

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080819\
 Data File : BG042110.D
 Acq On : 9 Aug 2019 20:10
 Operator : HP/JU
 Sample : K4041-20
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BENX3

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:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.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	3.147	5	10	33	rVB	1393401	2275806	100.00%	11.007%
2	3.412	52	55	59	rBV3	7014	7509	0.33%	0.036%
3	3.653	92	96	98	rBV3	1741	2667	0.12%	0.013%
4	3.935	141	144	150	rVB7	3823	5943	0.26%	0.029%
5	4.076	163	168	173	rBV	9013	15782	0.69%	0.076%
6	4.170	178	184	189	rBV5	2993	6674	0.29%	0.032%
7	7.184	689	697	712	rBV2	131728	297099	13.05%	1.437%
8	7.342	718	724	734	rBV	104341	186775	8.21%	0.903%
9	7.554	754	760	772	rVV	157205	286807	12.60%	1.387%
10	8.024	834	840	852	rBV	219757	388638	17.08%	1.880%
11	8.735	955	961	977	rBV	178221	343182	15.08%	1.660%
12	9.193	1031	1039	1049	rBV2	148822	273498	12.02%	1.323%
13	9.640	1110	1115	1121	rVB2	2808	4011	0.18%	0.019%
14	9.922	1156	1163	1177	rBV	130053	243998	10.72%	1.180%
15	10.468	1249	1256	1270	rBV	200234	403102	17.71%	1.950%
16	10.850	1312	1321	1331	rBV	320167	585508	25.73%	2.832%
17	10.979	1337	1343	1359	rBV	155363	329973	14.50%	1.596%
18	12.454	1587	1594	1599	rBV2	3828	6599	0.29%	0.032%
19	13.053	1693	1696	1703	rVB2	2144	3131	0.14%	0.015%
20	13.300	1731	1738	1745	rVB4	2740	5160	0.23%	0.025%
21	13.582	1780	1786	1791	rVB2	6510	10398	0.46%	0.050%
22	14.087	1865	1872	1876	rBV	434079	657613	28.90%	3.180%
23	14.134	1876	1880	1892	rVB	110335	170591	7.50%	0.825%
24	14.381	1911	1922	1932	rBV2	513490	828487	36.40%	4.007%
25	14.687	1967	1974	1982	rBV	565926	884142	38.85%	4.276%
26	14.887	2002	2008	2025	rBV2	54989	142076	6.24%	0.687%
27	15.098	2039	2044	2052	rVV3	12545	20146	0.89%	0.097%
28	15.492	2104	2111	2115	rBV4	3485	7454	0.33%	0.036%
29	15.680	2136	2143	2151	rBV	711369	1068845	46.97%	5.169%
30	15.797	2157	2163	2172	rBV	119875	199971	8.79%	0.967%
31	15.921	2180	2184	2192	rVB5	9584	13142	0.58%	0.064%
32	16.373	2255	2261	2262	rBV2	3620	5399	0.24%	0.026%
33	16.402	2262	2266	2268	rVB5	2919	4161	0.18%	0.020%
34	16.467	2273	2277	2282	rVB4	2534	4560	0.20%	0.022%

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35	16.784	2329	2331	2338	rVB5	2987	3294	0.14%	0.016%
36	17.031	2369	2373	2378	rVB3	2448	3862	0.17%	0.019%
37	17.196	2397	2401	2403	rBV4	2913	4080	0.18%	0.020%
38	17.348	2424	2427	2433	rVB4	2750	5150	0.23%	0.025%
39	17.436	2436	2442	2448	rBV2	673685	1063096	46.71%	5.142%
40	17.536	2453	2459	2470	rVB	732248	1146791	50.39%	5.546%
41	18.030	2539	2543	2549	rVB6	5561	11212	0.49%	0.054%
42	18.112	2553	2557	2560	rVB5	3636	5090	0.22%	0.025%
43	18.159	2560	2565	2568	rBV6	2982	5836	0.26%	0.028%
44	18.330	2588	2594	2598	rBV5	28215	50825	2.23%	0.246%
45	18.453	2611	2615	2619	rVB6	5177	6319	0.28%	0.031%
46	18.512	2620	2625	2629	rBV4	17974	29782	1.31%	0.144%
47	18.641	2639	2647	2649	rBV8	6577	14051	0.62%	0.068%
48	18.811	2669	2676	2682	rBV8	7937	19459	0.86%	0.094%
49	19.058	2713	2718	2723	rBV7	6461	12079	0.53%	0.058%
50	19.111	2724	2727	2730	rVB5	4119	5305	0.23%	0.026%
51	19.146	2730	2733	2737	rBV6	4737	7642	0.34%	0.037%
52	19.228	2742	2747	2750	rBV3	10808	20188	0.89%	0.098%
53	19.322	2761	2763	2766	rVB4	5143	3779	0.17%	0.018%
54	19.358	2766	2769	2773	rBV6	6404	9438	0.41%	0.046%
55	19.463	2784	2787	2788	rBV	11094	12542	0.55%	0.061%
56	19.493	2788	2792	2796	rVB	61342	86297	3.79%	0.417%
57	19.546	2799	2801	2806	rVB6	5911	6181	0.27%	0.030%
58	19.593	2807	2809	2812	rVB4	4851	5130	0.23%	0.025%
59	19.634	2814	2816	2821	rVB5	10410	11020	0.48%	0.053%
60	19.716	2822	2830	2833	rBV2	10054	24976	1.10%	0.121%
61	19.828	2843	2849	2863	rBV	983443	1543115	67.81%	7.463%
62	20.080	2888	2892	2894	rBV5	8681	9499	0.42%	0.046%
63	20.180	2905	2909	2910	rBV4	9729	12254	0.54%	0.059%
64	20.357	2933	2939	2943	rBV2	22509	45732	2.01%	0.221%
65	20.451	2951	2955	2959	rVB2	15754	20953	0.92%	0.101%
66	20.503	2959	2964	2967	rBV4	17235	33563	1.47%	0.162%
67	20.644	2985	2988	2992	rBV4	12375	19896	0.87%	0.096%
68	20.856	3019	3024	3029	rBV2	39285	63667	2.80%	0.308%
69	20.944	3035	3039	3043	rBV7	20971	30208	1.33%	0.146%
70	21.303	3097	3100	3107	rBV8	31727	76033	3.34%	0.368%
71	21.367	3107	3111	3120	rVB2	127088	234497	10.30%	1.134%

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72	21.543	3137	3141	3147	rBV3	26687	48759	2.14%	0.236%
73	21.614	3147	3153	3159	rVB	503898	727553	31.97%	3.519%
74	21.720	3164	3171	3177	rBV2	719747	1248019	54.84%	6.036%
75	21.925	3202	3206	3213	rVB4	51737	97510	4.28%	0.472%
76	22.013	3218	3221	3227	rVB	23000	34118	1.50%	0.165%
77	22.548	3306	3312	3315	rBV2	132348	229398	10.08%	1.109%
78	22.583	3315	3318	3328	rVB	113839	211491	9.29%	1.023%
79	22.777	3346	3351	3360	rVB	62690	127370	5.60%	0.616%
80	23.259	3429	3433	3440	rVB	64978	119977	5.27%	0.580%
81	23.952	3547	3551	3557	rBV3	20648	49856	2.19%	0.241%
82	24.082	3567	3573	3581	rVB	169783	369995	16.26%	1.789%
83	24.763	3680	3689	3701	rBV	498185	1409427	61.93%	6.816%
84	24.992	3718	3728	3737	rBV2	442842	1311180	57.61%	6.341%
85	26.220	3929	3937	3947	rVB3	73326	235850	10.36%	1.141%
86	27.613	4167	4174	4186	rVB7	30032	104564	4.59%	0.506%

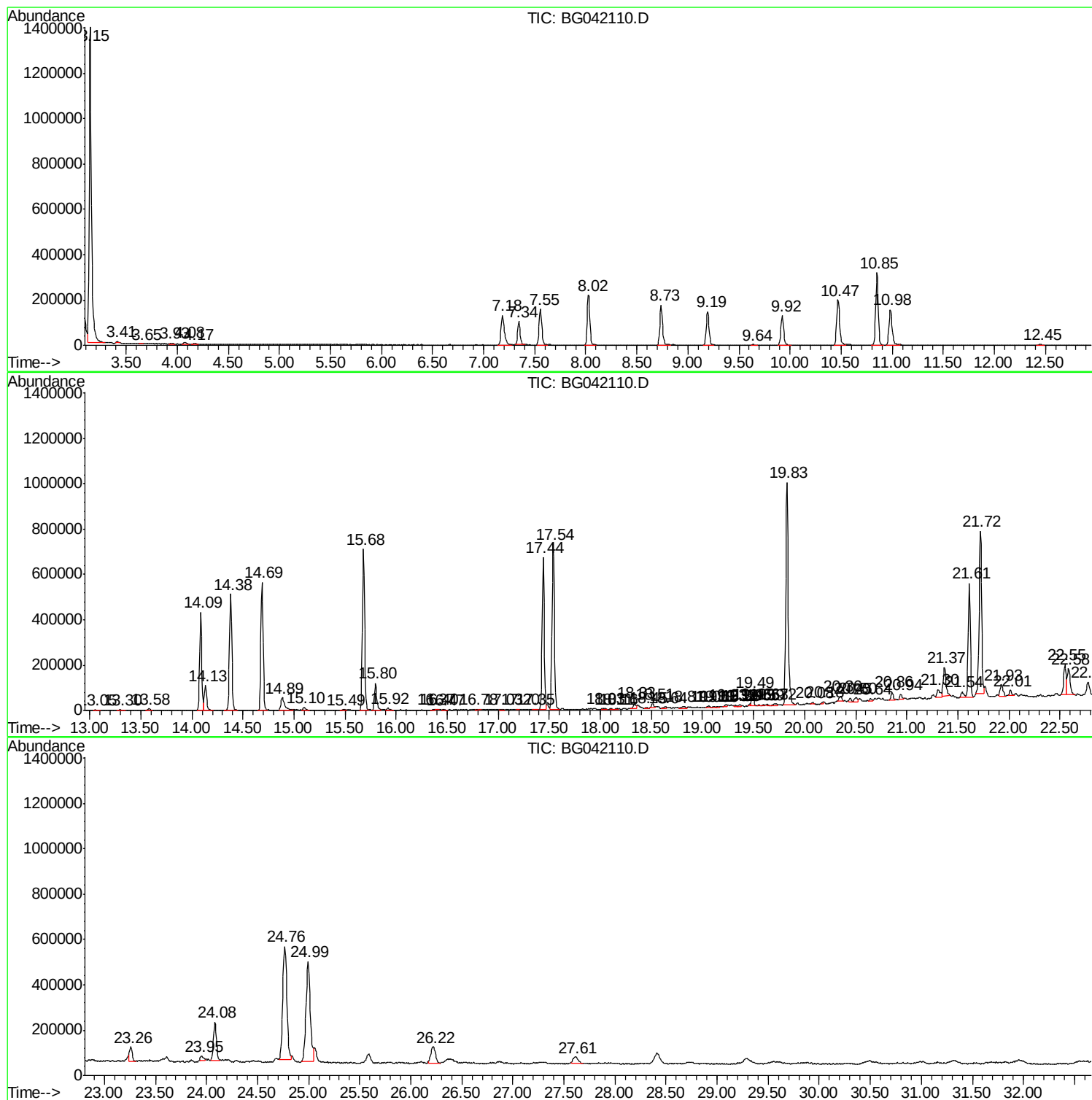
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
Quant Title : SVOA CALIBRATION

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



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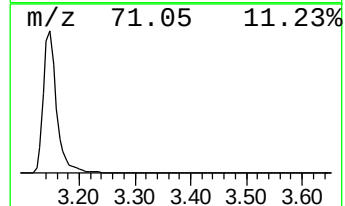
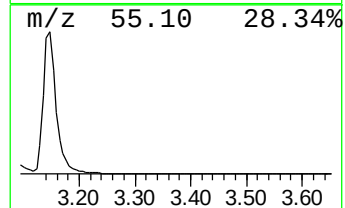
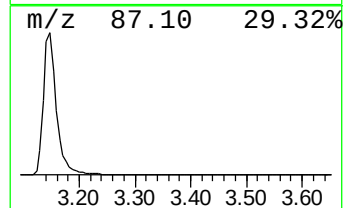
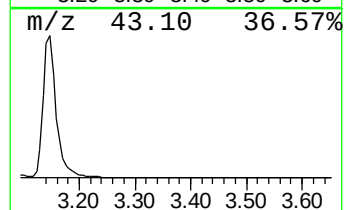
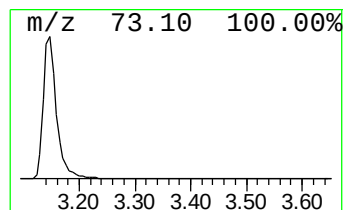
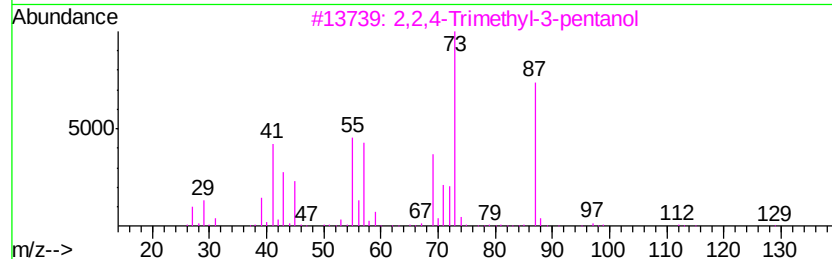
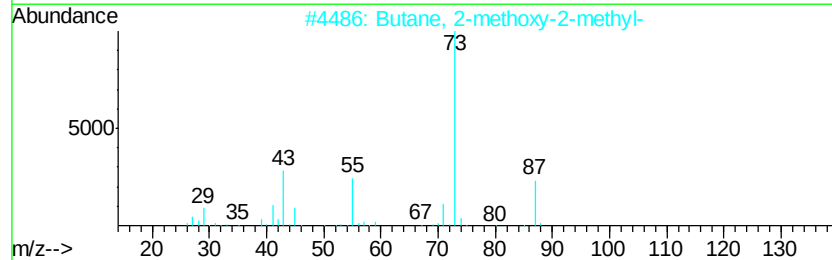
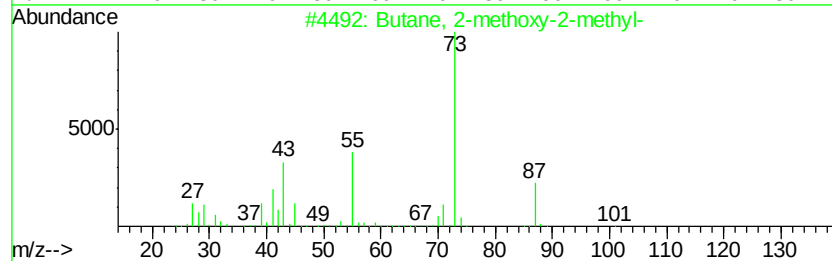
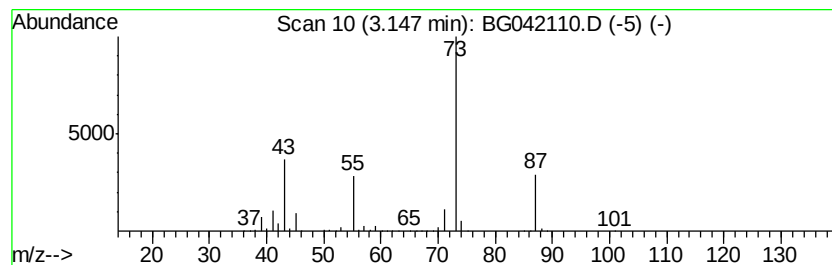
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.15	117.12 ng/ul	2275810	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	59
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	59
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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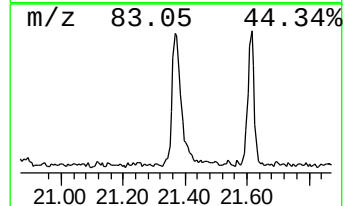
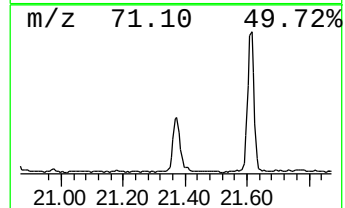
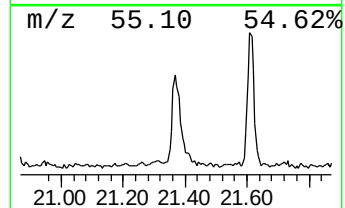
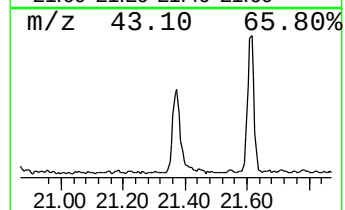
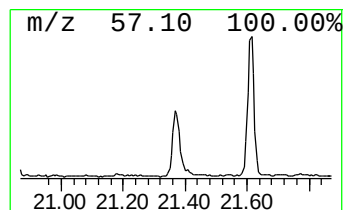
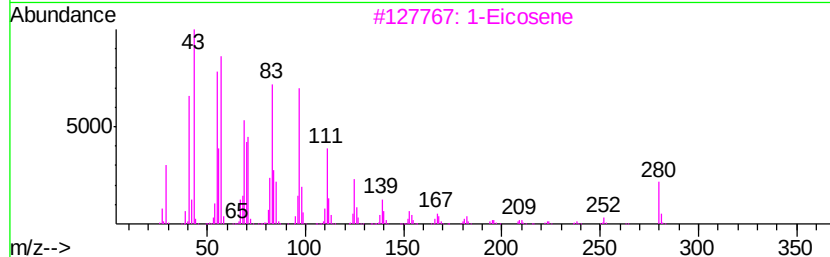
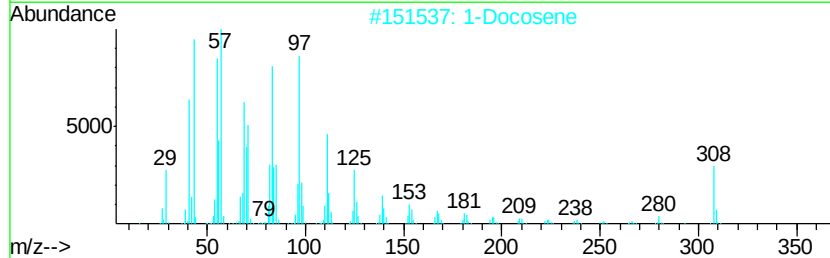
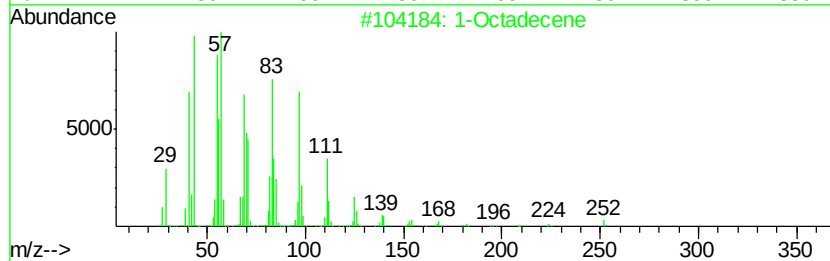
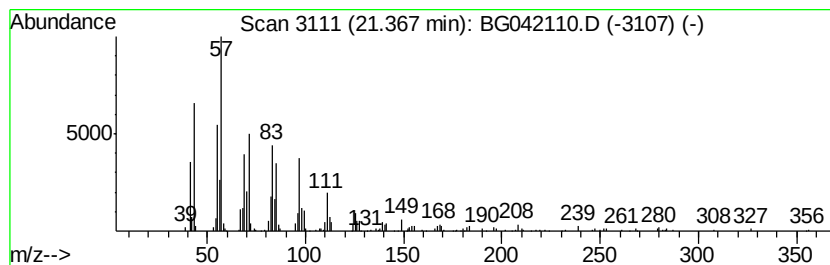
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Cyclic21.37 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.37	3.76 ng/ul	234497	Chrysene-d12	21.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Octadecene	252	C18H36	000112-88-9	96
2		1-Docosene	308	C22H44	001599-67-3	96
3		1-Eicosene	280	C20H40	003452-07-1	96
4		3-Eicosene, (E)-	280	C20H40	074685-33-9	95
5		1-Docosene	308	C22H44	001599-67-3	94



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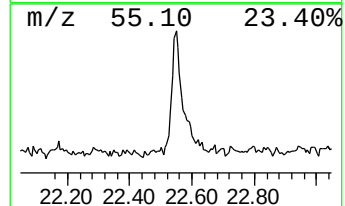
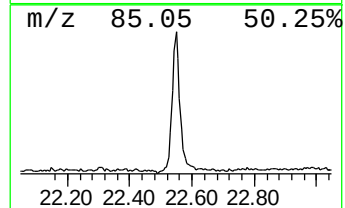
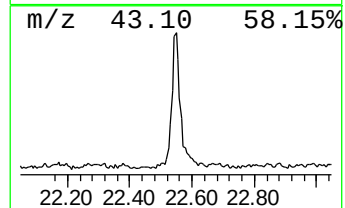
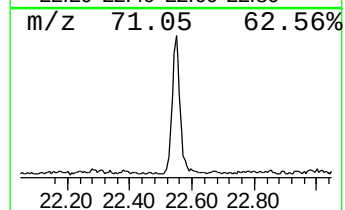
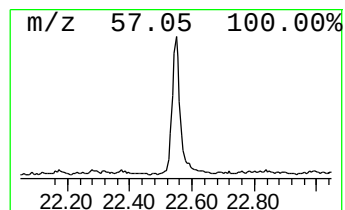
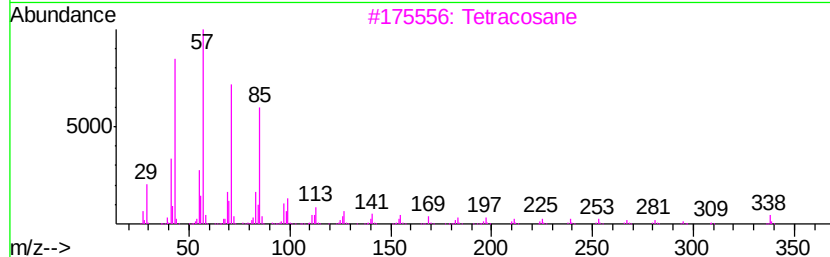
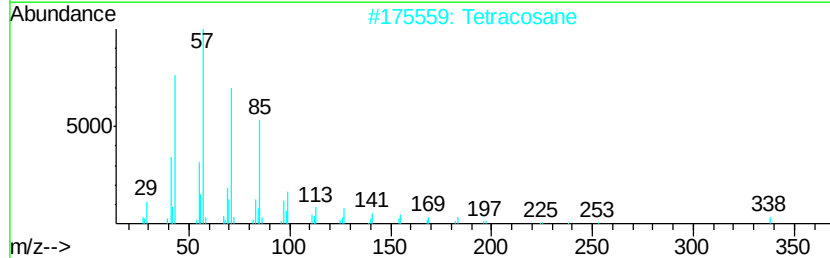
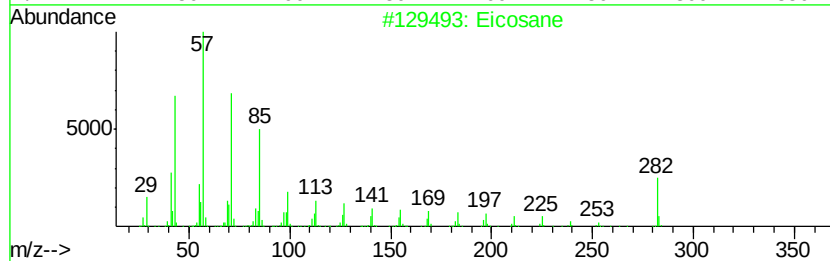
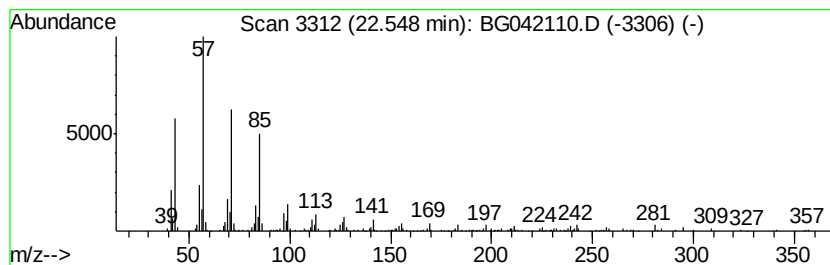
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.55	3.68 ng/ul	229398	Chrysene-d12	21.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	95
2		Tetracosane	338	C24H50	000646-31-1	93
3		Tetracosane	338	C24H50	000646-31-1	91
4		Heptadecane, 8-methyl-	254	C18H38	013287-23-5	90
5		Octacosane	394	C28H58	000630-02-4	90



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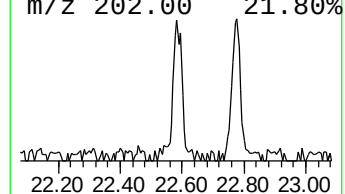
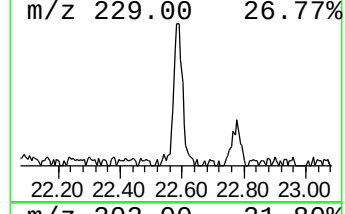
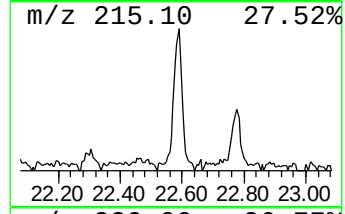
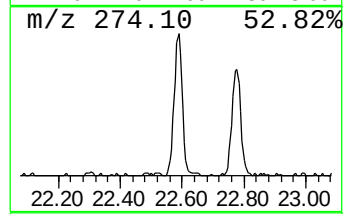
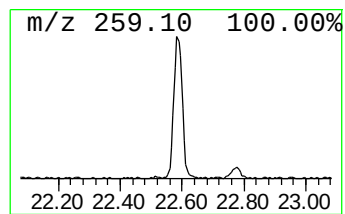
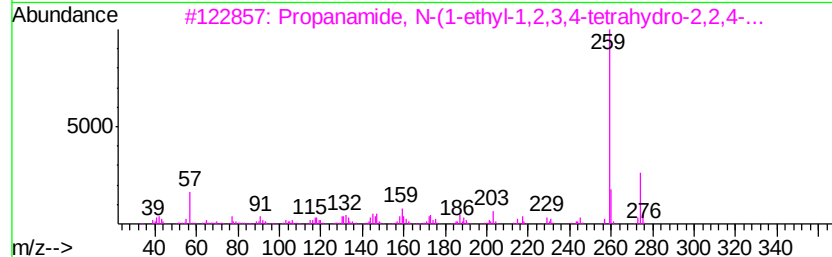
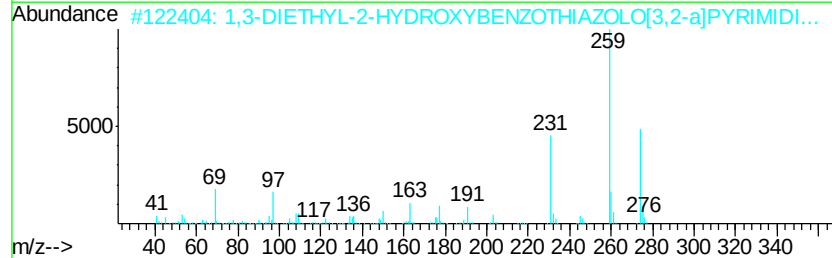
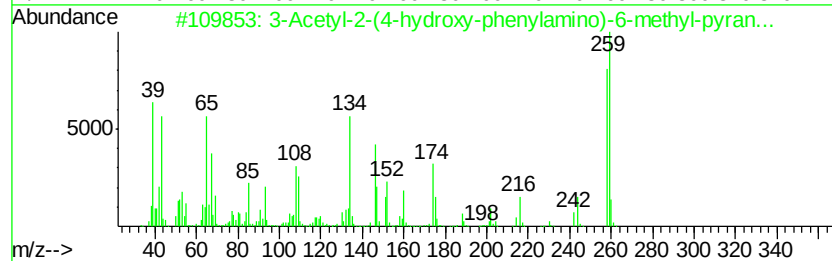
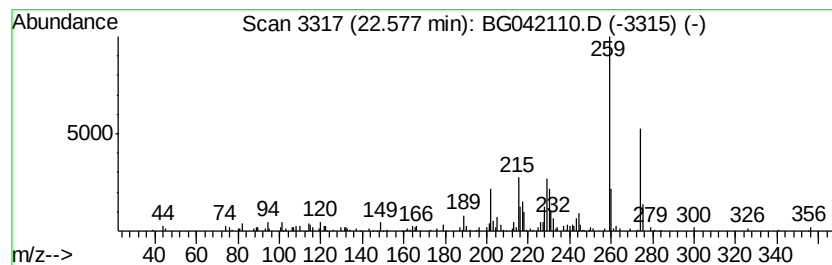
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BNA_G
ClientSampleId :
BENX3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 3-Acetyl-2-(4-hydroxy-pheny... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
22.58	3.39 ng/ul	211491	Chrysene-d12	21.72		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Acetyl-2-(4-hydroxy-phenylamin...	259	C14H13N04	1000300-99-5	86	
2	1,3-DIETHYL-2-HYDROXYBENZOTHAZO...	274	C14H14N2O2S	1000227-15-3	52	
3	Propanamide, N-(1-ethyl-1,2,3,4-...	274	C17H26N2O	063134-09-8	50	
4	Ethanone, 1-[5-(5-trifluoromethy...	274	C11H9F3N2O5	1000266-44-5	49	
5	10H-Phenoxaphosphine, 2-ethyl-10...	274	C15H15O3P	036360-91-5	47	



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080819\
Data File : BG042110.D
Acq On : 9 Aug 2019 20:10
Operator : HP/JU
Sample : K4041-20
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BENX3

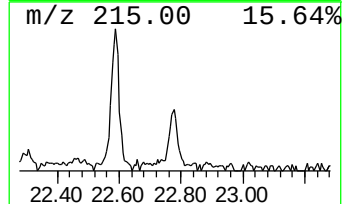
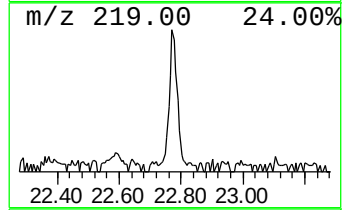
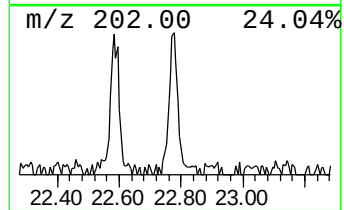
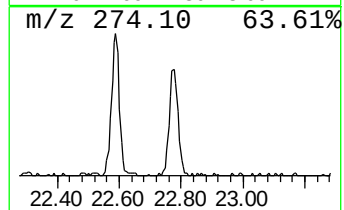
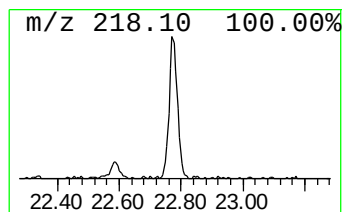
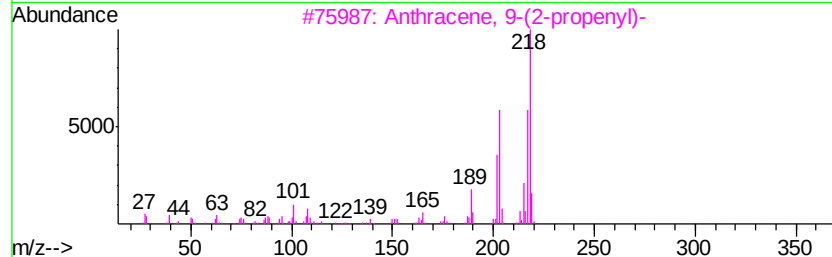
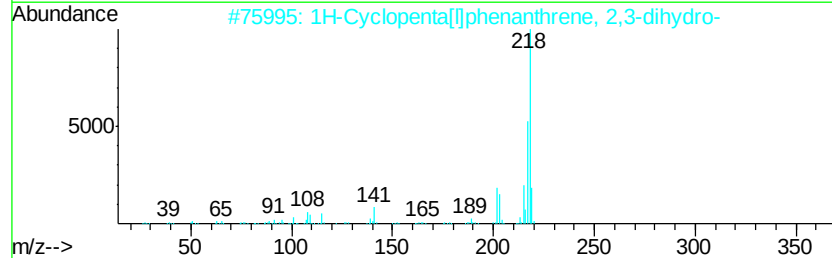
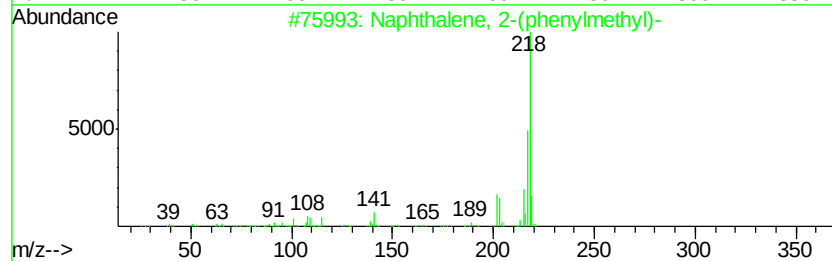
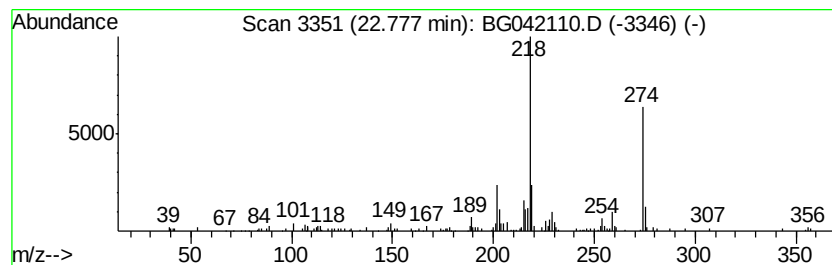
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.78	2.04 ng/ul	127370	Chrysene-d12	21.72

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2-(phenylmethyl)-	218	C17H14	000613-59-2	49
2		1H-Cyclopenta[<i>l</i>]phenanthrene, 2,...	218	C17H14	000723-98-8	49
3		Anthracene, 9-(2-propenyl)-	218	C17H14	023707-65-5	43
4		Isoquinoline, 1-(phenylmethyl)-	219	C16H13N	006907-59-1	43
5		Tetrachloro-o-benzoquinone	244	C6Cl4O2	002435-53-2	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080819\
Data File : BG042110.D
Acq On : 9 Aug 2019 20:10
Operator : HP/JU
Sample : K4041-20
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BENX3

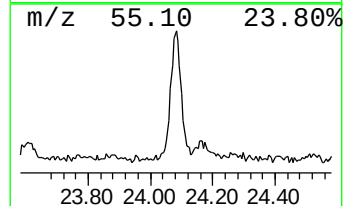
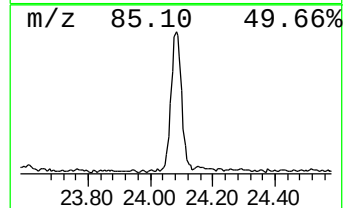
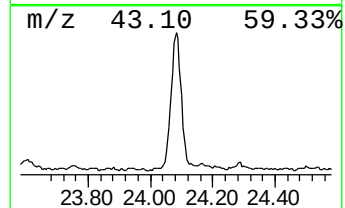
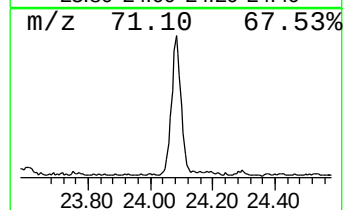
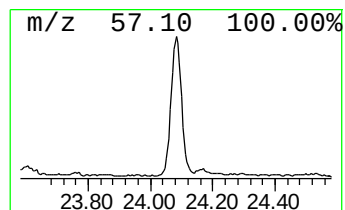
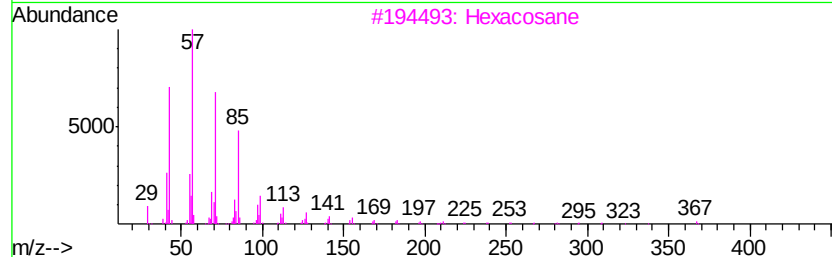
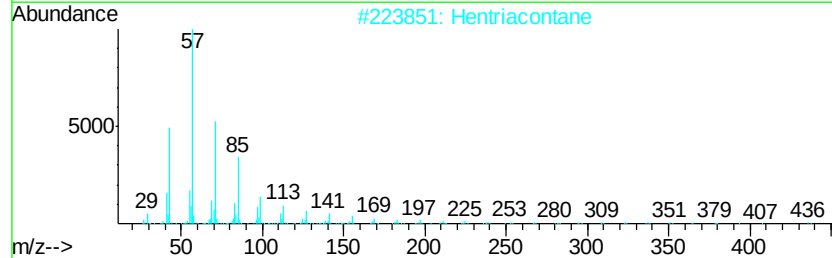
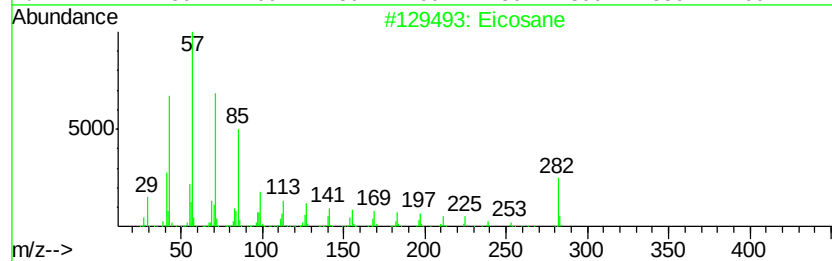
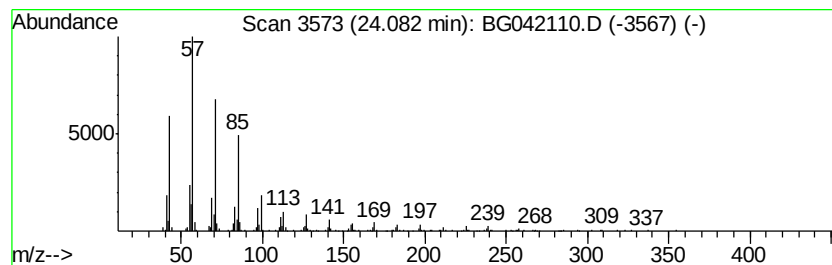
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.08	5.64 ng/ul	369995	Perylene-d12	24.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	96
2		Hentriacontane	437	C31H64	000630-04-6	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		Tetracosane, 11-decyl-	479	C34H70	055429-84-0	91
5		Tetratetracontane	619	C44H90	007098-22-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080819\
Data File : BG042110.D
Acq On : 9 Aug 2019 20:10
Operator : HP/JU
Sample : K4041-20
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BENX3

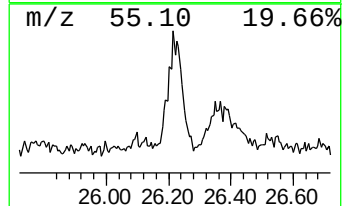
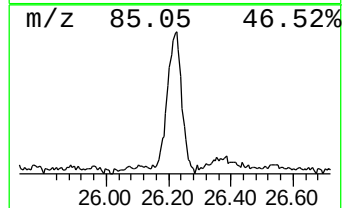
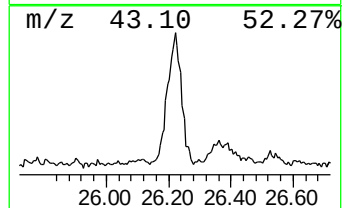
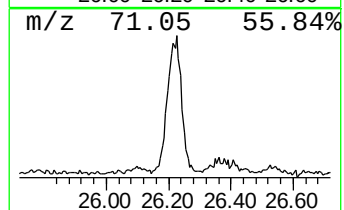
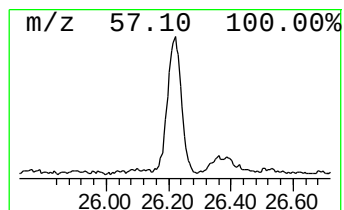
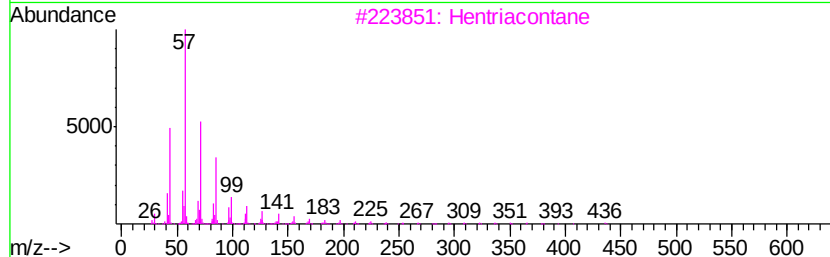
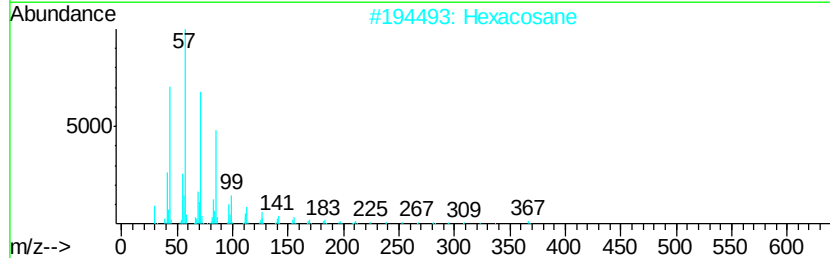
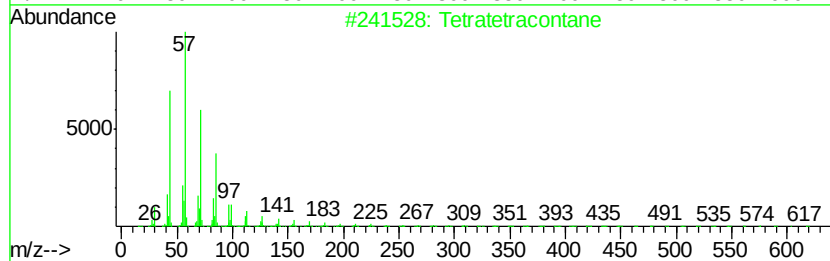
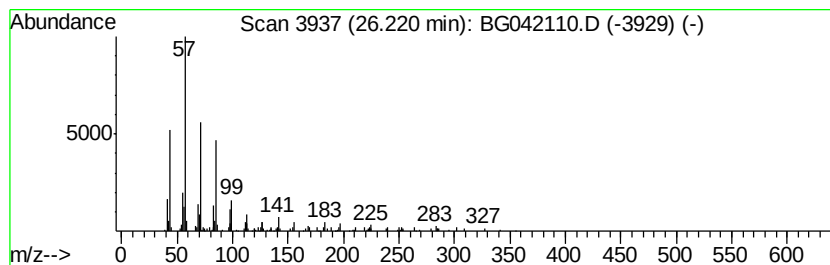
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.22	3.60 ng/ul	235850	Perylene-d12	24.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetratetracontane	619	C44H90	007098-22-8	90
2		Hexacosane	366	C26H54	000630-01-3	87
3		Hentriacontane	437	C31H64	000630-04-6	87
4		Octacosane	394	C28H58	000630-02-4	87
5		Hexacosane	366	C26H54	000630-01-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG080819\
Data File : BG042110.D
Acq On : 9 Aug 2019 20:10
Operator : HP/JU
Sample : K4041-20
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BENX3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.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
Butane, 2-methoxy...	3.15	117.1	ng/ul	2275810	1	8.03	388638	20.0
(DEL) Alkane: Cyc...	21.37	3.8	ng/ul	234497	5	21.72	1248020	20.0
(DEL) Alkane: Str...	22.55	3.7	ng/ul	229398	5	21.72	1248020	20.0
3-Acetyl-2-(4-hyd...	22.58	3.4	ng/ul	211491	5	21.72	1248020	20.0
unknown-01	22.78	2.0	ng/ul	127370	5	21.72	1248020	20.0
(DEL) Alkane: Str...	24.08	5.6	ng/ul	369995	6	24.99	1311180	20.0
(DEL) Alkane: Str...	26.22	3.6	ng/ul	235850	6	24.99	1311180	20.0