

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM110717\
 Data File : BM012800.D
 Acq On : 07 Nov 2017 17:56
 Operator : SJ/JU
 Sample : I6164-10
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ET0E0

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 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:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM110417MA.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	3.804	118	122	129	rVB	39430	52432	1.53%	0.121%
2	6.975	655	661	673	rVB2	288117	527058	15.43%	1.218%
3	7.134	681	688	696	rBV	235388	356210	10.43%	0.823%
4	7.334	716	722	733	rVB	302355	486877	14.25%	1.125%
5	7.798	789	801	809	rBV	351818	574460	16.81%	1.328%
6	8.510	916	922	930	rBV2	264707	440694	12.90%	1.019%
7	8.957	992	998	1004	rBV	295472	481436	14.09%	1.113%
8	9.022	1004	1009	1016	rVB2	28478	44836	1.31%	0.104%
9	9.675	1114	1120	1134	rBV	257170	441467	12.92%	1.020%
10	10.216	1206	1212	1226	rBV	394824	674937	19.75%	1.560%
11	10.592	1265	1276	1286	rBV	521510	933317	27.31%	2.157%
12	10.727	1291	1299	1311	rBV	205741	415638	12.16%	0.961%
13	11.610	1441	1449	1461	rBV4	15540	36264	1.06%	0.084%
14	11.910	1496	1500	1511	rVB	22456	39526	1.16%	0.091%
15	12.016	1511	1518	1525	rBV	51100	81547	2.39%	0.188%
16	12.121	1527	1536	1543	rBV2	31970	63416	1.86%	0.147%
17	13.263	1722	1730	1736	rBV	54726	89097	2.61%	0.206%
18	13.845	1823	1829	1834	rBV	827106	1168295	34.19%	2.700%
19	13.892	1834	1837	1846	rVB	391955	550668	16.12%	1.273%
20	14.133	1871	1878	1887	rVB	932703	1370368	40.11%	3.168%
21	14.339	1906	1913	1920	rBV	61695	83019	2.43%	0.192%
22	14.439	1920	1930	1937	rVV2	903628	1352798	39.59%	3.127%
23	14.504	1937	1941	1946	rVB	36109	49756	1.46%	0.115%
24	14.657	1961	1967	1980	rBV	212037	366087	10.71%	0.846%
25	14.839	1994	1998	2001	rBV	22164	35348	1.03%	0.082%
26	15.292	2070	2075	2079	rVB	50555	62785	1.84%	0.145%
27	15.433	2093	2099	2105	rBV	1275222	1771111	51.83%	4.094%
28	15.486	2105	2108	2116	rVB3	103481	148948	4.36%	0.344%
29	15.568	2116	2122	2133	rBV	261737	395388	11.57%	0.914%
30	15.698	2141	2144	2148	rVB	25332	34232	1.00%	0.079%
31	15.798	2155	2161	2167	rBV2	113403	166710	4.88%	0.385%
32	15.921	2176	2182	2188	rBV5	16442	34405	1.01%	0.080%
33	16.151	2216	2221	2224	rBV2	52104	84606	2.48%	0.196%
34	16.727	2311	2319	2323	rBV3	33070	56659	1.66%	0.131%

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35	16.845	2334	2339	2346	rVB	43103	63560	1.86%	0.147%
36	16.945	2351	2356	2359	rVB	39065	45757	1.34%	0.106%
37	17.004	2361	2366	2374	rVB	74204	117983	3.45%	0.273%
38	17.186	2387	2397	2400	rBV2	1192612	1726388	50.53%	3.990%
39	17.227	2400	2404	2409	rVB	1242803	1647368	48.21%	3.808%
40	17.286	2409	2414	2418	rBV	1267634	1831266	53.59%	4.233%
41	17.380	2426	2430	2434	rVB	35403	51474	1.51%	0.119%
42	17.586	2456	2465	2477	rVB2	128606	222998	6.53%	0.515%
43	17.680	2477	2481	2483	rBV2	39786	49727	1.46%	0.115%
44	17.962	2524	2529	2536	rBV6	35456	67391	1.97%	0.156%
45	18.074	2541	2548	2551	rVV2	236568	423250	12.39%	0.978%
46	18.109	2551	2554	2560	rVV	140543	198044	5.80%	0.458%
47	18.256	2573	2579	2583	rVV3	173036	287491	8.41%	0.665%
48	18.292	2583	2585	2591	rVB	55202	64861	1.90%	0.150%
49	18.368	2594	2598	2602	rBV3	43882	60442	1.77%	0.140%
50	18.562	2626	2631	2634	rBV	93377	126229	3.69%	0.292%
51	18.603	2634	2638	2643	rVB	257666	328797	9.62%	0.760%
52	18.980	2698	2702	2703	rBV2	37608	46996	1.38%	0.109%
53	18.998	2703	2705	2709	rVB4	34809	47759	1.40%	0.110%
54	19.121	2721	2726	2733	rBV2	80684	124511	3.64%	0.288%
55	19.245	2738	2747	2754	rVV	2580919	3416871	100.00%	7.898%
56	19.356	2760	2766	2776	rBV3	117979	211361	6.19%	0.489%
57	19.445	2776	2781	2782	rBV4	31090	47722	1.40%	0.110%
58	19.486	2785	2788	2792	rVB	64726	88518	2.59%	0.205%
59	19.574	2797	2803	2806	rBV2	1866340	2808455	82.19%	6.492%
60	19.603	2806	2808	2815	rVV	1801979	2276773	66.63%	5.263%
61	19.674	2817	2820	2825	rVB2	66826	95889	2.81%	0.222%
62	19.786	2835	2839	2845	rVB	88079	113938	3.33%	0.263%
63	19.845	2845	2849	2856	rVB	60709	103530	3.03%	0.239%
64	19.927	2858	2863	2864	rBV	73103	111134	3.25%	0.257%
65	19.986	2870	2873	2876	rBV	34089	35799	1.05%	0.083%
66	20.062	2881	2886	2888	rBV3	64595	110257	3.23%	0.255%
67	20.097	2888	2892	2899	rVB	186267	314159	9.19%	0.726%
68	20.197	2905	2909	2913	rBV	134154	165791	4.85%	0.383%
69	20.256	2913	2919	2923	rBV2	101132	166553	4.87%	0.385%
70	20.292	2923	2925	2929	rVB	48613	59328	1.74%	0.137%
71	20.339	2930	2933	2937	rVB4	30178	39632	1.16%	0.092%

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72	20.397	2940	2943	2946	rBV	50632	66681	1.95%	0.154%
73	20.444	2948	2951	2955	rVB2	45615	58090	1.70%	0.134%
74	20.844	3016	3019	3026	rVB	139141	182130	5.33%	0.421%
75	21.009	3042	3047	3051	rBV3	185346	295125	8.64%	0.682%
76	21.050	3051	3054	3055	rVV	144042	165906	4.86%	0.383%
77	21.074	3055	3058	3065	rVV3	229920	528518	15.47%	1.222%
78	21.144	3066	3070	3077	rVB2	154041	239956	7.02%	0.555%
79	21.362	3100	3107	3111	rBV2	1751332	3100855	90.75%	7.167%
80	21.403	3111	3114	3123	rVB	888080	1117148	32.70%	2.582%
81	21.515	3130	3133	3139	rBV	138835	189708	5.55%	0.439%
82	21.680	3158	3161	3166	rVB2	101300	145112	4.25%	0.335%
83	22.962	3372	3379	3384	rBV	654425	1578800	46.21%	3.649%
84	23.444	3457	3461	3465	rBV	301502	510537	14.94%	1.180%
85	23.497	3465	3470	3475	rVV2	871716	1657914	48.52%	3.832%
86	23.544	3475	3478	3487	rVB	476705	788009	23.06%	1.821%
87	23.644	3489	3495	3501	rBV	812394	1469930	43.02%	3.398%

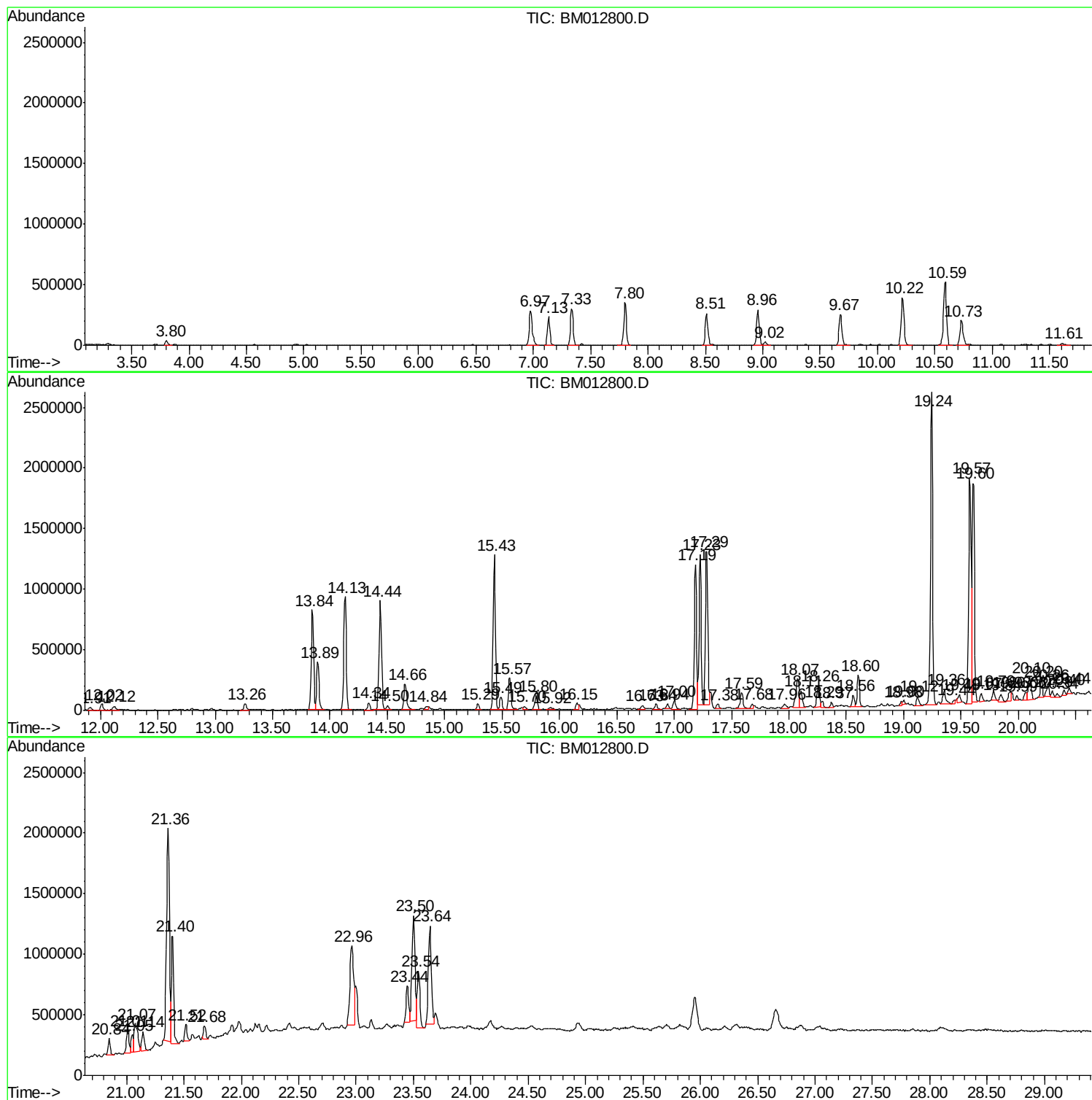
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Instrument :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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Instrument :
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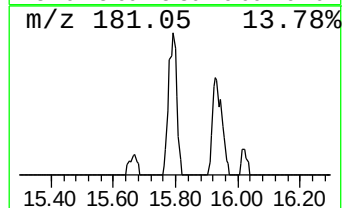
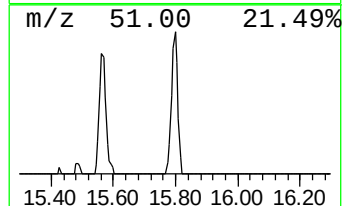
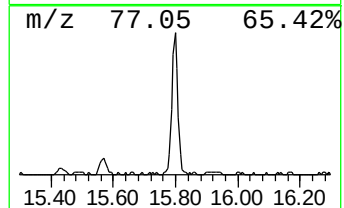
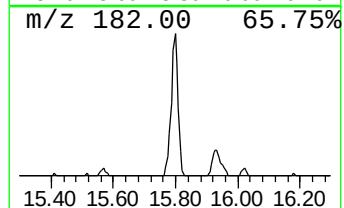
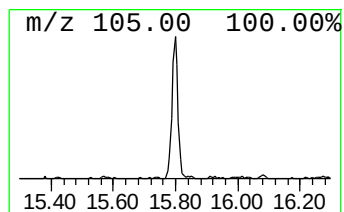
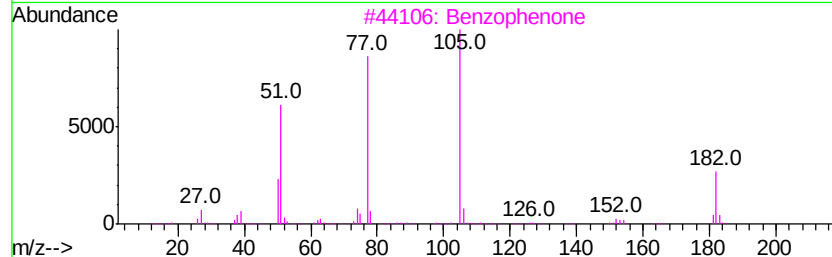
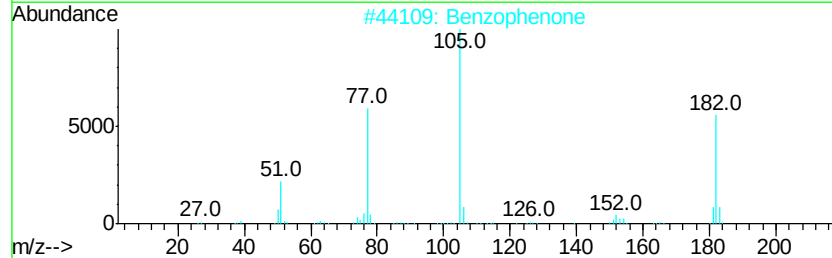
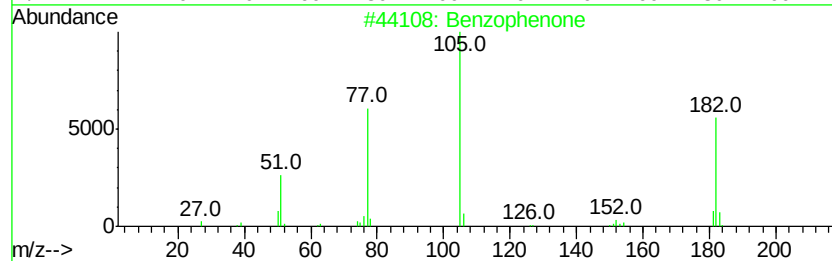
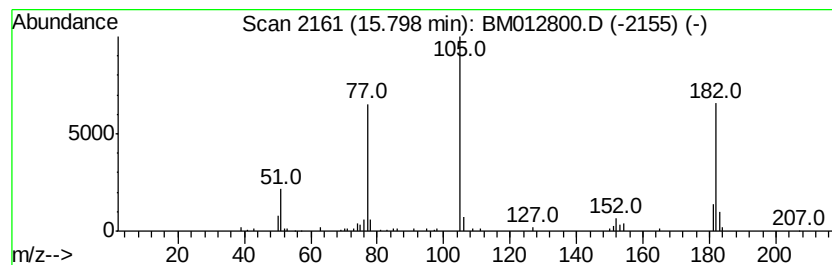
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzophenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.80	2.46 ng/ul	166710	Acenaphthene-d10	14.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzophenone	182	C13H10O	000119-61-9	96
2		Benzophenone	182	C13H10O	000119-61-9	96
3		Benzophenone	182	C13H10O	000119-61-9	95
4		Benzophenone	182	C13H10O	000119-61-9	95
5		Benzophenone	182	C13H10O	000119-61-9	91



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ALS Vial : 15 Sample Multiplier: 1

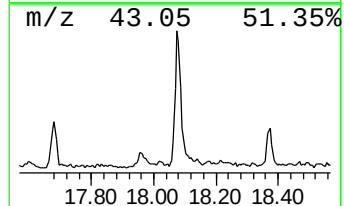
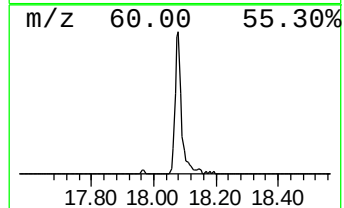
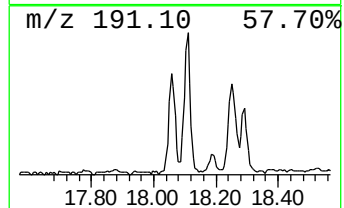
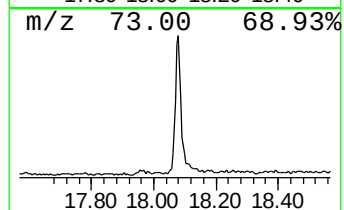
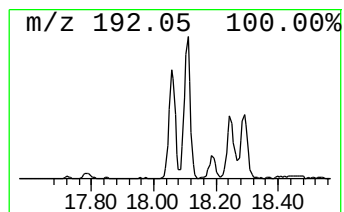
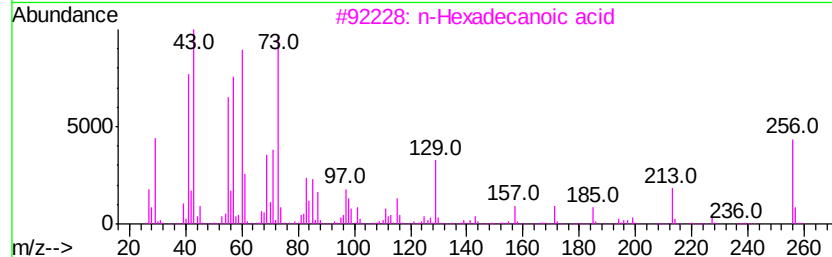
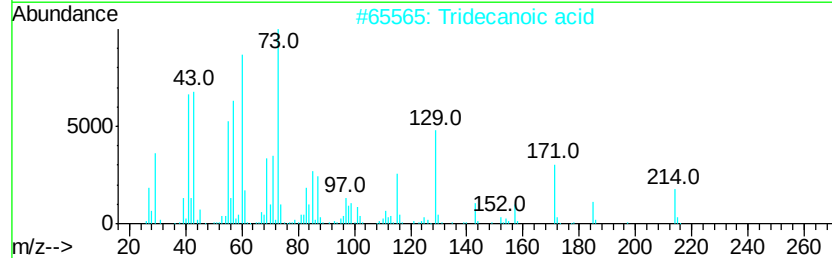
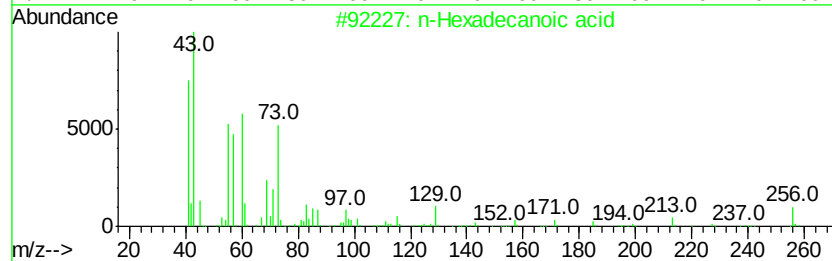
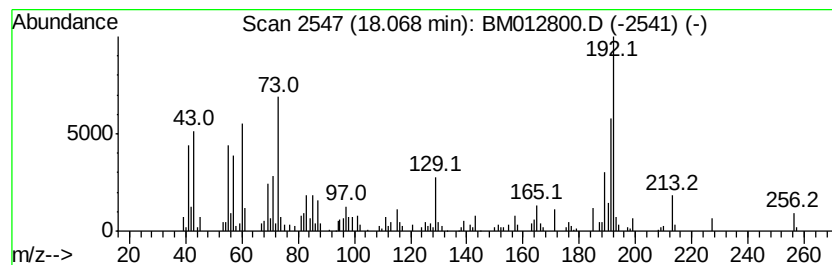
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.07	4.90 ng/ul	423250	Phenanthrene-d10		17.19	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	98
2	Tridecanoic acid		214	C13H26O2	000638-53-9	95
3	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	95
4	Tridecanoic acid		214	C13H26O2	000638-53-9	91
5	Pentadecanoic acid		242	C15H30O2	001002-84-2	74



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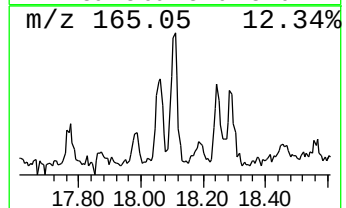
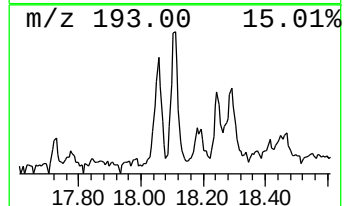
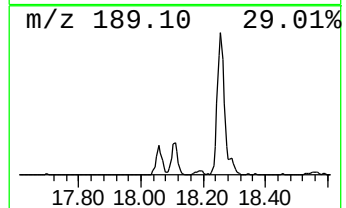
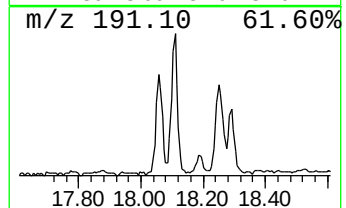
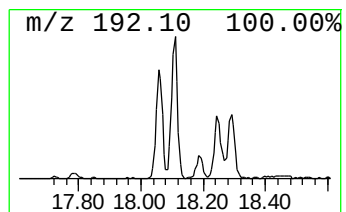
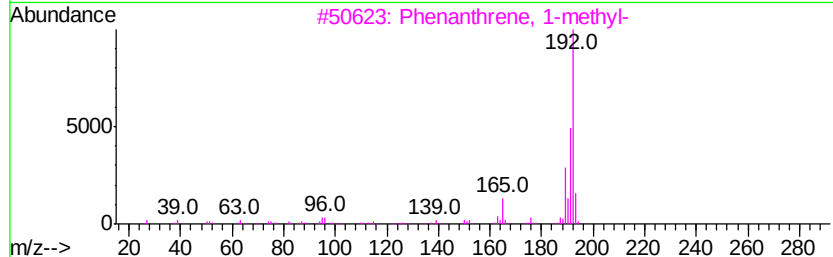
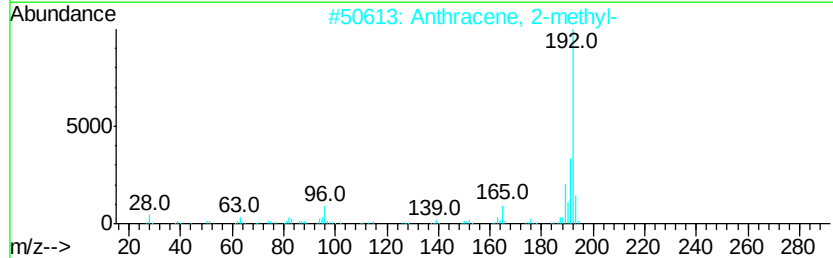
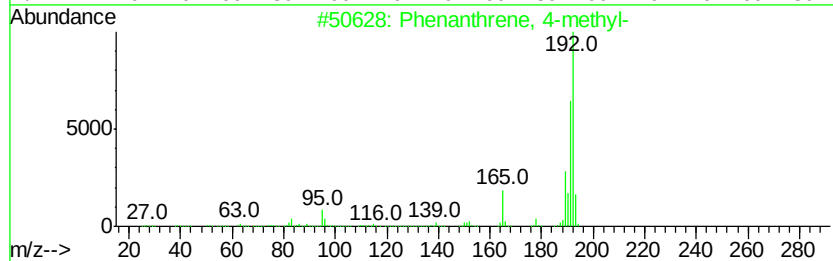
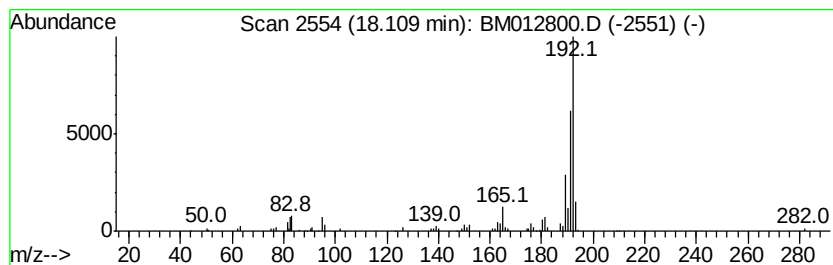
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Phenanthrene, 4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.11	2.29 ng/ul	198044	Phenanthrene-d10	17.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	93
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
4		5H-Dibenzofa,d]cycloheptene	192	C15H12	000256-81-5	91
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91



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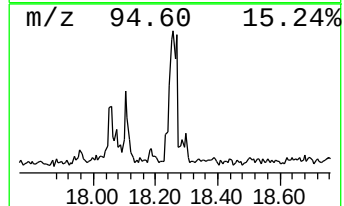
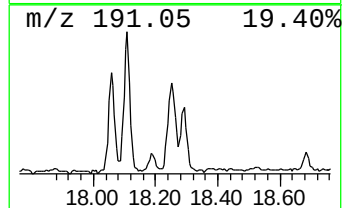
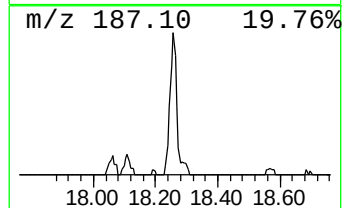
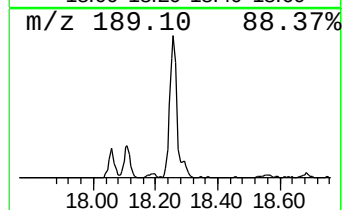
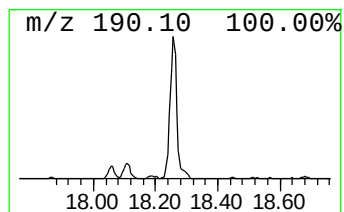
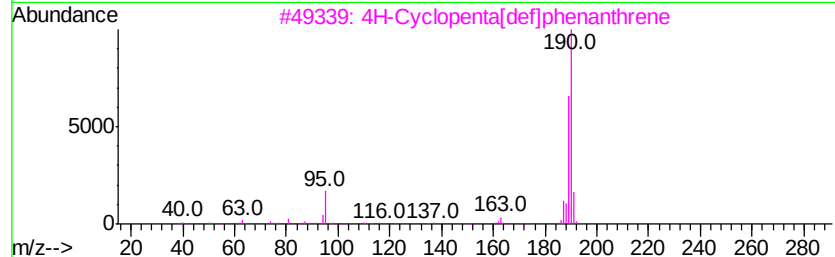
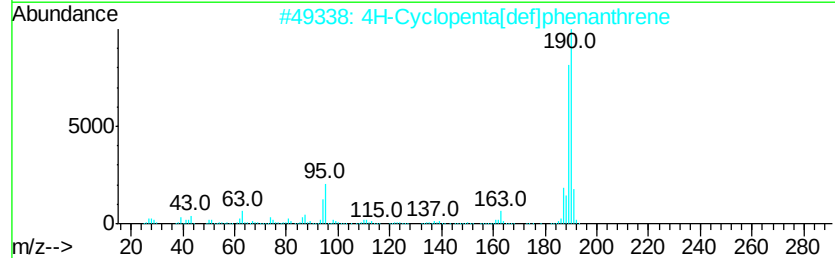
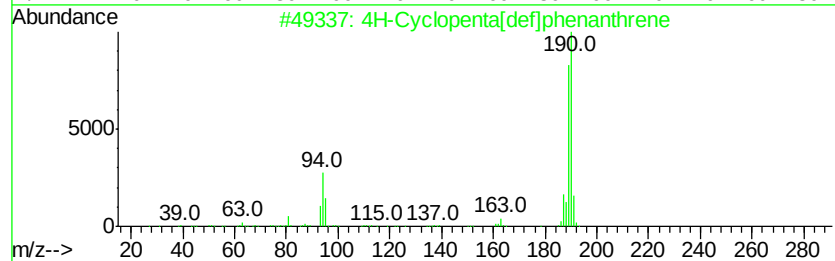
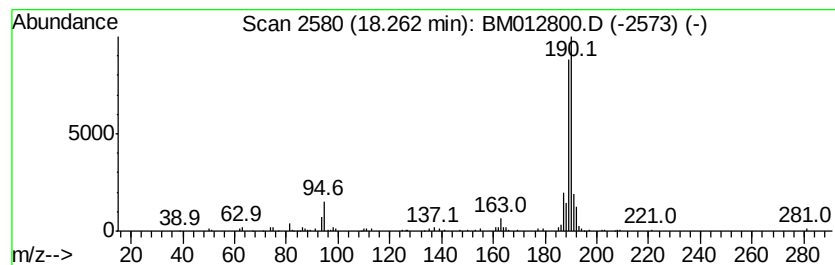
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.
18.26	3.33 ng/ul	287491	Phenanthrene-d10		17.19
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	81
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	72
4	6H-Cyclobuta[flk]phenanthrene	190	C15H10	083469-43-6	56
5	Methyl diselenide	190	C2H6Se2	007101-31-7	55



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Data File : BM012800.D
Acq On : 07 Nov 2017 17:56
Operator : SJ/JU
Sample : I6164-10
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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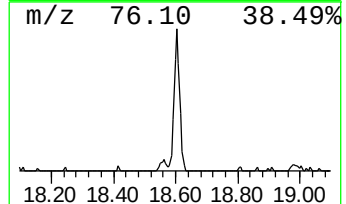
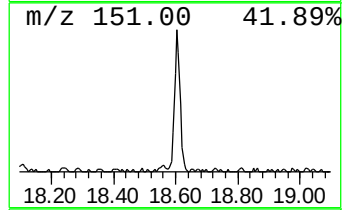
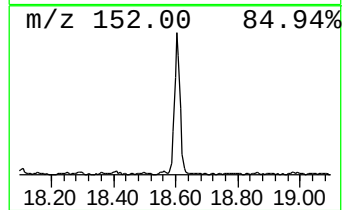
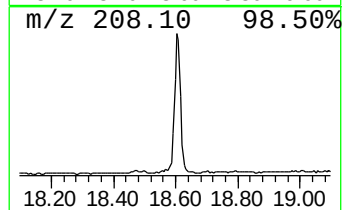
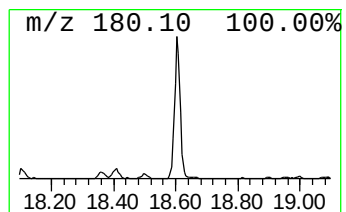
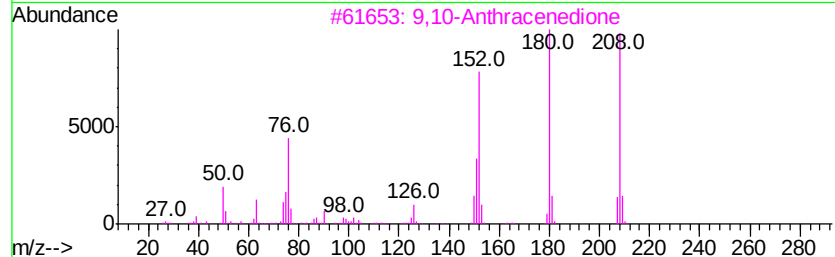
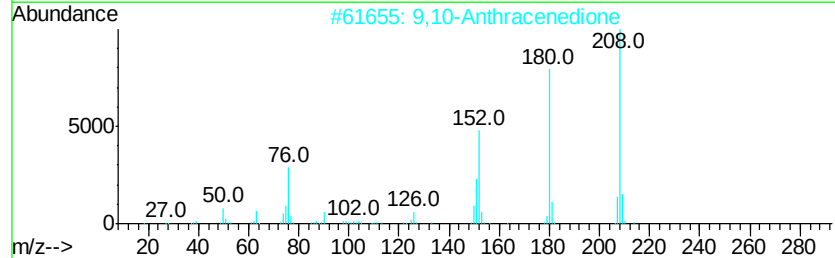
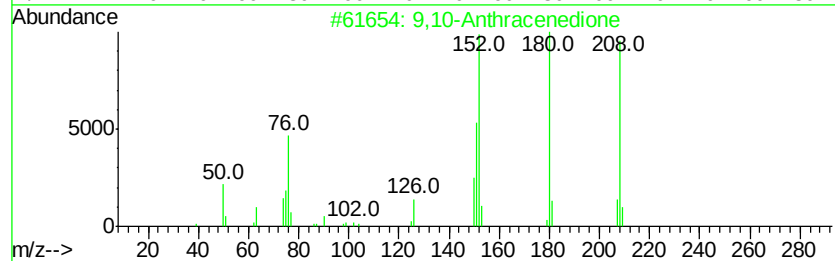
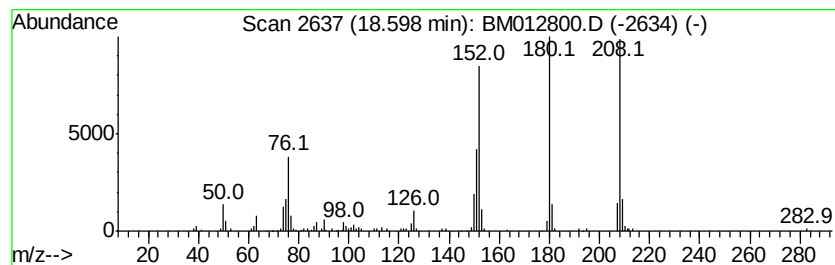
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 9,10-Anthracenedione Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.60	3.81 ng/ul	328797	Phenanthrene-d10	17.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Anthracenedione	208	C14H8O2	000084-65-1	97
2		9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
3		9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
4		Benzo[cl]cinoline	180	C12H8N2	000230-17-1	86
5		9H-Fluoren-9-one	180	C13H8O	000486-25-9	83



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA M\DATA\BM110717\
Data File : BM012800.D
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Sample : I6164-10
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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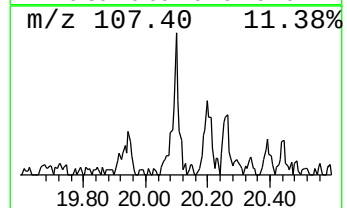
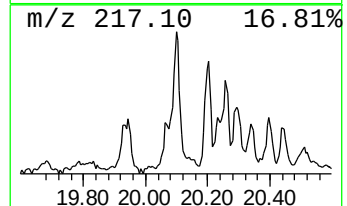
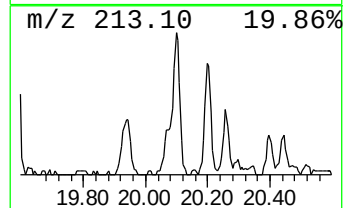
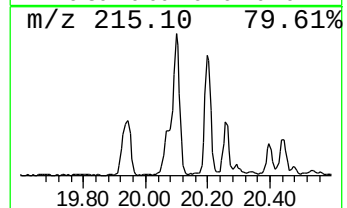
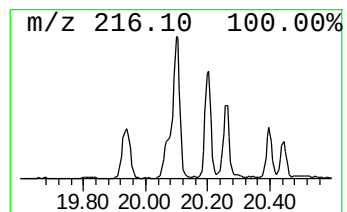
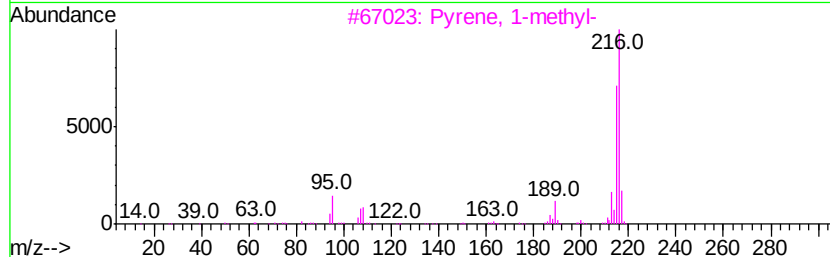
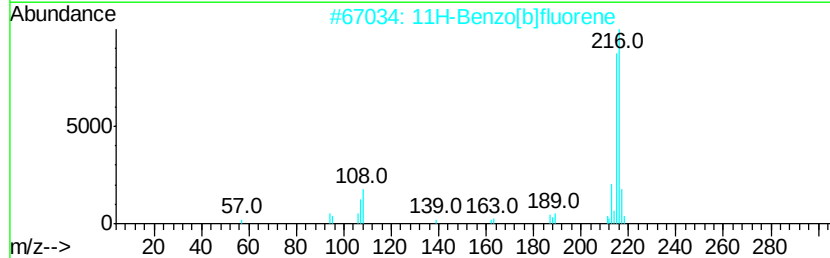
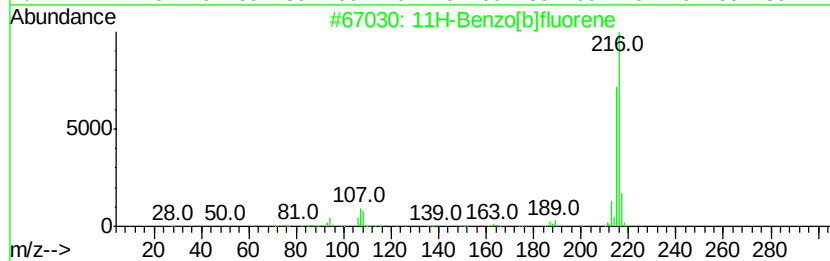
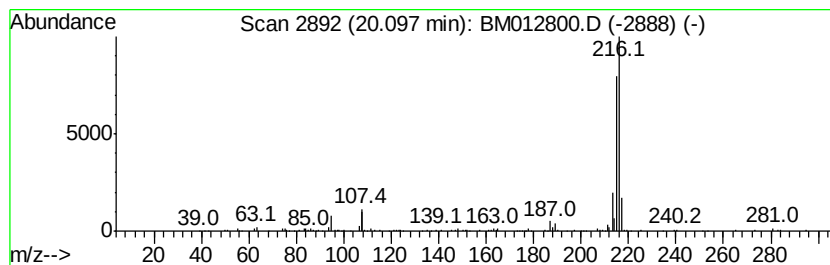
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 11H-Benzo[b]fluorene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.10	2.03 ng/ul	314159	Chrysene-d12	21.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
3		Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
4		Pyrene, 2-methyl-	216	C17H12	003442-78-2	90
5		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA M\DATA\BM110717\
Data File : BM012800.D
Acq On : 07 Nov 2017 17:56
Operator : SJ/JU
Sample : I6164-10
Misc :
ALS Vial : 15 Sample Multiplier: 1

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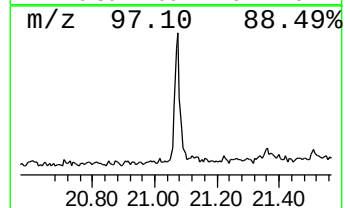
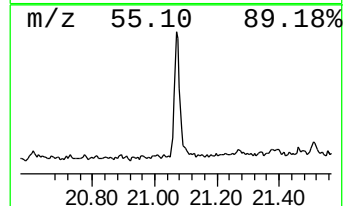
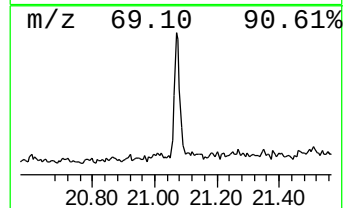
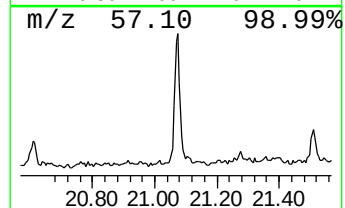
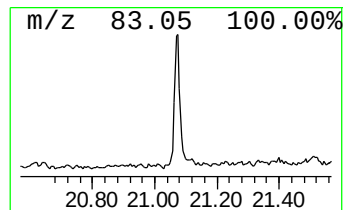
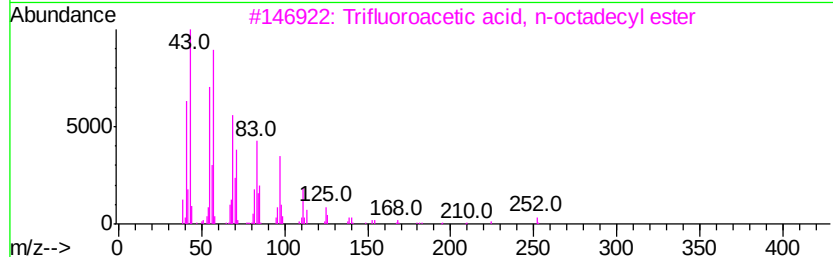
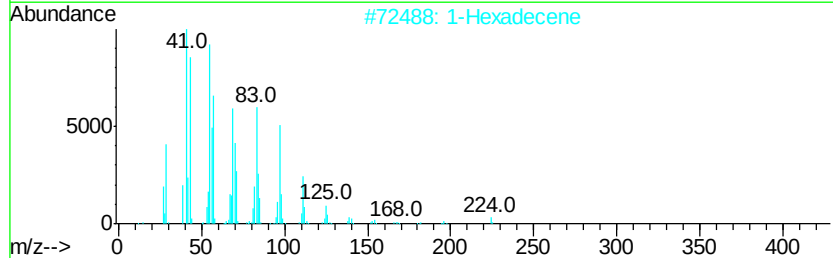
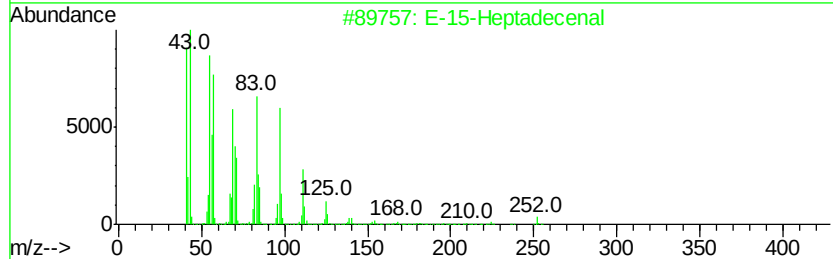
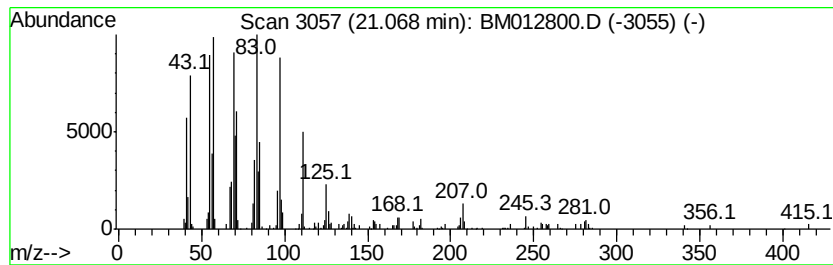
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 E-15-Heptadecenal Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.07	3.41 ng/ul	528518	Chrysene-d12	21.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	E-15-Heptadecenal			252	C17H32O	1000130-97-9	93
2	1-Hexadecene			224	C16H32	000629-73-2	91
3	Trifluoroacetic acid, n-octadecyl...			366	C20H37F3O2	079392-43-1	91
4	17-Pentatriacontene			491	C35H70	006971-40-0	91
5	Heptafluorobutyric acid, n-octadecyl...			466	C22H37F7O2	000400-57-7	91



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA M\DATA\BM110717\
Data File : BM012800.D
Acq On : 07 Nov 2017 17:56
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Sample : I6164-10
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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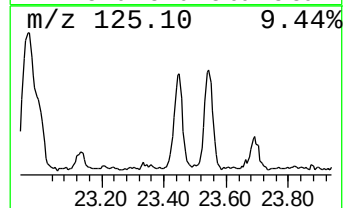
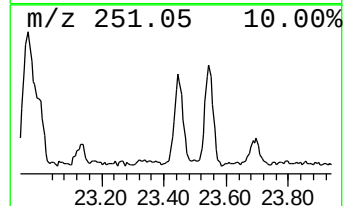
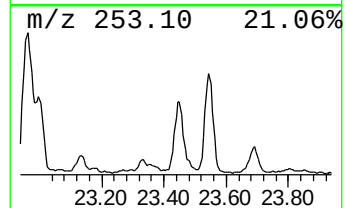
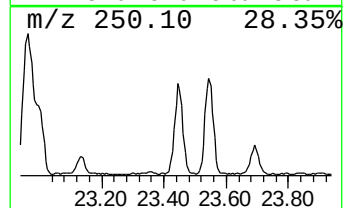
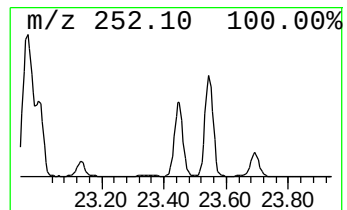
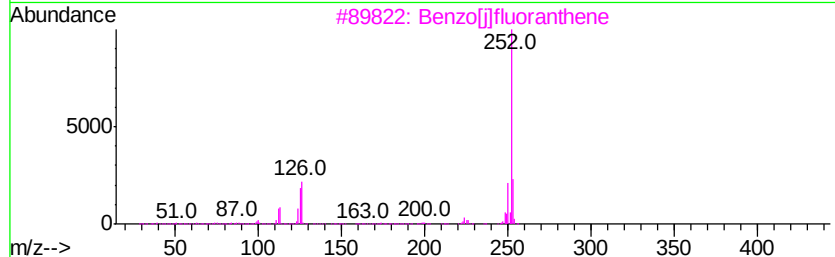
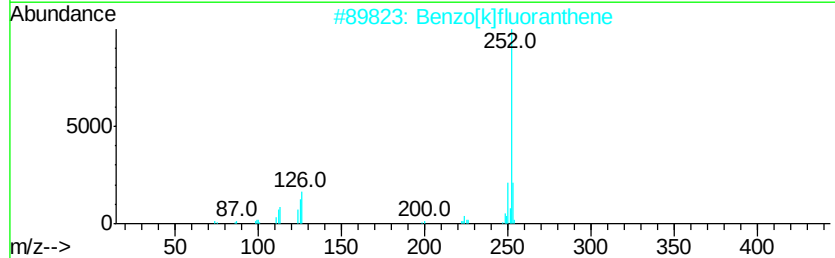
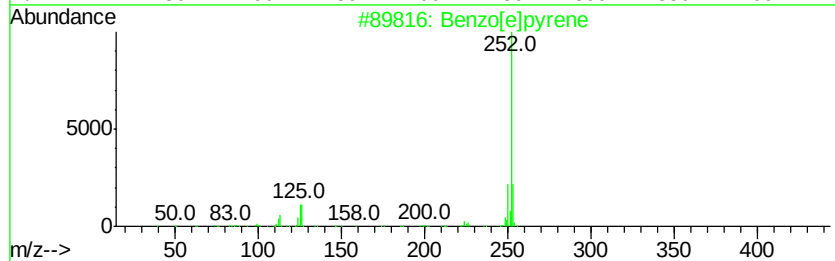
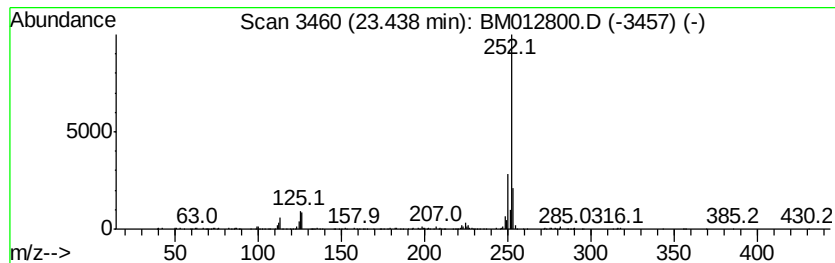
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.44	6.95 ng/ul	510537	Perylene-d12	23.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	95
2		Benzo[k]fluoranthene	252	C20H12	000207-08-9	94
3		Benzo[i]fluoranthene	252	C20H12	000205-82-3	94
4		Perylene	252	C20H12	000198-55-0	94
5		Benzo[a]pyrene	252	C20H12	000050-32-8	93



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_M\DATA\BM110717\
Data File : BM012800.D
Acq On : 07 Nov 2017 17:56
Operator : SJ/JU
Sample : I6164-10
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM110417MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	15.80	2.5	ng/ul	166710	3	14.44	1352800	20.0
n-Hexadecanoic acid	18.07	4.9	ng/ul	423250	4	17.19	1726390	20.0
Phenanthrene, 4-m...	18.11	2.3	ng/ul	198044	4	17.19	1726390	20.0
4H-Cyclopenta[def...	18.26	3.3	ng/ul	287491	4	17.19	1726390	20.0
9,10-Anthracenedione	18.60	3.8	ng/ul	328797	4	17.19	1726390	20.0
11H-Benzo[b]fluorene	20.10	2.0	ng/ul	314159	5	21.36	3100860	20.0
E-15-Heptadecenal	21.07	3.4	ng/ul	528518	5	21.36	3100860	20.0
Benzo[e]pyrene	23.44	7.0	ng/ul	510537	6	23.64	1469930	20.0