

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101415\
 Data File : BG019197.D
 Acq On : 14 Oct 2015 2:36
 Operator : UM/NP
 Sample : G3996-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 E5LG3

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 0.1 % 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_G\METHODS\SOM02.2-EPA-BG100415.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	3.143	48	52	60	rBV	51599	79121	18.40%	1.432%
2	3.220	60	65	80	rVB	295929	413044	96.06%	7.477%
3	3.995	194	197	202	rVB4	2712	3770	0.88%	0.068%
4	4.224	233	236	244	rVB3	5024	8496	1.98%	0.154%
5	5.141	386	392	396	rBV2	6707	10536	2.45%	0.191%
6	7.191	735	741	748	rBV	44788	74930	17.43%	1.356%
7	7.356	763	769	777	rVB	29819	49036	11.40%	0.888%
8	7.562	799	804	813	rVB	39016	65701	15.28%	1.189%
9	8.032	878	884	892	rVB	67566	108367	25.20%	1.962%
10	8.737	998	1004	1011	rBV	56920	95008	22.10%	1.720%
11	9.124	1065	1070	1075	rVV2	5634	7981	1.86%	0.144%
12	9.189	1075	1081	1088	rVB	45642	77971	18.13%	1.412%
13	9.912	1197	1204	1212	rBV	38601	67267	15.64%	1.218%
14	10.452	1290	1296	1304	rBV	55121	95566	22.23%	1.730%
15	10.734	1341	1344	1350	rBV3	3817	6449	1.50%	0.117%
16	10.834	1353	1361	1367	rVV	102389	180490	41.98%	3.267%
17	10.969	1378	1384	1390	rBV2	7381	13861	3.22%	0.251%
18	11.245	1427	1431	1439	rVB4	1938	3639	0.85%	0.066%
19	11.998	1555	1559	1567	rVB2	3529	5642	1.31%	0.102%
20	12.250	1594	1602	1608	rBV5	2537	4032	0.94%	0.073%
21	12.485	1639	1642	1648	rVB3	2865	3659	0.85%	0.066%
22	12.544	1648	1652	1657	rBV4	4427	7730	1.80%	0.140%
23	12.661	1668	1672	1676	rVB3	2663	4510	1.05%	0.082%
24	12.703	1676	1679	1686	rVB6	2650	4397	1.02%	0.080%
25	13.020	1727	1733	1738	rBV4	2144	3471	0.81%	0.063%
26	13.149	1752	1755	1759	rVB2	2562	3157	0.73%	0.057%
27	13.267	1771	1775	1779	rVB3	3939	5868	1.36%	0.106%
28	13.396	1793	1797	1800	rBV3	2214	3431	0.80%	0.062%
29	13.478	1809	1811	1817	rVB2	6476	8919	2.07%	0.161%
30	13.549	1817	1823	1827	rBV4	3058	6559	1.53%	0.119%
31	13.825	1865	1870	1875	rBV6	2416	5096	1.19%	0.092%
32	13.983	1891	1897	1900	rBV2	4258	6593	1.53%	0.119%
33	14.060	1902	1910	1914	rVV	119640	175915	40.91%	3.185%
34	14.107	1914	1918	1924	rVV	61175	88998	20.70%	1.611%

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35	14.213	1932	1936	1940	rBV4	2680	3964	0.92%	0.072%
36	14.348	1953	1959	1970	rVB	115475	186999	43.49%	3.385%
37	14.471	1978	1980	1984	rBV4	3380	3969	0.92%	0.072%
38	14.553	1989	1994	2004	rVB	13994	22773	5.30%	0.412%
39	14.659	2004	2012	2020	rBV2	169860	286742	66.69%	5.191%
40	14.853	2039	2045	2057	rVB2	46664	81096	18.86%	1.468%
41	15.053	2074	2079	2083	rBV4	3620	5688	1.32%	0.103%
42	15.129	2087	2092	2095	rBV3	2400	3919	0.91%	0.071%
43	15.446	2138	2146	2147	rBV6	7023	14803	3.44%	0.268%
44	15.499	2152	2155	2160	rVB	24305	28229	6.57%	0.511%
45	15.546	2160	2163	2168	rVB5	1982	3213	0.75%	0.058%
46	15.646	2174	2180	2186	rBV	163338	245641	57.13%	4.447%
47	15.758	2194	2199	2208	rBV	44192	70838	16.47%	1.282%
48	15.916	2223	2226	2230	rVB3	3416	5146	1.20%	0.093%
49	15.975	2234	2236	2240	rBV3	2560	3173	0.74%	0.057%
50	16.193	2270	2273	2278	rBV5	1684	2912	0.68%	0.053%
51	16.304	2290	2292	2298	rVB5	5927	9075	2.11%	0.164%
52	16.363	2298	2302	2311	rBV	34115	48899	11.37%	0.885%
53	16.475	2318	2321	2324	rBV4	3265	3948	0.92%	0.071%
54	16.539	2329	2332	2337	rVB4	5285	7290	1.70%	0.132%
55	16.586	2337	2340	2346	rBV3	3835	5590	1.30%	0.101%
56	16.686	2352	2357	2366	rVB4	8642	17489	4.07%	0.317%
57	16.945	2395	2401	2404	rBV6	3838	6112	1.42%	0.111%
58	17.109	2426	2429	2431	rBV4	7006	8289	1.93%	0.150%
59	17.156	2433	2437	2440	rVB	34131	38391	8.93%	0.695%
60	17.250	2449	2453	2457	rVB3	4638	7070	1.64%	0.128%
61	17.403	2472	2479	2490	rBV2	232480	352221	81.91%	6.376%
62	17.503	2490	2496	2503	rVV	174216	269745	62.73%	4.883%
63	17.779	2537	2543	2549	rBV3	24214	38630	8.98%	0.699%
64	17.855	2552	2556	2558	rVB4	6727	7278	1.69%	0.132%
65	17.897	2558	2563	2567	rBV	42663	50129	11.66%	0.908%
66	18.038	2585	2587	2590	rBV4	2363	3270	0.76%	0.059%
67	18.296	2624	2631	2635	rBV7	6637	14348	3.34%	0.260%
68	18.443	2653	2656	2660	rVV3	7770	10620	2.47%	0.192%
69	18.549	2671	2674	2677	rBV5	5709	5961	1.39%	0.108%
70	18.590	2677	2681	2684	rBV	37615	45418	10.56%	0.822%
71	18.713	2699	2702	2706	rVB2	11074	12411	2.89%	0.225%

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72	18.819	2717	2720	2725	rVB6	3528	4239	0.99%	0.077%
73	18.907	2733	2735	2739	rVB5	3167	3987	0.93%	0.072%
74	18.983	2746	2748	2750	rBV2	3910	4236	0.99%	0.077%
75	19.148	2772	2776	2779	rBV2	18545	26972	6.27%	0.488%
76	19.183	2779	2782	2784	rVV3	10881	14934	3.47%	0.270%
77	19.213	2784	2787	2794	rVB	43812	59313	13.79%	1.074%
78	19.342	2806	2809	2813	rBV5	3138	4609	1.07%	0.083%
79	19.671	2862	2865	2868	rBV4	3057	4231	0.98%	0.077%
80	19.706	2869	2871	2874	rVB4	5241	4554	1.06%	0.082%
81	19.788	2874	2885	2893	rBV	236218	364593	84.79%	6.600%
82	19.888	2898	2902	2905	rVB6	4948	7140	1.66%	0.129%
83	19.912	2905	2906	2911	rVB4	4211	5249	1.22%	0.095%
84	19.947	2911	2912	2914	rBV2	3717	3083	0.72%	0.056%
85	20.294	2969	2971	2973	rBV2	5936	5767	1.34%	0.104%
86	20.323	2973	2976	2981	rVB	26437	29073	6.76%	0.526%
87	20.799	3053	3057	3058	rBV2	28114	31889	7.42%	0.577%
88	21.316	3141	3145	3153	rBV2	45006	77507	18.03%	1.403%
89	21.674	3200	3206	3217	rVB	250840	429984	100.00%	7.784%
90	21.857	3234	3237	3238	rBV3	12570	13053	3.04%	0.236%
91	21.921	3245	3248	3254	rVB6	9890	15227	3.54%	0.276%
92	23.161	3455	3459	3460	rBV3	7078	9513	2.21%	0.172%
93	23.249	3470	3474	3480	rVB6	11398	21560	5.01%	0.390%
94	23.372	3492	3495	3501	rVB2	10778	17467	4.06%	0.316%
95	23.842	3571	3575	3576	rBV4	2974	3527	0.82%	0.064%
96	23.989	3597	3600	3601	rBV3	5540	5653	1.31%	0.102%
97	24.677	3708	3717	3727	rVB	106037	284191	66.09%	5.145%
98	24.900	3745	3755	3771	rBV	144054	419616	97.59%	7.596%
99	25.147	3794	3797	3799	rBV3	2616	3535	0.82%	0.064%
100	25.564	3862	3868	3875	rVB10	7895	18624	4.33%	0.337%

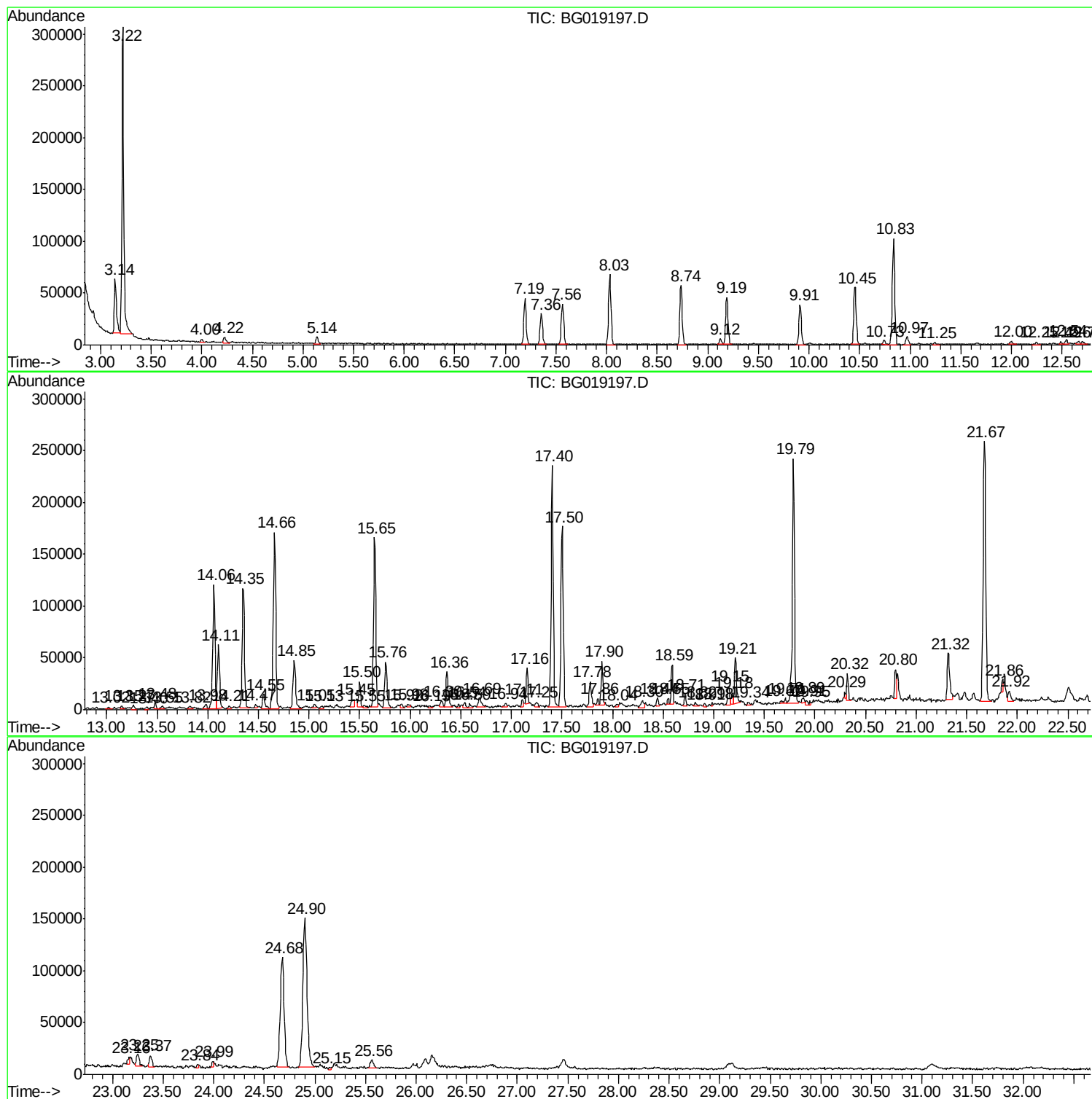
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG100415.M
Quant Title : SVOA CALIBRATION

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



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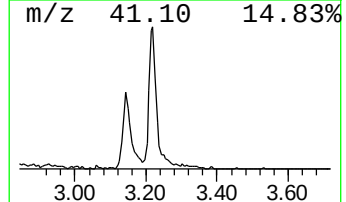
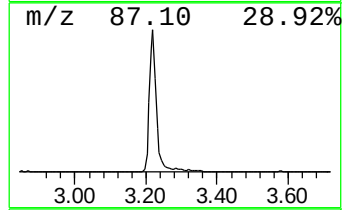
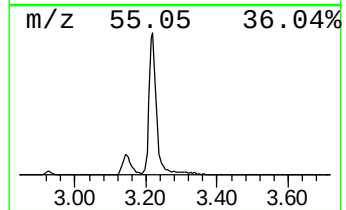
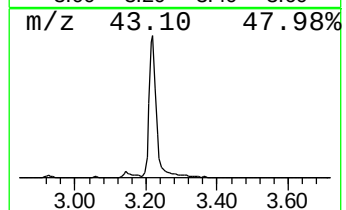
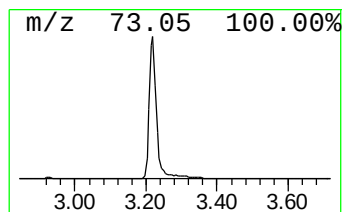
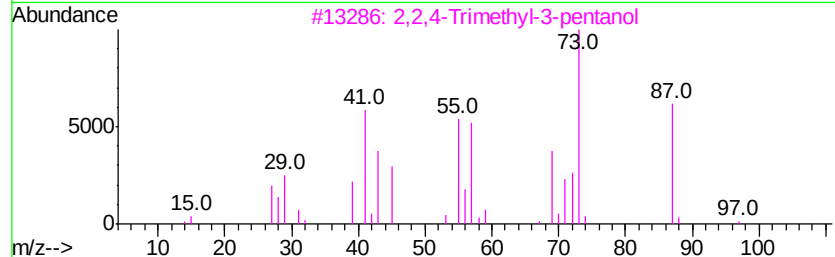
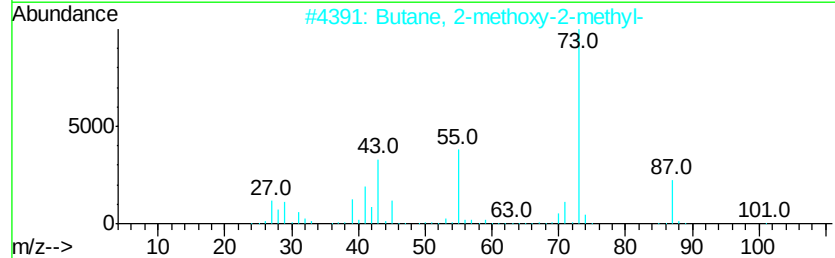
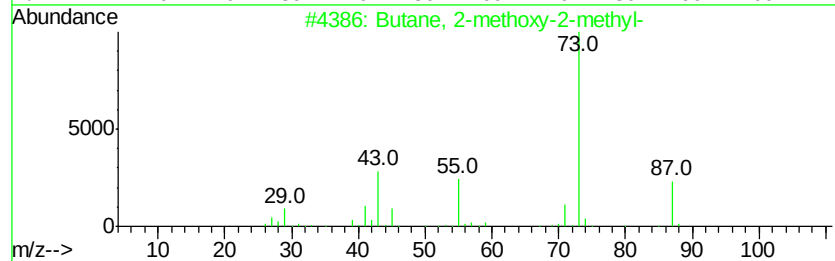
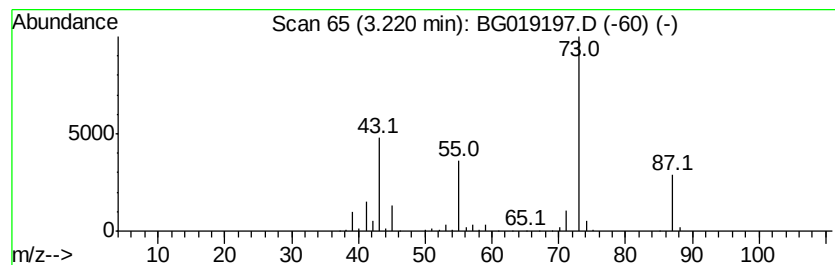
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG100415.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.22	76.23 ng/ul	413044	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	28
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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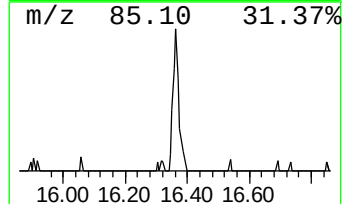
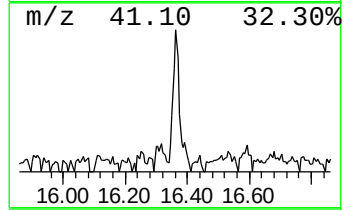
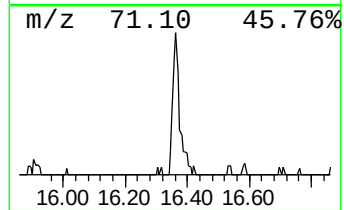
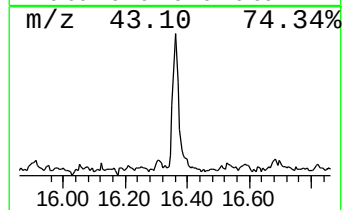
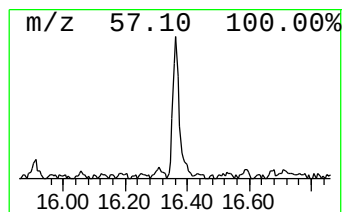
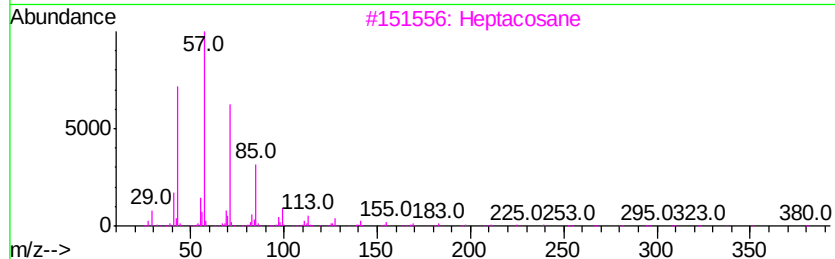
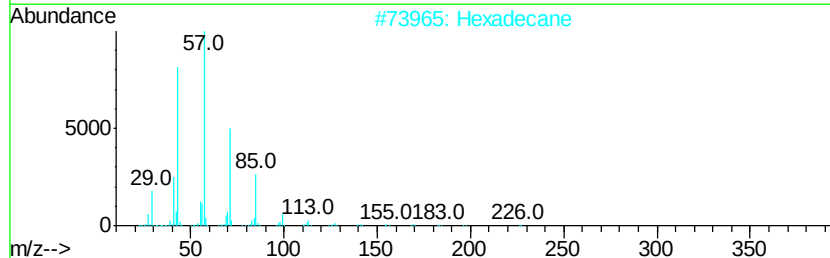
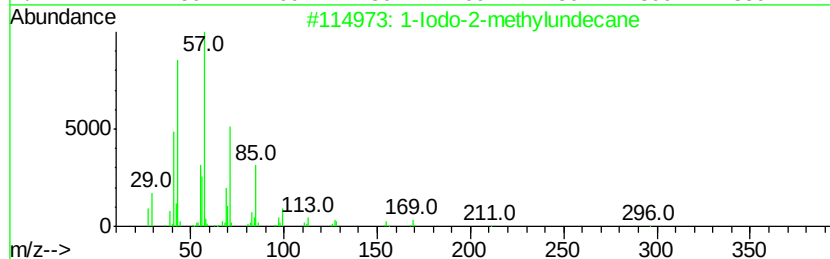
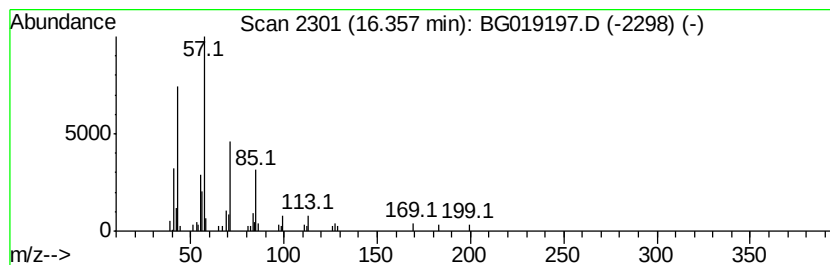
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG100415.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Iodo-2-methylundecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.36	2.78 ng/ul	48899	Phenanthrene-d10	17.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	86
2		Hexadecane	226	C16H34	000544-76-3	80
3		Heptacosane	380	C27H56	000593-49-7	80
4		Eicosane	282	C20H42	000112-95-8	74
5		Heptadecane	240	C17H36	000629-78-7	72



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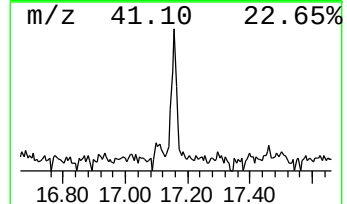
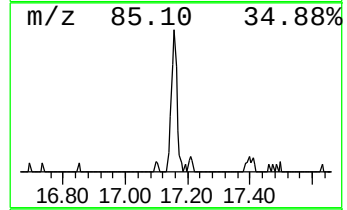
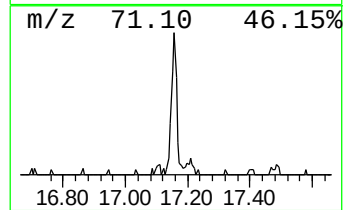
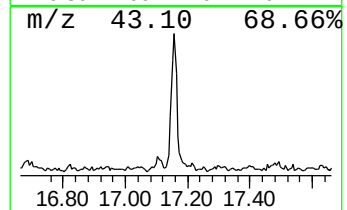
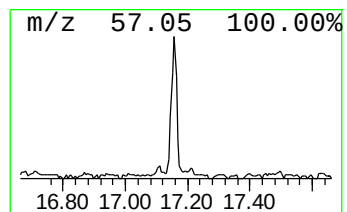
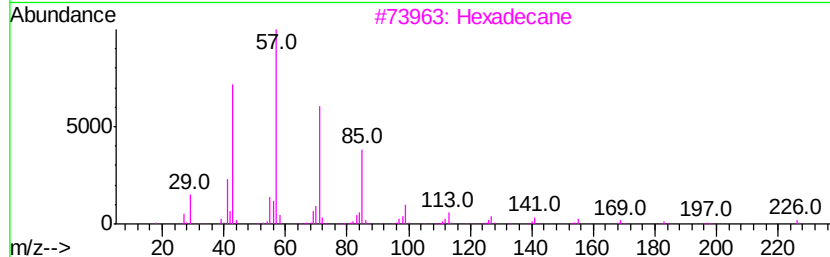
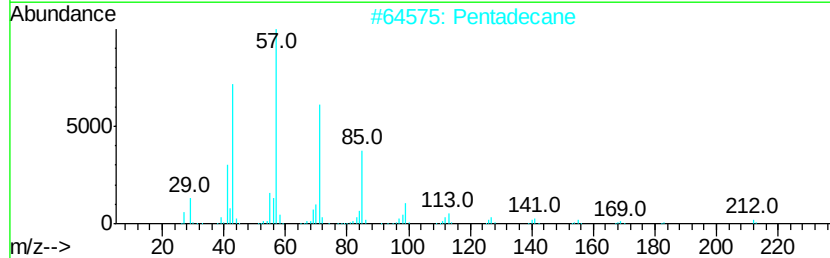
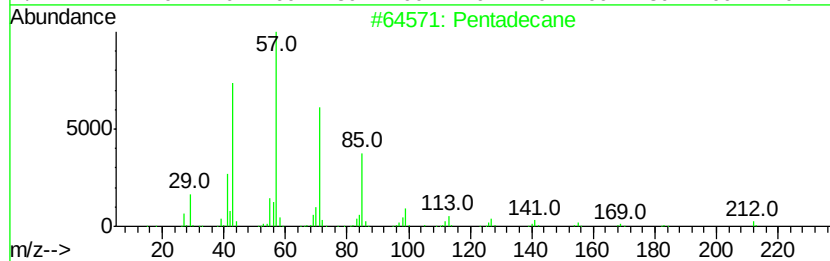
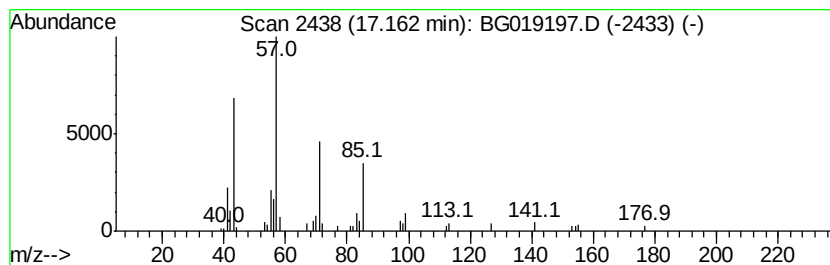
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TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.16	2.18 ng/ul	38391	Phenanthrene-d10	17.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane	212	C15H32	000629-62-9	90
2		Pentadecane	212	C15H32	000629-62-9	90
3		Hexadecane	226	C16H34	000544-76-3	90
4		Pentadecane	212	C15H32	000629-62-9	90
5		Tridecane	184	C13H28	000629-50-5	90



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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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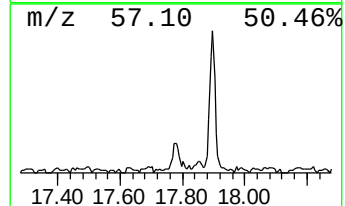
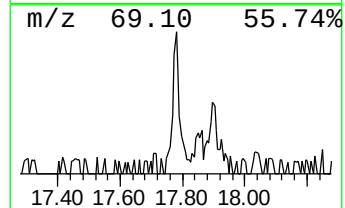
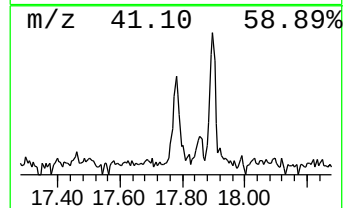
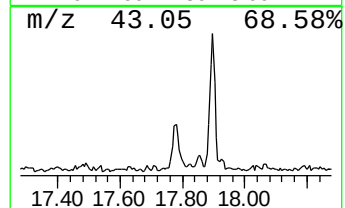
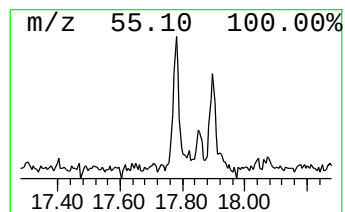
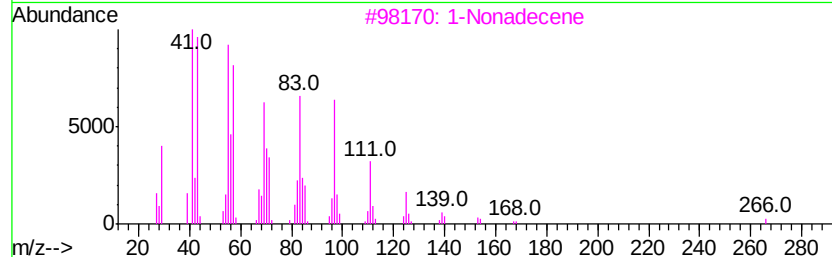
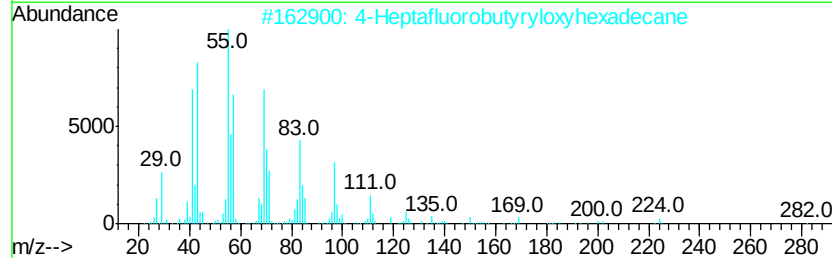
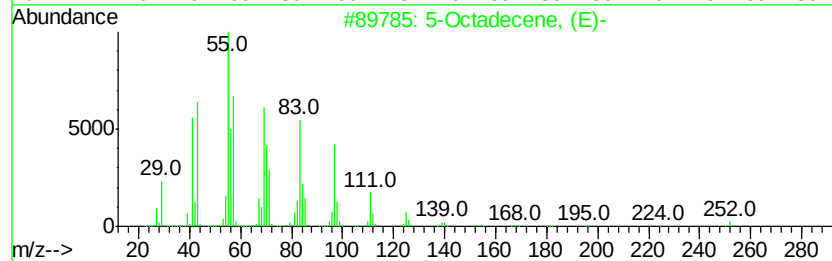
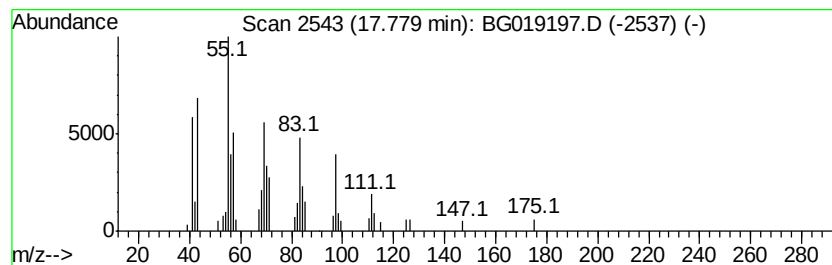
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Cyclic17.78 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.78	2.19 ng/ul	38630	Phenanthrene-d10	17.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Octadecene, (E)-	252	C18H36	007206-21-5	91
2		4-Heptafluorobutyryloxyhexadecane	438	C20H33F7O2	1000282-97-2	91
3		1-Nonadecene	266	C19H38	018435-45-5	91
4		5-Eicosene, (E)-	280	C20H40	074685-30-6	91
5		Trichloroacetic acid, dodecyl ester	330	C14H25Cl3O2	074339-50-7	87



Data Path : Z:\HPCHEM1\BNA G\DATA\BG101415\
Data File : BG019197.D
Acq On : 14 Oct 2015 2:36
Operator : UM/NP
Sample : G3996-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
E5LG3

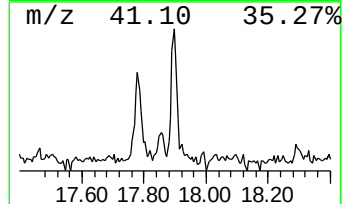
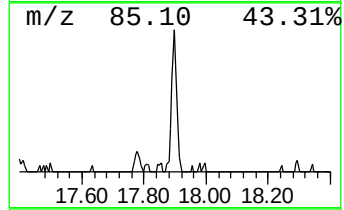
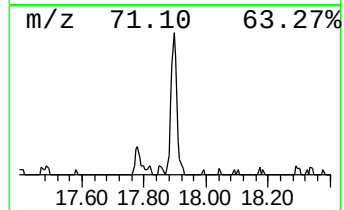
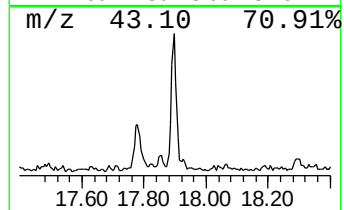
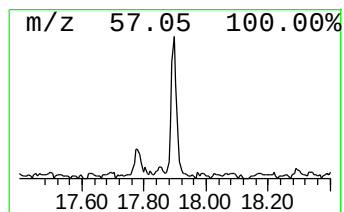
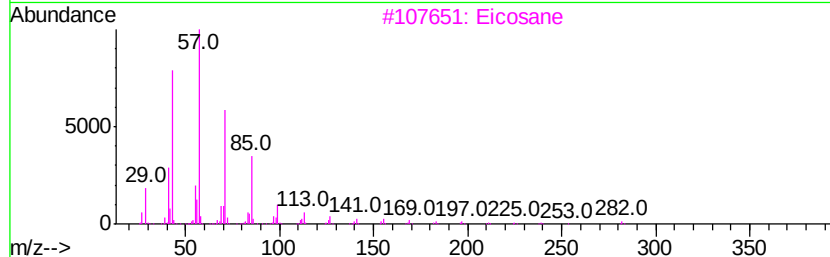
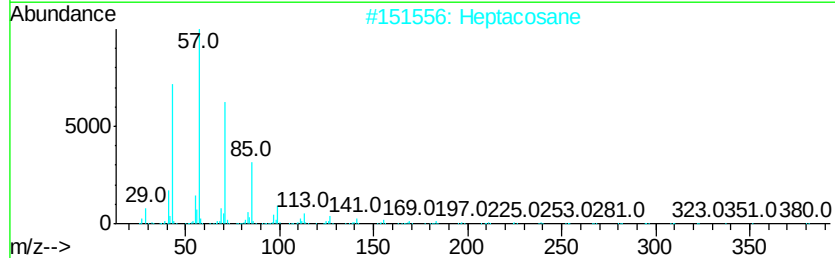
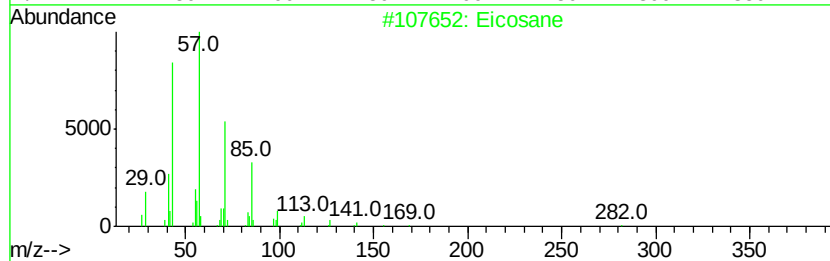
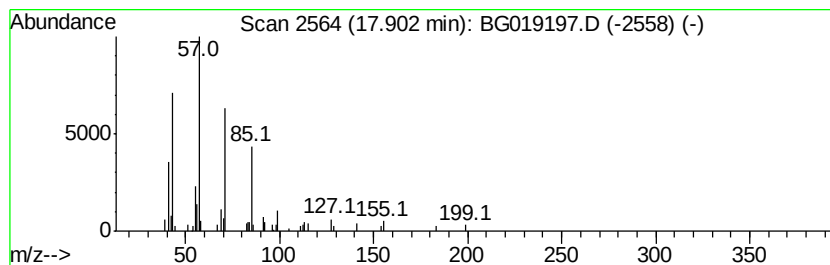
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.90	2.85 ng/ul	50129	Phenanthrene-d10	17.40

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	87
2			Heptacosane	380	C27H56	000593-49-7	86
3			Eicosane	282	C20H42	000112-95-8	80
4			Hexadecane	226	C16H34	000544-76-3	80
5			Tetradecane	198	C14H30	000629-59-4	72



Data Path : Z:\HPCHEM1\BNA G\DATA\BG101415\
Data File : BG019197.D
Acq On : 14 Oct 2015 2:36
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Sample : G3996-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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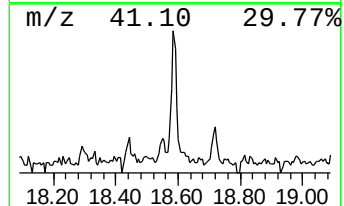
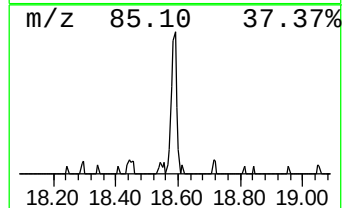
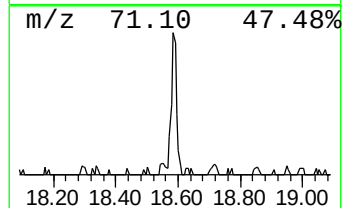
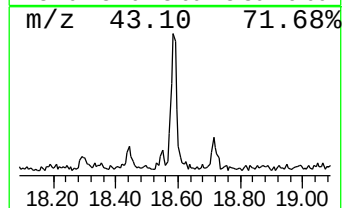
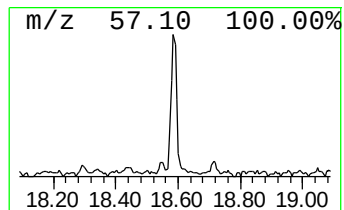
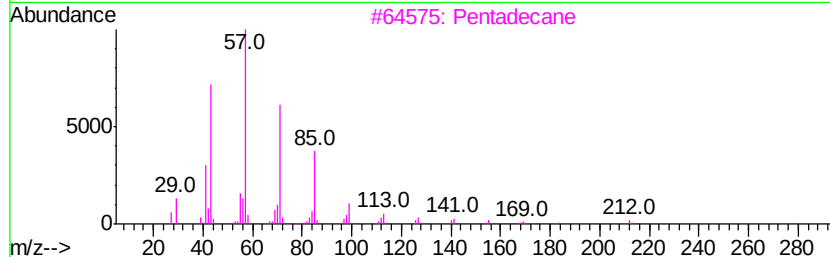
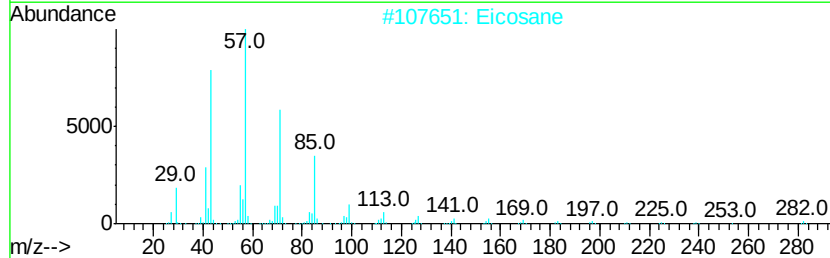
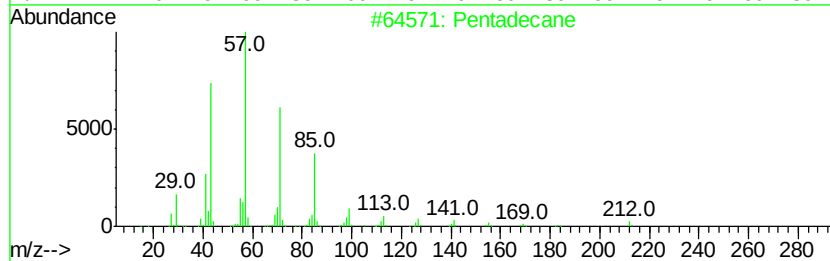
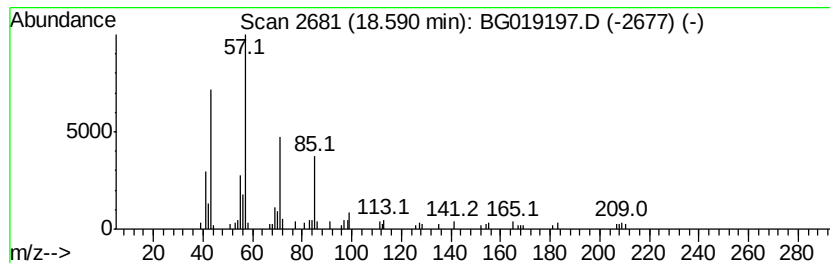
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Tridecane, 1-iodo- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.59	2.58 ng/ul	45418	Phenanthrene-d10	17.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane	212	C15H32	000629-62-9	80
2		Eicosane	282	C20H42	000112-95-8	80
3		Pentadecane	212	C15H32	000629-62-9	80
4		Pentadecane	212	C15H32	000629-62-9	80
5		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	80



Data Path : Z:\HPCHEM1\BNA G\DATA\BG101415\
Data File : BG019197.D
Acq On : 14 Oct 2015 2:36
Operator : UM/NP
Sample : G3996-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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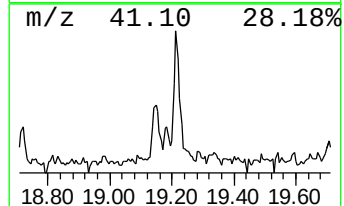
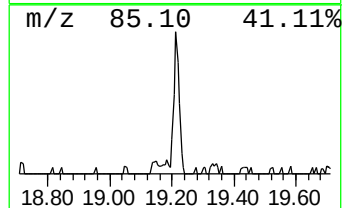
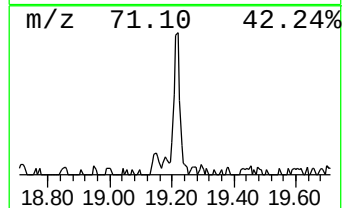
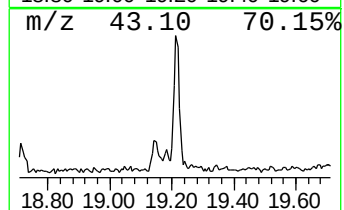
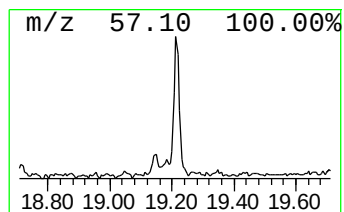
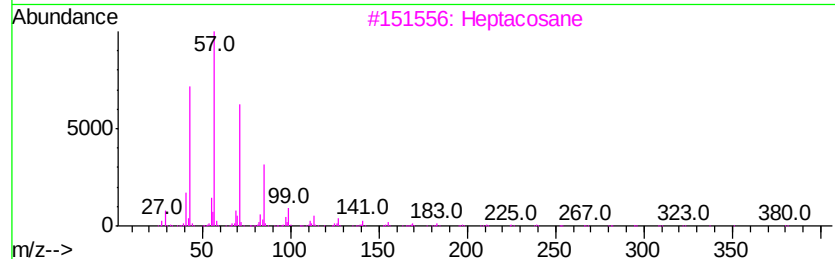
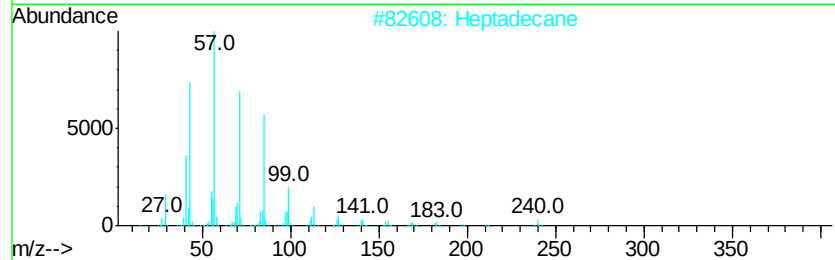
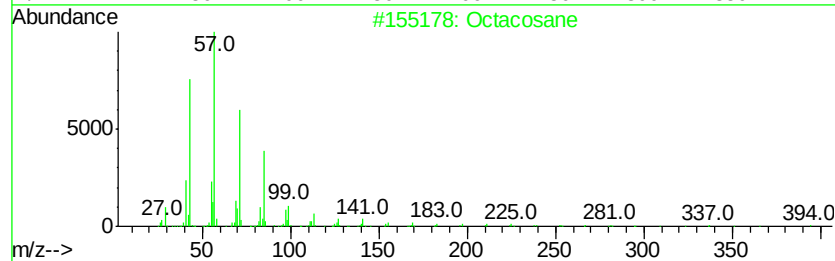
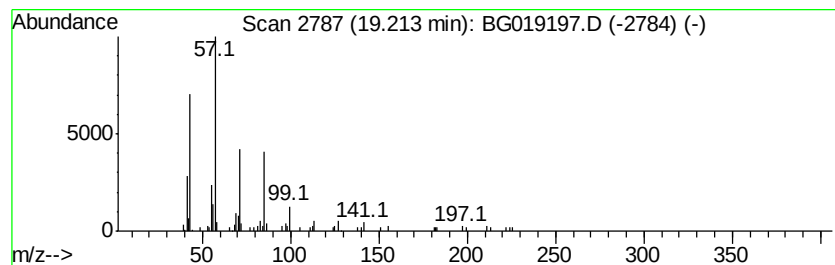
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.21	3.37 ng/ul	59313	Phenanthrene-d10	17.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	78
2		Heptadecane	240	C17H36	000629-78-7	72
3		Heptacosane	380	C27H56	000593-49-7	72
4		1-Iodoundecane	282	C11H23I	004282-44-4	72
5		triacontane	422	C30H62	000638-68-6	59



Data Path : Z:\HPCHEM1\BNA G\DATA\BG101415\
Data File : BG019197.D
Acq On : 14 Oct 2015 2:36
Operator : UM/NP
Sample : G3996-03
Misc :
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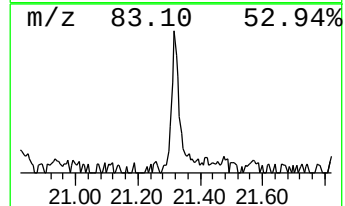
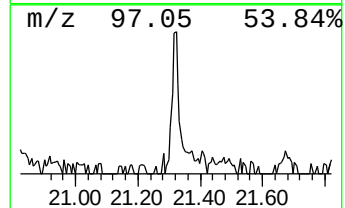
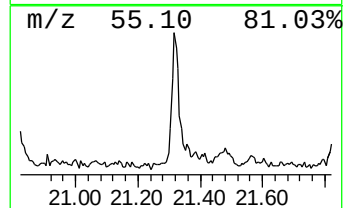
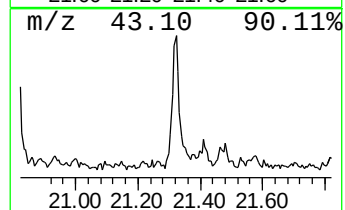
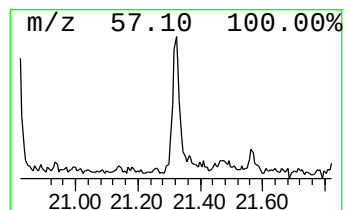
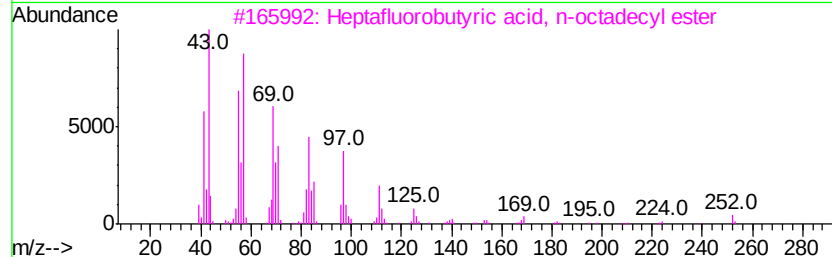
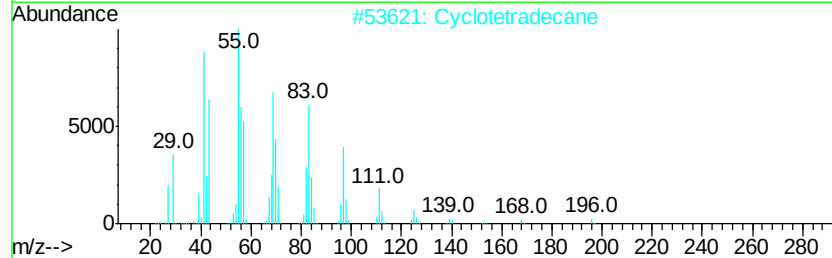
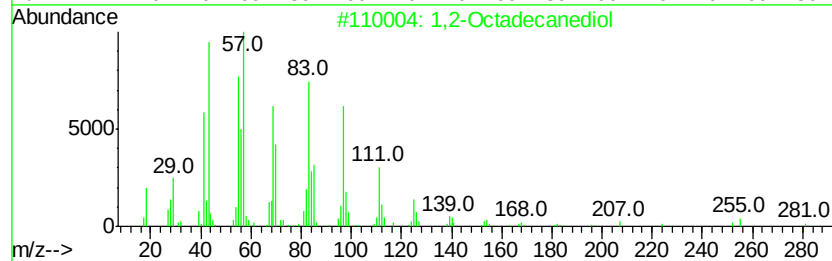
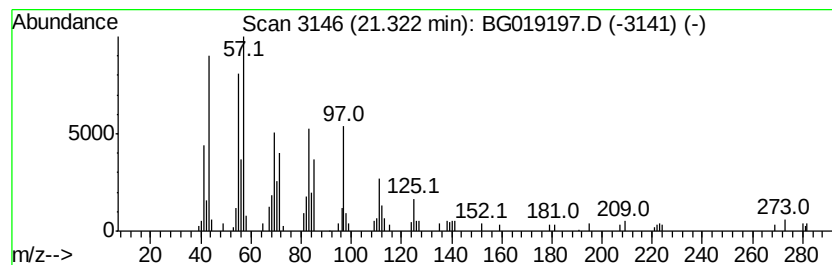
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG100415.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 1,2-Octadecanediol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
21.32	3.61 ng/ul	77507	Chrysene-d12	21.67		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,2-Octadecanediol	286	C18H38O2	020294-76-2	91
2		Cyclotetradecane	196	C14H28	000295-17-0	91
3		Heptafluorobutyric acid, n-octadecyl ester	466	C22H37F7O2	000400-57-7	91
4		Trifluoroacetic acid, n-octadecyl ester	366	C20H37F3O2	079392-43-1	91
5		1-Heptadecene	238	C17H34	006765-39-5	90



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG101415\
Data File : BG019197.D
Acq On : 14 Oct 2015 2:36
Operator : UM/NP
Sample : G3996-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG100415.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
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1-Iodo-2-methylun...	16.36	2.8	ng/ul	48899	4	17.40	352221	20.0
(DEL) Alkane: Str...	17.16	2.2	ng/ul	38391	4	17.40	352221	20.0
(DEL) Alkane: Cyc...	17.78	2.2	ng/ul	38630	4	17.40	352221	20.0
(DEL) Alkane: Str...	17.90	2.9	ng/ul	50129	4	17.40	352221	20.0
Tridecane, 1-iodo-	18.59	2.6	ng/ul	45418	4	17.40	352221	20.0
(DEL) Alkane: Str...	19.21	3.4	ng/ul	59313	4	17.40	352221	20.0
1,2-Octadecanediol	21.32	3.6	ng/ul	77507	5	21.67	429984	20.0