

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
 Data File : BG046418.D
 Acq On : 21 Aug 2020 18:01
 Operator : CG/JU
 Sample : PB131008BL
 Misc :
 ALS Vial : 82 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK08

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	6	10	24	rVB	50498	92933	3.93%	0.385%
2	3.399	49	53	61	rVB	20768	35506	1.50%	0.147%
3	4.145	177	180	187	rVB5	4199	7938	0.34%	0.033%
4	7.118	678	686	700	rBV	357398	631970	26.70%	2.616%
5	7.271	703	712	721	rBV	272147	461608	19.50%	1.911%
6	7.476	740	747	762	rVB	369722	653529	27.61%	2.705%
7	7.946	819	827	837	rVB	278080	488822	20.65%	2.024%
8	8.651	940	947	960	rBV	455106	791851	33.46%	3.278%
9	9.098	1015	1023	1036	rBV	334151	597627	25.25%	2.474%
10	9.821	1138	1146	1159	rBV	287886	517789	21.88%	2.144%
11	10.367	1231	1239	1253	rBV	469636	854542	36.11%	3.538%
12	10.737	1295	1302	1310	rBV	415669	745786	31.51%	3.087%
13	10.872	1317	1325	1334	rBV	420697	790535	33.40%	3.273%
14	12.324	1568	1572	1579	rVB3	6960	13045	0.55%	0.054%
15	13.975	1846	1853	1866	rBV	809382	1191206	50.33%	4.931%
16	14.262	1895	1902	1916	rBV	970842	1539663	65.05%	6.374%
17	14.574	1947	1955	1963	rBV	691442	1075757	45.45%	4.453%
18	14.774	1983	1989	2006	rBV	264593	483920	20.45%	2.003%
19	15.567	2116	2124	2137	rBV	1231710	1889574	79.84%	7.822%
20	15.684	2138	2144	2155	rVV	197292	311253	13.15%	1.289%
21	15.802	2158	2164	2170	rVB	15253	23460	0.99%	0.097%
22	16.342	2252	2256	2263	rVB4	5392	7787	0.33%	0.032%
23	17.000	2363	2368	2369	rBV3	2743	3409	0.14%	0.014%
24	17.094	2381	2384	2385	rBV2	3244	2660	0.11%	0.011%
25	17.312	2415	2421	2432	rBV	820396	1280432	54.10%	5.301%
26	17.412	2432	2438	2454	rBV2	1300650	2020097	85.35%	8.363%
27	17.564	2457	2464	2474	rBV3	21092	59955	2.53%	0.248%
28	17.688	2482	2485	2488	rBV5	2338	3261	0.14%	0.013%
29	17.905	2520	2522	2525	rVV4	3380	3056	0.13%	0.013%
30	17.929	2525	2526	2531	rVB5	2272	2430	0.10%	0.010%
31	18.264	2581	2583	2584	rBV2	3372	3094	0.13%	0.013%
32	18.416	2608	2609	2612	rBV3	2632	3085	0.13%	0.013%
33	18.499	2621	2623	2625	rVV3	4346	2544	0.11%	0.011%
34	18.599	2637	2640	2641	rBV3	4099	4315	0.18%	0.018%

LSC Area Percent Report

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35	18.628	2644	2645	2646	rBV	4955	2656	0.11%	0.011%
36	19.333	2762	2765	2768	rBV	18967	23435	0.99%	0.097%
37	19.591	2805	2809	2812	rBV	19095	33248	1.40%	0.138%
38	19.703	2821	2828	2834	rBV	1615131	2366747	100.00%	9.798%
39	21.566	3139	3145	3151	rVB2	854496	1377688	58.21%	5.703%
40	24.468	3630	3639	3653	rVB	876551	2316008	97.86%	9.588%
41	24.680	3666	3675	3688	rVB2	489766	1441664	60.91%	5.968%

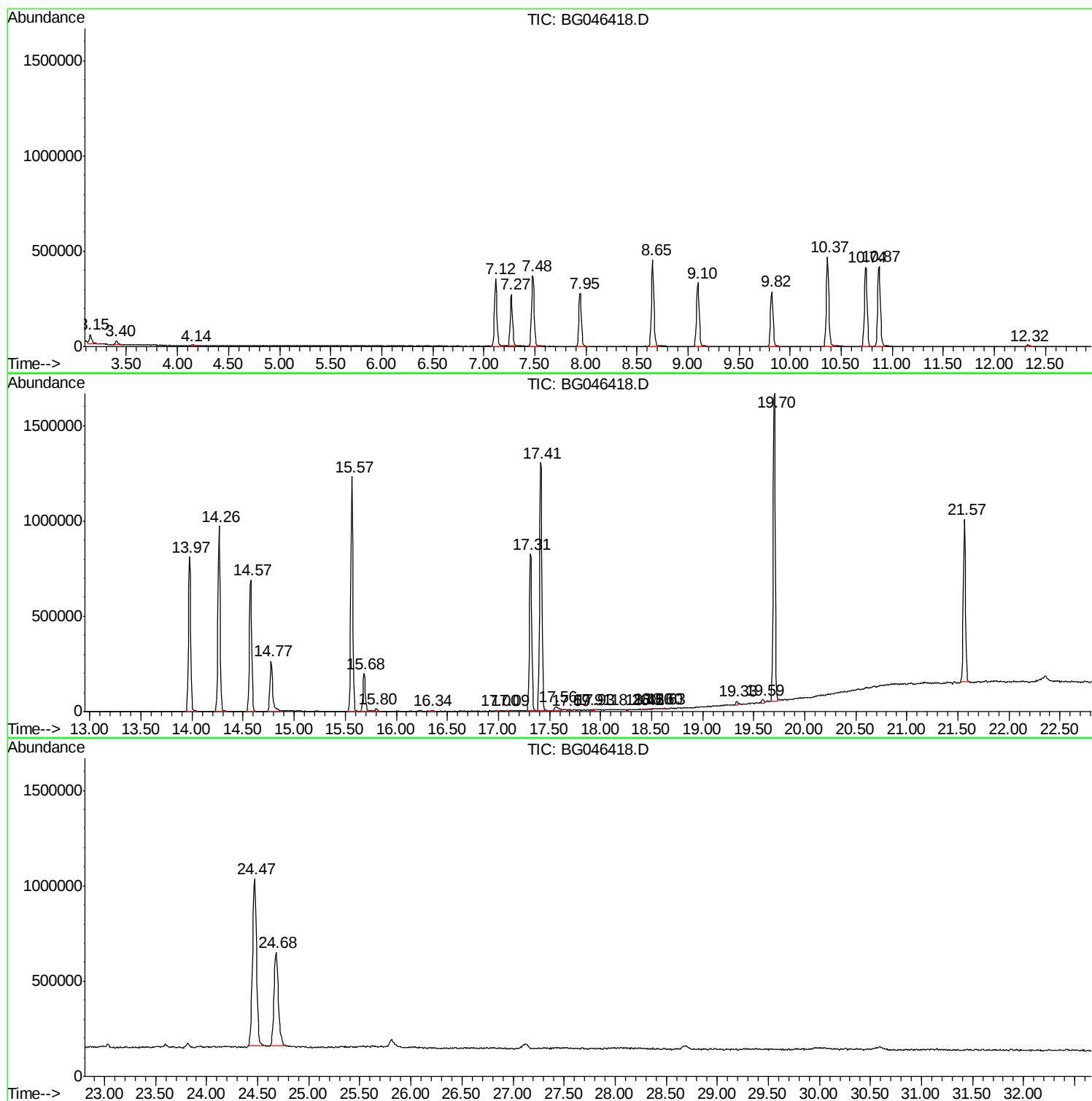
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG081320PAH.M
Quant Title : SVOA CALIBRATION

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



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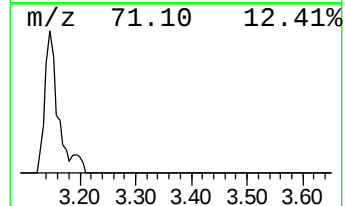
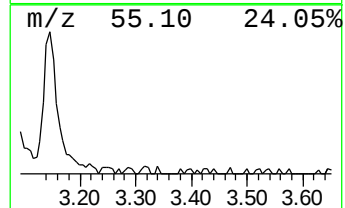
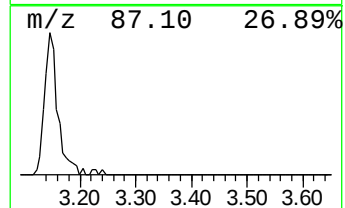
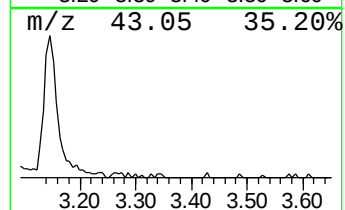
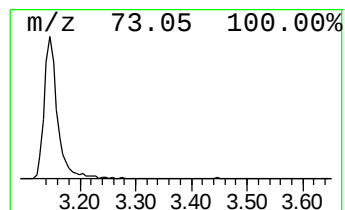
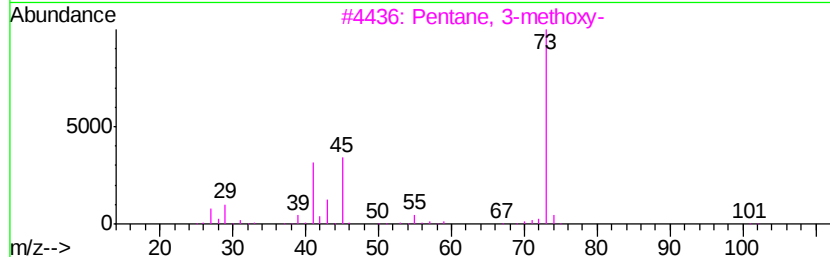
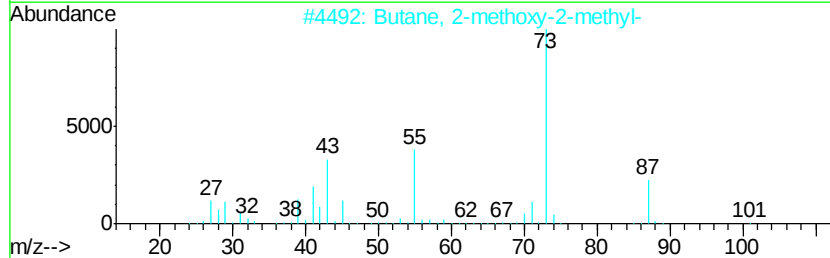
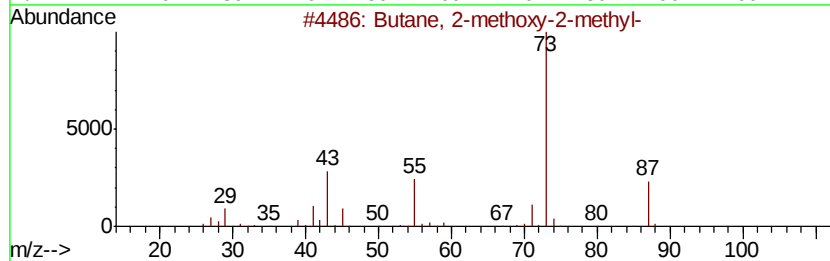
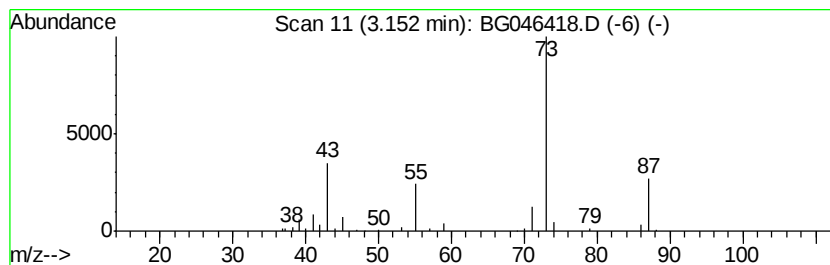
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.15	3.80 ng/ul	92933	1,4-Dichlorobenzene-d4	7.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	10
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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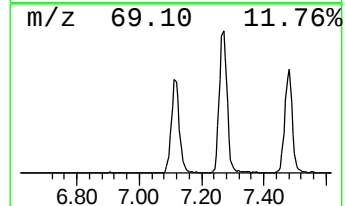
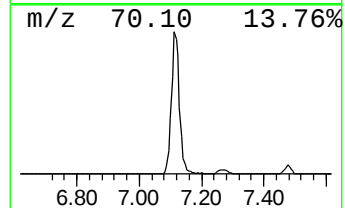
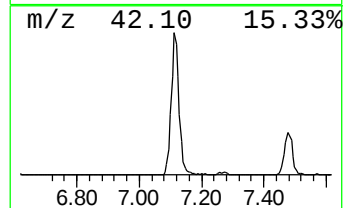
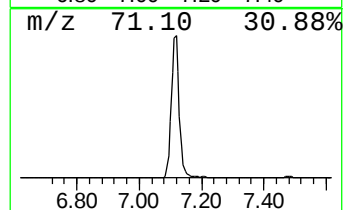
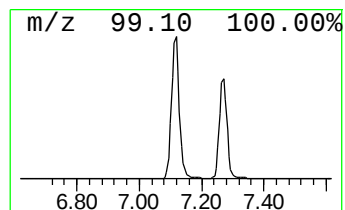
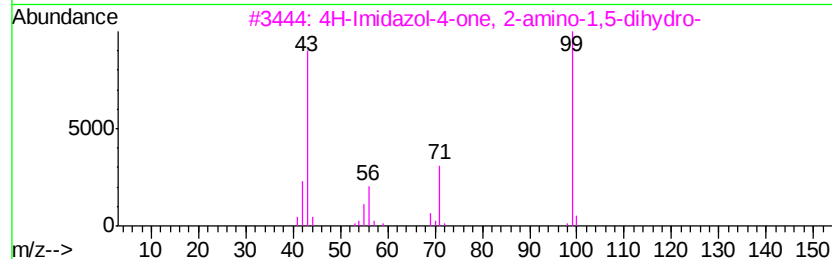
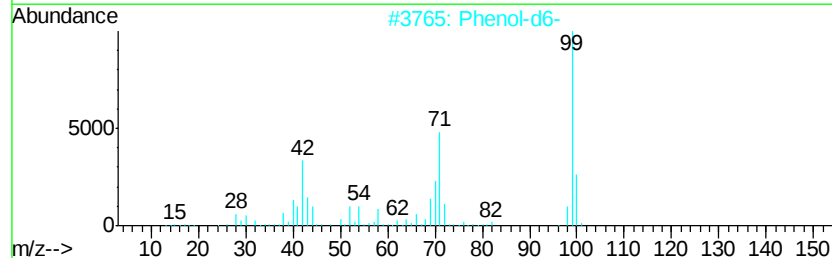
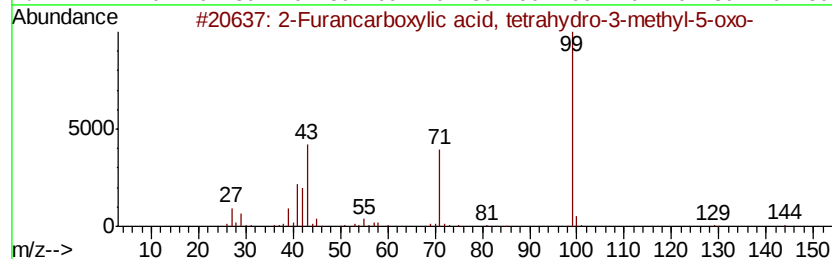
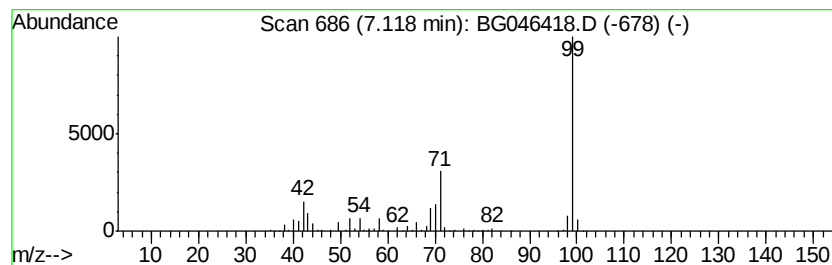
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 2-Furancarboxylic acid, tet... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.12	25.86 ng/ul	631970	1,4-Dichlorobenzene-d4	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Furancarboxylic acid, tetrahyd...	144	C6H8O4	022073-04-7	59
2		Phenol-d6-	100	C6D6O	013127-88-3	50
3		4H-Imidazol-4-one, 2-amino-1,5-d...	99	C3H5N3O	000503-86-6	45
4		n-Heptyl methylphosphonofluoridate	196	C8H18F02P	162085-82-7	42
5		Dihydropyrimidine-4,6-dione, 5-[...	373	C18H23N5O2S	1000304-28-2	40



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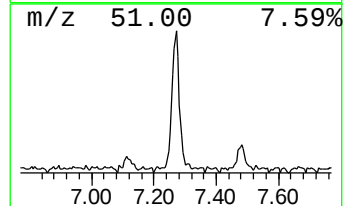
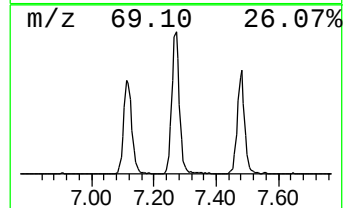
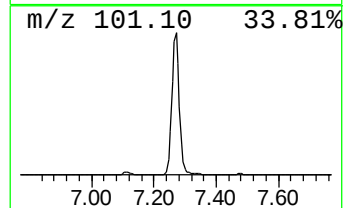
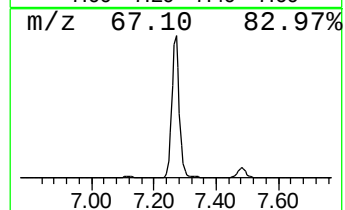
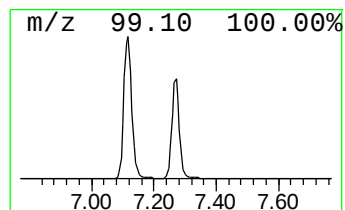
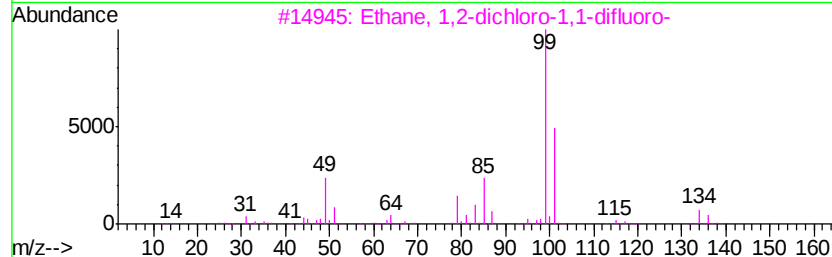
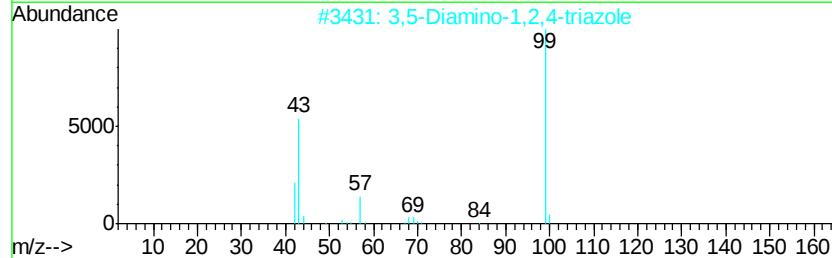
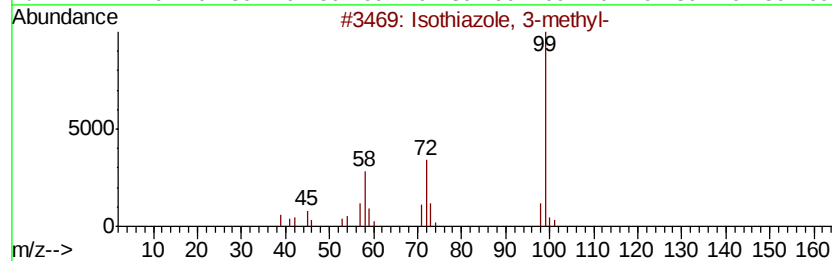
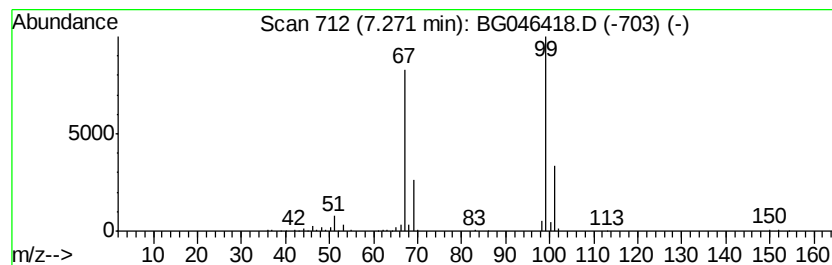
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.27	18.89 ng/ul	461608	1,4-Dichlorobenzene-d4	7.95

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Isothiazole, 3-methyl-	99	C4H5NS	000693-92-5	9
2		3,5-Diamino-1,2,4-triazole	99	C2H5N5	001455-77-2	9
3		Ethane, 1,2-dichloro-1,1-difluoro-	134	C2H2Cl2F2	001649-08-7	9
4		1-Butyne, 3,3-dimethyl-	82	C6H10	000917-92-0	9
5		1-Pentanamine, 5-(methylthio)-	133	C6H15NS	055021-78-8	9



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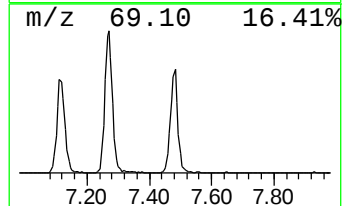
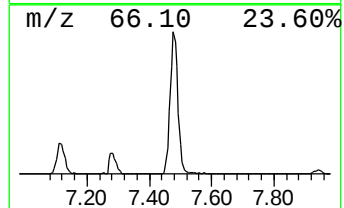
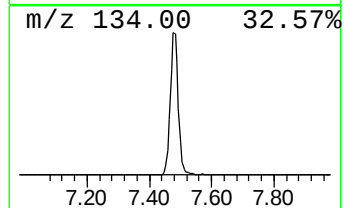
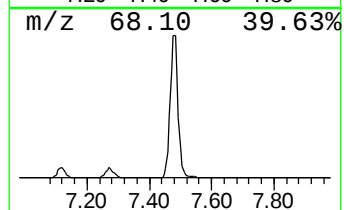
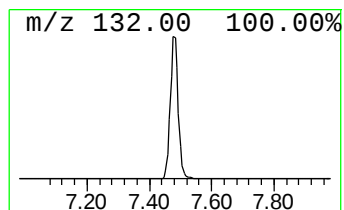
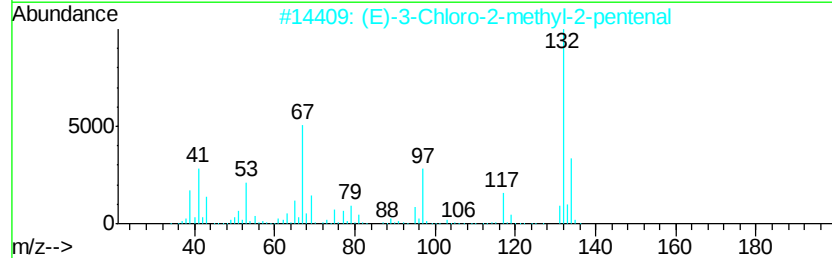
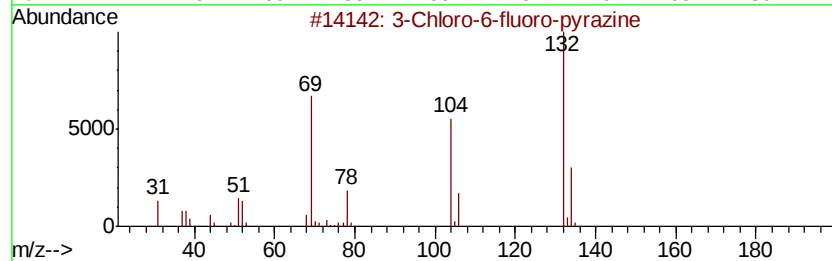
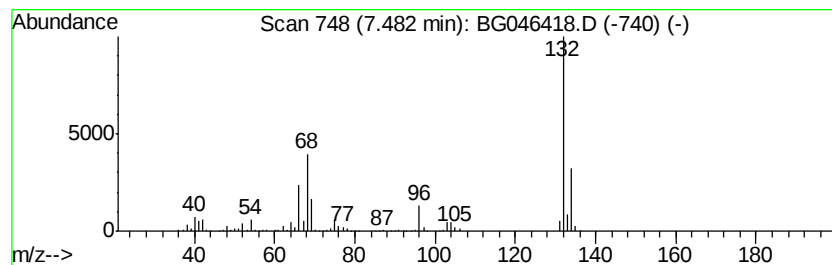
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Peak Number 4 unknown-02 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.48	26.74 ng/ul	653529	1,4-Dichlorobenzene-d4	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Tranvlcyproline-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



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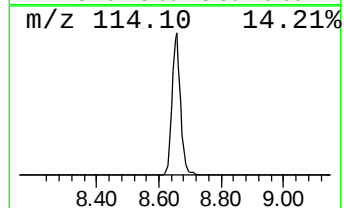
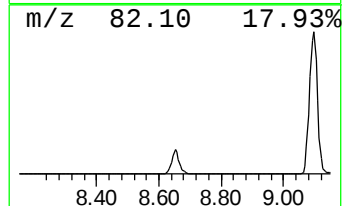
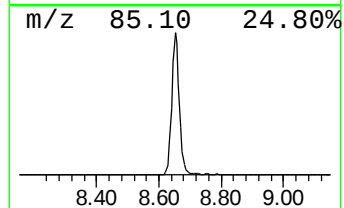
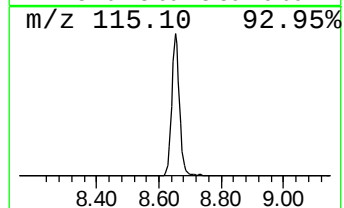
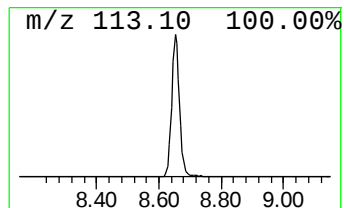
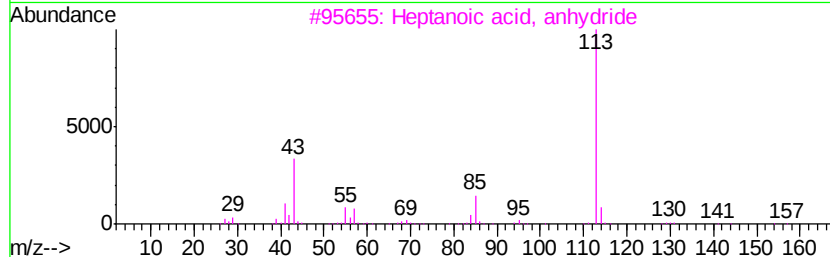
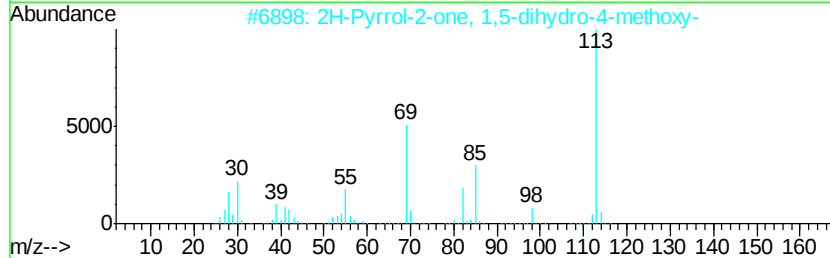
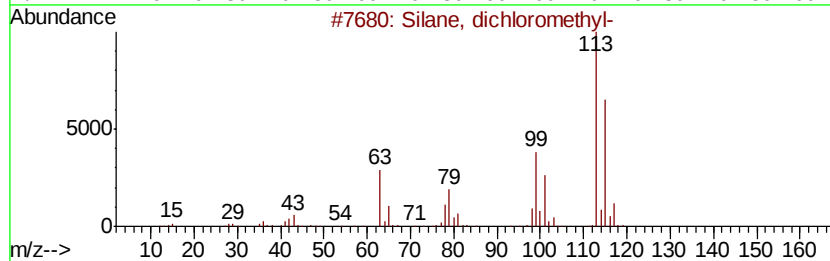
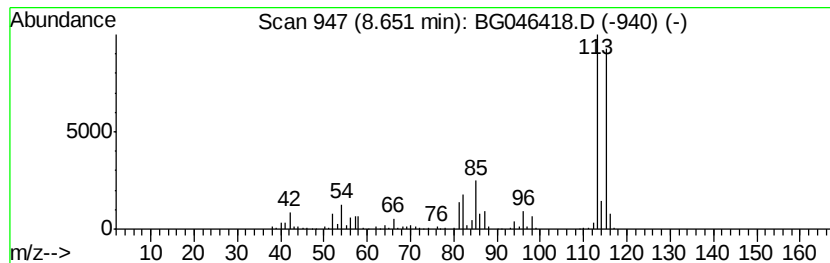
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Quant Title : SVOA CALIBRATION

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	32.40 ng/ul	791851	1,4-Dichlorobenzene-d4	7.95

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Silane, dichloromethyl-	114	CH4Cl2Si	000075-54-7	38
2		2H-Pyrrol-2-one, 1,5-dihydro-4-m...	113	C5H7NO2	069778-83-2	32
3		Heptanoic acid, anhydride	242	C14H26O3	000626-27-7	27
4		Heptanoic acid, anhydride	242	C14H26O3	000626-27-7	25
5		3,4-Methylpropylsuccinimide	155	C8H13NO2	1000248-78-0	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046418.D
Acq On : 21 Aug 2020 18:01
Operator : CG/JU
Sample : PB131008BL
Misc :
ALS Vial : 82 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK08

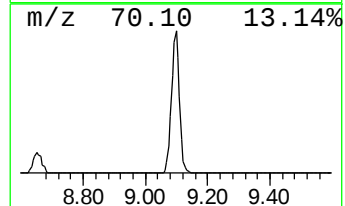
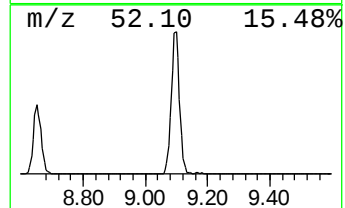
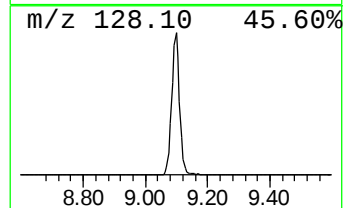
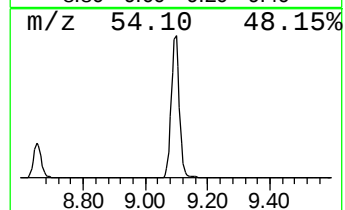
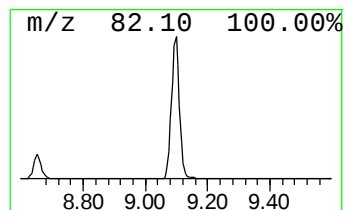
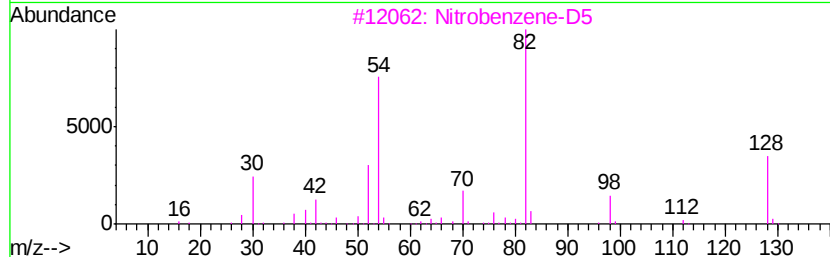
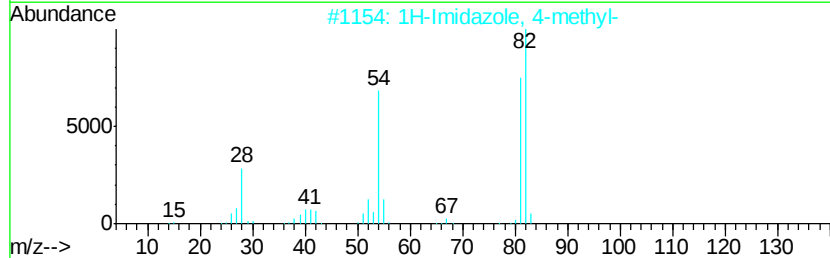
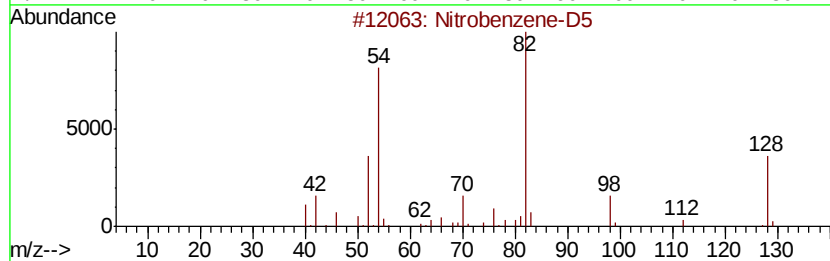
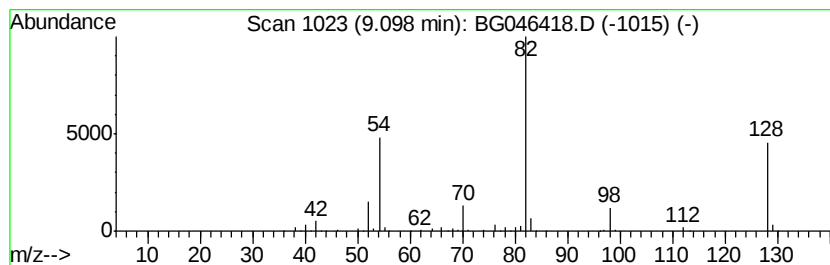
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Quant Title : SVOA CALIBRATION

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.
9.10	24.45 ng/ul	597627	1,4-Dichlorobenzene-d4	7.95

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	38
4		1H-Imidazole, 2-methyl-	82	C4H6N2	000693-98-1	38
5		1H-Imidazole, 4-methyl-	82	C4H6N2	000822-36-6	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046418.D
Acq On : 21 Aug 2020 18:01
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Sample : PB131008BL
Misc :
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Instrument :
BNA_G
ClientSampleId :
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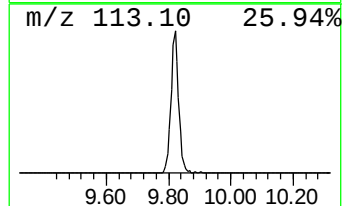
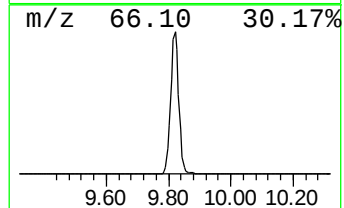
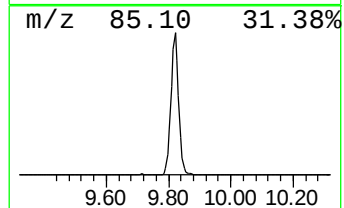
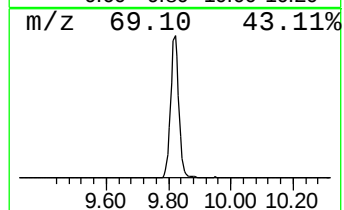
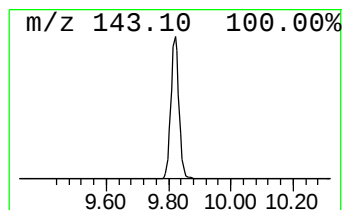
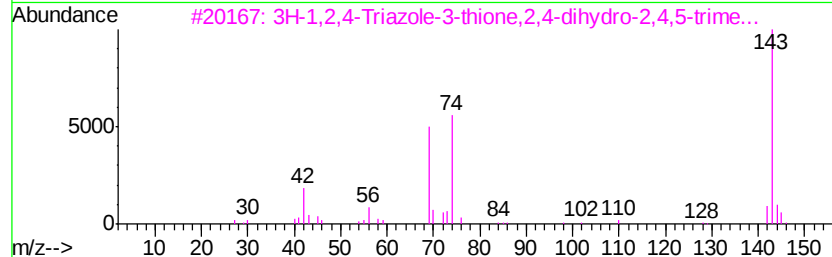
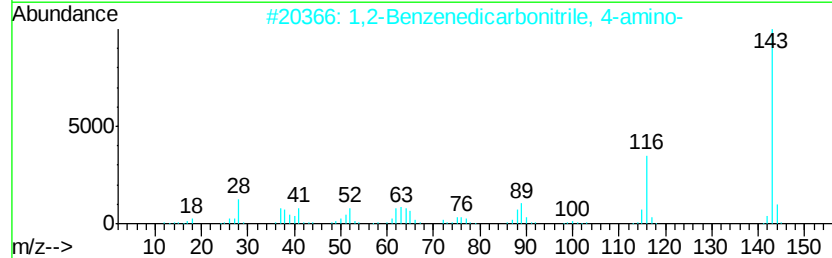
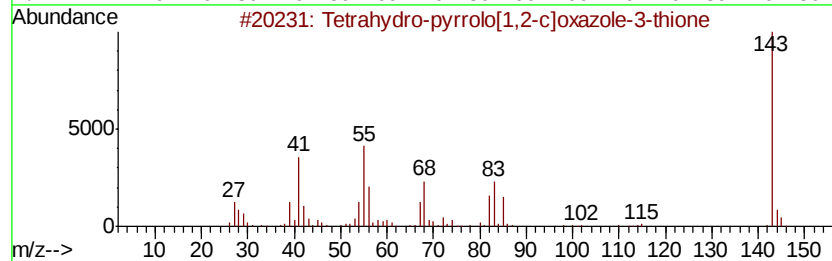
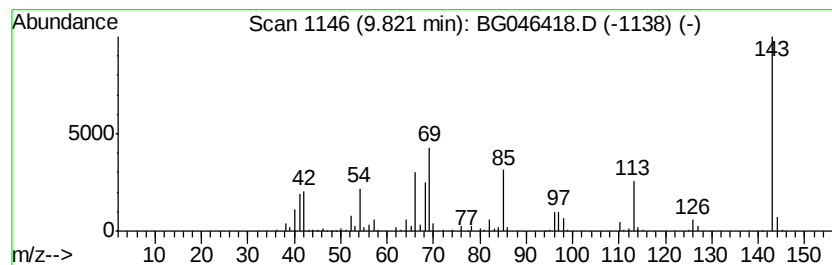
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-04 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.82	13.89 ng/ul	517789	Naphthalene-d8	10.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetrahydro-pyrrolo[1,2-c]oxazole...	143	C6H9NOS	1000190-53-6	17
2		1,2-Benzenedicarbonitrile, 4-amino-	143	C8H5N3	056765-79-8	14
3		3H-1,2,4-Triazole-3-thione,2,4-d...	143	C5H9N3S	037526-42-4	14
4		4(1H)-Pyrimidinone, 6-amino-2,3-...	143	C4H5N3OS	001004-40-6	12
5		1-(1-Butoxypropan-2-yloxy)propan...	274	C15H30O4	1000367-12-9	12



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046418.D
Acq On : 21 Aug 2020 18:01
Operator : CG/JU
Sample : PB131008BL
Misc :
ALS Vial : 82 Sample Multiplier: 1

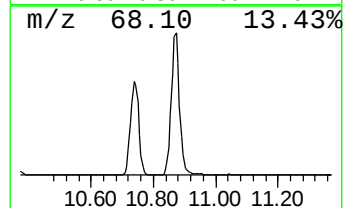
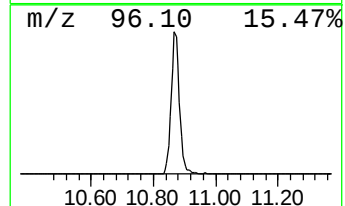
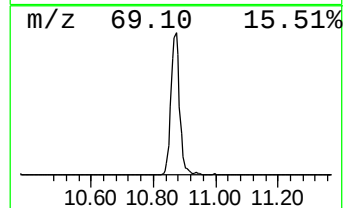
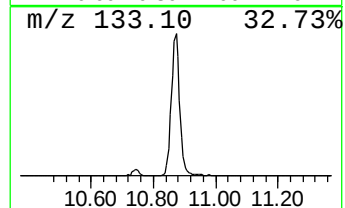
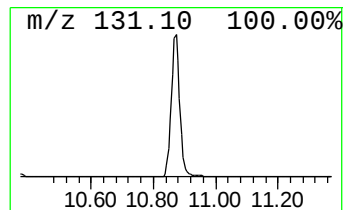
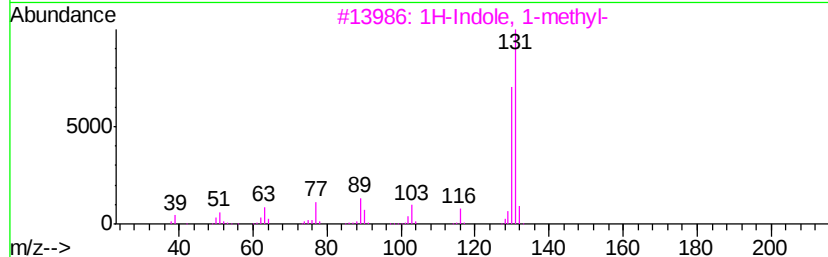
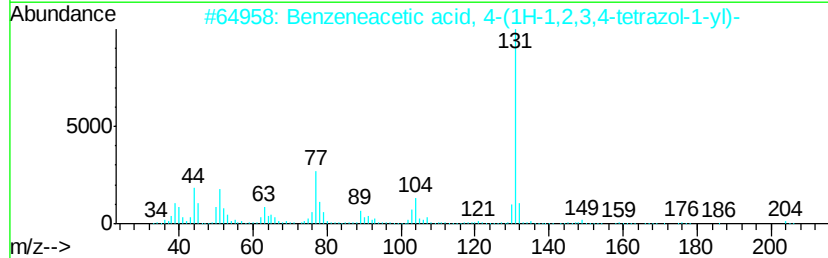
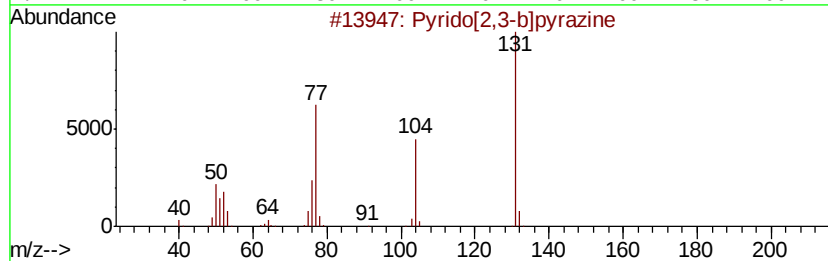
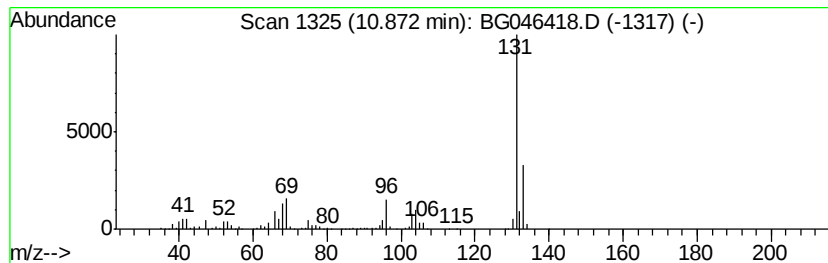
Instrument :
BNA_G
ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-05 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.87	21.20 ng/ul	790535	Naphthalene-d8	10.74
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Pyrido[2,3-b]pyrazine	131 C7H5N3	000322-46-3 40
2		Benzeneacetic acid, 4-(1H-1,2,3,...	204 C9H8N4O2	1000351-56-5 33
3		1H-Indole, 1-methyl-	131 C9H9N	000603-76-9 28
4		Pyrido[2,3-b]pyrazine	131 C7H5N3	000322-46-3 9
5		4-Aminobenzyl cyanide	132 C8H8N2	003544-25-0 9



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
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Operator : CG/JU
Sample : PB131008BL
Misc :
ALS Vial : 82 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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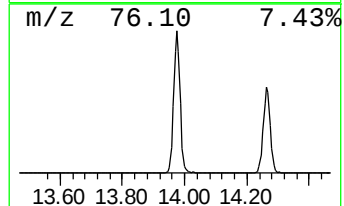
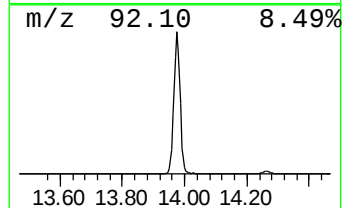
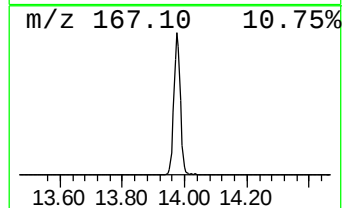
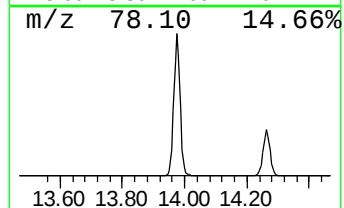
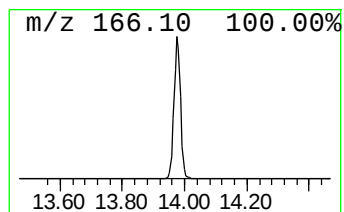
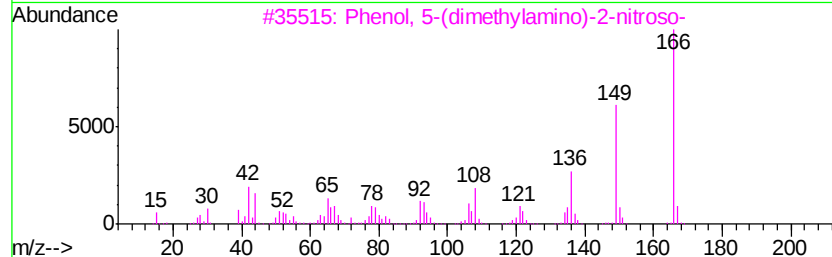
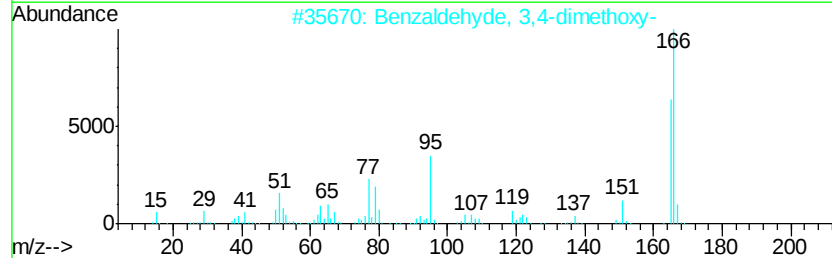
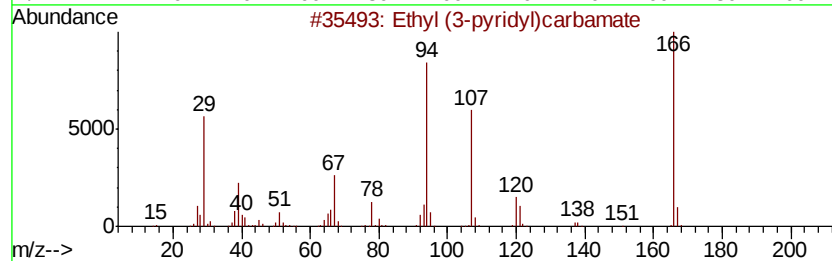
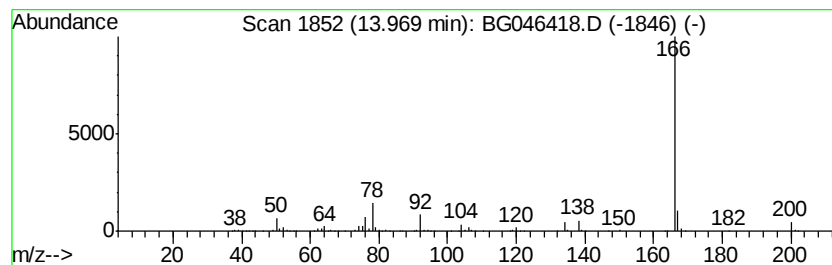
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-06 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	22.15 ng/ul	1191210	Acenaphthene-d10	14.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethyl (3-pyridyl)carbamate	166	C8H10N2O2	006276-11-5	39
2		Benzaldehyde, 3,4-dimethoxy-	166	C9H10O3	000120-14-9	38
3		Phenol, 5-(dimethylamino)-2-nitr...	166	C8H10N2O2	016761-04-9	38
4		Benzenamine, N,N-dimethyl-4-nitro-	166	C8H10N2O2	000100-23-2	38
5		3,5-Diamino-2-methylbenzoic acid	166	C8H10N2O2	090007-30-0	36



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046418.D
Acq On : 21 Aug 2020 18:01
Operator : CG/JU
Sample : PB131008BL
Misc :
ALS Vial : 82 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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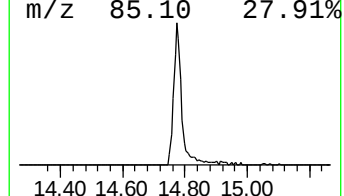
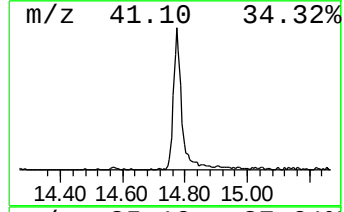
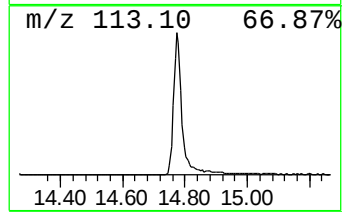
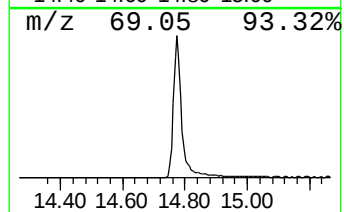
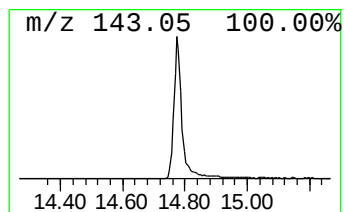
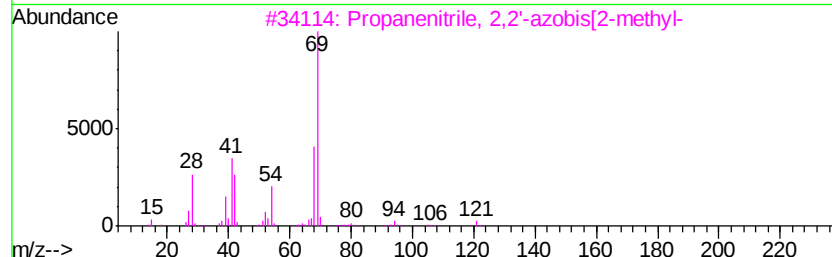
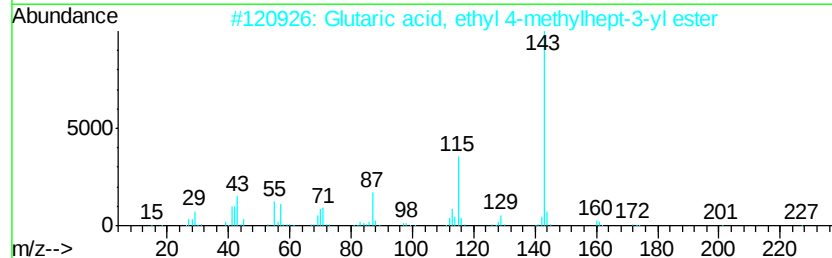
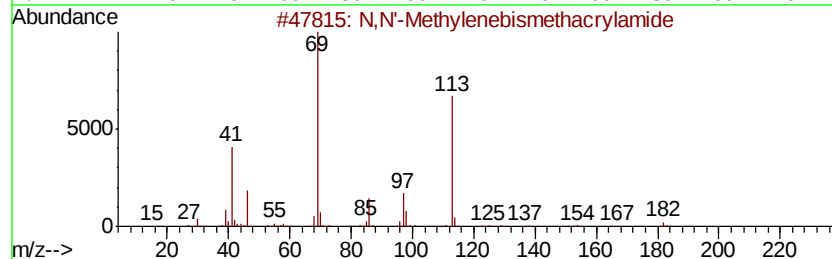
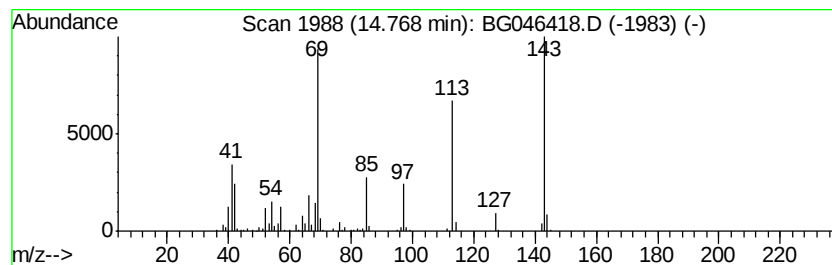
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-07 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.77	9.00 ng/ul	483920	Acenaphthene-d10	14.57

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N,N'-Methylenebismethacrylamide	182	C9H14N2O2	002359-15-1	25
2			Glutaric acid, ethyl 4-methylhept...	272	C15H28O4	1000377-14-9	22
3			Propanenitrile, 2,2'-azobis[2-me...	164	C8H12N4	000078-67-1	22
4			4-Amino-2-chloro-6-methylpyrimidine	143	C5H6ClN3	014394-60-6	16
5			1-Cyclopentylethanol, trifluoroa...	210	C9H13F3O2	1000364-11-9	12



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046418.D
Acq On : 21 Aug 2020 18:01
Operator : CG/JU
Sample : PB131008BL
Misc :
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Instrument :
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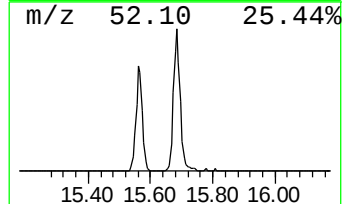
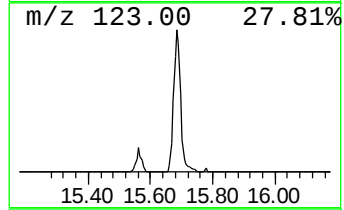
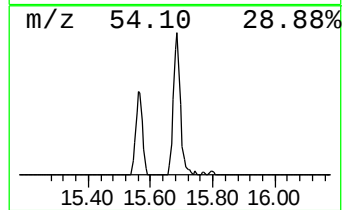
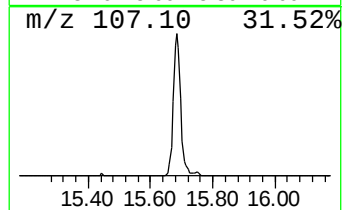
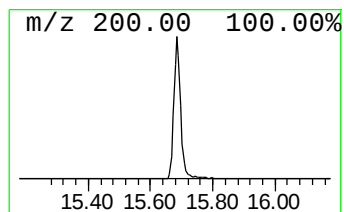
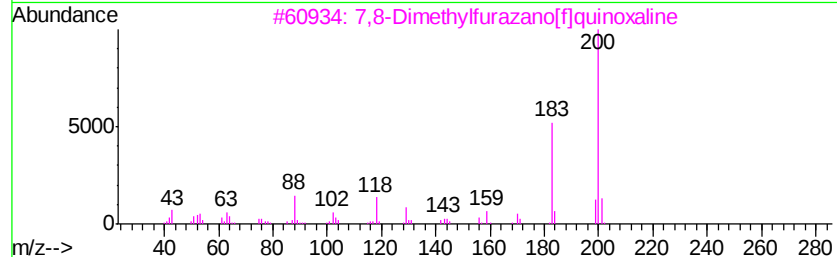
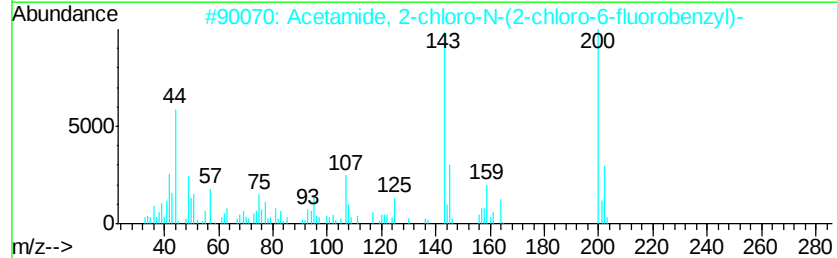
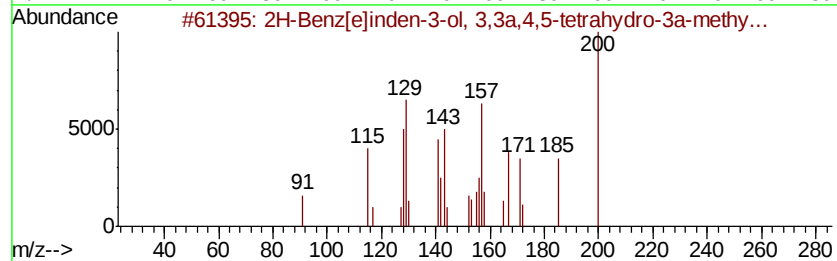
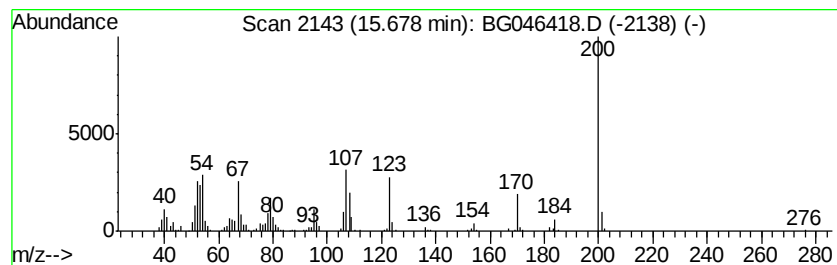
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Quant Title : SVOA CALIBRATION

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.
15.68	5.79 ng/ul	311253	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	38
2			Acetamide, 2-chloro-N-(2-chloro-...	235	C9H8Cl2FNO	1000268-05-5	16
3			7,8-Dimethylfurazano[f]quinoxaline	200	C10H8N4O	1000110-74-6	14
4			Benzene, 1-methoxy-4-phenoxy-	200	C13H12O2	001655-69-2	14
5			1-Hydroxy-4-methyl-2-oxo-1,2-dih...	200	C11H8N2O2	021771-74-4	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG081920\
Data File : BG046418.D
Acq On : 21 Aug 2020 18:01
Operator : CG/JU
Sample : PB131008BL
Misc :
ALS Vial : 82 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...	3.15	3.8	ng/ul	92933	1	7.95	488822	20.0
2-Furancarboxylic...	7.12	25.9	ng/ul	631970	1	7.95	488822	20.0
unknown-01	7.27	18.9	ng/ul	461608	1	7.95	488822	20.0
unknown-02	7.48	26.7	ng/ul	653529	1	7.95	488822	20.0
unknown-03	8.65	32.4	ng/ul	791851	1	7.95	488822	20.0
1H-Imidazole, 4-m...	9.10	24.4	ng/ul	597627	1	7.95	488822	20.0
unknown-04	9.82	13.9	ng/ul	517789	2	10.74	745786	20.0
unknown-05	10.87	21.2	ng/ul	790535	2	10.74	745786	20.0
unknown-06	13.97	22.1	ng/ul	1191210	3	14.57	1075760	20.0
unknown-07	14.77	9.0	ng/ul	483920	3	14.57	1075760	20.0
unknown-08	15.68	5.8	ng/ul	311253	3	14.57	1075760	20.0