LSC Area Percent Report

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA M\DATA\BM121117\
Data File : BM013317.D
Acq On : 11 Dec 2017 22:59
Operator : SJ/JU
Sample : I6759-03
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9A02

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-BM113017.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	4.852	295	300	305	rVB3	42154	65064	1.49%	0.091%
2	6.863	635	642	652	rBV2	514729	962928	22.00%	1.345%
3	7.028	664	670	678	rVB	386487	582584	13.31%	0.814%
4	7.222	695	703	713	rBV	521267	836537	19.11%	1.169%
5	7.687	776	782	790	rBV	695361	1103774	25.21%	1.542%
6	8.222	867	873	882	rBV2	53226	97288	2.22%	0.136%
7	8.393	894	902	911	rBV	659606	1079207	24.65%	1.508%
8	8.840	971	978	989	rBV	529185	861209	19.67%	1.203%
9	9.557	1093	1100	1107	rBV	444761	730438	16.68%	1.021%
10	10.093	1184	1191	1199	rBV	695323	1163620	26.58%	1.626%
11	10.469	1244	1255	1266	rBV	955550	1653108	37.76%	2.310%
12	10.610	1270	1279	1287	rBV	495625	878311	20.06%	1.227%
13	11.975	1507	1511	1521	rBV3	39437	82388	1.88%	0.115%
14	13.745	1805	1812	1816	rBV	1197036	1741050	39.77%	2.432%
15	13.786	1816	1819	1828	rVB	266859	389639	8.90%	0.544%
16	14.022	1848	1859	1863	rBV	1373439	2332773	53.28%	3.259%
17	14.233	1892	1895	1902	rVB3	54748	75120	1.72%	0.105%
18	14.328	1902	1911	1919	rBV2	1442416	2276540	52.00%	3.181%
19	14.539	1940	1947	1958	rBV	519250	880221	20.11%	1.230%
20	14.763	1982	1985	1994	rVB10	24718	46453	1.06%	0.065%
21	14.875	1997	2004	2009	rBV9	33771	75333	1.72%	0.105%
22	15.192	2054	2058	2065	rVB3	72711	104244	2.38%	0.146%
23	15.328	2074	2081	2087	rBV	1696494	2485947	56.78%	3.473%
24	15.386	2087	2091	2096	rBV8	27102	51594	1.18%	0.072%
25	15.451	2096	2102	2108	rVV	530343	770263	17.59%	1.076%
26	16.057	2200	2205	2213	rBV3	253041	466960	10.67%	0.652%
27	16.657	2305	2307	2311	rVB3	68746	75581	1.73%	0.106%
28	16.727	2314	2319	2326	rVV	962463	1339509	30.60%	1.871%
29	16.845	2336	2339	2343	rBV	109919	132798	3.03%	0.186%
30	16.898	2344	2348	2354	rVB2	158927	241905	5.53%	0.338%
31	17.080	2373	2379	2385	rBV	1992831	2902293	66.29%	4.055%
32	17.174	2389	2395	2399	rVV2	2051405	3058941	69.87%	4.274%
33	17.210	2399	2401	2407	rVV	860361	986976	22.54%	1.379%
34	17.269	2409	2411	2416	rVB6	61437	82074	1.87%	0.115%

LSC Area Percent Report

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Data File : BM013317.D
Acq On : 11 Dec 2017 22:59
Operator : SJ/JU
Sample : I6759-03
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9A02

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-BM113017.M
Title : SVOA CALIBRATION

35	17.486	2445	2448	2452	rVB2	84413	122453	2.80%	0.171%
36	17.586	2461	2465	2468	rBV3	239767	278466	6.36%	0.389%
37	17.751	2489	2493	2498	rVB2	137432	217015	4.96%	0.303%
38	17.980	2528	2532	2537	rBV2	419957	626817	14.32%	0.876%
39	18.080	2546	2549	2558	rBV3	158292	273708	6.25%	0.382%
40	18.280	2579	2583	2589	rBV	422030	544107	12.43%	0.760%
41	18.345	2591	2594	2598	rBV2	119171	164110	3.75%	0.229%
42	18.498	2617	2620	2624	rBV3	132018	186027	4.25%	0.260%
43	18.916	2688	2691	2695	rVB2	448523	489022	11.17%	0.683%
44	19.233	2742	2745	2748	rBV	269069	340319	7.77%	0.475%
45	19.263	2748	2750	2759	rVB3	204674	432120	9.87%	0.604%
46	19.363	2763	2767	2772	rVB	334589	446401	10.20%	0.624%
47	19.474	2781	2786	2793	rBV	2490485	3842196	87.76%	5.368%
48	19.533	2793	2796	2802	rVB3	175009	257448	5.88%	0.360%
49	20.027	2878	2880	2885	rVB	169072	182734	4.17%	0.255%
50	20.698	2991	2994	2999	rBV	246301	331982	7.58%	0.464%
51	20.986	3039	3043	3048	rBV2	1046103	1394597	31.86%	1.948%
52	21.262	3086	3090	3095	rBV	2857784	3684384	84.16%	5.148%
53	21.439	3116	3120	3126	rVB	572639	787147	17.98%	1.100%
54	21.568	3139	3142	3147	rBV2	198787	301832	6.89%	0.422%
55	22.327	3267	3271	3278	rVB4	319546	595711	13.61%	0.832%
56	22.680	3327	3331	3337	rBV2	276811	531115	12.13%	0.742%
57	22.739	3338	3341	3349	rVB	234004	444201	10.15%	0.621%
58	22.821	3350	3355	3367	rVB2	2123904	4377929	100.00%	6.116%
59	22.998	3380	3385	3390	rBV	549805	925497	21.14%	1.293%
60	23.303	3432	3437	3440	rBV	435440	788459	18.01%	1.102%
61	23.351	3440	3445	3450	rVV	1765259	3304793	75.49%	4.617%
62	23.392	3450	3452	3458	rVB	689157	1006653	22.99%	1.406%
63	23.492	3464	3469	3474	rBV	1463784	2508129	57.29%	3.504%
64	24.027	3557	3560	3569	rVB7	229901	431078	9.85%	0.602%
65	25.045	3726	3733	3741	rBV6	453283	1196243	27.32%	1.671%
66	25.492	3803	3809	3817	rVB3	536526	1303084	29.76%	1.821%
67	25.727	3841	3849	3861	rVB2	1481542	4277936	97.72%	5.977%
68	26.033	3896	3901	3917	rVB5	455386	1453961	33.21%	2.031%
69	26.409	3957	3965	3978	rVB2	1003661	2915513	66.60%	4.073%

Sum of corrected areas: 71575857

LSC Report - Integrated Chromatogram

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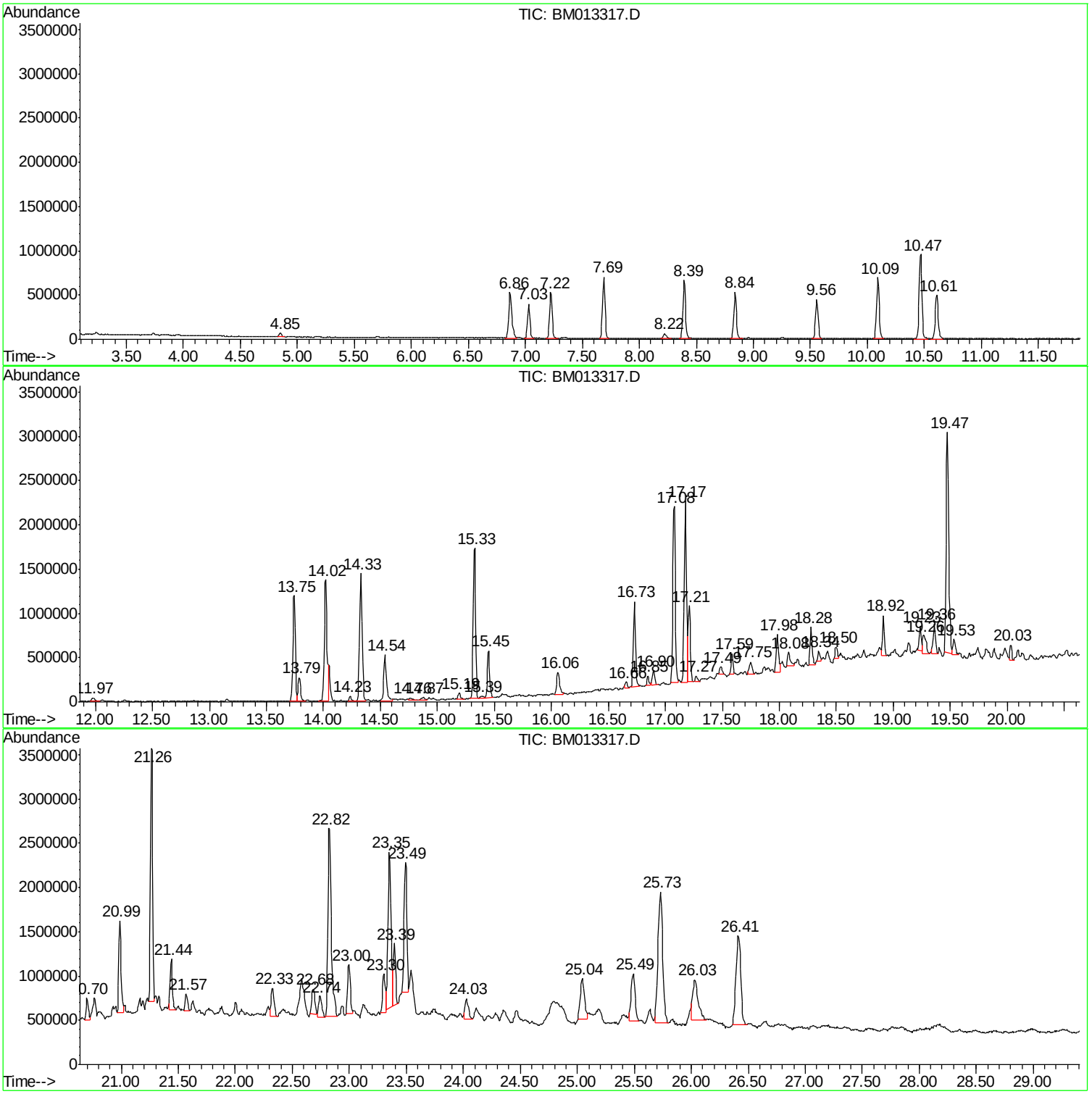
Data File : BM013317.D
Acq On : 11 Dec 2017 22:59
Operator : SJ/JU
Sample : I6759-03
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9A02

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

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Quant Title : SVOA CALIBRATION

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

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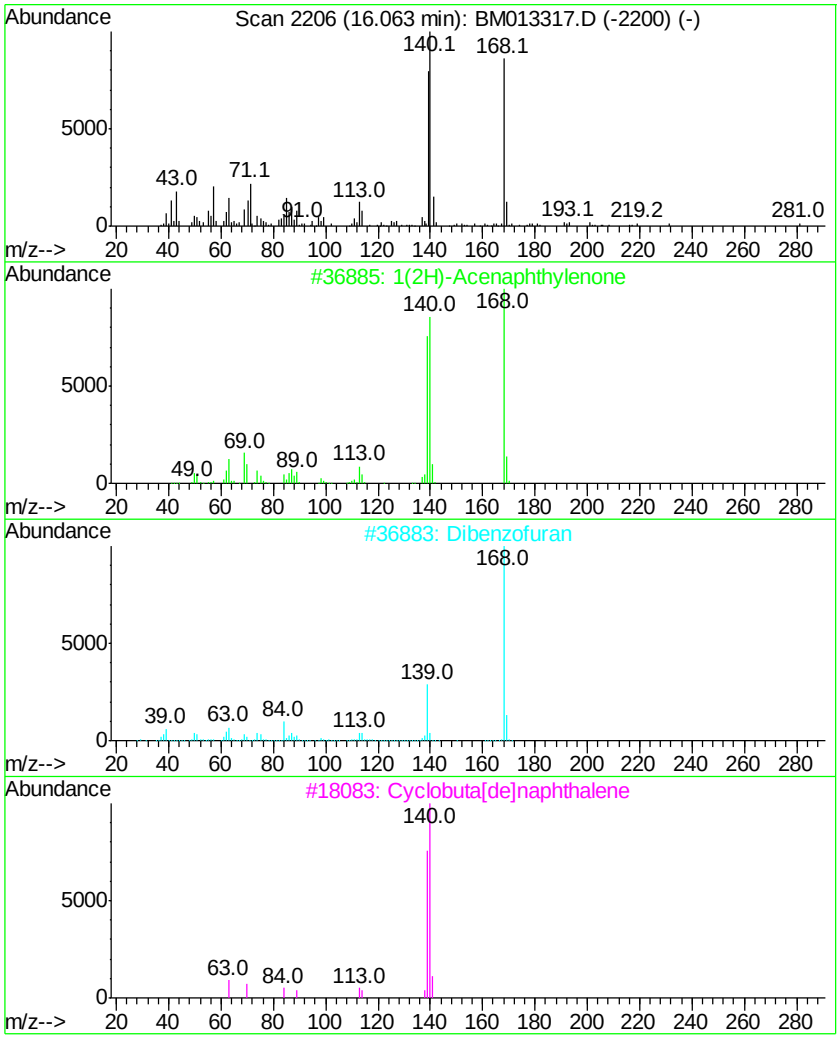
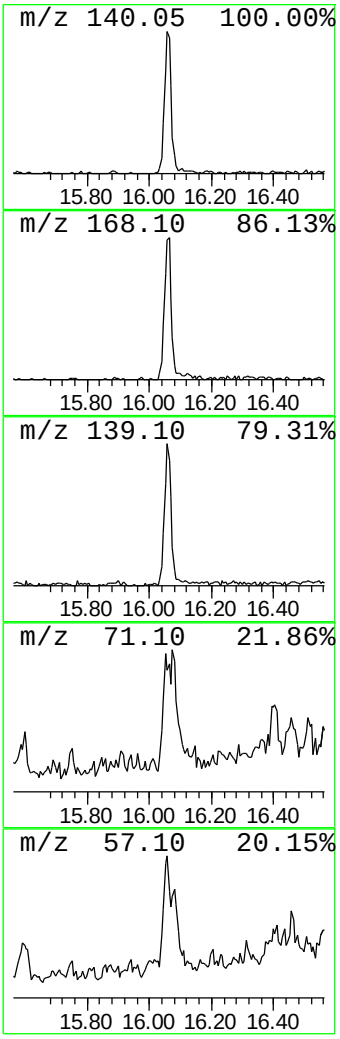
Peak Number 1 1(2H)-Acenaphthylenone Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.06	3.22 ng/ul	466960	Phenanthrene-d10	17.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(2H)-Acenaphthylenone	168	C12H8O	002235-15-6	94
2	Dibenzofuran	168	C12H8O	000132-64-9	52
3	Cyclobuta[de]naphthalene	140	C11H8	024973-91-9	49
4	3-Fluorosalicylaldehyde	140	C7H5FO2	000394-50-3	46
5	Benzaldehyde, 4-fluoro-3-hydroxy-	140	C7H5FO2	103438-85-3	46





Library Search Compound Report

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA M\DATA\BM121117\
Data File : BM013317.D
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Operator : SJ/JU
Sample : I6759-03
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9A02

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

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

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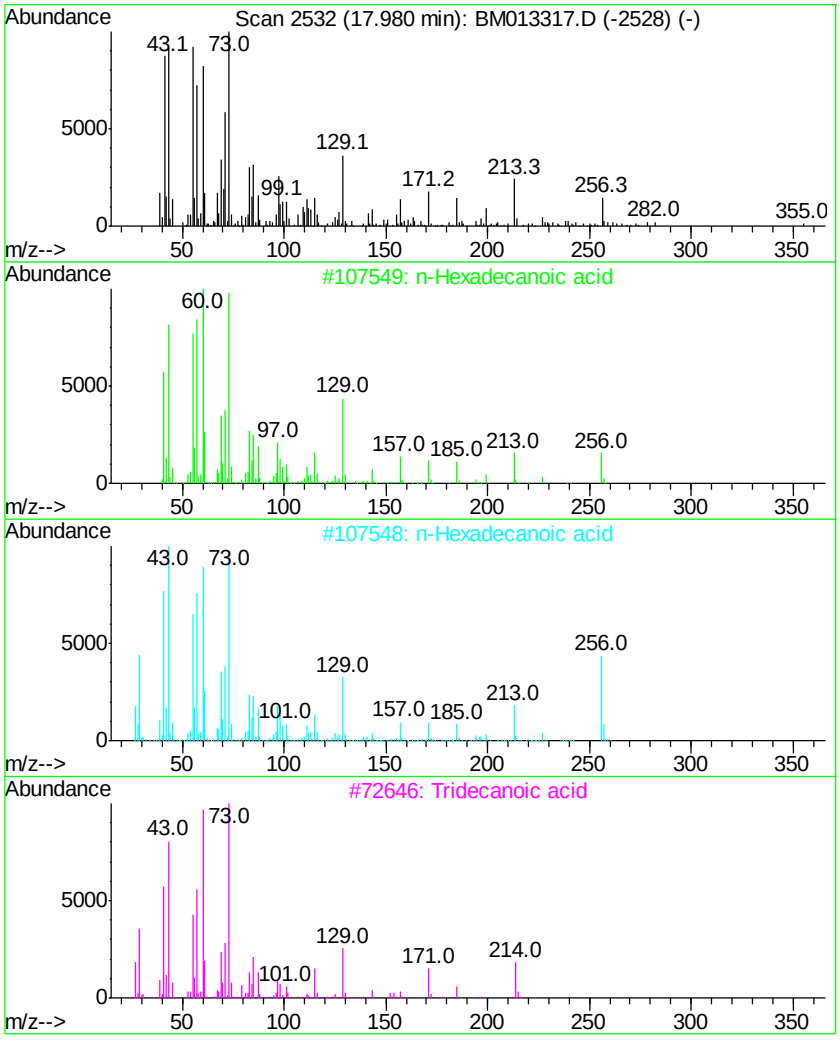
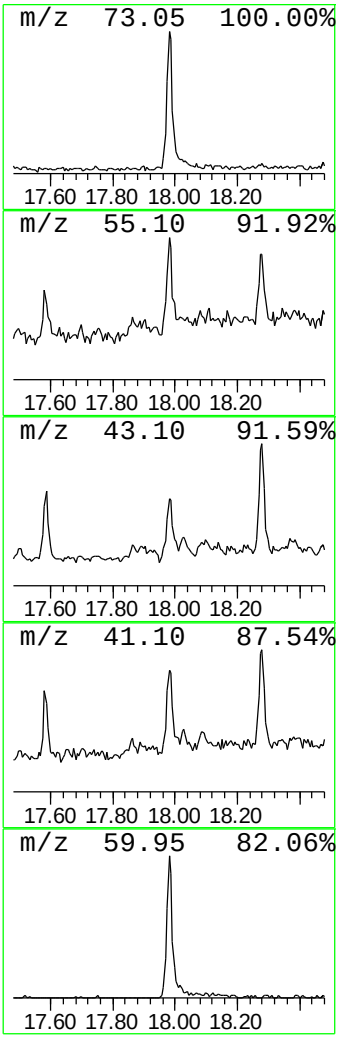
Peak Number 2 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.98	4.32 ng/ul	626817	Phenanthrene-d10	17.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
3	Tridecanoic acid	214	C13H26O2	000638-53-9	70
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	68
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	64





Library Search Compound Report

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Sample : I6759-03
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9A02

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

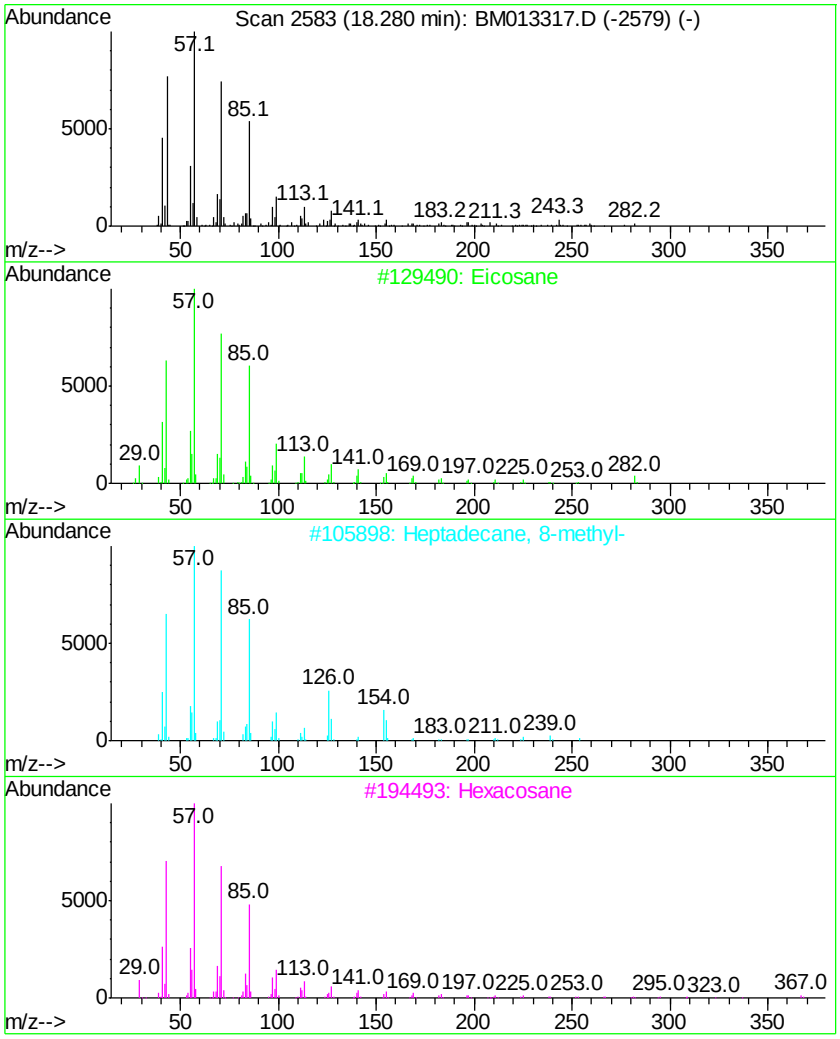
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

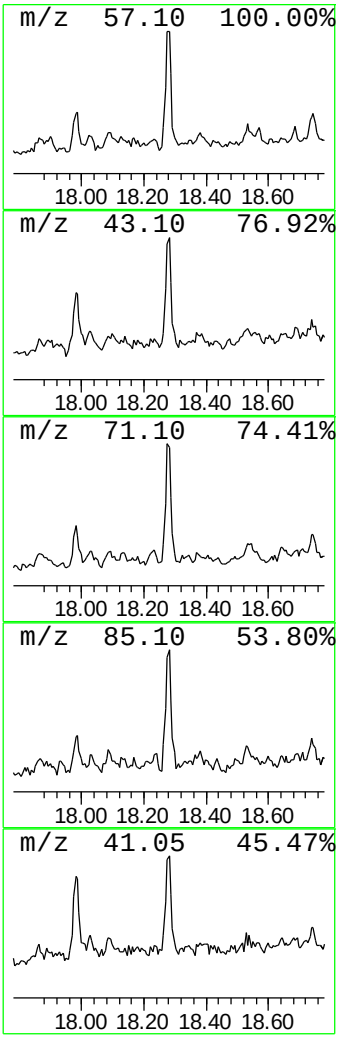
****

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.28	3.75 ng/ul	544107	Phenanthrene-d10	17.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	96
2	Heptadecane, 8-methyl-	254	C18H38	013287-23-5	95
3	Hexacosane	366	C26H54	000630-01-3	91
4	Tetracosane	338	C24H50	000646-31-1	91
5	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91





Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA M\DATA\BM121117\
Data File : BM013317.D
Acq On : 11 Dec 2017 22:59
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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9A02

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

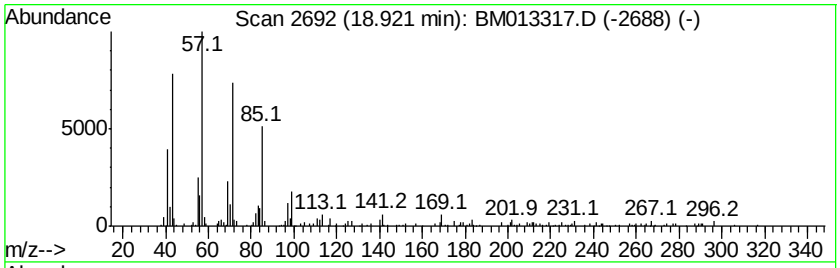
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

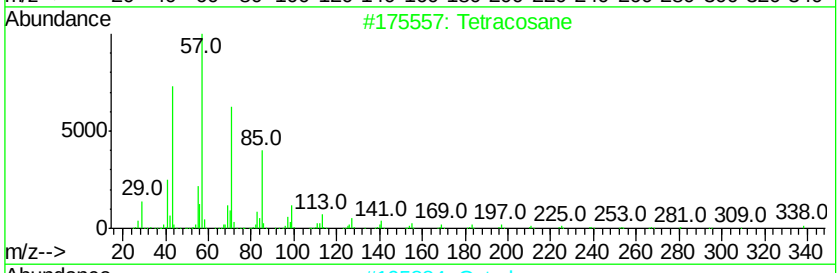
****

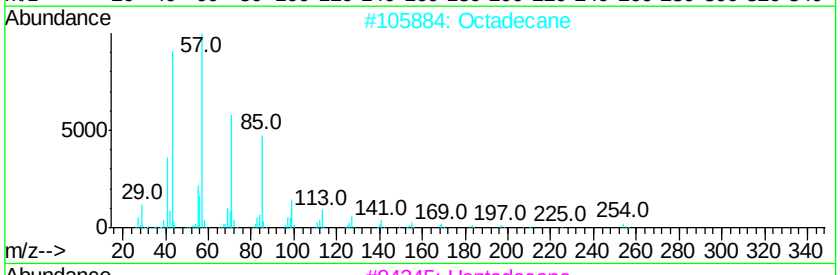
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 12

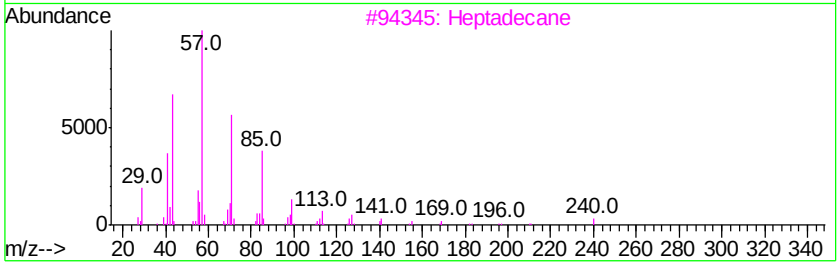
R.T.	EstConc	Area	Relative to ISTD	R.T.
18.92	3.37 ng/ul	489022	Phenanthrene-d10	17.08

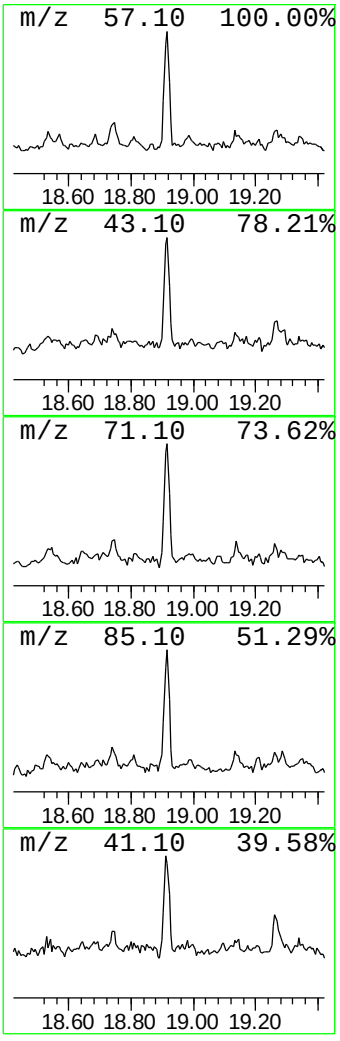
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	91
2	Octadecane	254	C18H38	000593-45-3	91
3	Heptadecane	240	C17H36	000629-78-7	90
4	Eicosane	282	C20H42	000112-95-8	90
5	Dodecane, 2-methyl-	184	C13H28	001560-97-0	90











Library Search Compound Report

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Sample : I6759-03
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9A02

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

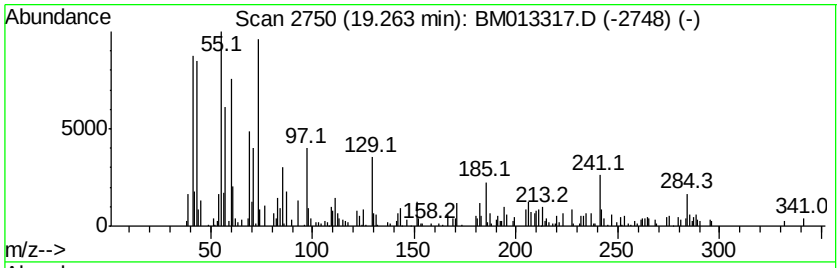
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

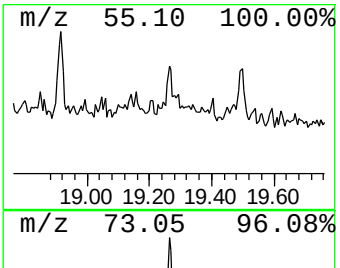
****

Peak Number 5 Octadecanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.26	2.35 ng/ul	432120	Chrysene-d12	21.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	92
2	Octadecanoic acid	284	C18H36O2	000057-11-4	90
3	Octadecanoic acid	284	C18H36O2	000057-11-4	86
4	Octadecanoic acid	284	C18H36O2	000057-11-4	70
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	55





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Data File : BM013317.D
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Operator : SJ/JU
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Misc :
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Instrument : BNA_M
ClientSampleId : F9A02

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

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

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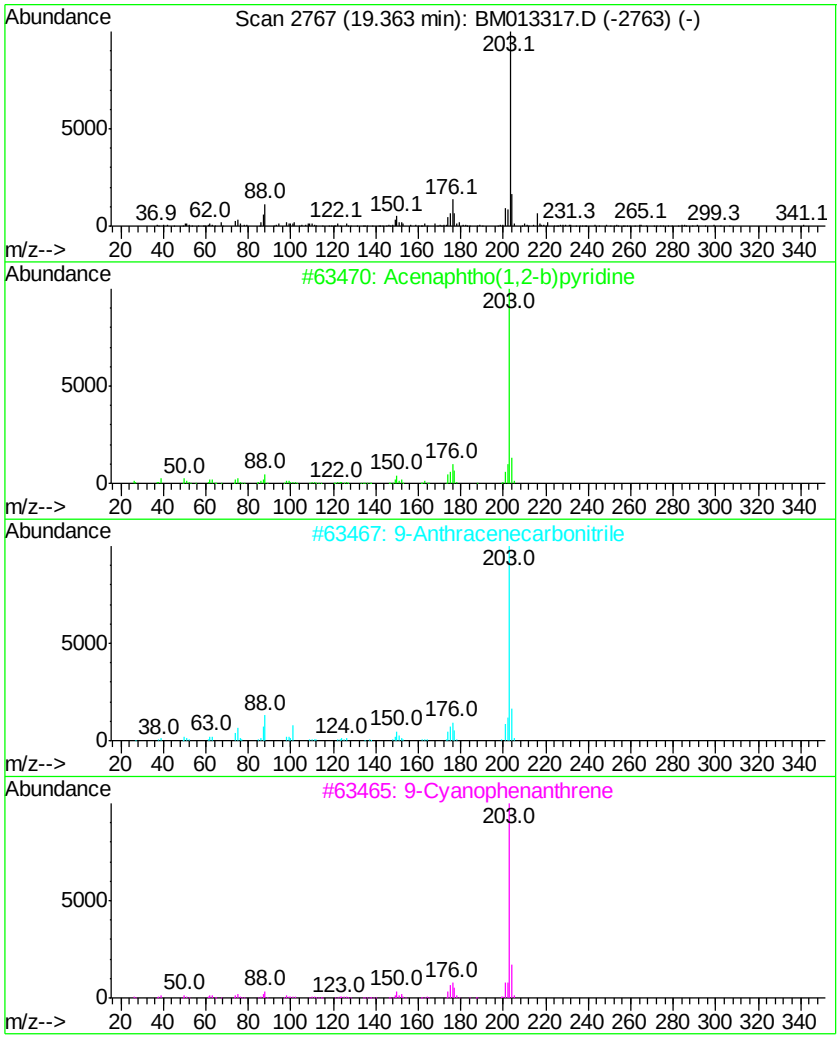
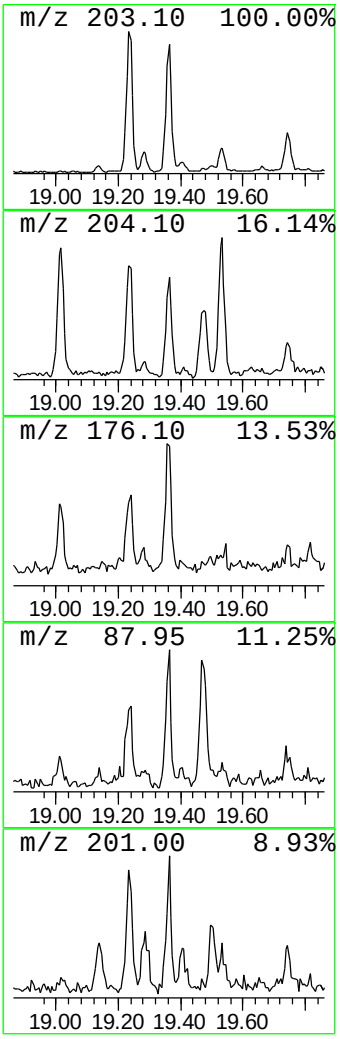
Peak Number 6 Acenaphtho(1,2-b)pyridine Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.36	2.42 ng/ul	446401	Chrysene-d12	21.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acenaphtho(1,2-b)pyridine	203	C15H9N	000206-49-5	95
2	9-Anthracenecarbonitrile	203	C15H9N	001210-12-4	95
3	9-Cyanophenanthrene	203	C15H9N	002510-55-6	94
4	Indeno(1,2,3-ij)isoquinoline	203	C15H9N	000206-56-4	94
5	9-(Cyanomethylene)fluorene	203	C15H9N	004425-74-5	94





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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9A02

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

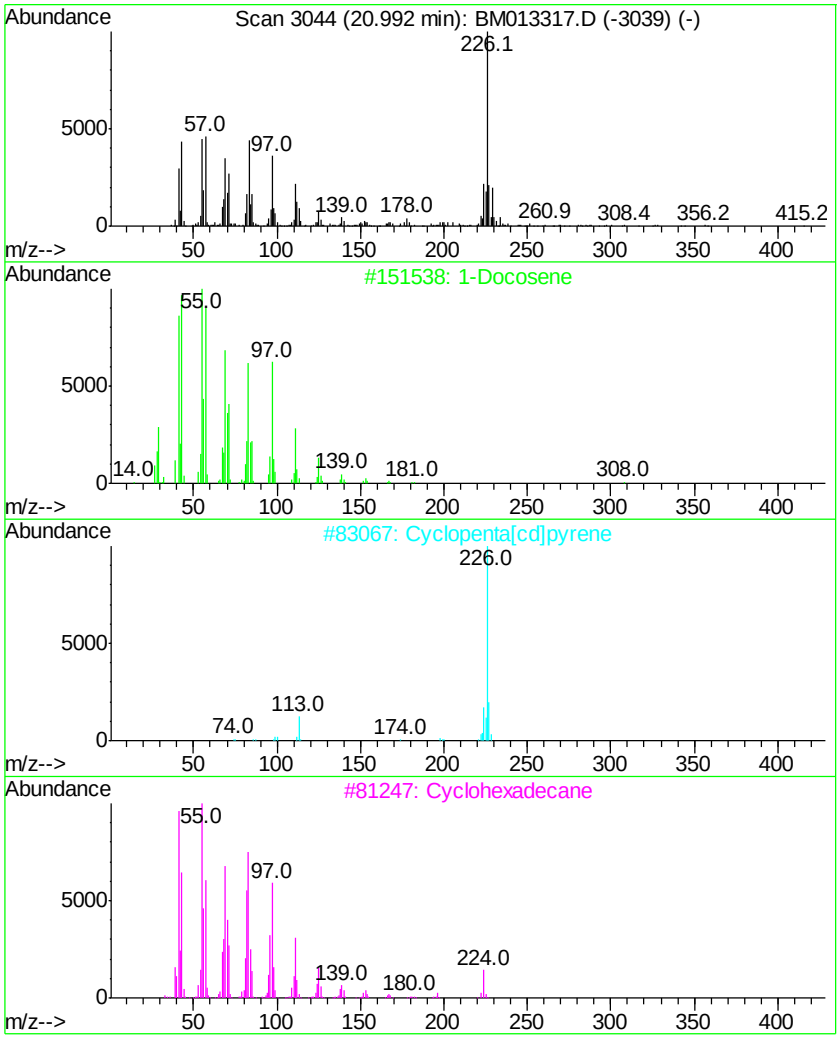
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

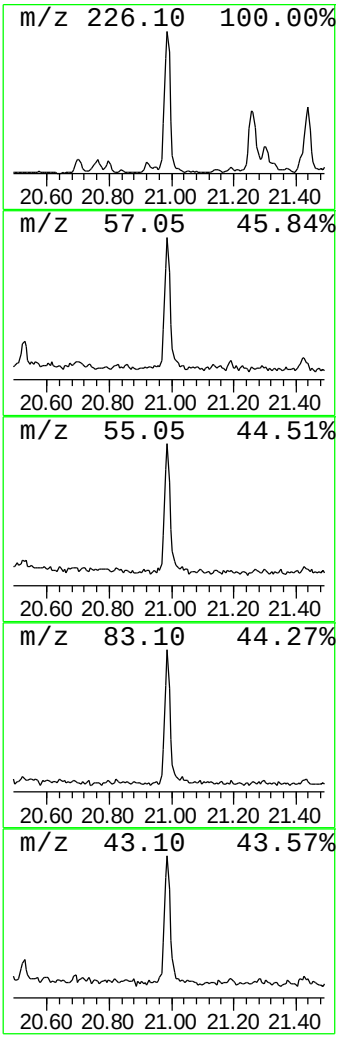
****

Peak Number 7 (DEL) Alkane: Cyclic20.99 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.99	7.57 ng/ul	1394600	Chrysene-d12	21.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	91
2	Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	50
3	Cyclohexadecane	224	C16H32	000295-65-8	43
4	n-Heptadecanol-1	256	C17H36O	001454-85-9	42
5	Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	38





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Acq On : 11 Dec 2017 22:59
Operator : SJ/JU
Sample : I6759-03
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9A02

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Quant Title : SVOA CALIBRATION

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

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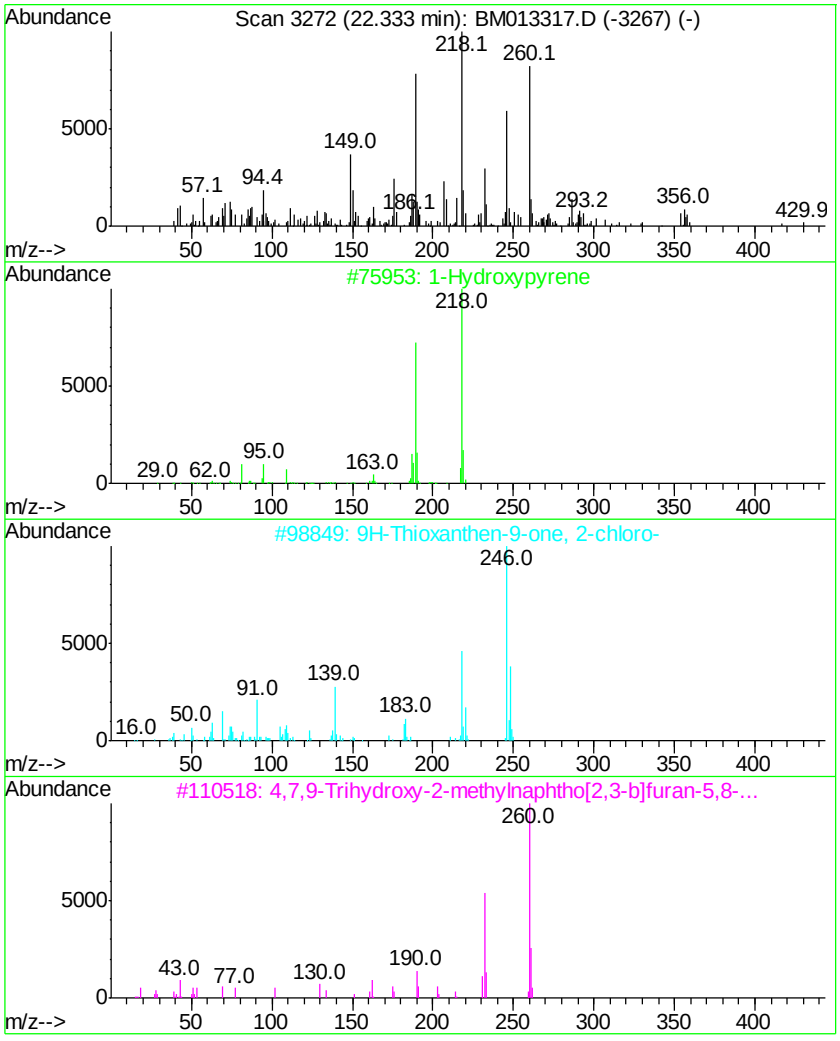
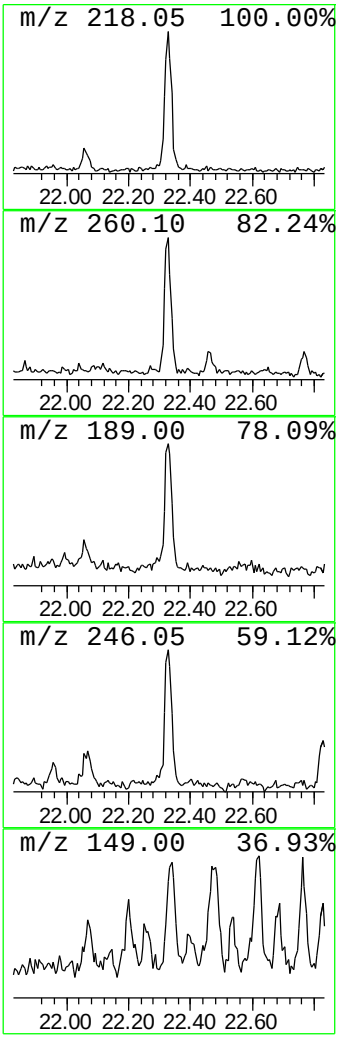
Peak Number 9 unknown-01 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.33	3.23 ng/ul	595711	Chrysene-d12	21.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hydroxypyrene	218	C16H10O	005315-79-7	38
2	9H-Thioxanthen-9-one, 2-chloro-	246	C13H7ClOS	000086-39-5	38
3	4,7,9-Trihydroxy-2-methylnaphtho...	260	C13H8O6	055293-73-7	38
4	5-Ethyl-5-p-tolylbarbituric acid	246	C13H14N2O3	052584-39-1	35
5	5-Ethyl-5-p-tolylbarbituric acid	246	C13H14N2O3	052584-39-1	35





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Instrument : BNA_M
ClientSampleId : F9A02

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Quant Title : SVOA CALIBRATION

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

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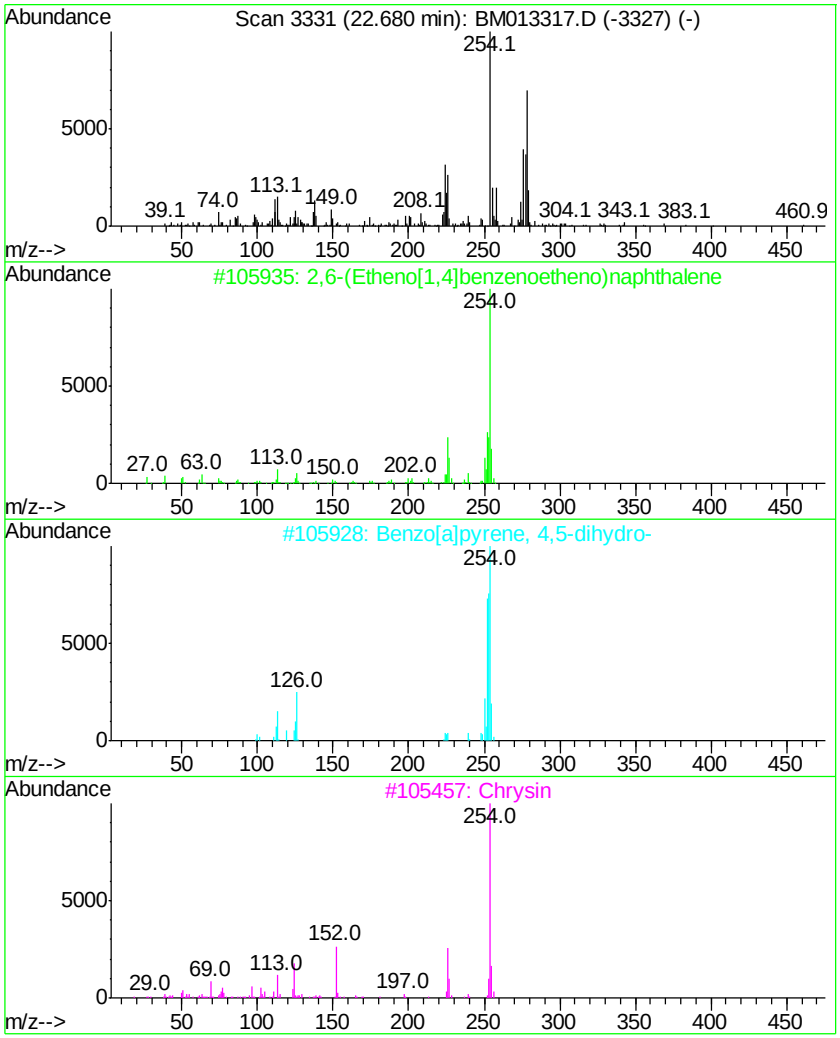
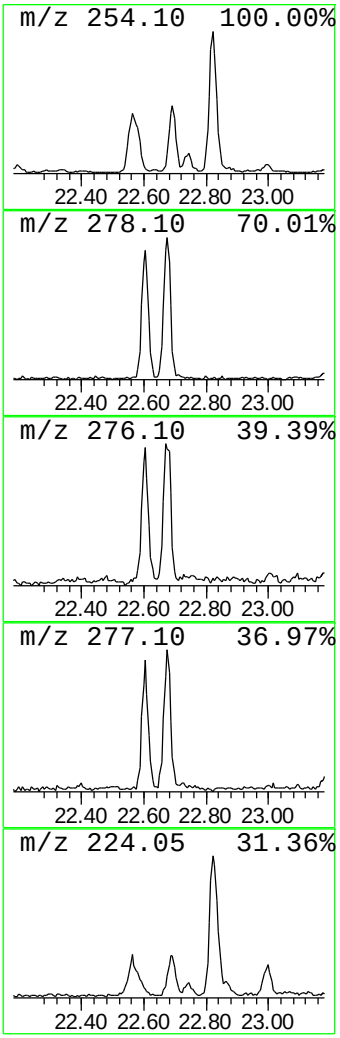
Peak Number 10 unknown-02 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.68	4.24 ng/ul	531115	Perylene-d12	23.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6-(Etheno[1,4]benzenoetheno)na...	254	C20H14	117054-70-3	38
2	Benzo[a]pyrene, 4,5-dihydro-	254	C20H14	057652-66-1	38
3	Chrysin	254	C15H10O4	000480-40-0	30
4	1,1'-Binaphthalene	254	C20H14	000604-53-5	30
5	Anthracene, 9-phenyl-	254	C20H14	000602-55-1	27





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Instrument : BNA_M
ClientSampleId : F9A02

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Quant Title : SVOA CALIBRATION

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

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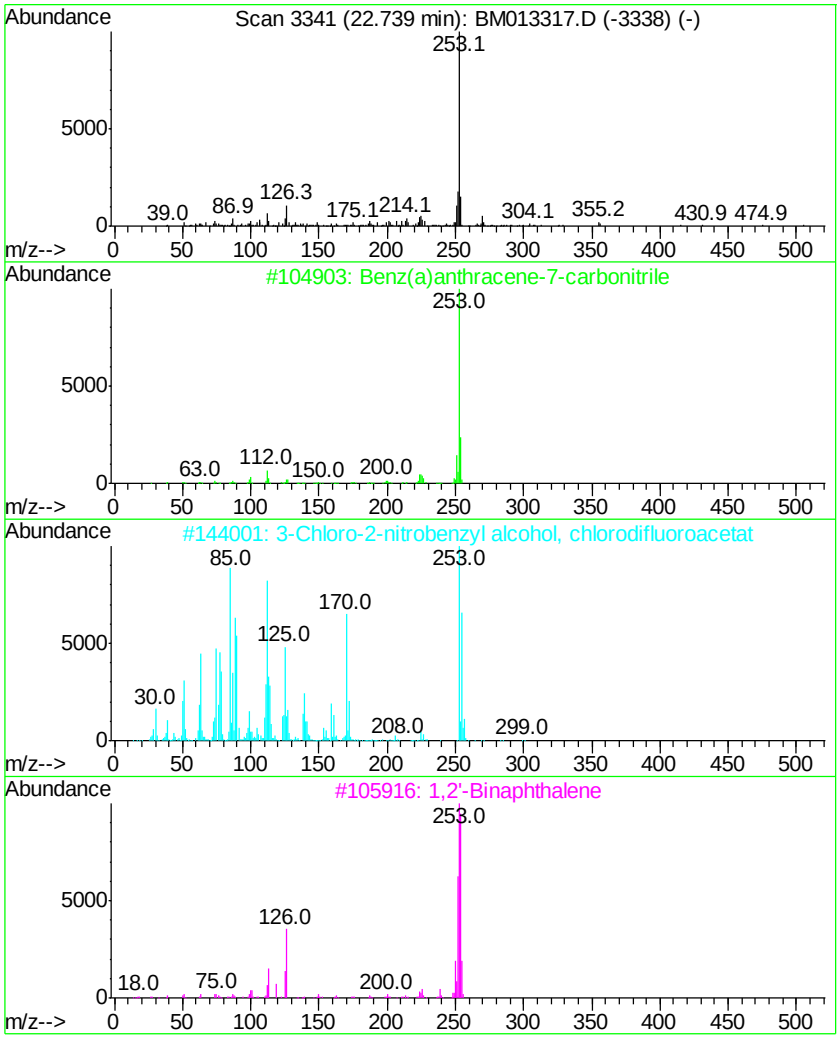
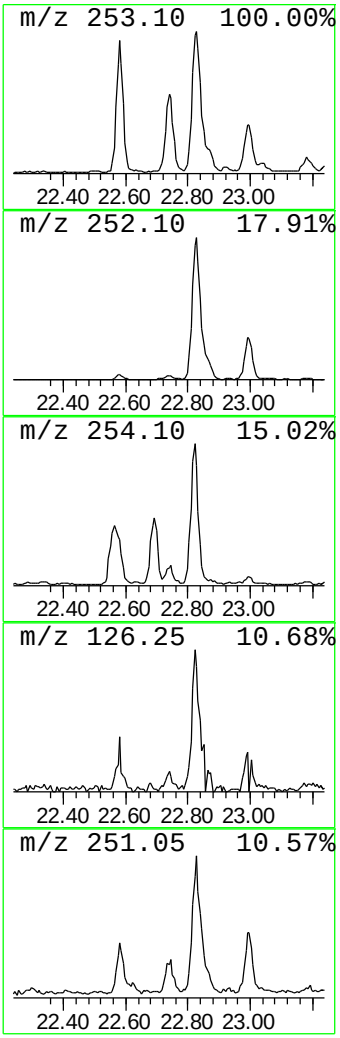
Peak Number 11 Benz(a)anthracene-7-carboni... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.74	3.54 ng/ul	444201	Perylene-d12	23.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benz(a)anthracene-7-carbonitrile	253	C19H11N	007476-08-6	64
2	3-Chloro-2-nitrobenzyl alcohol, ...	299	C9H5Cl2F2NO4	1000376-11-2	50
3	1,2'-Binaphthalene	254	C20H14	004325-74-0	42
4	Phthalic acid, 3,5-dimethylpheny...	376	C23H20O5	1000315-68-2	40
5	Phthalic acid, di(3-ethylphenyl)...	374	C24H22O4	1000357-08-4	40





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Instrument : BNA_M
ClientSampleId : F9A02

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Quant Title : SVOA CALIBRATION

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

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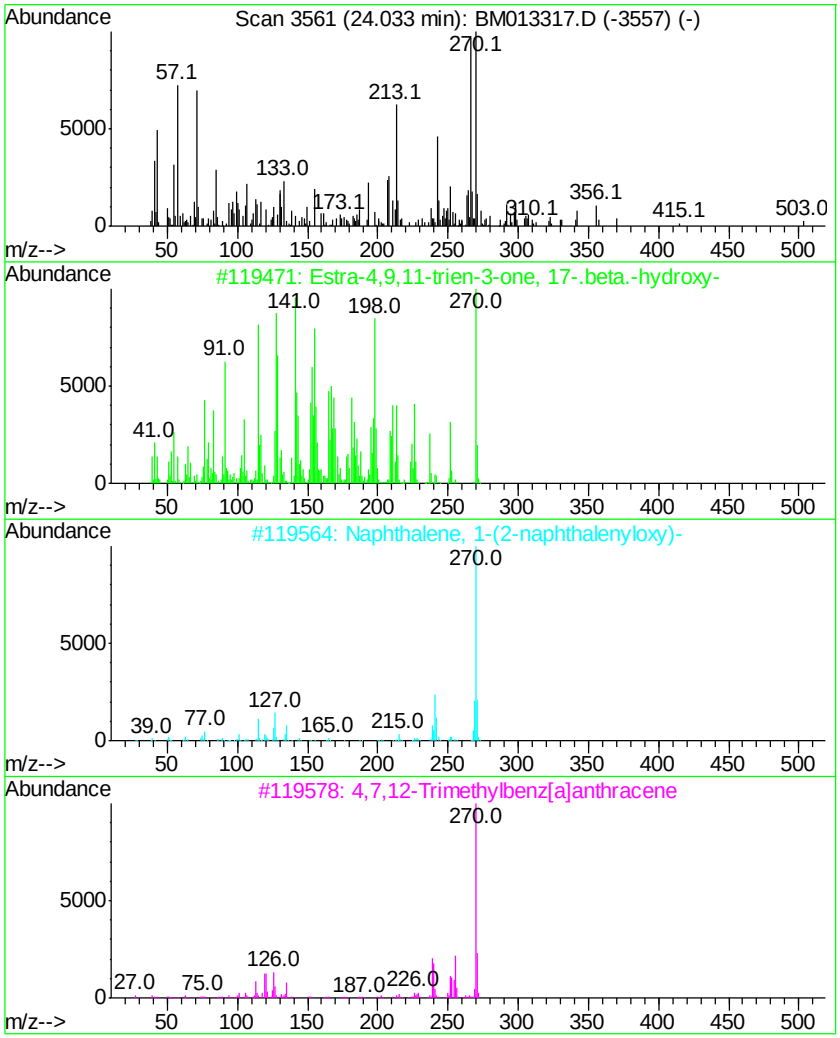
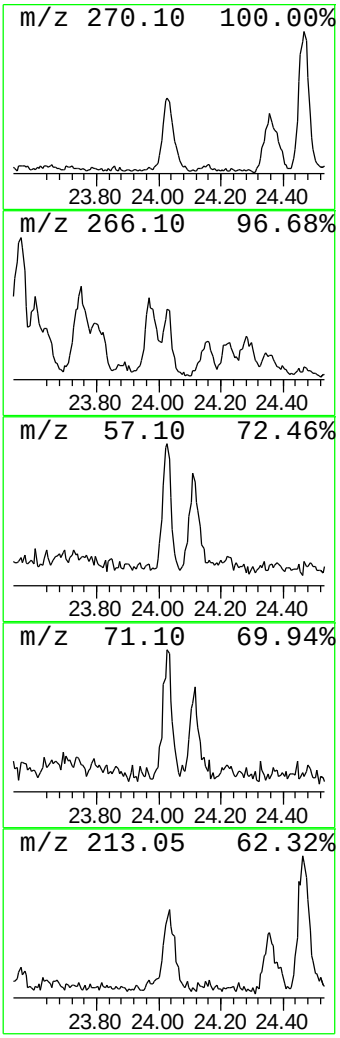
Peak Number 13 Estra-4,9,11-trien-3-one, 1... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.03	3.44 ng/ul	431078	Perylene-d12	23.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Estra-4,9,11-trien-3-one, 17-.be...	270	C18H22O2	010161-33-8	92
2	Naphthalene, 1-(2-naphthalenyloxy)-	270	C20H14O	000611-49-4	25
3	4,7,12-Trimethylbenz[a]anthracene	270	C21H18	035187-24-7	25
4	Benz(a)anthracene, 6,7,8-trimethyl-	270	C21H18	020627-32-1	25
5	4,4'-Di-tert-butyl-diphenylamine	281	C20H27N	1000370-31-1	20





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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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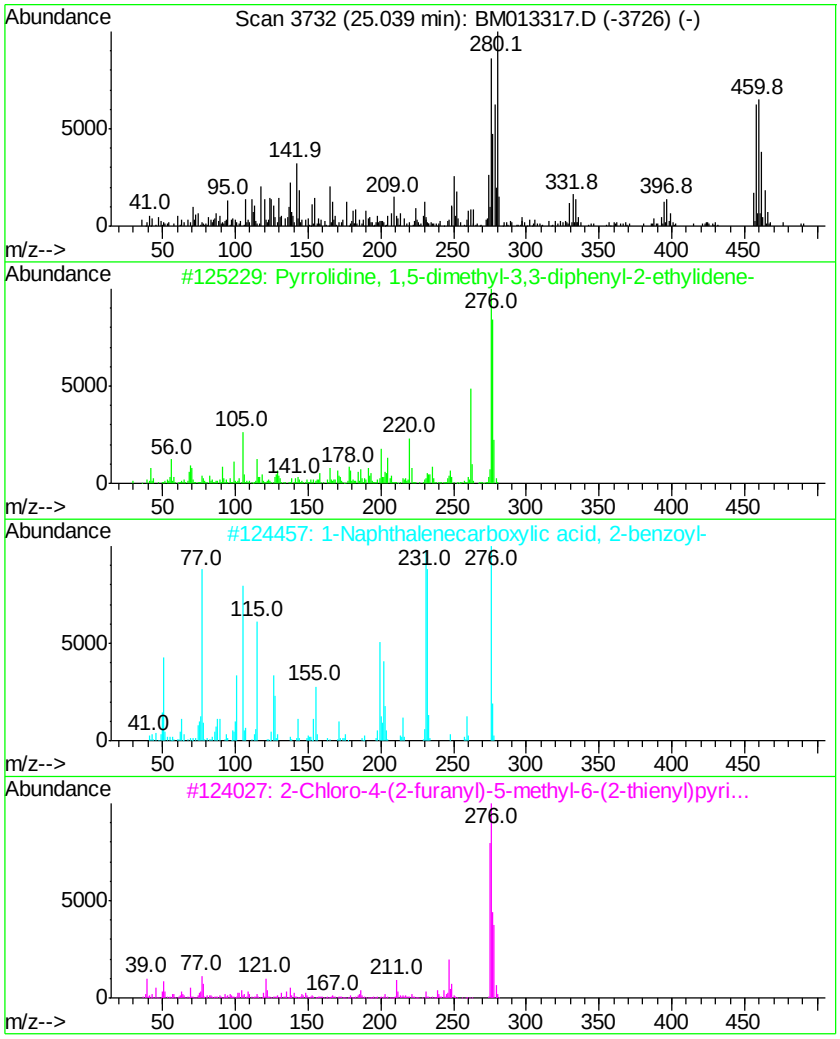
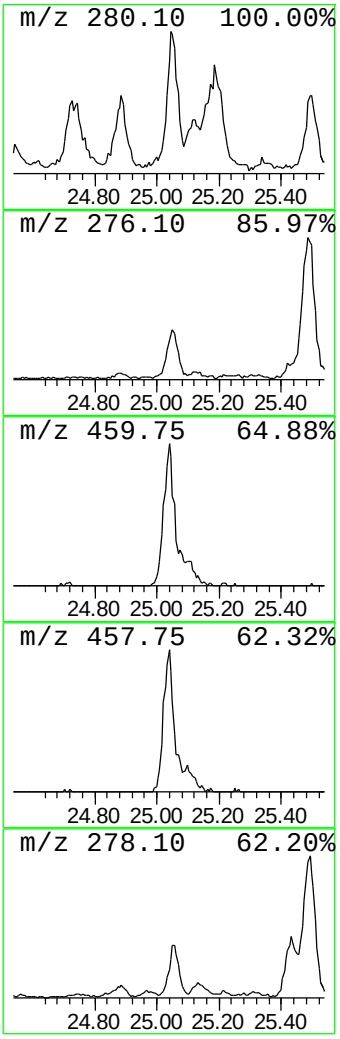
Peak Number 14 unknown-03 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.04	9.54 ng/ul	1196240	Perylene-d12	23.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrrolidine, 1,5-dimethyl-3,3-di...	277	C20H23N	030223-73-5	25
2	1-Naphthalenecarboxylic acid, 2-...	276	C18H12O3	038119-11-8	25
3	2-Chloro-4-(2-furanyl)-5-methyl-...	276	C13H9ClN2OS	131022-80-5	25
4	Dibenzo(a,c)fluoren-13-one	280	C21H12O	063041-47-4	25
5	3-Amino-6-[2-naphthylthio]picoli...	277	C16H11N3S	095095-92-4	25





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Sample : I6759-03
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Instrument : BNA_M
ClientSampleId : F9A02

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

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

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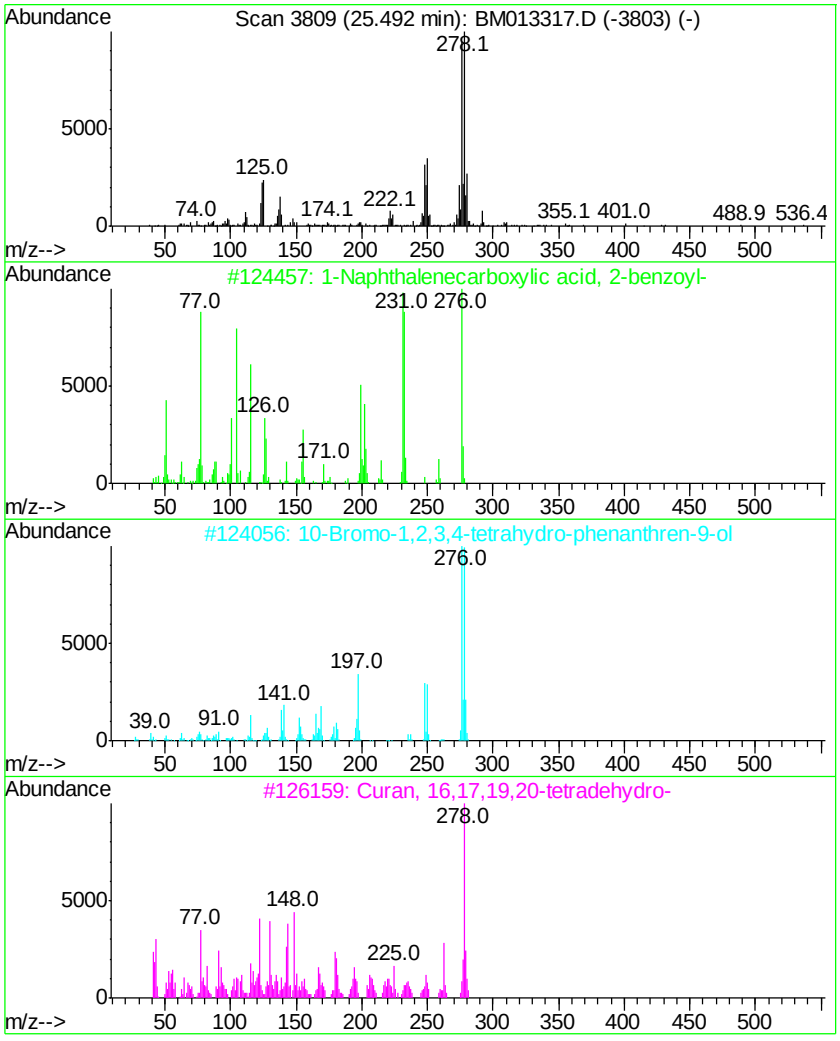
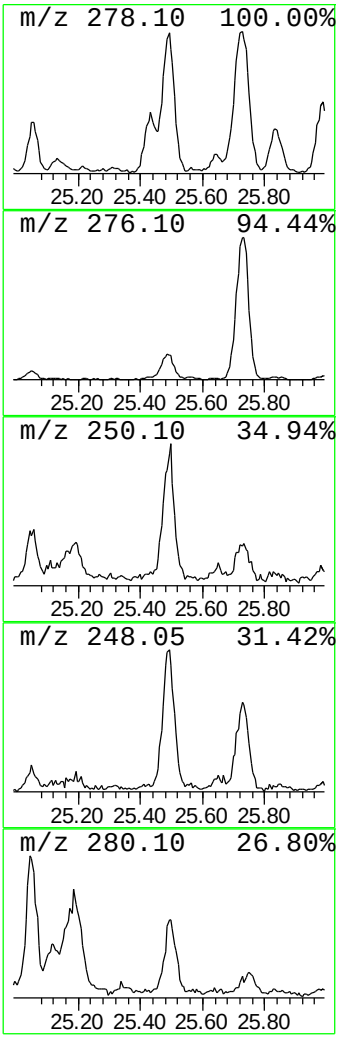
Peak Number 15 1-Naphthalenecarboxylic aci... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.49	10.39 ng/ul	1303080	Perylene-d12	23.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenecarboxylic acid, 2-...	276	C18H12O3	038119-11-8	68
2	10-Bromo-1,2,3,4-tetrahydro-phen...	276	C14H13BrO	1000190-54-9	64
3	Curan, 16,17,19,20-tetradehydro-	278	C19H22N2	056053-17-9	38
4	1,2,3,4-Phenazinetetrol, 7-chloro-	278	C12H7ClN2O4	023774-10-9	38
5	8-Methoxy-quinoline-5-sulfonic a...	278	C13H14N2O3S	1000274-64-1	35





Library Search Compound Report

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Operator : SJ/JU
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Misc :
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Instrument : BNA_M
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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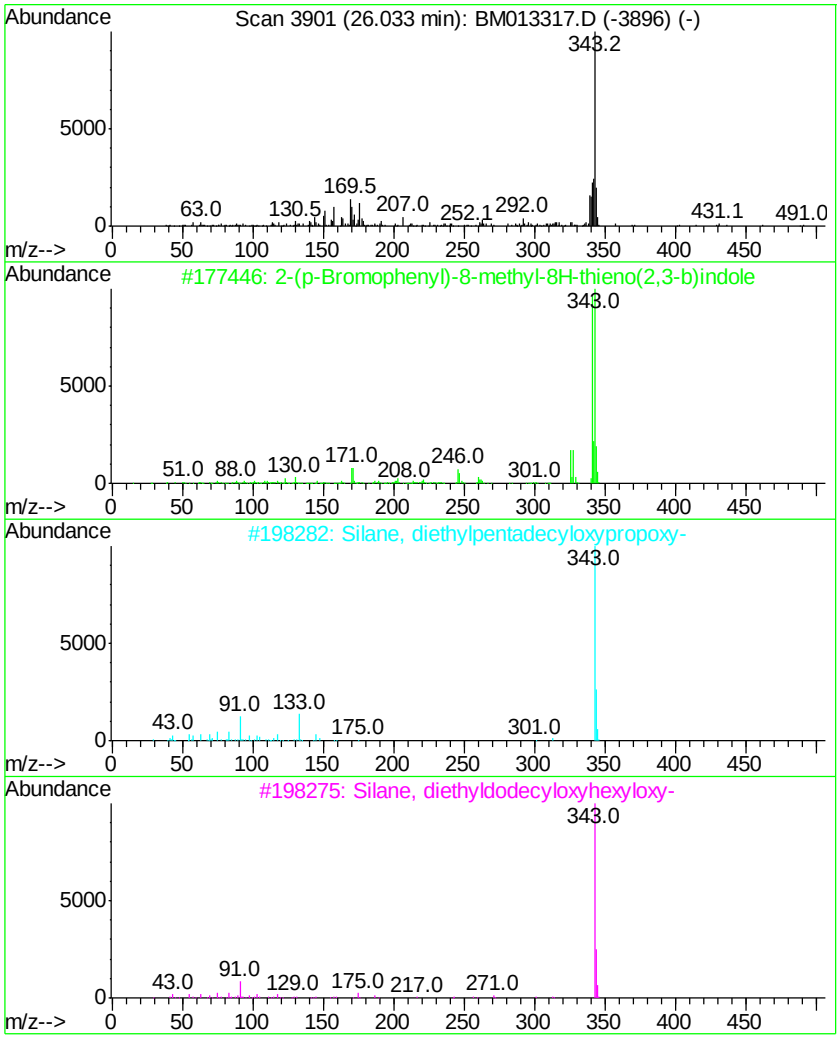
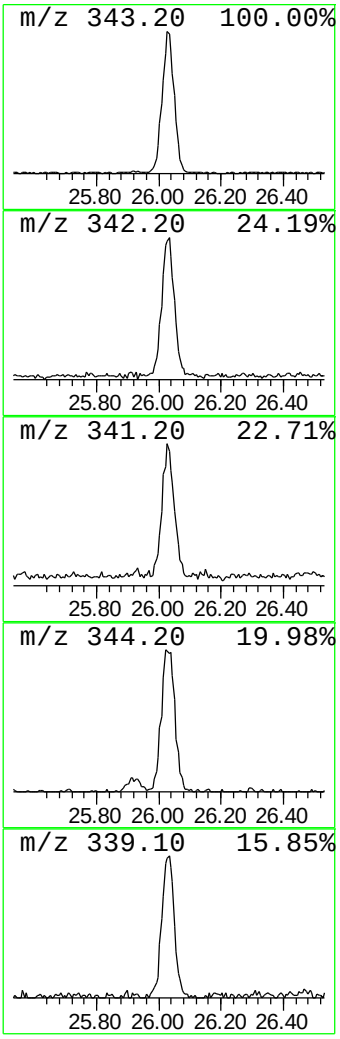
Peak Number 16 unknown-04 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.03	11.59 ng/ul	1453960	Perylene-d12	23.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(p-Bromophenyl)-8-methyl-8H-th...	341	C17H12BrNS	022315-11-3	45
2	Silane, diethylpentadecyloxyprop...	372	C22H48O2Si	1000363-96-2	43
3	Silane, diethyldodecyloxyhexyloxy-	372	C22H48O2Si	1000363-74-0	43
4	Benzamide, 2,3,4-trifluoro-N-(4-...	343	C19H12F3NO2	1000339-09-5	43
5	Benzoic acid, 4-(3,4,5-trimethox...	343	C19H21NO5	304683-21-4	38





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
1(2H)-Acenaphthyl...	16.06	3.2	ng/ul	466960	4	17.08	2902290	20.0
n-Hexadecanoic acid	17.98	4.3	ng/ul	626817	4	17.08	2902290	20.0
(DEL) Alkane: Str...	18.28	3.8	ng/ul	544107	4	17.08	2902290	20.0
(DEL) Alkane: Str...	18.92	3.4	ng/ul	489022	4	17.08	2902290	20.0
Octadecanoic acid	19.26	2.4	ng/ul	432120	5	21.27	3684380	20.0
Acenaphtho(1,2-b)...	19.36	2.4	ng/ul	446401	5	21.27	3684380	20.0
(DEL) Alkane: Cyc...	20.99	7.6	ng/ul	1394600	5	21.27	3684380	20.0
unknown-01	22.33	3.2	ng/ul	595711	5	21.27	3684380	20.0
unknown-02	22.68	4.2	ng/ul	531115	6	23.49	2508130	20.0
Benz(a)anthracene...	22.74	3.5	ng/ul	444201	6	23.49	2508130	20.0
Estra-4,9,11-trie...	24.03	3.4	ng/ul	431078	6	23.49	2508130	20.0
unknown-03	25.04	9.5	ng/ul	1196240	6	23.49	2508130	20.0
1-Naphthalenecarb...	25.49	10.4	ng/ul	1303080	6	23.49	2508130	20.0
unknown-04	26.03	11.6	ng/ul	1453960	6	23.49	2508130	20.0