LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112717\
Data File : BM013120.D
Acq On : 27 Nov 2017 20:49
Operator : SJ/JU
Sample : I6496-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AB4

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 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:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.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	4.863	297	302	306	rBV	65838	87527	2.12%	0.133%
2	6.893	636	647	661	rVB2	636536	1167659	28.28%	1.778%
3	7.051	665	674	684	rVV	472417	744838	18.04%	1.134%
4	7.246	700	707	714	rBV	637110	1018897	24.68%	1.552%
5	7.557	755	760	766	rVB	70181	103398	2.50%	0.157%
6	7.710	778	786	793	rBV	888351	1434640	34.75%	2.185%
7	7.846	802	809	814	rVB	28241	43331	1.05%	0.066%
8	8.422	900	907	915	rBV	758950	1197159	29.00%	1.823%
9	8.857	972	981	989	rBV	657635	1083031	26.23%	1.649%
10	9.281	1047	1053	1058	rVB2	31207	45388	1.10%	0.069%
11	9.575	1096	1103	1110	rBV	552809	913291	22.12%	1.391%
12	10.116	1187	1195	1203	rBV	863956	1457402	35.30%	2.220%
13	10.487	1249	1258	1265	rBV	1280906	2121605	51.39%	3.231%
14	10.628	1272	1282	1291	rBV	446068	798932	19.35%	1.217%
15	10.851	1315	1320	1326	rBV6	25884	55295	1.34%	0.084%
16	12.692	1623	1633	1640	rBV4	39631	66856	1.62%	0.102%
17	12.798	1646	1651	1657	rBV5	28681	47131	1.14%	0.072%
18	13.081	1692	1699	1707	rBV3	100704	168116	4.07%	0.256%
19	13.175	1711	1715	1719	rBV3	32573	47900	1.16%	0.073%
20	13.222	1719	1723	1728	rVV4	61532	86187	2.09%	0.131%
21	13.669	1794	1799	1806	rBV7	26898	44335	1.07%	0.068%
22	13.763	1809	1815	1819	rVV	1689661	2323074	56.27%	3.538%
23	13.810	1819	1823	1831	rVB	288346	398036	9.64%	0.606%
24	14.039	1851	1862	1872	rBV	1828122	2685390	65.04%	4.090%
25	14.257	1894	1899	1903	rVB	78526	101544	2.46%	0.155%
26	14.345	1906	1914	1923	rBV2	1921749	2959226	71.67%	4.507%
27	14.569	1946	1952	1967	rBV	756101	1310157	31.73%	1.995%
28	14.716	1972	1977	1982	rVB7	33558	59005	1.43%	0.090%
29	14.775	1982	1987	1991	rBV7	35782	52829	1.28%	0.080%
30	15.157	2048	2052	2058	rBV5	19916	44723	1.08%	0.068%
31	15.210	2058	2061	2067	rVB	84345	100091	2.42%	0.152%
32	15.339	2077	2083	2091	rVB	2229336	3197241	77.44%	4.869%
33	15.463	2097	2104	2111	rBV	661596	957984	23.20%	1.459%
34	15.580	2120	2124	2129	rBV6	30103	58612	1.42%	0.089%

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Operator : SJ/JU
Sample : I6496-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AB4

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 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:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Title : SVOA CALIBRATION

35	16.074	2204	2208	2215	rVB	82262	117630	2.85%	0.179%
36	16.863	2338	2342	2348	rBV2	50546	70928	1.72%	0.108%
37	17.004	2361	2366	2372	rBV8	47695	76420	1.85%	0.116%
38	17.092	2373	2381	2387	rBV2	2552244	3662968	88.72%	5.579%
39	17.192	2392	2398	2409	rVB	2598215	3583607	86.80%	5.458%
40	17.380	2427	2430	2436	rVB7	31770	49457	1.20%	0.075%
41	17.880	2509	2515	2521	rBV5	47572	103379	2.50%	0.157%
42	17.939	2521	2525	2530	rVV7	56034	83796	2.03%	0.128%
43	18.004	2530	2536	2542	rVV2	966697	1666747	40.37%	2.538%
44	18.051	2542	2544	2552	rVB2	320738	386972	9.37%	0.589%
45	18.168	2559	2564	2570	rVB9	38492	68134	1.65%	0.104%
46	18.298	2581	2586	2591	rBV4	38586	67762	1.64%	0.103%
47	18.445	2605	2611	2620	rBV8	97461	178923	4.33%	0.272%
48	18.880	2681	2685	2690	rVB8	72086	98445	2.38%	0.150%
49	18.962	2694	2699	2705	rVB	197795	261247	6.33%	0.398%
50	19.021	2705	2709	2714	rBV8	33411	41670	1.01%	0.063%
51	19.133	2718	2728	2729	rBV5	107494	227349	5.51%	0.346%
52	19.162	2729	2733	2736	rVV4	132477	239075	5.79%	0.364%
53	19.186	2736	2737	2744	rVB6	71983	119222	2.89%	0.182%
54	19.286	2749	2754	2765	rBV	738008	991208	24.01%	1.510%
55	19.486	2782	2788	2795	rBV	3050013	4128687	100.00%	6.288%
56	19.886	2850	2856	2863	rBV2	199219	321368	7.78%	0.489%
57	20.009	2874	2877	2881	rBV6	62124	79289	1.92%	0.121%
58	20.051	2881	2884	2891	rVB7	56565	73609	1.78%	0.112%
59	20.151	2898	2901	2909	rBV10	22527	65074	1.58%	0.099%
60	20.221	2909	2913	2920	rBV3	184916	276261	6.69%	0.421%
61	20.351	2930	2935	2941	rBV	654019	813278	19.70%	1.239%
62	20.580	2970	2974	2980	rBV7	38844	79068	1.92%	0.120%
63	20.945	3033	3036	3040	rBV5	49950	65825	1.59%	0.100%
64	20.998	3041	3045	3051	rVV	909954	1146233	27.76%	1.746%
65	21.245	3083	3087	3088	rBV2	125351	144647	3.50%	0.220%
66	21.274	3088	3092	3103	rVB	3145724	4010991	97.15%	6.109%
67	21.445	3118	3121	3127	rVB7	79645	114543	2.77%	0.174%
68	21.880	3190	3195	3205	rBV3	640296	1255597	30.41%	1.912%
69	22.080	3226	3229	3235	rVB2	160521	215176	5.21%	0.328%
70	22.333	3270	3272	3276	rVB3	105139	131563	3.19%	0.200%
71	22.562	3309	3311	3318	rVB4	147696	192882	4.67%	0.294%

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Instrument : BNA_M
ClientSampleId : C0AB4

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Title : SVOA CALIBRATION

72	22.845	3354	3359	3363	rBV	916742	1287687	31.19%	1.961%
73	22.886	3363	3366	3374	rVB	480386	731858	17.73%	1.115%
74	23.362	3440	3447	3453	rBV2	1712407	3157367	76.47%	4.809%
75	23.503	3464	3471	3479	rVB	1554669	2932714	71.03%	4.466%
76	23.721	3505	3508	3514	rVB4	145506	234336	5.68%	0.357%
77	24.056	3559	3565	3571	rVB2	419377	788102	19.09%	1.200%
78	24.139	3573	3579	3591	rVB2	522354	1118668	27.10%	1.704%
79	25.597	3820	3827	3834	rBV2	493262	1144753	27.73%	1.743%
80	26.544	3981	3988	4001	rVB6	431069	1326823	32.14%	2.021%
81	26.962	4055	4059	4069	rVB9	288648	748295	18.12%	1.140%

Sum of corrected areas: 65660463

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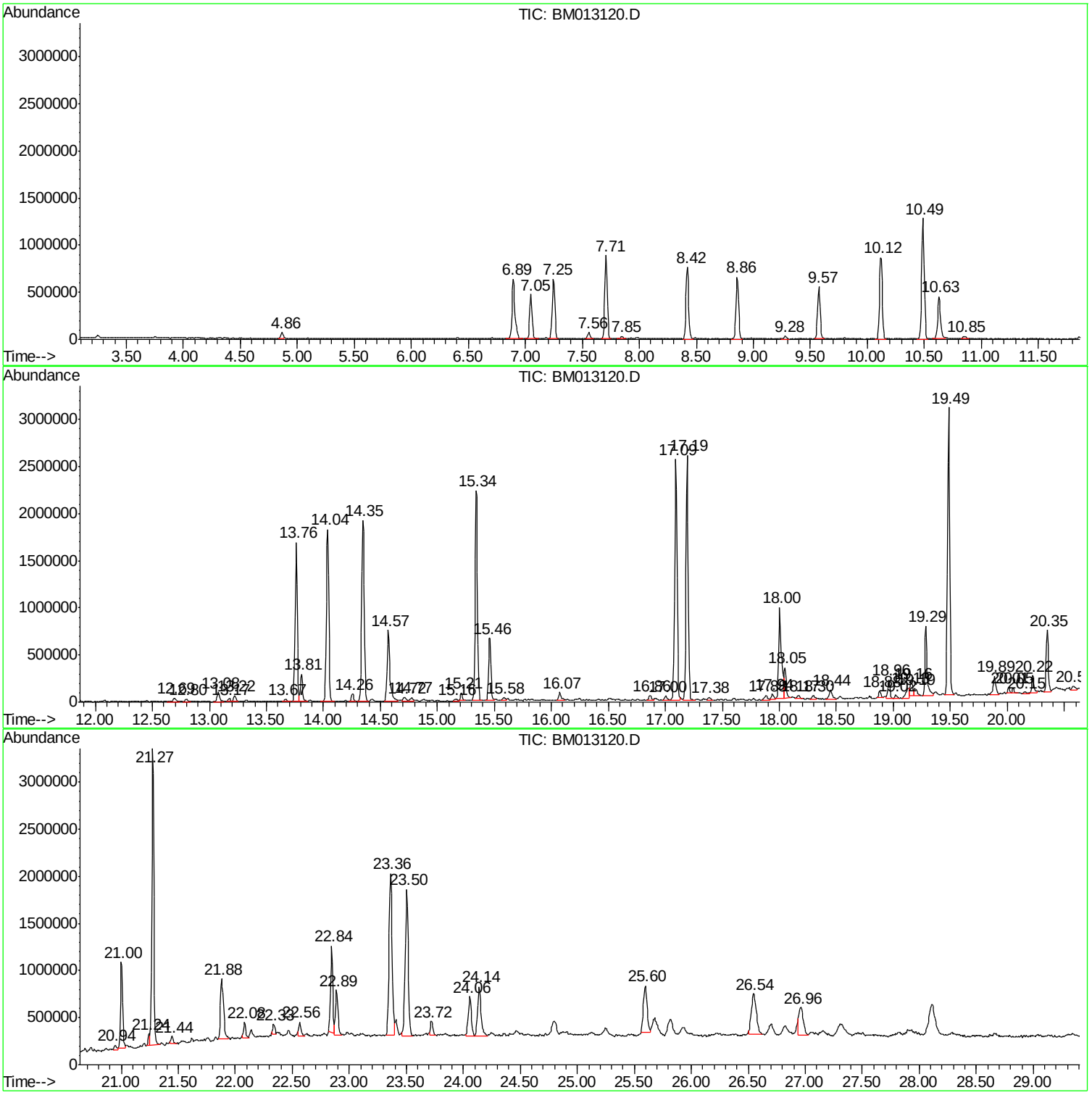
Instrument : BNA_M
ClientSampleId : C0AB4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112717\
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Acq On : 27 Nov 2017 20:49
Operator : SJ/JU
Sample : I6496-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AB4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

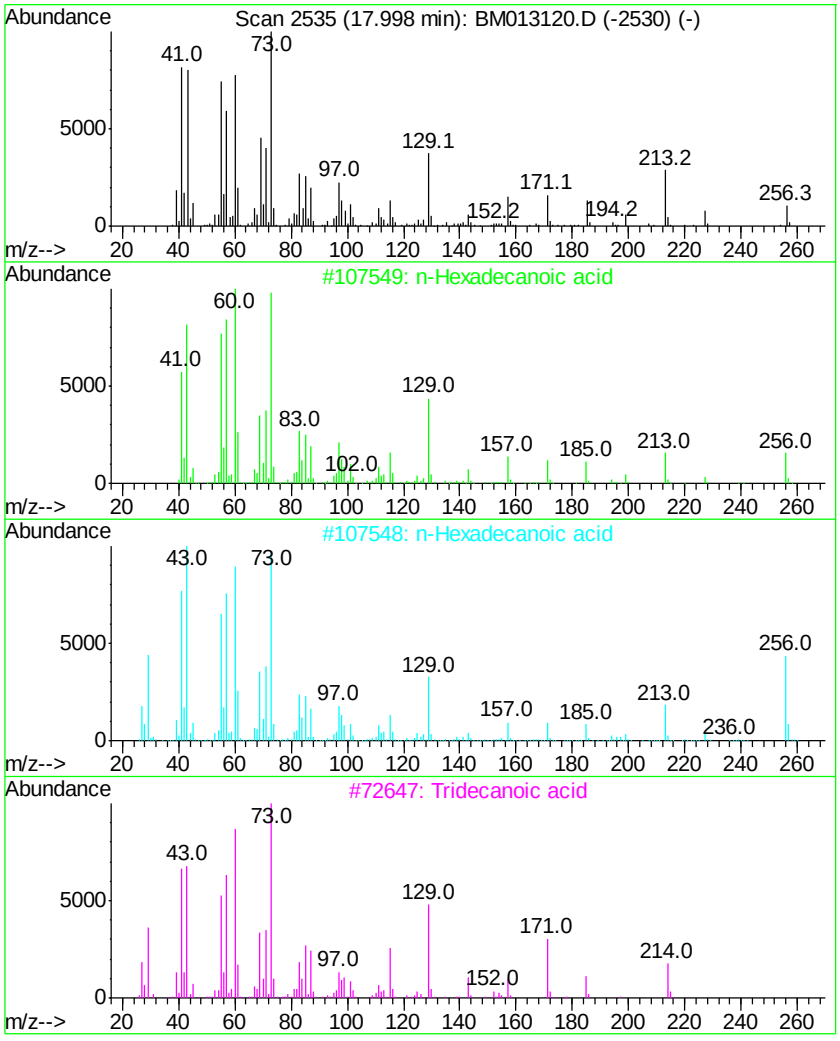
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

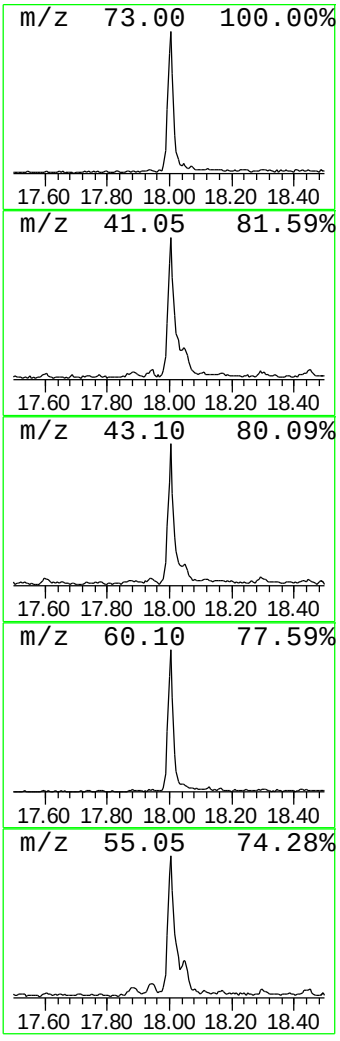
****

Peak Number 1 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.00	9.10 ng/ul	1666750	Phenanthrene-d10	17.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
3	Tridecanoic acid	214	C13H26O2	000638-53-9	95
4	Tridecanoic acid	214	C13H26O2	000638-53-9	94
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	94





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Sample : I6496-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AB4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

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

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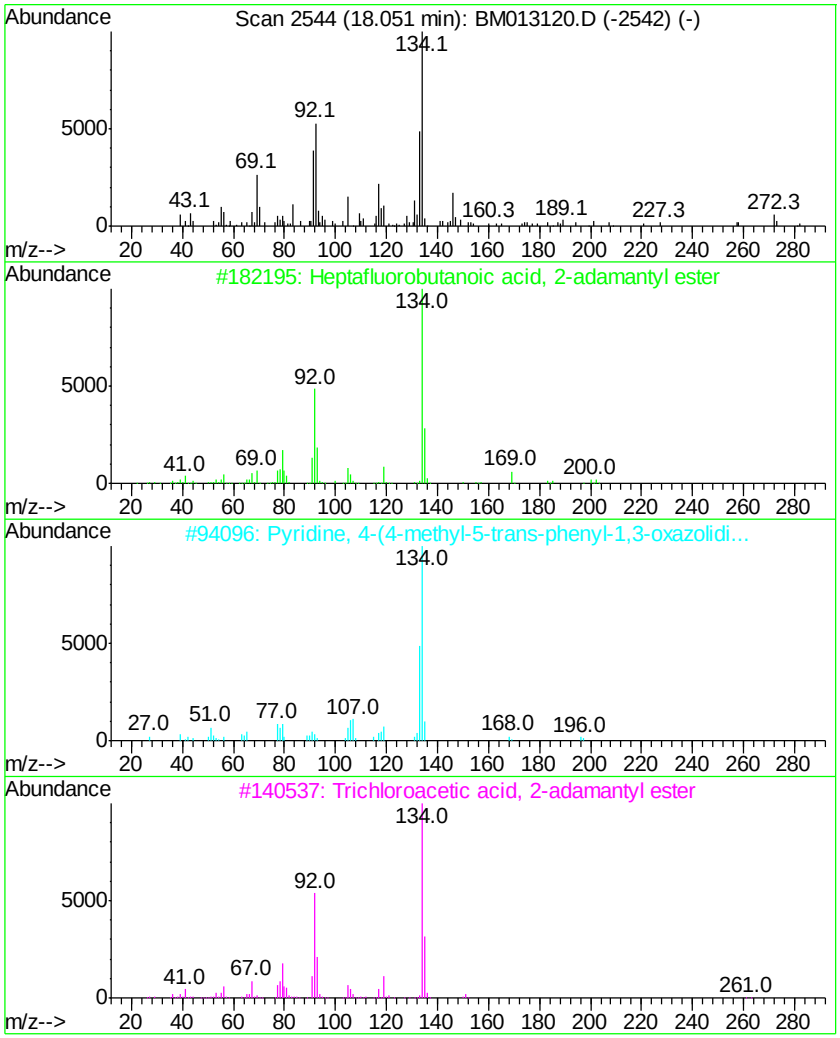
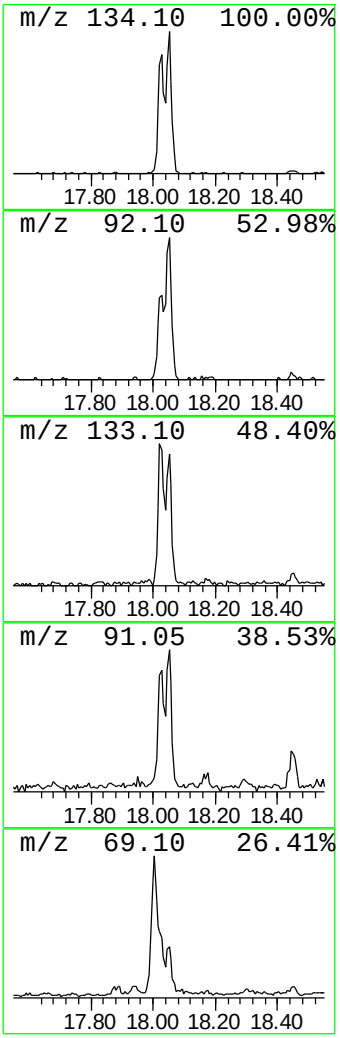
Peak Number 2 unknown-01 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.05	2.11 ng/ul	386972	Phenanthrene-d10	17.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptafluorobutanoic acid, 2-adam...	348	C14H15F7O2	1000282-96-8	43
2	Pyridine, 4-(4-methyl-5-trans-ph...	240	C15H16N2O	1000142-75-4	42
3	Trichloroacetic acid, 2-adamanty...	296	C12H15Cl3O2	1000282-91-9	38
4	Pentafluoropropionic acid, 2-ada...	298	C13H15F5O2	1000283-03-7	37
5	6-Bromohexanoic acid, 2-adamanty...	328	C16H25BrO2	1000282-90-1	37





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112717\
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Acq On : 27 Nov 2017 20:49
Operator : SJ/JU
Sample : I6496-04
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ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AB4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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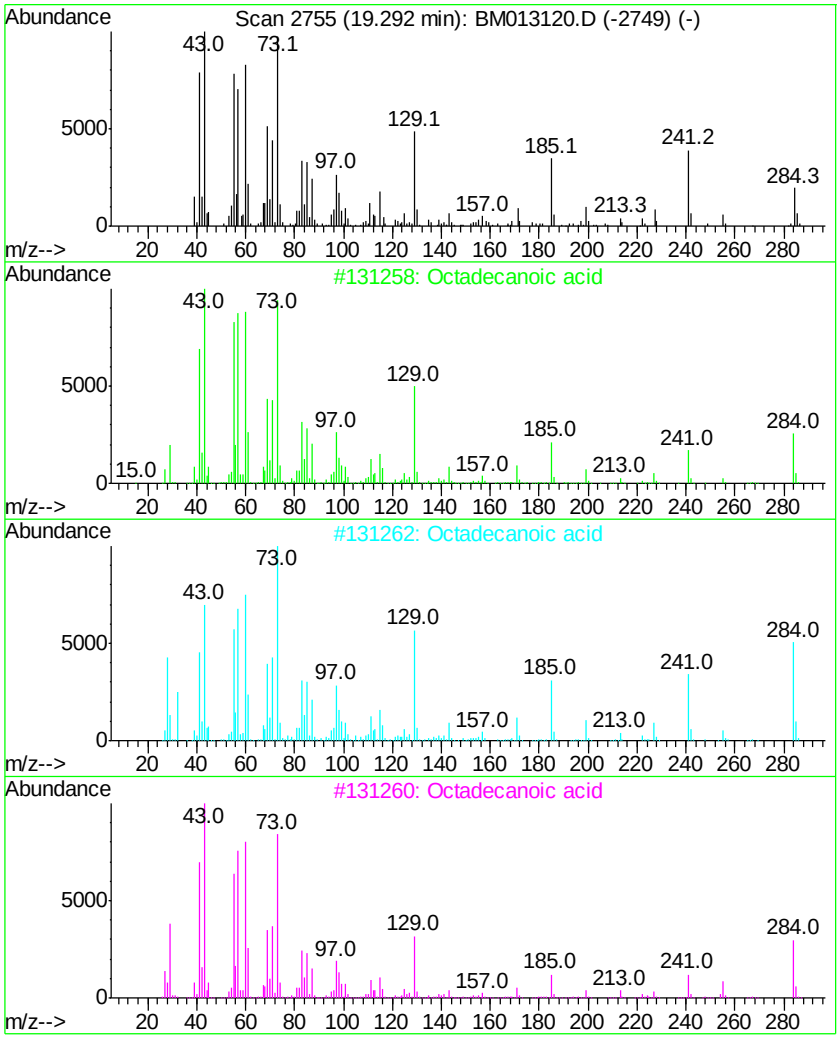
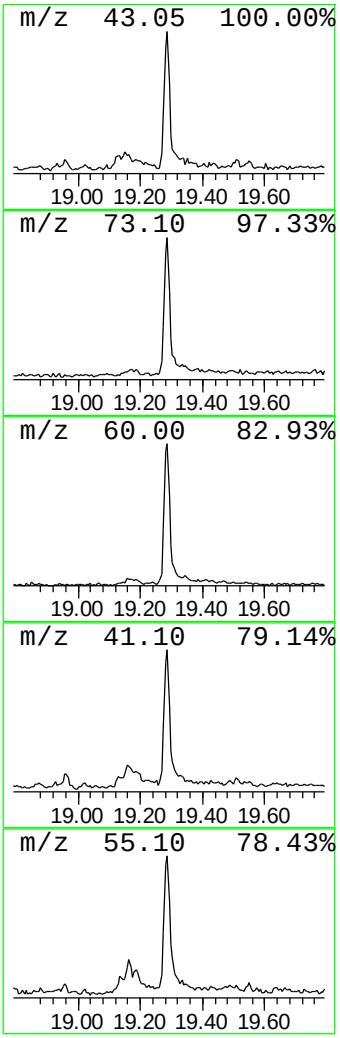
Peak Number 3 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.29	4.94 ng/ul	991208	Chrysene-d12	21.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	97
3	Octadecanoic acid	284	C18H36O2	000057-11-4	93
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
5	n-Decanoic acid	172	C10H20O2	000334-48-5	78





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Instrument : BNA_M
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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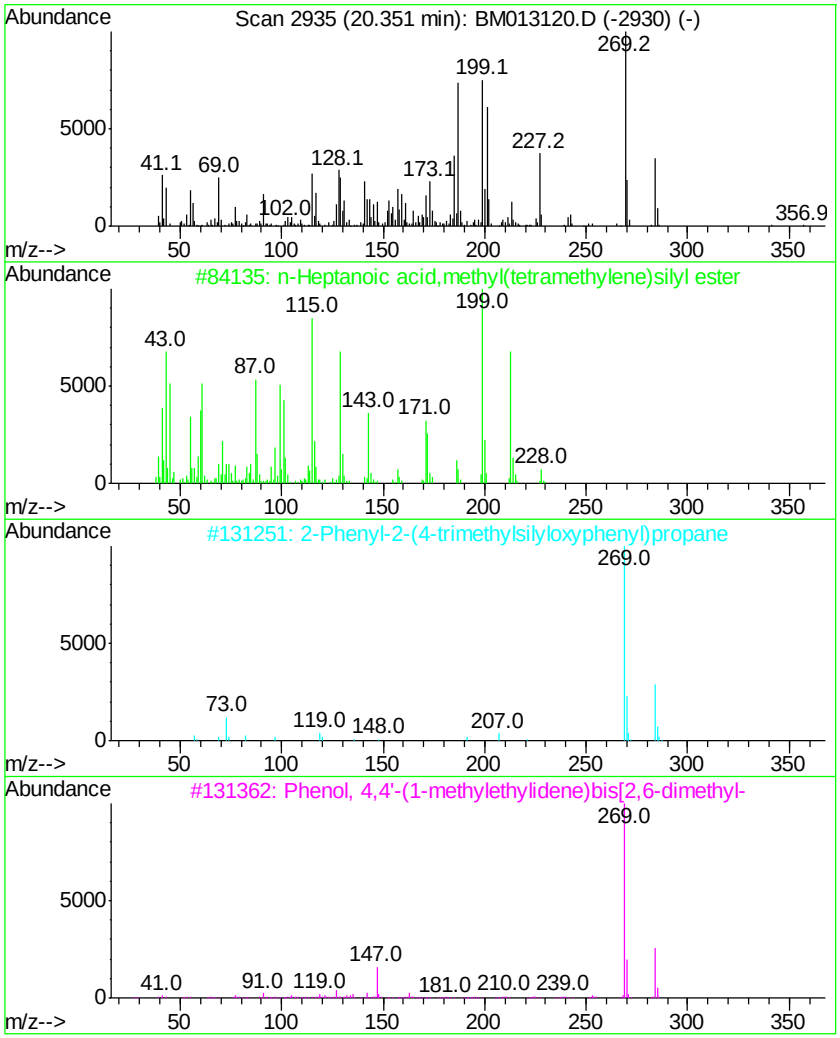
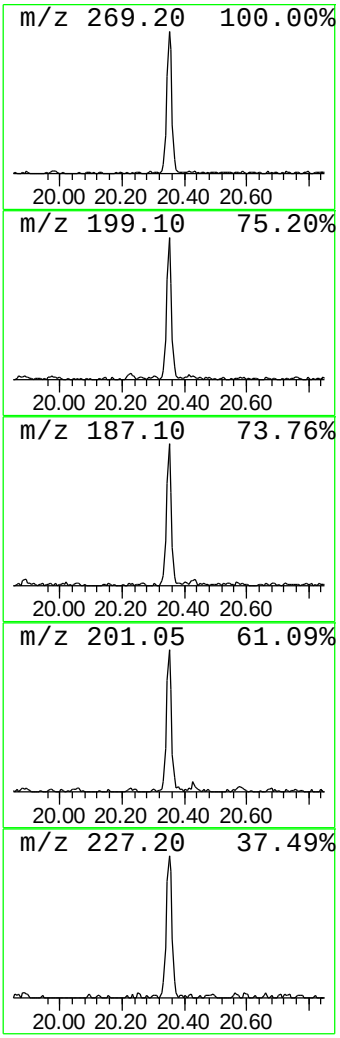
Peak Number 4 unknown-02 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.35	4.06 ng/ul	813278	Chrysene-d12	21.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Heptanoic acid,methyl(tetramet...	228	C12H24O2Si	1000217-03-6	25
2	2-Phenyl-2-(4-trimethylsilyloxyp...	284	C18H24OSi	261502-93-6	18
3	Phenol, 4,4'-(1-methylethylidene...	284	C19H24O2	005613-46-7	18
4	1,3-Dimethyl-5-(3-methyl-1H-indo...	269	C15H15N3O2	066723-75-9	15
5	(tert-Butyl)(4,6-diimidazol-1-yl...	284	C13H16N8	1000304-62-9	14





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Instrument : BNA_M
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

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

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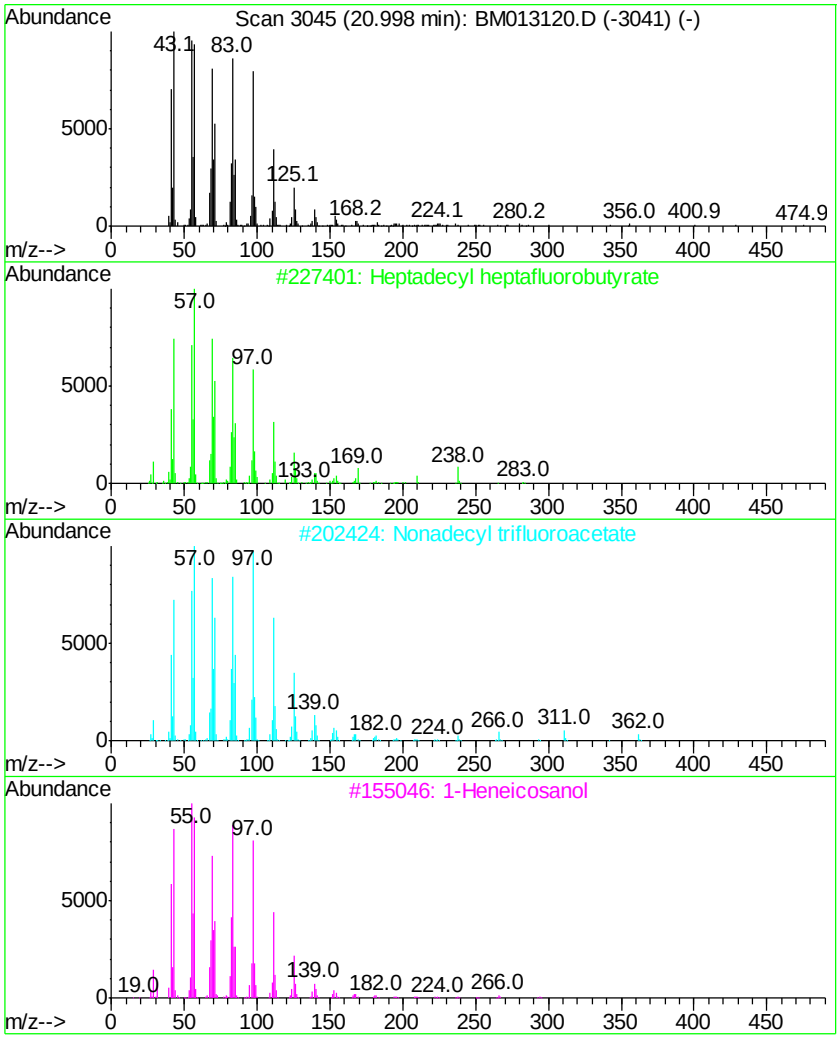
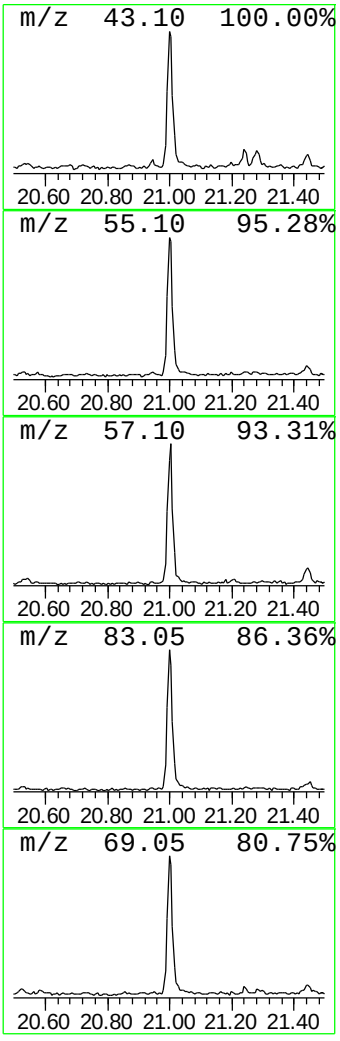
Peak Number 5 Heptadecyl heptafluorobutyrate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.00	5.72 ng/ul	1146230	Chrysene-d12	21.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
2	Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	94
3	1-Heneicosanol	312	C21H44O	015594-90-8	93
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	93
5	Eicosyl trifluoroacetate	394	C22H41F3O2	1000351-75-2	90





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Sample : I6496-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AB4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

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

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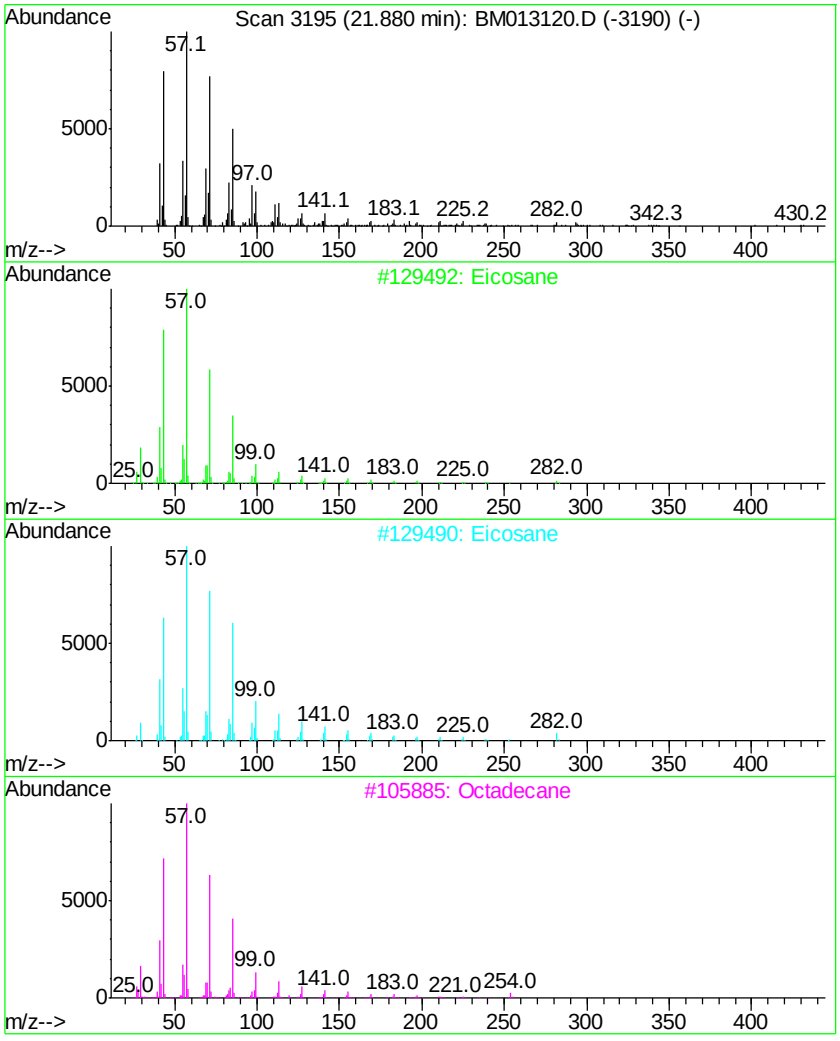
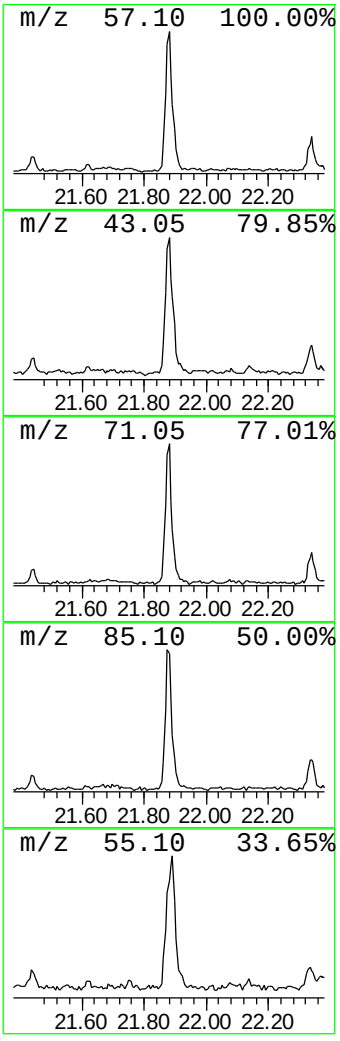
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.88	6.26 ng/ul	1255600	Chrysene-d12	21.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	96
2	Eicosane	282	C20H42	000112-95-8	95
3	Octadecane	254	C18H38	000593-45-3	95
4	Eicosane	282	C20H42	000112-95-8	94
5	Heneicosane	296	C21H44	000629-94-7	93





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112717\
Data File : BM013120.D
Acq On : 27 Nov 2017 20:49
Operator : SJ/JU
Sample : I6496-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AB4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

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

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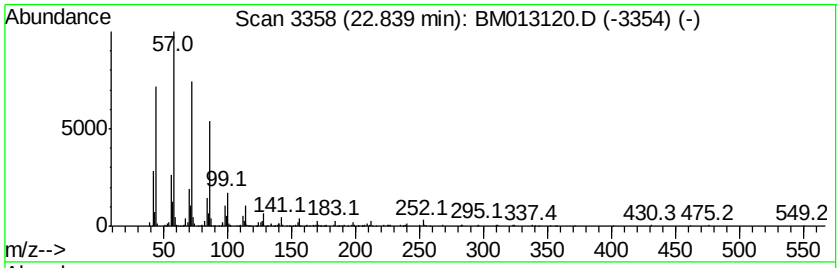
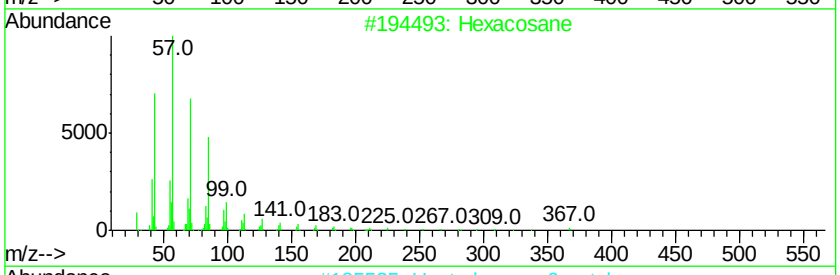
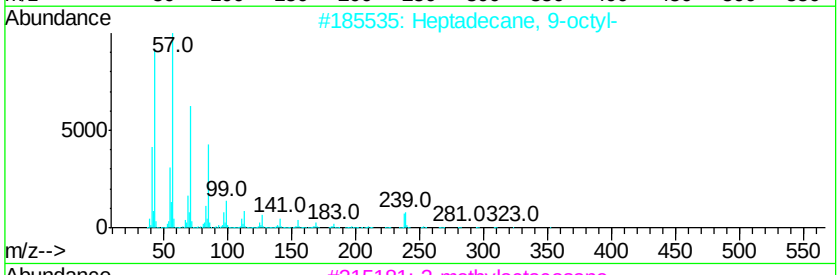
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 3

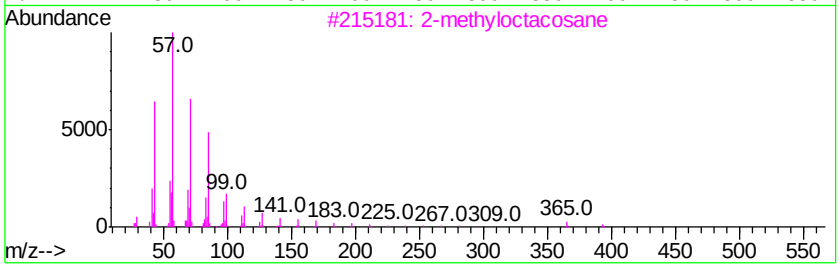
R.T.	EstConc	Area	Relative to ISTD	R.T.
22.84	8.78 ng/ul	1287690	Perylene-d12	23.50

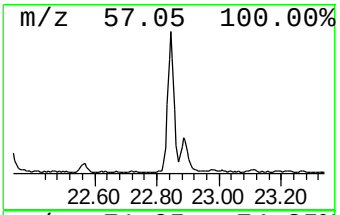
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	91
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3	2-methyloctacosane	408	C29H60	1000376-72-8	91
4	Heneicosane	296	C21H44	000629-94-7	91
5	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91

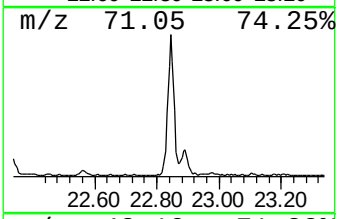


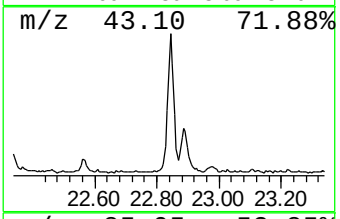


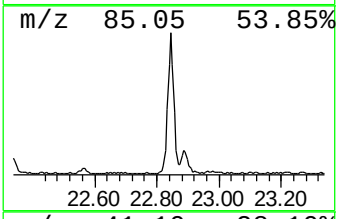


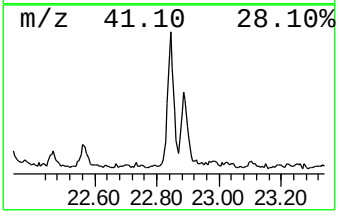












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Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AB4

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Quant Title : SVOA CALIBRATION

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

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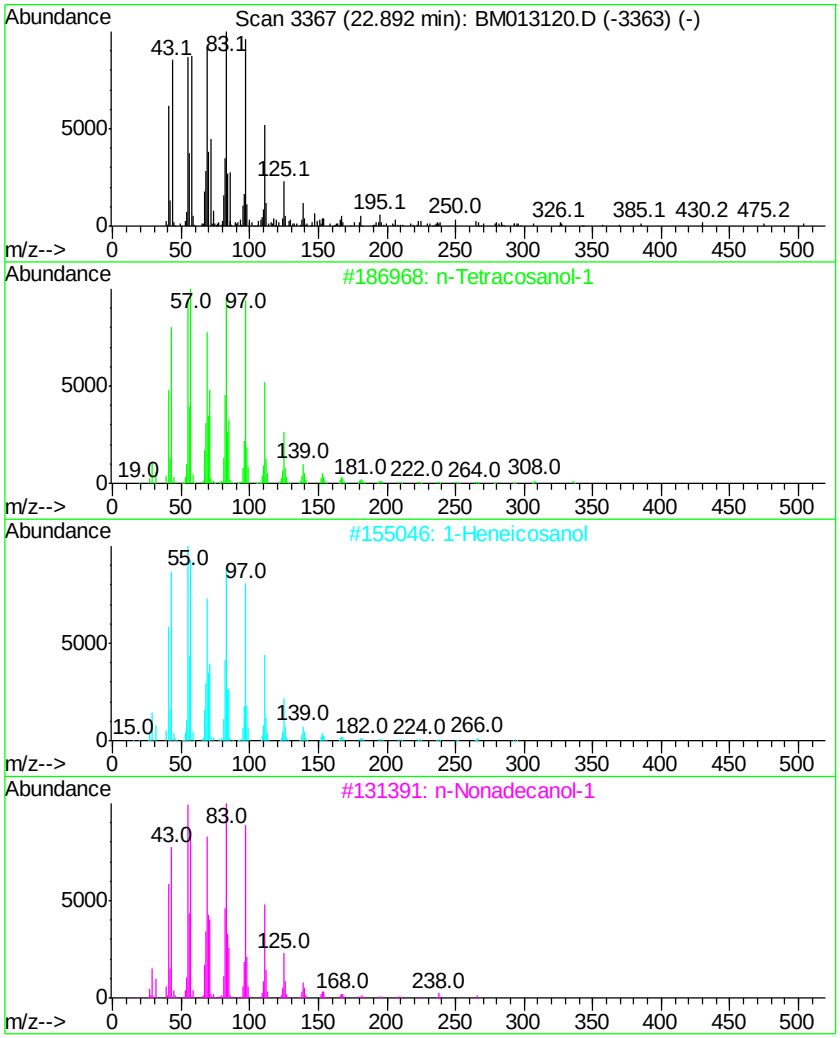
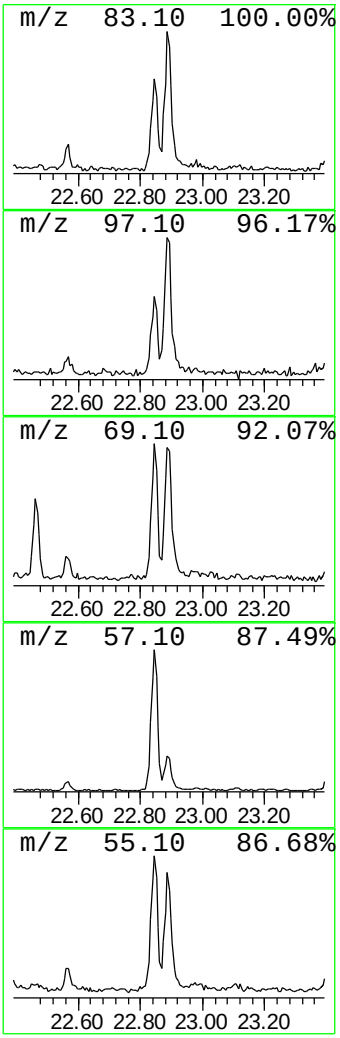
Peak Number 8 n-Tetracosanol-1 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.89	4.99 ng/ul	731858	Perylene-d12	23.50

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Tetracosanol-1	354	C24H50O	000506-51-4	95
2	1-Heneicosanol	312	C21H44O	015594-90-8	95
3	n-Nonadecanol-1	284	C19H40O	001454-84-8	94
4	Behenic alcohol	326	C22H46O	000661-19-8	91
5	Tetradecyl trifluoroacetate	310	C16H29F3O2	006222-02-2	91





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Instrument : BNA_M
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TIC Library : C:\DATABASE\NIST11.L
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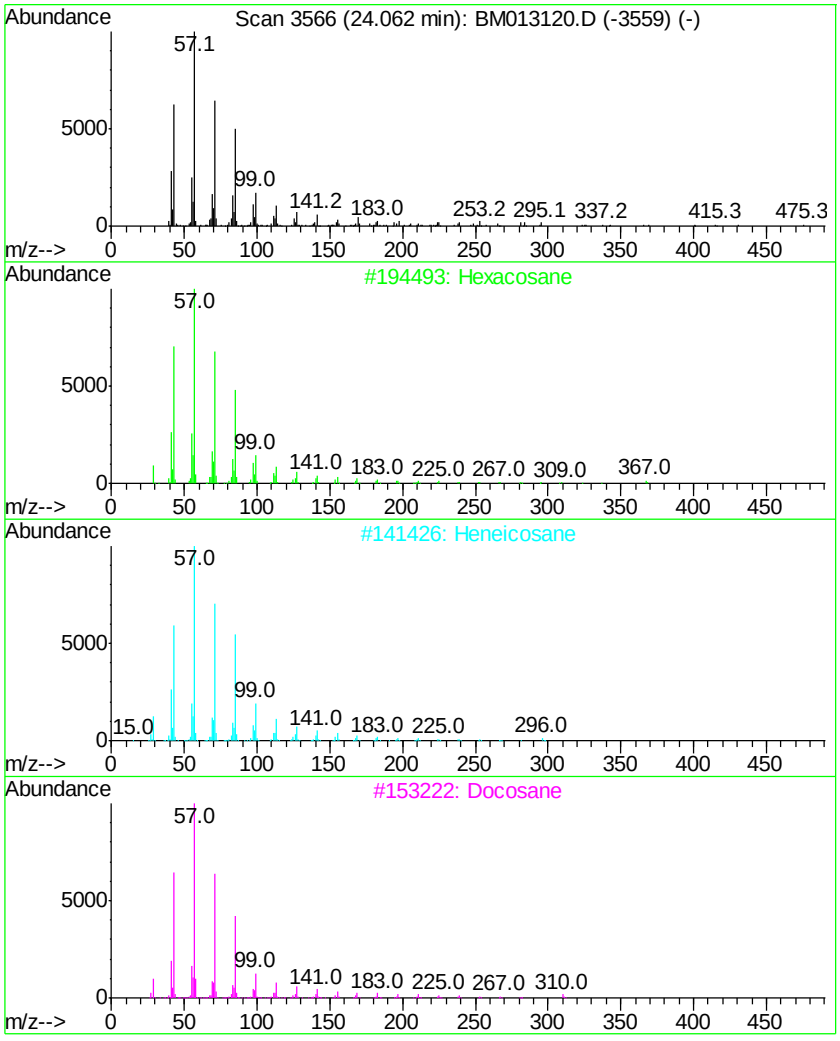
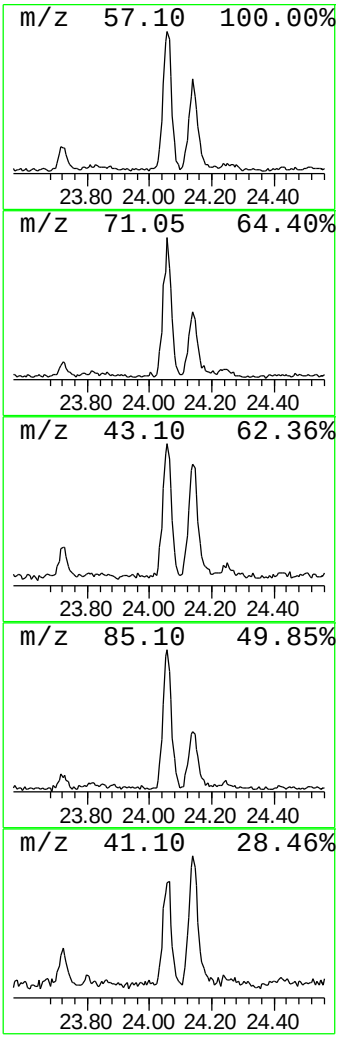
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.06	5.37 ng/ul	788102	Perylene-d12	23.50

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	91
2	Heneicosane	296	C21H44	000629-94-7	91
3	Docosane	310	C22H46	000629-97-0	91
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5	Octacosane	394	C28H58	000630-02-4	91





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Data File : BM013120.D
Acq On : 27 Nov 2017 20:49
Operator : SJ/JU
Sample : I6496-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AB4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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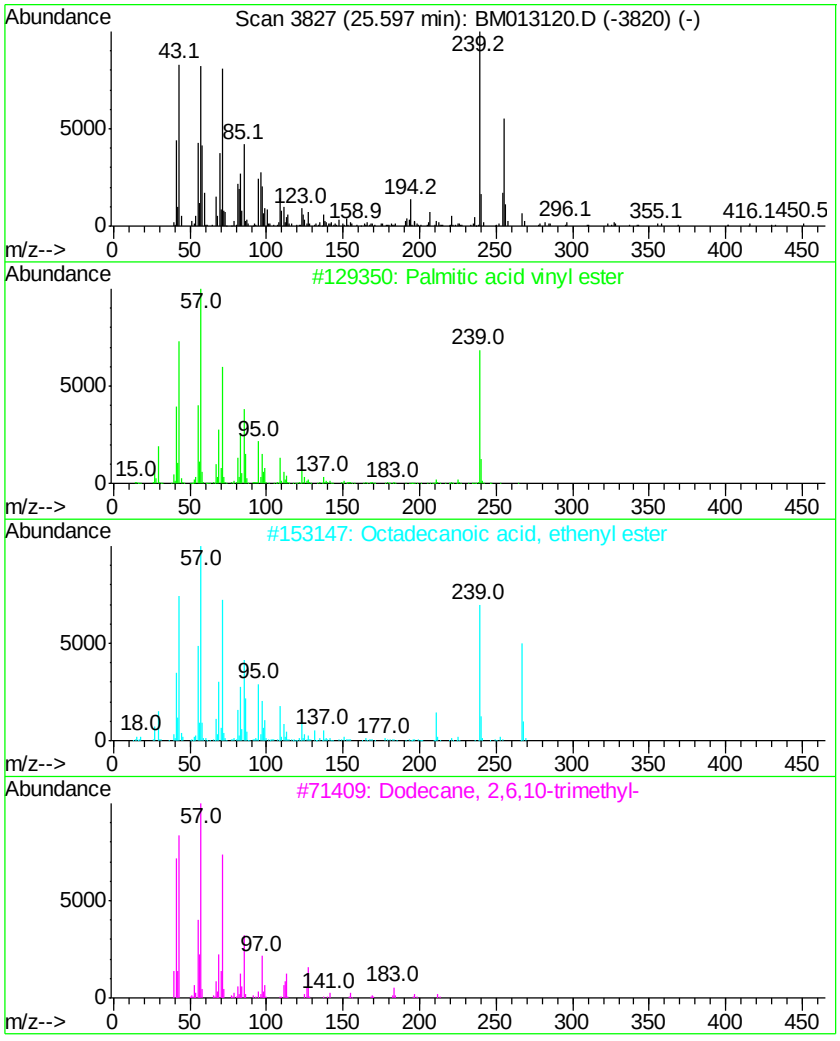
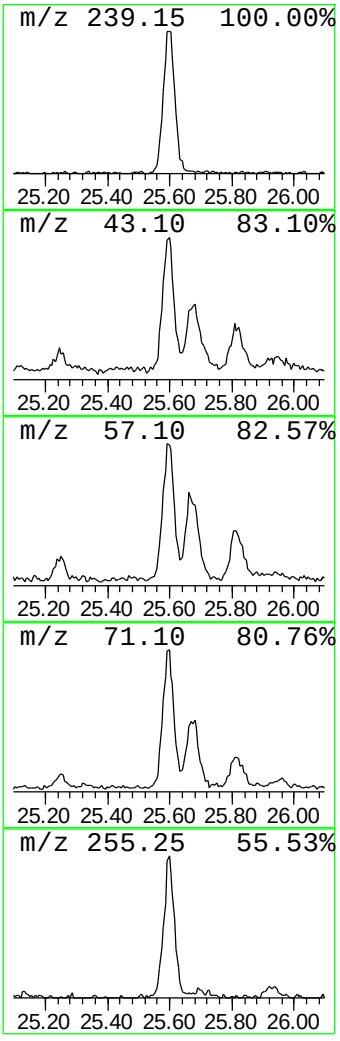
Peak Number 11 Palmitic acid vinyl ester Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.60	7.81 ng/ul	1144750	Perylene-d12	23.50

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Palmitic acid vinyl ester	282	C18H34O2	000693-38-9	68
2	Octadecanoic acid, ethenyl ester	310	C20H38O2	000111-63-7	53
3	Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	25
4	Nonadecane	268	C19H40	000629-92-5	25
5	1-Nonene, 4,6,8-trimethyl-	168	C12H24	054410-98-9	22





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Operator : SJ/JU
Sample : I6496-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AB4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

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

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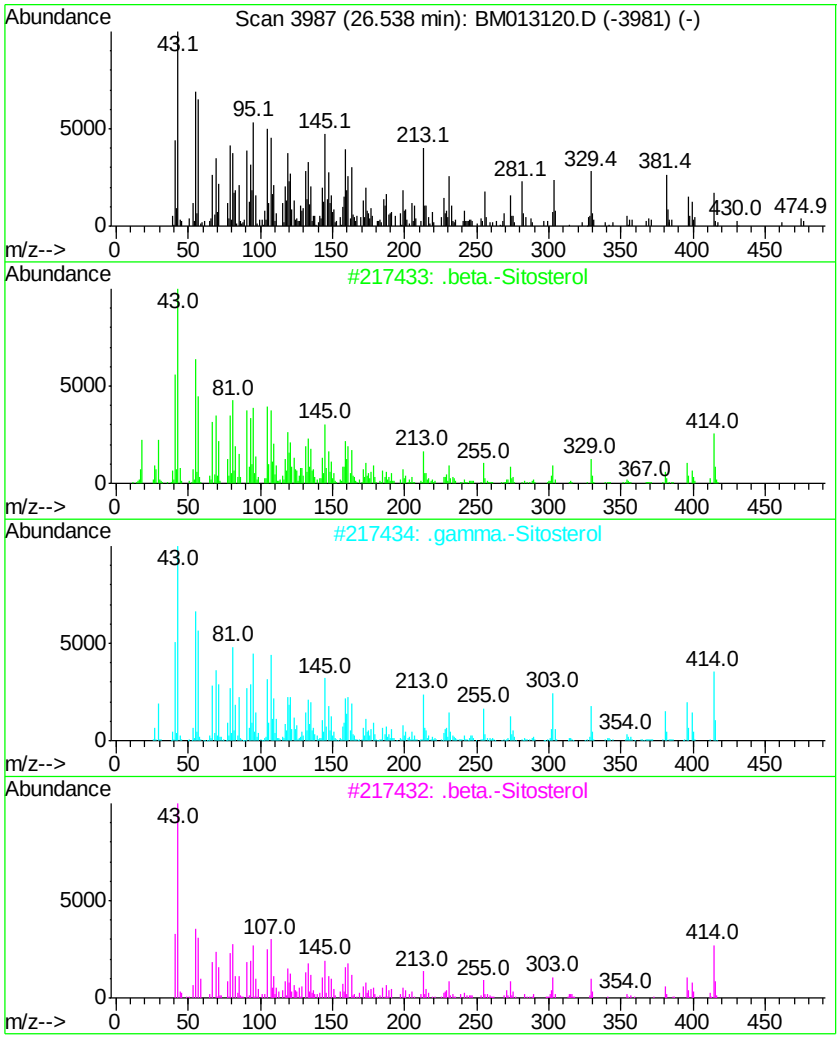
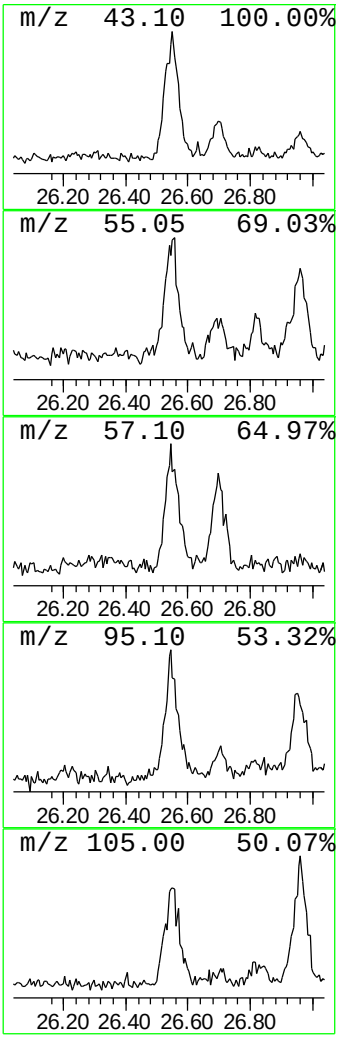
Peak Number 12 .beta.-Sitosterol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.54	9.05 ng/ul	1326820	Perylene-d12	23.50

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.beta.-Sitosterol	414	C29H50O	000083-46-5	97
2	.gamma.-Sitosterol	414	C29H50O	000083-47-6	92
3	.beta.-Sitosterol	414	C29H50O	000083-46-5	86
4	Cholest-5-ene, 3-ethoxy-, (3.bet...	414	C29H50O	000986-19-6	60
5	Ergost-7-en-3-ol, (3.beta.)-	400	C28H48O	026047-31-4	35





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Operator : SJ/JU
Sample : I6496-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AB4

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

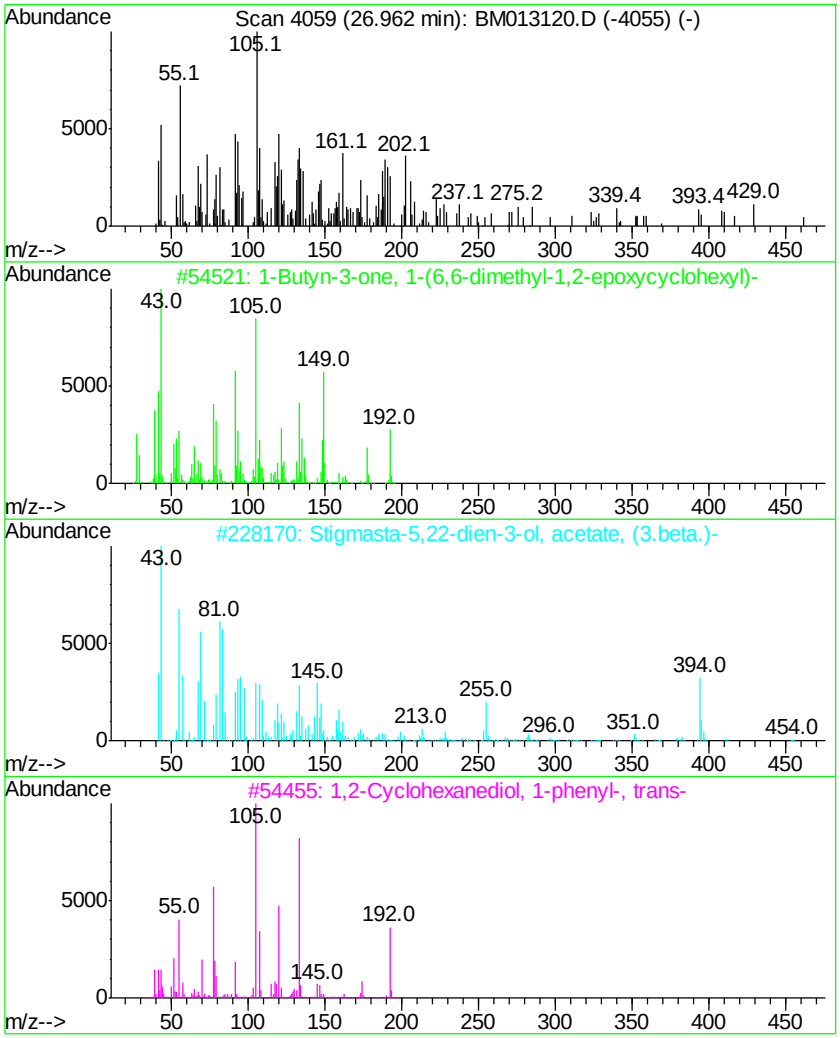
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

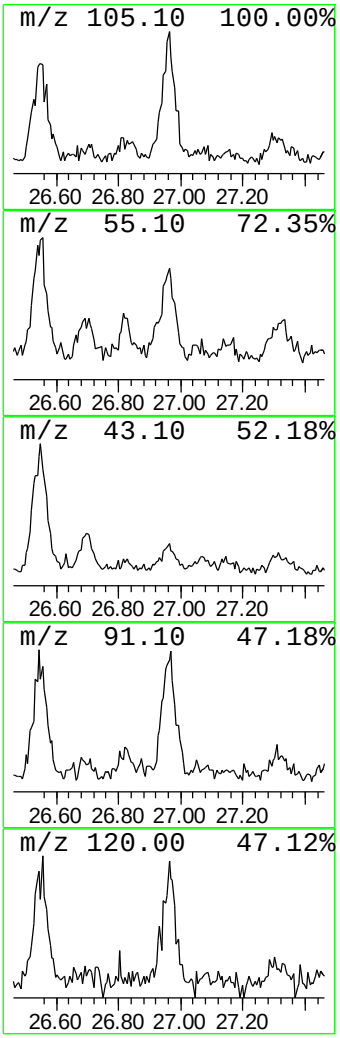
****

Peak Number 13 unknown-03 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.96	5.10 ng/ul	748295	Perylene-d12	23.50

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butyn-3-one, 1-(6,6-dimethyl-1...	192	C12H16O2	1000196-97-2	38
2	Stigmasta-5,22-dien-3-ol, acetat...	454	C31H50O2	004651-48-3	25
3	1,2-Cyclohexanediol, 1-phenyl-, ...	192	C12H16O2	027167-34-6	17
4	3-(4-Acetoxyphenyl)-2-[(2,4-dini...	416	C18H16N4O8	1000203-03-6	14
5	9,19-Cycloergost-24(28)-en-3-ol,...	468	C32H52O2	010376-42-8	12





Tentatively Identified Compound (LSC) summary

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Quant Title : SVOA CALIBRATION

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
n-Hexadecanoic acid	18.00	9.1	ng/ul	1666750	4	17.09	3662970	20.0
unknown-01	18.05	2.1	ng/ul	386972	4	17.09	3662970	20.0
Octadecanoic acid	19.29	4.9	ng/ul	991208	5	21.27	4010990	20.0
unknown-02	20.35	4.1	ng/ul	813278	5	21.27	4010990	20.0
Heptadecyl heptaf...	21.00	5.7	ng/ul	1146230	5	21.27	4010990	20.0
(DEL) Alkane: Str...	21.88	6.3	ng/ul	1255600	5	21.27	4010990	20.0
(DEL) Alkane: Str...	22.84	8.8	ng/ul	1287690	6	23.50	2932710	20.0
n-Tetracosanol-1	22.89	5.0	ng/ul	731858	6	23.50	2932710	20.0
(DEL) Alkane: Str...	24.06	5.4	ng/ul	788102	6	23.50	2932710	20.0
Palmitic acid vin...	25.60	7.8	ng/ul	1144750	6	23.50	2932710	20.0
.beta.-Sitosterol	26.54	9.1	ng/ul	1326820	6	23.50	2932710	20.0
unknown-03	26.96	5.1	ng/ul	748295	6	23.50	2932710	20.0