

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM081316\
 Data File : BM007101.D
 Acq On : 14 Aug 2016 11:01
 Operator : UM/SJ
 Sample : H4387-17
 Misc :
 ALS Vial : 66 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P003-SS004-1218-01

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\SOM02.2-EPA-BM081116.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	3.322	120	125	136	rVB	70847	99371	1.15%	0.091%
2	6.975	738	746	757	rBV	1274912	2174577	25.07%	1.987%
3	7.145	768	775	787	rVB	950871	1451828	16.74%	1.327%
4	7.345	801	809	819	rBV	1297083	2115224	24.39%	1.933%
5	7.810	880	888	896	rBV	1173647	1849760	21.33%	1.690%
6	8.504	998	1006	1016	rBV	1436557	2328334	26.84%	2.128%
7	8.963	1076	1084	1094	rBV	1222360	1992782	22.97%	1.821%
8	9.681	1198	1206	1214	rBV	1034381	1725402	19.89%	1.577%
9	10.210	1287	1296	1307	rBV	1597268	2709115	31.23%	2.476%
10	10.592	1354	1361	1368	rBV	1462713	2529798	29.17%	2.312%
11	10.728	1376	1384	1400	rBV	1252225	2230492	25.72%	2.038%
12	13.763	1891	1900	1904	rBV	64595	118219	1.36%	0.108%
13	13.851	1904	1915	1919	rVV	2668972	3916193	45.15%	3.579%
14	13.898	1919	1923	1932	rVB	2342043	3188346	36.76%	2.914%
15	14.133	1956	1963	1977	rVB	3048424	4857752	56.01%	4.439%
16	14.345	1994	1999	2004	rBV	75154	98596	1.14%	0.090%
17	14.439	2004	2015	2022	rBV2	2124516	3321057	38.29%	3.035%
18	14.627	2041	2047	2060	rBV	1253170	1958112	22.58%	1.789%
19	14.839	2079	2083	2089	rVB5	60963	106736	1.23%	0.098%
20	15.057	2112	2120	2124	rBV6	49283	126784	1.46%	0.116%
21	15.298	2157	2161	2167	rVB2	169599	236998	2.73%	0.217%
22	15.433	2175	2184	2190	rBV	4022940	5824431	67.15%	5.323%
23	15.545	2198	2203	2212	rVB	1026006	1411952	16.28%	1.290%
24	15.674	2220	2225	2228	rBV3	63165	112023	1.29%	0.102%
25	15.927	2260	2268	2277	rVB8	43454	116285	1.34%	0.106%
26	16.157	2302	2307	2315	rBV2	229999	393886	4.54%	0.360%
27	16.951	2438	2442	2447	rBV2	100852	140515	1.62%	0.128%
28	17.004	2447	2451	2459	rVB	111350	173803	2.00%	0.159%
29	17.180	2475	2481	2486	rBV2	2705545	4116886	47.46%	3.762%
30	17.221	2486	2488	2493	rVV	682068	899857	10.37%	0.822%
31	17.280	2493	2498	2508	rVV2	4297855	6334848	73.03%	5.789%
32	17.374	2508	2514	2519	rVB3	57954	117623	1.36%	0.107%
33	17.568	2543	2547	2558	rVB7	67857	130398	1.50%	0.119%
34	17.686	2563	2567	2573	rVV3	63628	116425	1.34%	0.106%

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35	17.762	2575	2580	2584	rVB	120789	169332	1.95%	0.155%
36	17.909	2600	2605	2609	rBV	74313	124901	1.44%	0.114%
37	18.080	2625	2634	2636	rVV2	573315	1353858	15.61%	1.237%
38	18.104	2636	2638	2644	rVV	526667	668861	7.71%	0.611%
39	18.186	2648	2652	2656	rVV2	106960	183655	2.12%	0.168%
40	18.245	2657	2662	2666	rVV3	410450	764734	8.82%	0.699%
41	18.286	2666	2669	2676	rVB	297185	421290	4.86%	0.385%
42	18.404	2686	2689	2693	rVV4	85805	138388	1.60%	0.126%
43	18.451	2693	2697	2702	rVB2	75978	126100	1.45%	0.115%
44	18.556	2711	2715	2718	rBV2	163423	225868	2.60%	0.206%
45	18.598	2718	2722	2730	rVB2	129858	249671	2.88%	0.228%
46	18.774	2749	2752	2754	rBV2	67045	88478	1.02%	0.081%
47	18.804	2754	2757	2761	rVV	144533	206077	2.38%	0.188%
48	18.856	2762	2766	2769	rVV	234185	336598	3.88%	0.308%
49	18.892	2769	2772	2776	rVB2	171393	226711	2.61%	0.207%
50	18.974	2781	2786	2791	rBV	500113	773674	8.92%	0.707%
51	19.033	2791	2796	2799	rVV3	198528	387532	4.47%	0.354%
52	19.068	2799	2802	2805	rVB	154895	190052	2.19%	0.174%
53	19.109	2805	2809	2816	rBV3	208616	430580	4.96%	0.393%
54	19.239	2826	2831	2837	rVB	1207422	1596552	18.41%	1.459%
55	19.304	2838	2842	2845	rBV2	111164	155609	1.79%	0.142%
56	19.356	2845	2851	2854	rBV3	188765	311002	3.59%	0.284%
57	19.392	2854	2857	2860	rVB	128052	144740	1.67%	0.132%
58	19.574	2882	2888	2892	rBV	5580276	8673756	100.00%	7.926%
59	19.680	2901	2906	2911	rBV2	274697	414162	4.77%	0.378%
60	19.839	2930	2933	2939	rVB4	138079	224668	2.59%	0.205%
61	19.921	2943	2947	2959	rBV5	238302	621309	7.16%	0.568%
62	20.009	2959	2962	2966	rBV4	81919	123583	1.42%	0.113%
63	20.062	2966	2971	2973	rVV4	198078	323188	3.73%	0.295%
64	20.092	2973	2976	2980	rVB	318880	473675	5.46%	0.433%
65	20.168	2986	2989	2991	rBV3	63637	89795	1.04%	0.082%
66	20.198	2991	2994	2998	rVB	154031	191113	2.20%	0.175%
67	20.256	2999	3004	3009	rVV	381095	547897	6.32%	0.501%
68	20.392	3024	3027	3031	rBV	324370	391241	4.51%	0.358%
69	20.439	3032	3035	3039	rVB	291140	362540	4.18%	0.331%
70	20.480	3039	3042	3050	rVB3	97986	154834	1.79%	0.141%
71	20.556	3052	3055	3059	rBV3	71591	104975	1.21%	0.096%

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72	20.845	3100	3104	3110	rVB	240461	385147	4.44%	0.352%
73	20.933	3116	3119	3123	rVB	112544	122249	1.41%	0.112%
74	21.009	3125	3132	3136	rVV3	264927	548410	6.32%	0.501%
75	21.050	3136	3139	3140	rVV	119047	122295	1.41%	0.112%
76	21.080	3140	3144	3150	rVB4	648197	872300	10.06%	0.797%
77	21.362	3185	3192	3196	rBV2	4631598	6936735	79.97%	6.339%
78	21.397	3196	3198	3208	rVB	879555	1107100	12.76%	1.012%
79	21.480	3209	3212	3215	rVB3	131135	152434	1.76%	0.139%
80	21.521	3216	3219	3223	rBV3	116800	187151	2.16%	0.171%
81	21.921	3284	3287	3291	rVB	294766	363815	4.19%	0.332%
82	21.974	3292	3296	3302	rBV2	386841	599238	6.91%	0.548%
83	22.433	3371	3374	3385	rVB7	202860	433376	5.00%	0.396%
84	22.950	3457	3462	3467	rBV	796814	1635642	18.86%	1.495%
85	22.991	3467	3469	3475	rVB	531322	768206	8.86%	0.702%
86	23.439	3541	3545	3548	rBV	343448	578517	6.67%	0.529%
87	23.491	3548	3554	3560	rBV2	3963974	7580979	87.40%	6.928%
88	23.639	3572	3579	3585	rBV	2746377	5433384	62.64%	4.965%
89	24.191	3669	3673	3680	rVB3	337413	606778	7.00%	0.554%

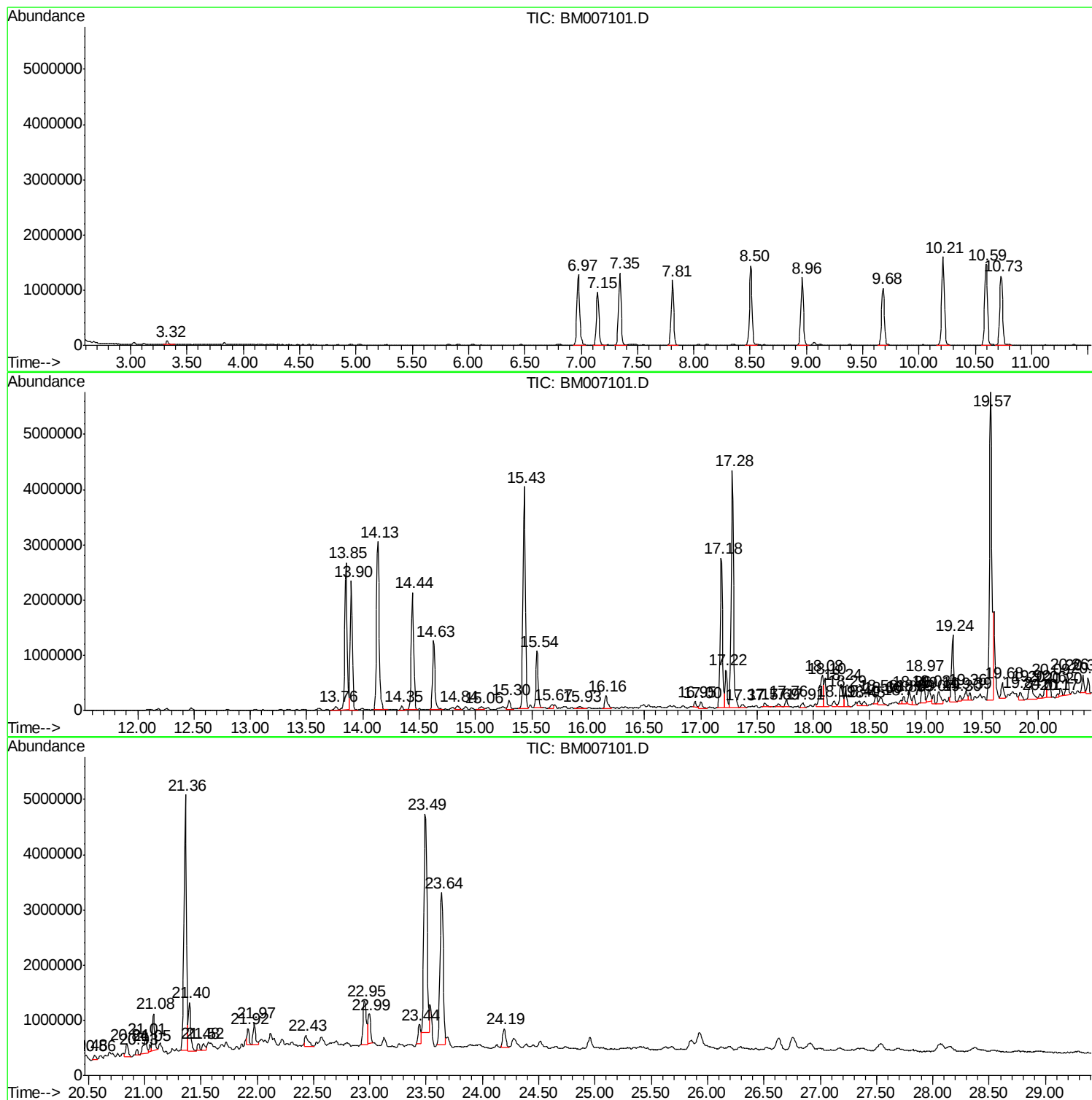
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM081116.M
Quant Title : SVOA CALIBRATION

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



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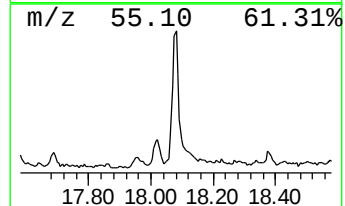
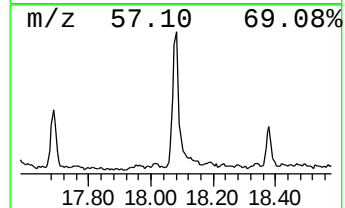
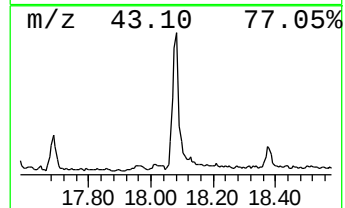
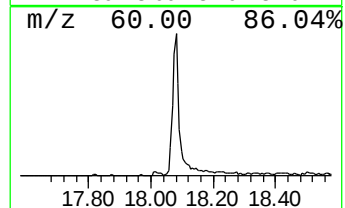
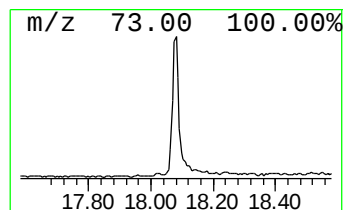
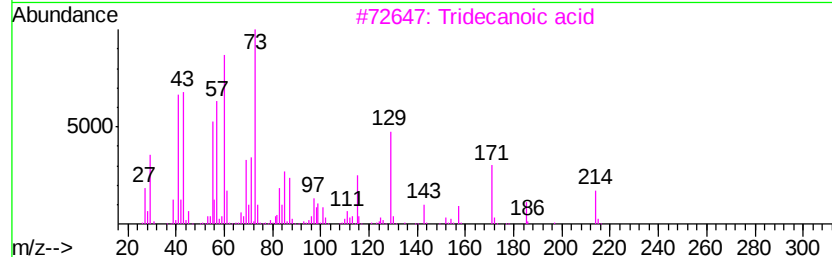
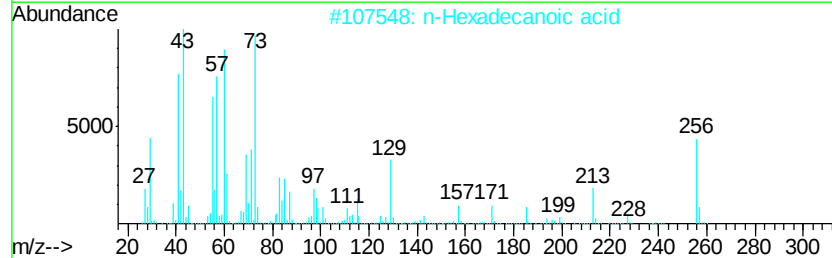
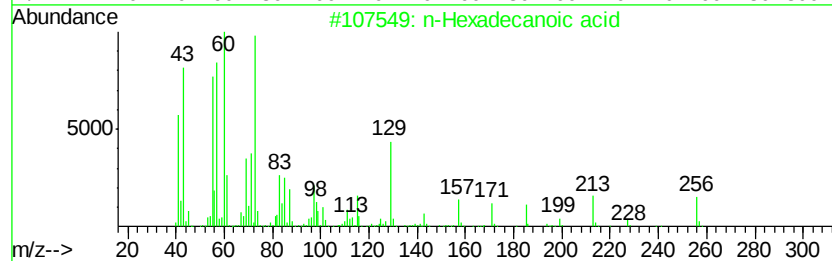
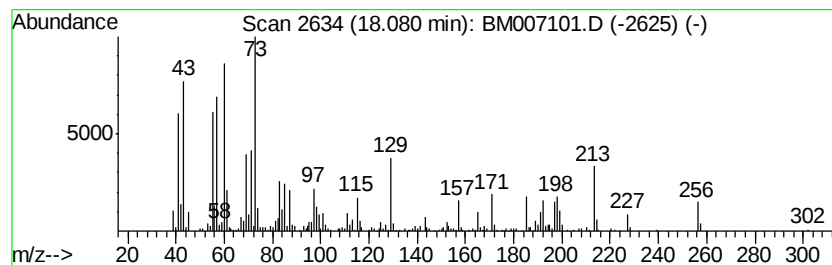
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM081116.M
Quant Title : SVOA CALIBRATION

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.		
18.08	6.58 ng/ul	1353860	Phenanthrene-d10	17.18		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99	
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98	
3	Tridecanoic acid	214	C13H26O2	000638-53-9	95	
4	Tridecanoic acid	214	C13H26O2	000638-53-9	94	
5	Tridecanoic acid	214	C13H26O2	000638-53-9	89	



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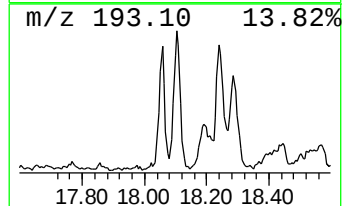
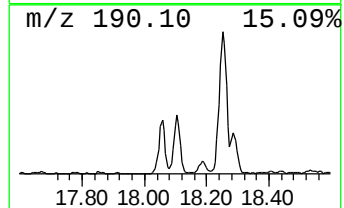
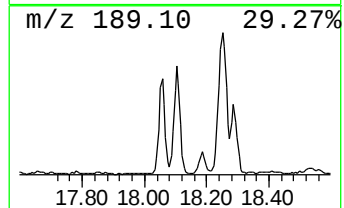
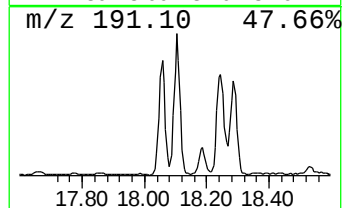
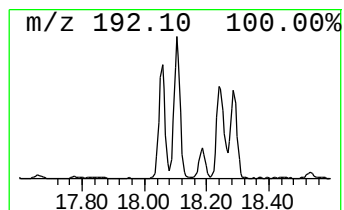
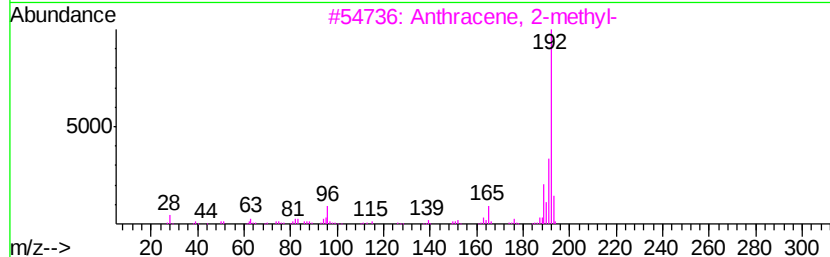
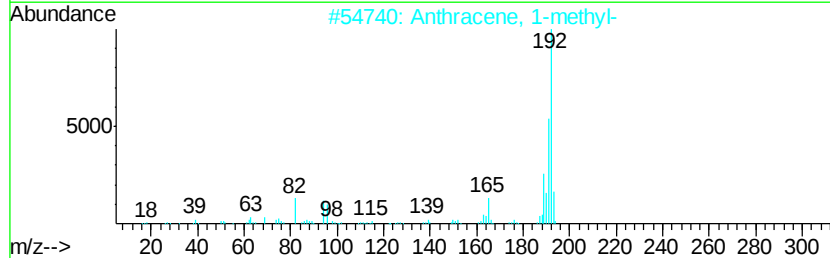
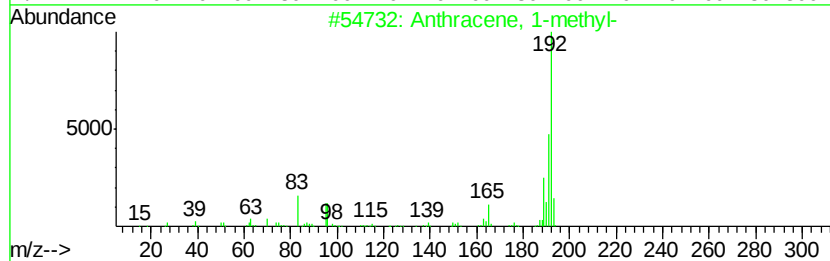
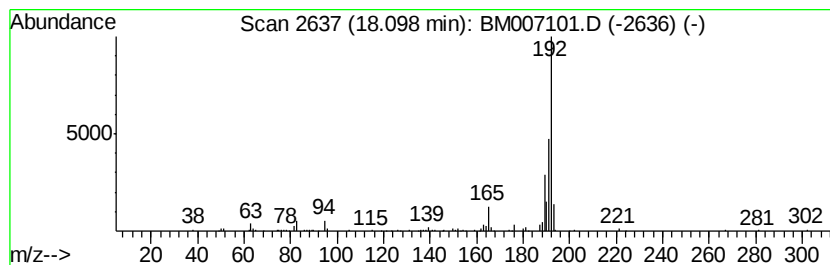
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Anthracene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.10	3.25 ng/ul	668861	Phenanthrene-d10	17.18		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
2		Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	89
4		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	87
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	87



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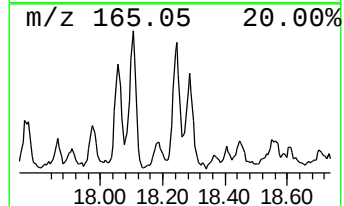
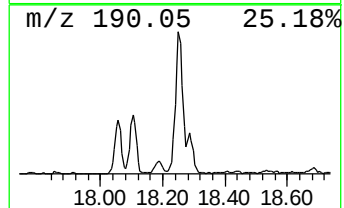
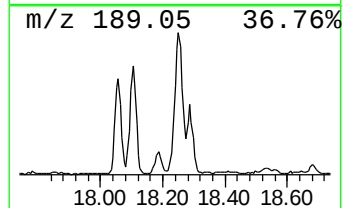
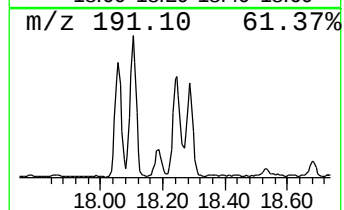
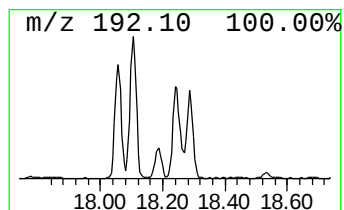
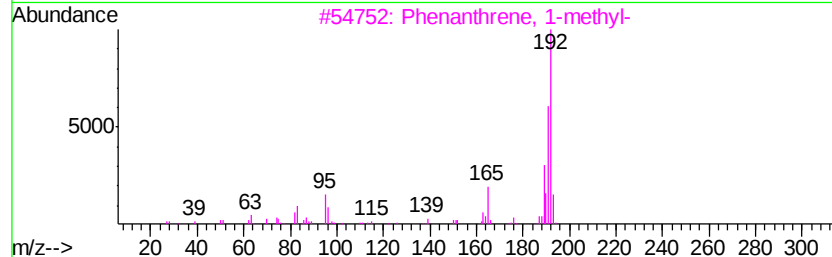
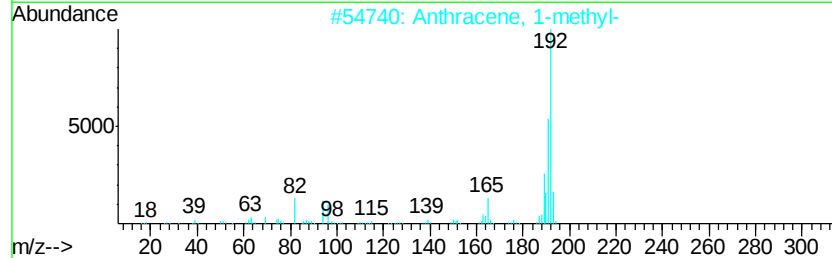
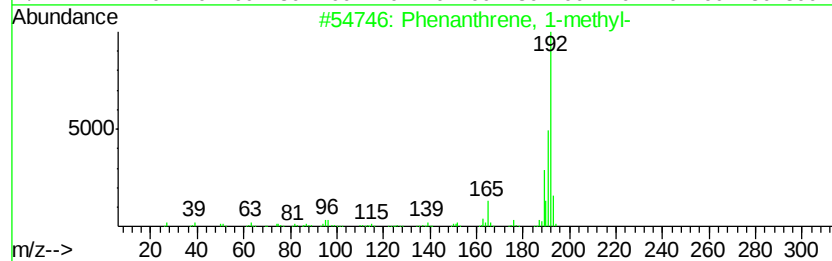
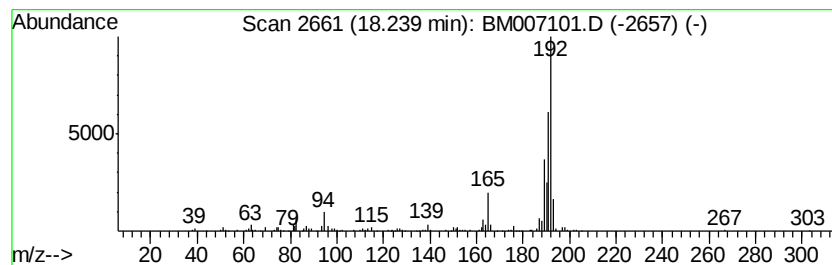
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Phenanthrene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.24	3.72 ng/ul	764734	Phenanthrene-d10	17.18		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
2		Anthracene, 1-methyl-	192	C15H12	000610-48-0	78
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	78
4		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	64
5		Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	64



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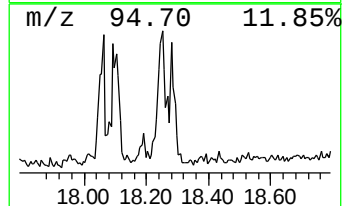
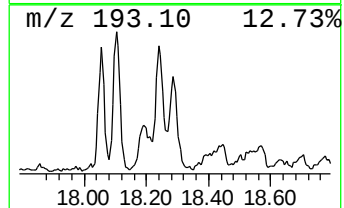
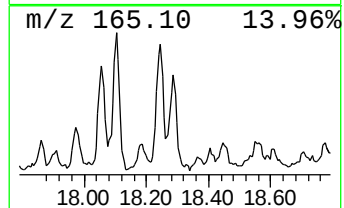
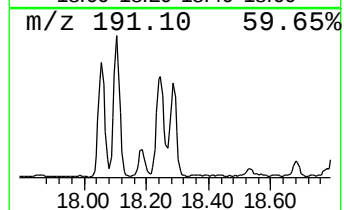
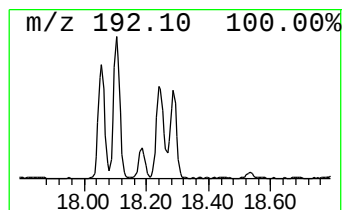
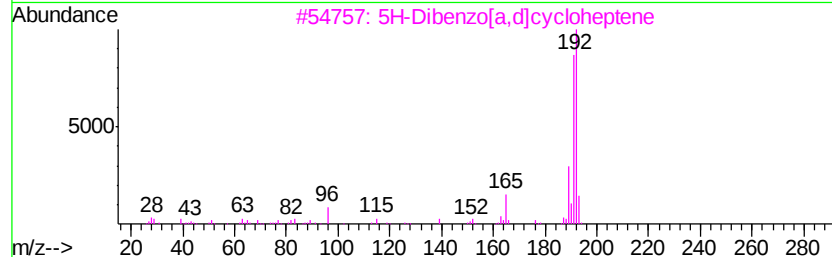
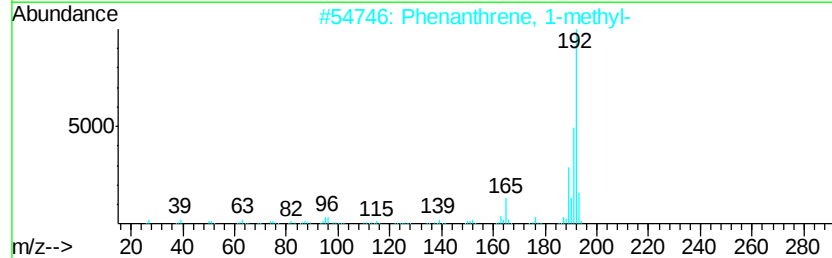
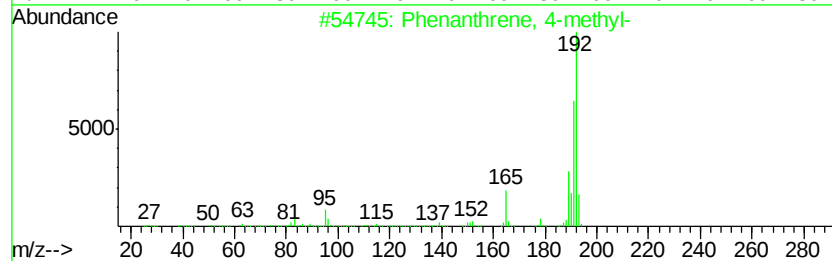
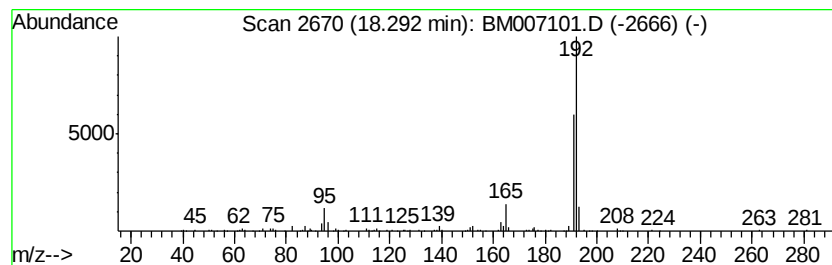
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenanthrene, 4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.29	2.05 ng/ul	421290	Phenanthrene-d10	17.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	90
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	87
3		5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	81
4		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	74
5		Propyne, 1,3-diphenyl-	192	C15H12	004980-70-5	74



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM081316\
Data File : BM007101.D
Acq On : 14 Aug 2016 11:01
Operator : UM/SJ
Sample : H4387-17
Misc :
ALS Vial : 66 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
P003-SS004-1218-01

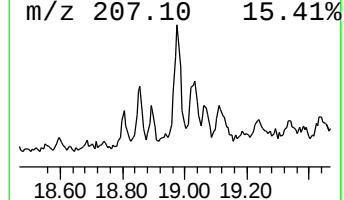
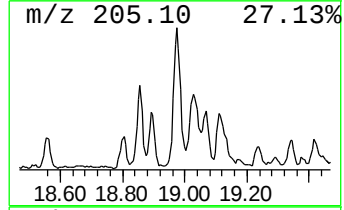
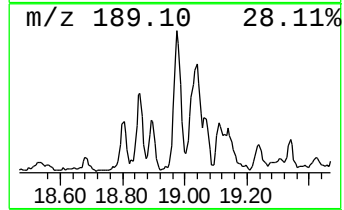
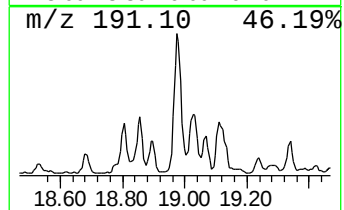
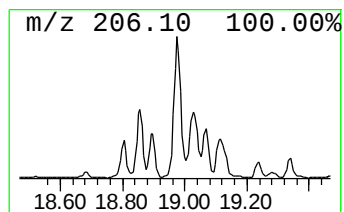
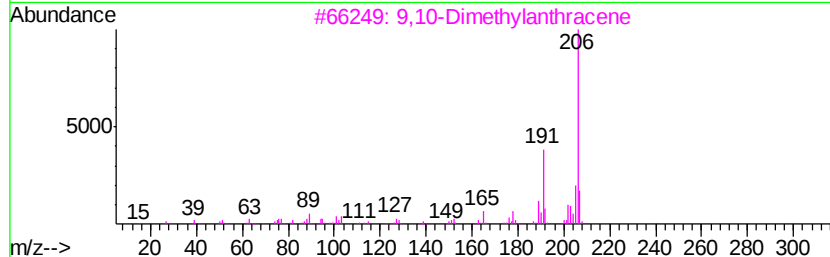
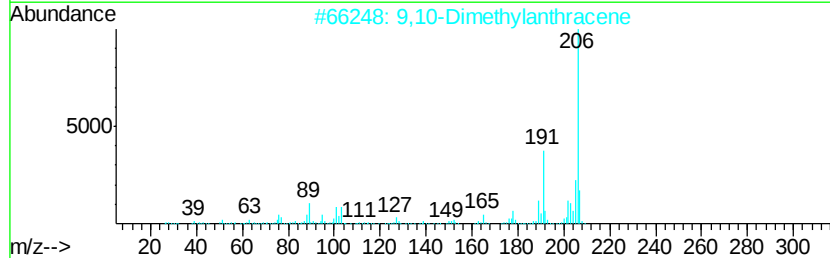
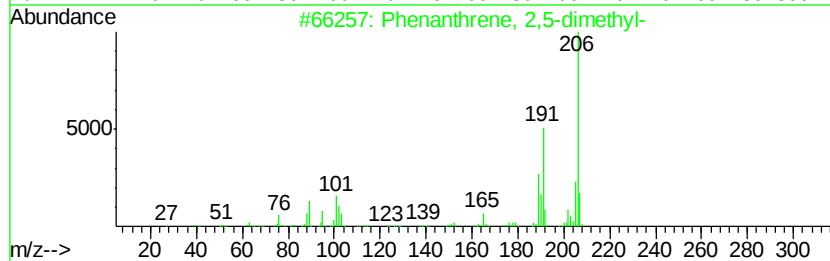
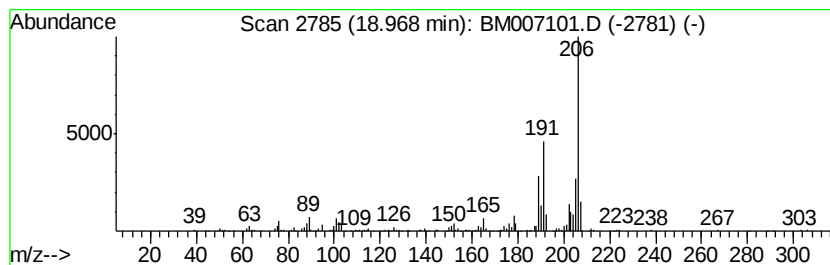
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Phenanthrene, 2,5-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.97	3.76 ng/ul	773674	Phenanthrene-d10	17.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2		9,10-Dimethylantracene	206	C16H14	000781-43-1	91
3		9,10-Dimethylantracene	206	C16H14	000781-43-1	91
4		Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	90
5		di-p-Tolylacetylene	206	C16H14	002789-88-0	87



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM081316\
Data File : BM007101.D
Acq On : 14 Aug 2016 11:01
Operator : UM/SJ
Sample : H4387-17
Misc :
ALS Vial : 66 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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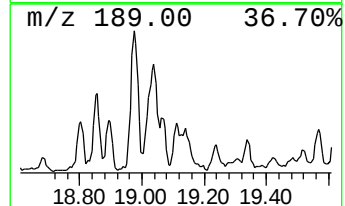
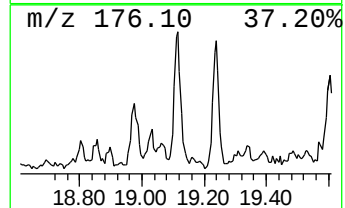
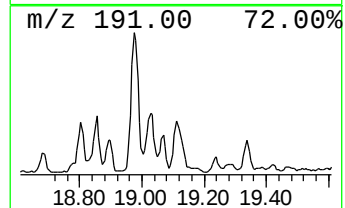
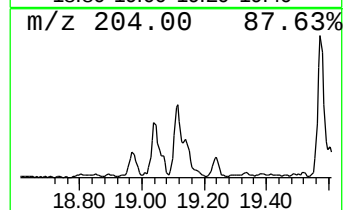
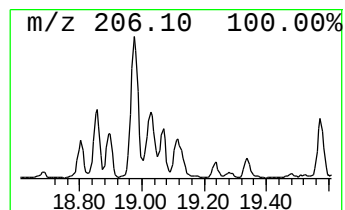
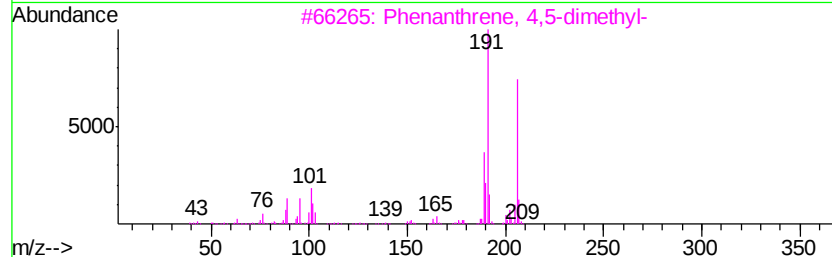
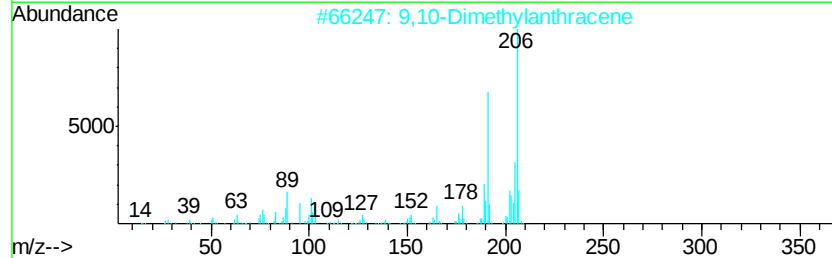
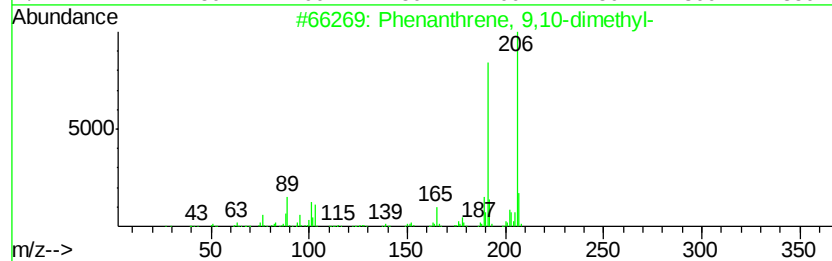
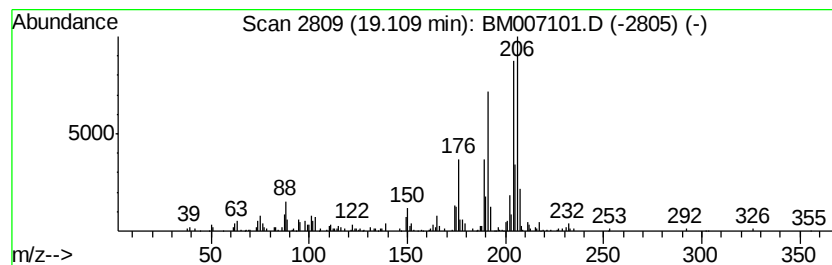
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Phenanthrene, 9,10-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.11	2.09 ng/ul	430580	Phenanthrene-d10	17.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 9,10-dimethyl-	206	C16H14	000604-83-1	70
2		9,10-Dimethylantracene	206	C16H14	000781-43-1	64
3		Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	60
4		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	60
5		Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	60



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM081316\
Data File : BM007101.D
Acq On : 14 Aug 2016 11:01
Operator : UM/SJ
Sample : H4387-17
Misc :
ALS Vial : 66 Sample Multiplier: 1

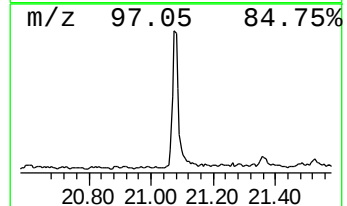
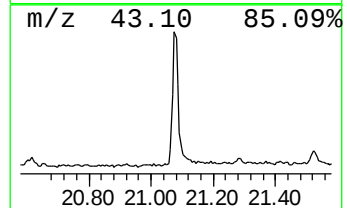
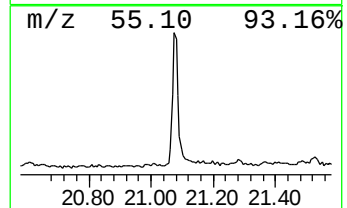
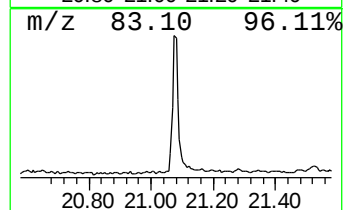
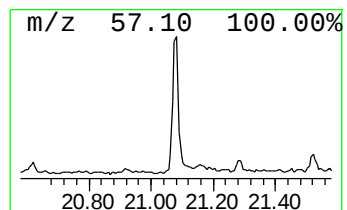
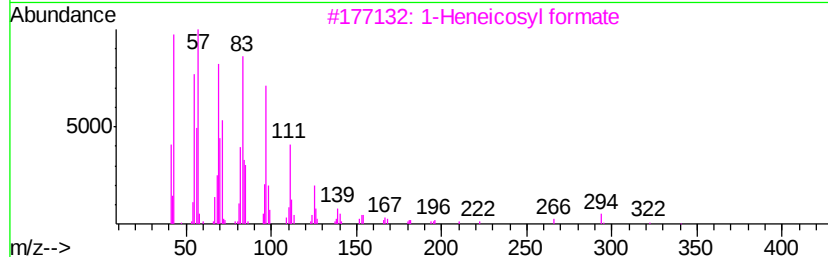
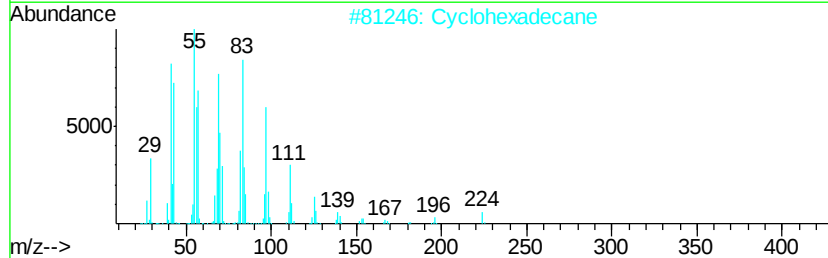
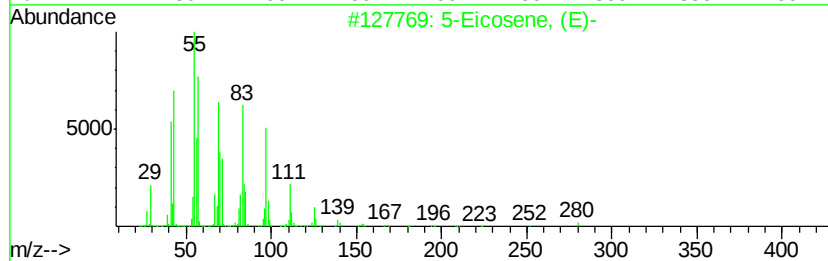
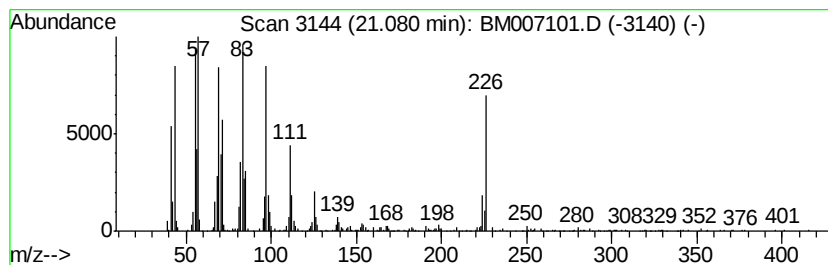
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Cyclic21.08 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.08	2.52 ng/ul	872300	Chrysene-d12	21.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-	280	C20H40	074685-30-6	98
2	Cyclohexadecane	224	C16H32	000295-65-8	95
3	1-Heneicosyl formate	340	C22H44O2	077899-03-7	91
4	Pentafluoropropionic acid, tetra...	360	C17H29F5O2	006222-06-6	90
5	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	90



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM081316\
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Misc :
ALS Vial : 66 Sample Multiplier: 1

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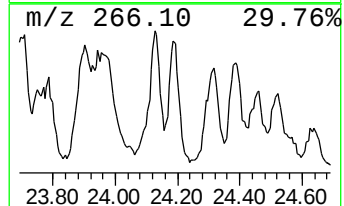
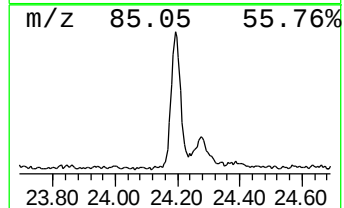
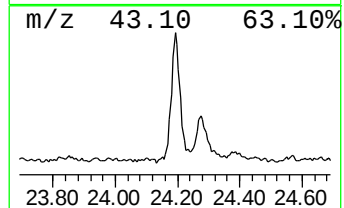
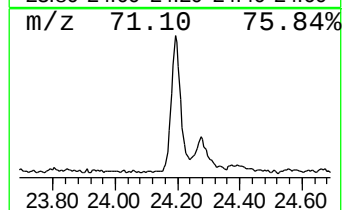
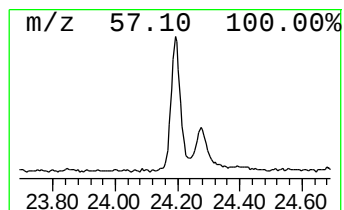
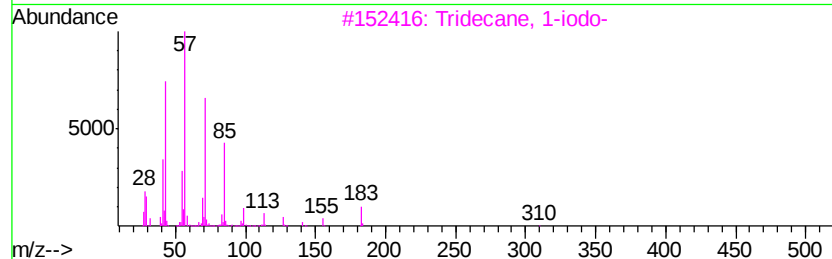
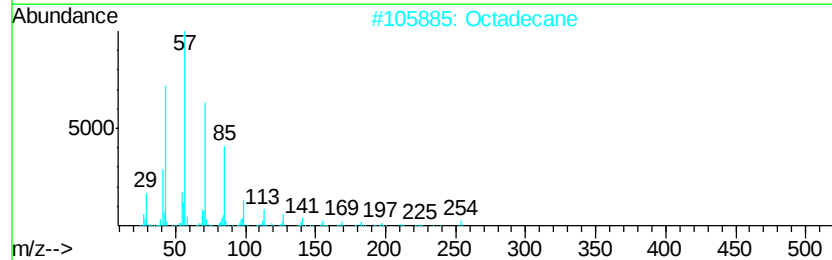
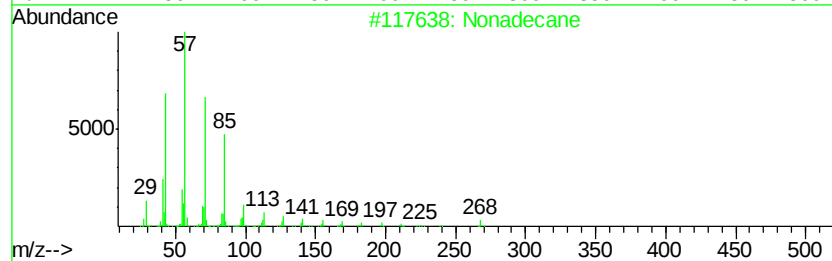
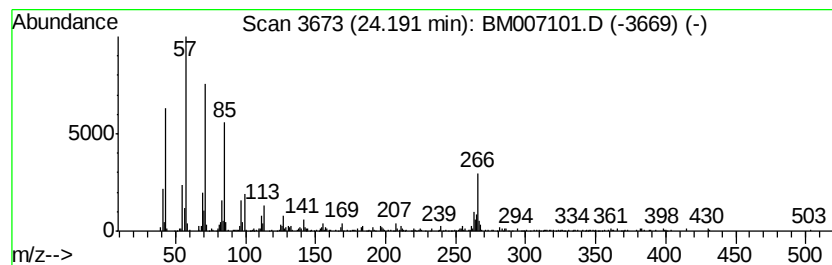
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Quant Title : SVOA CALIBRATION

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.
24.19	2.23 ng/ul	606778	Perylene-d12	23.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	96
2		Octadecane	254	C18H38	000593-45-3	92
3		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	70
4		triacontane	422	C30H62	000638-68-6	70
5		Hexacosane	366	C26H54	000630-01-3	70



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM081316\
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Acq On : 14 Aug 2016 11:01
Operator : UM/SJ
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Misc :
ALS Vial : 66 Sample Multiplier: 1

Instrument :
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ClientSampleId :
P003-SS004-1218-01

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM081116.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
n-Hexadecanoic acid	18.08	6.6	ng/ul	1353860	4	17.18	4116890	20.0
Anthracene, 1-met...	18.10	3.3	ng/ul	668861	4	17.18	4116890	20.0
Phenanthrene, 1-m...	18.24	3.7	ng/ul	764734	4	17.18	4116890	20.0
Phenanthrene, 4-m...	18.29	2.0	ng/ul	421290	4	17.18	4116890	20.0
Phenanthrene, 2,5...	18.97	3.8	ng/ul	773674	4	17.18	4116890	20.0
Phenanthrene, 9,1...	19.11	2.1	ng/ul	430580	4	17.18	4116890	20.0
(DEL) Alkane: Cyc...	21.08	2.5	ng/ul	872300	5	21.36	6936740	20.0
(DEL) Alkane: Str...	24.19	2.2	ng/ul	606778	6	23.64	5433380	20.0