

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081720\
 Data File : BG046336.D
 Acq On : 18 Aug 2020 20:33
 Operator : CG/JU
 Sample : L3636-16
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BFMN7

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	3.141	4	9	36	rVB	1291103	2092331	100.00%	15.450%
2	3.917	138	141	147	rVB6	4580	5566	0.27%	0.041%
3	4.052	159	164	174	rVB	12270	21925	1.05%	0.162%
4	4.140	176	179	184	rVB3	4381	5455	0.26%	0.040%
5	5.051	329	334	339	rBV	10739	16940	0.81%	0.125%
6	7.107	678	684	694	rBV2	31935	66445	3.18%	0.491%
7	7.266	704	711	721	rBV	35166	58560	2.80%	0.432%
8	7.477	741	747	755	rBV	38488	67960	3.25%	0.502%
9	7.583	758	765	768	rVB5	2897	6031	0.29%	0.045%
10	7.941	818	826	836	rBV	298719	505878	24.18%	3.735%
11	8.646	938	946	954	rBV2	36723	69084	3.30%	0.510%
12	9.093	1014	1022	1032	rBV	37088	70019	3.35%	0.517%
13	9.816	1139	1145	1157	rVB2	28043	53899	2.58%	0.398%
14	10.362	1231	1238	1247	rBV	38528	74915	3.58%	0.553%
15	10.738	1293	1302	1315	rBV	419704	760205	36.33%	5.613%
16	10.867	1317	1324	1335	rVB2	34085	67776	3.24%	0.500%
17	12.612	1616	1621	1624	rBV2	2403	3662	0.18%	0.027%
18	12.924	1672	1674	1680	rBV4	1995	3379	0.16%	0.025%
19	13.182	1712	1718	1721	rBV5	2413	3742	0.18%	0.028%
20	13.470	1761	1767	1776	rVV	90207	142346	6.80%	1.051%
21	13.970	1845	1852	1856	rBV	101917	151407	7.24%	1.118%
22	14.017	1856	1860	1871	rVB	51868	76871	3.67%	0.568%
23	14.257	1894	1901	1911	rBV	106462	167828	8.02%	1.239%
24	14.569	1946	1954	1961	rVV	710048	1111406	53.12%	8.207%
25	14.628	1961	1964	1970	rVB3	10298	16587	0.79%	0.122%
26	14.775	1984	1989	1994	rBV3	9160	19685	0.94%	0.145%
27	14.880	2003	2007	2012	rVB4	2327	3093	0.15%	0.023%
28	14.963	2017	2021	2027	rVB2	6836	11617	0.56%	0.086%
29	15.386	2086	2093	2094	rBV4	2593	4190	0.20%	0.031%
30	15.556	2116	2122	2128	rBV	152184	232587	11.12%	1.717%
31	15.615	2128	2132	2138	rVB2	11067	14909	0.71%	0.110%
32	15.697	2138	2146	2151	rBV5	40729	82777	3.96%	0.611%
33	15.908	2178	2182	2189	rBV2	4630	8062	0.39%	0.060%
34	16.055	2204	2207	2214	rBV3	3994	8496	0.41%	0.063%

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 Title : SVOA CALIBRATION

35	16.602	2290	2300	2301	rBV6	2414	4573	0.22%	0.034%
36	16.614	2301	2302	2308	rVB4	2716	3190	0.15%	0.024%
37	16.854	2336	2343	2345	rBV4	2403	4707	0.22%	0.035%
38	16.890	2345	2349	2351	rBV2	2543	2952	0.14%	0.022%
39	16.966	2357	2362	2367	rBV	9614	13517	0.65%	0.100%
40	17.048	2370	2376	2377	rBV4	3405	5524	0.26%	0.041%
41	17.066	2377	2379	2381	rVV2	3520	3536	0.17%	0.026%
42	17.101	2381	2385	2386	rVV3	3856	5482	0.26%	0.040%
43	17.125	2386	2389	2393	rVB	5443	7565	0.36%	0.056%
44	17.307	2413	2420	2425	rBV	863226	1350510	64.55%	9.972%
45	17.348	2425	2427	2432	rVV	132485	178104	8.51%	1.315%
46	17.407	2432	2437	2441	rVV2	154853	230791	11.03%	1.704%
47	17.436	2441	2442	2450	rVB	22128	26638	1.27%	0.197%
48	17.495	2450	2452	2457	rBV3	5120	6675	0.32%	0.049%
49	17.624	2470	2474	2479	rVB6	10352	13310	0.64%	0.098%
50	17.706	2479	2488	2492	rBV2	8822	18857	0.90%	0.139%
51	17.830	2505	2509	2515	rVB6	4371	8655	0.41%	0.064%
52	17.894	2515	2520	2524	rBV6	4756	8812	0.42%	0.065%
53	17.988	2533	2536	2540	rBV4	2643	3306	0.16%	0.024%
54	18.112	2552	2557	2558	rVB3	3206	3846	0.18%	0.028%
55	18.188	2564	2570	2573	rBV2	18487	31410	1.50%	0.232%
56	18.235	2573	2578	2582	rVV	29189	45726	2.19%	0.338%
57	18.276	2582	2585	2587	rVV	4388	5664	0.27%	0.042%
58	18.312	2589	2591	2597	rVB2	8217	10867	0.52%	0.080%
59	18.382	2597	2603	2606	rBV2	23368	44644	2.13%	0.330%
60	18.417	2607	2609	2617	rVB	15495	17424	0.83%	0.129%
61	18.629	2641	2645	2649	rVB6	2659	3910	0.19%	0.029%
62	18.682	2649	2654	2657	rBV	17688	25761	1.23%	0.190%
63	18.717	2657	2660	2665	rVB2	19717	27227	1.30%	0.201%
64	18.776	2668	2670	2673	rBV4	2474	3495	0.17%	0.026%
65	18.934	2692	2697	2700	rBV6	6960	10925	0.52%	0.081%
66	18.981	2700	2705	2708	rVB5	6187	8823	0.42%	0.065%
67	19.011	2708	2710	2715	rBV5	5355	7861	0.38%	0.058%
68	19.064	2715	2719	2722	rBV4	17293	27347	1.31%	0.202%
69	19.093	2722	2724	2729	rVB4	13235	19897	0.95%	0.147%
70	19.199	2738	2742	2744	rVB4	5599	6359	0.30%	0.047%
71	19.234	2744	2748	2753	rBV	16494	24791	1.18%	0.183%

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72	19.304	2757	2760	2762	rBV3	4905	5992	0.29%	0.044%
73	19.357	2762	2769	2778	rBV	210504	307124	14.68%	2.268%
74	19.457	2784	2786	2790	rVB5	7255	7764	0.37%	0.057%
75	19.510	2791	2795	2798	rBV4	6260	9711	0.46%	0.072%
76	19.557	2798	2803	2804	rBV5	8121	10748	0.51%	0.079%
77	19.598	2808	2810	2813	rVB2	5339	6557	0.31%	0.048%
78	19.692	2820	2826	2828	rBV2	227453	325599	15.56%	2.404%
79	19.716	2828	2830	2839	rVB	166429	233149	11.14%	1.722%
80	19.798	2839	2844	2847	rBV4	12627	25992	1.24%	0.192%
81	19.898	2857	2861	2866	rBV	14616	25826	1.23%	0.191%
82	19.957	2868	2871	2876	rVB3	11574	17484	0.84%	0.129%
83	20.027	2879	2883	2892	rVV5	46940	87474	4.18%	0.646%
84	20.180	2905	2909	2910	rBV3	12696	16880	0.81%	0.125%
85	20.209	2911	2914	2918	rVB	30532	44866	2.14%	0.331%
86	20.309	2928	2931	2935	rBV3	13440	21128	1.01%	0.156%
87	20.368	2936	2941	2946	rVV4	20136	39003	1.86%	0.288%
88	20.515	2963	2966	2968	rVB4	12358	13398	0.64%	0.099%
89	20.562	2970	2974	2978	rVV4	14368	21596	1.03%	0.159%
90	20.961	3039	3042	3049	rVB	21873	38869	1.86%	0.287%
91	21.226	3083	3087	3094	rBV5	55209	112168	5.36%	0.828%
92	21.555	3135	3143	3148	rBV	970717	1664391	79.55%	12.290%
93	21.596	3148	3150	3159	rVB	83985	134037	6.41%	0.990%
94	23.676	3501	3504	3511	rVB	28724	59354	2.84%	0.438%
95	23.805	3521	3526	3533	rVB4	38838	84273	4.03%	0.622%
96	24.440	3628	3634	3641	rBV	82540	185971	8.89%	1.373%
97	24.663	3662	3672	3679	rBV2	573438	1555324	74.33%	11.485%
98	25.803	3859	3866	3874	rVB2	51628	134416	6.42%	0.993%
99	27.107	4084	4088	4097	rVB4	44426	118999	5.69%	0.879%
100	30.738	4704	4706	4707	rBV2	9304	6383	0.31%	0.047%

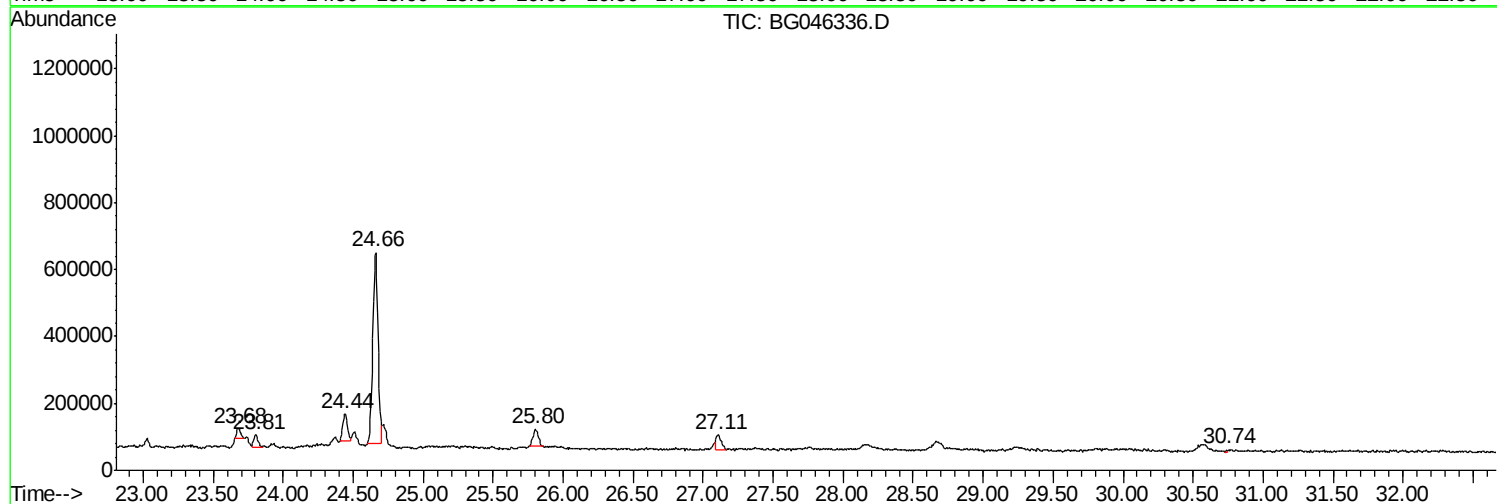
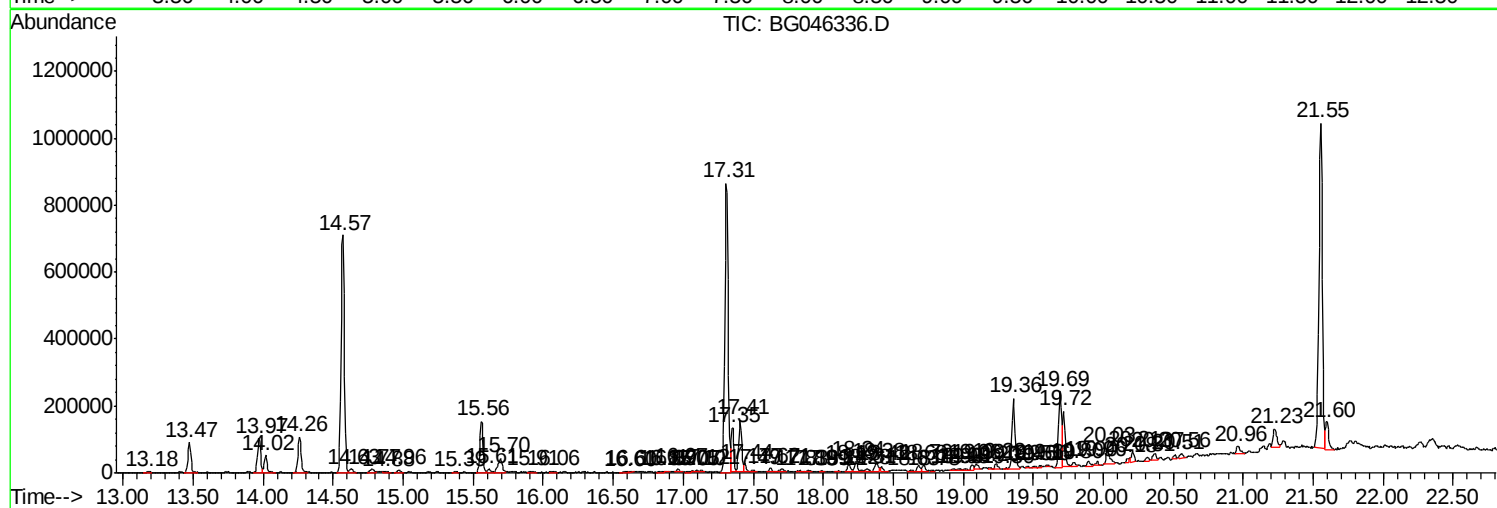
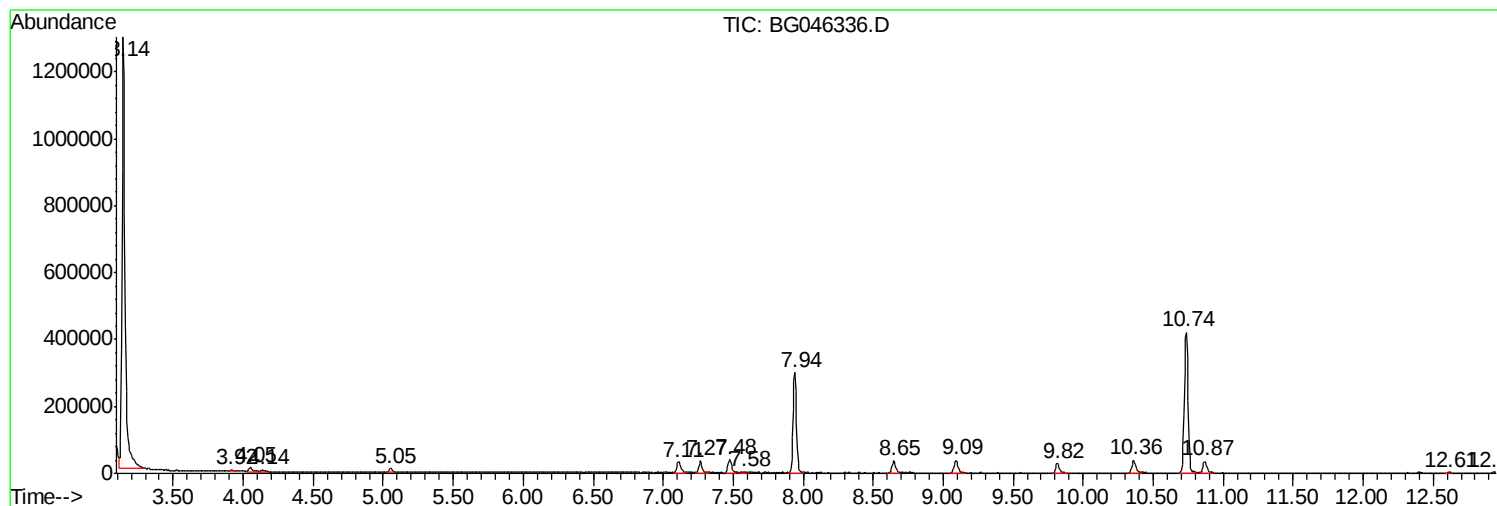
Sum of corrected areas: 13542790

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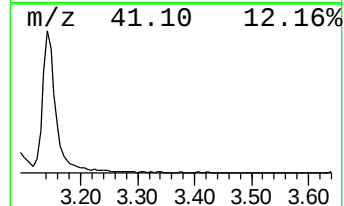
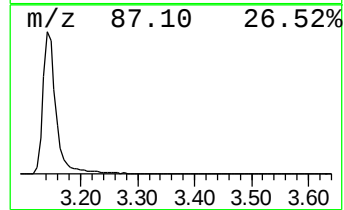
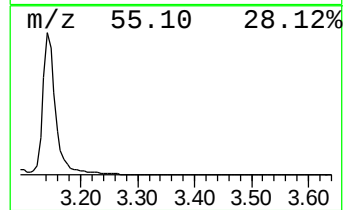
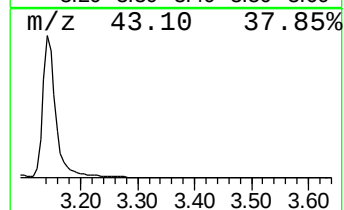
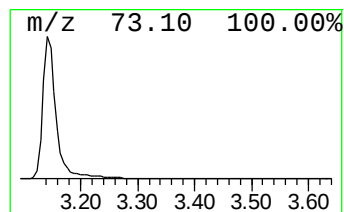
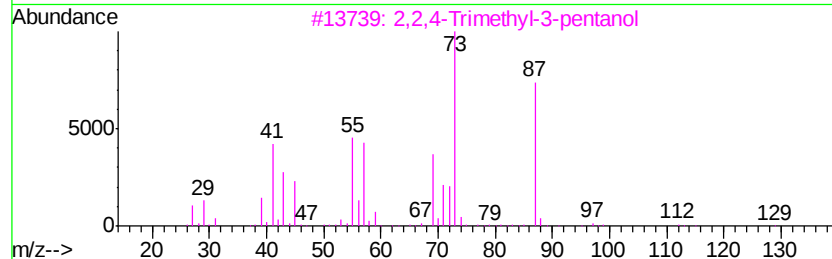
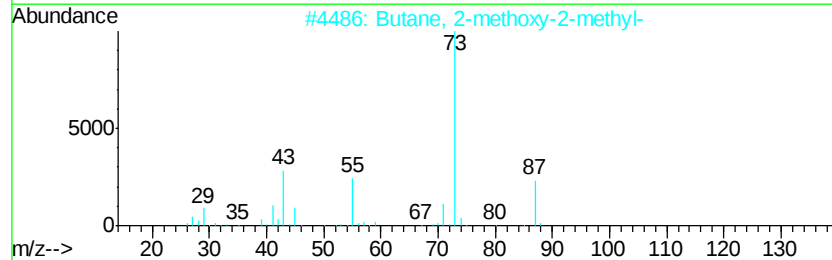
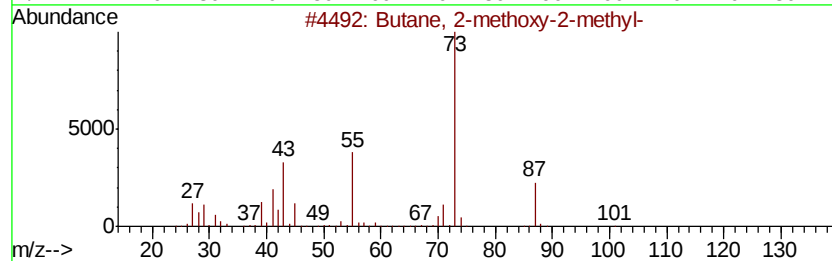
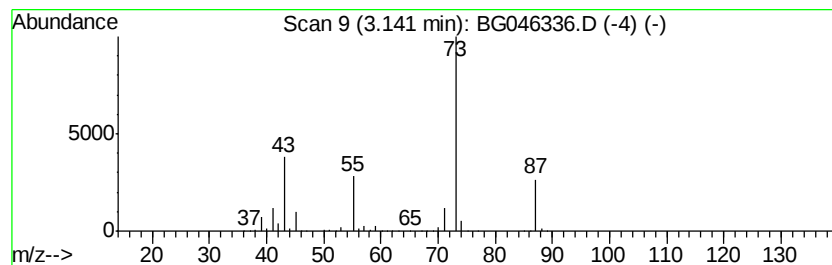
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG081320PAH.M
Quant Title : SVOA CALIBRATION

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.
3.14	82.72 ng/ul	2092330	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		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25



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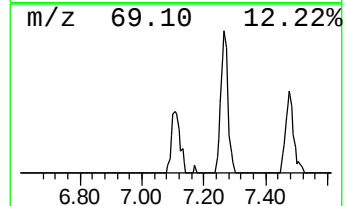
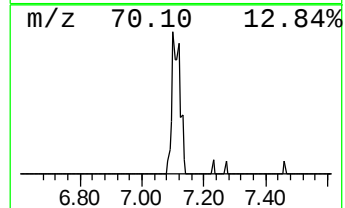
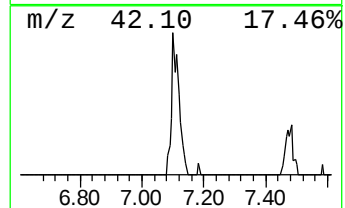
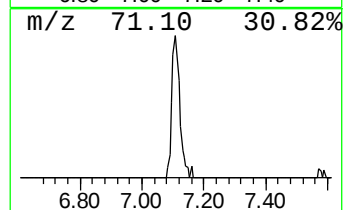
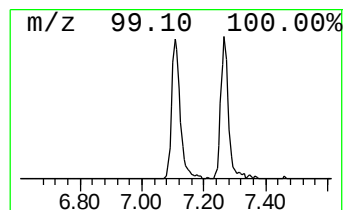
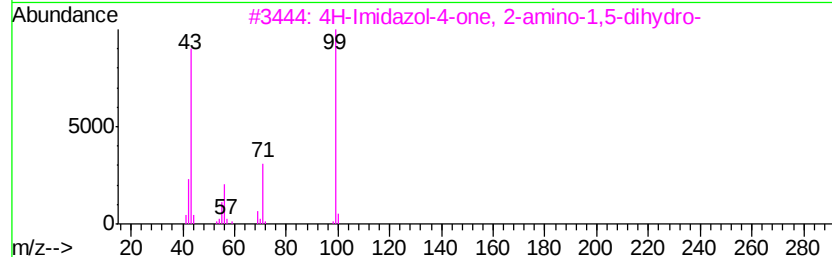
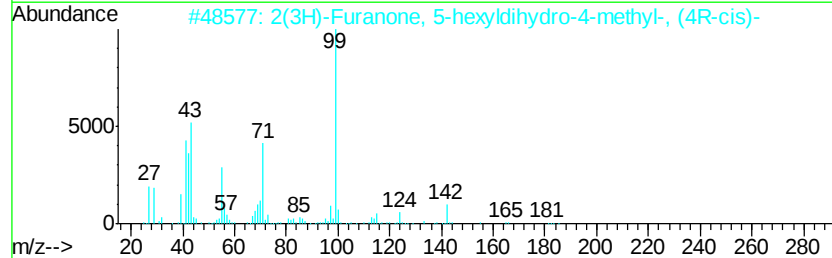
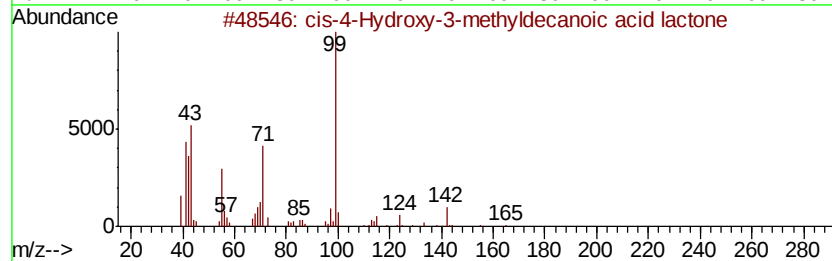
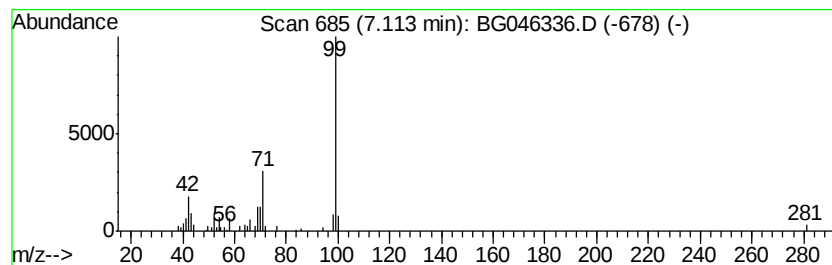
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 cis-4-Hydroxy-3-methyldecan... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.11	2.63 ng/ul	66445	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	cis-4-Hydroxy-3-methyldecanoic a...	184	C11H20O2	147254-32-8	64
2		2(3H)-Furanone, 5-hexyldihydro-4...	184	C11H20O2	121644-12-0	64
3		4H-Imidazol-4-one, 2-amino-1,5-d...	99	C3H5N3O	000503-86-6	59
4		Hexanoic acid, 5-tridecyl ester	298	C19H38O2	1000279-29-4	56
5		2(3H)-Furanone, 5-hexyldihydro-4...	184	C11H20O2	147254-33-9	56



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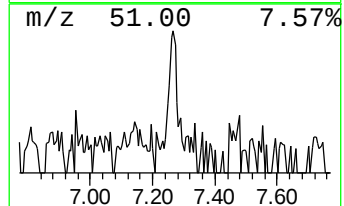
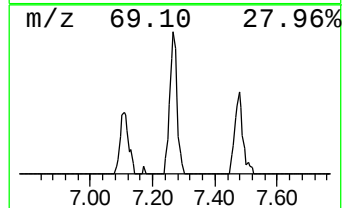
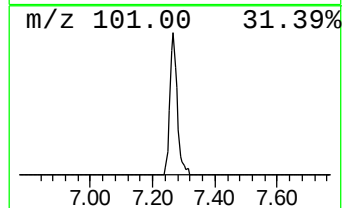
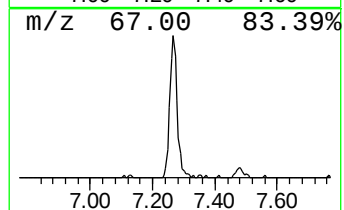
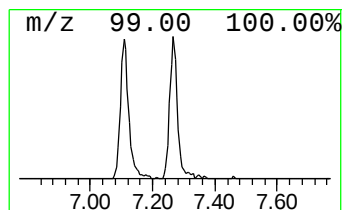
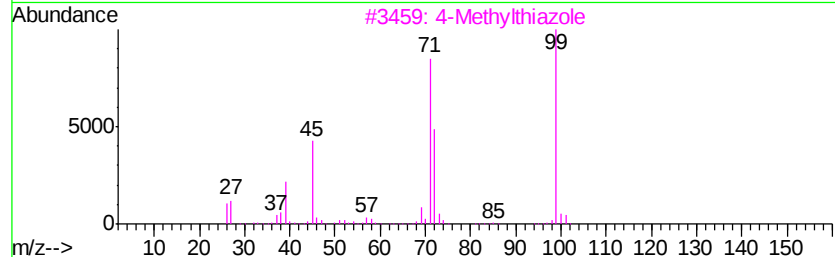
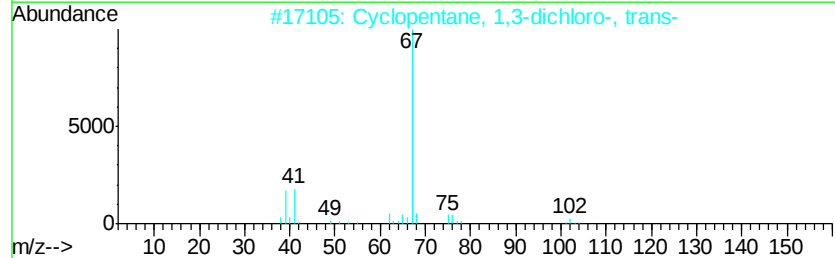
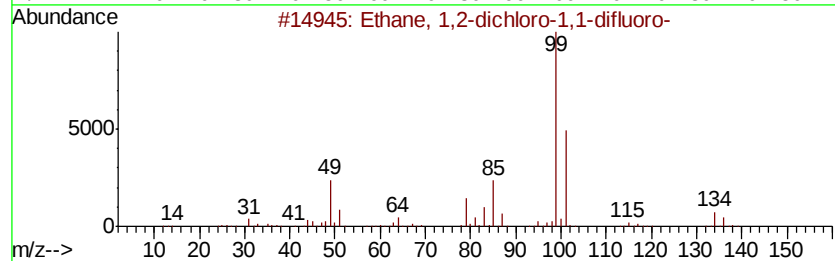
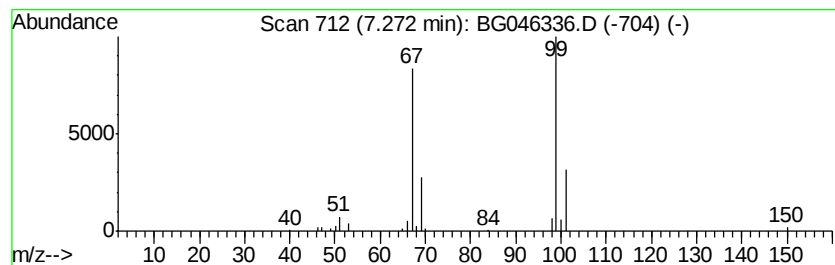
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.27	2.32 ng/ul	58560	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethane, 1,2-dichloro-1,1-difluoro-	134	C2H2Cl2F2	001649-08-7	25
2		Cyclopentane, 1,3-dichloro-, trans-	138	C5H8Cl2	026688-50-6	12
3		4-Methylthiazole	99	C4H5NS	000693-95-8	9
4		Methyl 2-butynoate	98	C5H6O2	023326-27-4	9
5		4-Methylthiazole	99	C4H5NS	000693-95-8	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081720\
Data File : BG046336.D
Acq On : 18 Aug 2020 20:33
Operator : CG/JU
Sample : L3636-16
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BFMN7

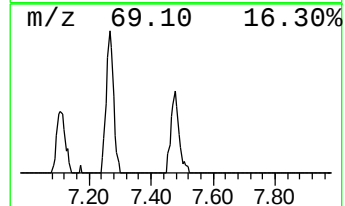
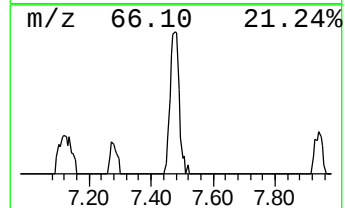
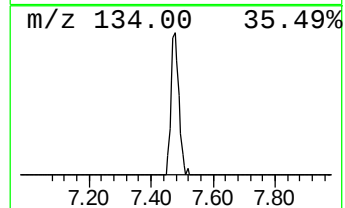
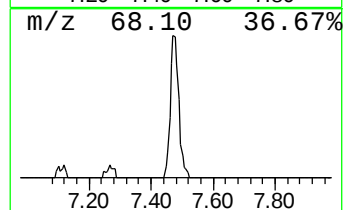
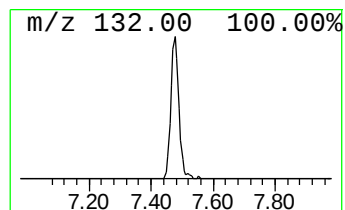
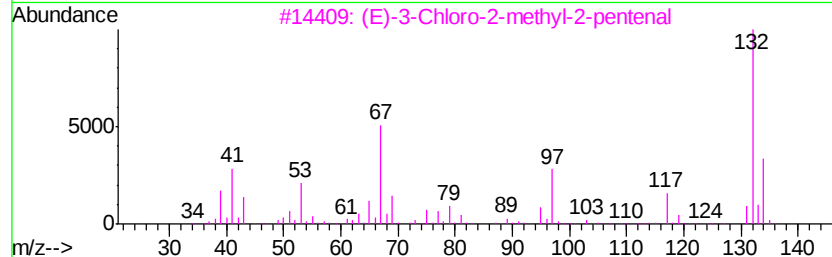
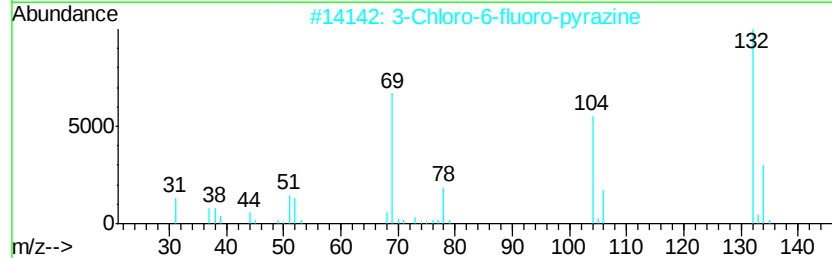
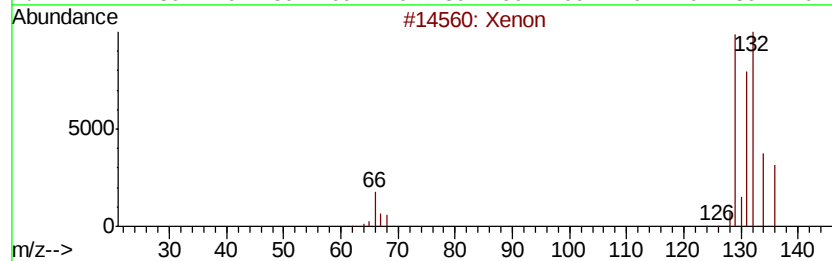
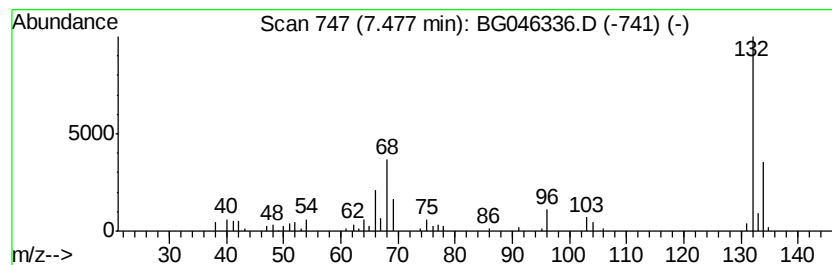
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Xenon Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.48	2.69 ng/ul	67960	1,4-Dichlorobenzene-d4	7.94

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Xenon		132	Xe	007440-63-3	59
2	3-Chloro-6-fluoro-pyrazine		132	C4H2ClFN2	1000146-10-7	35
3	(E)-3-Chloro-2-methyl-2-pentenal		132	C6H9ClO	031357-76-3	35
4	1,3-Cyclopentanedione, 2-chloro-		132	C5H5ClO2	014203-19-1	32
5	5-Fluoro-2-chloropyrimidine		132	C4H2ClFN2	062802-42-0	17



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081720\
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Acq On : 18 Aug 2020 20:33
Operator : CG/JU
Sample : L3636-16
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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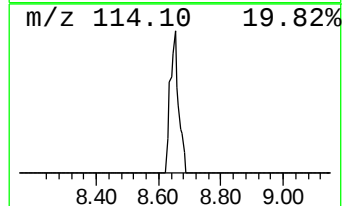
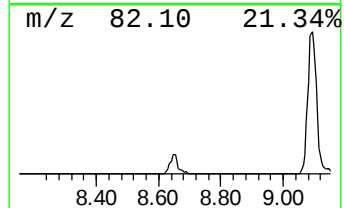
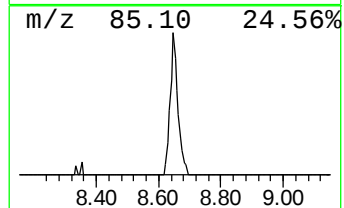
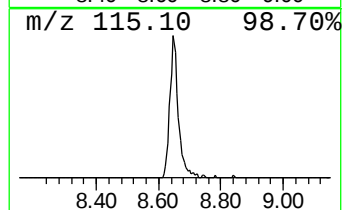
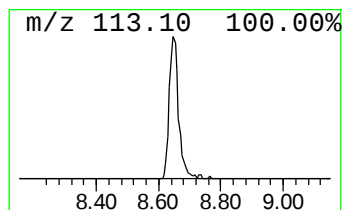
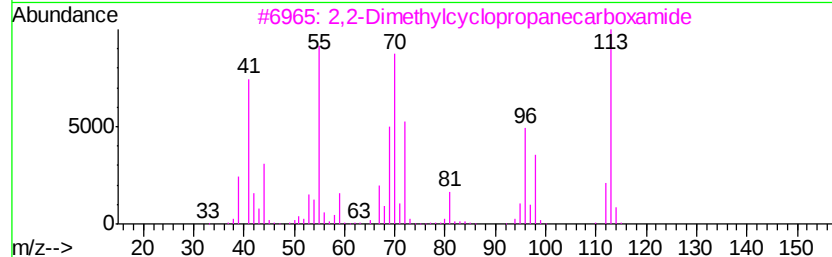
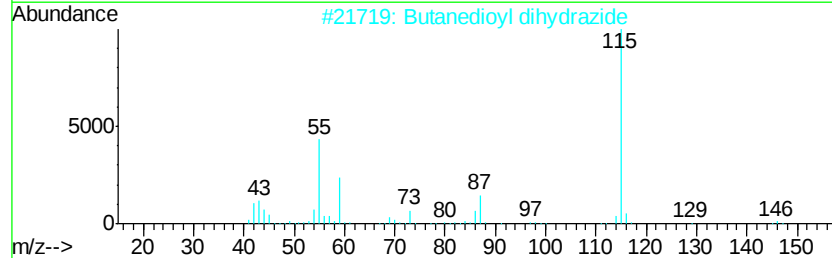
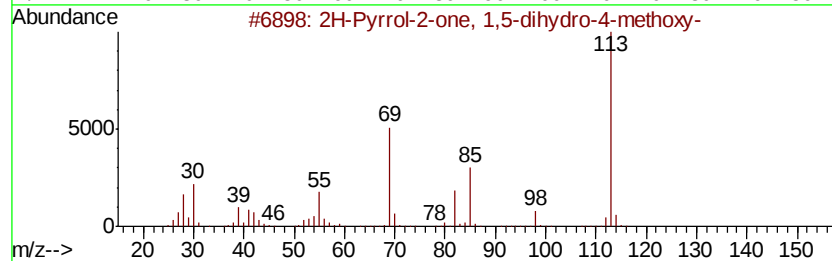
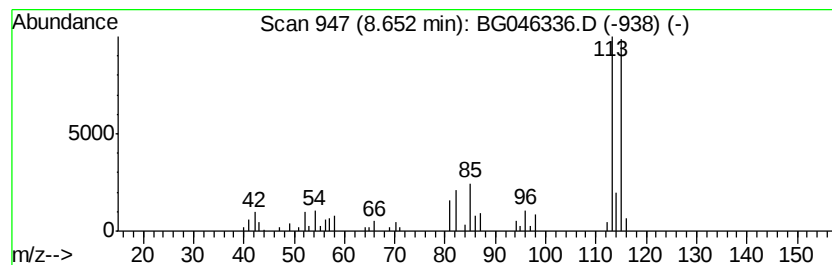
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-02 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.65	2.73 ng/ul	69084	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2H-Pyrrol-2-one, 1,5-dihydro-4-m...	113	C5H7NO2	069778-83-2	38
2		Butanedioyl dihydrazide	146	C4H10N4O2	004146-43-4	35
3		2,2-Dimethylcyclopropanecarboxamide	113	C6H11NO	075885-58-4	27
4		.alpha.-Methyl-.alpha.-propylsuc...	155	C8H13NO2	001497-19-4	25
5		Furazan-3-amine, 4-methoxy-	115	C3H5N3O2	078350-48-8	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081720\
Data File : BG046336.D
Acq On : 18 Aug 2020 20:33
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Sample : L3636-16
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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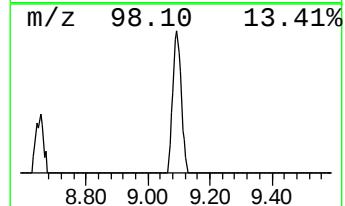
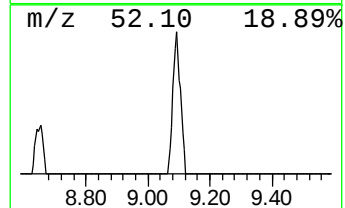
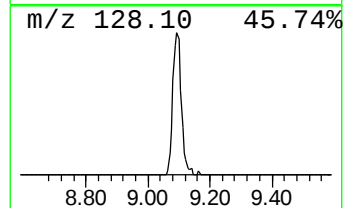
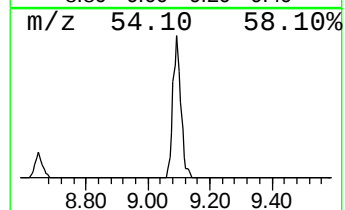
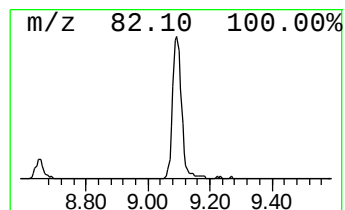
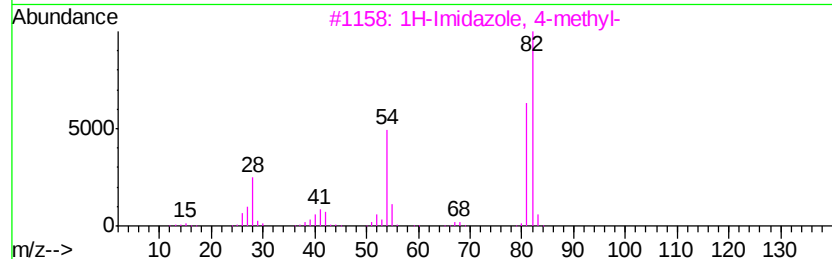
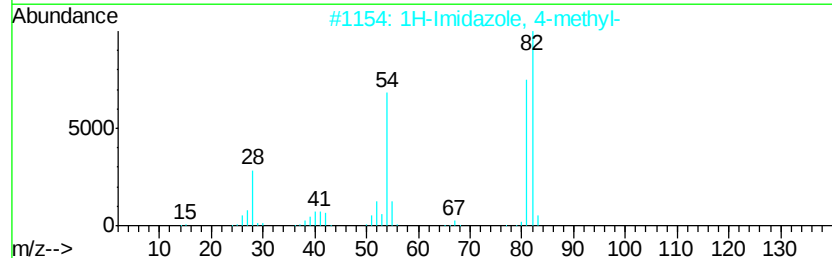
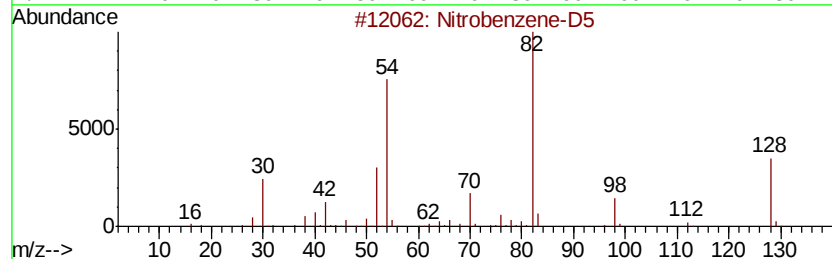
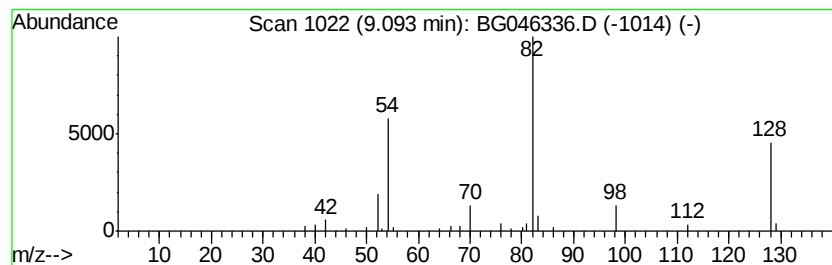
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-03 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.09	2.77 ng/ul	70019	1,4-Dichlorobenzene-d4	7.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nitrobenzene-D5	128	C6D5NO2	004165-60-0	91
2			1H-Imidazole, 4-methyl-	82	C4H6N2	000822-36-6	47
3			1H-Imidazole, 4-methyl-	82	C4H6N2	000822-36-6	47
4			Nitrobenzene-D5	128	C6D5NO2	004165-60-0	37
5			Butyramide, 2-cyano-2-ethyl-	140	C7H12N2O	018705-38-9	17



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081720\
Data File : BG046336.D
Acq On : 18 Aug 2020 20:33
Operator : CG/JU
Sample : L3636-16
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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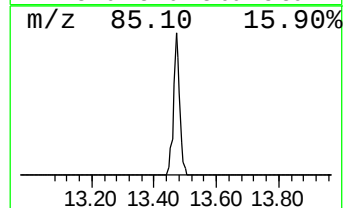
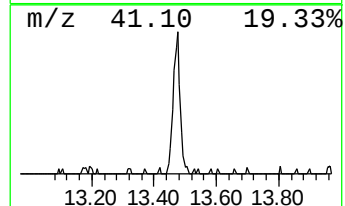
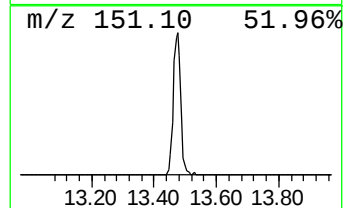
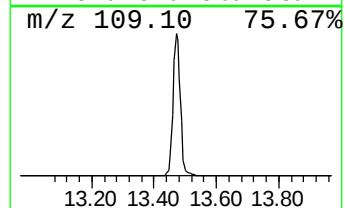
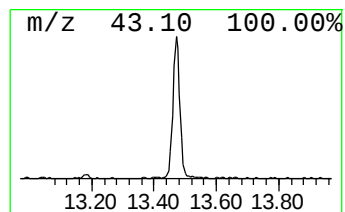
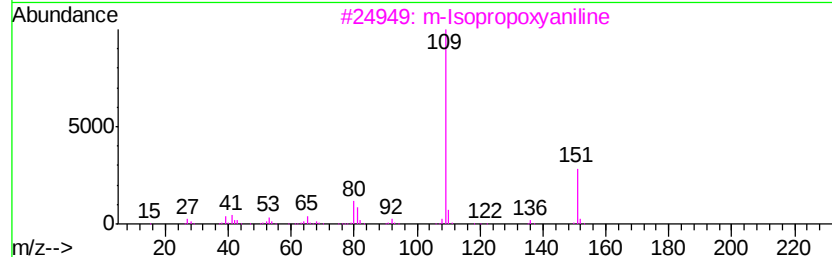
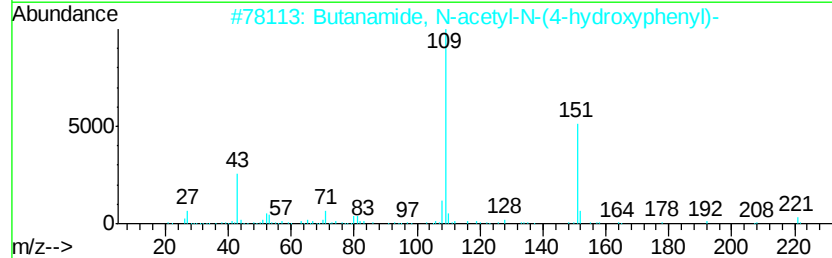
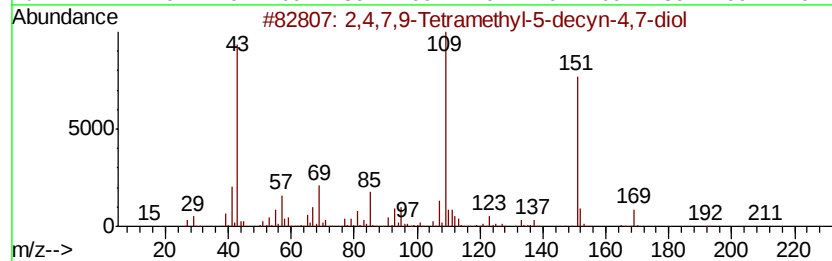
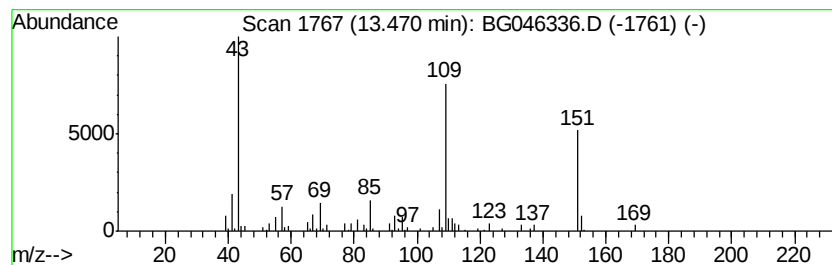
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	2.56 ng/ul	142346	Acenaphthene-d10	14.57

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91		
2	Butanamide, N-acetyl-N-(4-hydrox...	221	C12H15NO3	1000107-08-7	59		
3	m-Isopropoxvaniline	151	C9H13NO	041406-00-2	59		
4	Metacetamol	151	C8H9NO2	000621-42-1	59		
5	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	59		



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081720\
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Sample : L3636-16
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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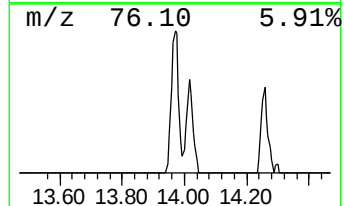
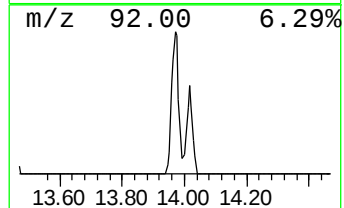
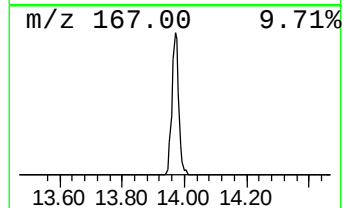
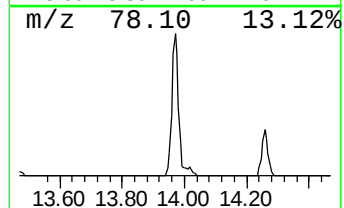
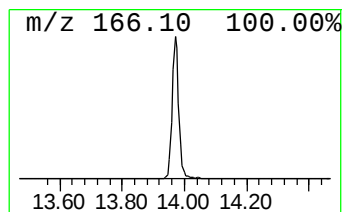
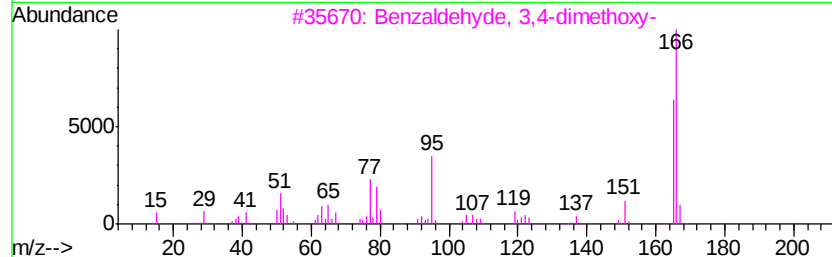
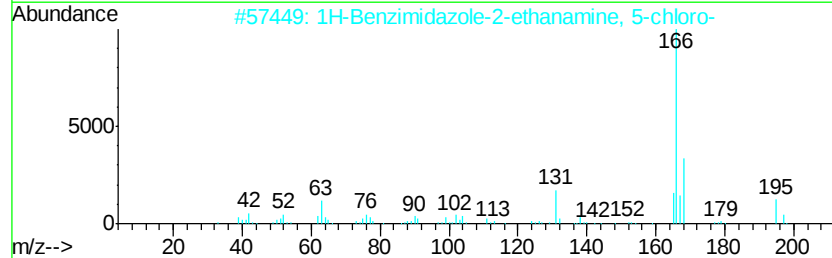
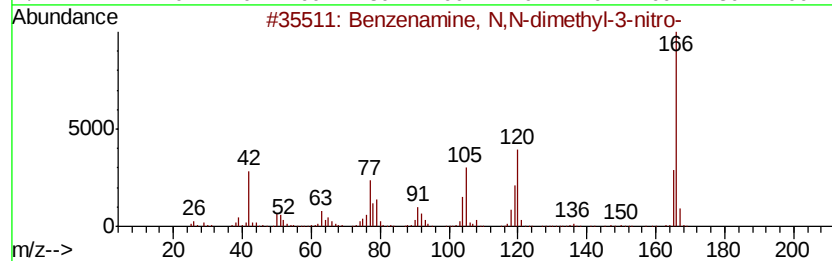
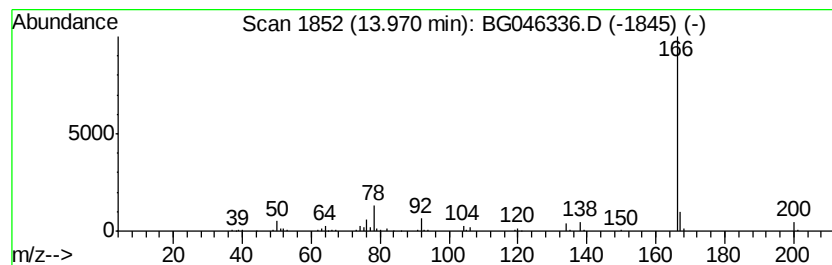
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzenamine, N,N-dimethyl-3... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	2.72 ng/ul	151407	Acenaphthene-d10	14.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenamine, N,N-dimethyl-3-nitro-	166	C8H10N2O2	000619-31-8	72
2		1H-Benzimidazole-2-ethanamine, 5...	195	C9H10ClN3	135875-16-0	38
3		Benzaldehyde, 3,4-dimethoxy-	166	C9H10O3	000120-14-9	38
4		Benzenamine, N,N-dimethyl-4-nitro-	166	C8H10N2O2	000100-23-2	38
5		3-Pyridineacetic acid, 1,4-dihyd...	167	C9H13NO2	039998-23-7	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG081720\
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Sample : L3636-16
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
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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...	3.14	82.7	ng/ul	2092330	1	7.94	505878	20.0
cis-4-Hydroxy-3-m...	7.11	2.6	ng/ul	66445	1	7.94	505878	20.0
unknown-01	7.27	2.3	ng/ul	58560	1	7.94	505878	20.0
Xenon	7.48	2.7	ng/ul	67960	1	7.94	505878	20.0
unknown-02	8.65	2.7	ng/ul	69084	1	7.94	505878	20.0
unknown-03	9.09	2.8	ng/ul	70019	1	7.94	505878	20.0
2,4,7,9-Tetrameth...	13.47	2.6	ng/ul	142346	3	14.57	1111410	20.0
Benzenamine, N,N-...	13.97	2.7	ng/ul	151407	3	14.57	1111410	20.0