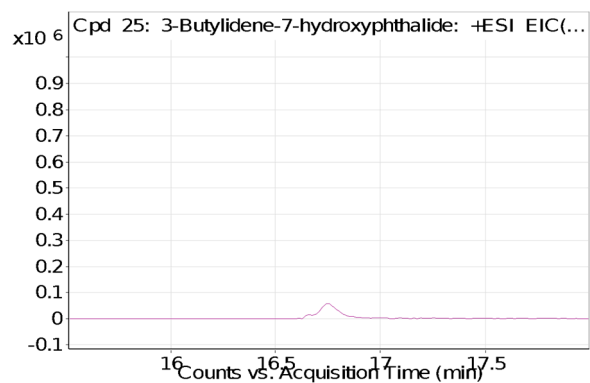
### MFE MS Spectrum



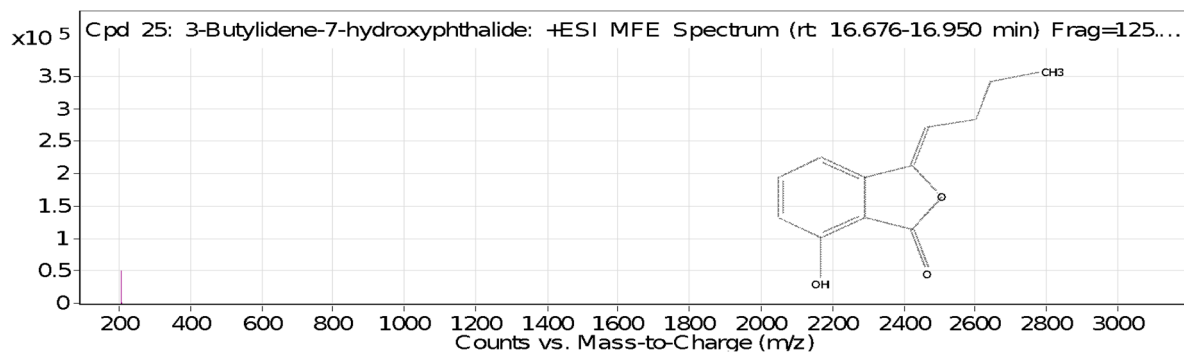
### Compound xii: 3-Butylidene-7-hydroxyphthalide

Compound Label	Name	<i>m/z</i>	RT	Algorithm	Mass
Cpd 25: 3-Butylidene-7-hydroxyphthalide	3-Butylidene-7-hydroxyphthalide	205.0858	16.755	Find by Molecular Feature	204.0786

### Compound Chromatograms



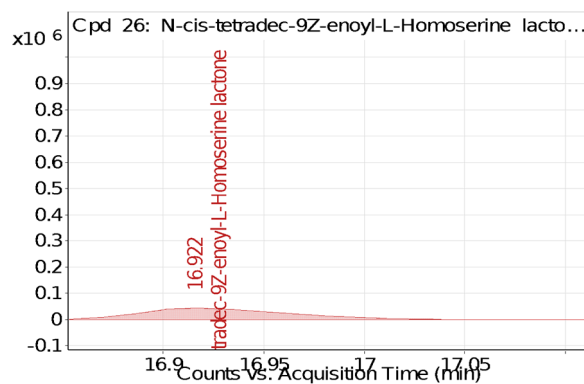
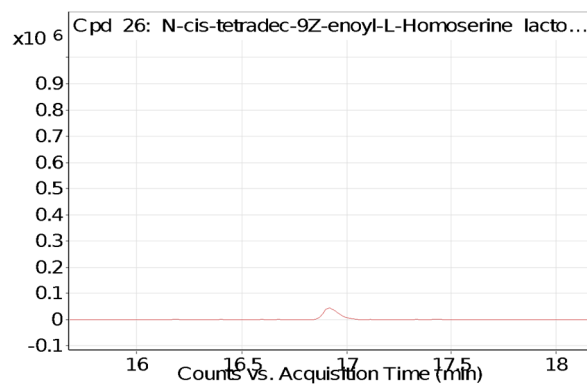
### MFE MS Spectrum



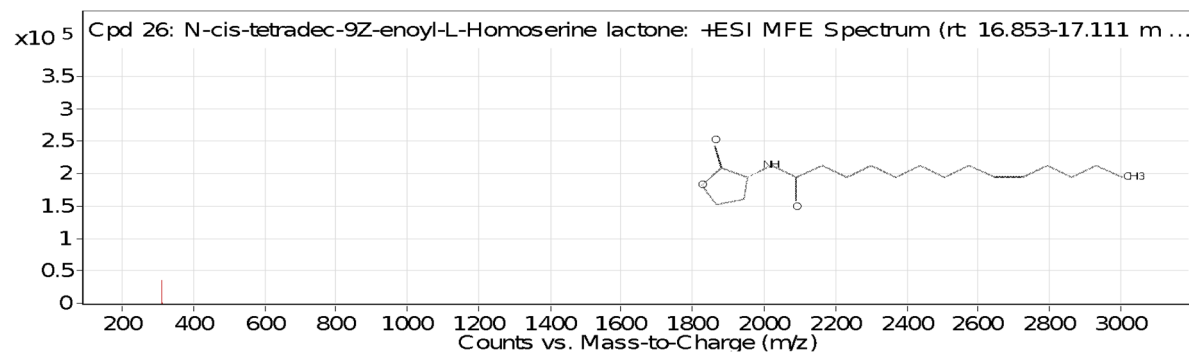
**Compound xiii: N-cis-tetradec-9Z-enoyl-L-Homoserine lactone**

Compound Label	Name	<i>m/z</i>	RT	Algorithm	Mass
Cpd 26: N-cis-tetradec-9Z-enoyl-L-Homoserine lactone	<b>N-cis-tetradec-9Z-enoyl-L-Homoserine lactone</b>	310.2374	16.922	Find by Molecular Feature	309.2303

**Compound Chromatograms**



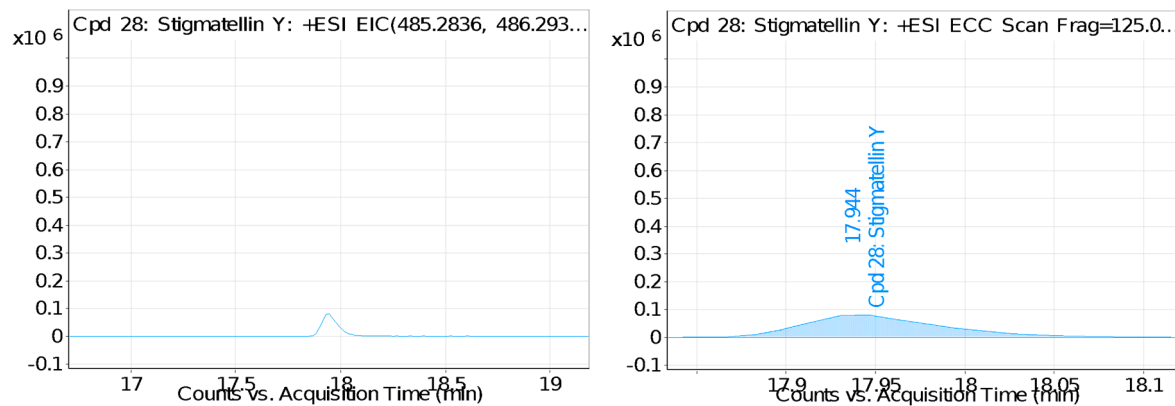
**MFE MS Spectrum**



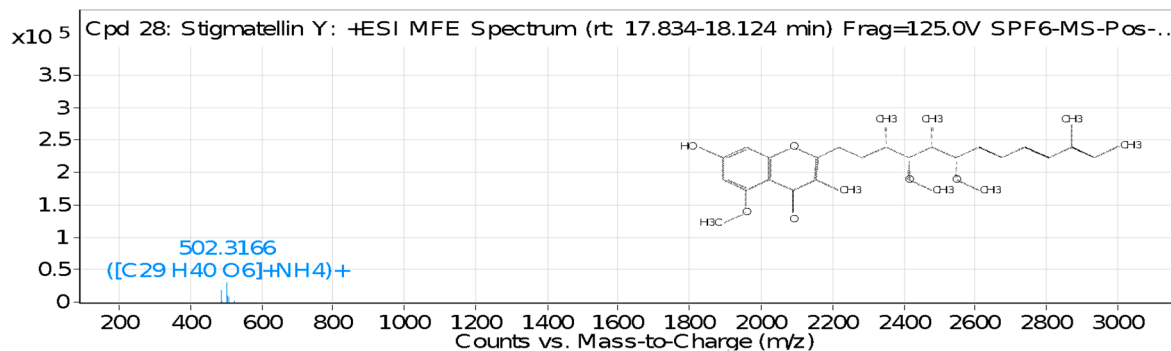
### Compound xiv: Stigmatellin Y

Compound Label	Name	<i>m/z</i>	RT	Algorithm	Mass
Cpd 28: Stigmatellin Y	Stigmatellin Y	502.3166	17.944	Find by Molecular Feature	484.2826

### Compound Chromatograms



### MFE MS Spectrum

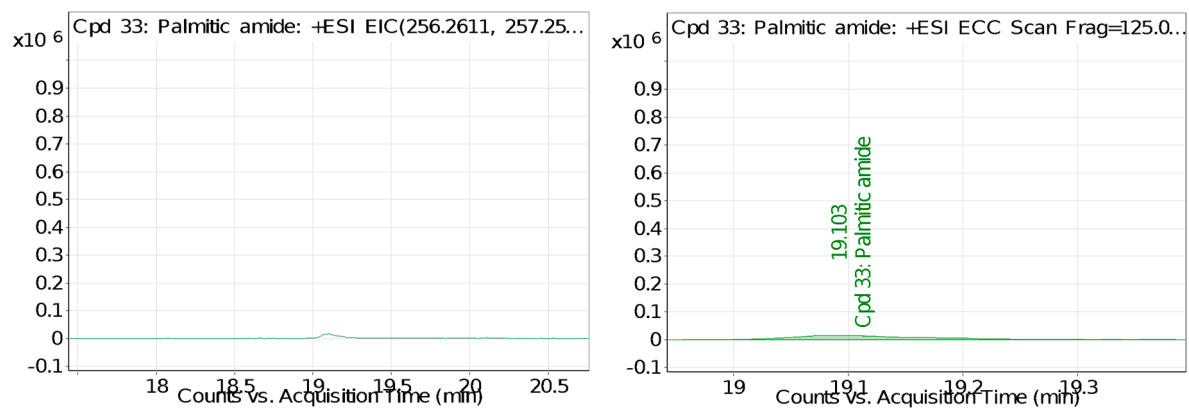




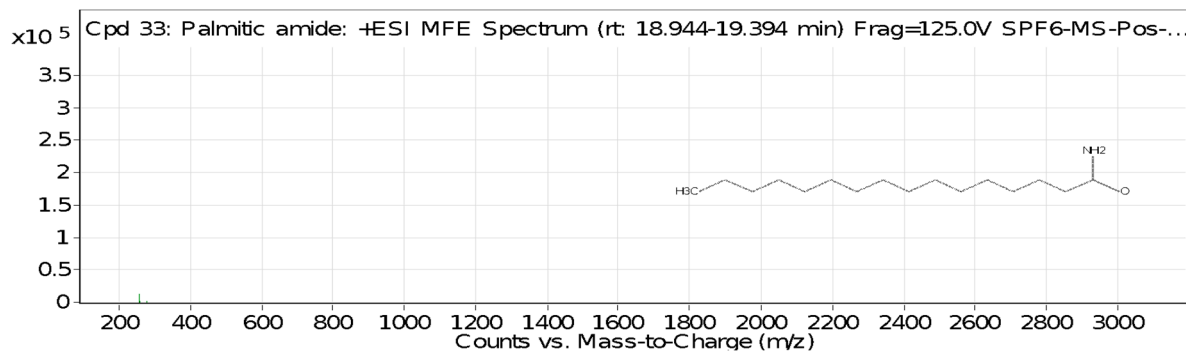
**Compound xv: Palmitic amide**

Compound Label	Name	<i>m/z</i>	RT	Algorithm	Mass
Cpd 33: Palmitic amide	Palmitic amide	256.2636	19.103	Find by Molecular Feature	255.2562

**Compound Chromatograms**



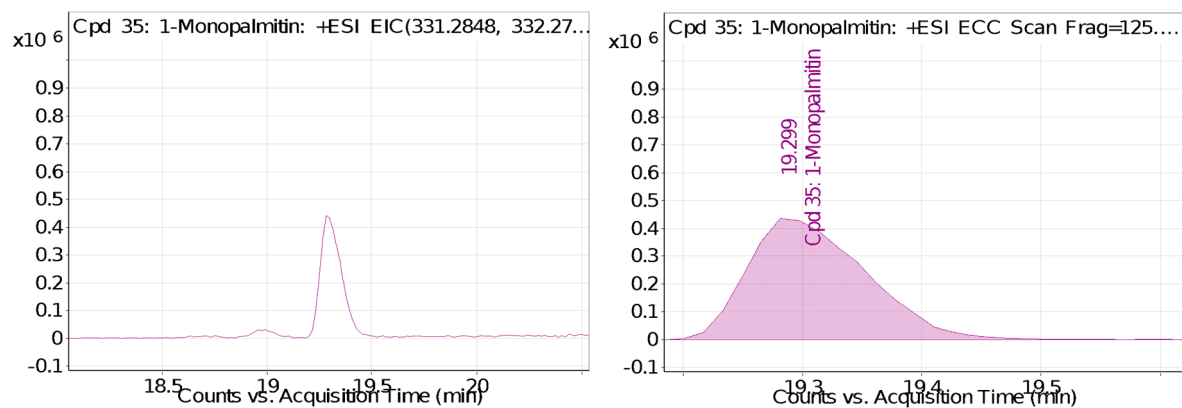
**MFE MS Spectrum**



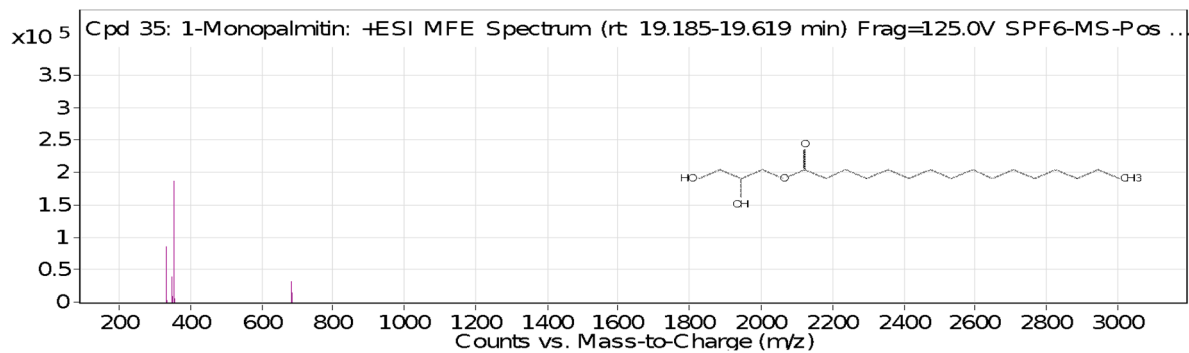
### Compound xvi: 1-monopalmitin

Compound Label	Name	<i>m/z</i>	RT	Algorithm	Mass
Cpd 35: 1-monopalmitin	1-monopalmitin	353.2669	19.299	Find by Molecular Feature	330.2773

### Compound Chromatograms



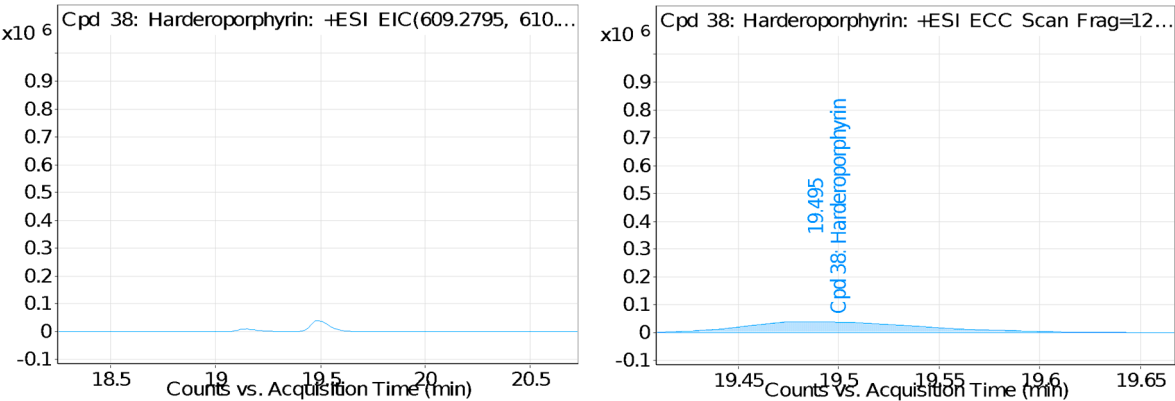
### MFE MS Spectrum



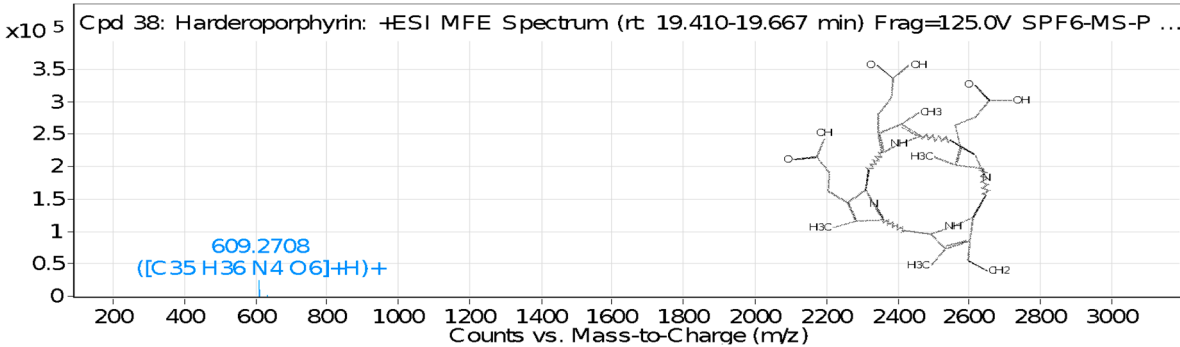
Compound xvii: Harderoporphyrin

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 38: Harderoporphyrin	Harderoporphyrin	609.2708	19.495	Find by Molecular Feature	608.2636

Compound Chromatograms



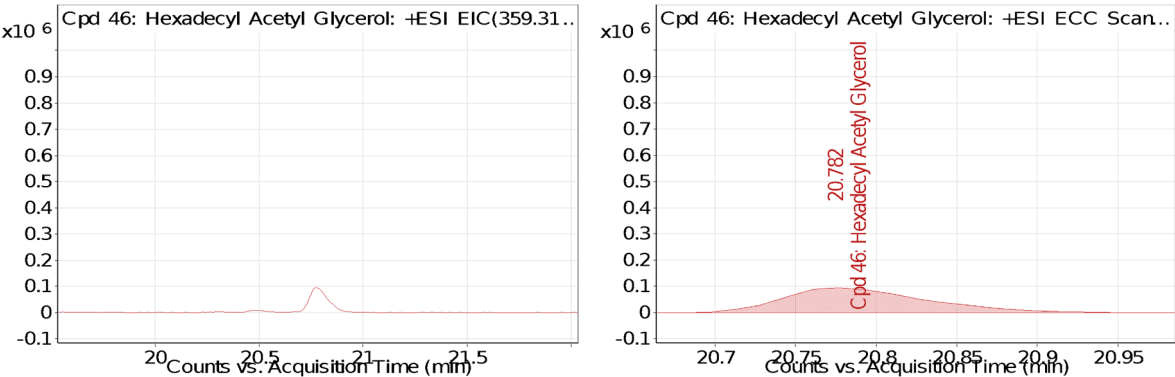
MFE MS Spectrum



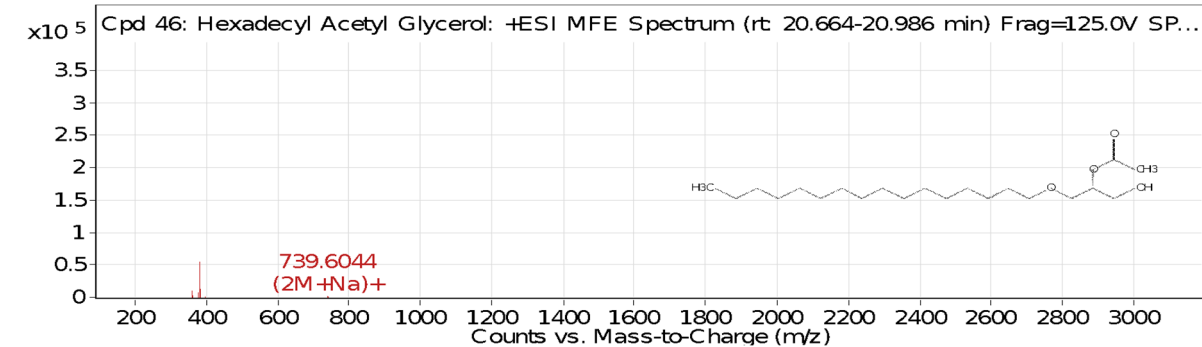
Compound xviii: Hexadecyl acetyl glycerol

Compound Label	Name	m/z	RT	Algorithm	Mass
Cpd 46: Hexadecyl acetyl glycerol	Hexadecyl acetyl glycerol	381.2983	20.782	Find by Molecular Feature	358.3092

Compound Chromatograms



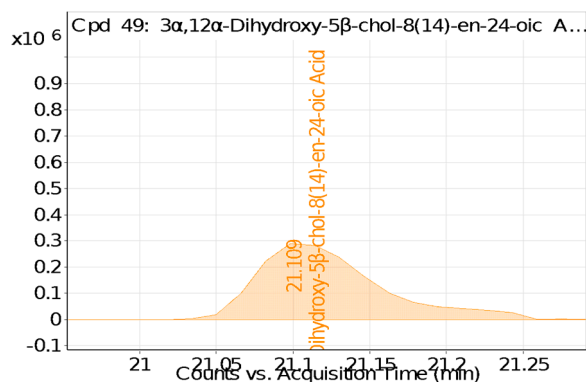
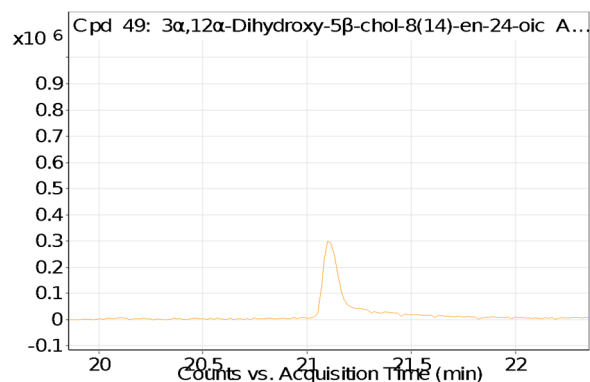
MFE MS Spectrum



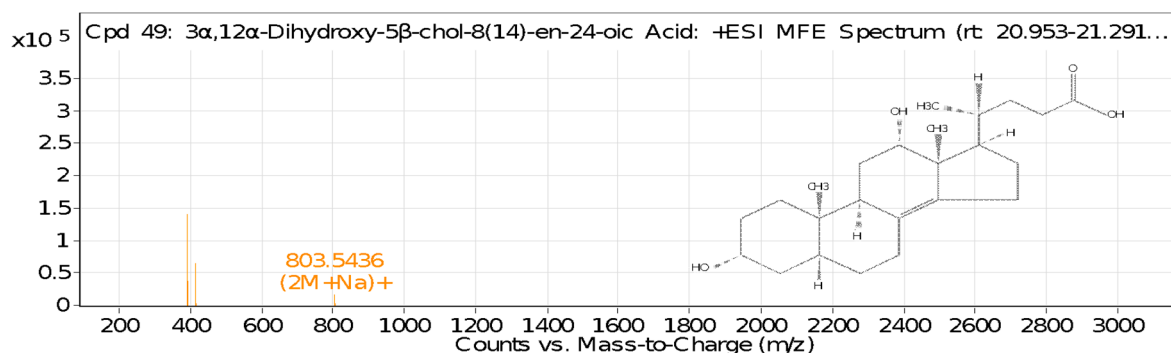
**Compound xix: 3 $\alpha$ ,12 $\alpha$ -Dihydroxy-5 $\beta$ -chol-8(14)-en-24-oic acid**

Compound Label	Name	<i>m/z</i>	RT	Algorithm	Mass
Cpd 49: 3 $\alpha$ ,12 $\alpha$ -Dihydroxy-5 $\beta$ -chol-8(14)-en-24-oic acid	3 $\alpha$ ,12 $\alpha$ -Dihydroxy-5 $\beta$ -chol-8(14)-en-24-oic acid	391.2854	21.109	Find by Molecular Feature	390.278

**Compound Chromatograms**



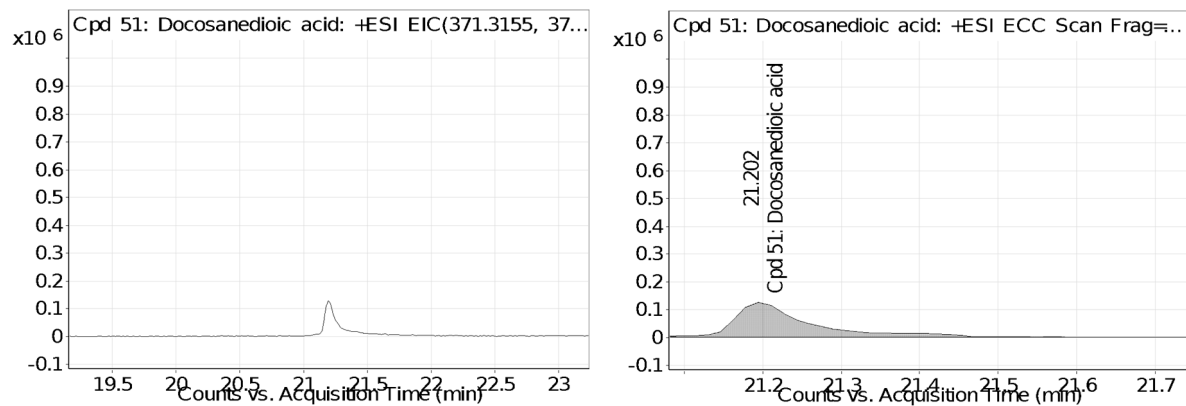
**MFE MS Spectrum**



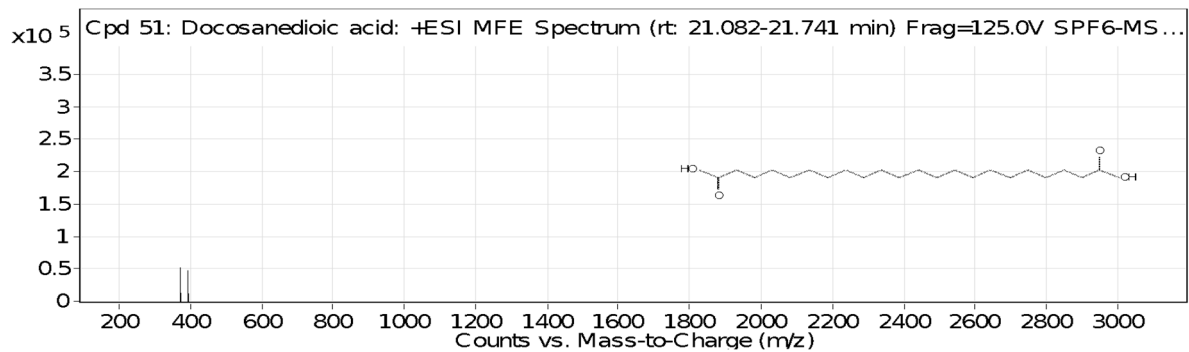
### Compound xx: Docosanedioic acid

Compound Label	Name	<i>m/z</i>	RT	Algorithm	Mass
Cpd 51: Docosanedioic acid	Docosanedioic acid	371.3151	21.202	Find by Molecular Feature	370.3077

### Compound Chromatograms



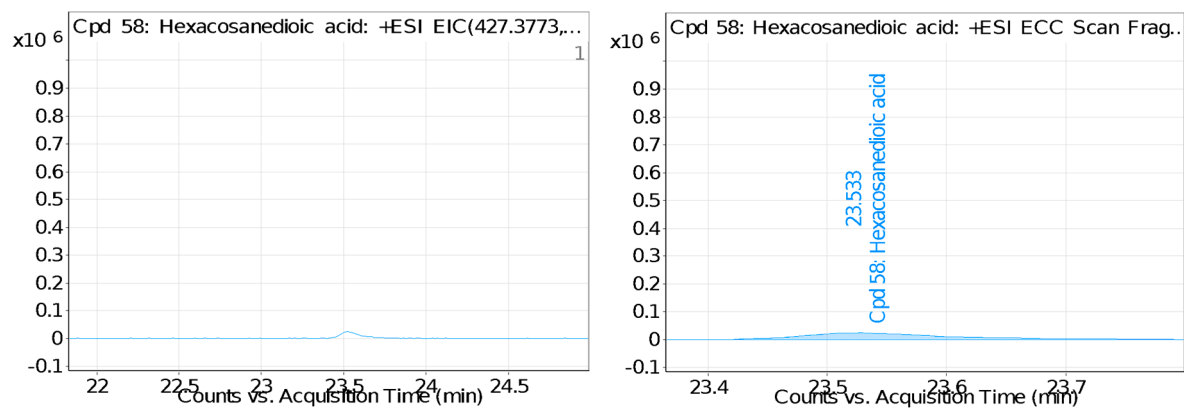
### MFE MS Spectrum



**Compound xxi: Hexacosanedioic acid**

Compound Label	Name	<i>m/z</i>	RT	Algorithm	Mass
Cpd 58: Hexacosanedioic acid	Hexacosanedioic acid	449.3598	23.533	Find by Molecular Feature	426.3705

## Compound Chromatograms



## MFE MS Spectrum

