

Data Path : Z:\HPCHEM1\BNA M\DATA\BM030118\
 Data File : BM014443.D
 Acq On : 01 Mar 2018 21:22
 Operator : SJ/JU
 Sample : J1678-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BE5T2

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-BM021418.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	4.669	281	286	288	rBV2	44162	67938	1.21%	0.092%
2	6.722	629	635	650	rVB	672923	1237241	22.11%	1.684%
3	6.863	654	659	670	rVV	484435	761706	13.61%	1.036%
4	7.051	685	691	699	rBV	734037	1189469	21.25%	1.619%
5	7.510	763	769	778	rBV	458531	741905	13.26%	1.010%
6	8.239	886	893	907	rBV	834497	1506141	26.91%	2.049%
7	8.675	960	967	978	rBV	742481	1247184	22.29%	1.697%
8	9.386	1080	1088	1101	rBV	652202	1194214	21.34%	1.625%
9	9.928	1173	1180	1199	rBV	834371	1679094	30.00%	2.285%
10	10.280	1231	1240	1246	rBV	683142	1141999	20.41%	1.554%
11	10.333	1246	1249	1254	rVB2	39974	60009	1.07%	0.082%
12	10.445	1260	1268	1288	rBV	683282	1383654	24.72%	1.883%
13	13.598	1798	1804	1808	rBV	1715697	2362707	42.22%	3.215%
14	13.645	1808	1812	1818	rVB	469931	650965	11.63%	0.886%
15	13.863	1842	1849	1862	rBV	1982347	2979540	53.24%	4.054%
16	14.068	1878	1884	1890	rBV2	87780	139618	2.49%	0.190%
17	14.174	1895	1902	1908	rVV2	1009691	1578416	28.21%	2.148%
18	14.239	1908	1913	1921	rVB3	183051	317038	5.67%	0.431%
19	14.463	1945	1951	1967	rBV	407969	1113136	19.89%	1.515%
20	14.580	1967	1971	1977	rVV	80754	139726	2.50%	0.190%
21	14.980	2032	2039	2043	rBV	64762	84954	1.52%	0.116%
22	15.033	2043	2048	2052	rVB2	101143	134238	2.40%	0.183%
23	15.180	2066	2073	2079	rBV	2297502	3341948	59.72%	4.547%
24	15.233	2079	2082	2088	rVB	169057	226513	4.05%	0.308%
25	15.345	2095	2101	2107	rBV2	291833	500638	8.95%	0.681%
26	15.539	2129	2134	2140	rBV4	44285	77749	1.39%	0.106%
27	15.898	2190	2195	2197	rBV2	124133	165493	2.96%	0.225%
28	15.933	2197	2201	2210	rVB2	172944	267208	4.77%	0.364%
29	16.145	2232	2237	2243	rBV7	39762	72364	1.29%	0.098%
30	16.239	2247	2253	2259	rVV9	37720	81976	1.46%	0.112%
31	16.604	2310	2315	2318	rBV4	63646	104135	1.86%	0.142%
32	16.692	2321	2330	2335	rVV7	127275	321292	5.74%	0.437%
33	16.751	2337	2340	2347	rVB4	88228	150082	2.68%	0.204%
34	16.851	2353	2357	2364	rVB10	42636	65440	1.17%	0.089%

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35	16.939	2364	2372	2375	rBV	1278151	1846151	32.99%	2.512%
36	16.980	2375	2379	2383	rVV	1579750	2102661	37.57%	2.861%
37	17.039	2383	2389	2393	rVV	2558827	3720168	66.48%	5.062%
38	17.068	2393	2394	2400	rVB	345299	350379	6.26%	0.477%
39	17.245	2420	2424	2429	rBV8	51384	103014	1.84%	0.140%
40	17.327	2434	2438	2440	rVV4	91674	130925	2.34%	0.178%
41	17.356	2440	2443	2452	rVV	187545	336987	6.02%	0.459%
42	17.433	2452	2456	2458	rVV4	83108	113782	2.03%	0.155%
43	17.468	2458	2462	2468	rVV7	44885	106996	1.91%	0.146%
44	17.533	2468	2473	2479	rVB10	55501	107921	1.93%	0.147%
45	17.727	2503	2506	2513	rBV8	48436	77922	1.39%	0.106%
46	17.815	2514	2521	2523	rBV	156491	240077	4.29%	0.327%
47	17.851	2523	2527	2536	rVV4	542275	1064504	19.02%	1.448%
48	17.945	2540	2543	2547	rVV	72015	117862	2.11%	0.160%
49	18.009	2549	2554	2558	rVV3	267762	436793	7.81%	0.594%
50	18.045	2558	2560	2565	rVB2	96154	111693	2.00%	0.152%
51	18.133	2571	2575	2579	rBV7	70063	118789	2.12%	0.162%
52	18.321	2603	2607	2613	rVB	108082	163017	2.91%	0.222%
53	18.380	2613	2617	2623	rBV2	187408	276078	4.93%	0.376%
54	18.886	2699	2703	2709	rVB2	232219	348546	6.23%	0.474%
55	19.009	2719	2724	2729	rVB	2261477	3110150	55.58%	4.232%
56	19.139	2743	2746	2754	rVB5	147932	232320	4.15%	0.316%
57	19.350	2776	2782	2785	rBV	3451165	5059233	90.40%	6.884%
58	19.380	2785	2787	2795	rVB	1839745	1940088	34.67%	2.640%
59	19.562	2815	2818	2824	rVB	128216	153833	2.75%	0.209%
60	19.880	2869	2872	2878	rVB2	325197	454762	8.13%	0.619%
61	19.980	2886	2889	2893	rBV	189222	211905	3.79%	0.288%
62	20.791	3025	3027	3031	rBV	205142	209334	3.74%	0.285%
63	20.862	3036	3039	3048	rVB5	566444	842915	15.06%	1.147%
64	20.980	3056	3059	3064	rVB	964960	1077292	19.25%	1.466%
65	21.086	3070	3077	3082	rBV2	3308592	5596201	100.00%	7.615%
66	21.150	3082	3088	3092	rVV3	2228140	4241338	75.79%	5.771%
67	21.191	3092	3095	3102	rVB	1046561	1355395	24.22%	1.844%
68	21.433	3133	3136	3140	rVB2	374996	421869	7.54%	0.574%
69	21.750	3187	3190	3194	rVV5	322475	374883	6.70%	0.510%
70	21.797	3194	3198	3210	rVB	978891	1724478	30.82%	2.347%
71	22.656	3340	3344	3347	rBV	1202059	1757252	31.40%	2.391%

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72	23.197	3431	3436	3440	rBV2	1698092	2745929	49.07%	3.736%
73	23.327	3454	3458	3464	rBV2	734489	1244825	22.24%	1.694%
74	23.803	3534	3539	3545	rVB2	975559	1777050	31.75%	2.418%

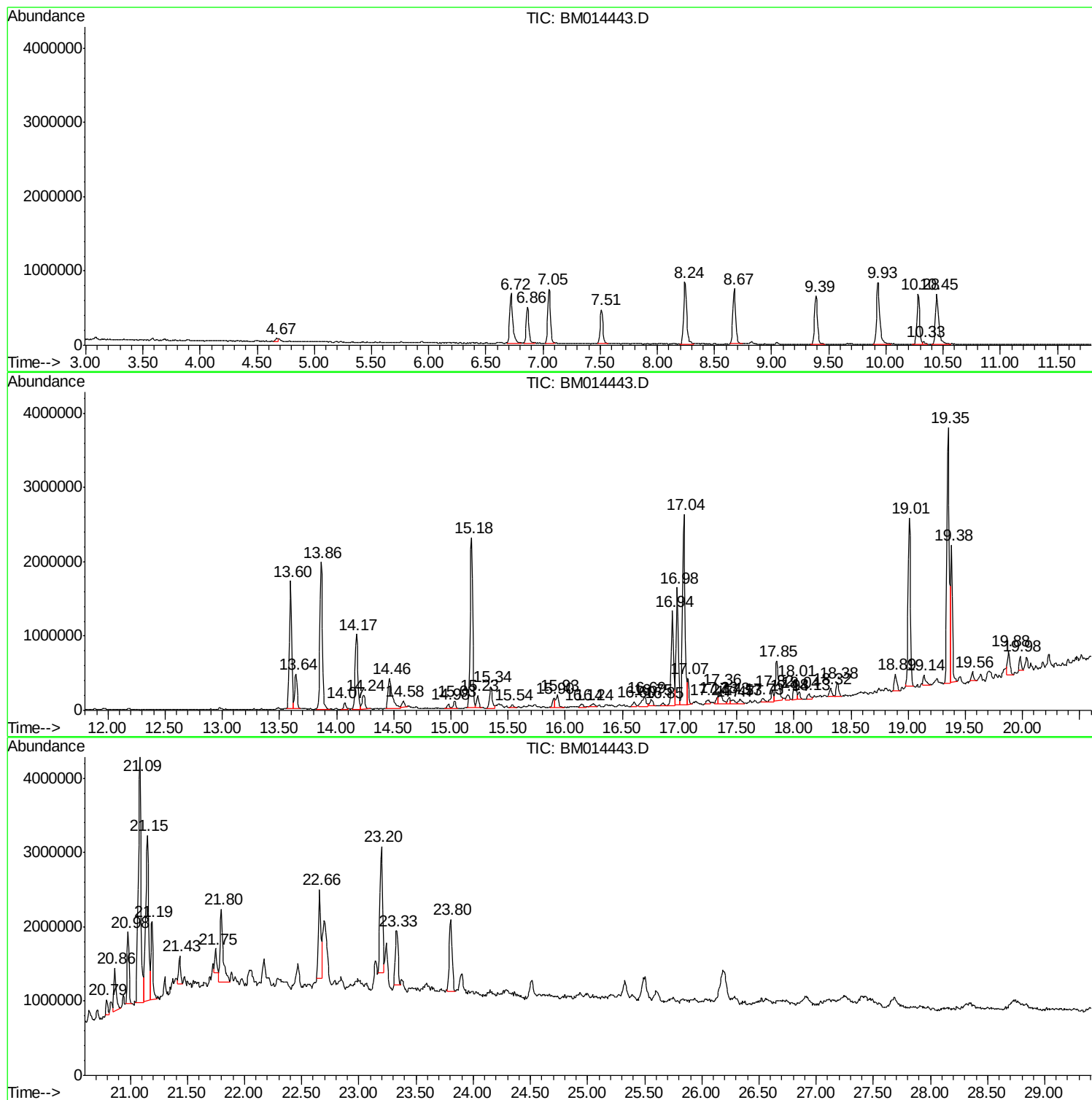
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

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



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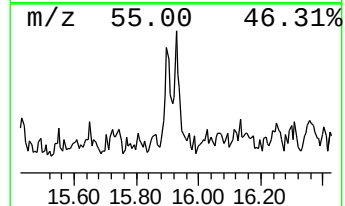
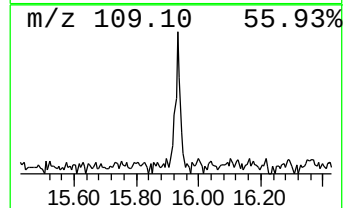
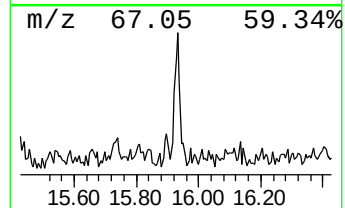
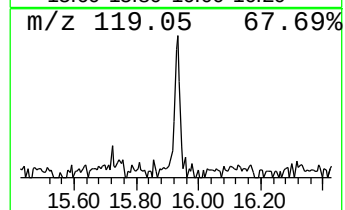
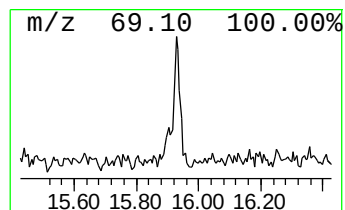
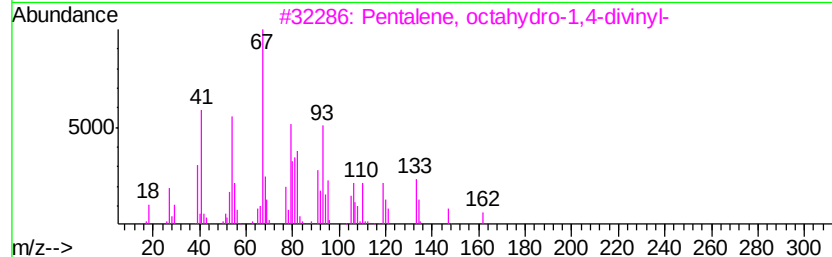
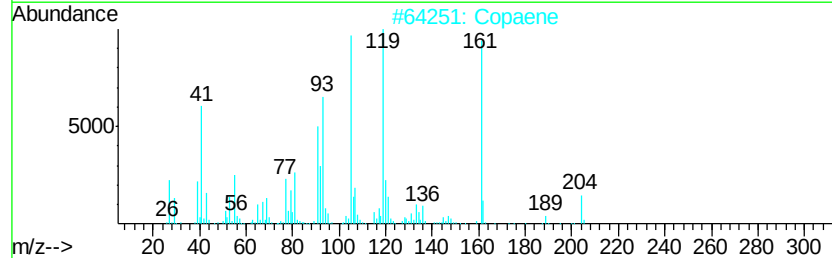
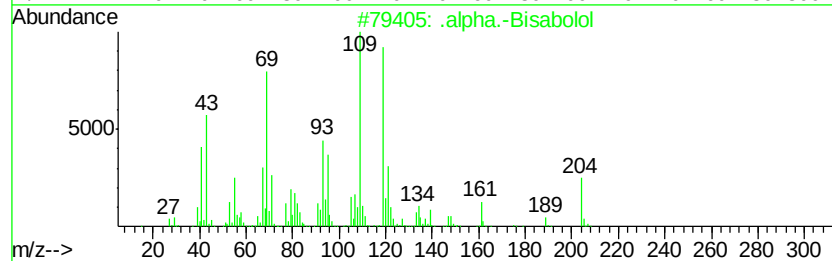
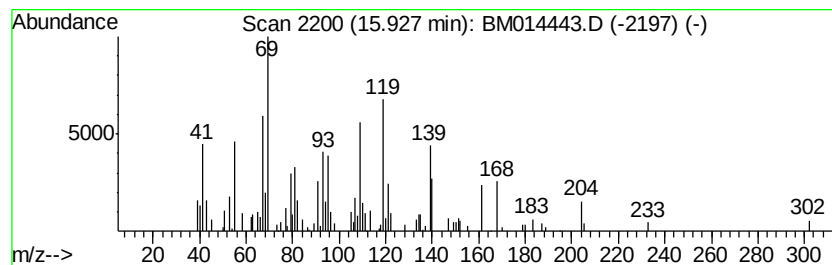
Instrument :
BNA_M
ClientSampled :
BE5T2

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.93	2.89 ng/ul	267208	Phenanthrene-d10	16.94	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Bisabolol	222	C15H26O	000515-69-5	25
2	Copaene	204	C15H24	003856-25-5	25
3	Pentalene, octahydro-1,4-divinyl-	162	C12H18	017572-84-8	25
4	(1,3-Dimethyl-2-methylene-cyclop...	140	C9H16O	1000190-16-2	22
5	3-Methyl-3-buten-1-ol, pentaflu...	232	C8H9F5O2	1000352-29-2	14



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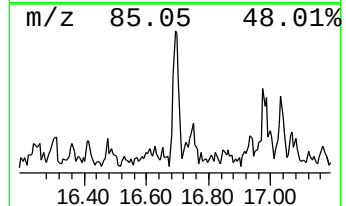
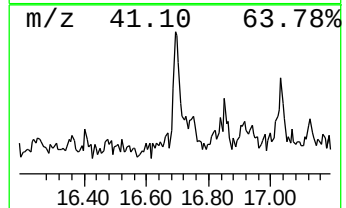
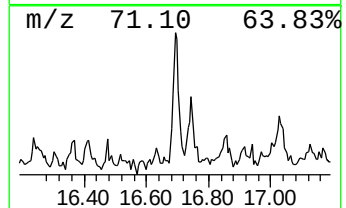
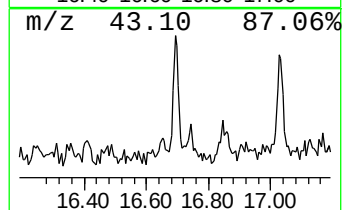
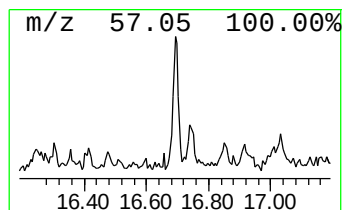
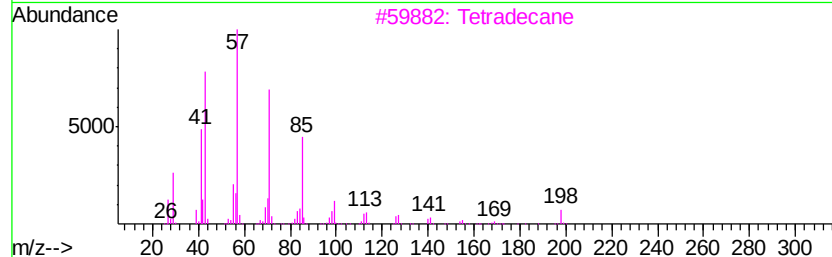
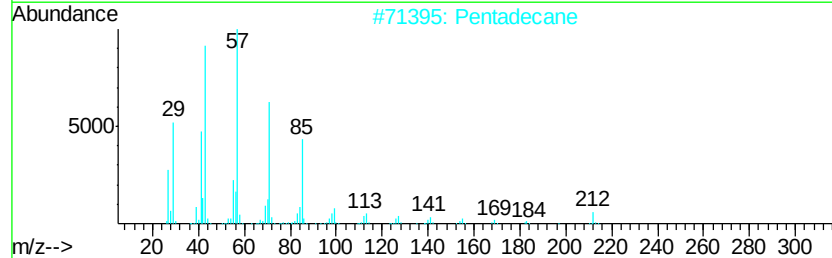
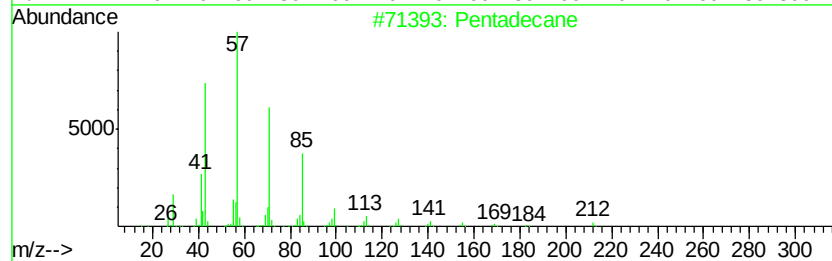
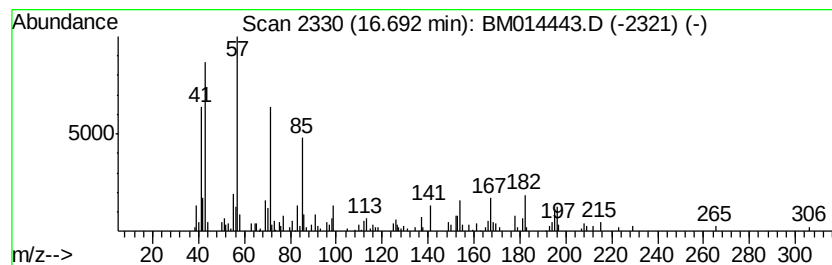
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.69	3.48 ng/ul	321292	Phenanthrene-d10	16.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane	212	C15H32	000629-62-9	70
2			Pentadecane	212	C15H32	000629-62-9	55
3			Tetradecane	198	C14H30	000629-59-4	49
4			Octacosane	394	C28H58	000630-02-4	49
5			Undecane, 5-ethyl-5-propyl-	226	C16H34	002755-07-9	49



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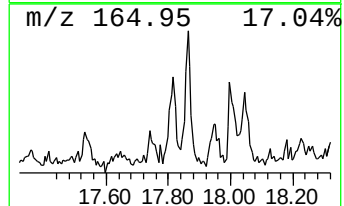
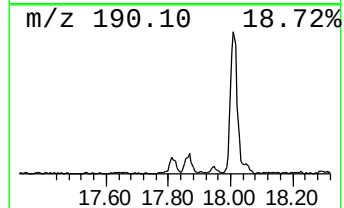
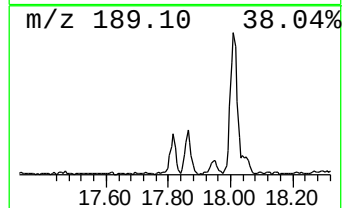
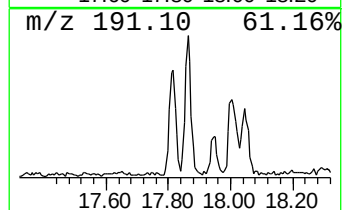
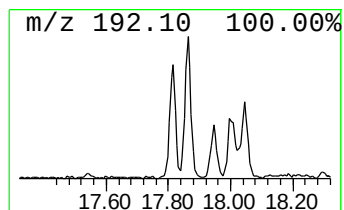
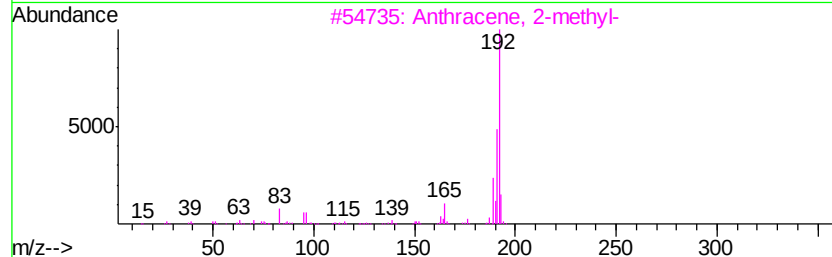
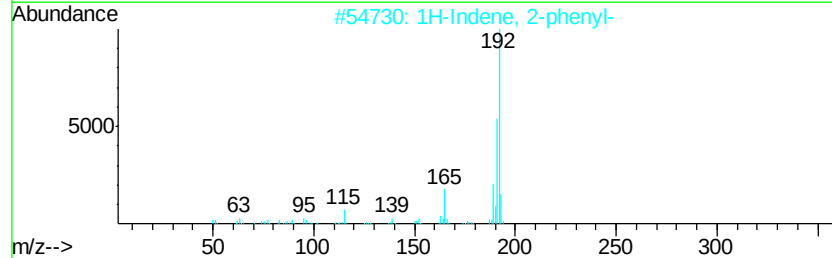
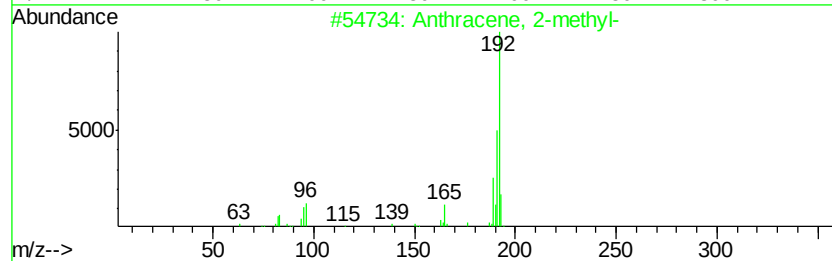
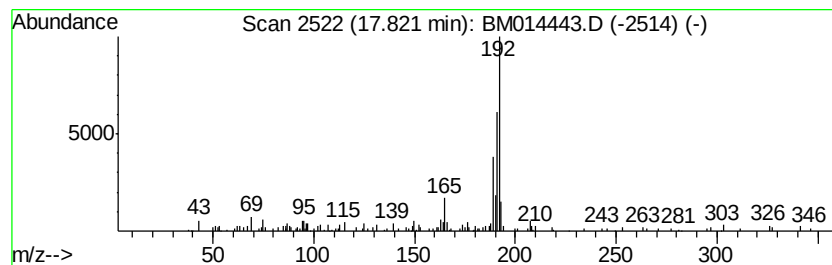
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Anthracene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.82	2.60 ng/ul	240077	Phenanthrene-d10	16.94	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
2	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	93
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	91



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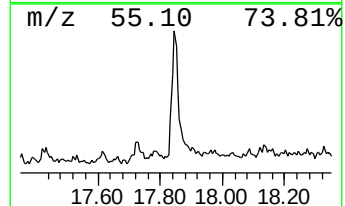
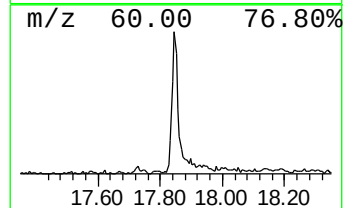
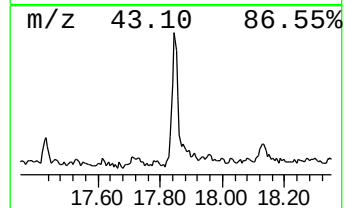
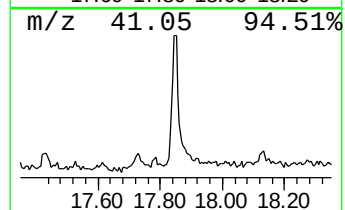
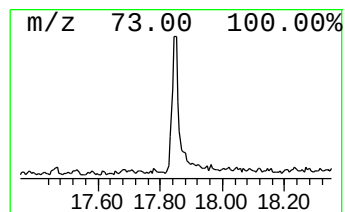
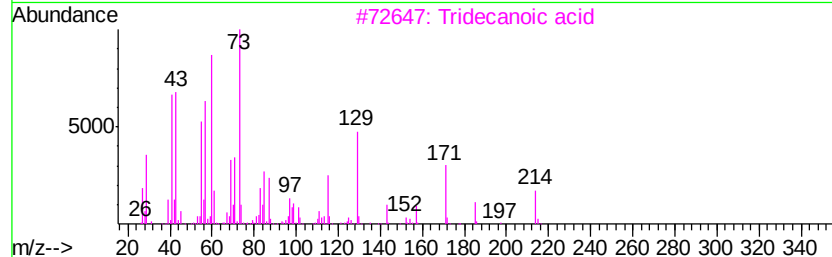
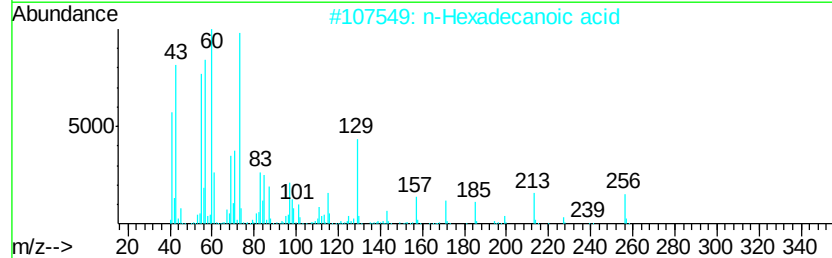
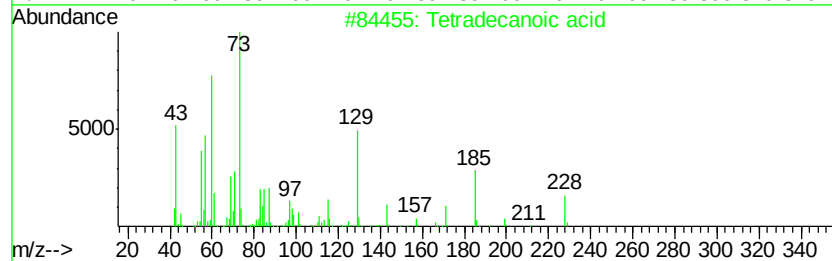
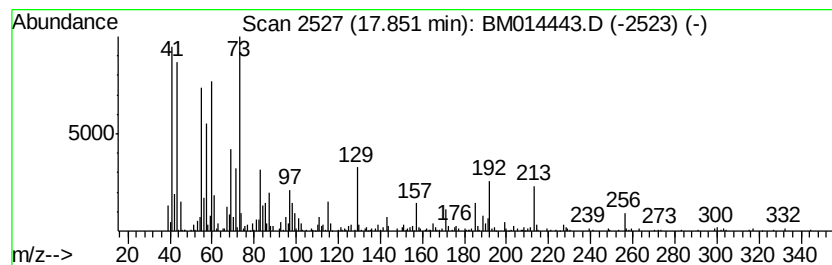
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Tetradecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
17.85	11.53 ng/ul	1064500	Phenanthrene-d10	16.94		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecanoic acid	228	C14H28O2	000544-63-8	93
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	92
3		Tridecanoic acid	214	C13H26O2	000638-53-9	70
4		Tridecanoic acid	214	C13H26O2	000638-53-9	62
5		Tridecanoic acid	214	C13H26O2	000638-53-9	58



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030118\
Data File : BM014443.D
Acq On : 01 Mar 2018 21:22
Operator : SJ/JU
Sample : J1678-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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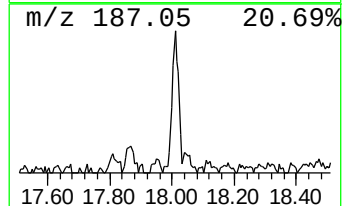
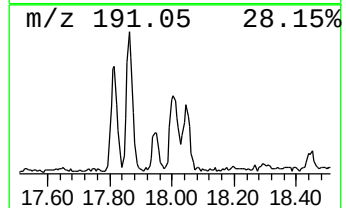
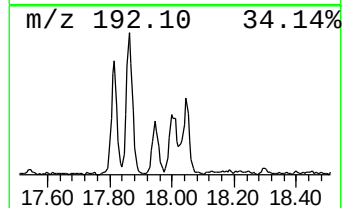
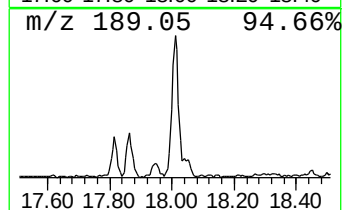
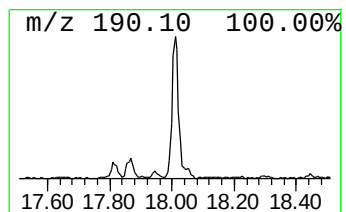
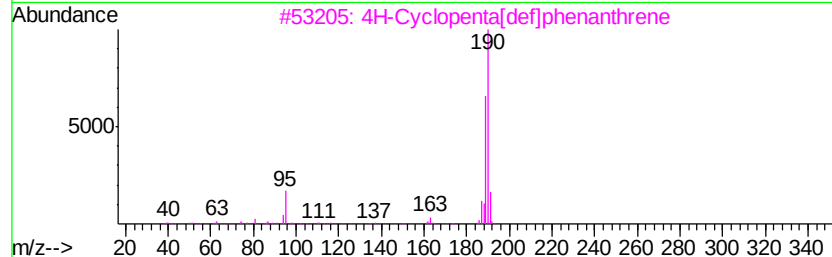
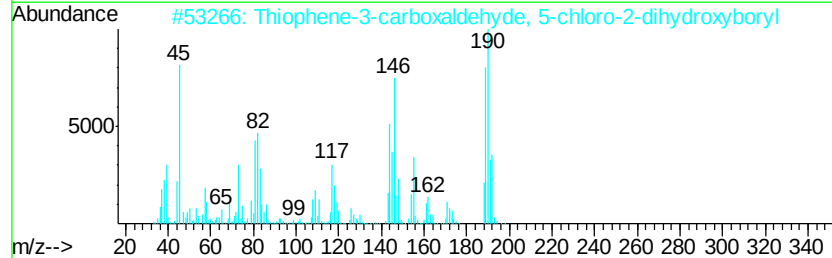
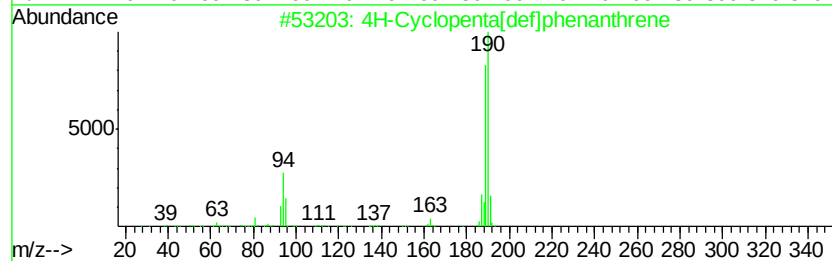
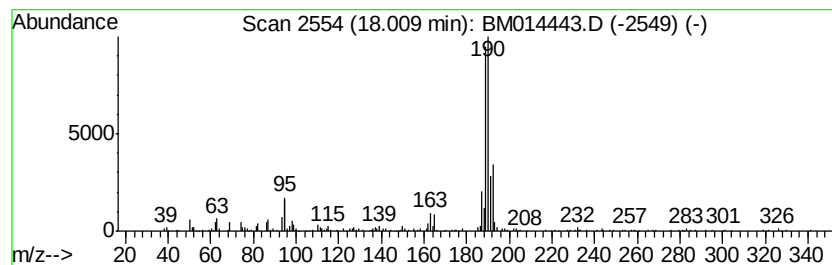
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.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.
18.01	4.73 ng/ul	436793	Phenanthrene-d10	16.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2		Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	72
3		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	72
4		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
5		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	52



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030118\
Data File : BM014443.D
Acq On : 01 Mar 2018 21:22
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Sample : J1678-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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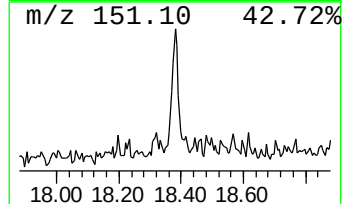
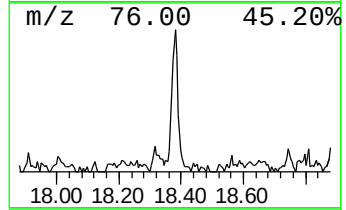
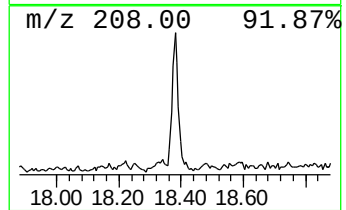
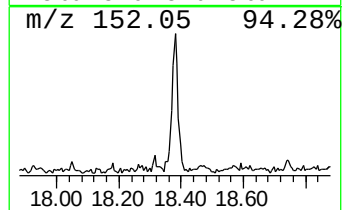
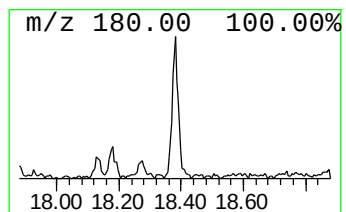
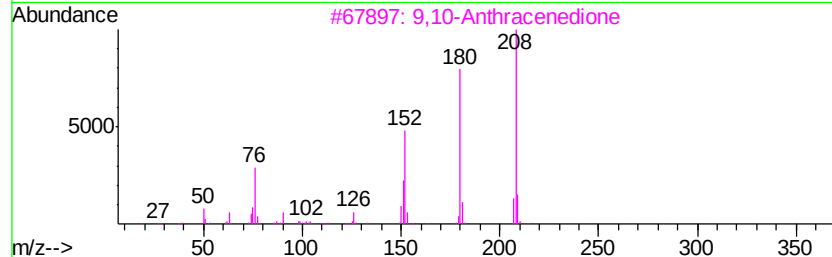
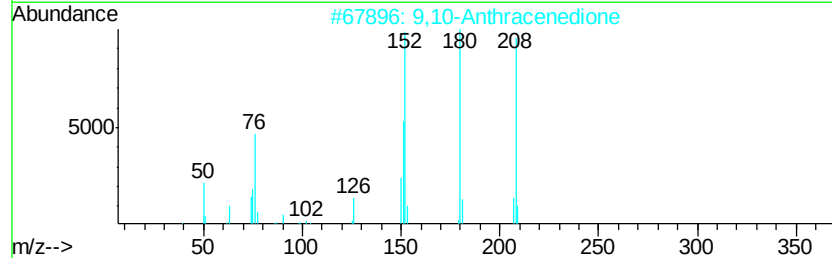
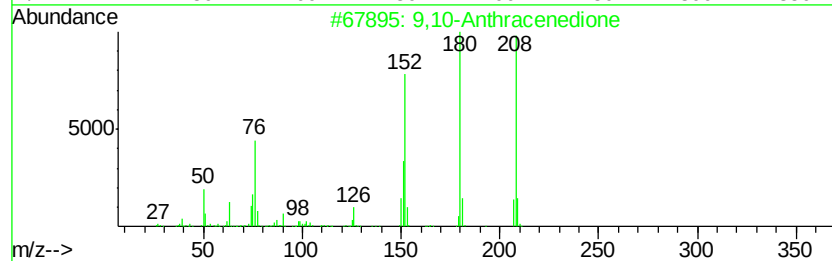
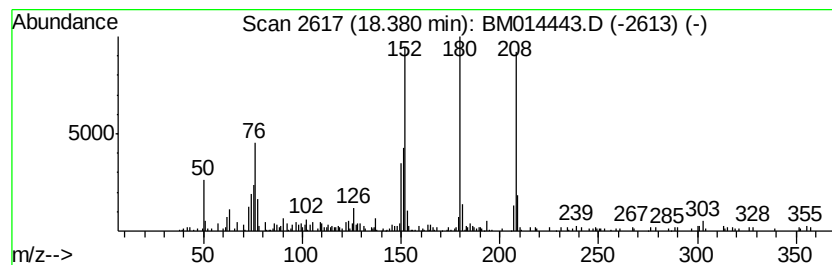
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 9,10-Anthracenedione Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.38	2.99 ng/ul	276078	Phenanthrene-d10	16.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Anthracenedione	208	C14H8O2	000084-65-1	99
2		9,10-Anthracenedione	208	C14H8O2	000084-65-1	97
3		9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
4		Benzo[cl]cinoline	180	C12H8N2	000230-17-1	90
5		9H-Fluoren-9-one	180	C13H8O	000486-25-9	86



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030118\
Data File : BM014443.D
Acq On : 01 Mar 2018 21:22
Operator : SJ/JU
Sample : J1678-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

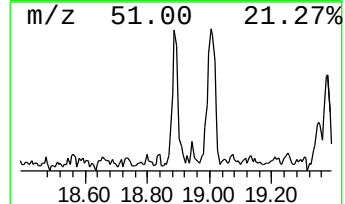
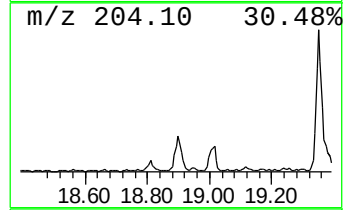
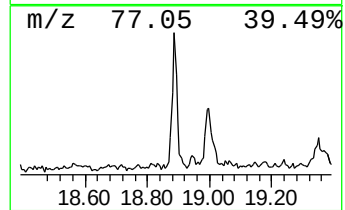
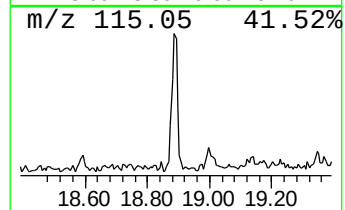
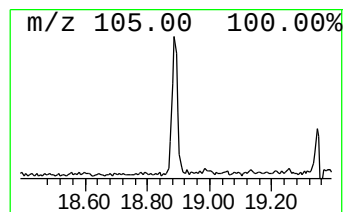
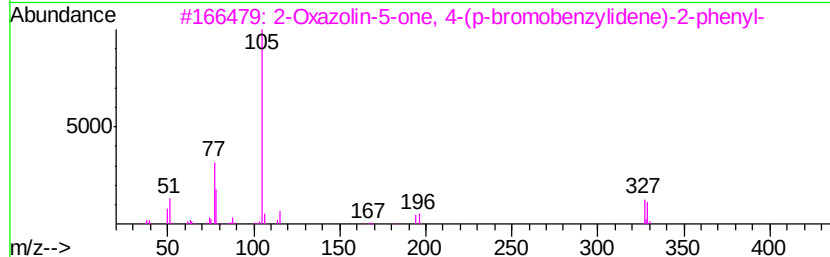
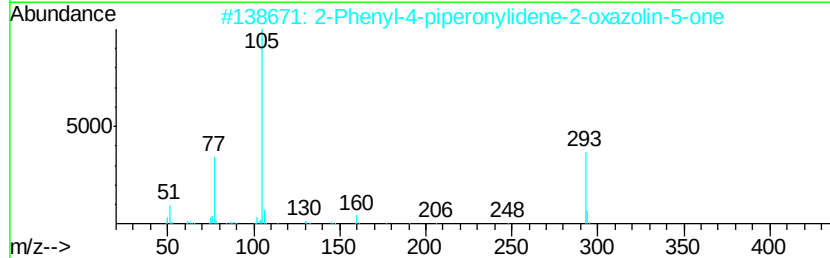
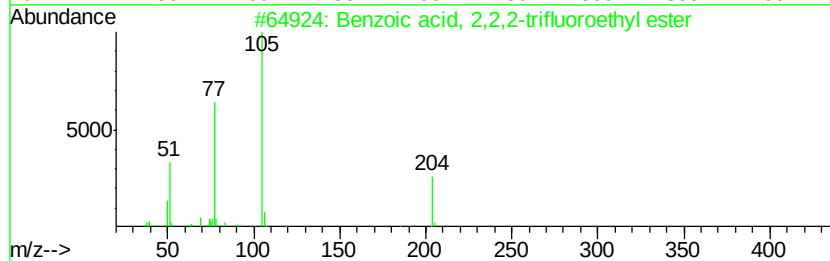
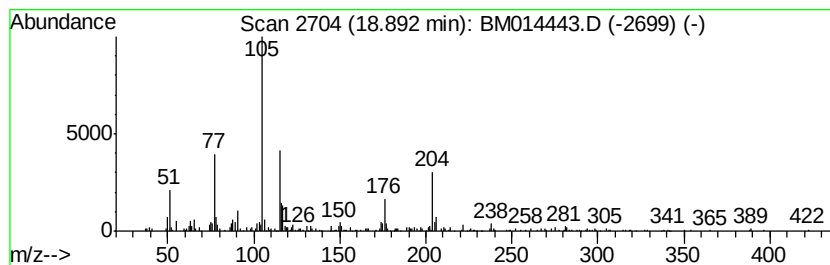
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzoic acid, 2,2,2-trifluo... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.89	3.78 ng/ul	348546	Phenanthrene-d10	16.94		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 2,2,2-trifluoroeth...	204	C9H7F3O2	1000376-54-4	53	
2	2-Phenyl-4-piperonylidene-2-oxaz...	293	C17H11N04	006412-89-1	46	
3	2-Oxazolin-5-one, 4-(p-bromobenz...	327	C16H10BrN02	020345-16-8	46	
4	Benzenepentanoic acid, .delta.-oxo-	192	C11H12O3	001501-05-9	43	
5	Tetrazole, 1-(1-phenylpyrazol-3-...	288	C16H12N6	1000259-06-8	43	



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030118\
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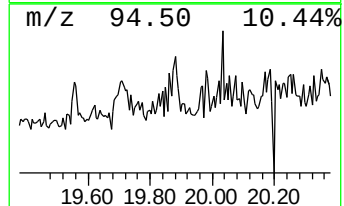
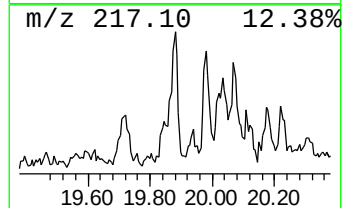
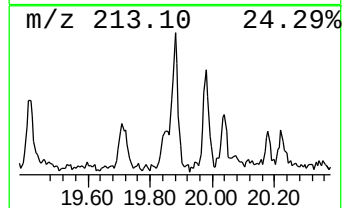
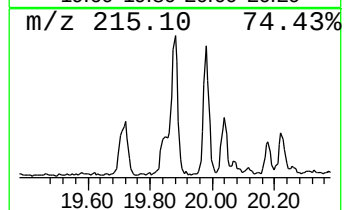
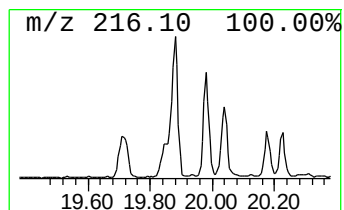
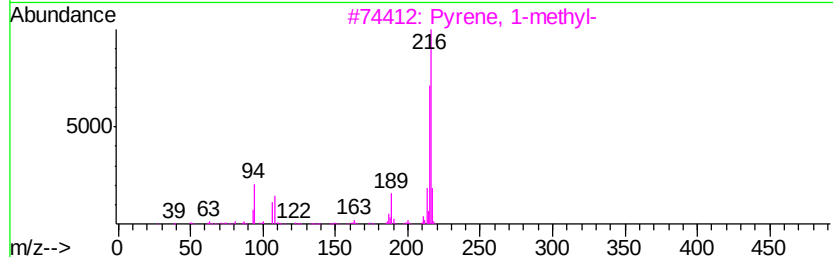
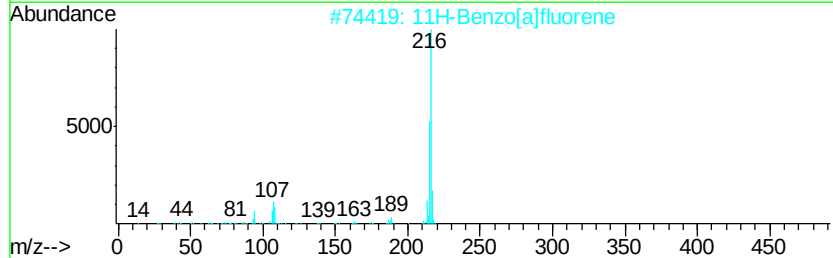
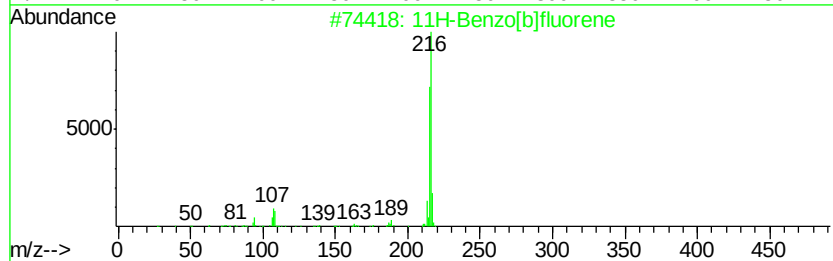
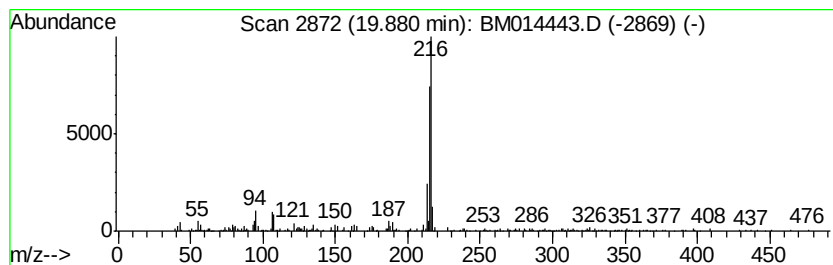
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 11H-Benzo[b]fluorene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.88	2.14 ng/ul	454762	Chrysene-d12	21.15
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		11H-Benzo[b]fluorene	216 C17H12	000243-17-4 83
2		11H-Benzo[a]fluorene	216 C17H12	000238-84-6 80
3		Pvrene, 1-methyl-	216 C17H12	002381-21-7 70
4		Fluoranthene, 2-methyl-	216 C17H12	033543-31-6 68
5		Fluoranthene, 2-methyl-	216 C17H12	033543-31-6 64



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030118\
Data File : BM014443.D
Acq On : 01 Mar 2018 21:22
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Misc :
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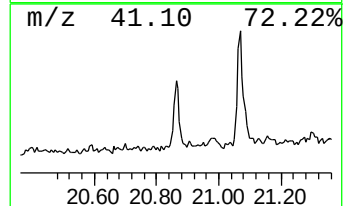
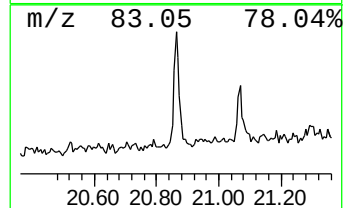
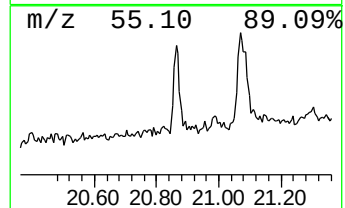
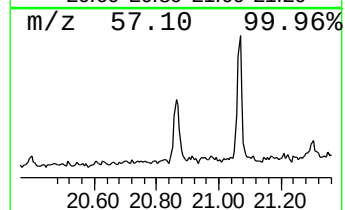
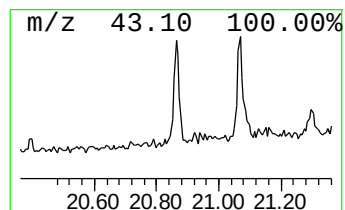
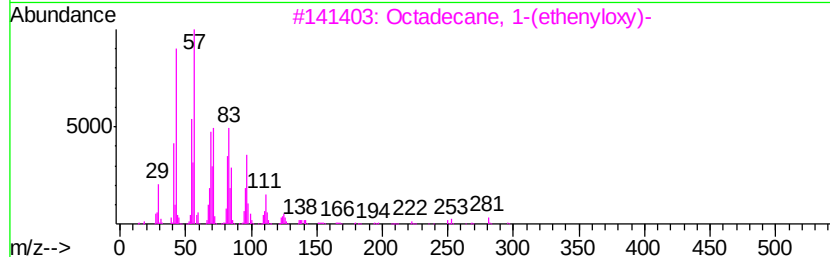
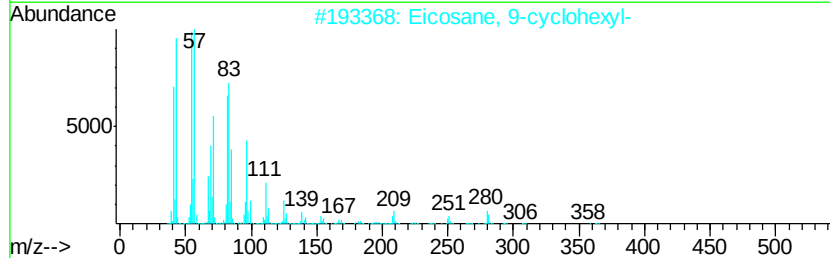
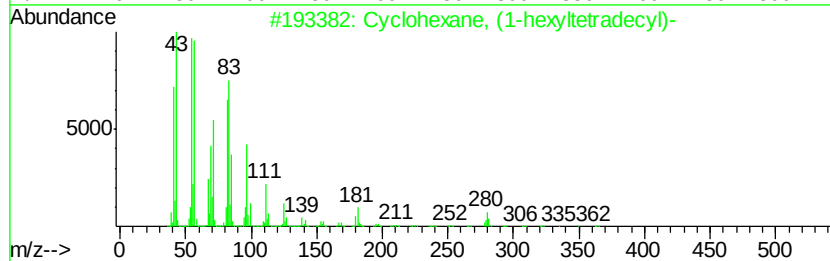
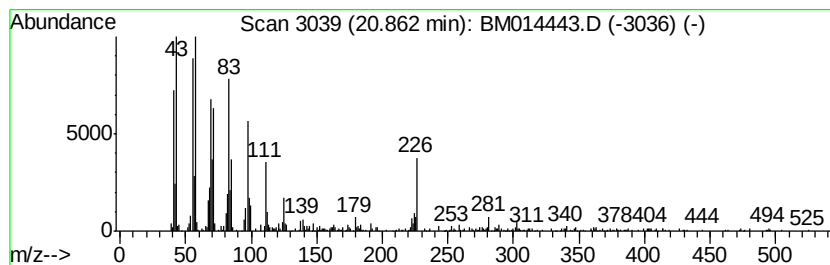
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.86	3.97 ng/ul	842915	Chrysene-d12	21.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (1-hexyltetradecyl)-	364	C26H52	004443-60-1	90
2			Eicosane, 9-cyclohexyl-	364	C26H52	004443-61-2	90
3			Octadecane, 1-(ethenvloxy)-	296	C20H40	000930-02-9	90
4			Z-5-Nonadecene	266	C19H38	1000131-11-8	89
5			9-Eicosene, (E)-	280	C20H40	074685-29-3	89



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030118\
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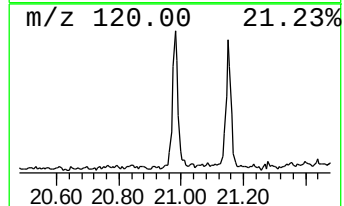
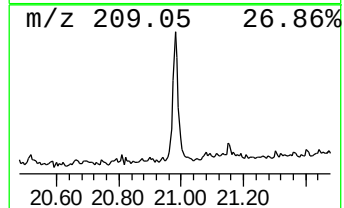
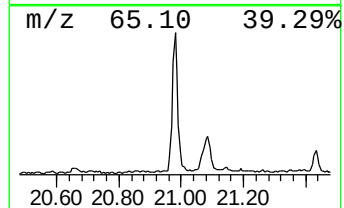
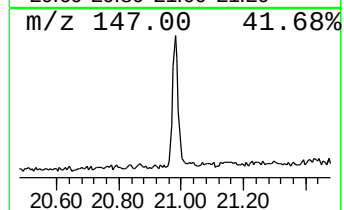
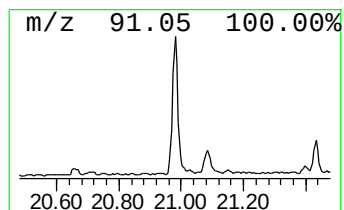
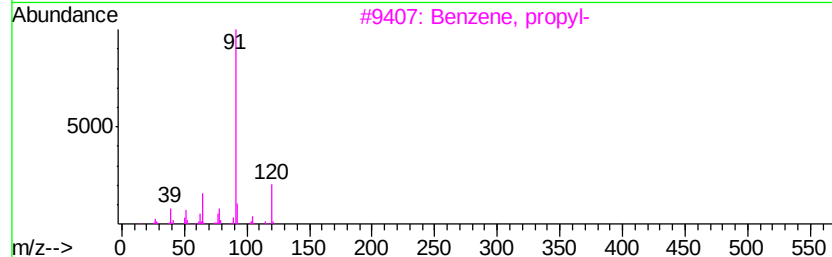
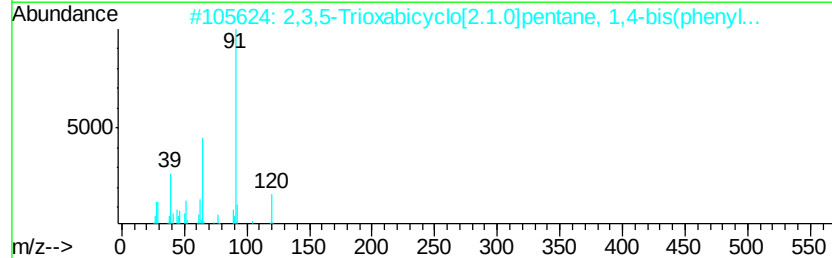
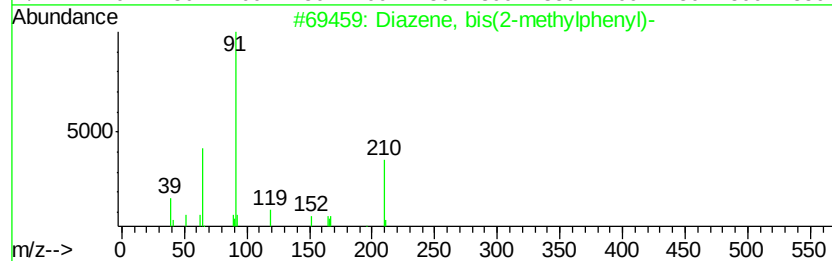
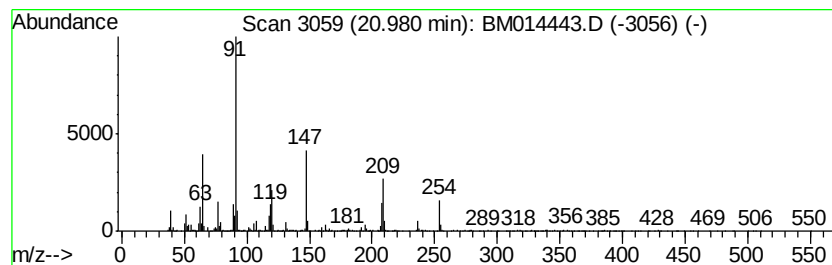
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Diazene, bis(2-methylphenyl)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.98	5.08 ng/ul	1077290	Chrysene-d12	21.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diazene, bis(2-methylphenyl)-	210	C14H14N2	000584-90-7	53
2		2,3,5-Trioxabicyclo[2.1.0]pentan...	254	C16H14O3	056247-48-4	50
3		Benzene, propyl-	120	C9H12	000103-65-1	49
4		Benzaldehyde, 4-methyl-	120	C8H8O	000104-87-0	47
5		Benzene, 1-methyl-4-nitroso-	121	C7H7NO	000623-11-0	43



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030118\
Data File : BM014443.D
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ClientSampleId :
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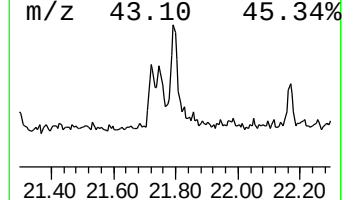
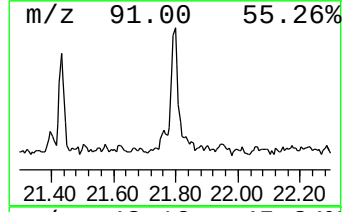
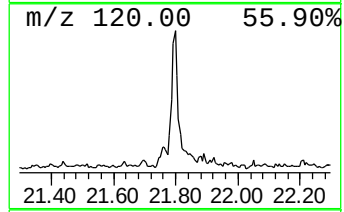
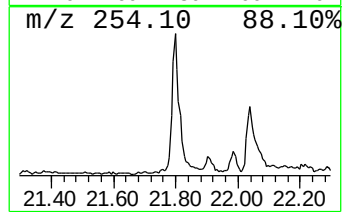
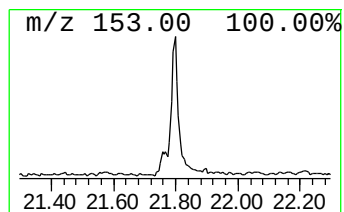
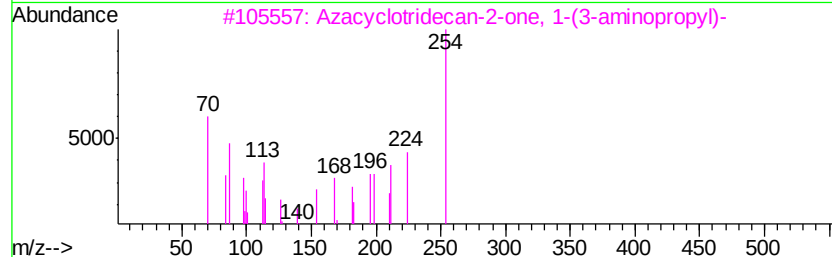
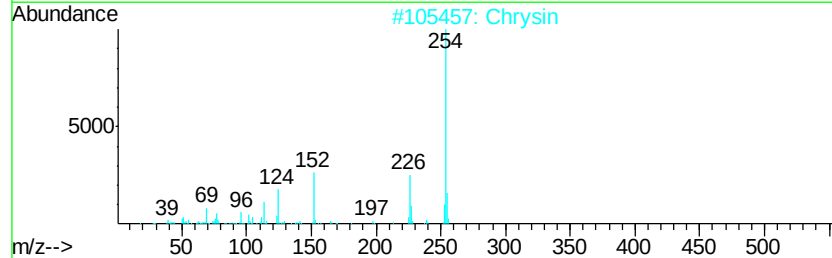
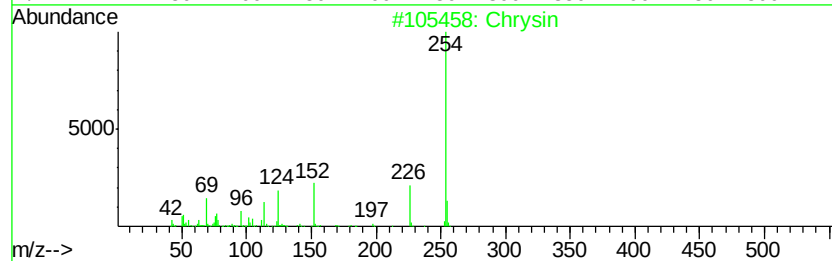
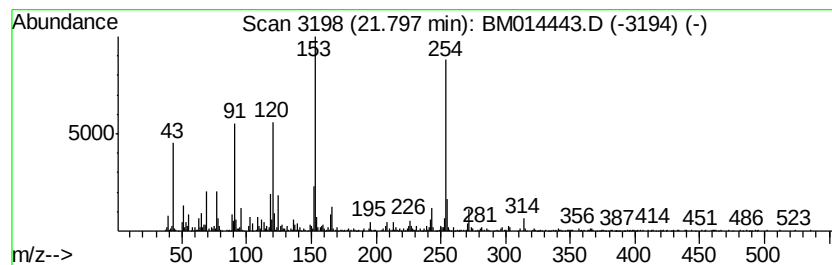
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-02 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.80	8.13 ng/ul	1724480	Chrysene-d12	21.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Chrysin	254	C15H10O4	000480-40-0	38
2		Chrysin	254	C15H10O4	000480-40-0	38
3		Azacyclotridecan-2-one, 1-(3-ami...	254	C15H30N2O	064414-61-5	35
4		3-Pyridinemethanol, 5-hydroxy-4,...	153	C8H11NO2	000061-67-6	27
5		Chrysin	254	C15H10O4	000480-40-0	22



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030118\
Data File : BM014443.D
Acq On : 01 Mar 2018 21:22
Operator : SJ/JU
Sample : J1678-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BE5T2

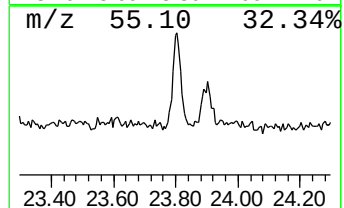
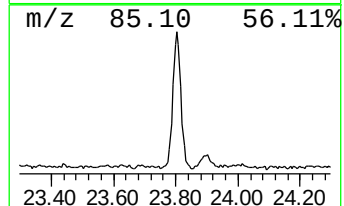
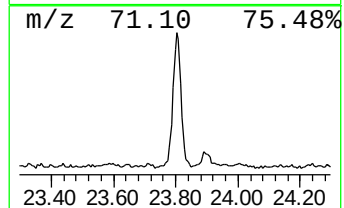
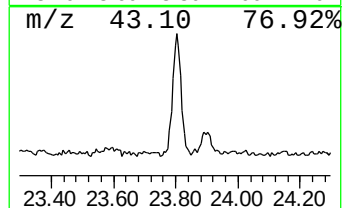
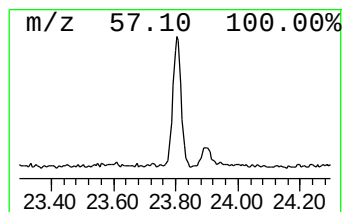
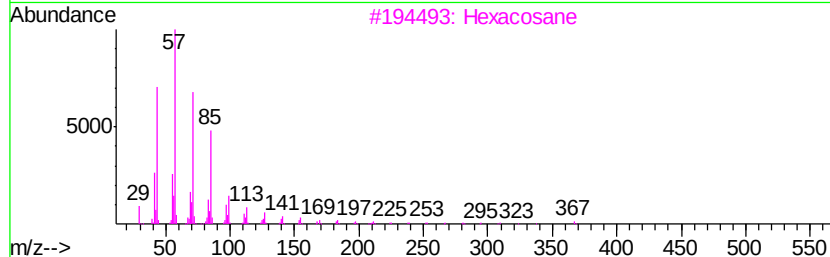
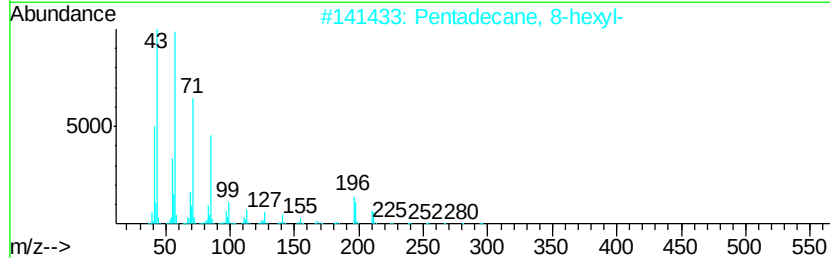
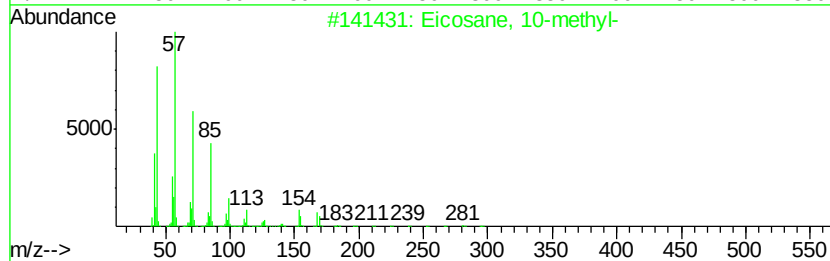
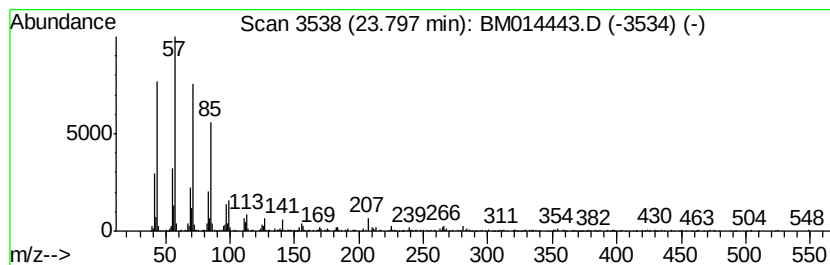
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Branched23.80 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.80	28.55 ng/ul	1777050	Perylene-d12	23.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93
3		Hexacosane	366	C26H54	000630-01-3	90
4		Heptadecane	240	C17H36	000629-78-7	87
5		Docosane	310	C22H46	000629-97-0	87



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030118\
 Data File : BM014443.D
 Acq On : 01 Mar 2018 21:22
 Operator : SJ/JU
 Sample : J1678-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BE5T2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	15.93	2.9	ng/ul	267208	4	16.94	1846150	20.0
(DEL) Alkane: Str...	16.69	3.5	ng/ul	321292	4	16.94	1846150	20.0
Anthracene, 2-met...	17.82	2.6	ng/ul	240077	4	16.94	1846150	20.0
Tetradecanoic acid	17.85	11.5	ng/ul	1064500	4	16.94	1846150	20.0
4H-Cyclopenta[def...	18.01	4.7	ng/ul	436793	4	16.94	1846150	20.0
9,10-Anthracenedione	18.38	3.0	ng/ul	276078	4	16.94	1846150	20.0
Benzoic acid, 2,2...	18.89	3.8	ng/ul	348546	4	16.94	1846150	20.0
11H-Benzo[b]fluorene	19.88	2.1	ng/ul	454762	5	21.15	4241340	20.0
(DEL) Alkane: Str...	20.86	4.0	ng/ul	842915	5	21.15	4241340	20.0
Diazene, bis(2-me...	20.98	5.1	ng/ul	1077290	5	21.15	4241340	20.0
unknown-02	21.80	8.1	ng/ul	1724480	5	21.15	4241340	20.0
(DEL) Alkane: Bra...	23.80	28.6	ng/ul	1777050	6	23.33	1244830	20.0