

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101720\
 Data File : BG046949.D
 Acq On : 17 Oct 2020 19:39
 Operator : CG/JU
 Sample : L4201-20
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AM3

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 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-BG101520MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.479	19	24	31	rVB3	20981	36420	1.06%	0.110%
2	3.755	67	71	75	rBV5	6069	9204	0.27%	0.028%
3	4.002	110	113	116	rBV3	2887	3826	0.11%	0.012%
4	4.237	148	153	155	rBV4	3294	6217	0.18%	0.019%
5	4.302	162	164	170	rBV4	2471	3639	0.11%	0.011%
6	7.222	653	661	673	rBV	468791	836358	24.29%	2.526%
7	7.392	681	690	700	rBV	333622	537802	15.62%	1.624%
8	7.598	718	725	732	rBV	450689	761579	22.12%	2.300%
9	7.657	732	735	741	rVB	11068	17310	0.50%	0.052%
10	8.068	796	805	812	rBV	206031	351677	10.22%	1.062%
11	8.127	812	815	819	rVB3	2876	4488	0.13%	0.014%
12	8.773	917	925	936	rBV	576339	1011830	29.39%	3.056%
13	9.231	995	1003	1011	rBV	455572	797376	23.16%	2.408%
14	9.648	1070	1074	1079	rVB4	2737	3831	0.11%	0.012%
15	9.954	1117	1126	1137	rBV	425745	752223	21.85%	2.272%
16	10.495	1211	1218	1230	rBV	627786	1103184	32.05%	3.332%
17	10.882	1276	1284	1293	rVB	298547	514205	14.94%	1.553%
18	11.012	1298	1306	1316	rBV	565282	1030126	29.92%	3.111%
19	11.358	1359	1365	1368	rBV3	4545	7868	0.23%	0.024%
20	12.457	1546	1552	1560	rVB3	11399	19727	0.57%	0.060%
21	13.520	1730	1733	1736	rVB3	3342	3712	0.11%	0.011%
22	13.585	1740	1744	1752	rVB5	9080	15048	0.44%	0.045%
23	14.102	1825	1832	1836	rBV	1210341	1799041	52.26%	5.434%
24	14.143	1836	1839	1847	rVB	207320	280661	8.15%	0.848%
25	14.396	1874	1882	1891	rVB	1294892	2048023	59.49%	6.186%
26	14.584	1912	1914	1921	rVB3	2719	4040	0.12%	0.012%
27	14.695	1926	1933	1940	rVV	507328	790784	22.97%	2.388%
28	14.883	1958	1965	1980	rBV	608359	985643	28.63%	2.977%
29	15.095	1998	2001	2005	rBV3	4184	5776	0.17%	0.017%
30	15.489	2064	2068	2071	rBV4	3954	6381	0.19%	0.019%
31	15.688	2095	2102	2110	rVB	1685019	2612619	75.89%	7.891%
32	15.800	2113	2121	2132	rBV	663767	946545	27.50%	2.859%
33	15.929	2138	2143	2148	rBV	25887	36832	1.07%	0.111%
34	16.399	2219	2223	2226	rVV4	2879	4515	0.13%	0.014%

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35	16.470	2230	2235	2239	rVV6	9436	13702	0.40%	0.041%
36	16.834	2293	2297	2300	rBV3	3934	4918	0.14%	0.015%
37	16.975	2318	2321	2325	rVB2	5329	5945	0.17%	0.018%
38	17.122	2342	2346	2350	rVB5	4646	6545	0.19%	0.020%
39	17.234	2360	2365	2368	rVB3	5232	9274	0.27%	0.028%
40	17.339	2378	2383	2384	rBV	2585	4155	0.12%	0.013%
41	17.439	2394	2400	2407	rBV	667372	1005738	29.22%	3.038%
42	17.545	2410	2418	2426	rBV	1792161	2789637	81.03%	8.425%
43	17.710	2440	2446	2450	rBV	26394	33687	0.98%	0.102%
44	17.757	2450	2454	2458	rBV	49809	59451	1.73%	0.180%
45	17.980	2487	2492	2495	rBV2	2610	5168	0.15%	0.016%
46	18.103	2510	2513	2517	rVB4	4814	5482	0.16%	0.017%
47	18.321	2545	2550	2558	rBV	239790	365978	10.63%	1.105%
48	18.650	2604	2606	2612	rBV6	4198	9009	0.26%	0.027%
49	18.808	2632	2633	2635	rBV2	4508	3496	0.10%	0.011%
50	18.990	2660	2664	2665	rBV4	3722	4283	0.12%	0.013%
51	19.020	2665	2669	2672	rVV	26876	31636	0.92%	0.096%
52	19.061	2672	2676	2681	rVB	51253	61041	1.77%	0.184%
53	19.108	2682	2684	2690	rVV6	2647	4868	0.14%	0.015%
54	19.220	2699	2703	2707	rBV6	21559	25714	0.75%	0.078%
55	19.290	2712	2715	2716	rBV3	4592	4196	0.12%	0.013%
56	19.390	2730	2732	2736	rVB5	6195	6891	0.20%	0.021%
57	19.455	2738	2743	2748	rBV	27360	47558	1.38%	0.144%
58	19.590	2762	2766	2773	rBV2	86449	135595	3.94%	0.410%
59	19.678	2779	2781	2783	rBV3	3566	3510	0.10%	0.011%
60	19.707	2783	2786	2790	rVV	19204	28147	0.82%	0.085%
61	19.742	2790	2792	2795	rVB2	10324	10423	0.30%	0.031%
62	19.825	2798	2806	2816	rBV	2325769	3442538	100.00%	10.397%
63	20.048	2840	2844	2847	rVV6	7547	10731	0.31%	0.032%
64	20.136	2854	2859	2862	rVV	214549	253534	7.36%	0.766%
65	20.171	2862	2865	2871	rVB	436940	469086	13.63%	1.417%
66	20.318	2886	2890	2892	rBV5	14157	21096	0.61%	0.064%
67	20.406	2903	2905	2910	rVB6	6684	7165	0.21%	0.022%
68	20.489	2917	2919	2923	rBV5	5609	4118	0.12%	0.012%
69	20.724	2953	2959	2965	rBV	96125	121653	3.53%	0.367%
70	20.841	2974	2979	2983	rVB8	16568	26098	0.76%	0.079%
71	21.123	3023	3027	3030	rBV2	80241	106389	3.09%	0.321%

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72	21.164	3030	3034	3038	rVB	145408	183538	5.33%	0.554%
73	21.346	3059	3065	3076	rBV	371772	592447	17.21%	1.789%
74	21.711	3120	3127	3136	rBV	656445	1049555	30.49%	3.170%
75	21.905	3157	3160	3164	rVB5	22509	22808	0.66%	0.069%
76	22.228	3211	3215	3219	rBV3	30426	44718	1.30%	0.135%
77	22.275	3219	3223	3228	rVB3	60283	89993	2.61%	0.272%
78	22.439	3249	3251	3252	rBV2	8524	4443	0.13%	0.013%
79	22.510	3260	3263	3264	rBV2	23953	23970	0.70%	0.072%
80	23.015	3347	3349	3350	rBV2	8009	5650	0.16%	0.017%
81	23.209	3379	3382	3390	rVB8	31106	57861	1.68%	0.175%
82	23.638	3451	3455	3459	rBV6	29100	50238	1.46%	0.152%
83	23.697	3461	3465	3473	rVB4	52106	107594	3.13%	0.325%
84	24.020	3516	3520	3527	rVB7	31157	70282	2.04%	0.212%
85	24.737	3630	3642	3657	rVB	1115395	3050486	88.61%	9.213%
86	24.954	3669	3679	3691	rVB2	360468	1104269	32.08%	3.335%
87	25.148	3710	3712	3715	rBV3	8372	10803	0.31%	0.033%
88	26.253	3894	3900	3913	rBV9	35529	111949	3.25%	0.338%
89	26.423	3927	3929	3930	rBV2	8870	3824	0.11%	0.012%
90	27.445	4101	4103	4104	rBV2	10447	8048	0.23%	0.024%
91	28.121	4216	4218	4220	rBV3	8416	7609	0.22%	0.023%
92	32.005	4877	4879	4880	rBV2	7637	4795	0.14%	0.014%
93	32.316	4930	4932	4934	rBV3	8355	7994	0.23%	0.024%

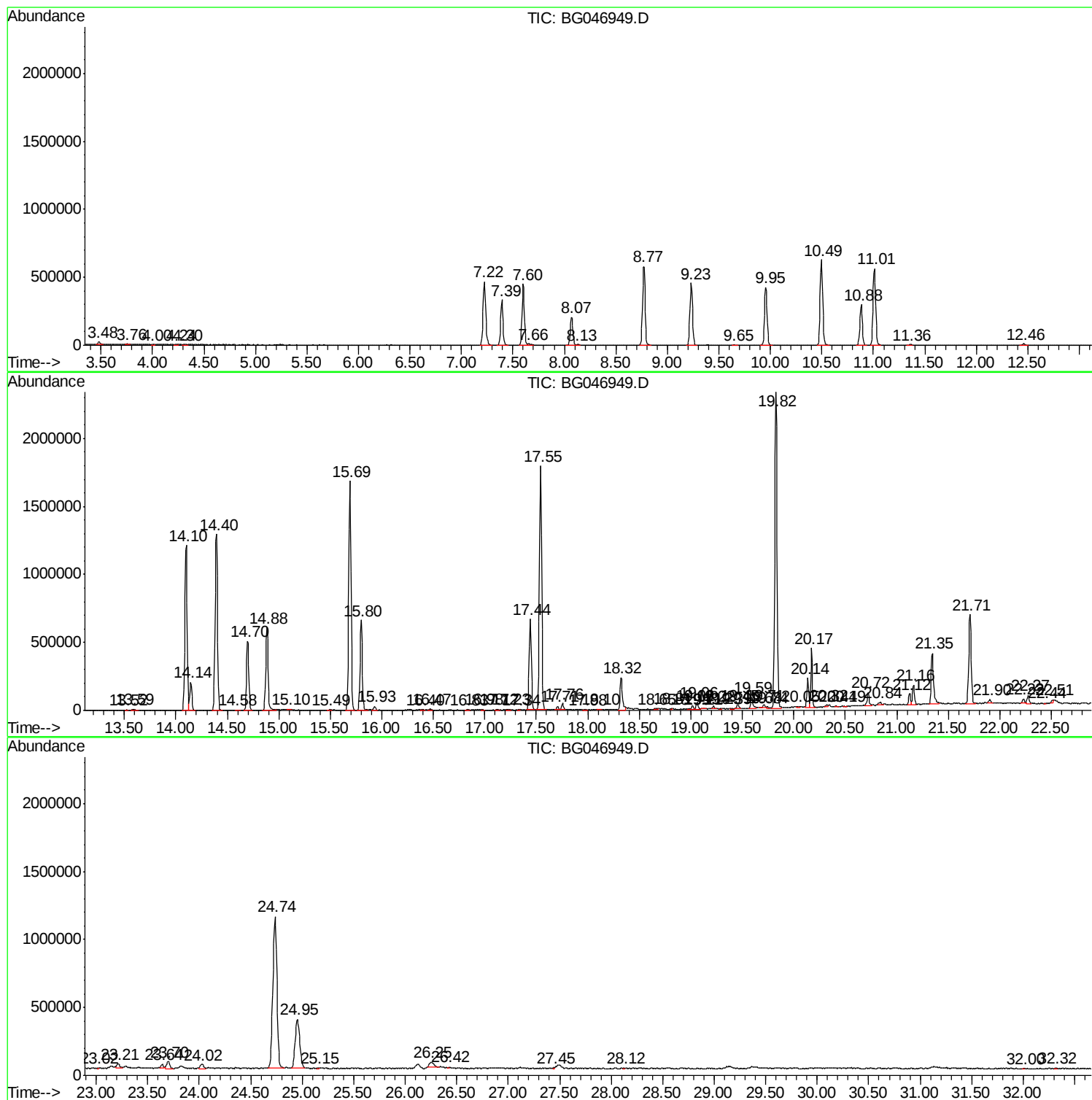
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG101520MA.M
Quant Title : SVOA CALIBRATION

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



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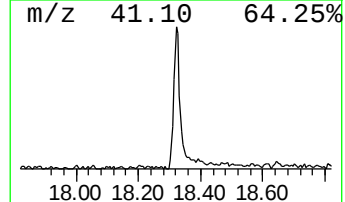
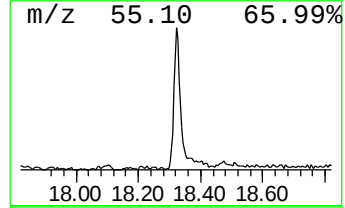
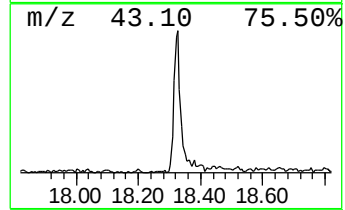
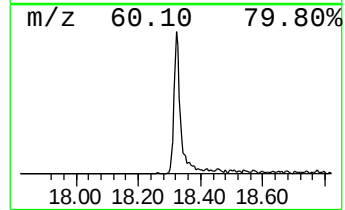
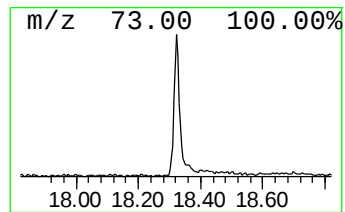
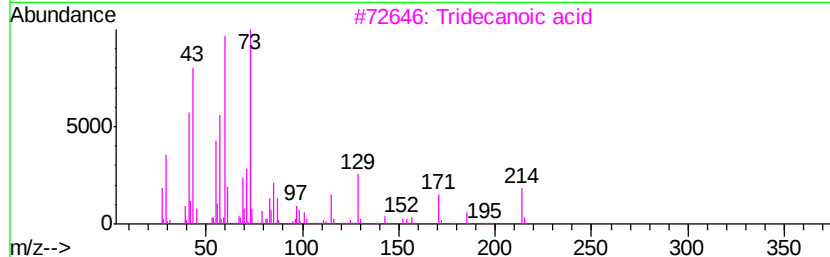
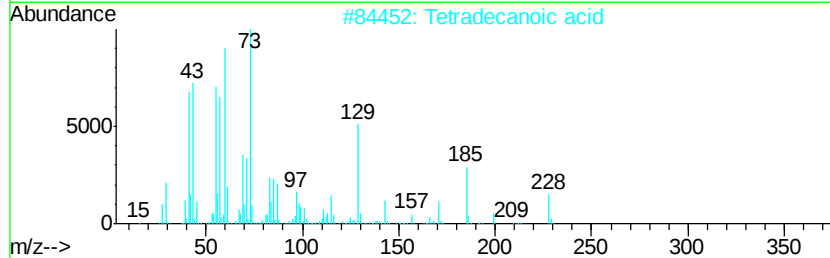
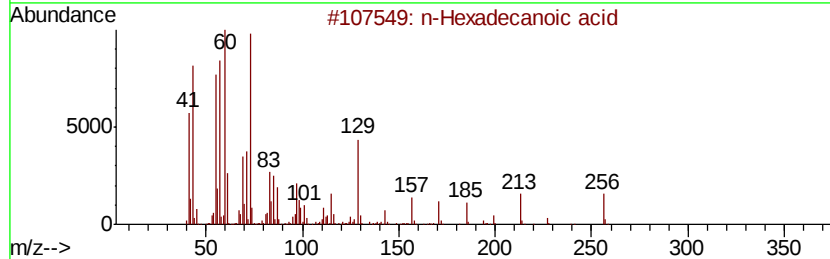
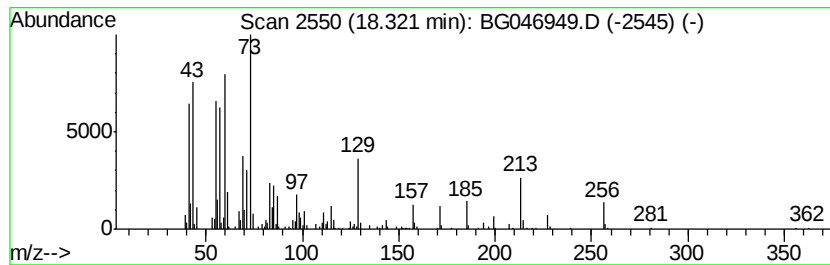
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG101520MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.32	7.28 ng/ul	365978	Phenanthrene-d10	17.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	93
3		Tridecanoic acid	214	C13H26O2	000638-53-9	86
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	76
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	72



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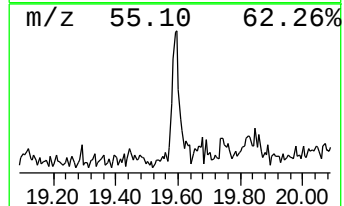
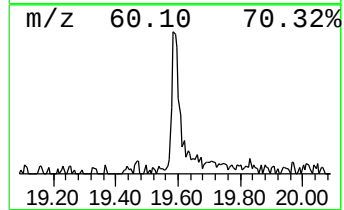
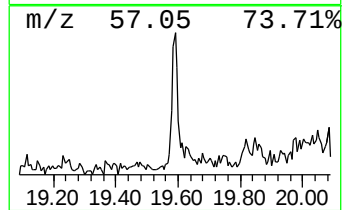
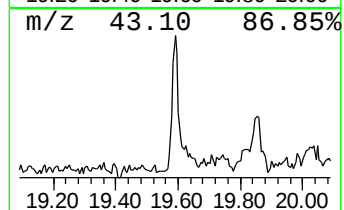
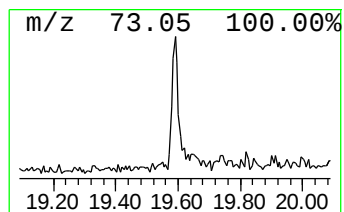
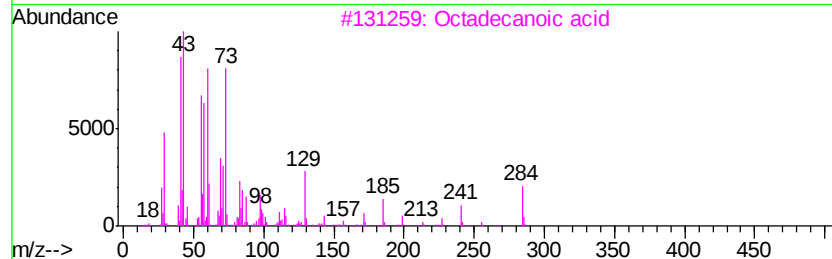
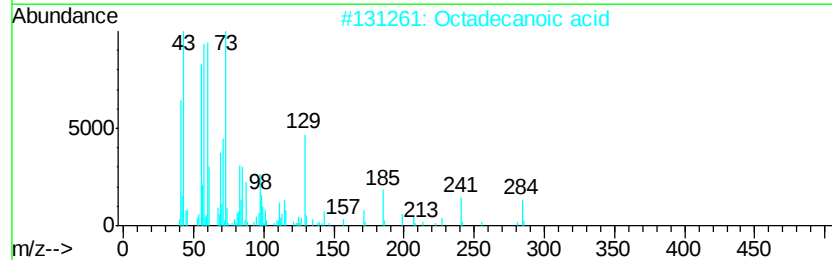
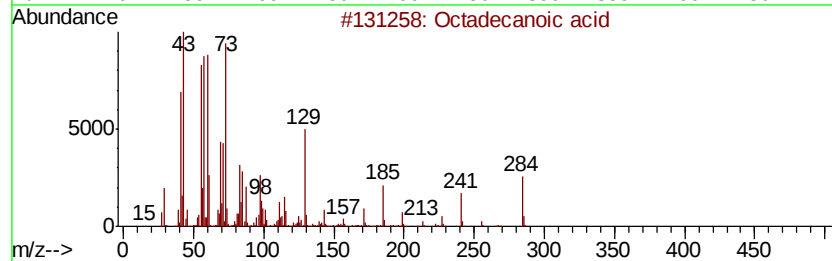
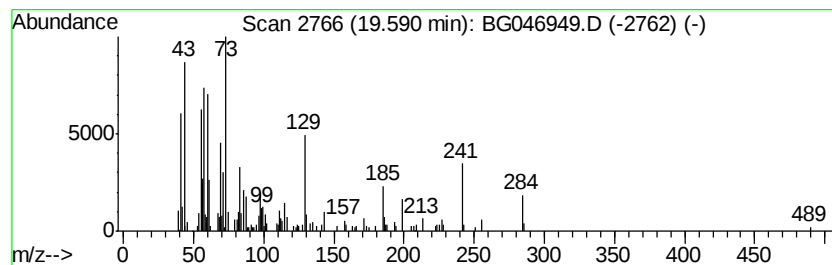
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.59	2.58 ng/ul	135595	Chrysene-d12	21.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Octadecanoic acid	284	C18H36O2	000057-11-4	98
3		Octadecanoic acid	284	C18H36O2	000057-11-4	96
4		Octadecanoic acid	284	C18H36O2	000057-11-4	95
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	89



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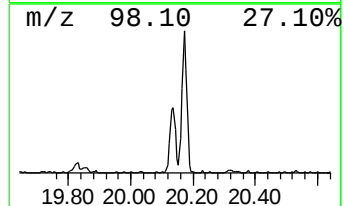
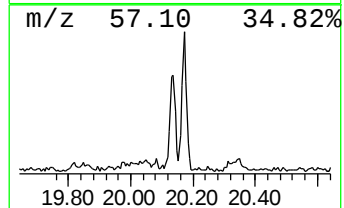
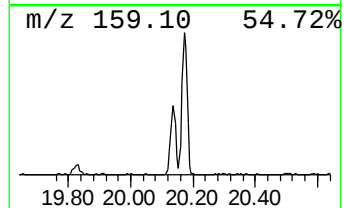
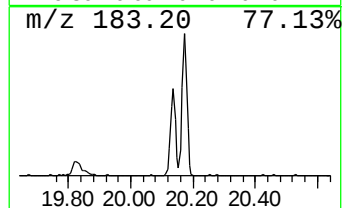
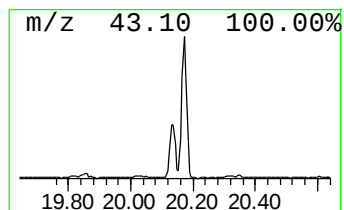
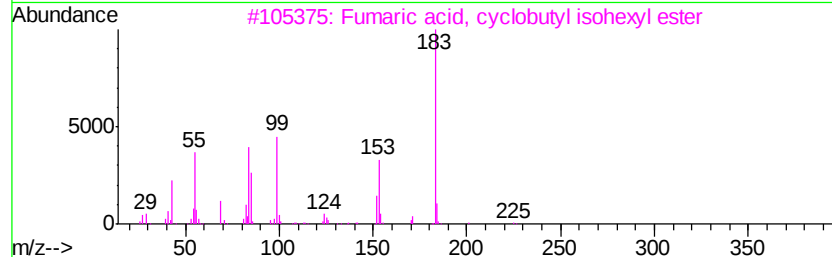
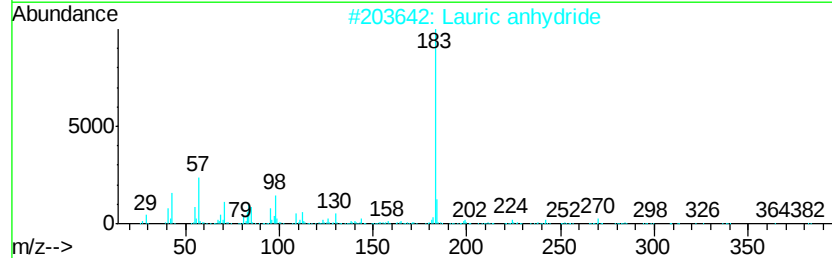
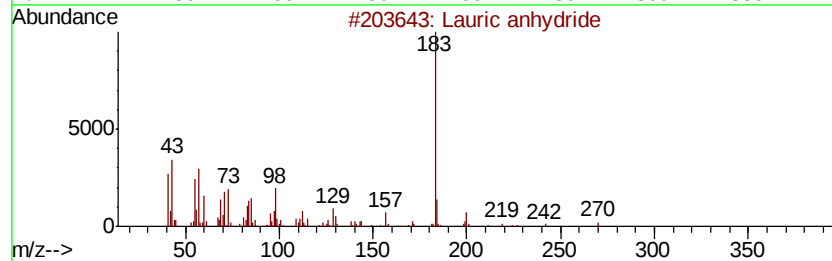
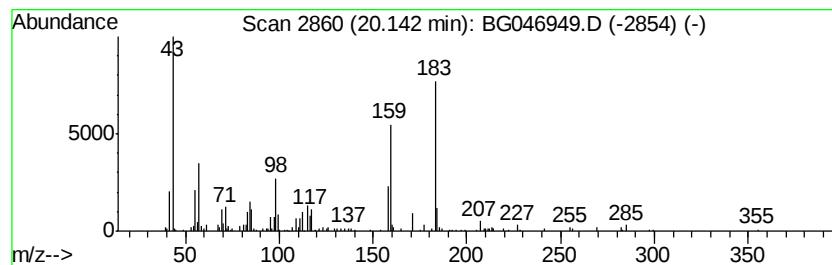
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.14	4.83 ng/ul	253534	Chrysene-d12	21.71
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Lauric anhydride	382 C24H46O3	000645-66-9	49
2	Lauric anhydride	382 C24H46O3	000645-66-9	40
3	Fumaric acid, cyclobutyl isohexyl ester	254 C14H22O4	1000345-03-5	38
4	Fumaric acid, hexyl 2-pentyl ester	270 C15H26O4	1000330-41-6	38
5	Fumaric acid, isohexyl 4-methylp...	284 C16H28O4	1000348-31-9	38



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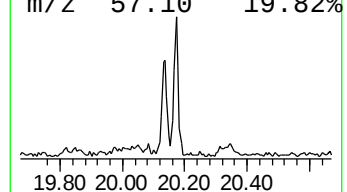
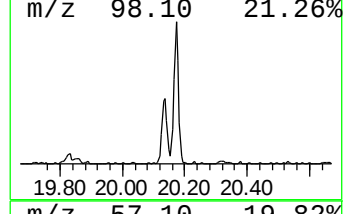
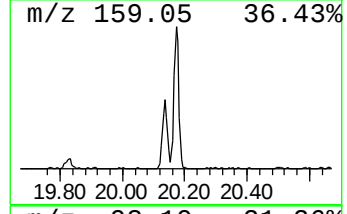
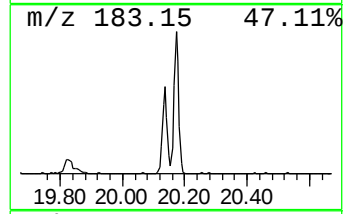
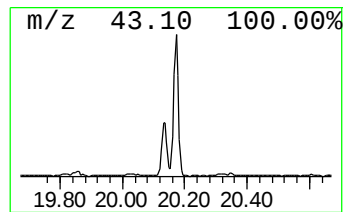
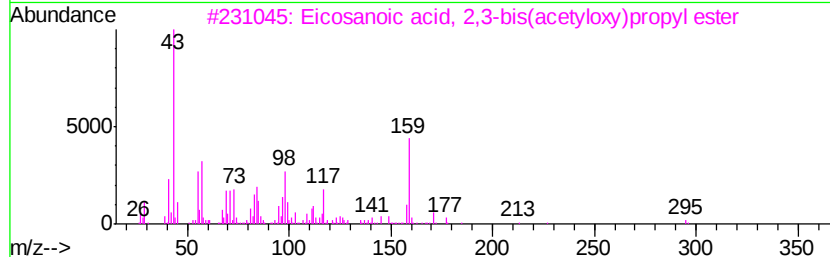
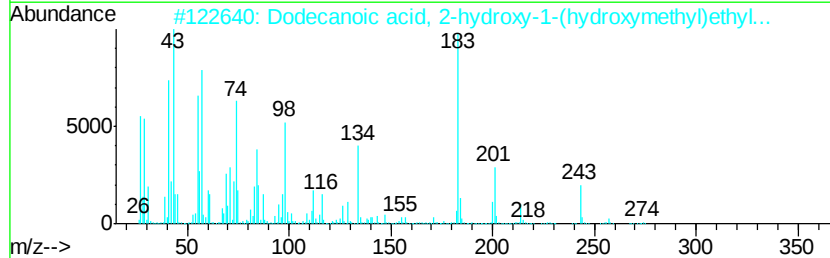
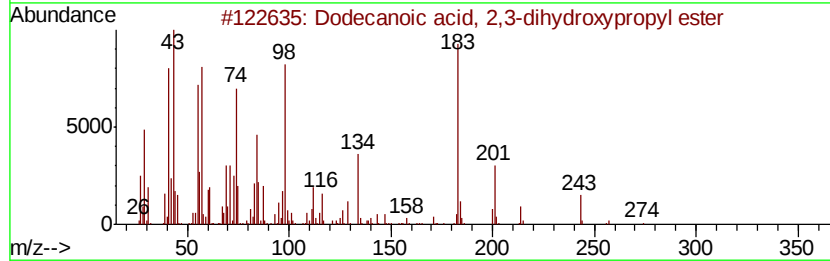
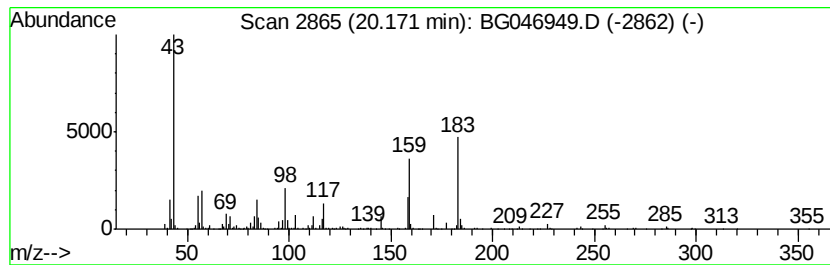
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-02 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.17	8.94 ng/ul	469086	Chrysene-d12	21.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecanoic acid, 2,3-dihydroxypr...	274	C15H30O4	000142-18-7	35
2		Dodecanoic acid, 2-hydroxy-1-(hy...	274	C15H30O4	001678-45-1	25
3		Eicosanoic acid, 2,3-bis(acetylo...	470	C27H50O6	055429-67-9	16
4		Dodecanoyl chloride	218	C12H23ClO	000112-16-3	10
5		Eicosanoic acid, 2-(acetyloxy)-1...	470	C27H50O6	055429-68-0	10



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101720\
Data File : BG046949.D
Acq On : 17 Oct 2020 19:39
Operator : CG/JU
Sample : L4201-20
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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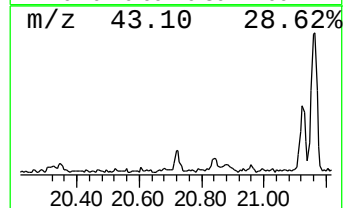
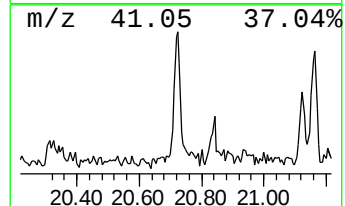
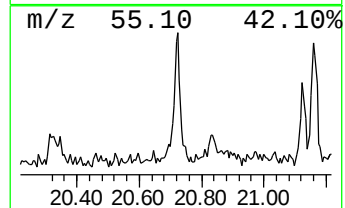
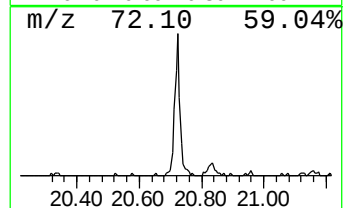
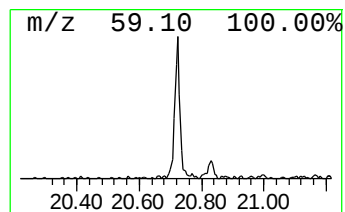
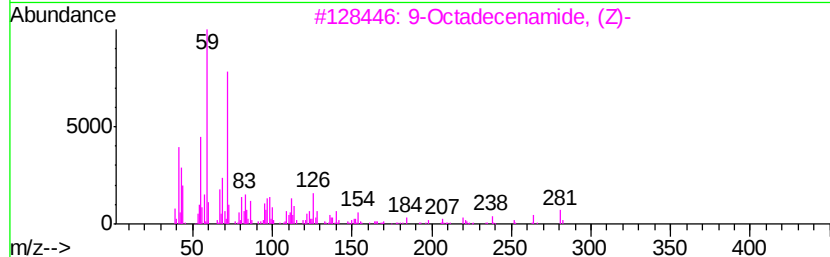
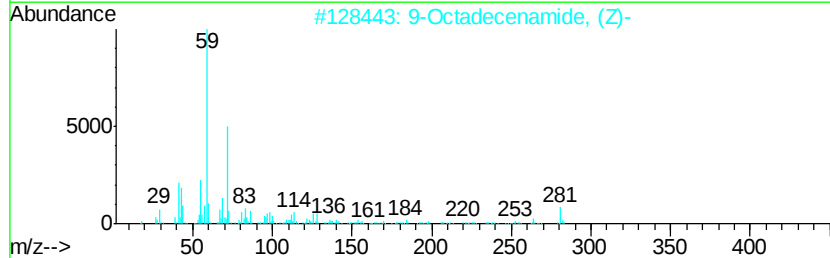
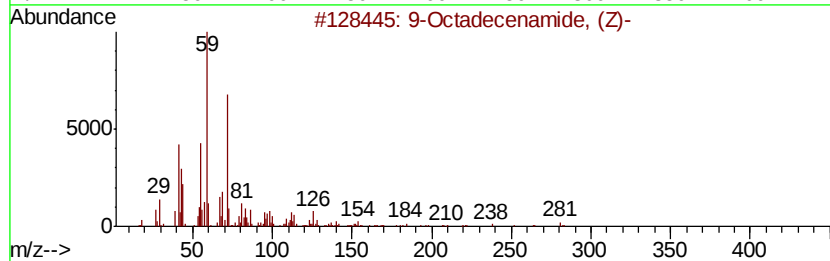
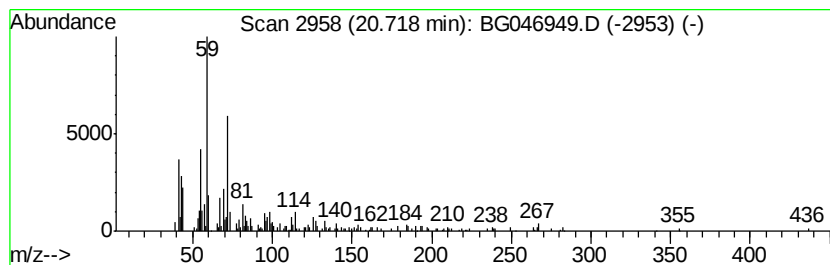
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 9-Octadecenamide, (Z)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.72	2.32 ng/ul	121653	Chrysene-d12	21.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	87
2	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	86
3	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	78
4	Decanamide-			171	C10H21NO	002319-29-1	78
5	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	78



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101720\
Data File : BG046949.D
Acq On : 17 Oct 2020 19:39
Operator : CG/JU
Sample : L4201-20
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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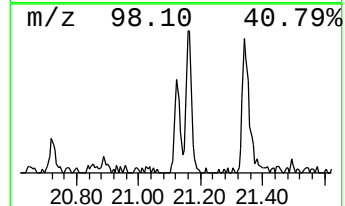
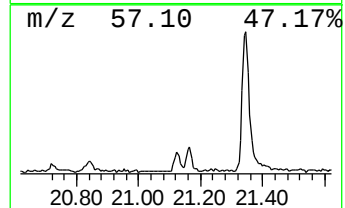
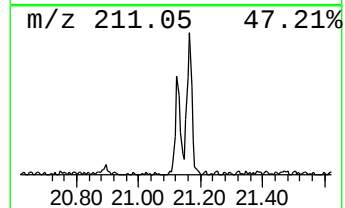
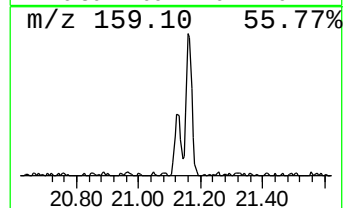
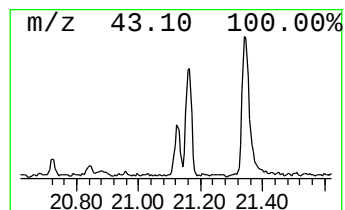
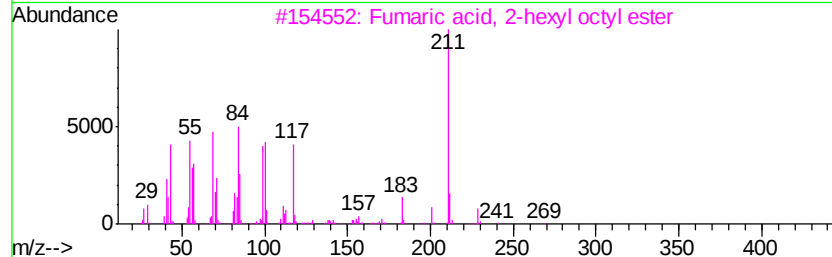
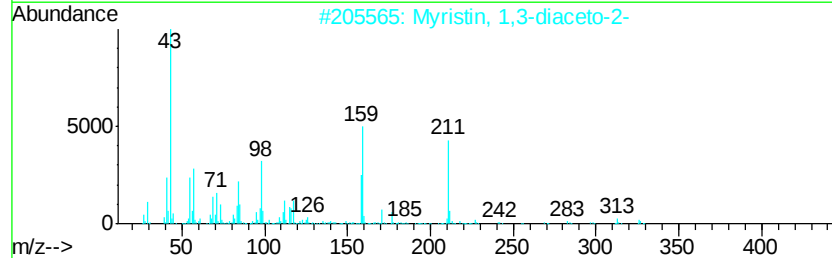
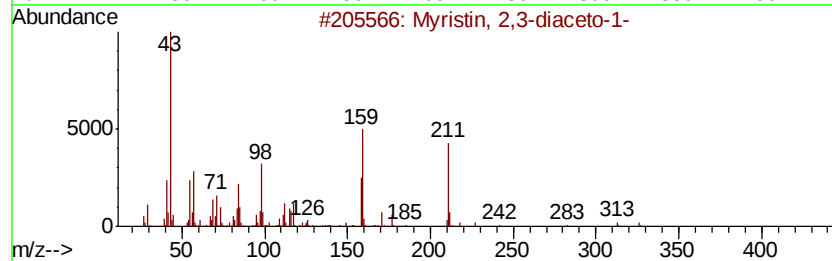
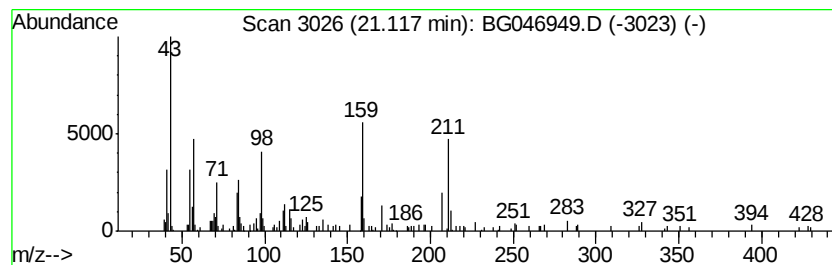
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Myristin, 2,3-diaceto-1- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.12	2.03 ng/ul	106389	Chrysene-d12	21.71

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Myristin, 2,3-diaceto-1-	386	C21H38O6	014473-55-3	90	
2	Myristin, 1,3-diaceto-2-	386	C21H38O6	014290-23-4	38	
3	Fumaric acid, 2-hexyl octyl ester	312	C18H32O4	1000348-74-6	37	
4	Fumaric acid, 4-octyl octyl ester	340	C20H36O4	1000339-21-9	35	
5	2,4-Dimethoxyphenylamine, N-meth...	211	C10H13NO4	1000343-98-3	27	



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101720\
Data File : BG046949.D
Acq On : 17 Oct 2020 19:39
Operator : CG/JU
Sample : L4201-20
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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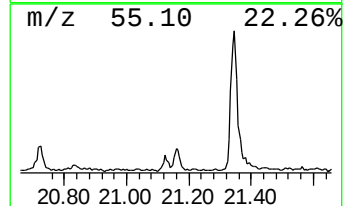
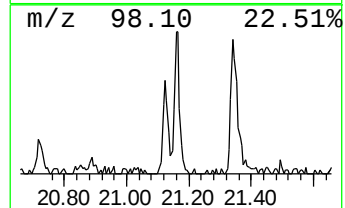
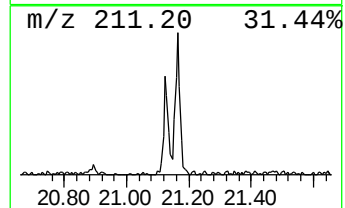
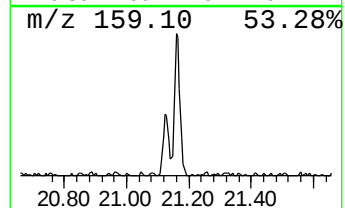
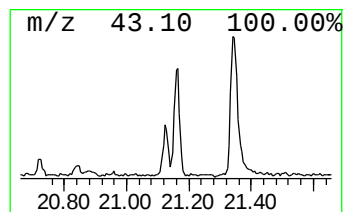
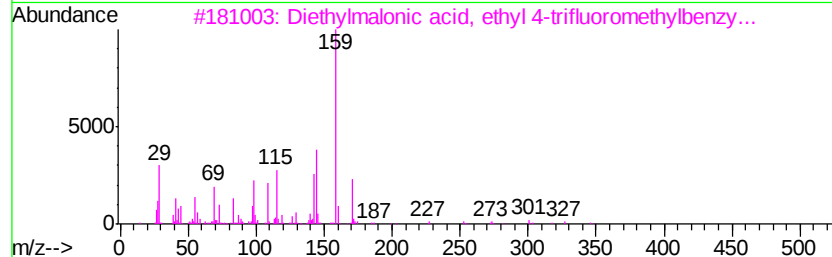
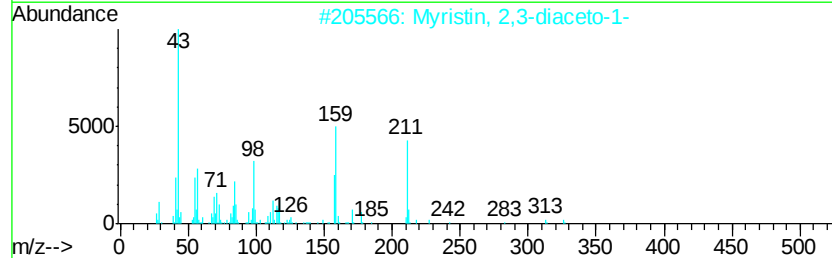
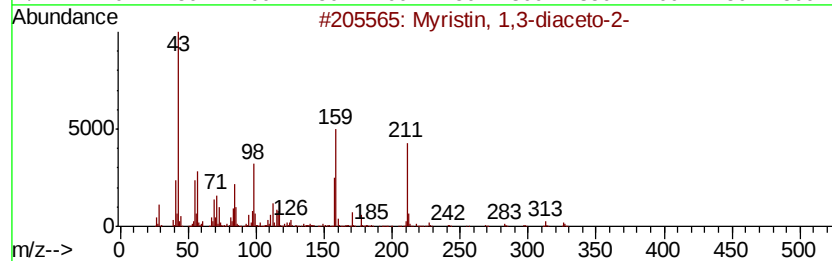
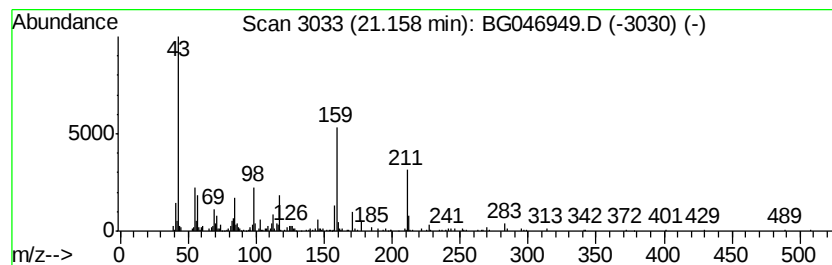
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-03 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.16	3.50 ng/ul	183538	Chrysene-d12	21.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Myristin, 1,3-diaceto-2-	386	C21H38O6	014290-23-4	38
2		Myristin, 2,3-diaceto-1-	386	C21H38O6	014473-55-3	27
3		Diethylmalonic acid, ethyl 4-tri...	346	C17H21F3O4	1000368-39-9	16
4		2,4-Difluorobenzoic acid, octade...	410	C25H40F2O2	1000338-79-6	14
5		2,6-Difluorobenzoic acid, octade...	410	C25H40F2O2	1000292-39-7	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101720\
Data File : BG046949.D
Acq On : 17 Oct 2020 19:39
Operator : CG/JU
Sample : L4201-20
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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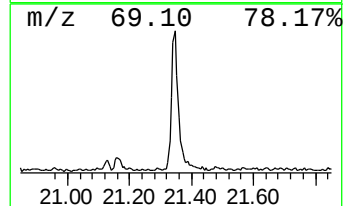
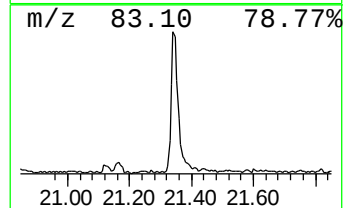
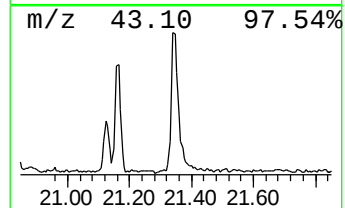
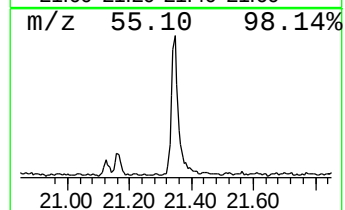
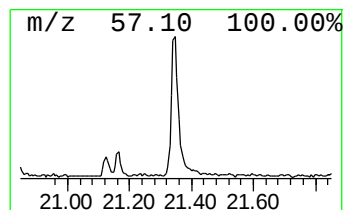
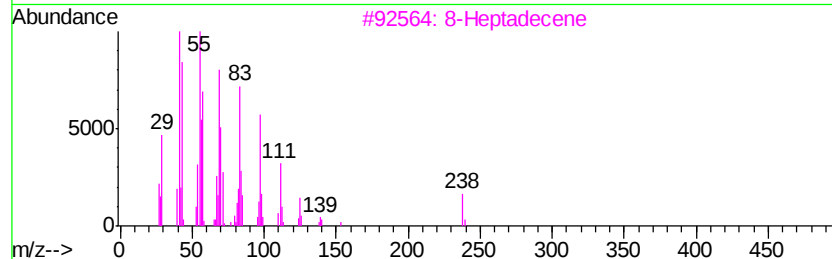
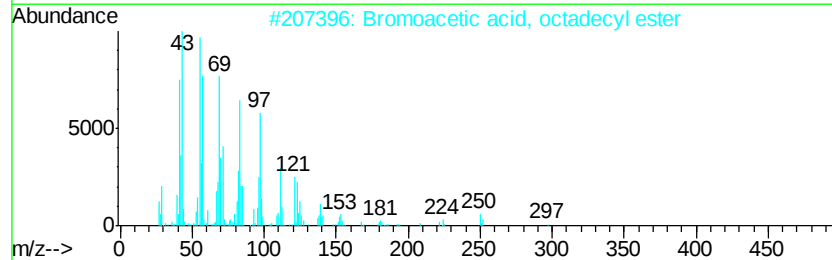
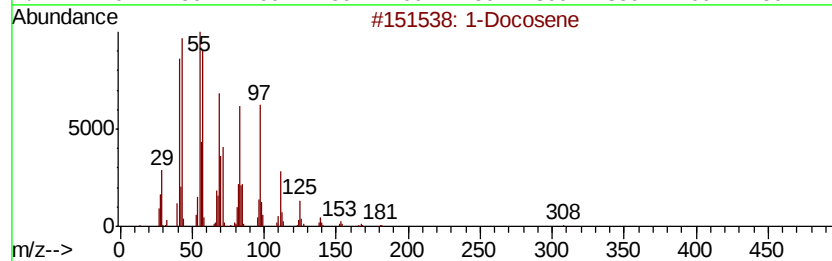
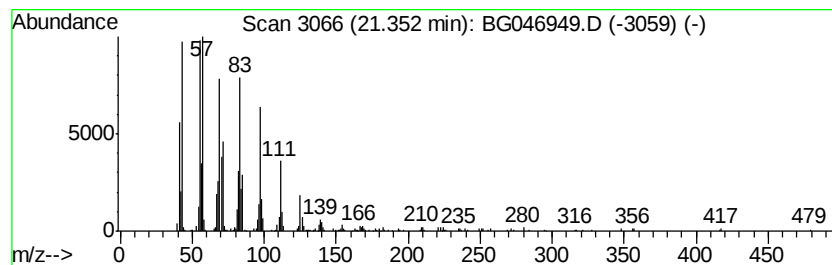
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.35	11.29 ng/ul	592447	Chrysene-d12	21.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Docosene	308	C22H44	001599-67-3	99
2		Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	94
3		8-Heptadecene	238	C17H34	002579-04-6	93
4		3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	91
5		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101720\
Data File : BG046949.D
Acq On : 17 Oct 2020 19:39
Operator : CG/JU
Sample : L4201-20
Misc :
ALS Vial : 15 Sample Multiplier: 1

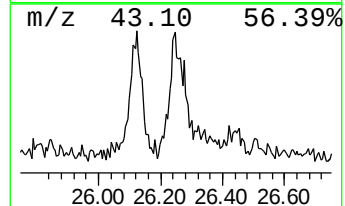
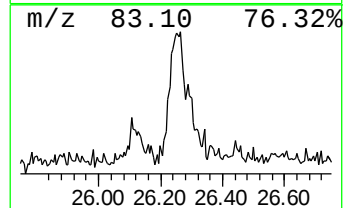
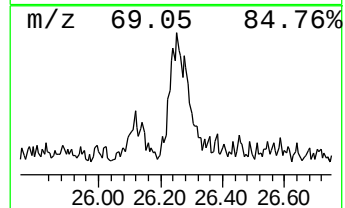
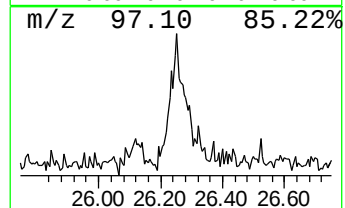
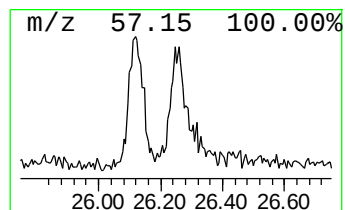
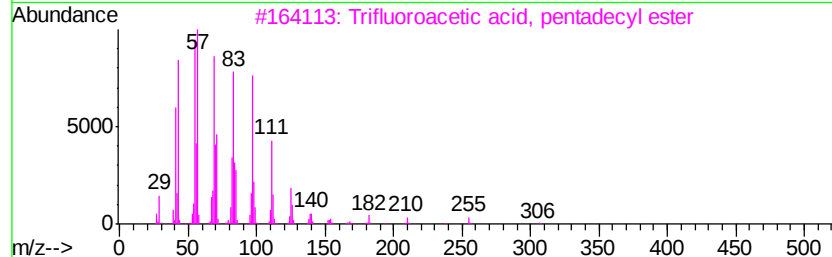
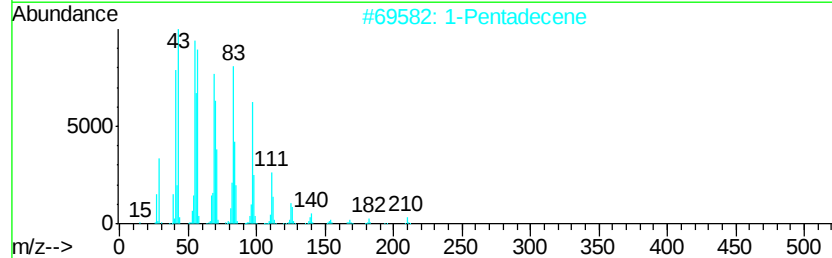
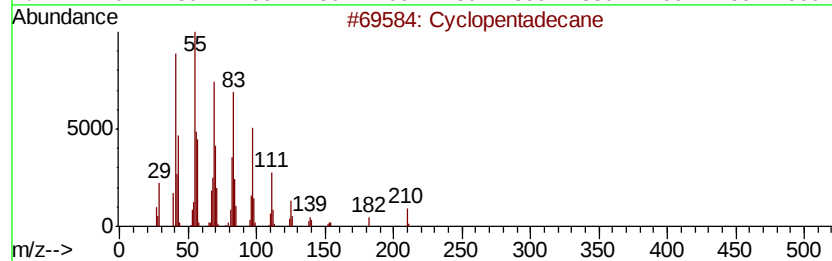
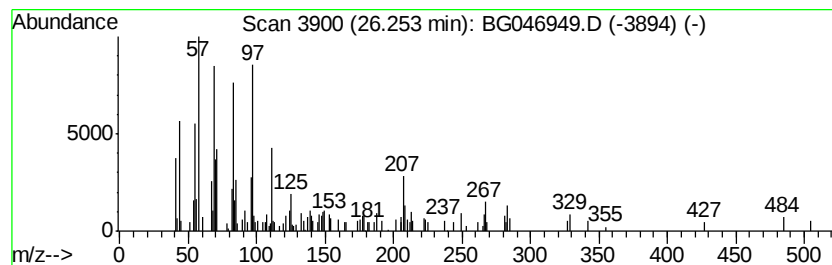
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG101520MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Cyclic26.25 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.25	2.03 ng/ul	111949	Perylene-d12	24.95
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclopentadecane	210 C15H30	000295-48-7 81
2		1-Pentadecene	210 C15H30	013360-61-7 76
3		Trifluoroacetic acid, pentadecyl...	324 C17H31F3O2	959010-23-2 74
4		Trifluoroacetoxv hexadecane	338 C18H33F3O2	006222-03-3 72
5		Z-5-Nonadecene	266 C19H38	1000131-11-8 60



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG101720\
Data File : BG046949.D
Acq On : 17 Oct 2020 19:39
Operator : CG/JU
Sample : L4201-20
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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
n-Hexadecanoic acid	18.32	7.3	ng/ul	365978	4	17.44	1005740	20.0
Octadecanoic acid	19.59	2.6	ng/ul	135595	5	21.71	1049560	20.0
unknown-01	20.14	4.8	ng/ul	253534	5	21.71	1049560	20.0
unknown-02	20.17	8.9	ng/ul	469086	5	21.71	1049560	20.0
9-Octadecenamide,...	20.72	2.3	ng/ul	121653	5	21.71	1049560	20.0
Myristin, 2,3-dia...	21.12	2.0	ng/ul	106389	5	21.71	1049560	20.0
unknown-03	21.16	3.5	ng/ul	183538	5	21.71	1049560	20.0
(DEL) Alkane: Str...	21.35	11.3	ng/ul	592447	5	21.71	1049560	20.0
(DEL) Alkane: Cyc...	26.25	2.0	ng/ul	111949	6	24.95	1104270	20.0