

Data Path : Z:\HPCHEM1\BNA G\DATA\BG020916\
 Data File : BG020825.D
 Acq On : 9 Feb 2016 20:03
 Operator : SJ/UM
 Sample : H1371-16
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C04J2

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-BG020816.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	3.256	65	71	80	rBV	67653	127572	23.64%	2.161%
2	3.332	80	84	90	rVB	66937	91084	16.88%	1.543%
3	3.966	188	192	196	rVV	4662	6258	1.16%	0.106%
4	4.148	220	223	228	rVB2	2524	3448	0.64%	0.058%
5	4.383	259	263	268	rVB	4256	6137	1.14%	0.104%
6	5.875	511	517	529	rBV2	17592	40183	7.45%	0.681%
7	6.257	575	582	587	rVB	15174	23414	4.34%	0.397%
8	7.344	763	767	772	rVV	6172	8935	1.66%	0.151%
9	7.409	772	778	786	rVV	47197	81661	15.13%	1.383%
10	7.479	786	790	795	rVB	6002	8682	1.61%	0.147%
11	7.585	801	808	814	rBV	37332	60919	11.29%	1.032%
12	7.655	814	820	826	rVV	4468	7512	1.39%	0.127%
13	7.796	837	844	853	rVV	42611	73508	13.62%	1.245%
14	7.879	853	858	860	rVV2	7595	11610	2.15%	0.197%
15	7.914	860	864	869	rVV	24162	36079	6.69%	0.611%
16	8.272	913	925	932	rBV	81665	153688	28.48%	2.603%
17	8.390	939	945	952	rBV3	12370	27437	5.08%	0.465%
18	8.824	1011	1019	1028	rVB5	4919	12507	2.32%	0.212%
19	8.913	1028	1034	1038	rBV3	6220	10838	2.01%	0.184%
20	8.965	1038	1043	1049	rVB2	48855	81809	15.16%	1.386%
21	9.095	1060	1065	1072	rBV3	5441	13008	2.41%	0.220%
22	9.383	1106	1114	1118	rBV3	4922	9019	1.67%	0.153%
23	9.441	1118	1124	1130	rBV	39428	68036	12.61%	1.152%
24	9.494	1130	1133	1138	rVB2	3675	5077	0.94%	0.086%
25	10.170	1239	1248	1254	rBV2	30412	53097	9.84%	0.899%
26	10.493	1298	1303	1305	rBV2	2566	3561	0.66%	0.060%
27	10.516	1305	1307	1315	rVB3	2756	4214	0.78%	0.071%
28	10.710	1333	1340	1347	rVB	54545	91988	17.05%	1.558%
29	10.933	1372	1378	1387	rBV2	12700	23816	4.41%	0.403%
30	11.045	1393	1397	1400	rVV3	2245	3262	0.60%	0.055%
31	11.098	1400	1406	1412	rVV	130714	238352	44.17%	4.037%
32	11.145	1412	1414	1421	rVB	4428	6455	1.20%	0.109%
33	11.227	1421	1428	1437	rBV	46655	84585	15.68%	1.433%
34	11.521	1472	1478	1483	rVB	11721	18631	3.45%	0.316%

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35	11.662	1496	1502	1509	rBV	28096	46737	8.66%	0.792%
36	11.732	1509	1514	1520	rVV2	2988	4151	0.77%	0.070%
37	12.091	1570	1575	1579	rBV	2515	3792	0.70%	0.064%
38	12.478	1635	1641	1645	rBV2	2816	4088	0.76%	0.069%
39	12.737	1678	1685	1691	rBV	4973	8310	1.54%	0.141%
40	12.954	1717	1722	1727	rVB	3805	6062	1.12%	0.103%
41	13.336	1783	1787	1791	rBV2	2970	4951	0.92%	0.084%
42	13.412	1796	1800	1806	rVB4	2438	3750	0.69%	0.064%
43	13.694	1843	1848	1856	rBV2	6272	10492	1.94%	0.178%
44	14.205	1930	1935	1941	rBV2	16474	26812	4.97%	0.454%
45	14.282	1941	1948	1952	rBV	106852	157341	29.16%	2.665%
46	14.329	1952	1956	1963	rVB	44368	65911	12.22%	1.116%
47	14.581	1993	1999	2007	rBV	135843	210476	39.01%	3.565%
48	14.716	2015	2022	2025	rBV2	2913	4908	0.91%	0.083%
49	14.758	2025	2029	2031	rBV	15835	21102	3.91%	0.357%
50	14.799	2031	2036	2044	rVV	339157	539582	100.00%	9.139%
51	14.887	2044	2051	2058	rVB2	197708	320991	59.49%	5.437%
52	14.957	2058	2063	2069	rVB5	2989	5063	0.94%	0.086%
53	15.051	2072	2079	2086	rBV	34312	56042	10.39%	0.949%
54	15.192	2098	2103	2105	rBV2	3203	5116	0.95%	0.087%
55	15.210	2105	2106	2110	rVV3	3083	3256	0.60%	0.055%
56	15.275	2110	2117	2122	rVV3	20950	33323	6.18%	0.564%
57	15.369	2127	2133	2136	rBV6	4450	8489	1.57%	0.144%
58	15.545	2158	2163	2168	rBV3	4724	7397	1.37%	0.125%
59	15.604	2168	2173	2178	rVB5	6998	11102	2.06%	0.188%
60	15.668	2178	2184	2186	rBV4	5104	8140	1.51%	0.138%
61	15.703	2186	2190	2196	rVV	21920	28099	5.21%	0.476%
62	15.874	2210	2219	2225	rVV	169027	257684	47.76%	4.365%
63	15.921	2225	2227	2232	rVV4	4741	7437	1.38%	0.126%
64	15.974	2232	2236	2242	rVB2	21094	30739	5.70%	0.521%
65	16.109	2252	2259	2264	rBV2	8154	16237	3.01%	0.275%
66	16.209	2272	2276	2281	rVB5	2696	5424	1.01%	0.092%
67	16.326	2292	2296	2299	rBV4	3677	5870	1.09%	0.099%
68	16.561	2332	2336	2338	rBV	20560	25059	4.64%	0.424%
69	16.585	2338	2340	2346	rVB3	16435	21431	3.97%	0.363%
70	16.696	2355	2359	2362	rBV5	5219	7300	1.35%	0.124%
71	16.896	2388	2393	2398	rVV3	13010	23352	4.33%	0.396%

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72	16.966	2403	2405	2412	rVB6	4205	7922	1.47%	0.134%
73	17.066	2419	2422	2425	rBV3	3693	4666	0.86%	0.079%
74	17.096	2425	2427	2430	rVV2	4235	4071	0.75%	0.069%
75	17.219	2443	2448	2454	rVB	56948	83796	15.53%	1.419%
76	17.278	2454	2458	2461	rBV5	3915	6420	1.19%	0.109%
77	17.354	2467	2471	2474	rBV	13571	18890	3.50%	0.320%
78	17.407	2477	2480	2483	rVB4	6925	9910	1.84%	0.168%
79	17.519	2496	2499	2504	rVB4	3241	3871	0.72%	0.066%
80	17.624	2511	2517	2521	rBV	219071	338413	62.72%	5.732%
81	17.665	2521	2524	2528	rVV	41449	60621	11.23%	1.027%
82	17.718	2528	2533	2542	rVV2	171014	280608	52.00%	4.753%
83	17.795	2542	2546	2552	rVV7	7618	13517	2.51%	0.229%
84	18.088	2593	2596	2600	rVB4	8910	11590	2.15%	0.196%
85	18.165	2605	2609	2617	rVB	17667	29175	5.41%	0.494%
86	18.482	2658	2663	2668	rBV4	37336	61077	11.32%	1.035%
87	18.558	2670	2676	2680	rVB2	28847	54420	10.09%	0.922%
88	18.682	2693	2697	2701	rBV5	9616	16272	3.02%	0.276%
89	18.735	2701	2706	2710	rVB3	73626	105056	19.47%	1.779%
90	18.899	2730	2734	2743	rVB2	35740	54226	10.05%	0.918%
91	19.081	2760	2765	2773	rVB	210034	280150	51.92%	4.745%
92	19.216	2784	2788	2794	rVB	55039	73426	13.61%	1.244%
93	19.392	2816	2818	2823	rVB4	13510	15280	2.83%	0.259%
94	19.657	2858	2863	2867	rVB	54047	73820	13.68%	1.250%
95	19.704	2867	2871	2875	rBV	33340	44357	8.22%	0.751%
96	19.739	2875	2877	2882	rVB6	11560	14730	2.73%	0.249%
97	19.986	2914	2919	2923	rBV	179851	285235	52.86%	4.831%
98	20.421	2989	2993	2999	rBV	37312	47219	8.75%	0.800%
99	21.513	3176	3179	3186	rVB	50443	70214	13.01%	1.189%
100	21.907	3240	3246	3252	rBV	187075	332591	61.64%	5.633%

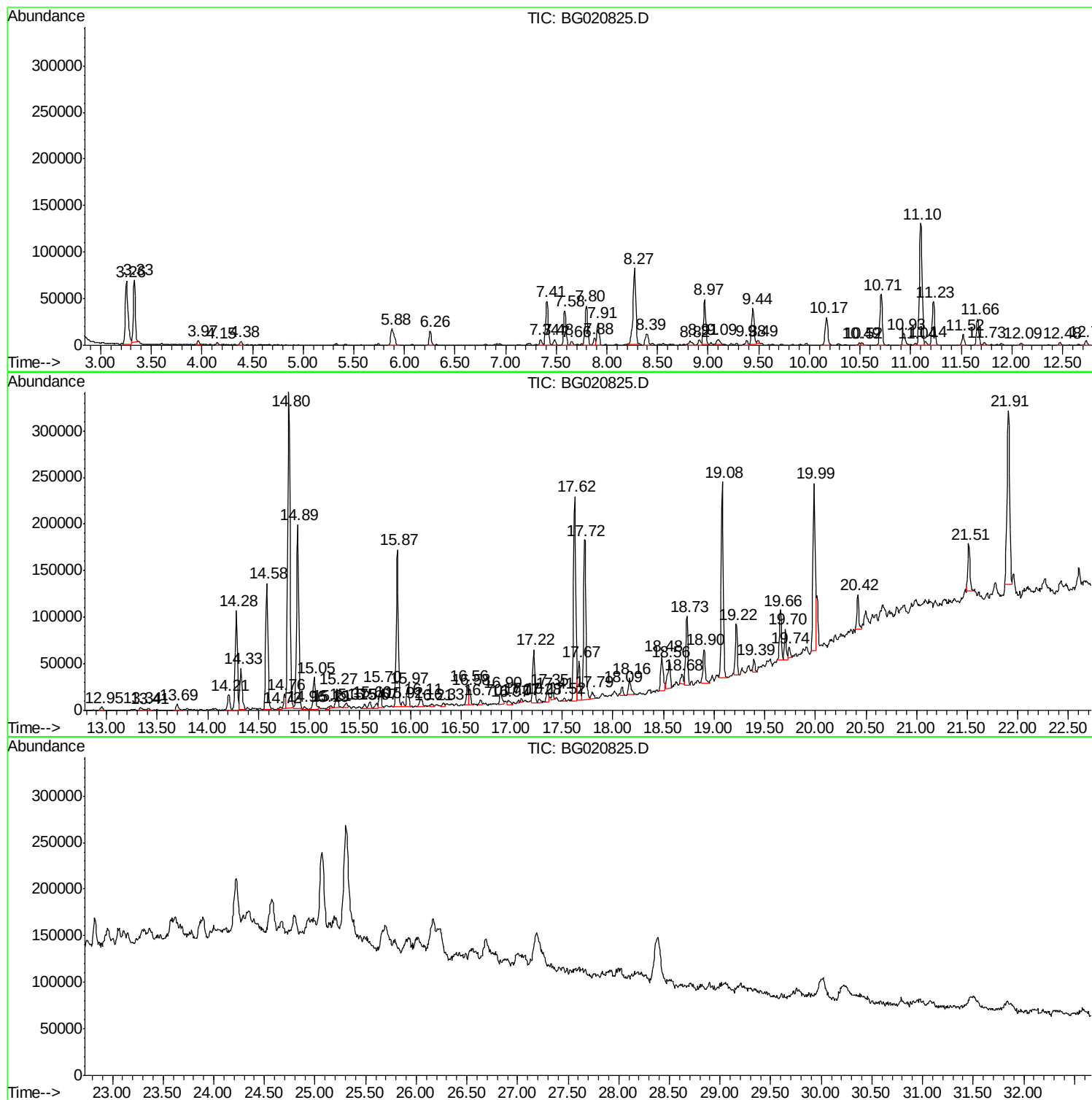
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG020816.M
Quant Title : SVOA CALIBRATION

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



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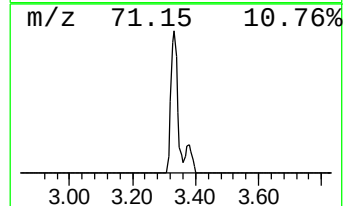
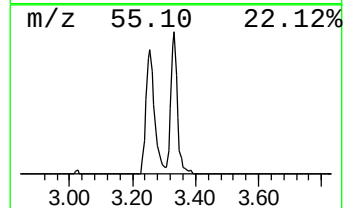
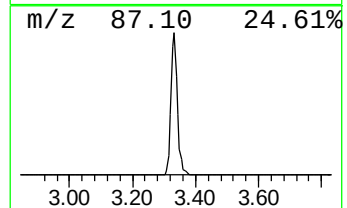
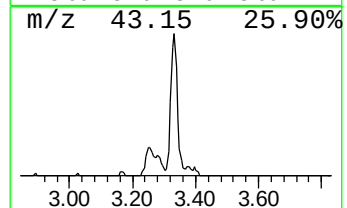
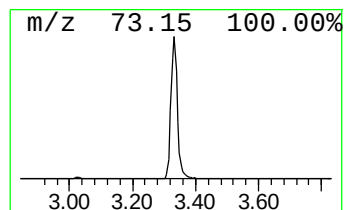
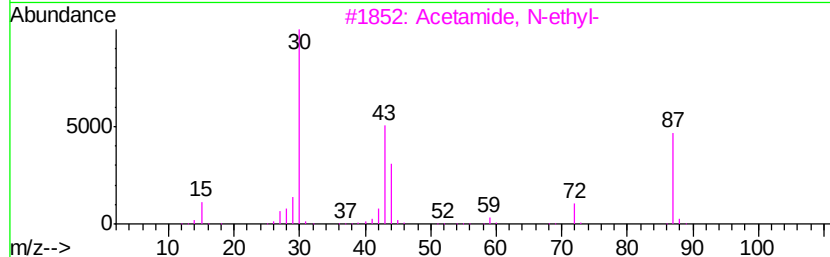
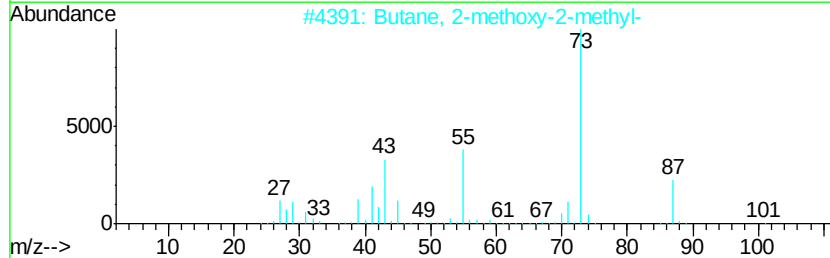
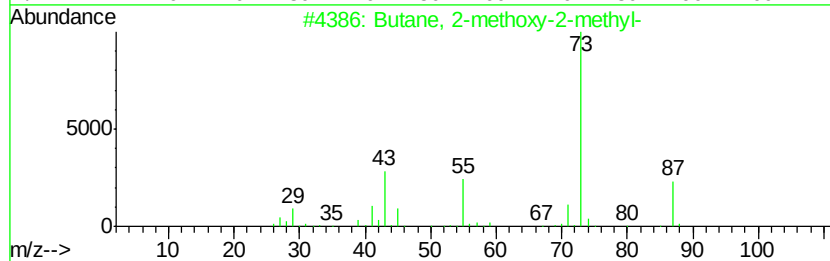
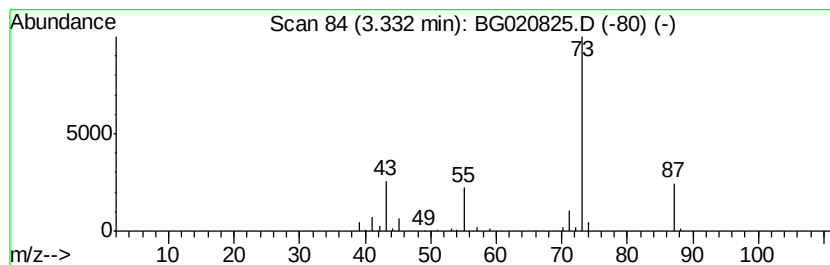
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG020816.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.33	11.85 ng/ul	91084	1,4-Dichlorobenzene-d4	8.27

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		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	9



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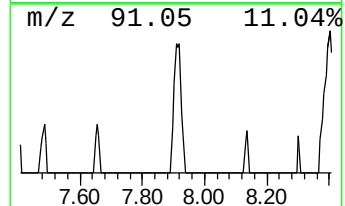
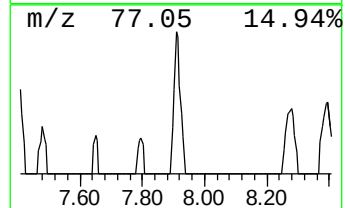
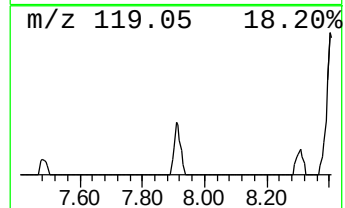
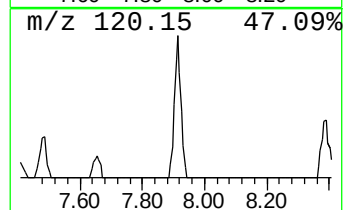
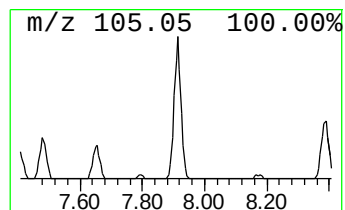
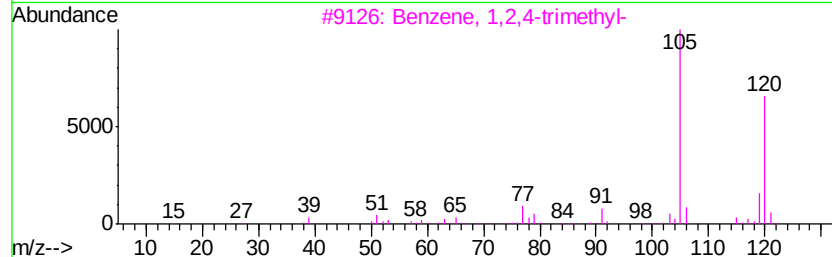
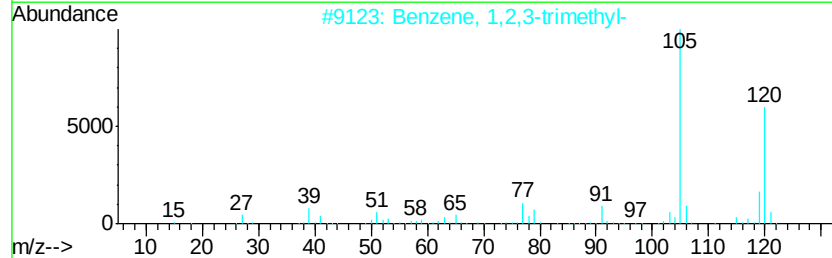
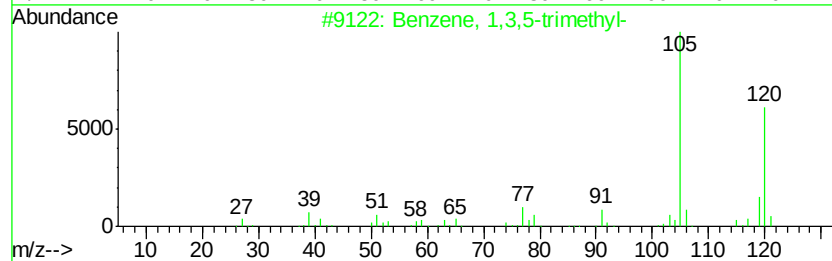
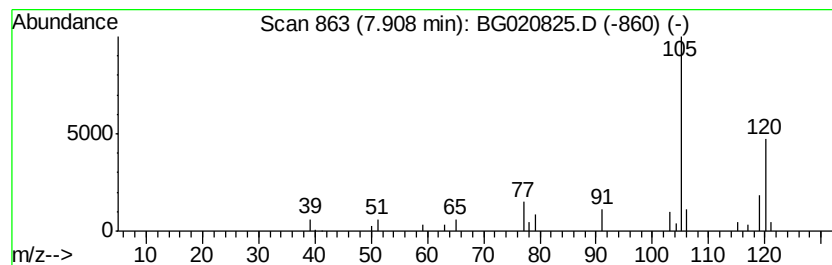
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG020816.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1,3,5-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.91	4.70 ng/ul	36079	1,4-Dichlorobenzene-d4	8.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	93
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90
4		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	90
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90



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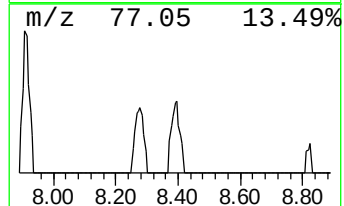
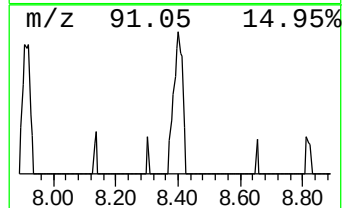
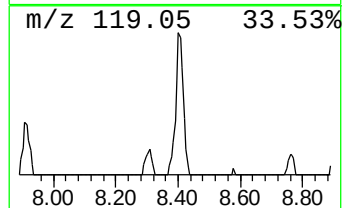
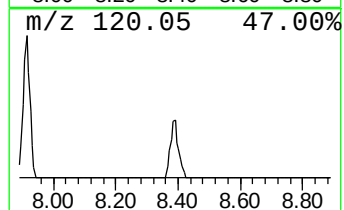
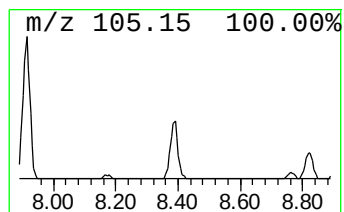
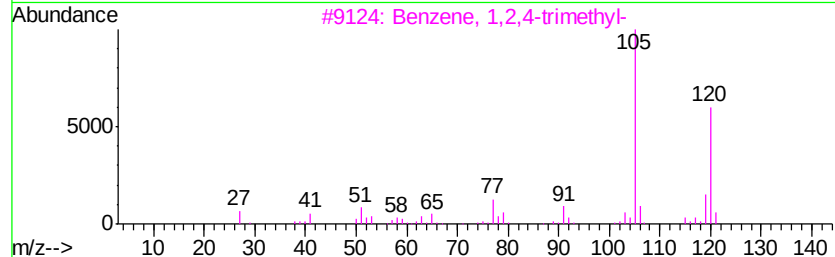
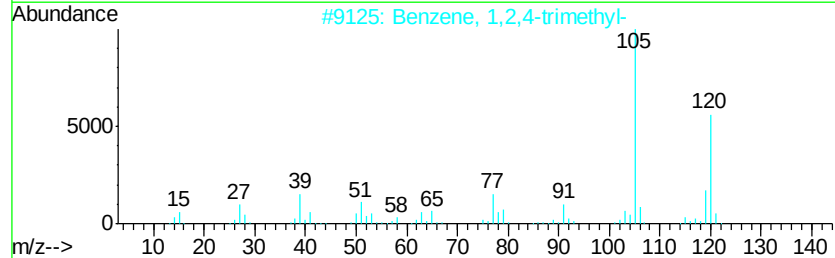
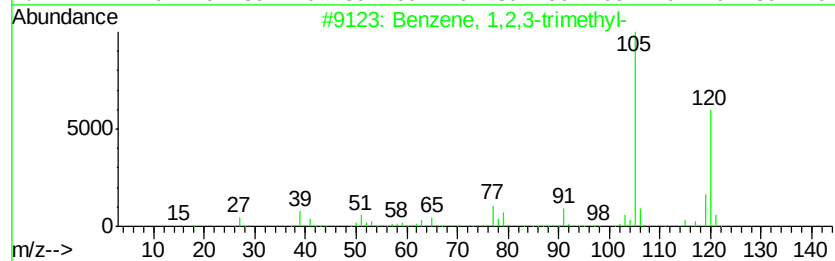
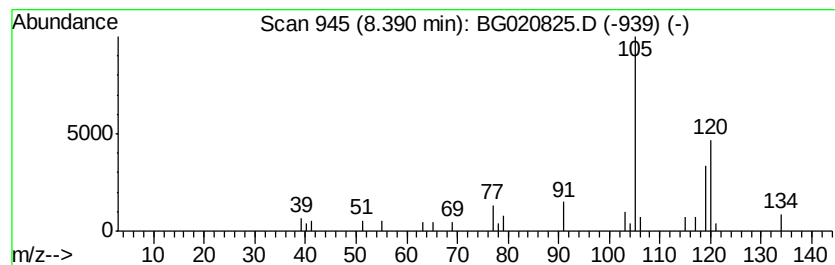
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.39	3.57 ng/ul	27437	1,4-Dichlorobenzene-d4	8.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	83
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	80
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	80
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	72
5		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	72



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Operator : SJ/UM
Sample : H1371-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

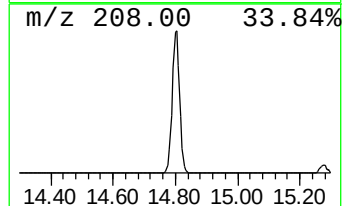
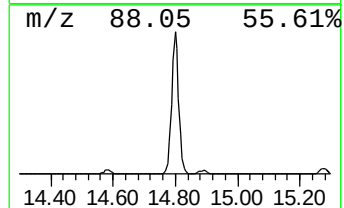
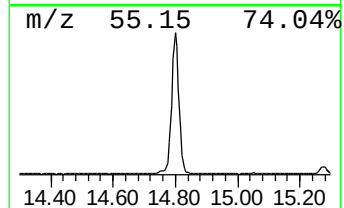
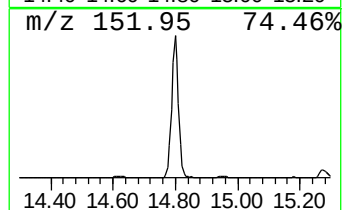
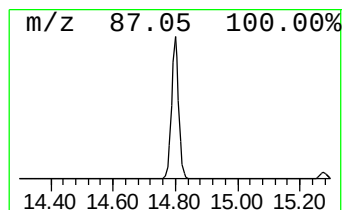
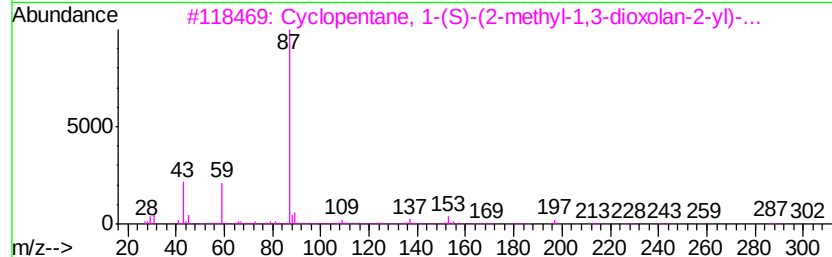
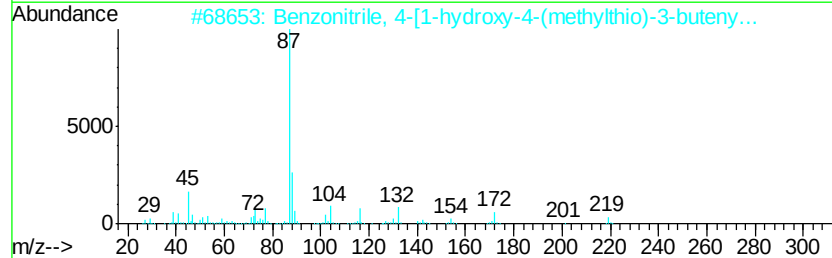
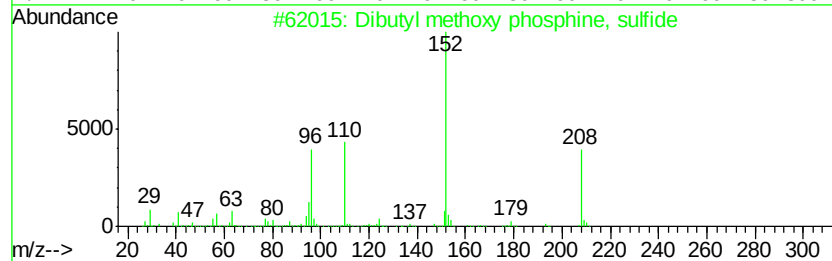
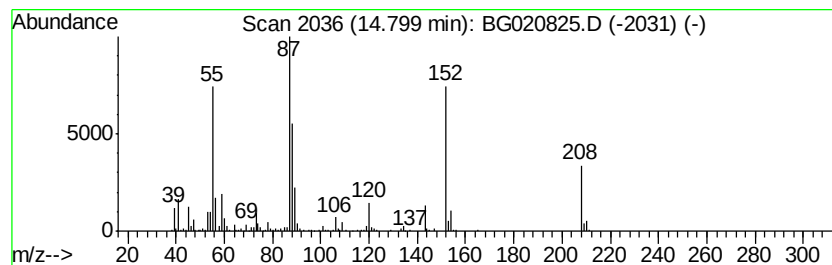
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG020816.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.80	33.62 ng/ul	539582	Acenaphthene-d10	14.89
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Dibutyl methoxy phosphine, sulfide	208 C9H210PS	098958-59-9 25
2		Benzonitrile, 4-[1-hydroxy-4-(me...	219 C12H13NOS	097369-82-9 10
3		Cyclopentane, 1-(S)-(2-methyl-1,...	302 C15H26O6	1000155-90-5 10
4		2-Methoxy-3-methyl-butylric acid,...	146 C7H14O3	109989-32-4 10
5		Hydrazine, 1,1-diethyl-2-pentyl-	158 C9H22N2	067398-41-8 10



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020916\
Data File : BG020825.D
Acq On : 9 Feb 2016 20:03
Operator : SJ/UM
Sample : H1371-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

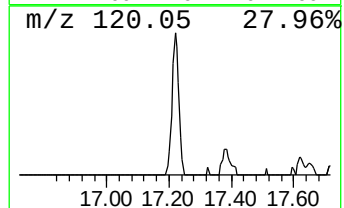
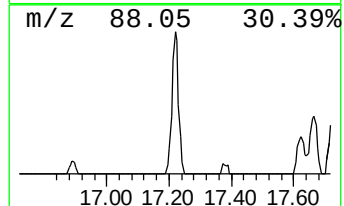
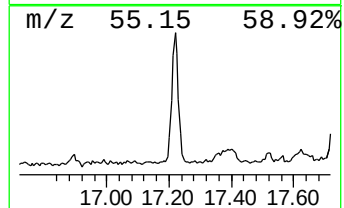
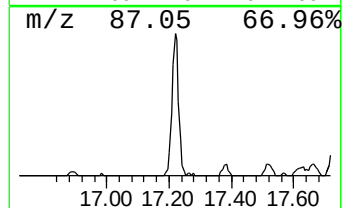
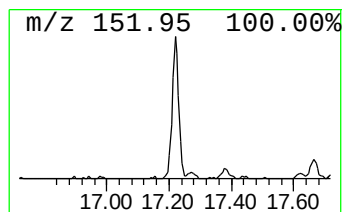
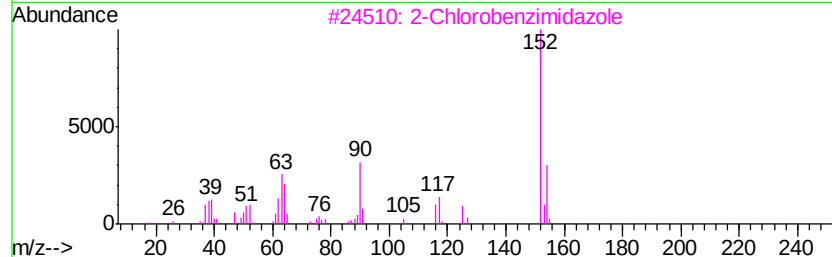
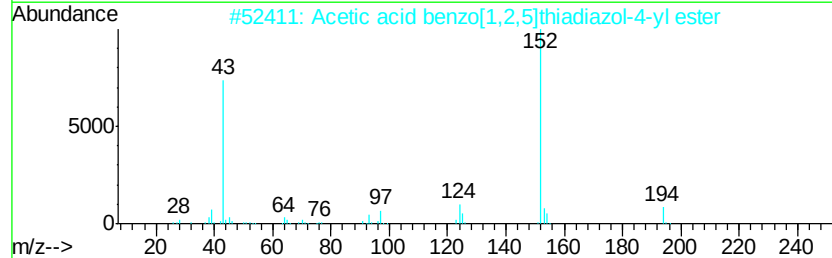
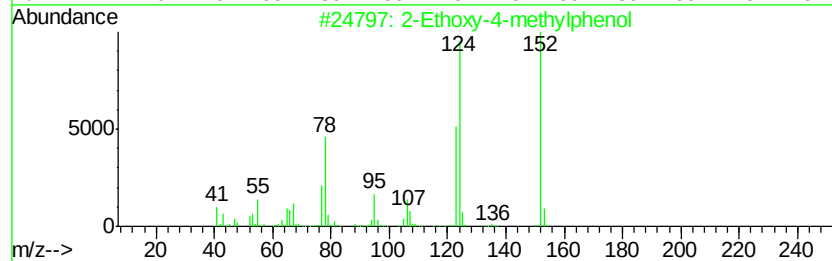
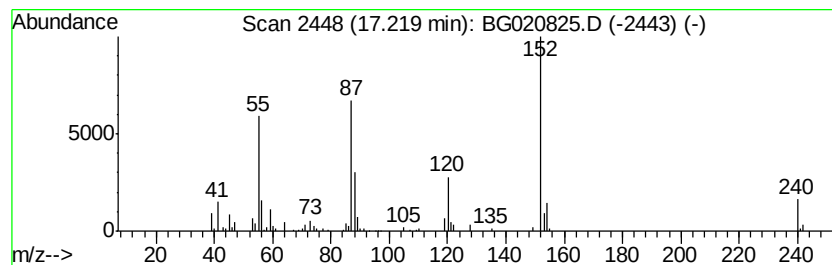
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG020816.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-02 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.22	4.95 ng/ul	83796	Phenanthrene-d10	17.62	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Ethoxy-4-methylphenol	152	C9H12O2	002563-07-7	25
2	Acetic acid benzo[1,2,5]thiadiaz...	194	C8H6N2O2S	001753-22-6	17
3	2-Chlorobenzimidazole	152	C7H5ClN2	004857-06-1	17
4	Thiophene, 2-butyltetrahydro-	144	C8H16S	001613-49-6	10
5	4-Amino-2-chlorobenzonitrile	152	C7H5ClN2	020925-27-3	9



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020916\
Data File : BG020825.D
Acq On : 9 Feb 2016 20:03
Operator : SJ/UM
Sample : H1371-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

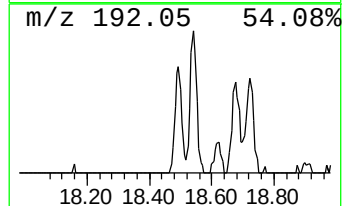
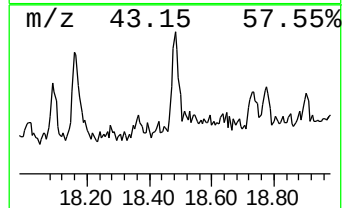
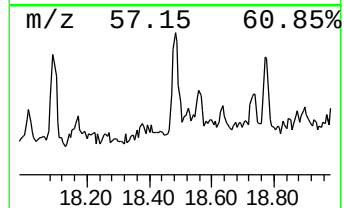
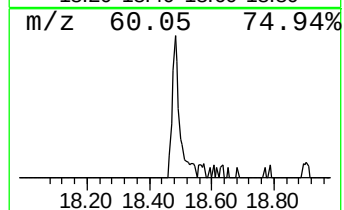
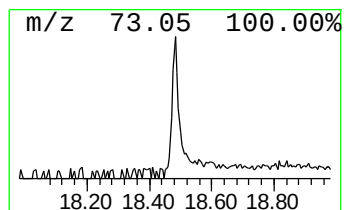
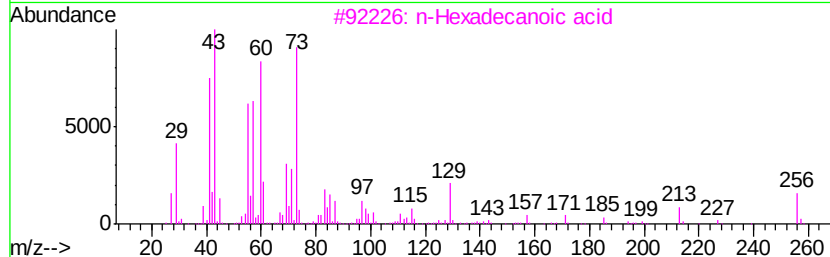
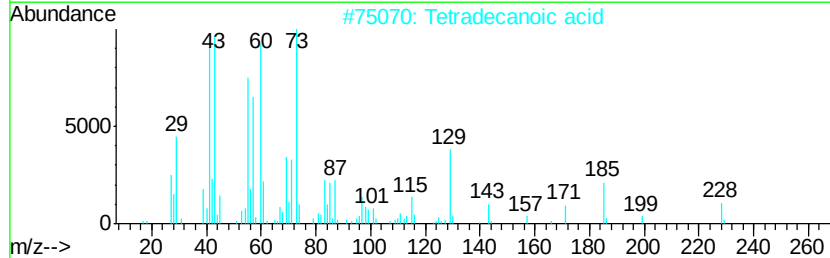
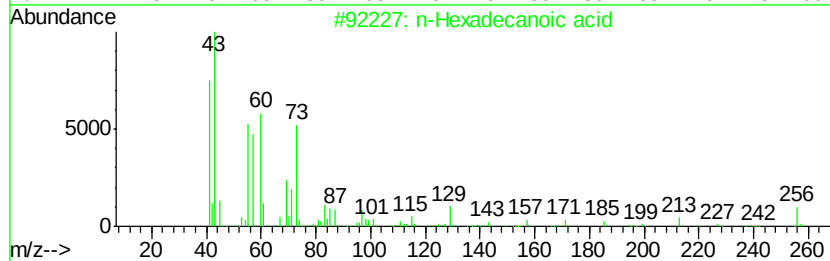
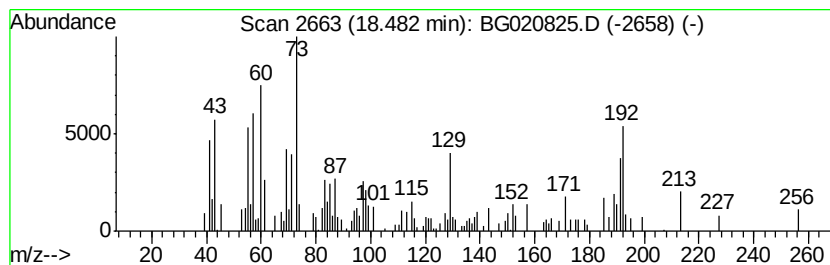
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG020816.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 n-Hexadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.48	3.61 ng/ul	61077	Phenanthrene-d10	17.62	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	43
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	43
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	38
5	.alpha.-D-Glucopyranoside, .alph...	342	C12H22O11	000099-20-7	35



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020916\
Data File : BG020825.D
Acq On : 9 Feb 2016 20:03
Operator : SJ/UM
Sample : H1371-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

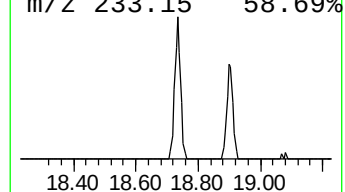
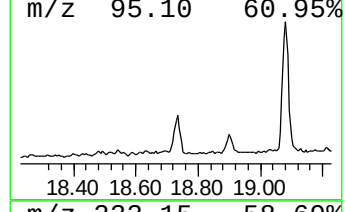
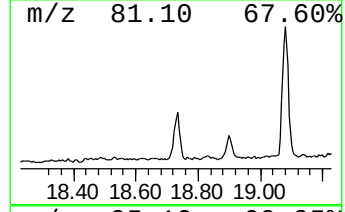
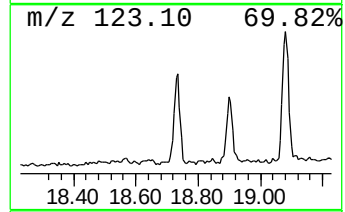
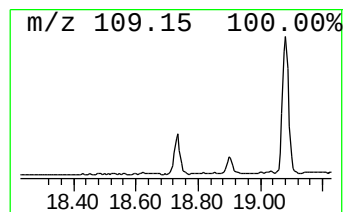
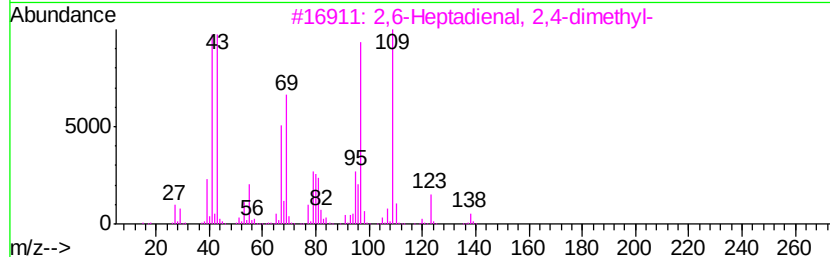
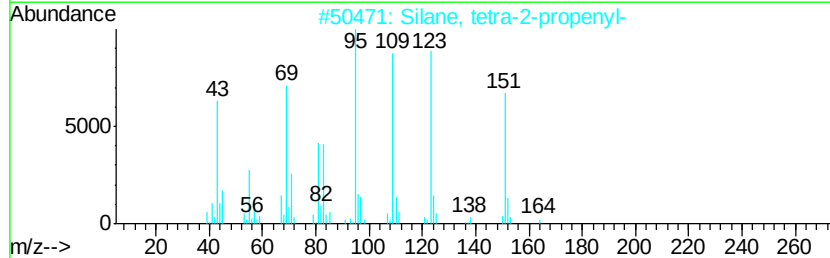
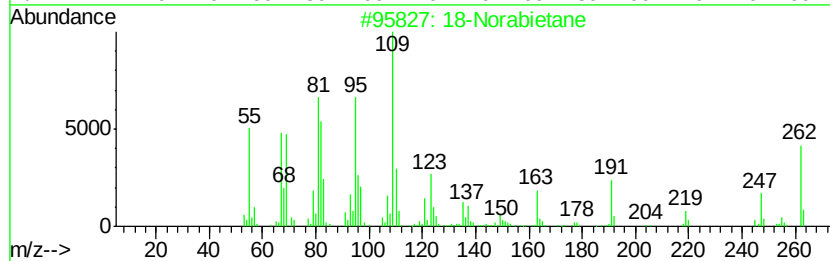
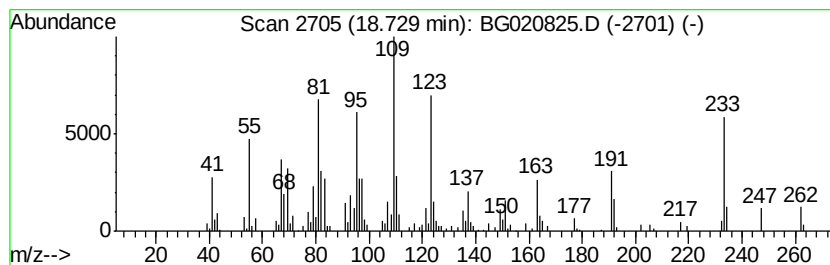
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG020816.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 18-Norabietane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.73	6.21 ng/ul	105056	Phenanthrene-d10	17.62	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	18-Norabietane	262	C19H34	1000293-16-6	83
2	Silane, tetra-2-propenyl-	192	C12H20Si	001112-66-9	35
3	2,6-Heptadienal, 2,4-dimethyl-	138	C9H14O	085136-08-9	22
4	1H-Pyrazole, 4-ethyl-3,5-dimethyl-	124	C7H12N2	007554-67-8	18
5	2,6-Bis(1,1-dimethylethyl)-4-(1-...	262	C17H26O2	014035-34-8	15



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020916\
Data File : BG020825.D
Acq On : 9 Feb 2016 20:03
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Sample : H1371-16
Misc :
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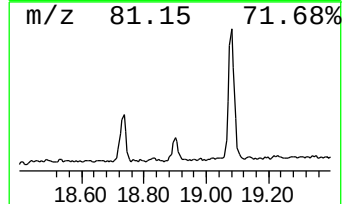
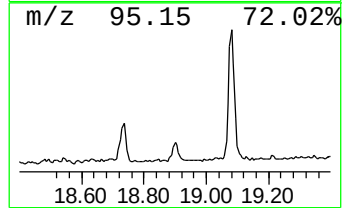
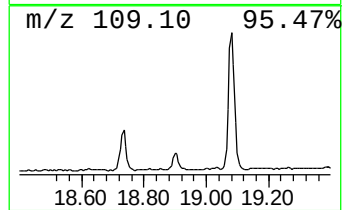
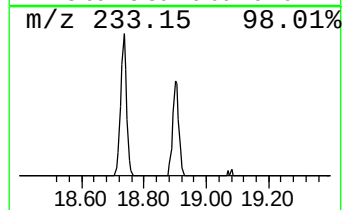
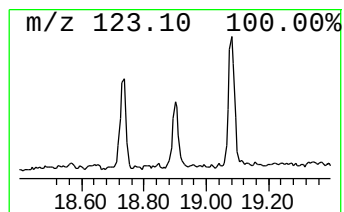
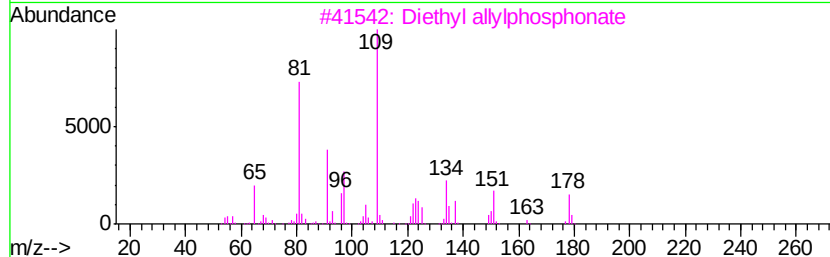
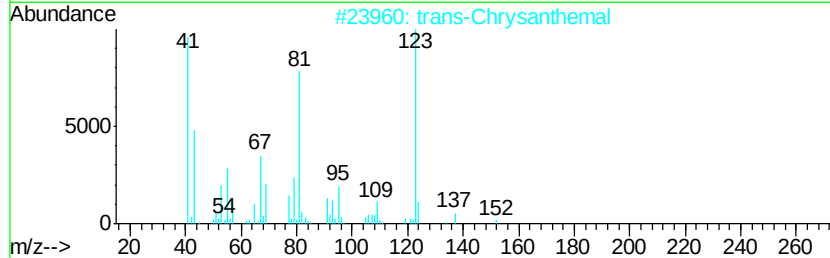
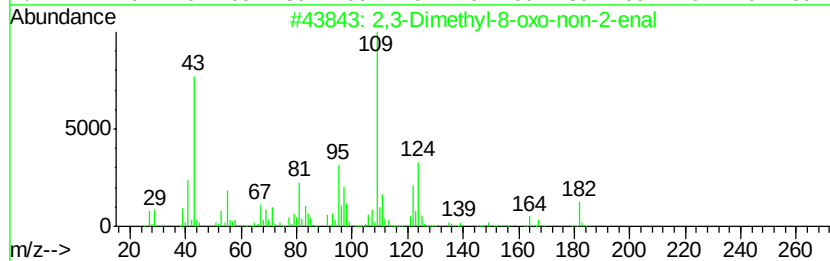
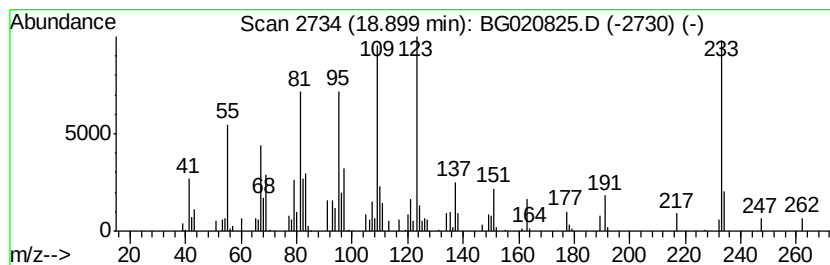
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 unknown-03 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.90	3.20 ng/ul	54226	Phenanthrene-d10	17.62

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3-Dimethyl-8-oxo-non-2-enal	182	C11H18O2	1000186-82-6	27		
2	trans-Chrysanthemal	152	C10H16O	020104-05-6	25		
3	Diethyl allvlphosphonate	178	C7H15O3P	001067-87-4	22		
4	Cyclohexene, 4-(4-ethylcyclohexyl)-	262	C19H34	301643-32-3	15		
5	4,4-Dimethoxy-2,5-cyclohexadien-	154	C8H10O3	000935-50-2	14		



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020916\
Data File : BG020825.D
Acq On : 9 Feb 2016 20:03
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Sample : H1371-16
Misc :
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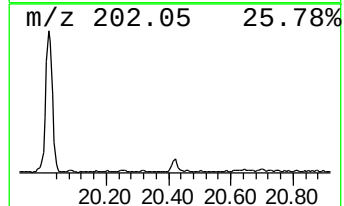
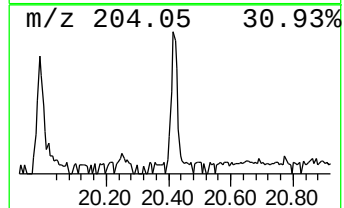
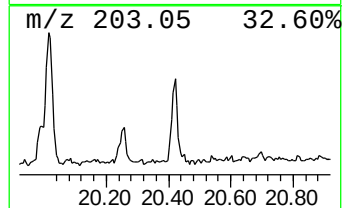
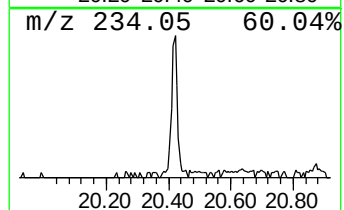
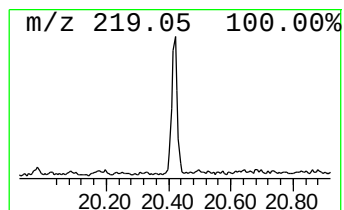
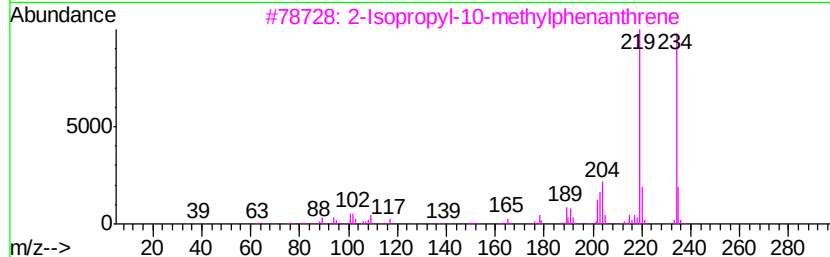
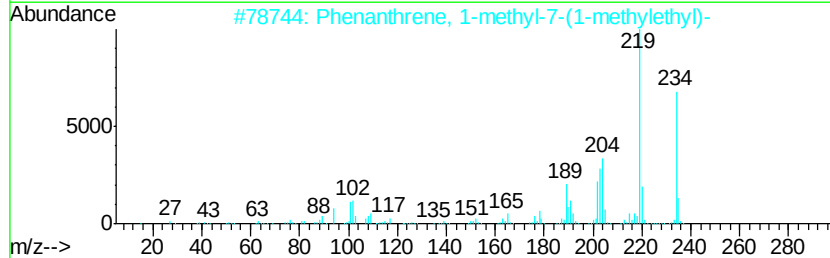
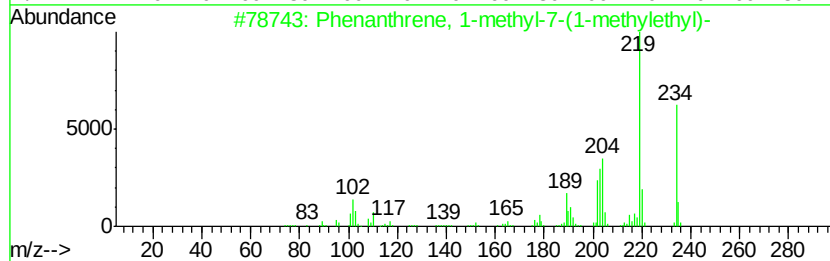
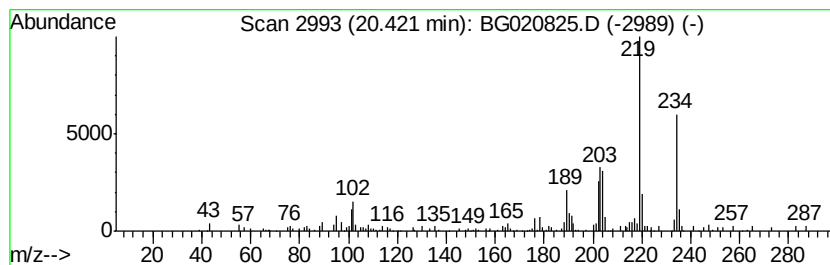
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Phenanthrene, 1-methyl-7-(1... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.42	2.84 ng/ul	47219	Chrysene-d12	21.91

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-7-(1-meth...	234	C18H18	000483-65-8	98
2			Phenanthrene, 1-methyl-7-(1-meth...	234	C18H18	000483-65-8	96
3			2-Isopropyl-10-methylphenanthrene	234	C18H18	066552-97-4	90
4			Phenanthrene, 2,4,5,7-tetramethyl-	234	C18H18	007396-38-5	64
5			1-(10-Methylantracen-9-yl)ethanone	234	C17H14O	036778-18-4	62



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020916\
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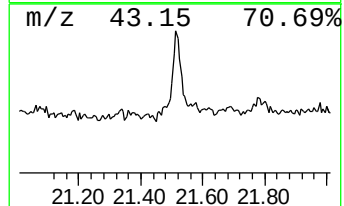
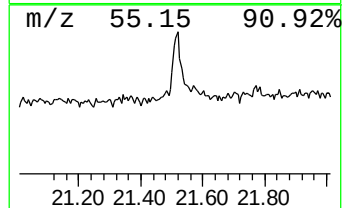
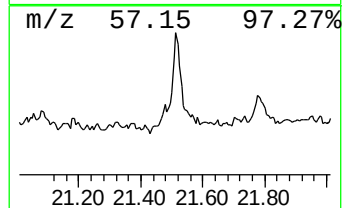
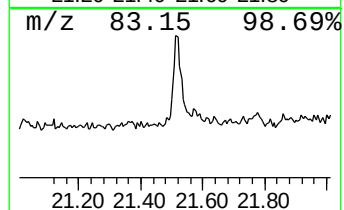
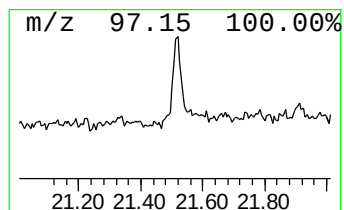
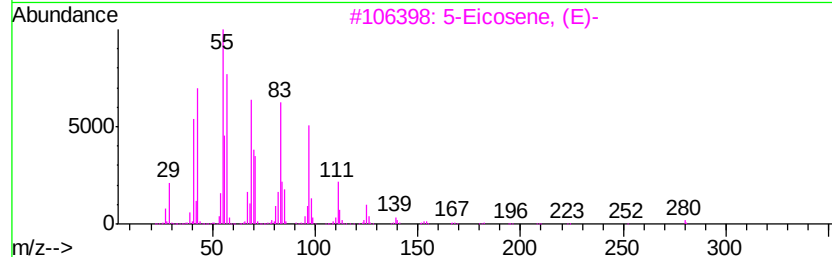
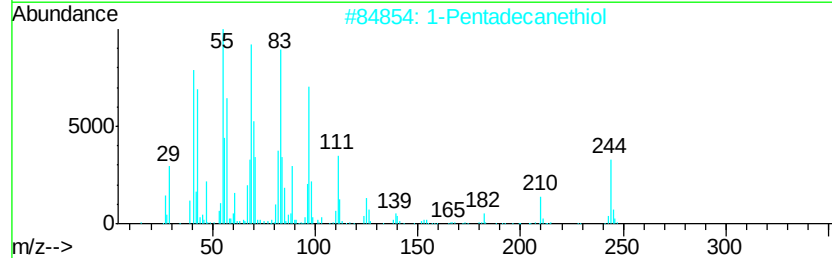
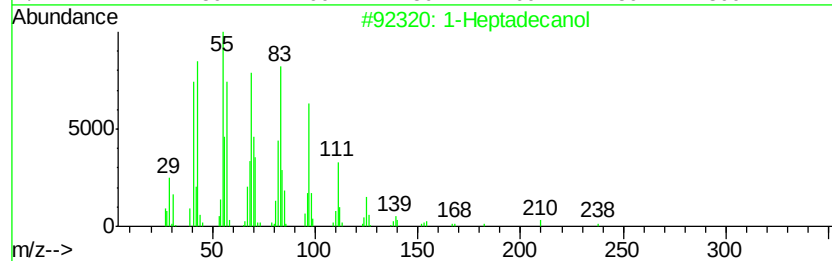
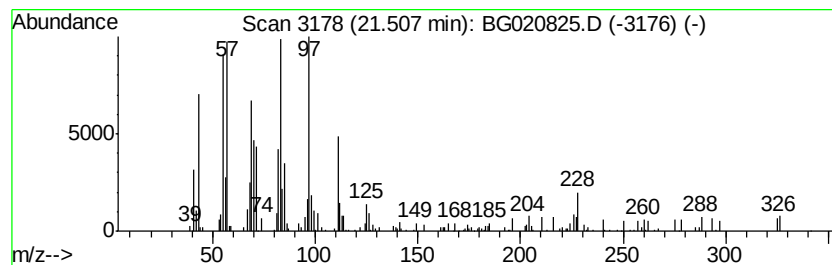
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 1-Heptadecanol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.51	4.22 ng/ul	70214	Chrysene-d12	21.91
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1-Heptadecanol	256 C17H36O	001454-85-9 90
2		1-Pentadecanethiol	244 C15H32S	025276-70-4 87
3		5-Eicosene, (E)-	280 C20H40	074685-30-6 83
4		Cycloeicosane	280 C20H40	000296-56-0 78
5		3-Eicosene, (E)-	280 C20H40	074685-33-9 78



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG020916\
Data File : BG020825.D
Acq On : 9 Feb 2016 20:03
Operator : SJ/UM
Sample : H1371-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C04J2

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG020816.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
Butane, 2-methoxy...	3.33	11.9	ng/ul	91084	1	8.27	153688	20.0
Benzene, 1,3,5-tr...	7.91	4.7	ng/ul	36079	1	8.27	153688	20.0
Benzene, 1,2,3-tr...	8.39	3.6	ng/ul	27437	1	8.27	153688	20.0
unknown-01	14.80	33.6	ng/ul	539582	3	14.89	320991	20.0
unknown-02	17.22	5.0	ng/ul	83796	4	17.62	338413	20.0
n-Hexadecanoic acid	18.48	3.6	ng/ul	61077	4	17.62	338413	20.0
18-Norabietane	18.73	6.2	ng/ul	105056	4	17.62	338413	20.0
unknown-03	18.90	3.2	ng/ul	54226	4	17.62	338413	20.0
Phenanthrene, 1-m...	20.42	2.8	ng/ul	47219	5	21.91	332591	20.0
1-Heptadecanol	21.51	4.2	ng/ul	70214	5	21.91	332591	20.0