

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
 Data File : BG046350.D
 Acq On : 19 Aug 2020 17:24
 Operator : CG/JU
 Sample : PB130882BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK82

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	3.142	5	9	15	rVB	36480	52967	2.16%	0.209%
2	3.400	49	53	59	rVB	21377	32204	1.32%	0.127%
3	4.146	174	180	185	rVV4	4431	7796	0.32%	0.031%
4	7.108	675	684	700	rBV	350923	612441	25.01%	2.417%
5	7.266	704	711	723	rVV	278200	463248	18.92%	1.829%
6	7.478	740	747	759	rBV	355436	625541	25.55%	2.469%
7	7.942	819	826	834	rBV	310054	528248	21.58%	2.085%
8	8.647	939	946	961	rBV	433271	765951	31.28%	3.023%
9	9.094	1013	1022	1031	rBV	323745	580859	23.72%	2.293%
10	9.816	1136	1145	1155	rBV	275541	500430	20.44%	1.975%
11	10.357	1230	1237	1255	rBV	447573	813868	33.24%	3.212%
12	10.739	1292	1302	1313	rBV	452708	800168	32.68%	3.158%
13	10.868	1316	1324	1337	rBV	404427	771053	31.49%	3.043%
14	12.319	1564	1571	1578	rBV4	6307	12412	0.51%	0.049%
15	13.970	1845	1852	1867	rBV	804043	1181762	48.27%	4.665%
16	14.258	1894	1901	1913	rVB	971378	1498965	61.22%	5.917%
17	14.569	1946	1954	1967	rVB	718039	1139433	46.54%	4.498%
18	14.763	1979	1987	2007	rBV	257152	493976	20.18%	1.950%
19	15.557	2115	2122	2134	rBV	1244061	1897523	77.50%	7.490%
20	15.674	2136	2142	2153	rBV	290553	428329	17.49%	1.691%
21	15.797	2159	2163	2169	rVB	14507	23159	0.95%	0.091%
22	16.344	2249	2256	2263	rVB4	4396	7791	0.32%	0.031%
23	16.996	2363	2367	2370	rBV3	2036	2912	0.12%	0.011%
24	17.096	2379	2384	2389	rBV5	2930	5088	0.21%	0.020%
25	17.307	2413	2420	2429	rBV	908978	1359934	55.54%	5.368%
26	17.407	2429	2437	2450	rVB2	1304772	2002273	81.78%	7.903%
27	17.584	2461	2467	2468	rBV3	2613	3468	0.14%	0.014%
28	18.535	2625	2629	2636	rVB4	2284	5399	0.22%	0.021%
29	19.047	2713	2716	2719	rBV4	2210	3020	0.12%	0.012%
30	19.264	2749	2753	2758	rBV5	2830	4453	0.18%	0.018%
31	19.329	2758	2764	2770	rVV	15738	25555	1.04%	0.101%
32	19.581	2803	2807	2812	rBV2	15068	22871	0.93%	0.090%
33	19.634	2812	2816	2819	rVB5	4546	5257	0.21%	0.021%
34	19.693	2819	2826	2837	rBV	1660502	2411128	98.48%	9.517%

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35	20.069	2889	2890	2892	rBV2	3281	2571	0.11%	0.010%
36	20.421	2948	2950	2951	rBV2	4543	3294	0.13%	0.013%
37	20.615	2980	2983	2984	rBV3	7706	9018	0.37%	0.036%
38	21.555	3136	3143	3155	rBV	989035	1554076	63.47%	6.134%
39	23.030	3389	3394	3399	rVB3	30259	57615	2.35%	0.227%
40	23.806	3521	3526	3533	rVB	47871	91339	3.73%	0.361%
41	24.446	3625	3635	3647	rBV	886579	2448350	100.00%	9.664%
42	24.658	3661	3671	3690	rVB2	570398	1738567	71.01%	6.862%
43	25.803	3859	3866	3876	rVB2	65244	177355	7.24%	0.700%
44	27.102	4082	4087	4097	rVB3	56633	162963	6.66%	0.643%

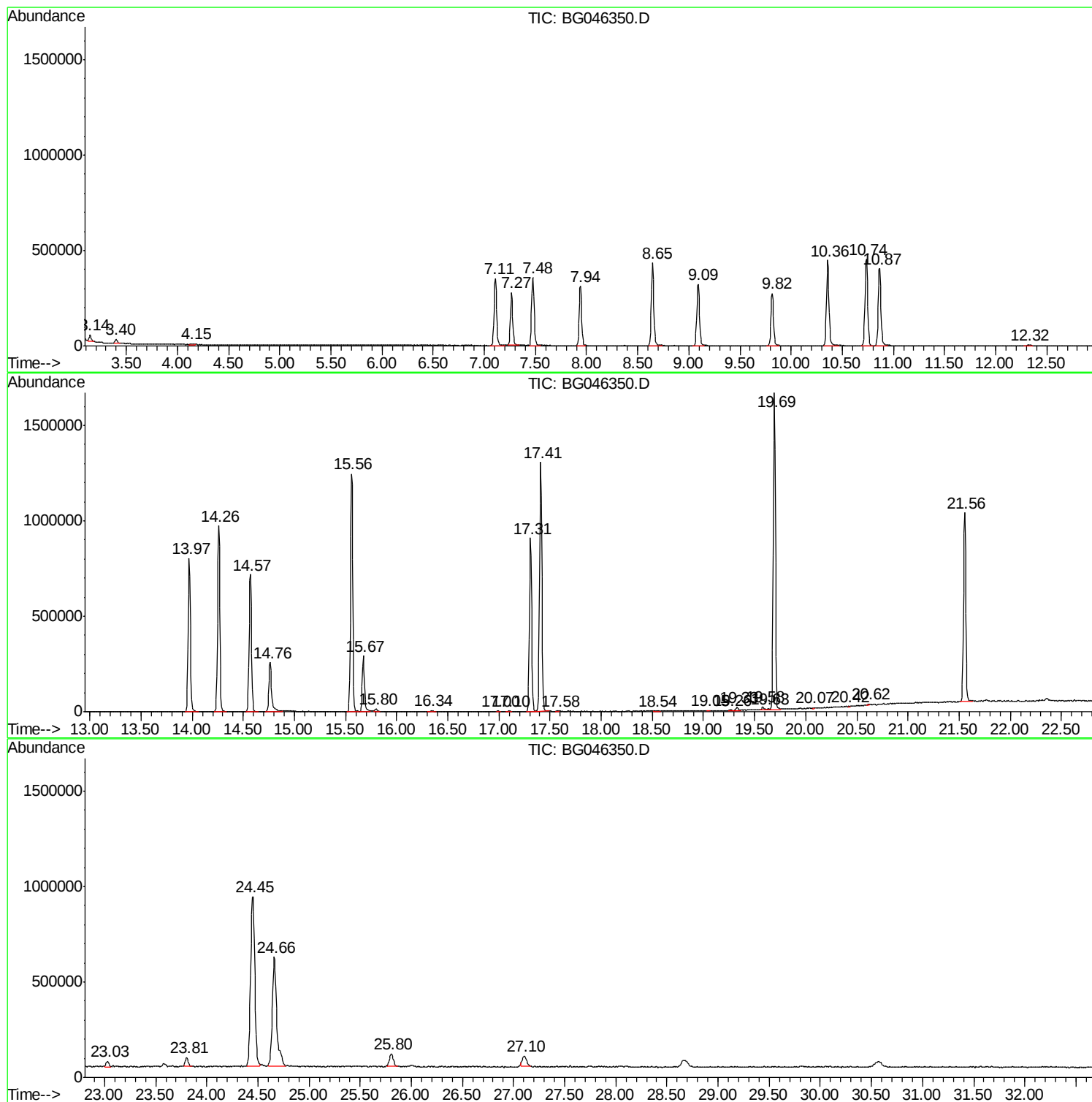
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG081320PAH.M
Quant Title : SVOA CALIBRATION

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



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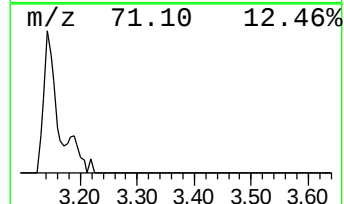
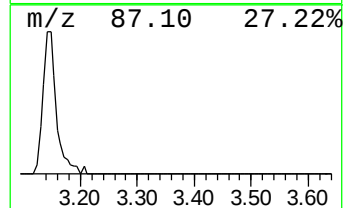
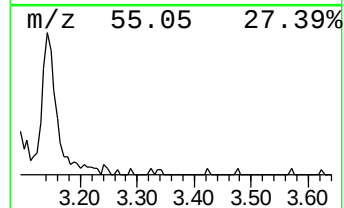
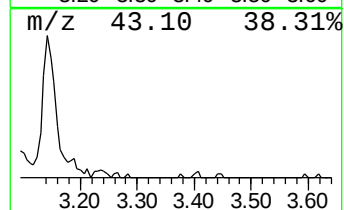
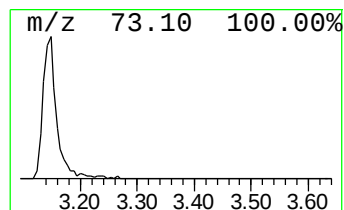
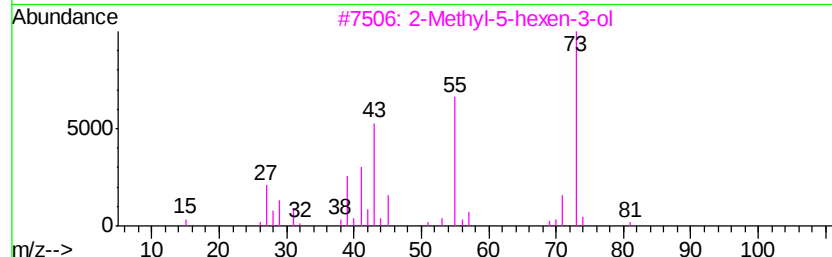
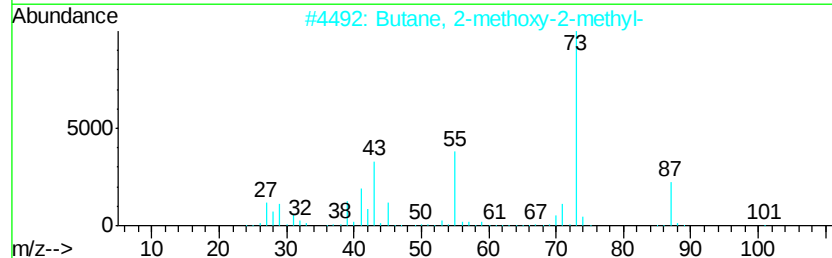
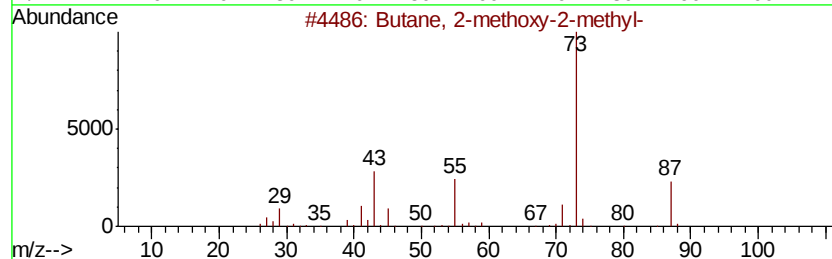
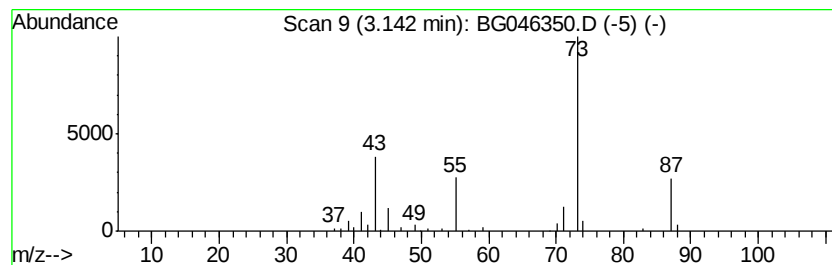
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.14	2.01 ng/ul	52967	1,4-Dichlorobenzene-d4	7.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	56
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	40
3			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	12
4			1,4-Butanediamine	88	C4H12N2	000110-60-1	9
5			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	9



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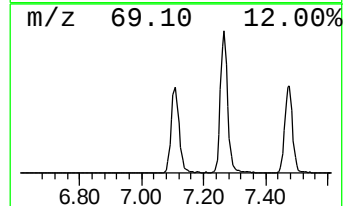
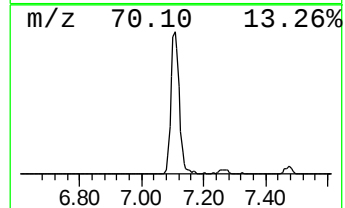
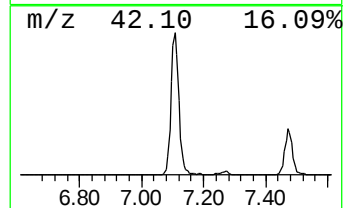
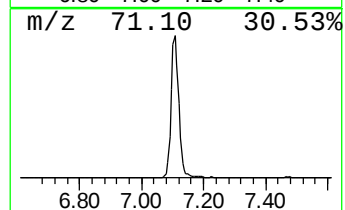
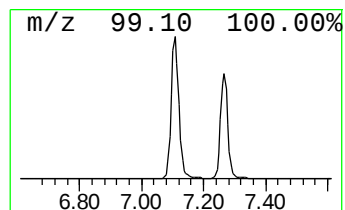
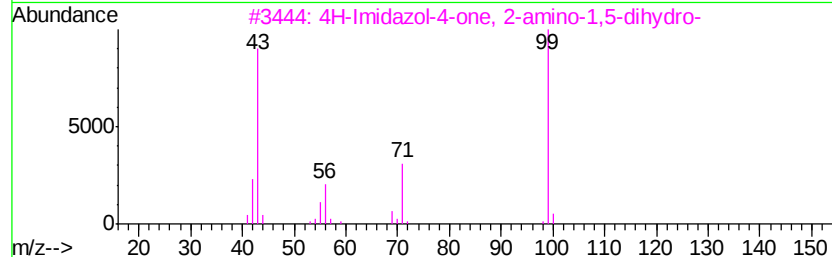
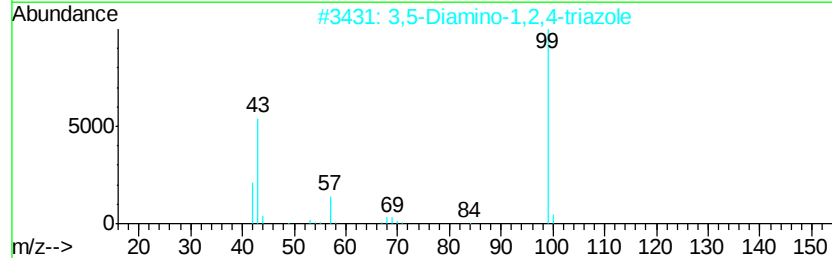
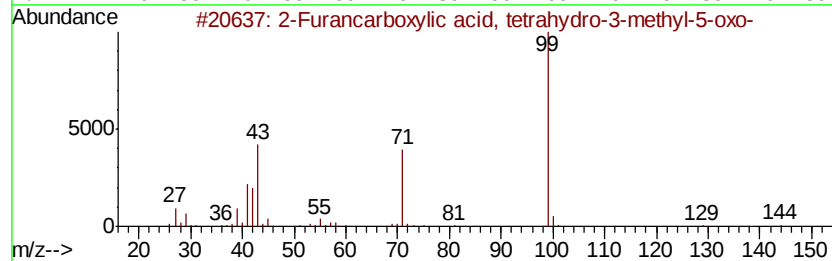
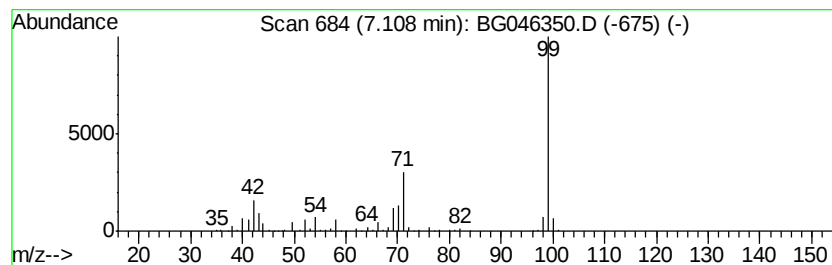
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 2-Furancarboxylic acid, tet... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.11	23.19 ng/ul	612441	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Furancarboxylic acid, tetrahyd...	144	C6H8O4	022073-04-7	59
2		3,5-Diamino-1,2,4-triazole	99	C2H5N5	001455-77-2	50
3		4H-Imidazol-4-one, 2-amino-1,5-d...	99	C3H5N3O	000503-86-6	45
4		Phenol-d6-	100	C6D6O	013127-88-3	40
5		2-Furancarboxylic acid, tetrahyd...	158	C7H10O4	085547-53-1	38



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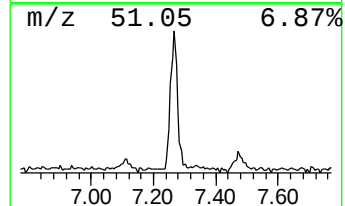
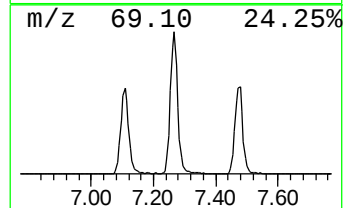
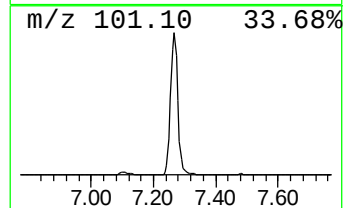
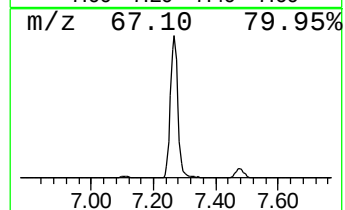
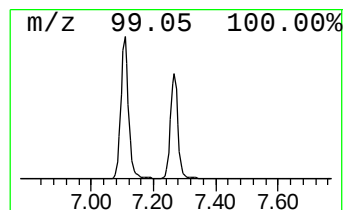
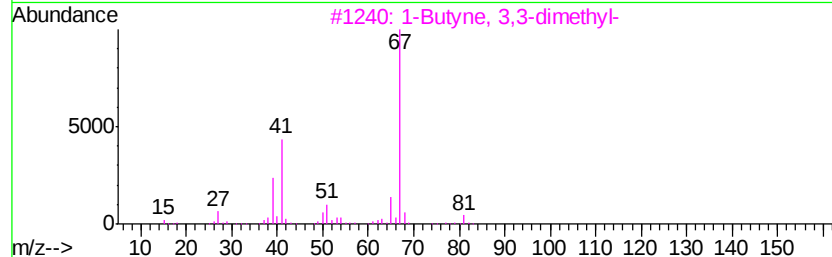
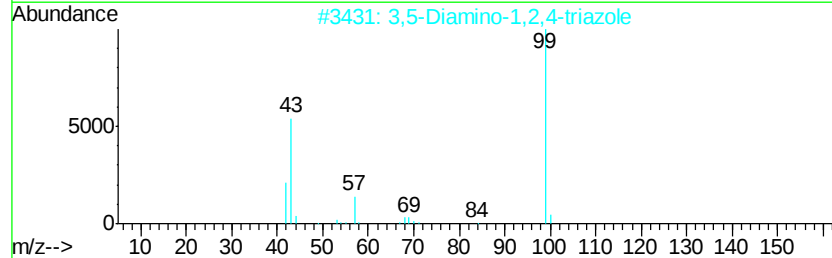
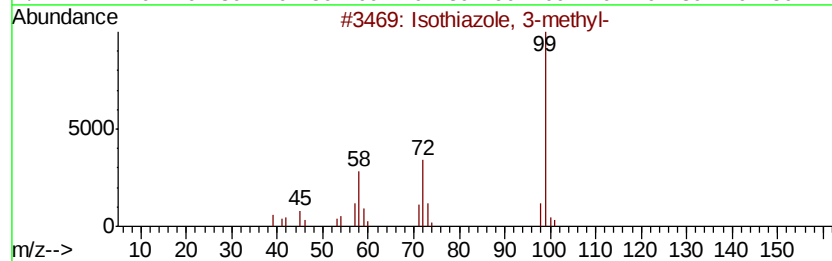
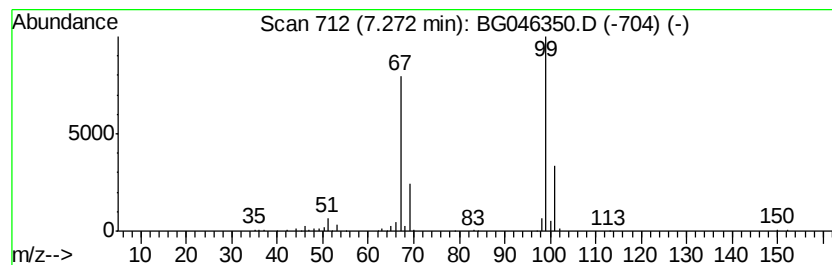
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.27	17.54 ng/ul	463248	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isothiazole, 3-methyl-	99	C4H5NS	000693-92-5	9
2		3,5-Diamino-1,2,4-triazole	99	C2H5N5	001455-77-2	9
3		1-Butyne, 3,3-dimethyl-	82	C6H10	000917-92-0	9
4		1-Pentanamine, 5-(methylthio)-	133	C6H15NS	055021-78-8	9
5		3,5-Diamino-1,2,4-triazole	99	C2H5N5	001455-77-2	7



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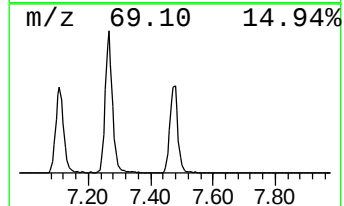
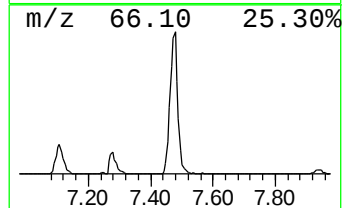
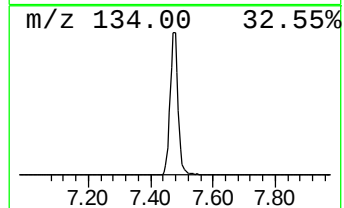
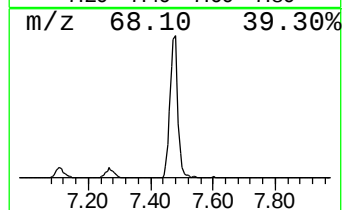
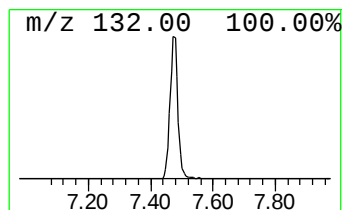
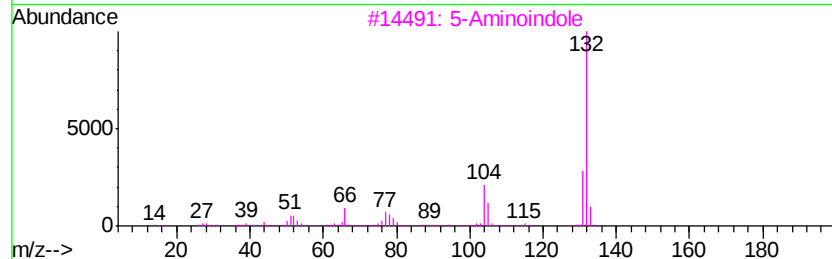
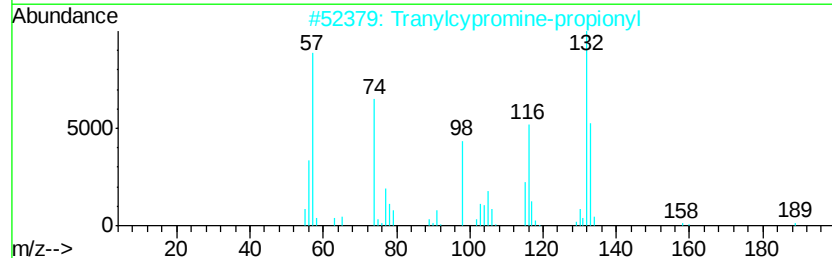
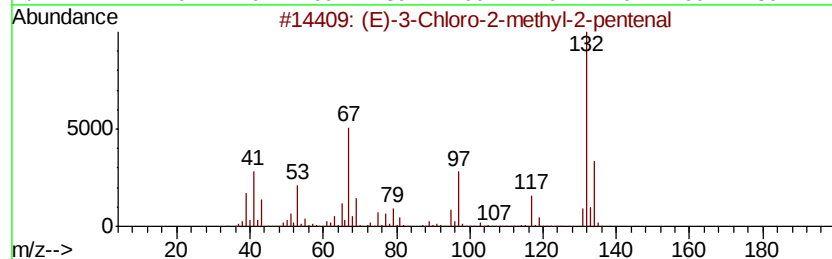
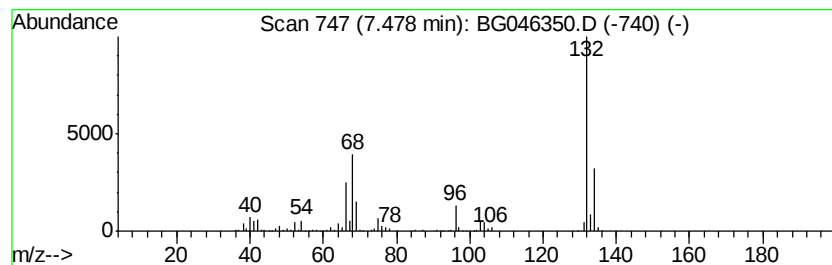
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown-02 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.48	23.68 ng/ul	625541	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		N-(-phenylisopropyl)benzalimine	223	C16H17N	1000379-01-3	10
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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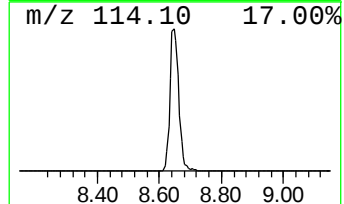
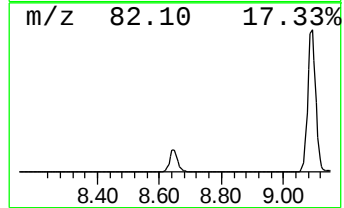
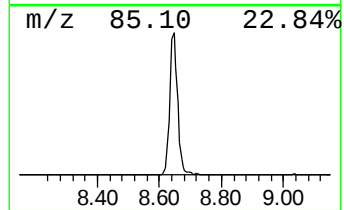
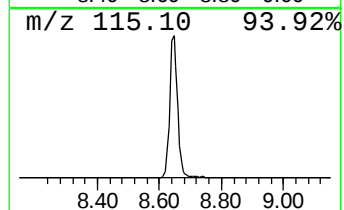
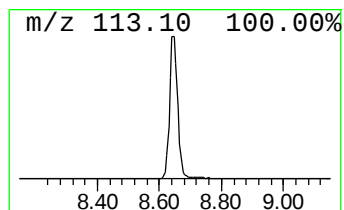
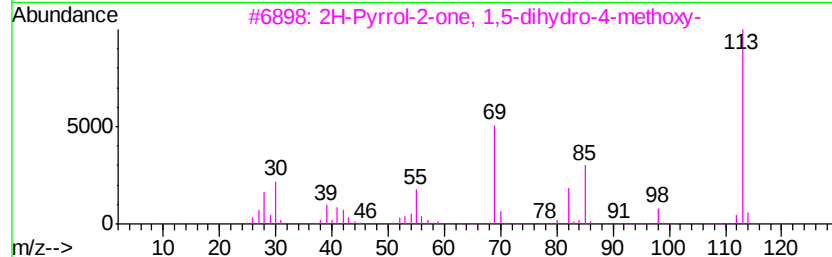
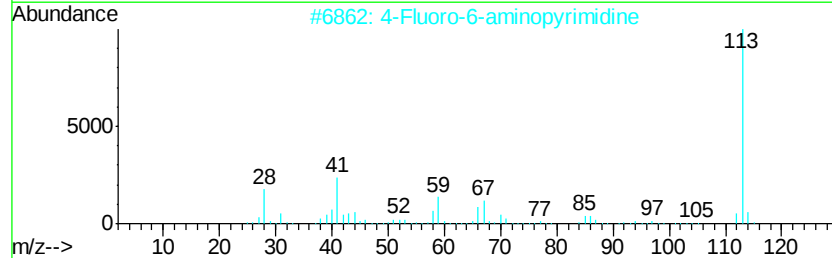
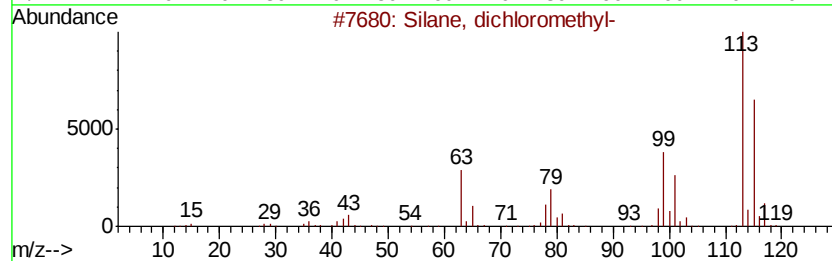
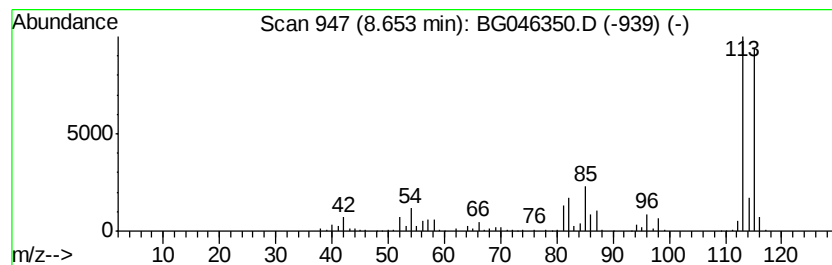
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-03 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.65	29.00 ng/ul	765951	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Silane, dichloromethyl-	114	CH4Cl2Si	000075-54-7	38
2		4-Fluoro-6-aminopyrimidine	113	C4H4FN3	051421-96-6	35
3		2H-Pyrrol-2-one, 1,5-dihydro-4-methoxy-	113	C5H7NO2	069778-83-2	32
4		Heptanoic acid, anhydride	242	C14H26O3	000626-27-7	27
5		2,2-Dimethylcyclopropanecarboxamide	113	C6H11NO	075885-58-4	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046350.D
Acq On : 19 Aug 2020 17:24
Operator : CG/JU
Sample : PB130882BL
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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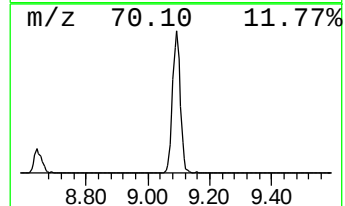
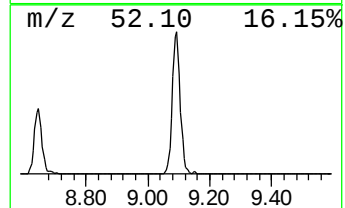
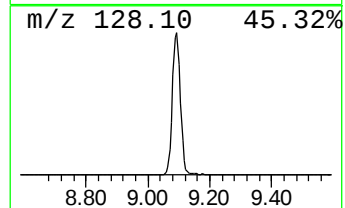
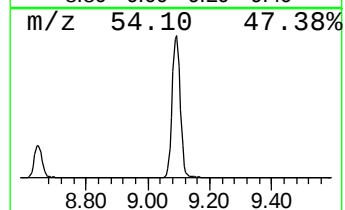
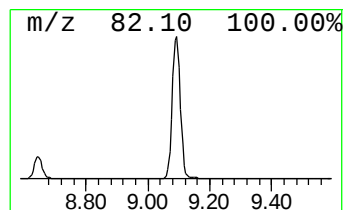
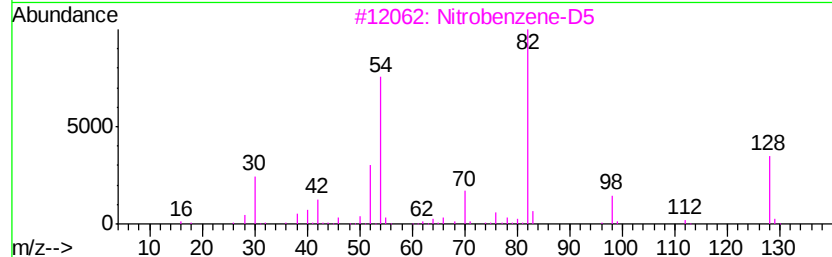
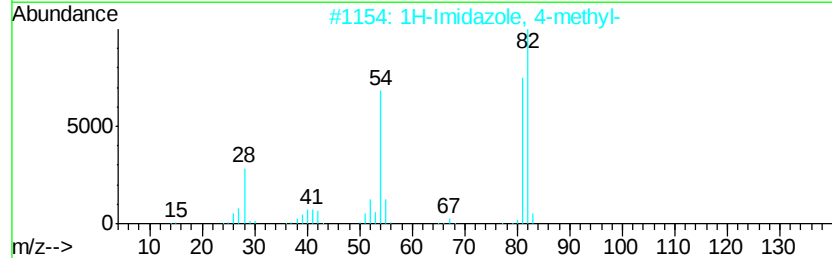
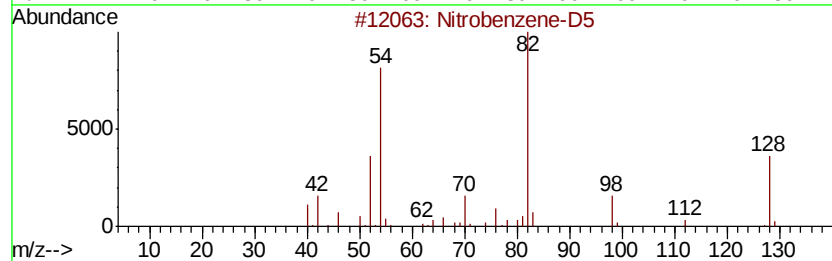
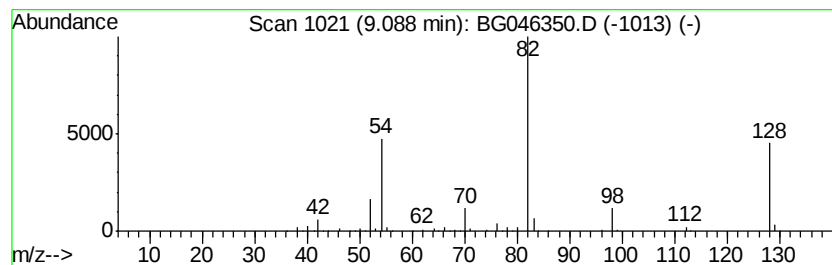
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1H-Imidazole, 4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.09	21.99 ng/ul	580859	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nitrobenzene-D5	128	C6D5NO2	004165-60-0	56
2		1H-Imidazole, 4-methyl-	82	C4H6N2	000822-36-6	50
3		Nitrobenzene-D5	128	C6D5NO2	004165-60-0	47
4		1H-Imidazole, 4-methyl-	82	C4H6N2	000822-36-6	38
5		1H-Imidazole, 2-methyl-	82	C4H6N2	000693-98-1	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046350.D
Acq On : 19 Aug 2020 17:24
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Sample : PB130882BL
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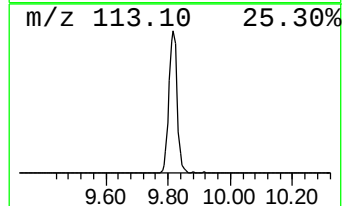
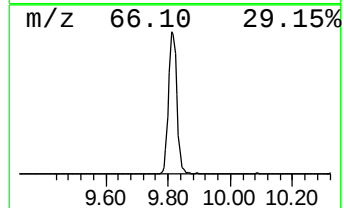
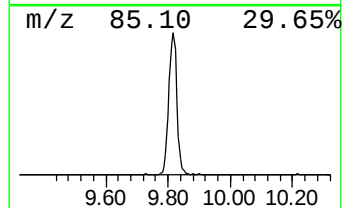
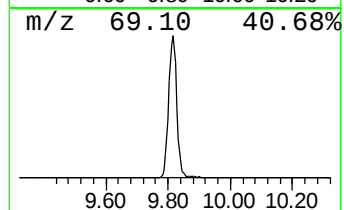
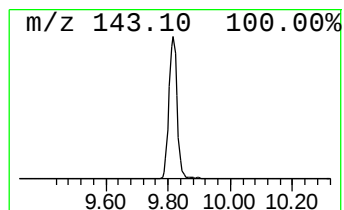
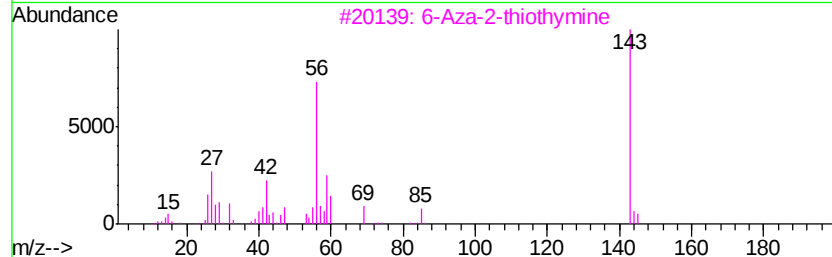
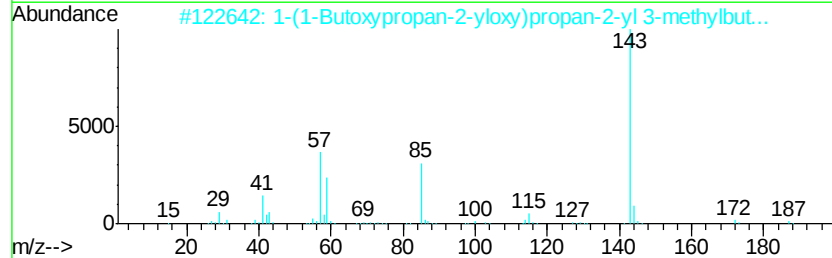
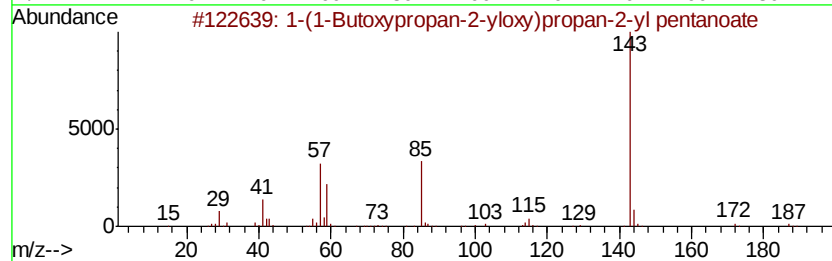
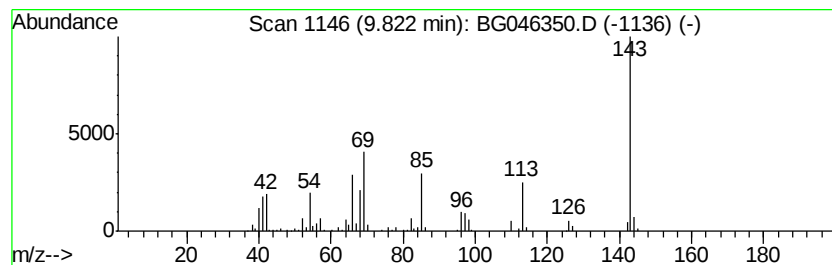
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-04 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.82	12.51 ng/ul	500430	Naphthalene-d8	10.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-(1-Butoxypropan-2-yloxy)propan...	274	C15H30O4	1000367-13-0	16
2		1-(1-Butoxypropan-2-yloxy)propan...	274	C15H30O4	1000367-12-9	16
3		6-Aza-2-thiothymine	143	C4H5N3OS	000615-76-9	10
4		3H-1,2,4-Triazole-3-thione,2,4-d...	143	C5H9N3S	037526-42-4	10
5		Quinoline, 7-methyl-	143	C10H9N	000612-60-2	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
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Misc :
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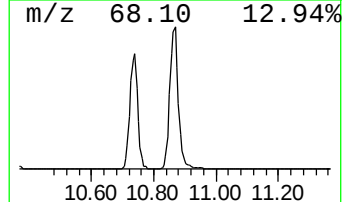
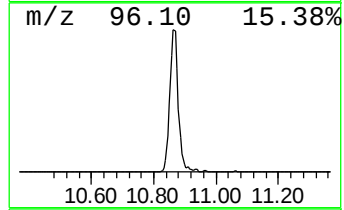
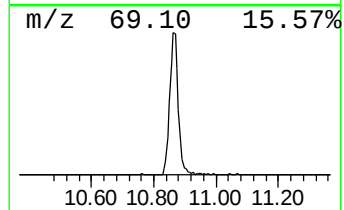
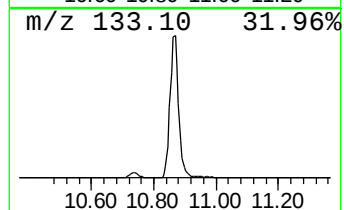
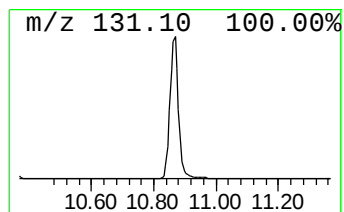
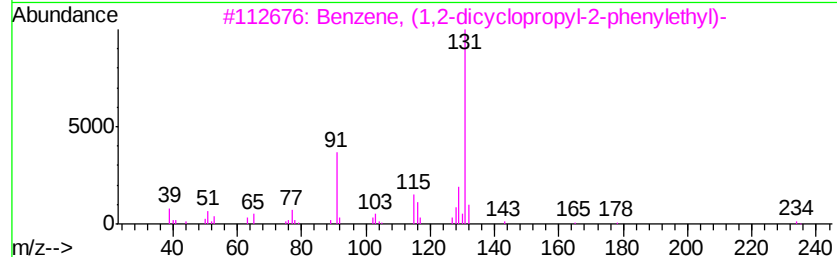
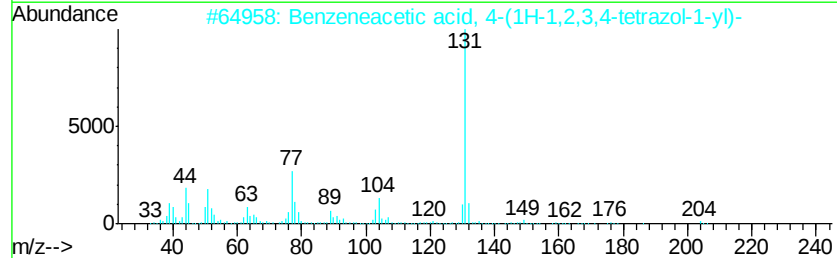
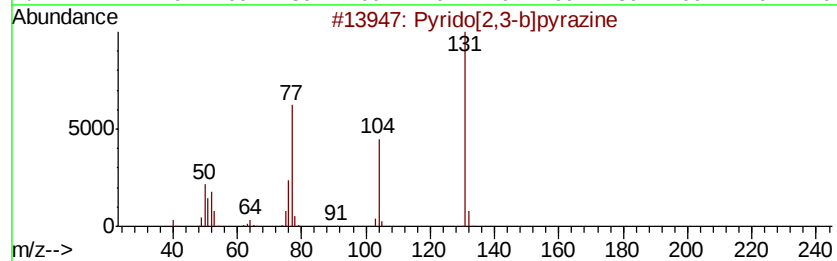
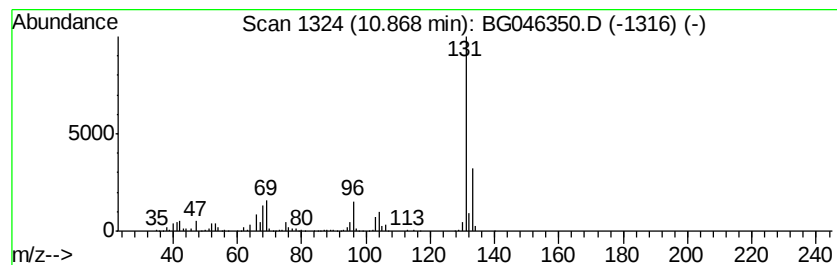
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-05 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.87	19.27 ng/ul	771053	Naphthalene-d8	10.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrido[2,3-b]pyrazine	131	C7H5N3	000322-46-3	38
2		Benzeneacetic acid, 4-(1H-1,2,3,...	204	C9H8N4O2	1000351-56-5	33
3		Benzene, (1,2-dicyclopropyl-2-ph...	262	C20H22	110330-90-0	33
4		1H-Indole, 1-methyl-	131	C9H9N	000603-76-9	28
5		Pyrido[2,3-b]pyrazine	131	C7H5N3	000322-46-3	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046350.D
Acq On : 19 Aug 2020 17:24
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Misc :
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ClientSampleId :
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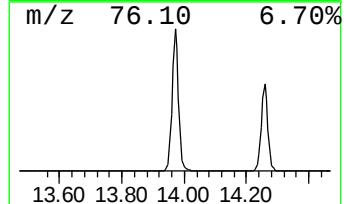
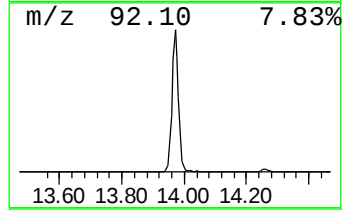
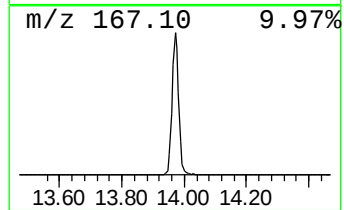
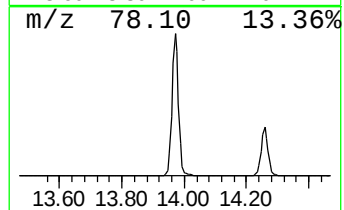
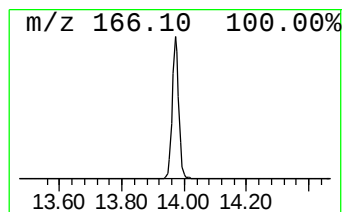
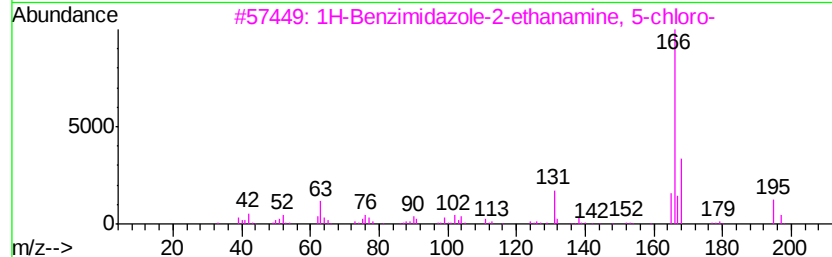
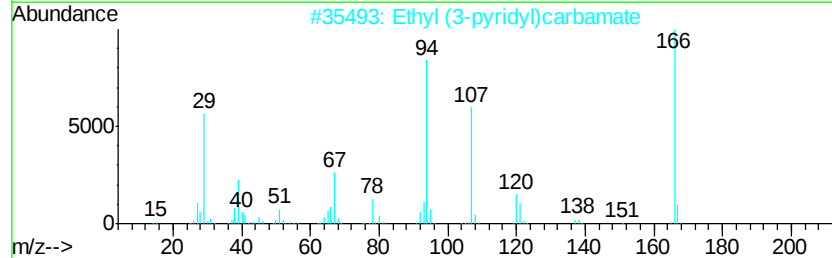
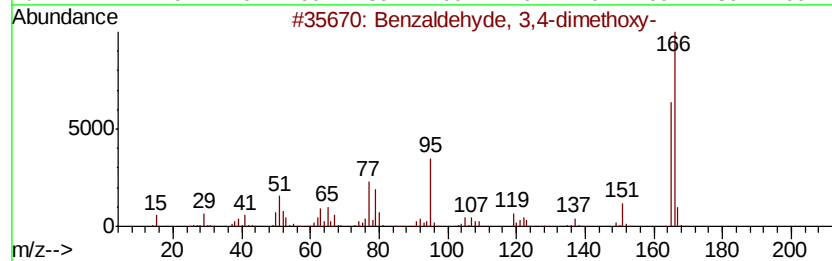
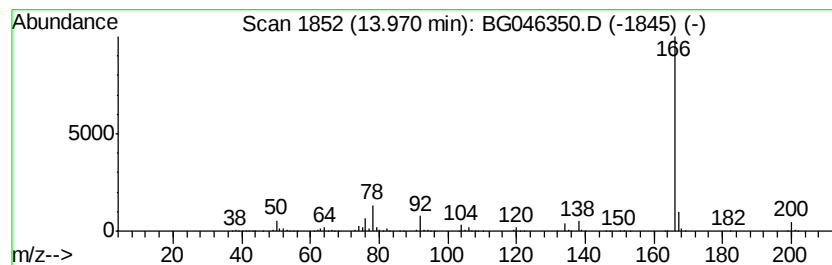
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-06 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	20.74 ng/ul	1181760	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		Ethyl (3-pyridyl)carbamate	166	C8H10N2O2	006276-11-5	39
3		1H-Benzimidazole-2-ethanamine, 5...	195	C9H10ClN3	135875-16-0	38
4		Benzenamine, N,N-dimethyl-4-nitro-	166	C8H10N2O2	000100-23-2	38
5		1H-Purine, 6-(methylthio)-	166	C6H6N4S	000050-66-8	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
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Acq On : 19 Aug 2020 17:24
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Sample : PB130882BL
Misc :
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Instrument :
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ClientSampleId :
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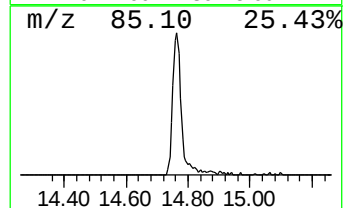
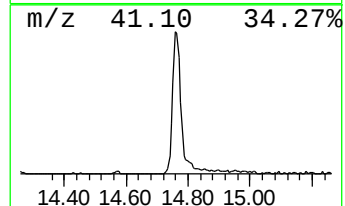
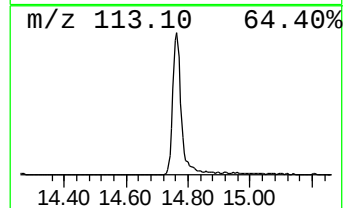
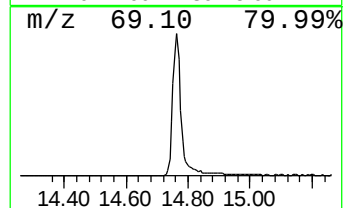
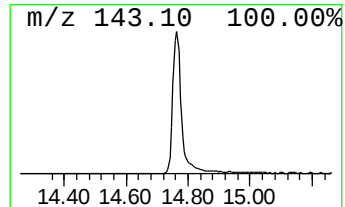
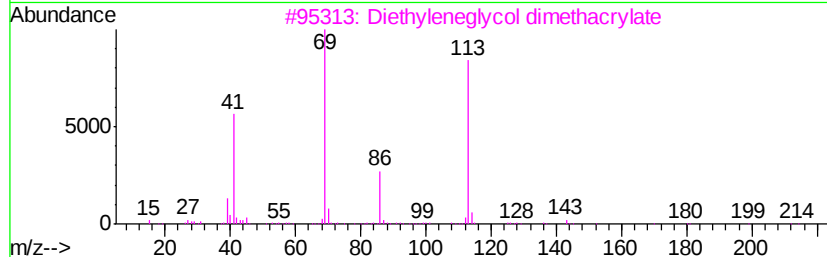
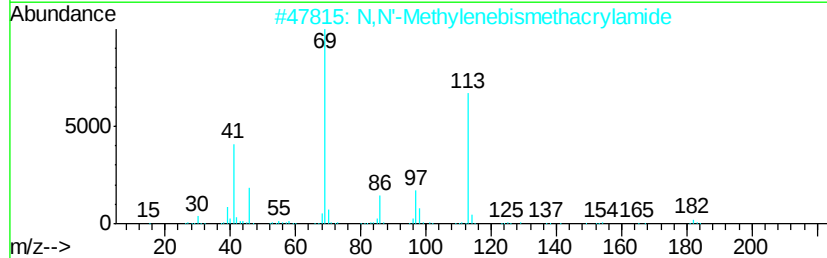
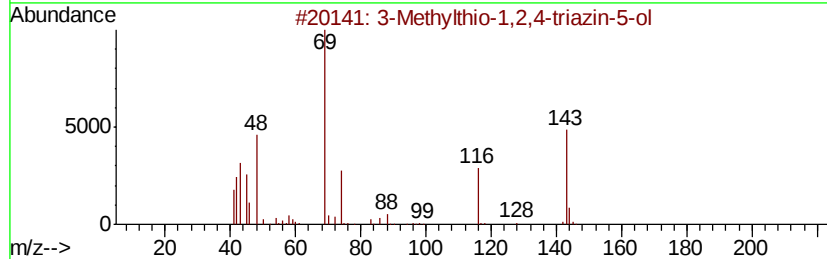
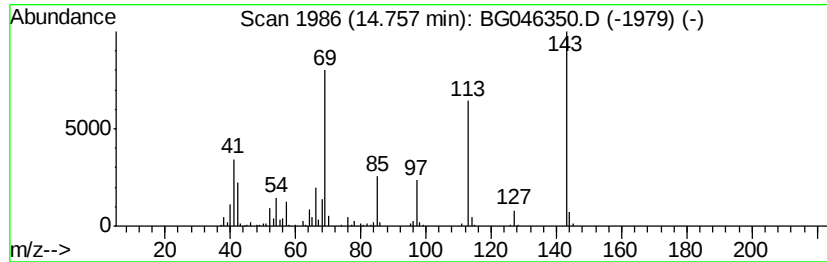
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-07 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.76	8.67 ng/ul	493976	Acenaphthene-d10	14.57

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Methylthio-1,2,4-triazin-5-ol	143	C4H5N3OS	018060-72-5	38
2			N,N'-Methylenebismethacrylamide	182	C9H14N2O2	002359-15-1	32
3			Diethyleneglycol dimethacrylate	242	C12H18O5	002358-84-1	25
4			4-Amino-2-chloro-6-methylpyrimidine	143	C5H6ClN3	014394-60-6	22
5			Glutaric acid, ethyl 4-methylhep...	272	C15H28O4	1000377-14-9	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
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Sample : PB130882BL
Misc :
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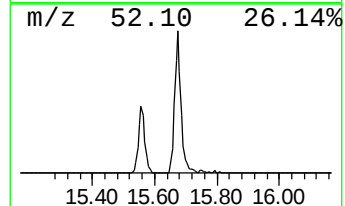
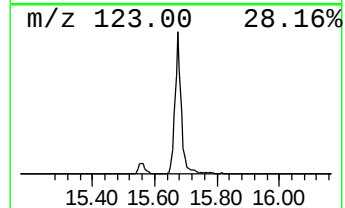
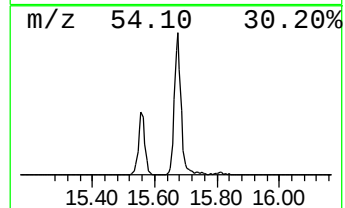
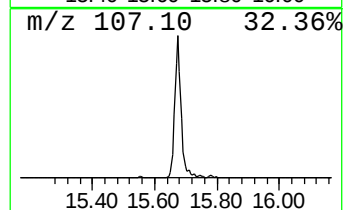
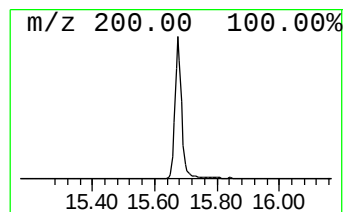
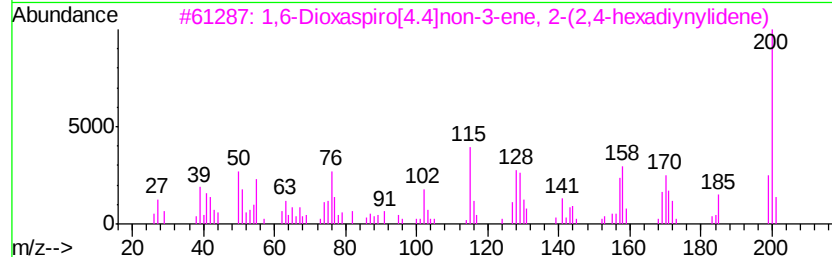
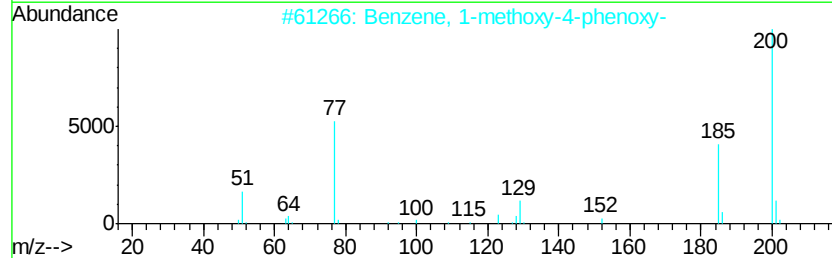
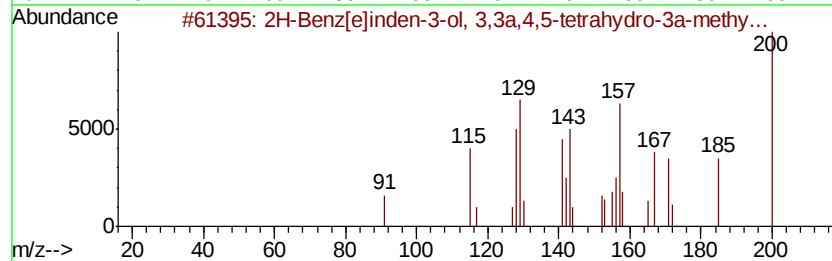
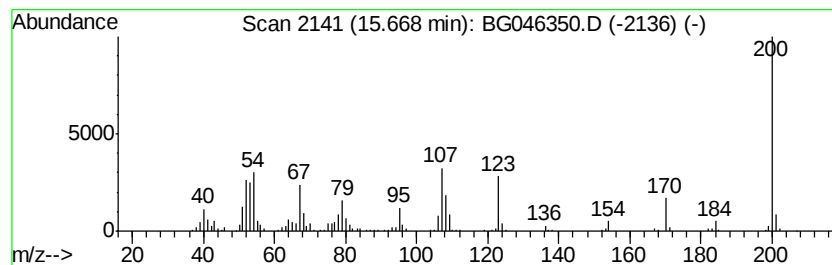
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-08 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.67	7.52 ng/ul	428329	Acenaphthene-d10	14.57

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2H-Benz[e]inden-3-ol, 3,3a,4,5-t...	200	C14H16O	071805-91-9	35
2			Benzene, 1-methoxy-4-phenoxy-	200	C13H12O2	001655-69-2	22
3			1,6-Dioxaspiro[4.4]non-3-ene, 2-...	200	C13H12O2	016863-61-9	16
4			2-Aminocaprylic acid, N-propoxyc...	469	C28H55NO4	1000325-43-2	14
5			1-Hydroxy-4-methyl-2-oxo-1,2-dih...	200	C11H8N2O2	021771-74-4	10



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
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Misc :
ALS Vial : 14 Sample Multiplier: 1

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ClientSampleId :
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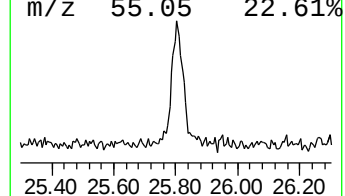
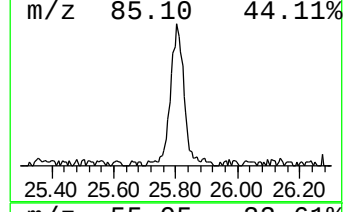
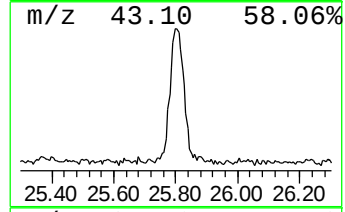
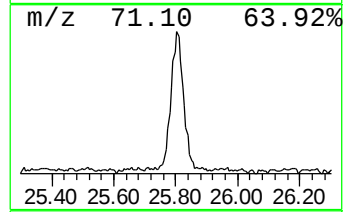
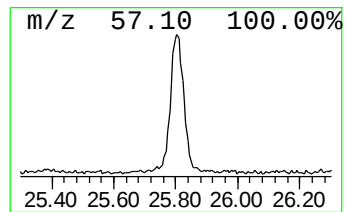
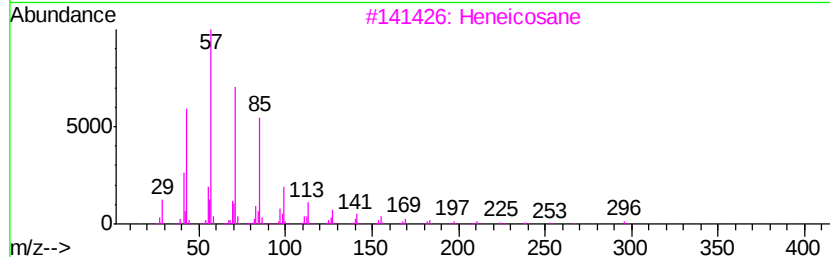
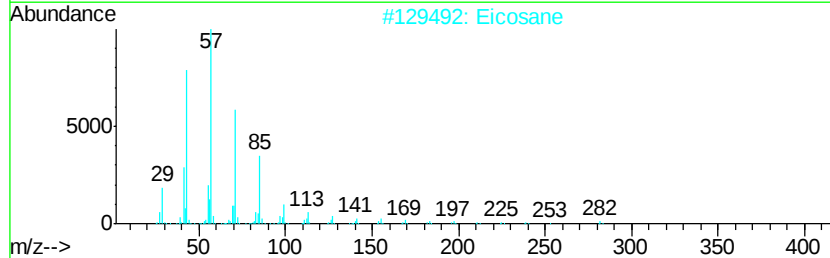
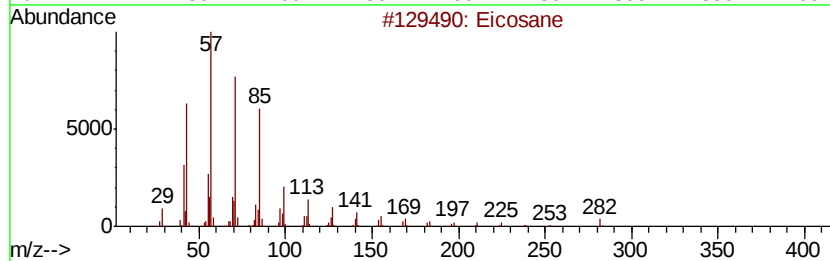
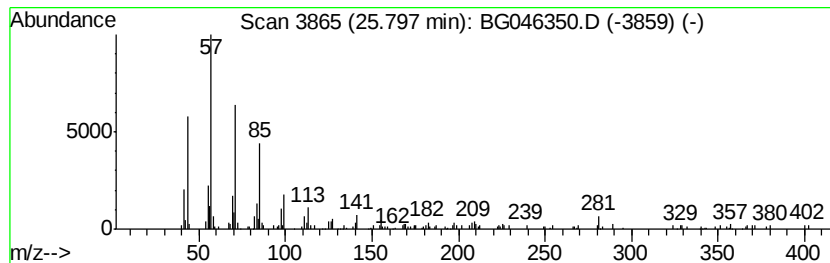
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.80	2.04 ng/ul	177355	Perylene-d12	24.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane			282	C20H42	000112-95-8	94
2	Eicosane			282	C20H42	000112-95-8	91
3	Heneicosane			296	C21H44	000629-94-7	87
4	Octacosane			394	C28H58	000630-02-4	87
5	Pentacosane			352	C25H52	000629-99-2	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG081920\
Data File : BG046350.D
Acq On : 19 Aug 2020 17:24
Operator : CG/JU
Sample : PB130882BL
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK82

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.14	2.0	ng/ul	52967	1	7.94	528248	20.0
2-Furancarboxylic...	7.11	23.2	ng/ul	612441	1	7.94	528248	20.0
unknown-01	7.27	17.5	ng/ul	463248	1	7.94	528248	20.0
unknown-02	7.48	23.7	ng/ul	625541	1	7.94	528248	20.0
unknown-03	8.65	29.0	ng/ul	765951	1	7.94	528248	20.0
1H-Imidazole, 4-m...	9.09	22.0	ng/ul	580859	1	7.94	528248	20.0
unknown-04	9.82	12.5	ng/ul	500430	2	10.74	800168	20.0
unknown-05	10.87	19.3	ng/ul	771053	2	10.74	800168	20.0
unknown-06	13.97	20.7	ng/ul	1181760	3	14.57	1139430	20.0
unknown-07	14.76	8.7	ng/ul	493976	3	14.57	1139430	20.0
unknown-08	15.67	7.5	ng/ul	428329	3	14.57	1139430	20.0
(DEL) Alkane: Str...	25.80	2.0	ng/ul	177355	6	24.66	1738570	20.0