

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
 Data File : BG046415.D
 Acq On : 21 Aug 2020 16:00
 Operator : CG/JU
 Sample : L3635-19DL 5X
 Misc :
 ALS Vial : 79 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BFMJ5DL

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-BG081320PAH.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	3.146	5	10	29	rVB	376479	637107	30.67%	3.814%
2	3.399	51	53	59	rVB4	3607	4565	0.22%	0.027%
3	3.922	137	142	149	rVB3	9794	16089	0.77%	0.096%
4	3.986	149	153	157	rVV5	1431	2646	0.13%	0.016%
5	4.045	161	163	171	rVB5	3846	6411	0.31%	0.038%
6	4.151	176	181	183	rVV2	2594	3814	0.18%	0.023%
7	5.056	331	335	341	rVB2	3255	4990	0.24%	0.030%
8	5.150	348	351	356	rVB4	2242	3535	0.17%	0.021%
9	7.118	679	686	694	rBV	54850	99983	4.81%	0.599%
10	7.271	706	712	722	rBV	43271	75502	3.63%	0.452%
11	7.482	741	748	758	rBV	58674	105382	5.07%	0.631%
12	7.941	819	826	836	rBV	330566	567675	27.32%	3.399%
13	8.651	940	947	958	rBV	69453	126963	6.11%	0.760%
14	8.763	962	966	972	rVB5	2256	4529	0.22%	0.027%
15	9.098	1016	1023	1032	rBV	54023	101196	4.87%	0.606%
16	9.821	1137	1146	1155	rBV	47273	84239	4.05%	0.504%
17	10.367	1232	1239	1251	rBV	74297	143138	6.89%	0.857%
18	10.737	1293	1302	1310	rBV	479822	860438	41.41%	5.151%
19	10.790	1310	1311	1317	rVB	7138	6111	0.29%	0.037%
20	10.872	1317	1325	1337	rBV3	39618	86060	4.14%	0.515%
21	12.165	1541	1545	1550	rVB5	1354	2313	0.11%	0.014%
22	12.329	1570	1573	1578	rVB5	1860	2123	0.10%	0.013%
23	12.400	1580	1585	1591	rVB5	3291	6138	0.30%	0.037%
24	12.617	1614	1622	1626	rBV2	3440	5055	0.24%	0.030%
25	12.941	1671	1677	1681	rVB5	1915	3341	0.16%	0.020%
26	13.181	1712	1718	1724	rBV5	1974	3917	0.19%	0.023%
27	13.405	1749	1756	1761	rBV7	3405	6956	0.33%	0.042%
28	13.751	1812	1815	1819	rVB4	1903	2486	0.12%	0.015%
29	13.892	1834	1839	1844	rBV5	4369	7785	0.37%	0.047%
30	13.975	1844	1853	1858	rBV	156065	226108	10.88%	1.354%
31	14.022	1858	1861	1866	rVB	25545	34754	1.67%	0.208%
32	14.133	1876	1880	1883	rVB2	2323	2872	0.14%	0.017%
33	14.263	1894	1902	1912	rBV	188425	313971	15.11%	1.880%
34	14.333	1912	1914	1918	rVB2	2496	2799	0.13%	0.017%

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35	14.480	1935	1939	1941	rBV3	2933	3529	0.17%	0.021%
36	14.574	1946	1955	1961	rBV	757863	1229091	59.16%	7.358%
37	14.633	1961	1965	1971	rVB2	23149	34820	1.68%	0.208%
38	14.780	1982	1990	1997	rBV2	30538	64799	3.12%	0.388%
39	14.968	2017	2022	2027	rVB3	9334	15851	0.76%	0.095%
40	15.050	2027	2036	2040	rVB6	3582	8040	0.39%	0.048%
41	15.197	2057	2061	2066	rVB5	2791	4666	0.22%	0.028%
42	15.244	2066	2069	2075	rBV5	2884	4660	0.22%	0.028%
43	15.338	2083	2085	2086	rBV	2249	2121	0.10%	0.013%
44	15.367	2087	2090	2093	rVV4	3265	4951	0.24%	0.030%
45	15.432	2097	2101	2105	rBV6	3900	5906	0.28%	0.035%
46	15.561	2116	2123	2129	rBV	241302	371003	17.86%	2.221%
47	15.614	2129	2132	2137	rVV3	16713	26650	1.28%	0.160%
48	15.690	2139	2145	2148	rVV3	7835	14049	0.68%	0.084%
49	15.743	2150	2154	2157	rVV5	7109	10360	0.50%	0.062%
50	15.784	2157	2161	2166	rVV7	7508	13607	0.65%	0.081%
51	15.837	2166	2170	2172	rVV5	5571	8032	0.39%	0.048%
52	15.867	2173	2175	2179	rVV5	3786	4597	0.22%	0.028%
53	15.908	2179	2182	2188	rVB5	6416	12605	0.61%	0.075%
54	15.996	2195	2197	2199	rVB2	2608	2438	0.12%	0.015%
55	16.025	2199	2202	2204	rBV3	3068	4871	0.23%	0.029%
56	16.060	2204	2208	2215	rVB3	7804	17364	0.84%	0.104%
57	16.125	2215	2219	2220	rBV4	3593	4019	0.19%	0.024%
58	16.295	2243	2248	2250	rVV4	7739	11831	0.57%	0.071%
59	16.384	2261	2263	2266	rBV4	2688	2080	0.10%	0.012%
60	16.431	2268	2271	2272	rBV3	3503	2955	0.14%	0.018%
61	16.495	2280	2282	2285	rBV4	2177	2549	0.12%	0.015%
62	16.530	2285	2288	2291	rBV5	2923	3905	0.19%	0.023%
63	16.572	2293	2295	2296	rBV2	2909	2888	0.14%	0.017%
64	16.666	2309	2311	2312	rBV2	4119	3250	0.16%	0.019%
65	16.830	2335	2339	2340	rBV3	6063	7298	0.35%	0.044%
66	16.842	2340	2341	2343	rVV2	4982	4853	0.23%	0.029%
67	16.895	2346	2350	2354	rBV7	10375	16510	0.79%	0.099%
68	16.965	2358	2362	2371	rVB	25214	42912	2.07%	0.257%
69	17.047	2371	2376	2378	rBV5	10732	15969	0.77%	0.096%
70	17.318	2415	2422	2426	rBV	901065	1435655	69.10%	8.595%
71	17.359	2426	2429	2433	rVB	275030	352749	16.98%	2.112%

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72	17.412	2433	2438	2443	rBV2	238377	366823	17.66%	2.196%
73	17.500	2450	2453	2459	rBV4	16030	30935	1.49%	0.185%
74	17.717	2487	2490	2494	rVB	18685	22535	1.08%	0.135%
75	17.835	2506	2510	2515	rVB4	19318	31891	1.53%	0.191%
76	18.193	2567	2571	2575	rBV	43365	61223	2.95%	0.367%
77	18.240	2575	2579	2582	rVV	56964	78082	3.76%	0.467%
78	18.281	2582	2586	2590	rVB	55638	65492	3.15%	0.392%
79	18.381	2599	2603	2608	rBV2	60813	120315	5.79%	0.720%
80	18.422	2608	2610	2614	rVB2	35800	39373	1.90%	0.236%
81	18.687	2651	2655	2658	rBV	48700	68517	3.30%	0.410%
82	18.728	2659	2662	2673	rVB2	58009	93153	4.48%	0.558%
83	19.104	2723	2726	2731	rBV4	44603	69447	3.34%	0.416%
84	19.368	2766	2771	2776	rVB	578606	771593	37.14%	4.619%
85	19.697	2822	2827	2829	rBV2	400510	573040	27.58%	3.431%
86	19.727	2829	2832	2841	rVB	604942	878157	42.27%	5.257%
87	21.566	3138	3145	3150	rBV2	1032463	2077601	100.00%	12.438%
88	21.607	3150	3152	3163	rVB	346144	541916	26.08%	3.244%
89	23.699	3502	3508	3516	rBV	323164	1066033	51.31%	6.382%
90	24.539	3646	3651	3662	rVB	434432	1037381	49.93%	6.211%
91	24.686	3669	3676	3685	rBV2	517563	1375137	66.19%	8.233%

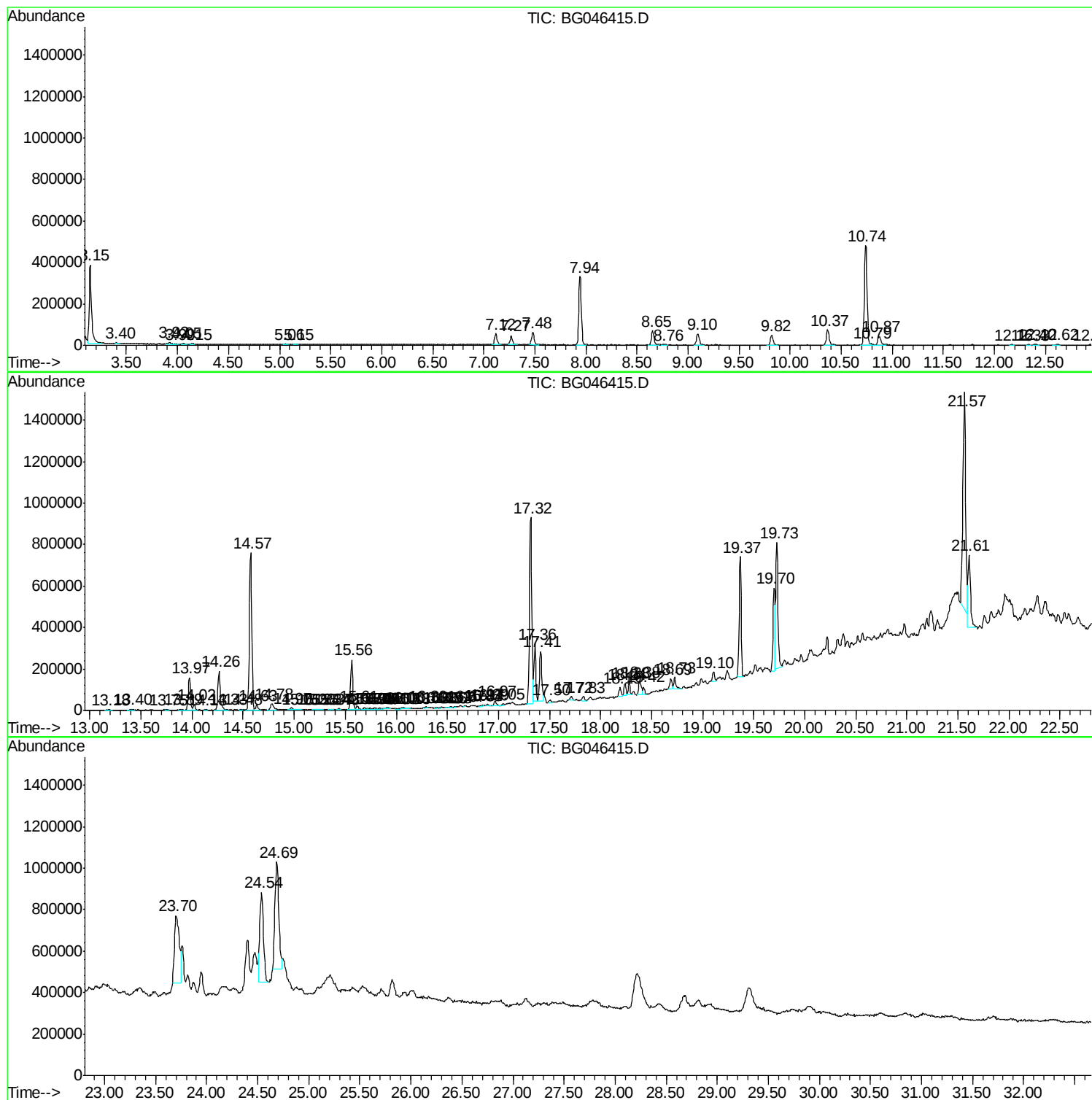
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG081320PAH.M
Quant Title : SVOA CALIBRATION

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



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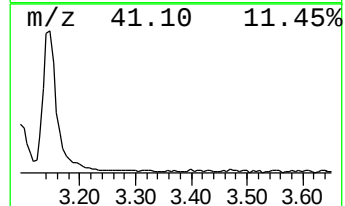
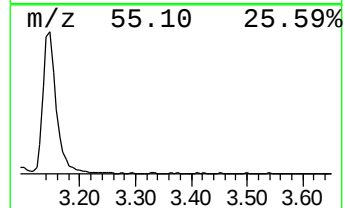
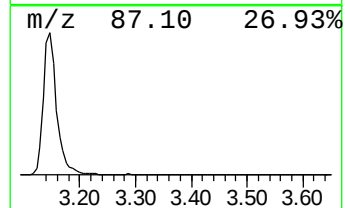
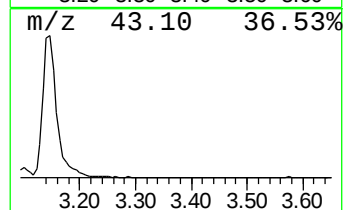
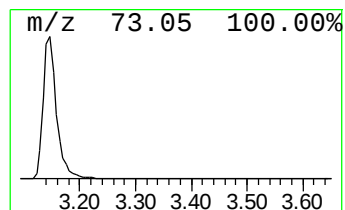
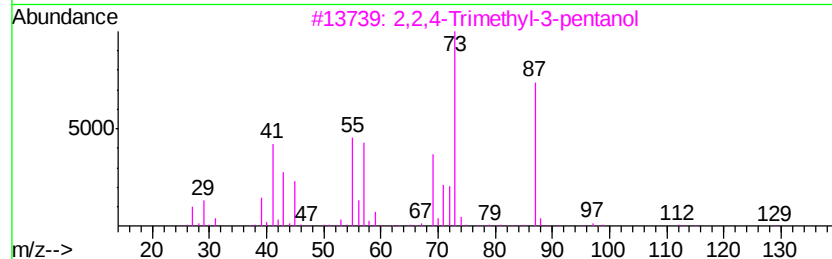
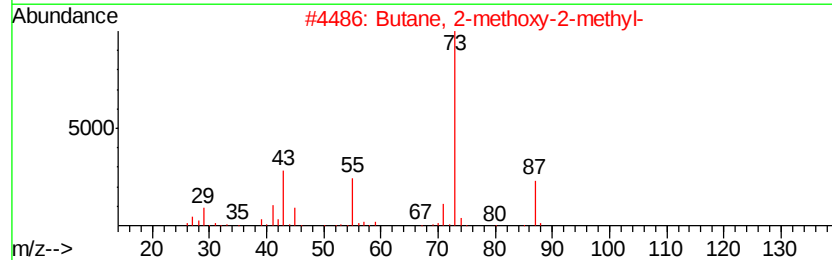
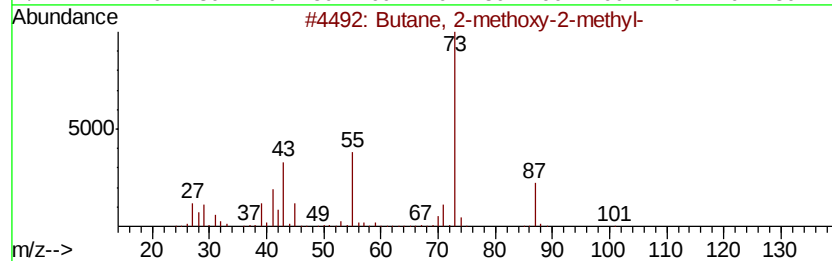
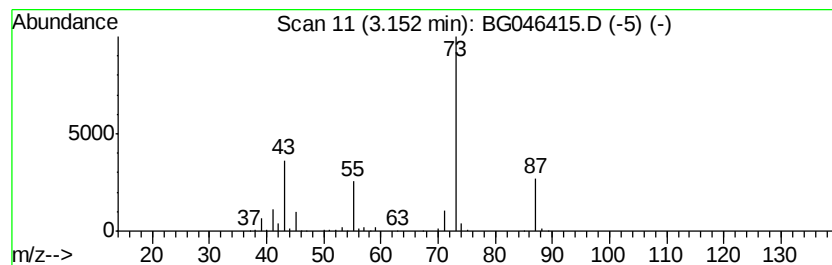
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG081320PAH.M
Quant Title : SVOA CALIBRATION

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	22.45 ng/ul	637107	1,4-Dichlorobenzene-d4	7.94

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	78
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	27
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17



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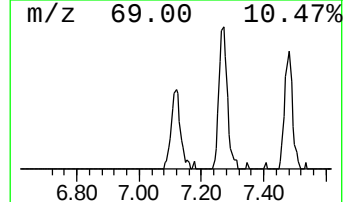
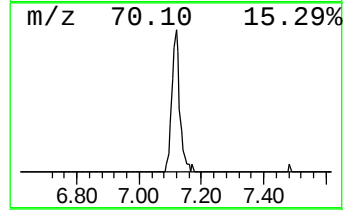
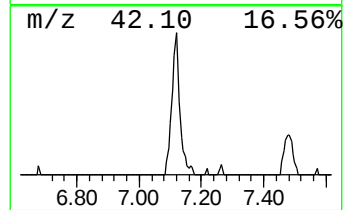
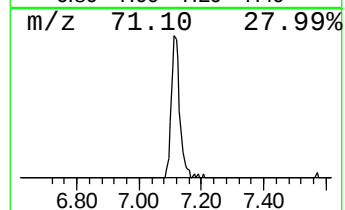
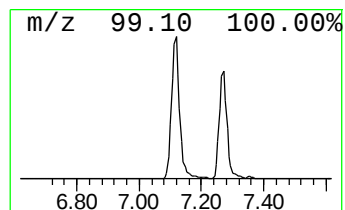
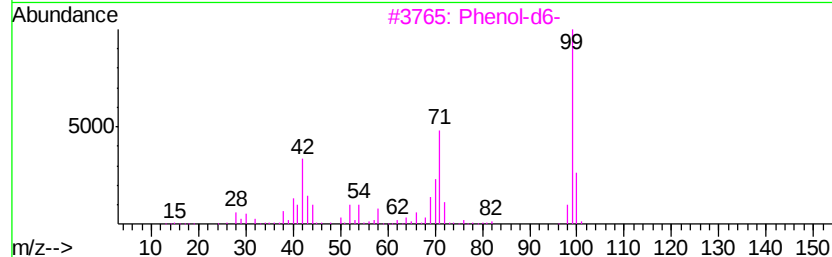
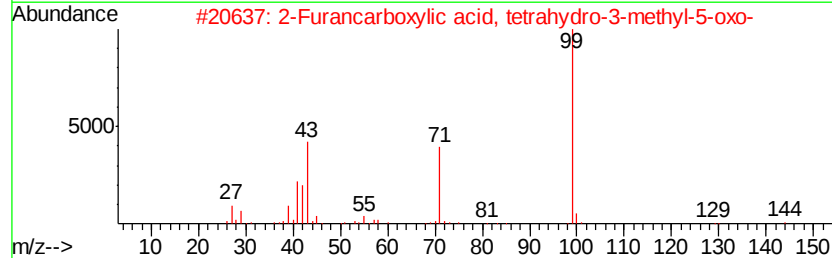
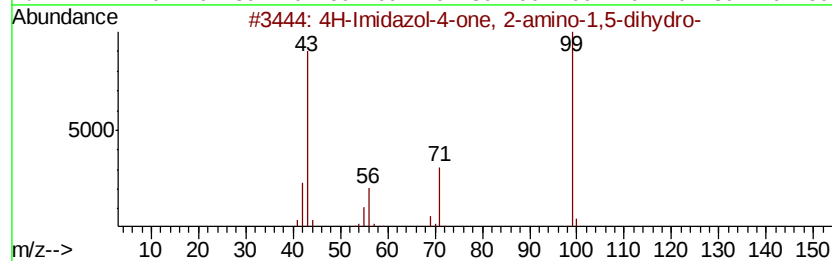
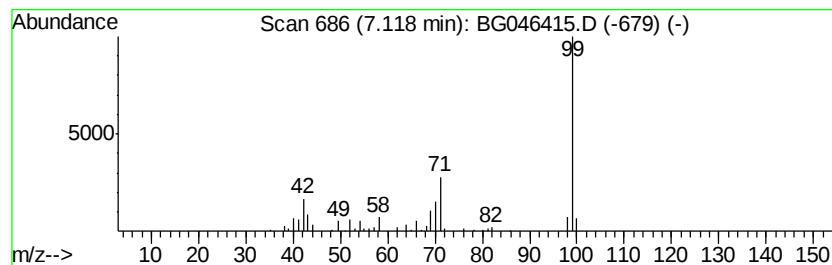
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG081320PAH.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 4H-Imidazol-4-one, 2-amino-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.12	3.52 ng/ul	99983	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Imidazol-4-one, 2-amino-1,5-d...	99	C3H5N3O	000503-86-6	72
2		2-Furancarboxylic acid, tetrahyd...	144	C6H8O4	022073-04-7	59
3		Phenol-d6-	100	C6D6O	013127-88-3	56
4		n-Heptyl methylphosphonofluoridate	196	C8H18F02P	162085-82-7	53
5		Hexanoic acid, anhydride	214	C12H22O3	002051-49-2	50



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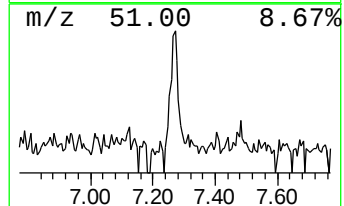
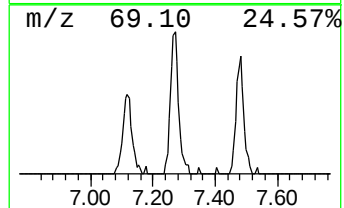
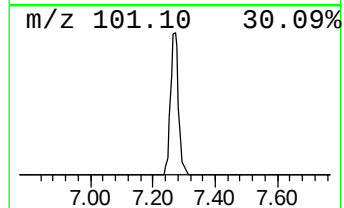
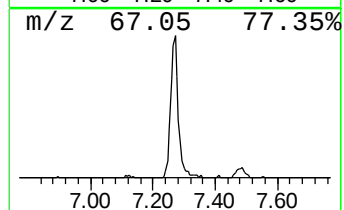
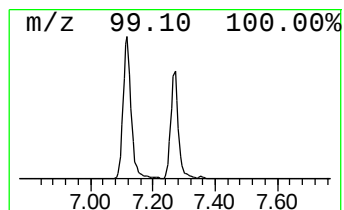
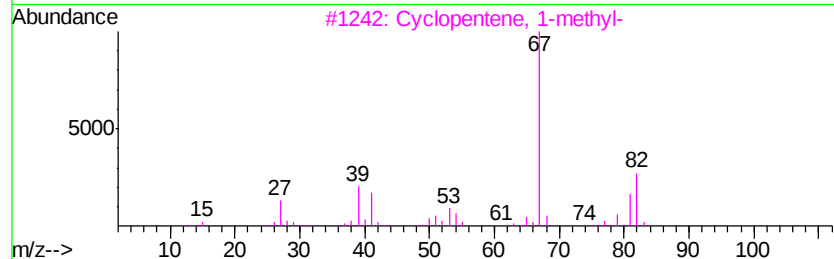
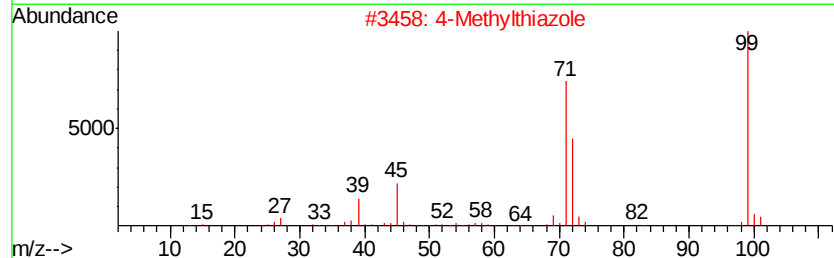
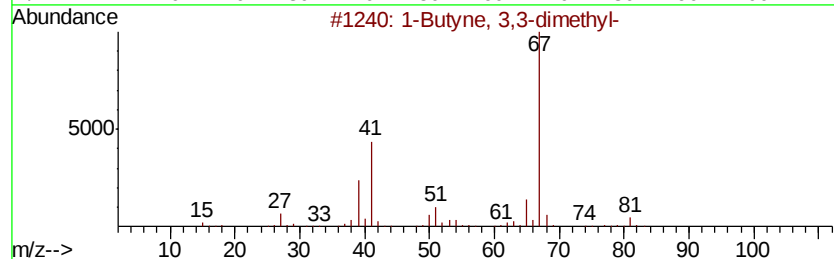
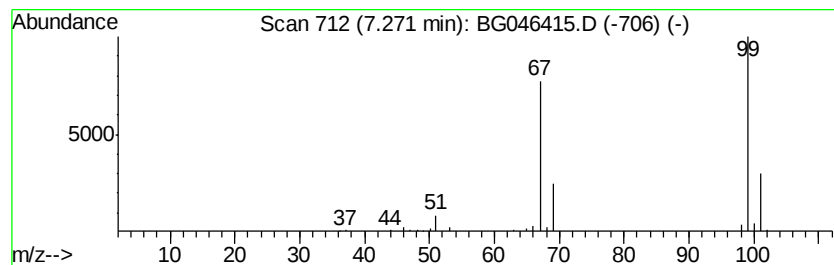
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG081320PAH.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.27	2.66 ng/ul	75502	1,4-Dichlorobenzene-d4	7.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Butyne, 3,3-dimethyl-	82	C6H10	000917-92-0	12
2			4-Methylthiazole	99	C4H5NS	000693-95-8	9
3			Cyclopentene, 1-methyl-	82	C6H10	000693-89-0	9
4			Cyclopentene, 3-methyl-	82	C6H10	001120-62-3	9
5			Thiazole, 5-methyl-	99	C4H5NS	003581-89-3	7



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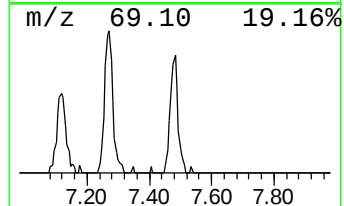
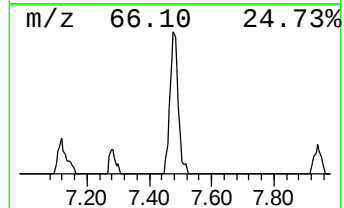
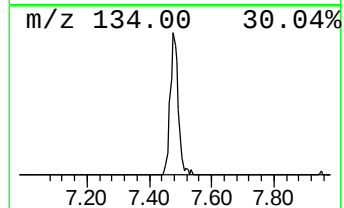
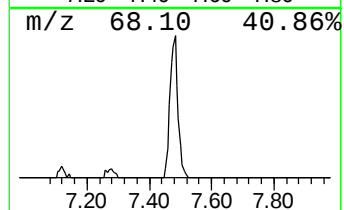
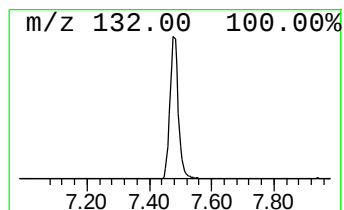
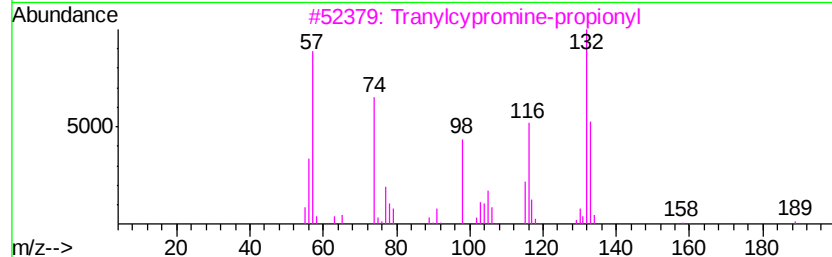
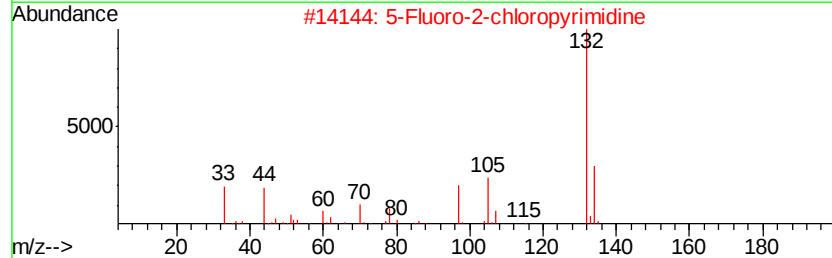
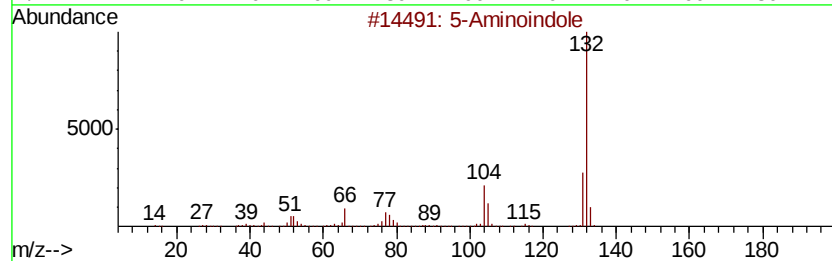
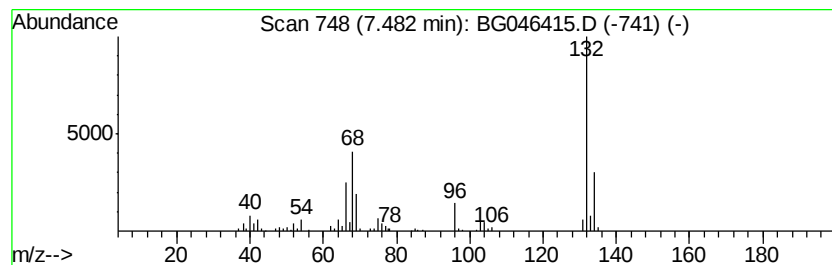
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-02 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.48	3.71 ng/ul	105382	1,4-Dichlorobenzene-d4	7.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Aminoindole	132	C8H8N2	005192-03-0	22
2			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3			Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5			Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046415.D
Acq On : 21 Aug 2020 16:00
Operator : CG/JU
Sample : L3635-19DL 5X
Misc :
ALS Vial : 79 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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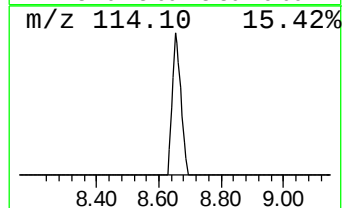
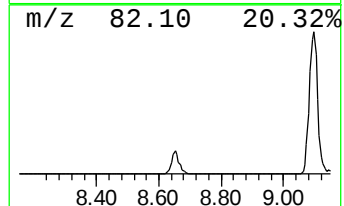
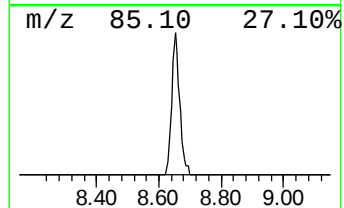
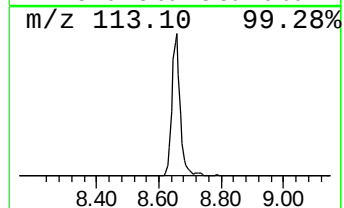
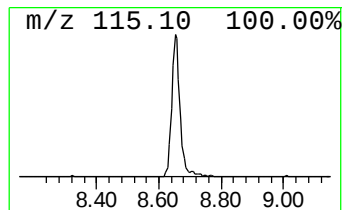
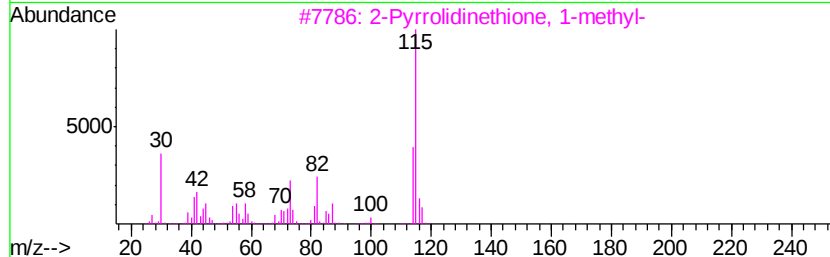
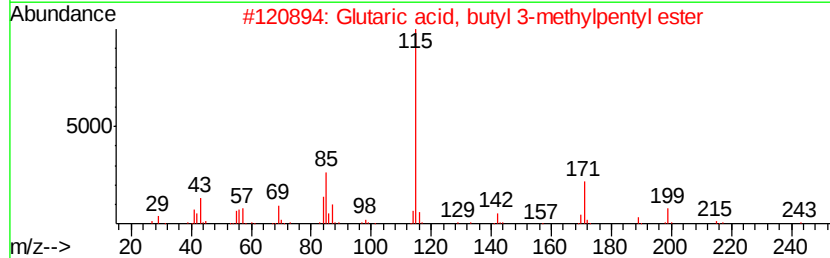
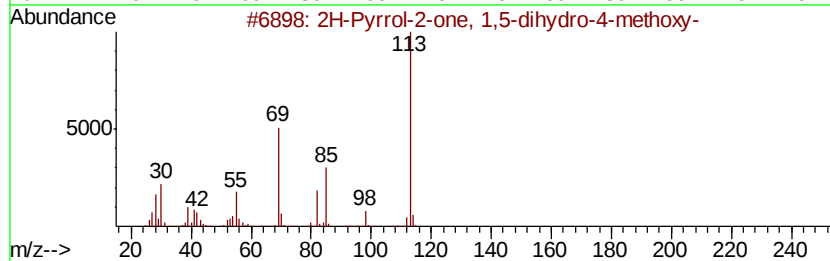
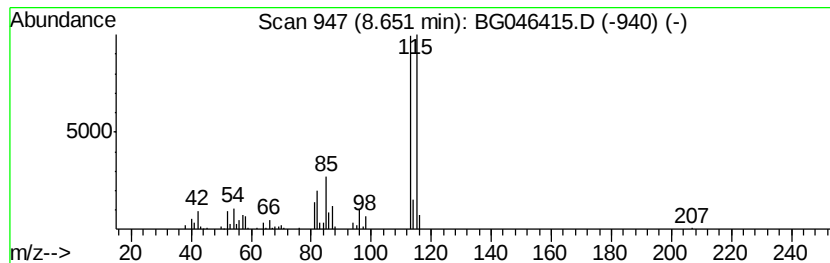
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-03 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.65	4.47 ng/ul	126963	1,4-Dichlorobenzene-d4	7.94

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2H-Pyrrol-2-one, 1,5-dihydro-4-m...	113	C5H7NO2	069778-83-2	38
2		Glutaric acid, butyl 3-methylpen...	272	C15H28O4	1000360-05-7	38
3		2-Pyrrolidinethione, 1-methyl-	115	C5H9NS	010441-57-3	38
4		Glutaric acid, 3,3-dimethylbut-2...	272	C15H28O4	1000359-74-9	38
5		Glutaric acid, di(2-methylpentyl...	300	C17H32O4	1000358-42-8	37



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046415.D
Acq On : 21 Aug 2020 16:00
Operator : CG/JU
Sample : L3635-19DL 5X
Misc :
ALS Vial : 79 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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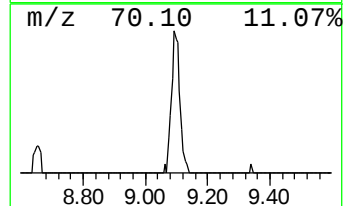
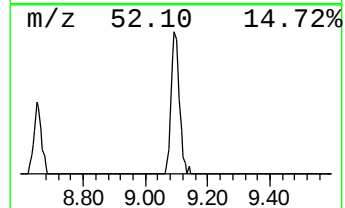
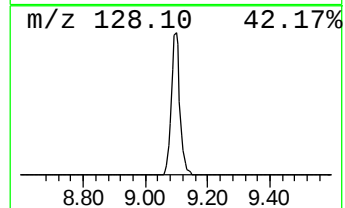
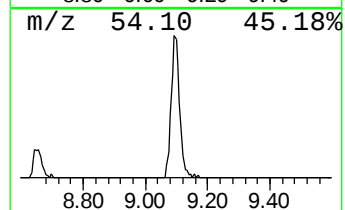
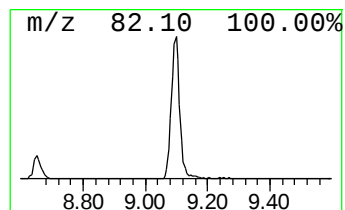
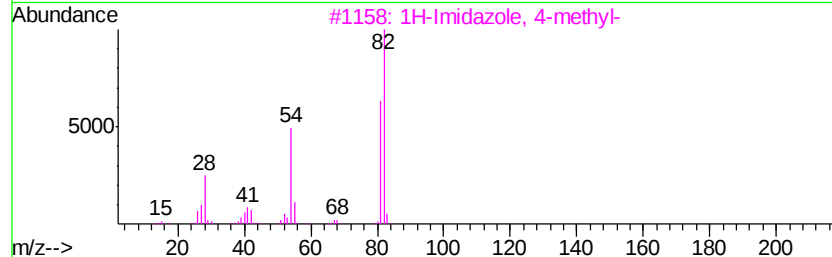
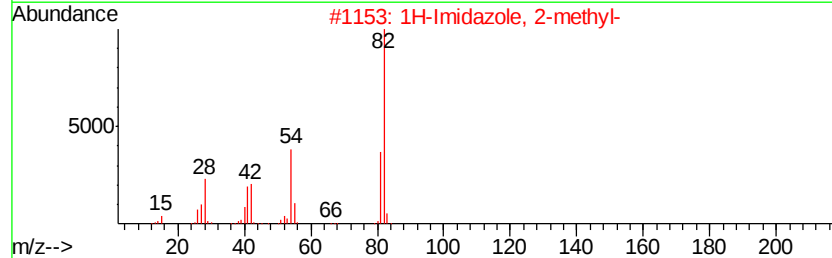
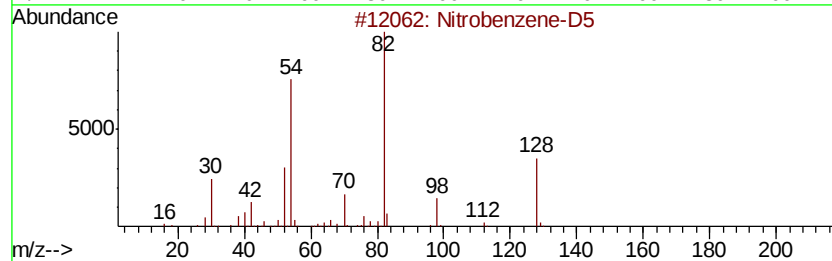
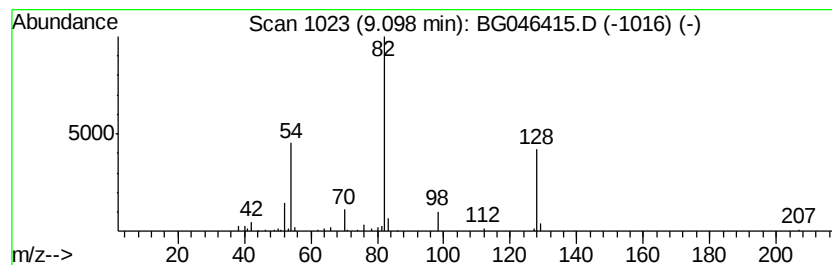
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-04 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.10	3.57 ng/ul	101196	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nitrobenzene-D5	128	C6D5NO2	004165-60-0	47
2		1H-Imidazole, 2-methyl-	82	C4H6N2	000693-98-1	38
3		1H-Imidazole, 4-methyl-	82	C4H6N2	000822-36-6	38
4		1H-Imidazole, 2-methyl-	82	C4H6N2	000693-98-1	28
5		2-Cyclohexen-1-one, 3-methyl-	110	C7H10O	001193-18-6	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046415.D
Acq On : 21 Aug 2020 16:00
Operator : CG/JU
Sample : L3635-19DL 5X
Misc :
ALS Vial : 79 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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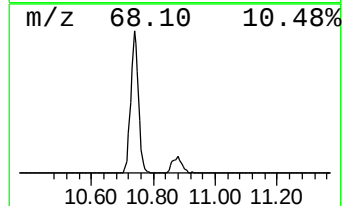
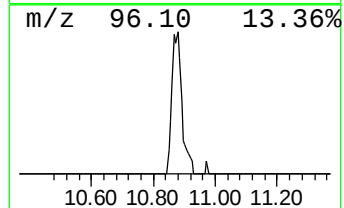
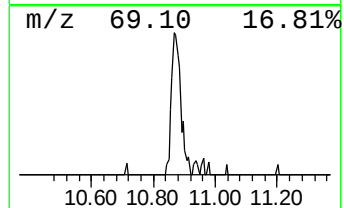
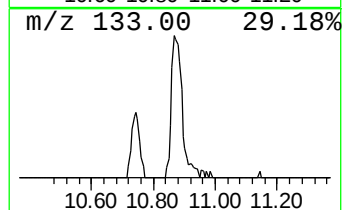
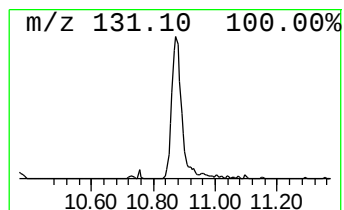
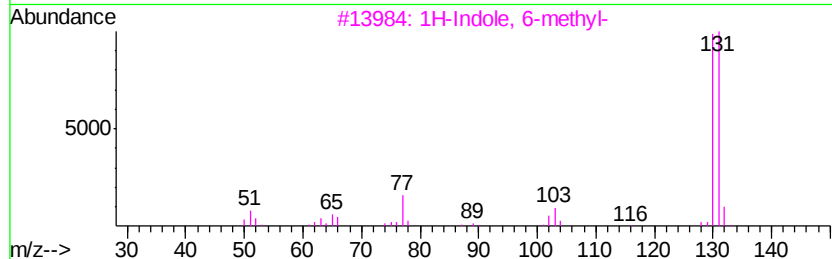
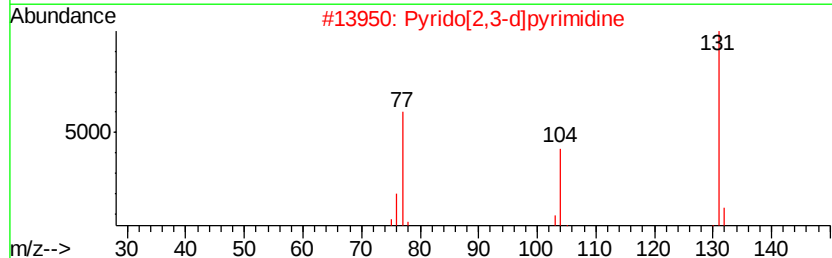
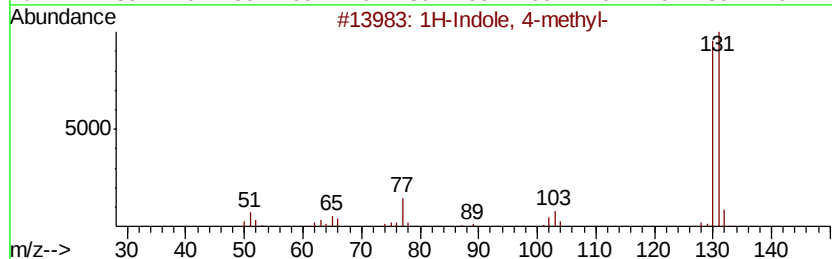
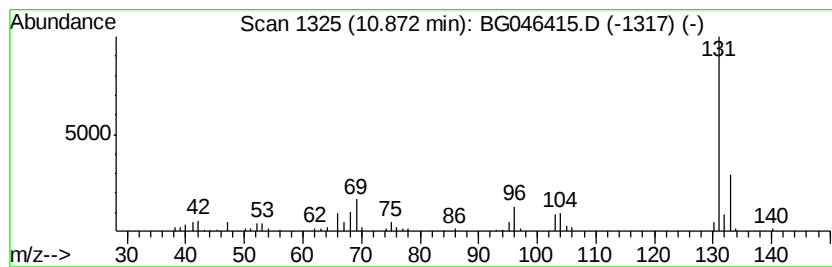
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-05 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.87	2.00 ng/ul	86060	Naphthalene-d8	10.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indole, 4-methyl-	131	C9H9N	016096-32-5	40
2		Pyrido[2,3-d]pyrimidine	131	C7H5N3	000254-61-5	40
3		1H-Indole, 6-methyl-	131	C9H9N	003420-02-8	33
4		1H-Indole, 5-methyl-	131	C9H9N	000614-96-0	33
5		1H-Indole, 1-methyl-	131	C9H9N	000603-76-9	28



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046415.D
Acq On : 21 Aug 2020 16:00
Operator : CG/JU
Sample : L3635-19DL 5X
Misc :
ALS Vial : 79 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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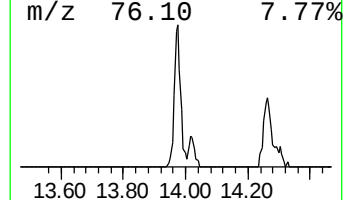
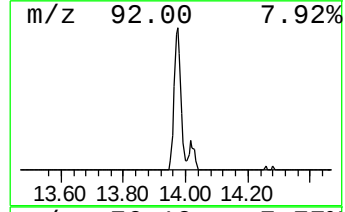
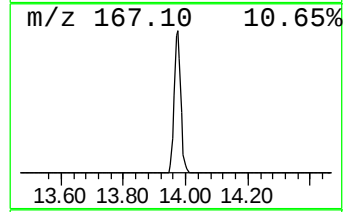
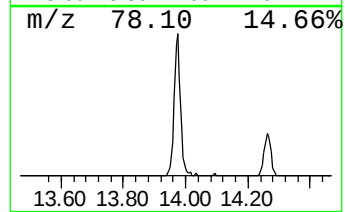
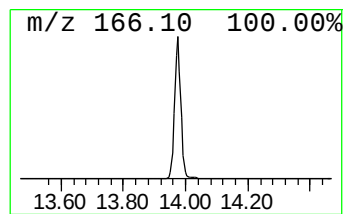
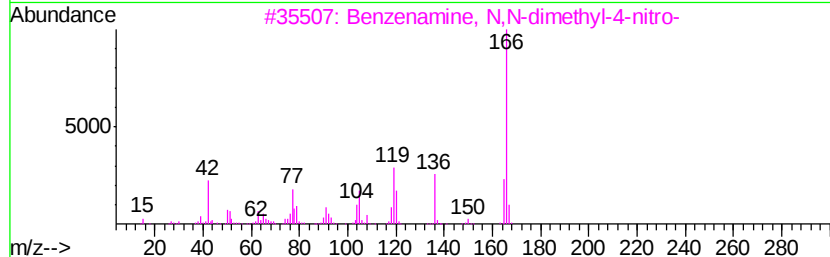
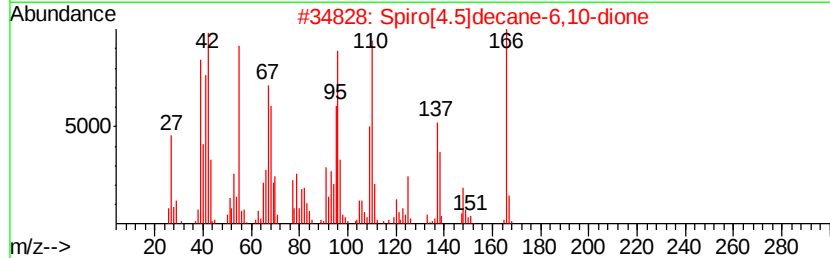
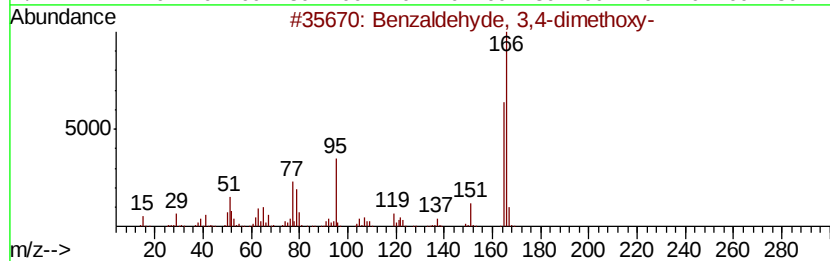
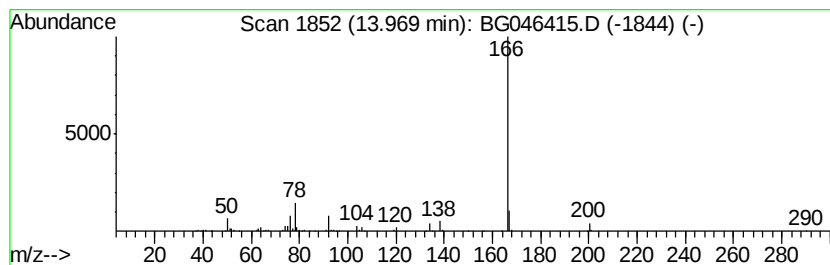
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-06 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	3.68 ng/ul	226108	Acenaphthene-d10	14.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzaldehyde, 3,4-dimethoxy-	166	C9H10O3	000120-14-9	45
2		Spiro[4.5]decane-6,10-dione	166	C10H14O2	006684-66-8	40
3		Benzenamine, N,N-dimethyl-4-nitro-	166	C8H10N2O2	000100-23-2	38
4		6H-Purine-6-thione, 1,7-dihydro-...	166	C6H6N4S	001126-23-4	9
5		Benzo[b]tetrahydrofuran-3-one, 5...	166	C8H6O4	014771-00-7	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG081920\
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Misc :
ALS Vial : 79 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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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	22.4	ng/ul	637107	1	7.94	567675	20.0
4H-Imidazol-4-one...	7.12	3.5	ng/ul	99983	1	7.94	567675	20.0
unknown-01	7.27	2.7	ng/ul	75502	1	7.94	567675	20.0
unknown-02	7.48	3.7	ng/ul	105382	1	7.94	567675	20.0
unknown-03	8.65	4.5	ng/ul	126963	1	7.94	567675	20.0
unknown-04	9.10	3.6	ng/ul	101196	1	7.94	567675	20.0
unknown-05	10.87	2.0	ng/ul	86060	2	10.74	860438	20.0
unknown-06	13.97	3.7	ng/ul	226108	3	14.57	1229090	20.0