

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100220\
 Data File : BG046771.D
 Acq On : 3 Oct 2020 14:30
 Operator : CG/JU
 Sample : L4150-21
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CB7M2

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG046771.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.532	29	33	38	rVB3	12710	17731	0.63%	0.067%
2	4.178	138	143	149	rVB7	8402	16449	0.59%	0.062%
3	4.290	156	162	167	rBV4	7765	13159	0.47%	0.050%
4	4.496	192	197	203	rBV2	30002	43553	1.55%	0.165%
5	4.648	218	223	226	rBV5	8510	13911	0.49%	0.053%
6	4.913	263	268	274	rVB	24957	34890	1.24%	0.132%
7	5.207	313	318	322	rVB2	3771	7267	0.26%	0.028%
8	5.289	328	332	337	rBV4	3529	6622	0.24%	0.025%
9	5.412	349	353	355	rBV4	4001	5728	0.20%	0.022%
10	5.589	377	383	389	rBV2	29546	43264	1.54%	0.164%
11	6.153	476	479	485	rVB2	6424	10140	0.36%	0.038%
12	7.057	630	633	637	rBV4	4014	5265	0.19%	0.020%
13	7.198	651	657	660	rBV2	24063	35548	1.26%	0.135%
14	7.257	660	667	677	rVB	222944	405128	14.41%	1.536%
15	7.433	690	697	705	rBV	153322	260924	9.28%	0.989%
16	7.639	725	732	739	rBV	218497	373427	13.28%	1.416%
17	7.739	744	749	752	rVV2	4845	6676	0.24%	0.025%
18	7.786	754	757	764	rVB4	3169	6120	0.22%	0.023%
19	8.109	805	812	819	rBV	275128	466914	16.61%	1.770%
20	8.162	819	821	825	rVB2	5562	6942	0.25%	0.026%
21	8.338	847	851	858	rVB5	6434	12650	0.45%	0.048%
22	8.708	908	914	919	rBV5	5643	9642	0.34%	0.037%
23	8.802	923	930	943	rVB	282449	493564	17.56%	1.871%
24	8.932	947	952	957	rBV3	17523	30985	1.10%	0.117%
25	9.214	996	1000	1003	rBV2	4130	6977	0.25%	0.026%
26	9.267	1003	1009	1019	rVB	217189	390965	13.91%	1.482%
27	9.431	1031	1037	1044	rBV4	14762	22255	0.79%	0.084%
28	9.607	1062	1067	1072	rBV4	3355	6099	0.22%	0.023%
29	9.701	1077	1083	1089	rVB2	12924	20455	0.73%	0.078%
30	9.995	1125	1133	1140	rBV2	185663	333541	11.87%	1.264%
31	10.312	1183	1187	1191	rBV5	4977	8761	0.31%	0.033%
32	10.430	1200	1207	1217	rBV2	58684	98816	3.52%	0.375%
33	10.536	1217	1225	1235	rVB	293981	530888	18.89%	2.012%
34	10.923	1283	1291	1300	rVB	382844	676409	24.06%	2.564%
35	11.047	1305	1312	1320	rBV	202188	359814	12.80%	1.364%
36	11.576	1397	1402	1406	rVB4	3749	6788	0.24%	0.026%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

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37	11.693	1417	1422	1429	rVB5	5264	9198	0.33%	0.035%
38	11.852	1445	1449	1452	rVV3	6821	8096	0.29%	0.031%
39	11.934	1459	1463	1465	rBV4	8945	10825	0.39%	0.041%
40	12.193	1502	1507	1511	rBV2	7058	10649	0.38%	0.040%

41	12.498	1555	1559	1563	rVB2	6516	9786	0.35%	0.037%
42	12.645	1578	1584	1589	rVB5	6096	11384	0.40%	0.043%
43	13.186	1672	1676	1680	rBV4	6118	9508	0.34%	0.036%
44	13.350	1696	1704	1705	rBV4	7887	12666	0.45%	0.048%
45	13.368	1705	1707	1711	rVV2	6361	7312	0.26%	0.028%

46	13.562	1734	1740	1742	rBV3	4317	7037	0.25%	0.027%
47	13.597	1744	1746	1750	rVV5	8786	8936	0.32%	0.034%
48	14.131	1831	1837	1842	rVV	555717	807065	28.71%	3.059%
49	14.178	1842	1845	1850	rVB2	25065	35964	1.28%	0.136%
50	14.255	1857	1858	1864	rVB4	4001	6188	0.22%	0.023%

51	14.372	1872	1878	1882	rBV6	15723	34512	1.23%	0.131%
52	14.425	1882	1887	1897	rVB	582919	901660	32.08%	3.418%
53	14.590	1911	1915	1919	rBV2	18748	23094	0.82%	0.088%
54	14.625	1919	1921	1926	rVB5	6867	6663	0.24%	0.025%
55	14.737	1930	1940	1948	rBV	629979	1037553	36.91%	3.933%

56	14.901	1962	1968	1978	rBV	288824	443379	15.77%	1.681%
57	15.054	1992	1994	1998	rVB4	6483	6786	0.24%	0.026%
58	15.130	2003	2007	2013	rBV2	36834	52098	1.85%	0.197%
59	15.383	2043	2050	2054	rBV5	4877	11693	0.42%	0.044%
60	15.512	2069	2072	2073	rBV2	5652	5389	0.19%	0.020%

61	15.577	2079	2083	2087	rVB3	5914	7756	0.28%	0.029%
62	15.724	2101	2108	2115	rBV	772436	1160512	41.28%	4.399%
63	15.824	2118	2125	2134	rBV2	185731	286735	10.20%	1.087%
64	15.959	2141	2148	2156	rVB6	14480	28938	1.03%	0.110%
65	16.129	2173	2177	2178	rBV4	4243	5388	0.19%	0.020%

66	16.258	2196	2199	2202	rBV4	8804	12684	0.45%	0.048%
67	16.441	2225	2230	2233	rBV3	21987	31626	1.13%	0.120%
68	16.505	2239	2241	2245	rVB4	7068	9134	0.32%	0.035%
69	16.793	2284	2290	2297	rBV6	22834	45726	1.63%	0.173%
70	16.864	2299	2302	2307	rVV6	11757	18505	0.66%	0.070%

71	16.934	2309	2314	2315	rBV3	7201	9006	0.32%	0.034%
72	17.081	2332	2339	2342	rBV7	10447	20648	0.73%	0.078%
73	17.222	2361	2363	2367	rBV3	8868	9521	0.34%	0.036%
74	17.357	2382	2386	2389	rBV6	12263	18558	0.66%	0.070%
75	17.475	2400	2406	2412	rVV	802617	1228233	43.69%	4.656%

76	17.575	2416	2423	2431	rVB	784374	1256861	44.71%	4.764%
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 Stop Thrs : 0 Peak Location: TOP

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77	17.780	2456	2458	2465	rVB7	11864	20256	0.72%	0.077%
78	17.851	2466	2470	2472	rBV4	20825	28281	1.01%	0.107%
79	17.968	2488	2490	2495	rVB5	10062	12442	0.44%	0.047%
80	18.238	2532	2536	2542	rBV5	42426	62830	2.24%	0.238%
81	18.297	2542	2546	2552	rBV3	90958	138583	4.93%	0.525%
82	18.362	2552	2557	2568	rVV	320433	482857	17.18%	1.830%
83	18.691	2611	2613	2618	rVB6	18992	24567	0.87%	0.093%
84	19.085	2677	2680	2686	rVB2	108583	143238	5.10%	0.543%
85	19.372	2726	2729	2732	rBV3	33246	36110	1.28%	0.137%
86	19.484	2745	2748	2749	rBV3	37567	42399	1.51%	0.161%
87	19.513	2749	2753	2760	rVV3	129837	267579	9.52%	1.014%
88	19.578	2760	2764	2768	rVB4	84954	123193	4.38%	0.467%
89	19.625	2769	2772	2777	rVV	94442	108415	3.86%	0.411%
90	19.854	2806	2811	2823	rVB	1022451	1465998	52.15%	5.557%
91	20.295	2883	2886	2890	rVB	51251	60302	2.15%	0.229%
92	21.394	3068	3073	3083	rVB	411543	676838	24.08%	2.566%
93	21.752	3128	3134	3141	rBV	805657	1291514	45.94%	4.896%
94	22.598	3272	3278	3295	rVB2	452258	1078285	38.36%	4.087%
95	23.644	3450	3456	3466	rVB	516971	1029057	36.61%	3.901%
96	24.126	3532	3538	3541	rBV2	304864	611399	21.75%	2.318%
97	24.179	3542	3547	3562	rVB	1183182	2811034	100.00%	10.656%
98	24.795	3645	3652	3663	rVB	444187	1217670	43.32%	4.616%
99	25.025	3684	3691	3700	rBV2	406718	1062383	37.79%	4.027%
100	26.264	3895	3902	3912	rVB2	391846	1169346	41.60%	4.433%

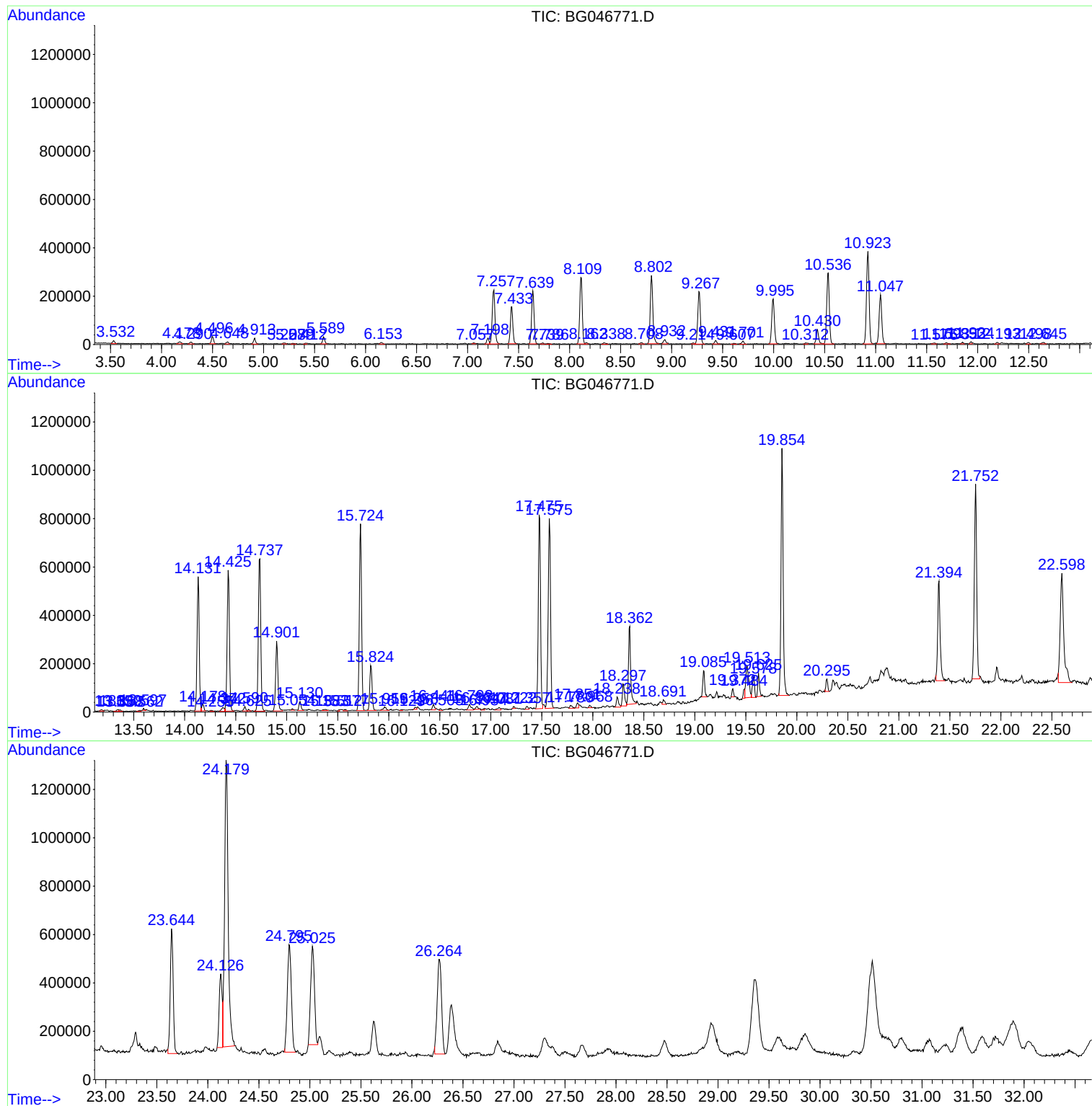
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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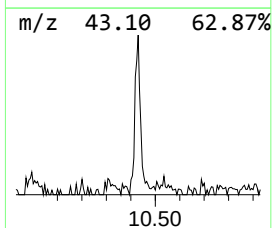
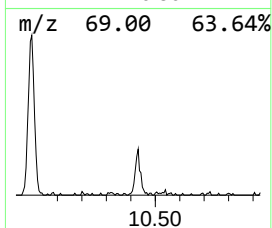
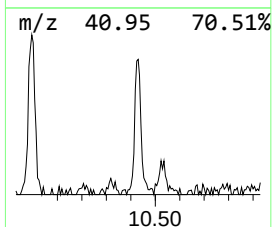
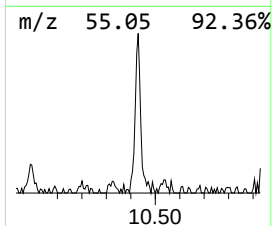
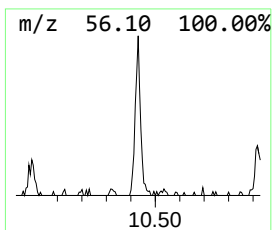
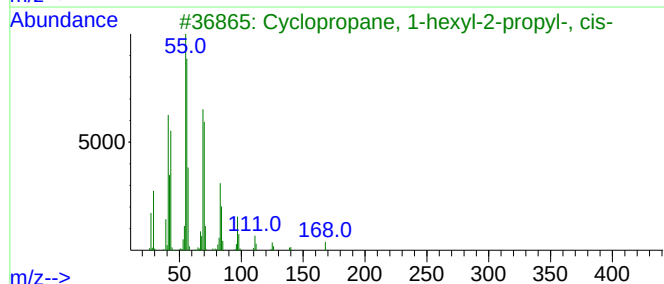
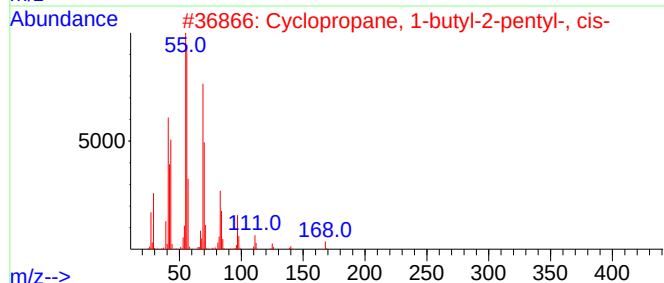
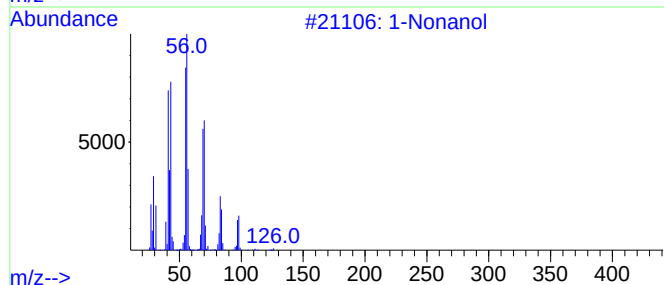
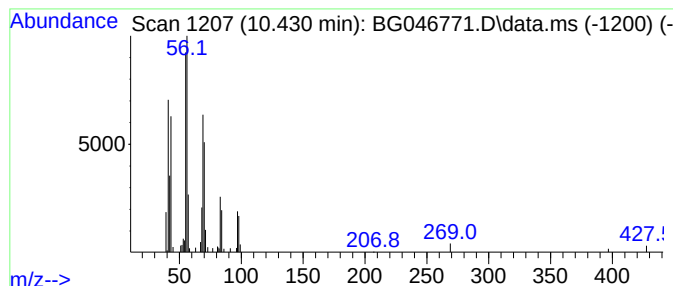
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 1-Nonanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.430	2.92 ng/ul	98816	Naphthalene-d8	10.924

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonanol			144	C9H20O	000143-08-8	86
2	Cyclopropane, 1-butyl-2-pentyl-,...			168	C12H24	074663-88-0	78
3	Cyclopropane, 1-hexyl-2-propyl-,...			168	C12H24	074630-58-3	78
4	Cyclopropane, pentyl-			112	C8H16	002511-91-3	72
5	3-Dodecene, (E)-			168	C12H24	007206-14-6	53



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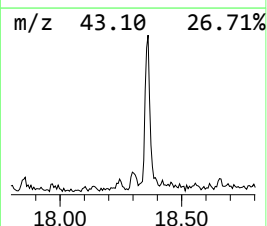
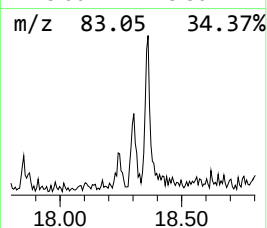
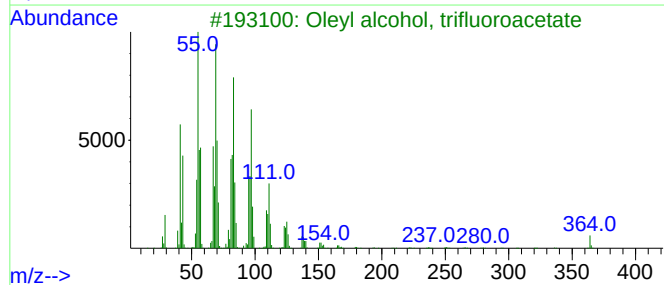
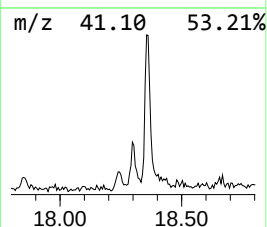
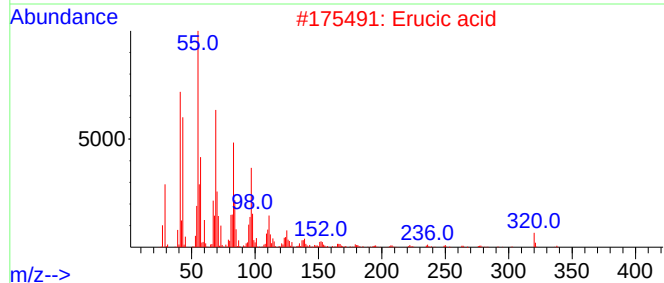
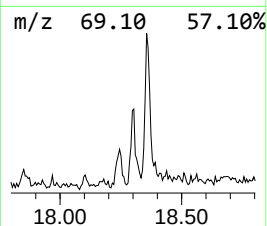
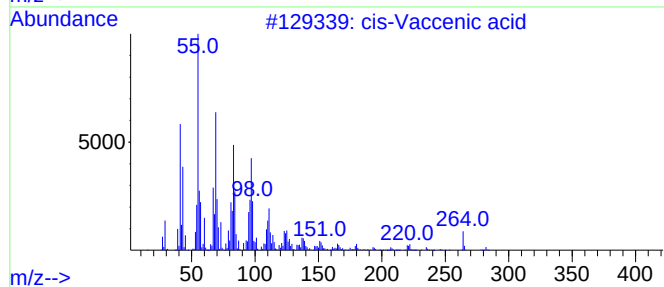
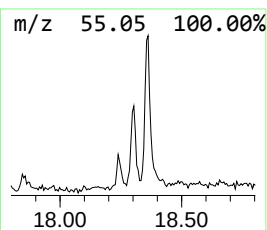
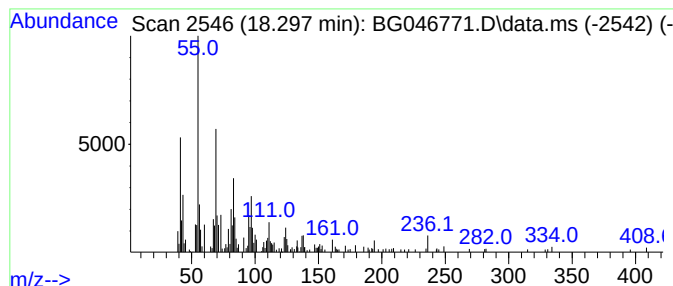
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 cis-Vaccenic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.300	2.26 ng/ul	138583	Phenanthrene-d10	17.475

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-Vaccenic acid	282	C18H34O2	000506-17-2	68
2			Erucic acid	338	C22H42O2	000112-86-7	64
3			Oleyl alcohol, trifluoroacetate	364	C20H35F3O2	1000352-68-4	62
4			22-Tricosenoic acid	352	C23H44O2	065119-95-1	58
5			Z-11-Tetradecenoic acid	226	C14H26O2	1000130-83-3	55



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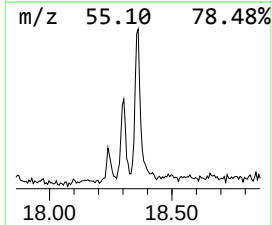
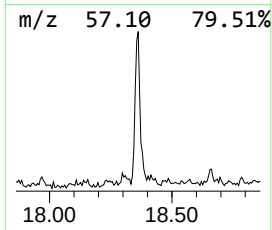
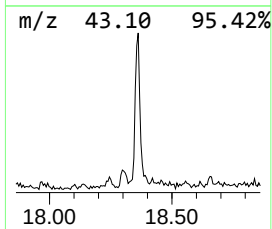
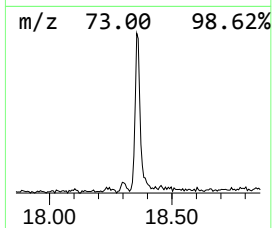
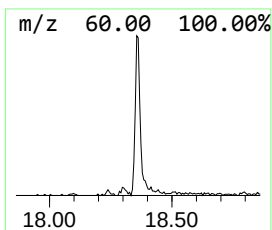
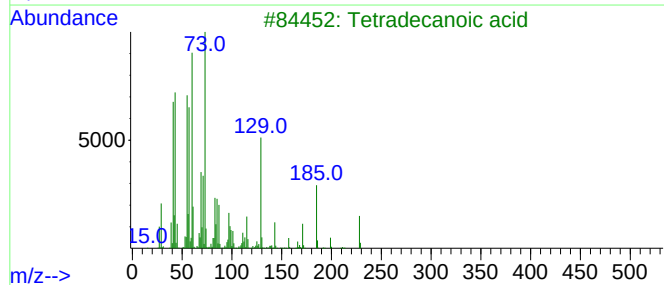
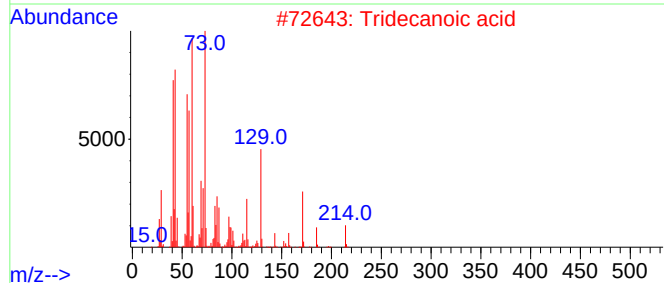
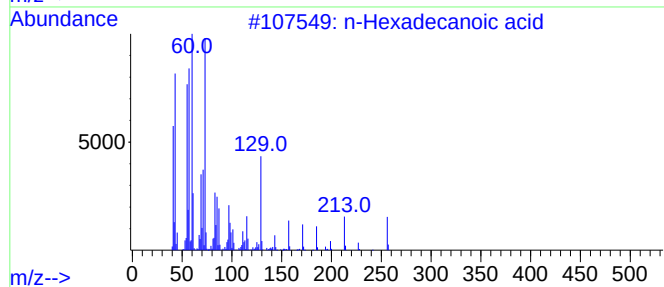
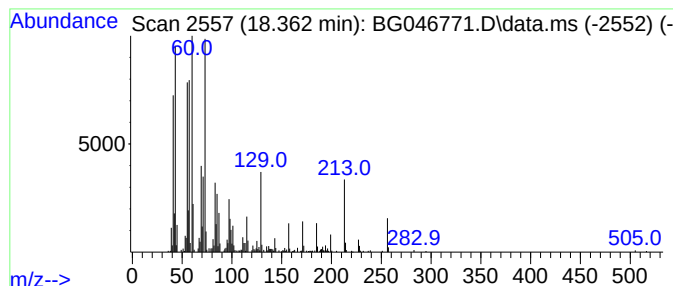
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.360	7.86 ng/ul	482857	Phenanthrene-d10	17.475

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Tridecanoic acid	214	C13H26O2	000638-53-9	95
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	94
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	81
5			Dodecanoic acid	200	C12H24O2	000143-07-7	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100220\
Data File : BG046771.D
Acq On : 3 Oct 2020 14:30
Operator : CG/JU
Sample : L4150-21
Misc :
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Instrument :
BNA_G
ClientSampleId :
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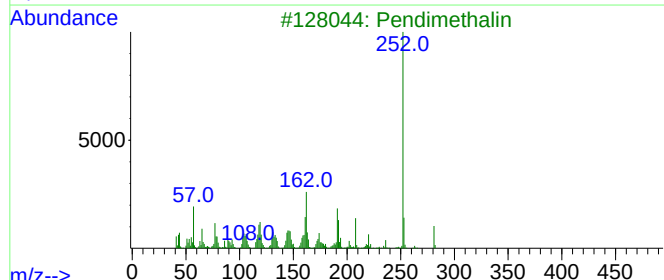
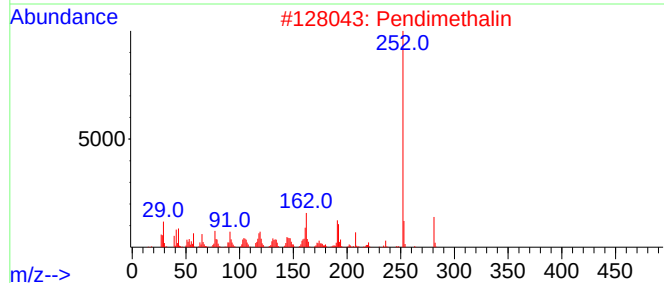
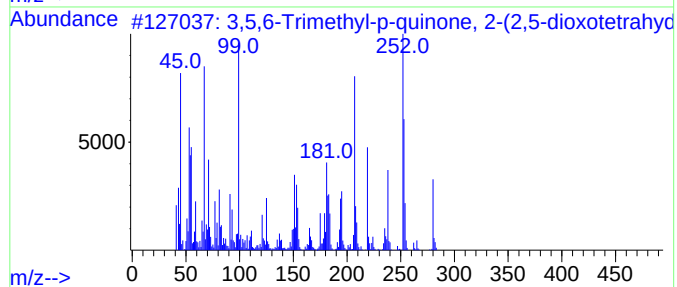
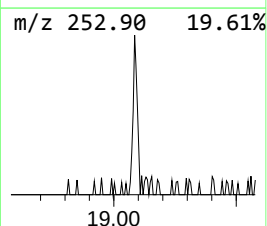
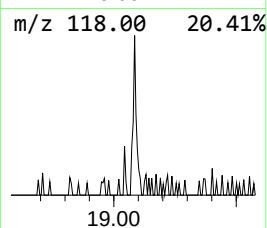
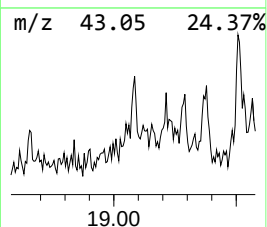
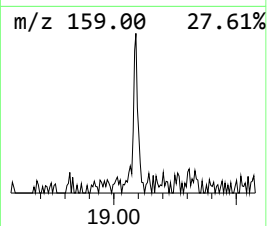
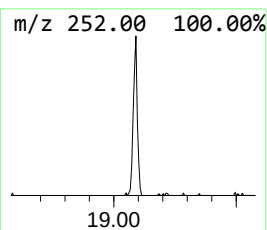
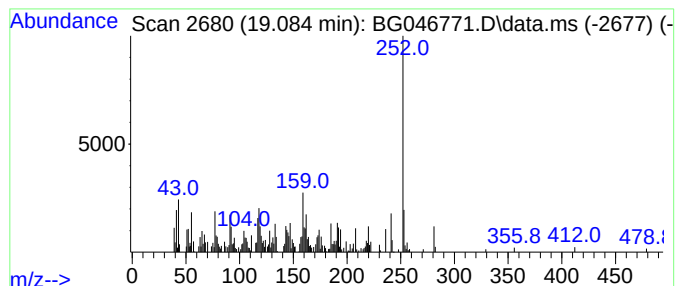
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 3,5,6-Trimethyl-p-quinone, ... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.080	2.33 ng/ul	143238	Phenanthrene-d10	17.475

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,5,6-Trimethyl-p-quinone, 2-(2,...	280	C13H12O5S	126848-01-9	95
2			Pendimethalin	281	C13H19N3O4	040487-42-1	78
3			Pendimethalin	281	C13H19N3O4	040487-42-1	70
4			Pendimethalin	281	C13H19N3O4	040487-42-1	64
5			Pyridine-3-carbonitrile, 2-amino...	281	C17H19N3O	1000264-87-9	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100220\
Data File : BG046771.D
Acq On : 3 Oct 2020 14:30
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Sample : L4150-21
Misc :
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ClientSampleId :
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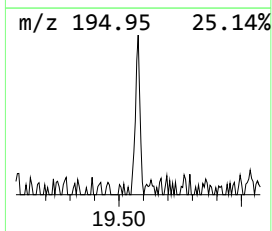
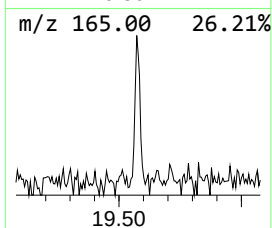
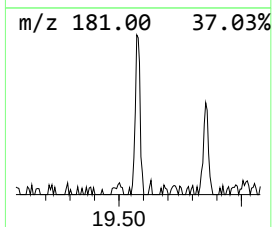
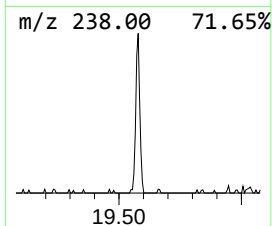
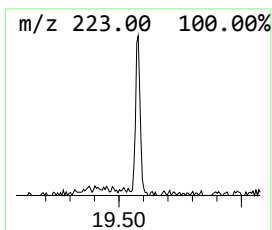
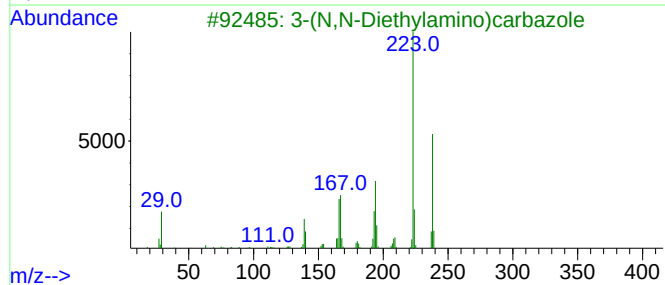
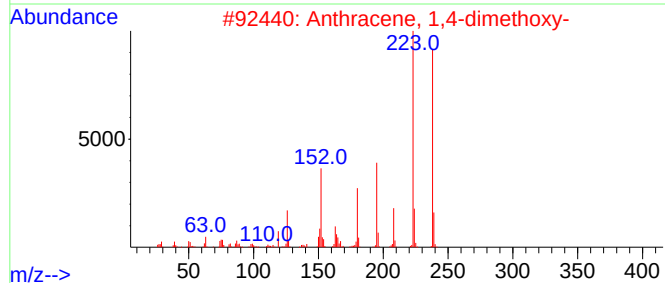
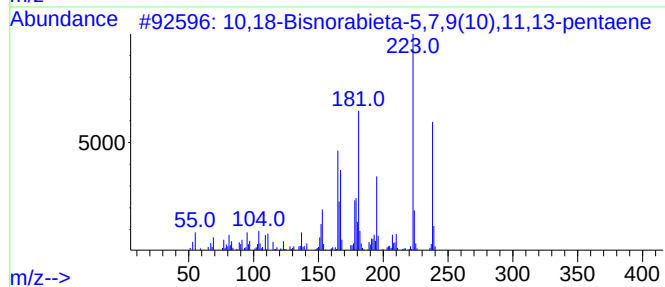
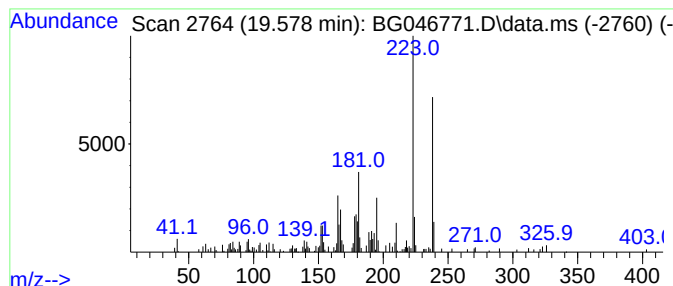
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.580	2.01 ng/ul	123193	Phenanthrene-d10	17.475

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	95		
2	Anthracene, 1,4-dimethoxy-	238	C16H14O2	013076-29-4	70		
3	3-(N,N-Diethylamino)carbazole	238	C16H18N2	054994-31-9	58		
4	9,10-Anthracenedione, 1-amino-	223	C14H9NO2	000082-45-1	49		
5	Anthraquinone, 2-amino-	223	C14H9NO2	000117-79-3	49		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100220\
Data File : BG046771.D
Acq On : 3 Oct 2020 14:30
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Sample : L4150-21
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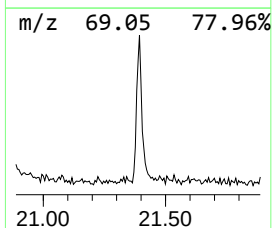
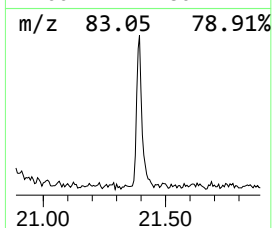
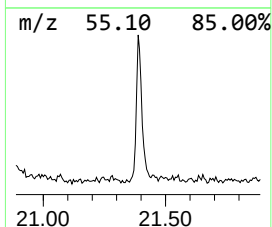
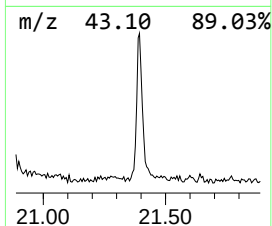
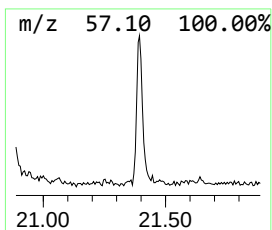
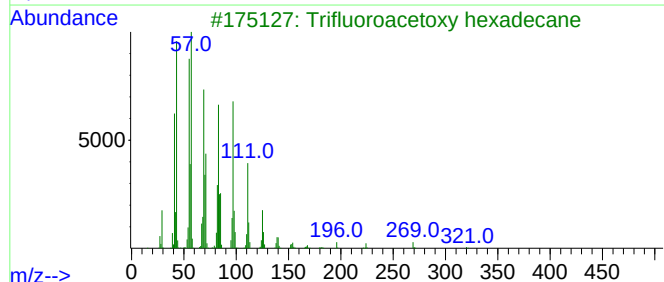
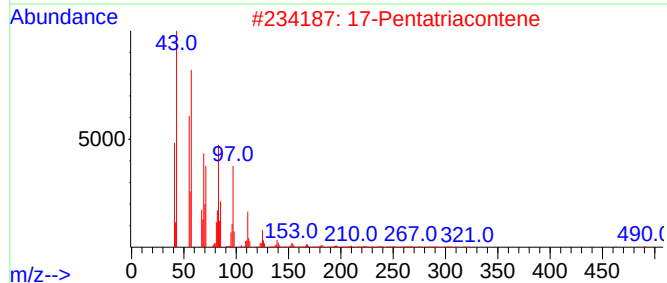
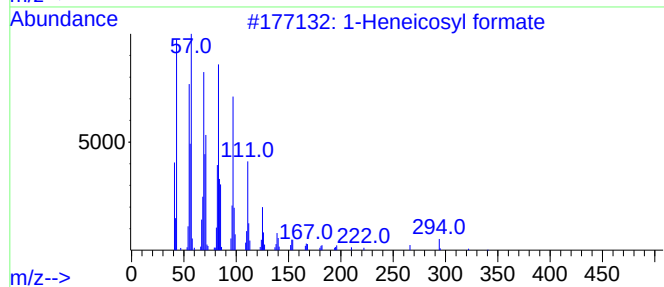
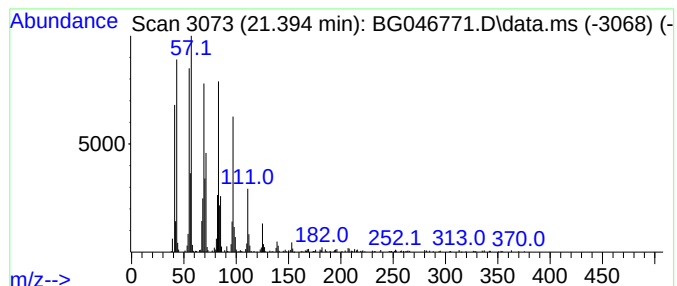
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 1-Heneicosyl formate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.390	10.48 ng/ul	676838	Chrysene-d12	21.752

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosyl formate	340	C22H44O2	077899-03-7	91
2			17-Pentatriacontene	491	C35H70	006971-40-0	91
3			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	91
4			3-Eicosene, (E)-	280	C20H40	074685-33-9	87
5			Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100220\
Data File : BG046771.D
Acq On : 3 Oct 2020 14:30
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Sample : L4150-21
Misc :
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Instrument :
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ClientSampleId :
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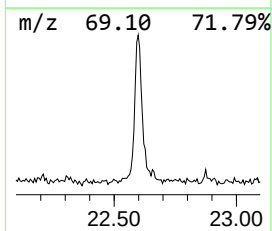
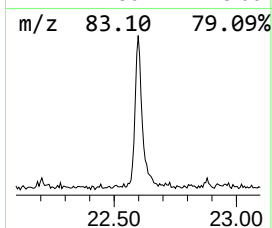
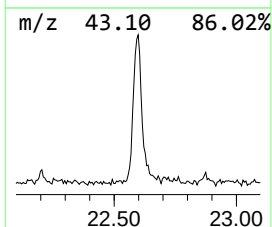
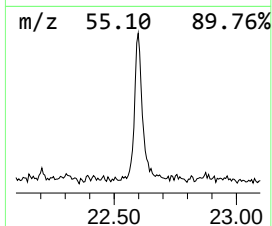
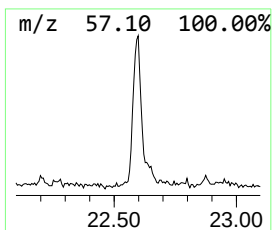
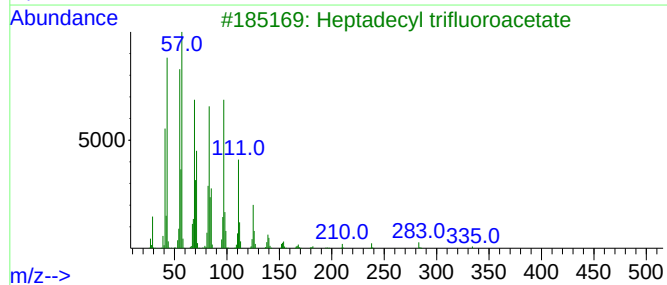
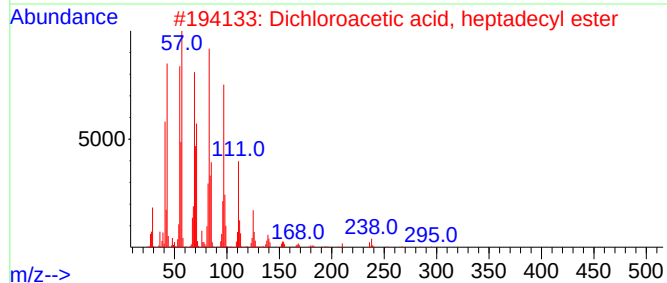
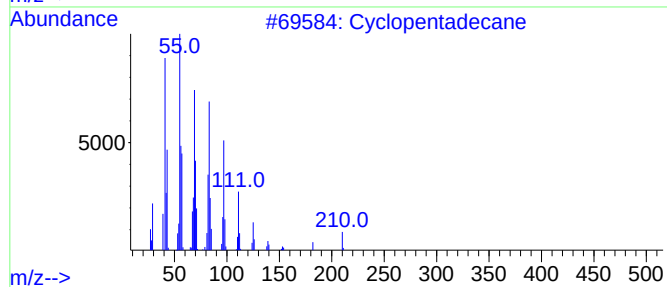
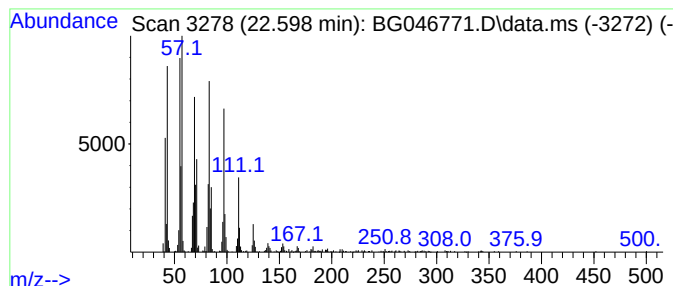
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Cyclic22.60 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.600	16.70 ng/ul	1078290	Chrysene-d12	21.752

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentadecane	210	C15H30	000295-48-7	98
2			Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	94
3			Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	91
4			1-Heneicosanol	312	C21H44O	015594-90-8	91
5			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100220\
Data File : BG046771.D
Acq On : 3 Oct 2020 14:30
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Sample : L4150-21
Misc :
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Instrument :
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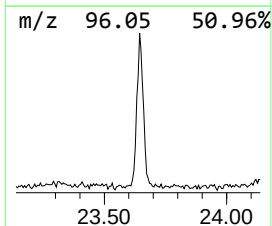
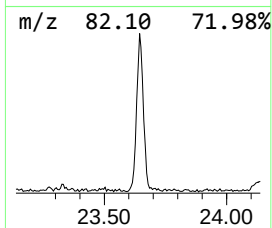
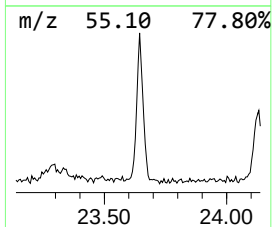
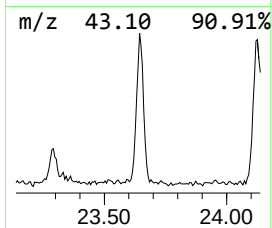
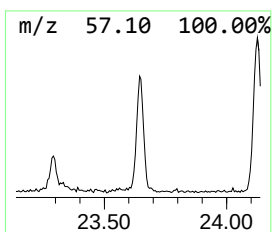
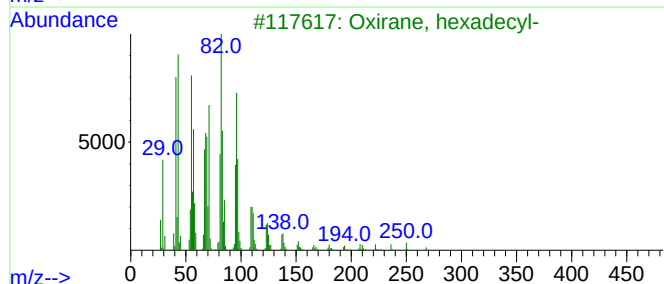
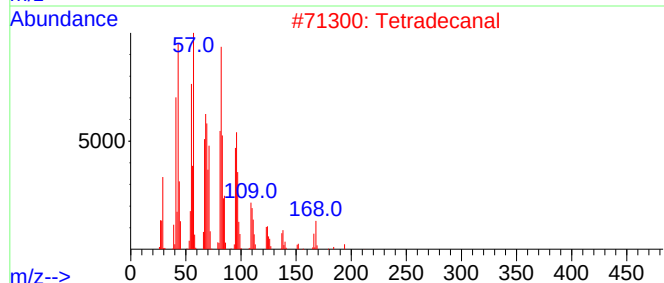
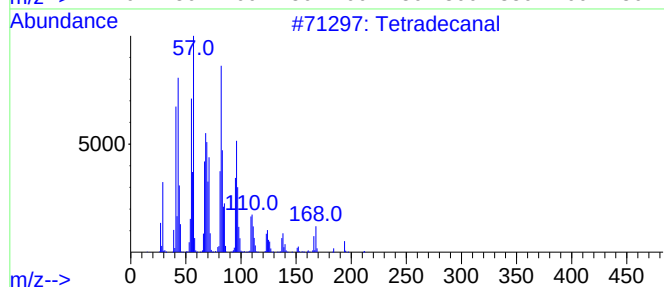
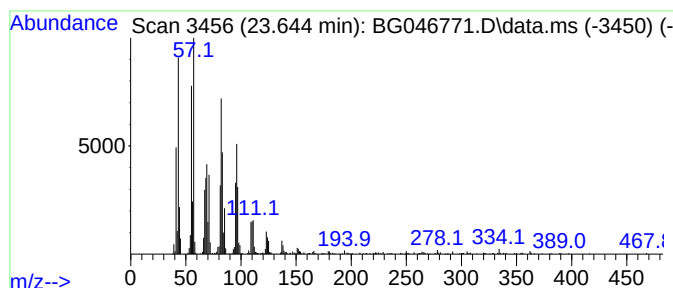
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Tetradecanal Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.640	19.37 ng/ul	1029060	Perylene-d12	25.025

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecanal	212	C14H28O	000124-25-4	94
2			Tetradecanal	212	C14H28O	000124-25-4	91
3			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
4			Octadecanal	268	C18H36O	000638-66-4	91
5			Pentadecanal	226	C15H30O	002765-11-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100220\
Data File : BG046771.D
Acq On : 3 Oct 2020 14:30
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Sample : L4150-21
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Instrument :
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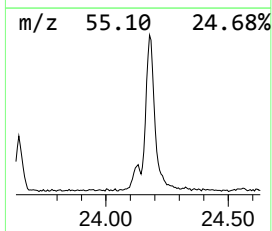
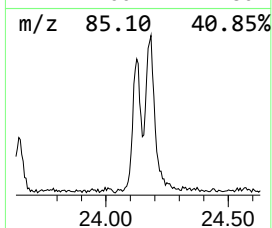
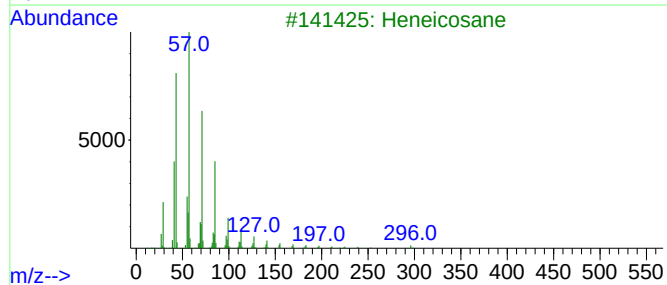
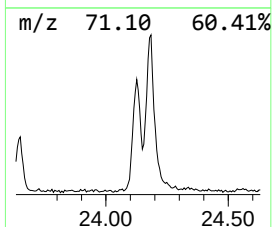
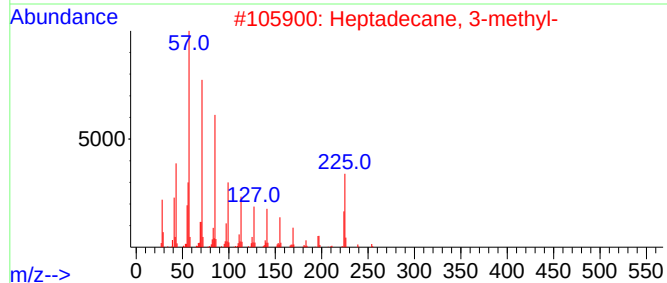
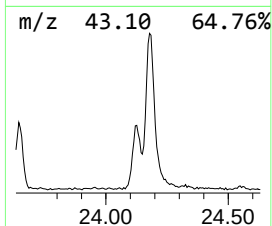
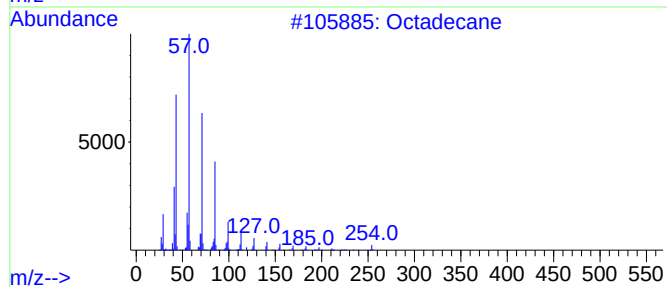
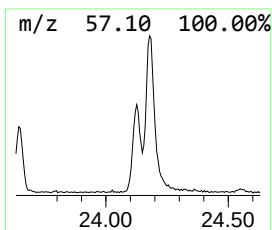
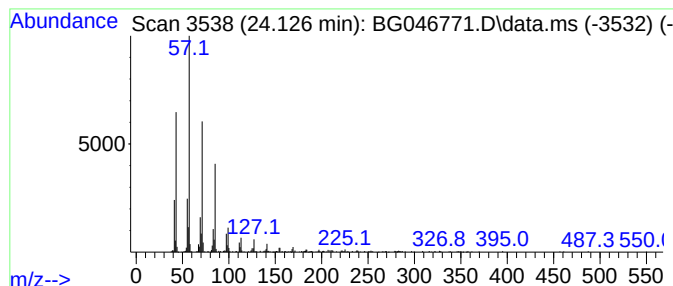
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.130	11.51 ng/ul	611399	Perylene-d12	25.025

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecane	254	C18H38	000593-45-3	94
2		Heptadecane, 3-methyl-	254	C18H38	006418-44-6	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Heptacosane	380	C27H56	000593-49-7	90
5		Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100220\
Data File : BG046771.D
Acq On : 3 Oct 2020 14:30
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Sample : L4150-21
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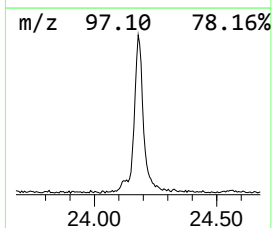
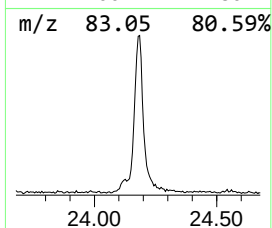
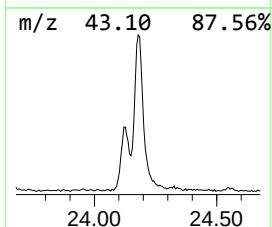
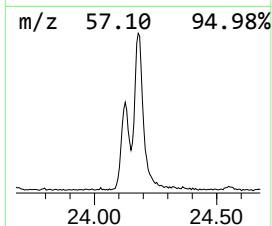
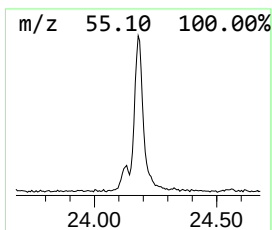
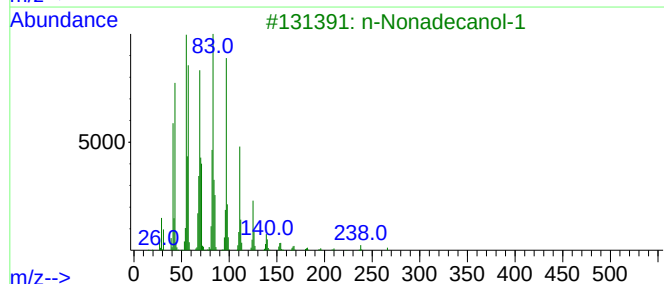
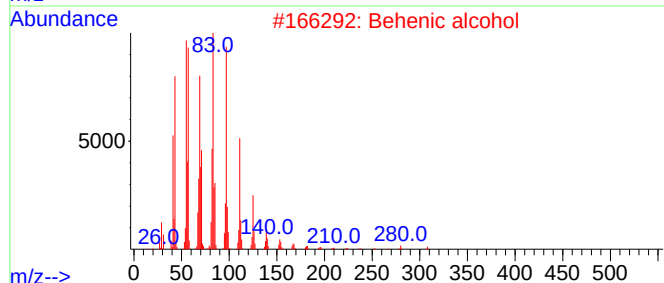
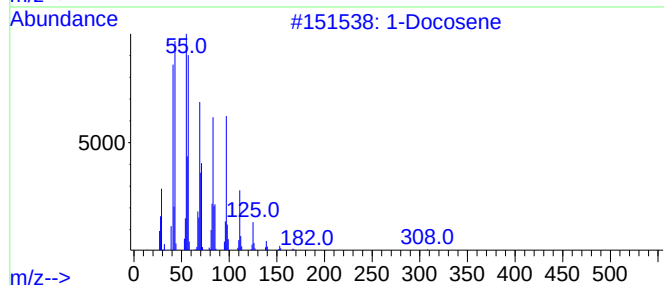
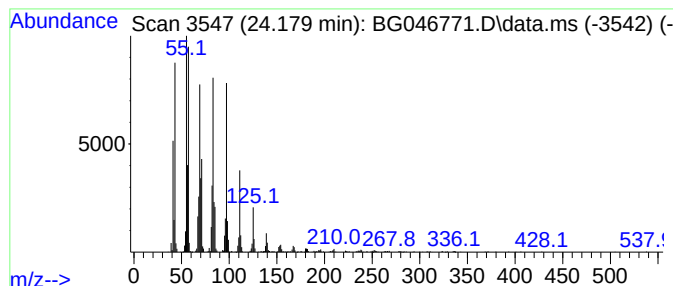
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.180	52.92 ng/ul	2811030	Perylene-d12	25.025

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene			308	C22H44	001599-67-3	99
2	Behenic alcohol			326	C22H46O	000661-19-8	91
3	n-Nonadecanol-1			284	C19H40O	001454-84-8	91
4	Trifluoroacetoxy hexadecane			338	C18H33F3O2	006222-03-3	90
5	3-Eicosene, (E)-			280	C20H40	074685-33-9	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100220\
Data File : BG046771.D
Acq On : 3 Oct 2020 14:30
Operator : CG/JU
Sample : L4150-21
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CB7M2

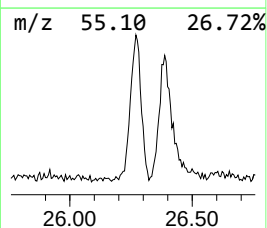
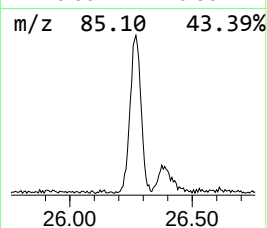
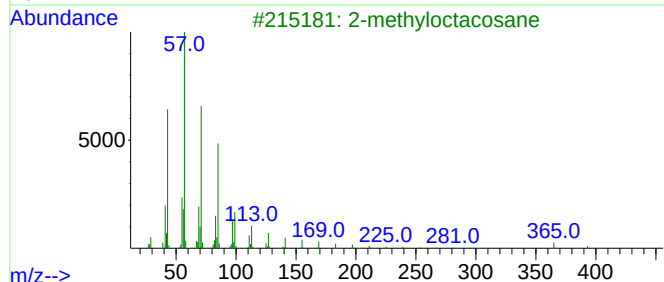
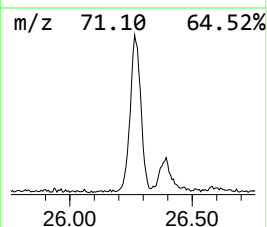
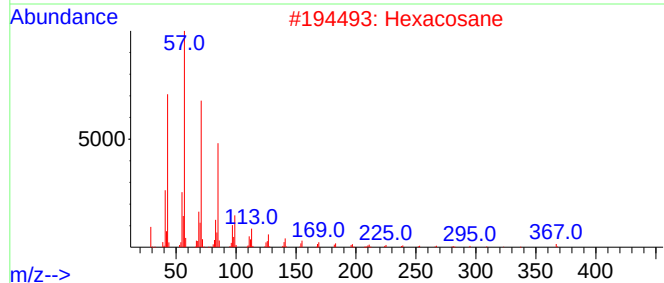
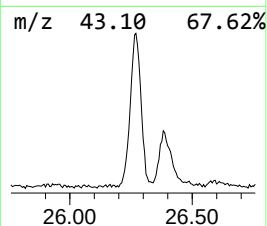
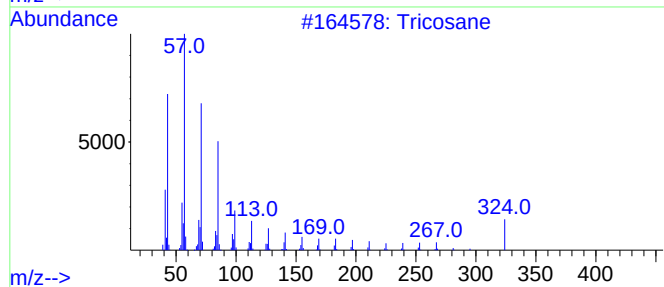
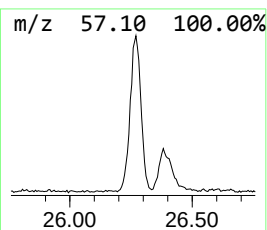
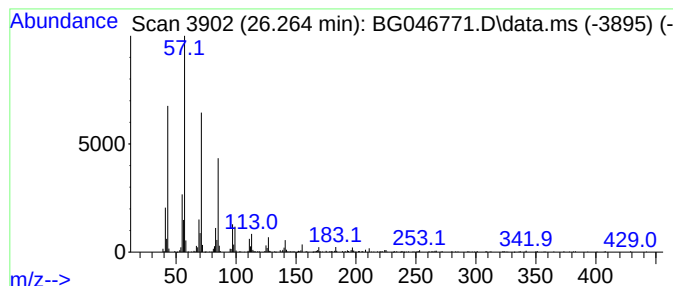
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.260	22.01 ng/ul	1169350	Perylene-d12	25.025

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tricosane	324	C23H48	000638-67-5	95
2		Hexacosane	366	C26H54	000630-01-3	93
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Heptacosane	380	C27H56	000593-49-7	91
5		Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100220\
Data File : BG046771.D
Acq On : 3 Oct 2020 14:30
Operator : CG/JU
Sample : L4150-21
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CB7M2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Nonanol	10.430	2.9	ng/ul	98816	2	10.924	676409	20.0
cis-Vaccenic acid	18.300	2.3	ng/ul	138583	4	17.475	1228230	20.0
n-Hexadecanoic ...	18.360	7.9	ng/ul	482857	4	17.475	1228230	20.0
3,5,6-Trimethyl...	19.080	2.3	ng/ul	143238	4	17.475	1228230	20.0
10,18-Bisnorabi...	19.580	2.0	ng/ul	123193	4	17.475	1228230	20.0
1-Heneicosyl fo...	21.390	10.5	ng/ul	676838	5	21.752	1291510	20.0
(DEL) Alkane: C...	22.600	16.7	ng/ul	1078290	5	21.752	1291510	20.0
Tetradecanal	23.640	19.4	ng/ul	1029060	6	25.025	1062380	20.0
(DEL) Alkane: S...	24.130	11.5	ng/ul	611399	6	25.025	1062380	20.0
(DEL) Alkane: S...	24.180	52.9	ng/ul	2811030	6	25.025	1062380	20.0
(DEL) Alkane: S...	26.260	22.0	ng/ul	1169350	6	25.025	1062380	20.0