

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\
 Data File : BG046250.D
 Acq On : 13 Aug 2020 15:39
 Operator : CG/JU
 Sample : PB130901BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK01

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	3.164	9	13	24	rVB	42680	77085	2.41%	0.268%
2	3.299	34	36	39	rVB3	3544	3673	0.11%	0.013%
3	3.417	52	56	63	rVB	20474	32294	1.01%	0.112%
4	4.157	178	182	187	rBV4	3714	6189	0.19%	0.022%
5	7.112	675	685	702	rBV	375013	651168	20.33%	2.265%
6	7.277	706	713	726	rVV	312793	529864	16.54%	1.843%
7	7.483	741	748	758	rBV	385472	655538	20.47%	2.280%
8	7.947	821	827	840	rVB	283125	485664	15.16%	1.689%
9	8.652	940	947	963	rBV	467116	812697	25.38%	2.827%
10	9.098	1016	1023	1036	rBV	363739	640753	20.01%	2.229%
11	9.821	1139	1146	1160	rBV	303429	543279	16.96%	1.890%
12	10.367	1231	1239	1250	rBV	460170	844508	26.37%	2.938%
13	10.743	1295	1303	1311	rBV	404132	709427	22.15%	2.468%
14	10.873	1317	1325	1343	rBV	474732	884201	27.61%	3.076%
15	12.330	1566	1573	1580	rVB2	7805	13526	0.42%	0.047%
16	13.981	1847	1854	1863	rBV	1002924	1448641	45.23%	5.039%
17	14.269	1895	1903	1913	rBV	1116049	1778238	55.52%	6.186%
18	14.574	1948	1955	1963	rBV	669126	1056761	33.00%	3.676%
19	14.762	1980	1987	2007	rBV	306067	601869	18.79%	2.094%
20	15.567	2117	2124	2135	rBV	1513626	2270790	70.90%	7.899%
21	15.673	2137	2142	2153	rBV	395713	610328	19.06%	2.123%
22	15.808	2160	2165	2170	rVB2	18383	26145	0.82%	0.091%
23	16.349	2253	2257	2262	rVB3	5698	7983	0.25%	0.028%
24	16.842	2339	2341	2347	rVB4	2963	4178	0.13%	0.015%
25	16.995	2360	2367	2372	rVB5	2738	4350	0.14%	0.015%
26	17.101	2383	2385	2393	rVB4	3310	4651	0.15%	0.016%
27	17.312	2412	2421	2431	rBV	864671	1345198	42.00%	4.679%
28	17.412	2431	2438	2456	rVV	1638773	2479264	77.41%	8.624%
29	19.333	2760	2765	2770	rBV2	20777	28997	0.91%	0.101%
30	19.586	2803	2808	2812	rVB	18091	23239	0.73%	0.081%
31	19.698	2820	2827	2841	rBV	2134484	3124189	97.55%	10.868%
32	20.767	3007	3009	3011	rBV3	4300	3518	0.11%	0.012%
33	21.560	3137	3144	3156	rBV	1005946	1595269	49.81%	5.549%
34	21.778	3178	3181	3186	rVB2	22221	29545	0.92%	0.103%

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Integration Parameters: LSCINT.P

Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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Title : SVOA CALIBRATION

35	22.371	3277	3282	3290	rVB2	49855	87506	2.73%	0.304%
36	23.041	3391	3396	3403	rVB	59526	106199	3.32%	0.369%
37	23.816	3523	3528	3537	rVB	64745	129113	4.03%	0.449%
38	24.457	3626	3637	3654	rBV	1230588	3202626	100.00%	11.141%
39	24.662	3662	3672	3681	rBV2	597503	1651976	51.58%	5.747%
40	25.832	3863	3871	3880	rVB2	45941	132641	4.14%	0.461%
41	27.130	4086	4092	4102	rVB6	35332	104006	3.25%	0.362%

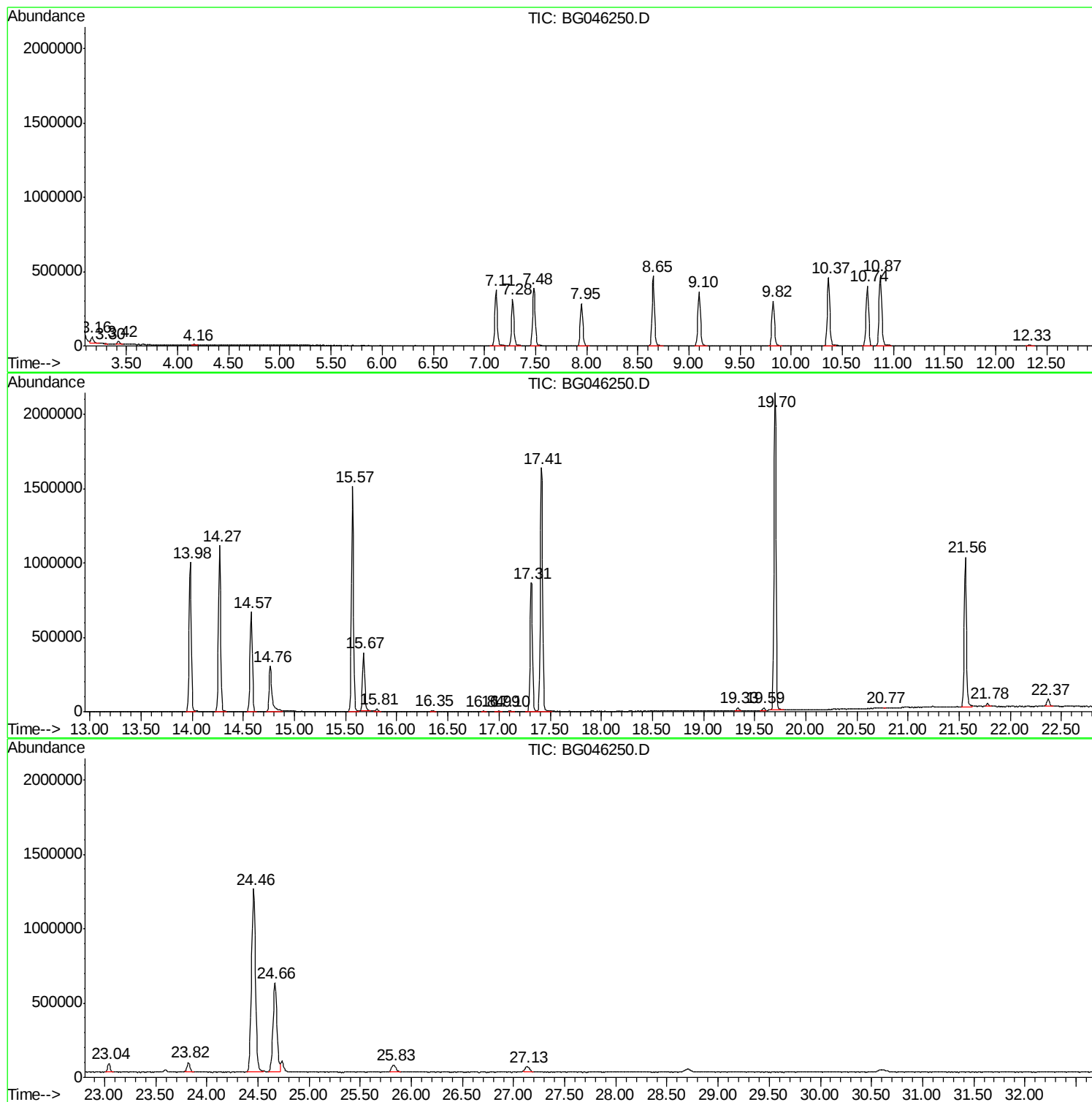
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG081320PAH.M
Quant Title : SVOA CALIBRATION

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



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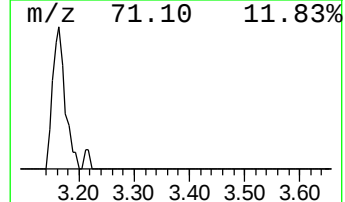
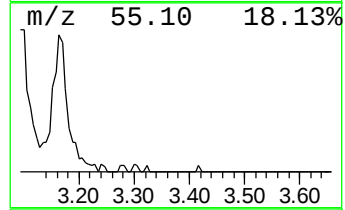
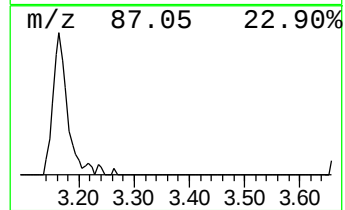
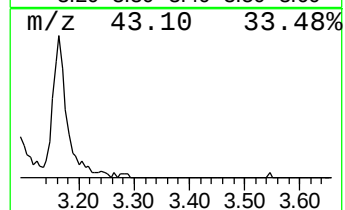
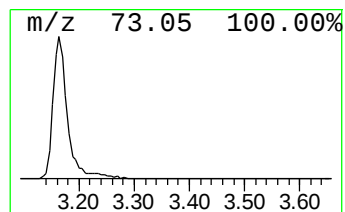
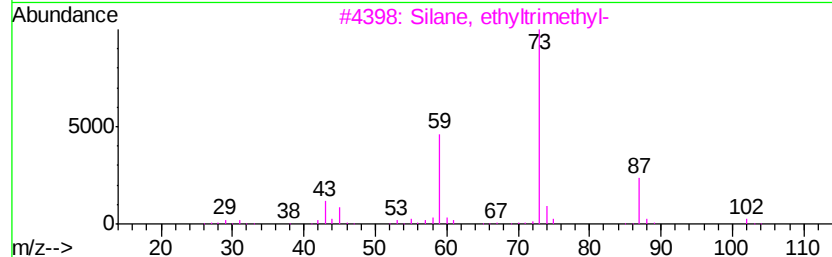
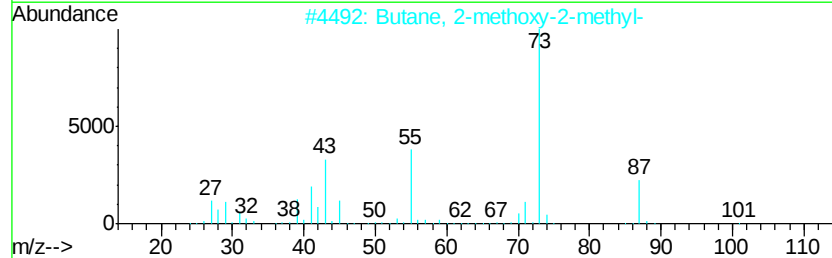
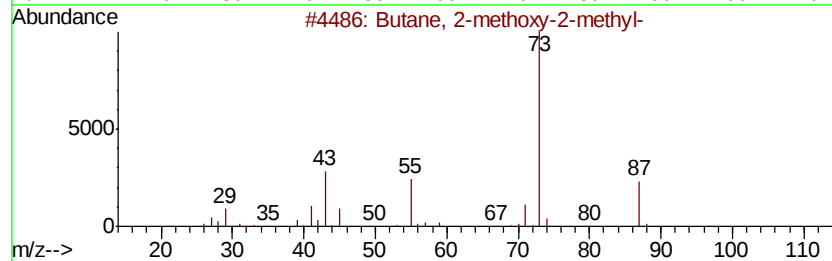
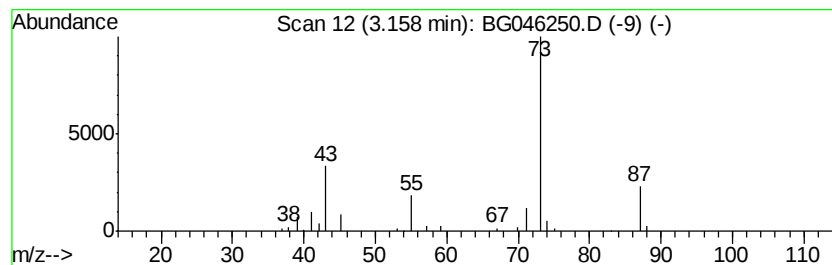
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG081320PAH.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.16	3.17 ng/ul	77085	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	56
3			Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	36
4			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	36
5			1,4-Butanediamine	88	C4H12N2	000110-60-1	9



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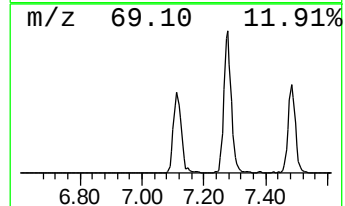
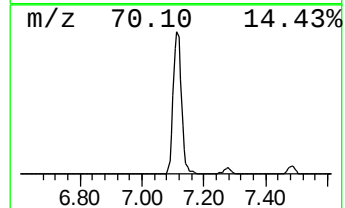
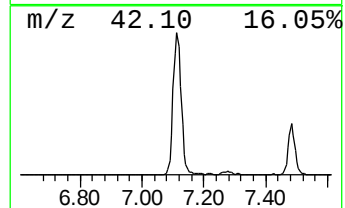
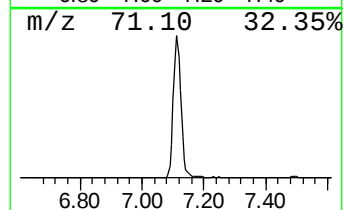
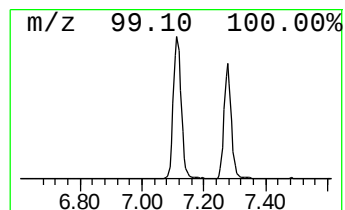
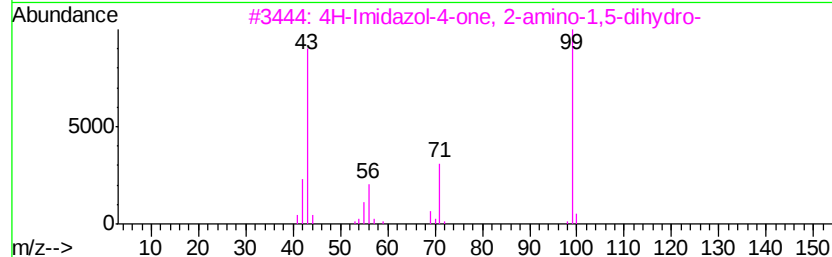
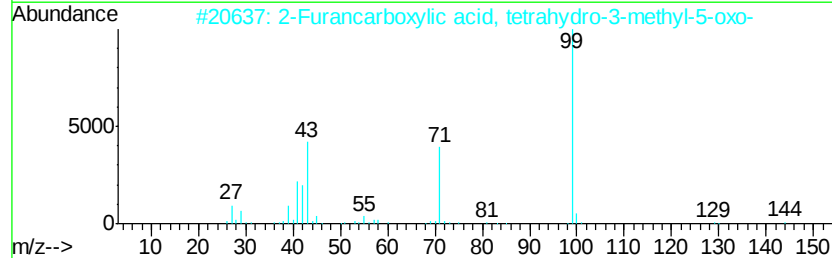
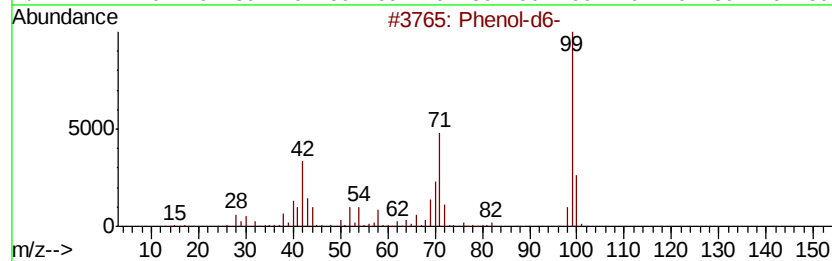
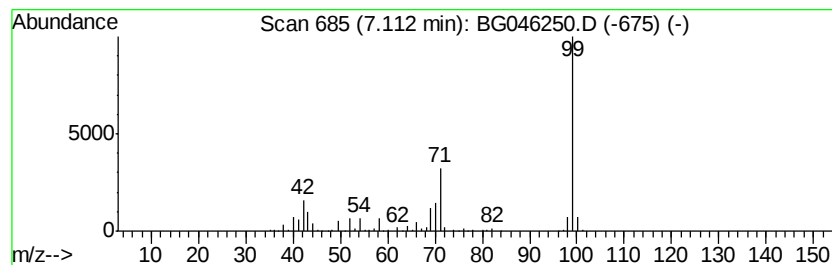
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Peak Number 2 2-Furancarboxylic acid, tet... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.11	26.82 ng/ul	651168	1,4-Dichlorobenzene-d4	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol-d6-	100	C6D6O	013127-88-3	64
2		2-Furancarboxylic acid, tetrahyd...	144	C6H8O4	022073-04-7	59
3		4H-Imidazol-4-one, 2-amino-1,5-d...	99	C3H5N3O	000503-86-6	59
4		2-Pyrrolidinone, 4-methyl-	99	C5H9NO	1000197-00-3	53
5		Hexanoic acid, anhydride	214	C12H22O3	002051-49-2	38



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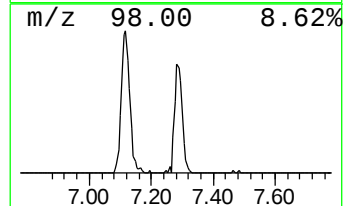
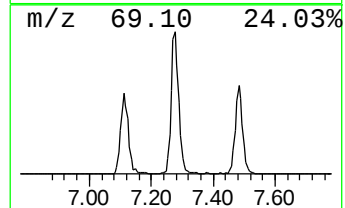
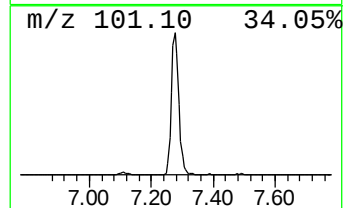
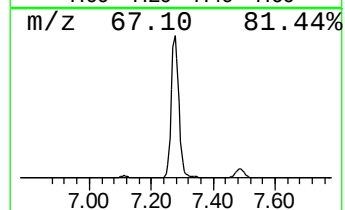
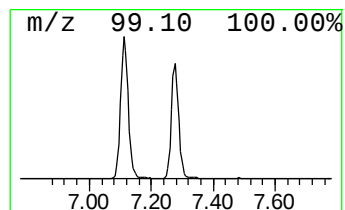
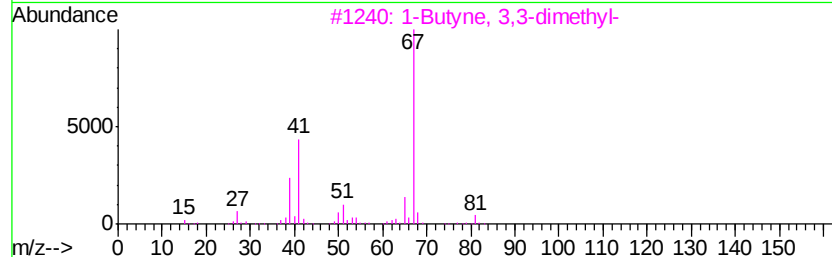
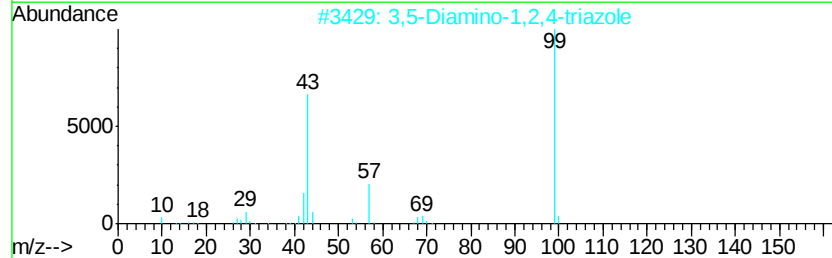
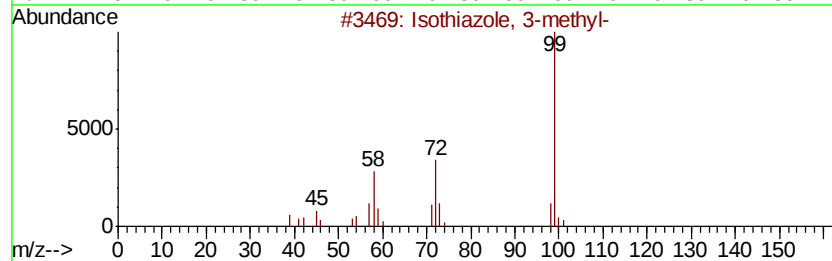
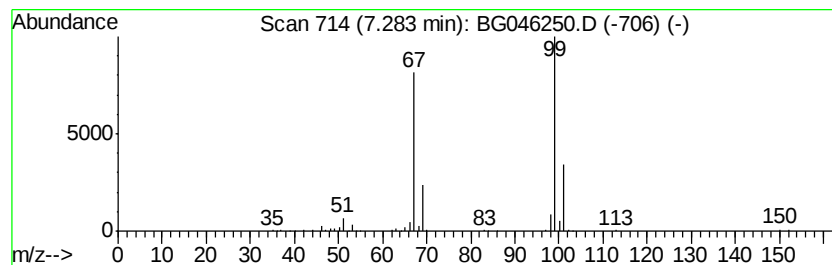
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG081320PAH.M
Quant Title : SVOA CALIBRATION

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.
7.28	21.82 ng/ul	529864	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			1-Butyne, 3,3-dimethyl-	82	C6H10	000917-92-0	9
4			Allyl Isothiocyanate	99	C4H5NS	000057-06-7	7
5			Cyclopropane, isothiocyanato-	99	C4H5NS	056601-42-4	7



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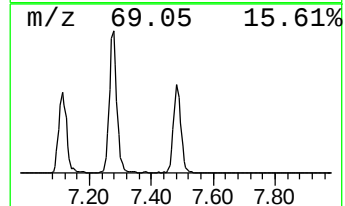
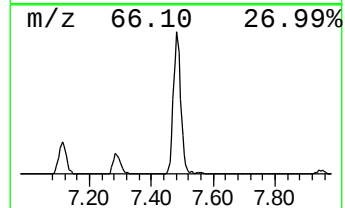
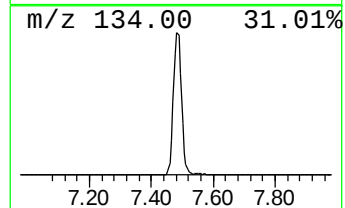
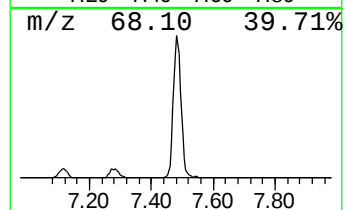
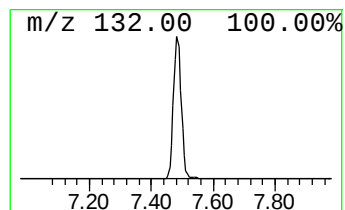
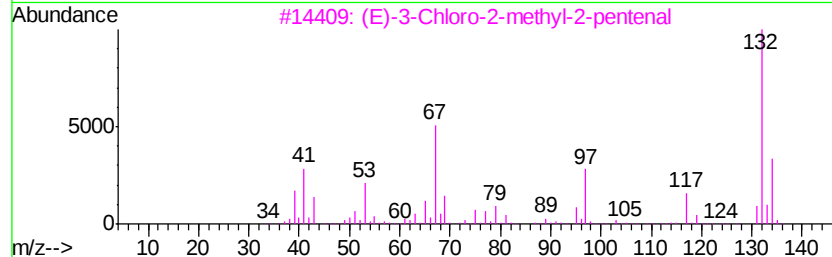
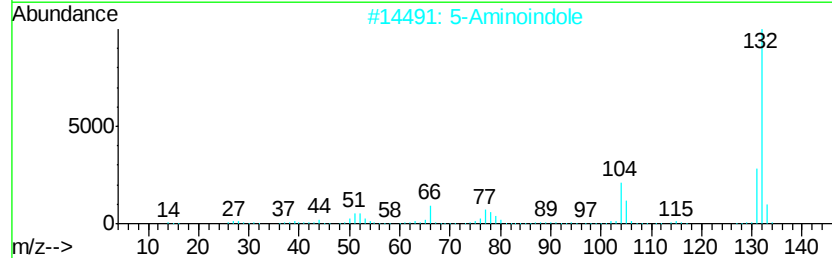
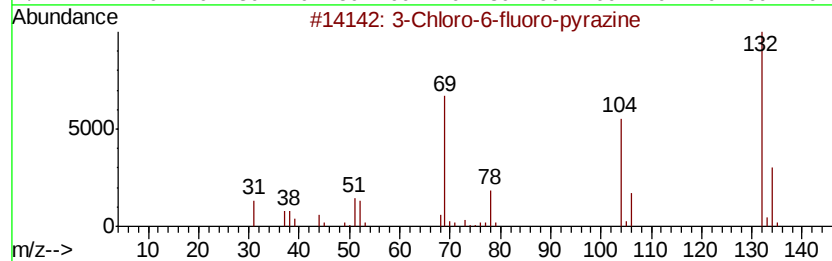
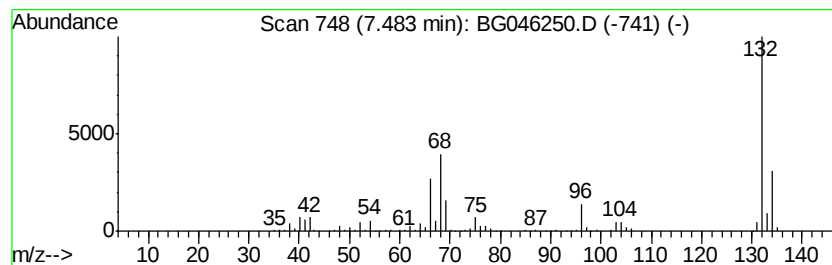
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown-02 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.48	27.00 ng/ul	655538	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		5-Aminoindole	132	C8H8N2	005192-03-0	22
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
4		Amphetaminil	250	C17H18N2	017590-01-1	12
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12



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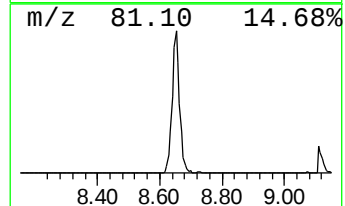
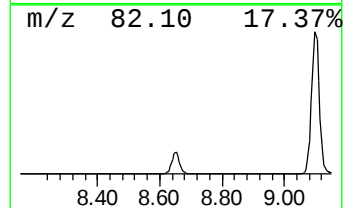
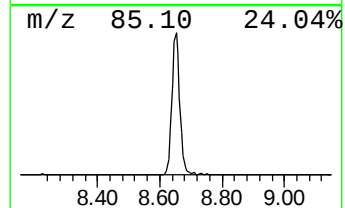
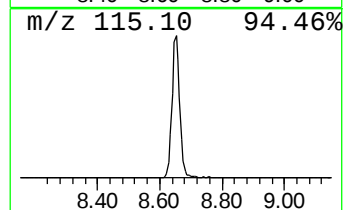
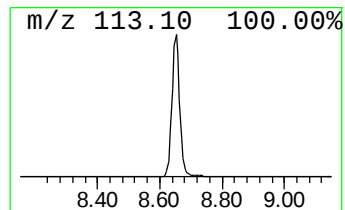
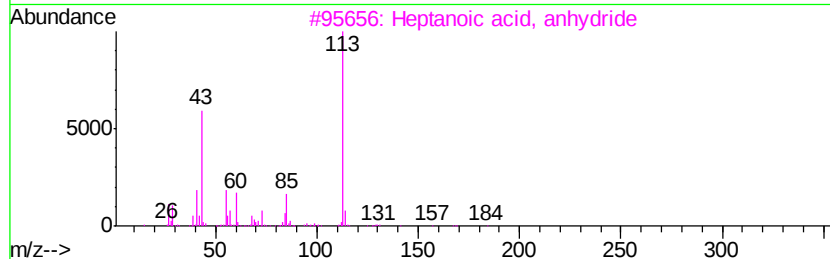
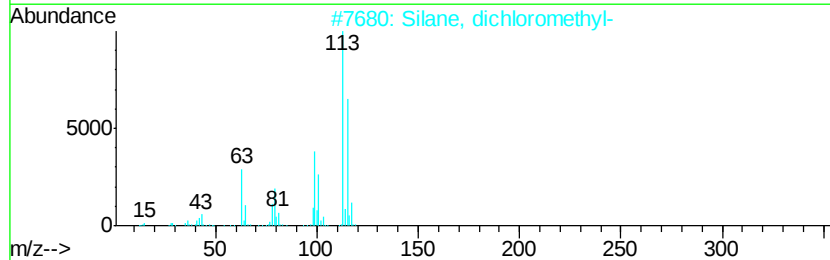
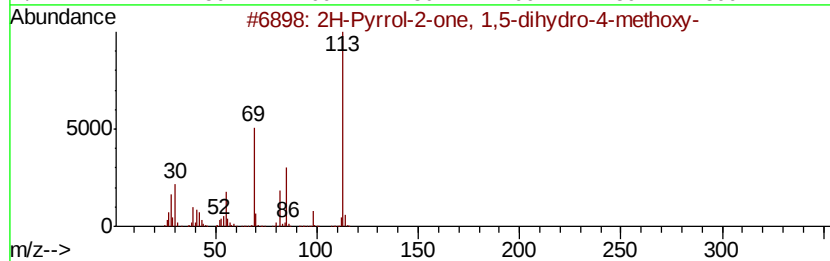
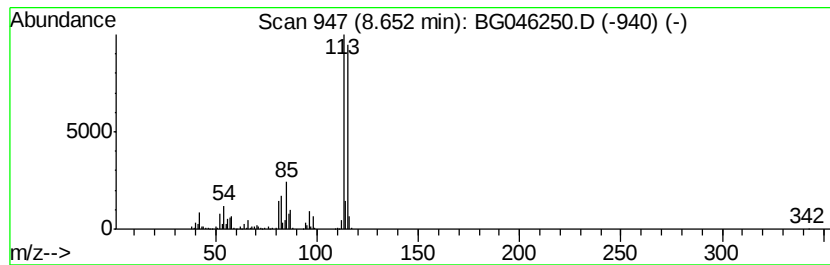
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Quant Title : SVOA CALIBRATION

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	33.47 ng/ul	812697	1,4-Dichlorobenzene-d4	7.95

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		Silane, dichloromethyl-	114	CH4Cl2Si	000075-54-7	38
3		Heptanoic acid, anhydride	242	C14H26O3	000626-27-7	32
4		Butanedioyl dihydrazide	146	C4H10N4O2	004146-43-4	25
5		Furazan-3-amine, 4-methoxy-	115	C3H5N3O2	078350-48-8	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\
Data File : BG046250.D
Acq On : 13 Aug 2020 15:39
Operator : CG/JU
Sample : PB130901BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK01

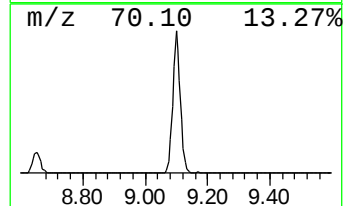
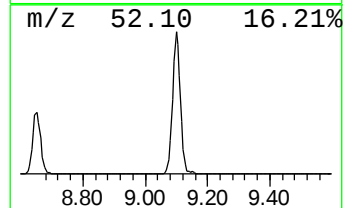
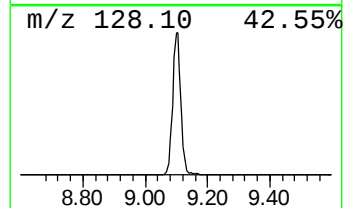
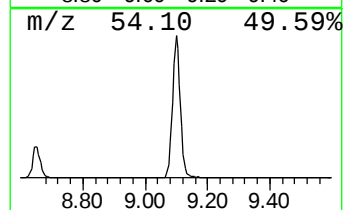
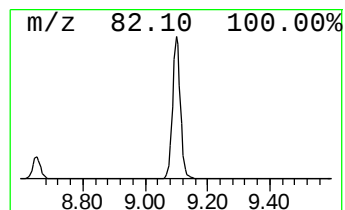
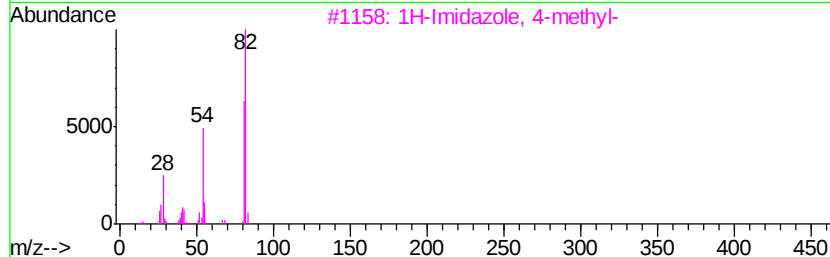
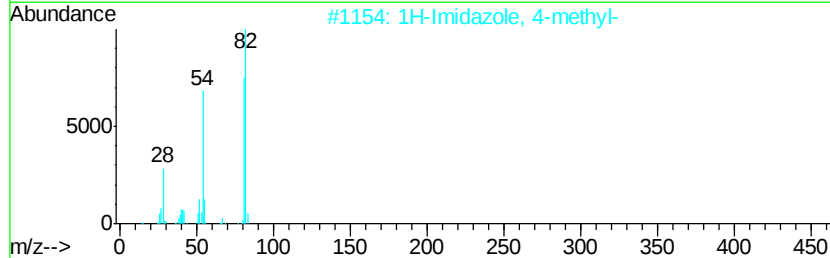
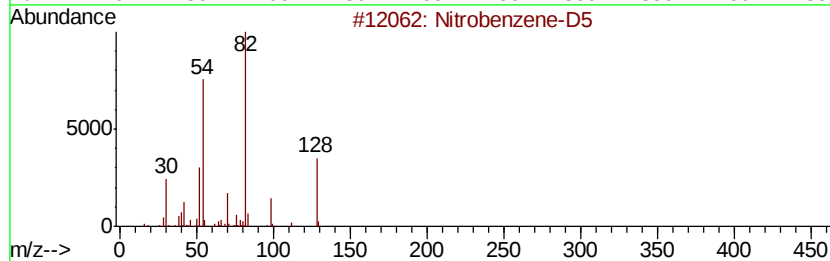
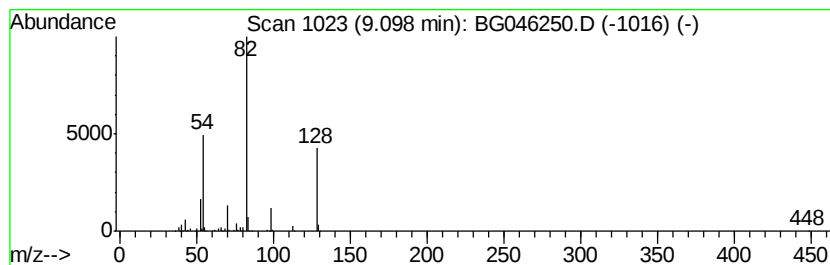
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1H-Imidazole, 4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.10	26.39 ng/ul	640753	1,4-Dichlorobenzene-d4	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nitrobenzene-D5	128	C6D5N02	004165-60-0	64
2		1H-Imidazole, 4-methyl-	82	C4H6N2	000822-36-6	50
3		1H-Imidazole, 4-methyl-	82	C4H6N2	000822-36-6	38
4		Nitrobenzene-D5	128	C6D5N02	004165-60-0	32
5		1H-Imidazole, 2-methyl-	82	C4H6N2	000693-98-1	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\
Data File : BG046250.D
Acq On : 13 Aug 2020 15:39
Operator : CG/JU
Sample : PB130901BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK01

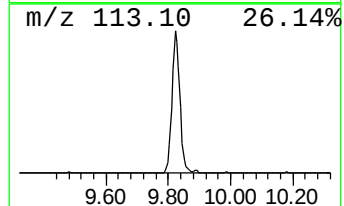
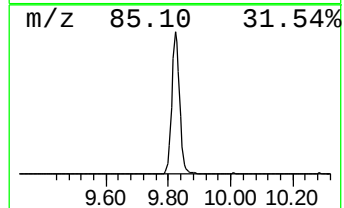
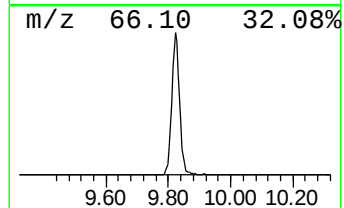
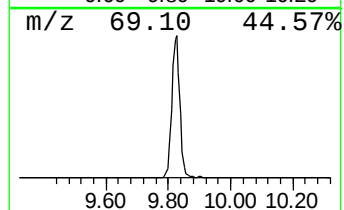
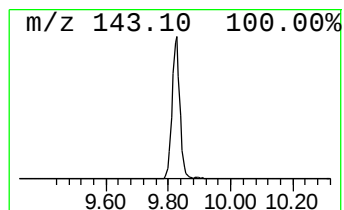
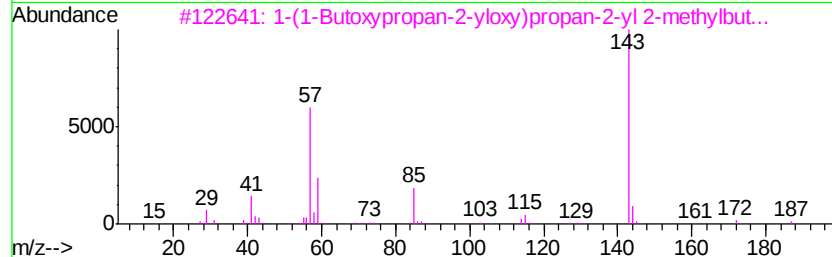
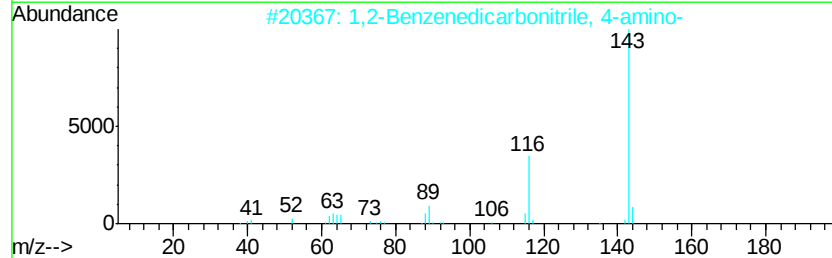
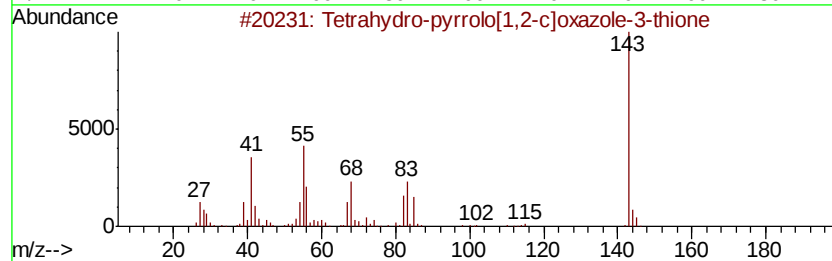
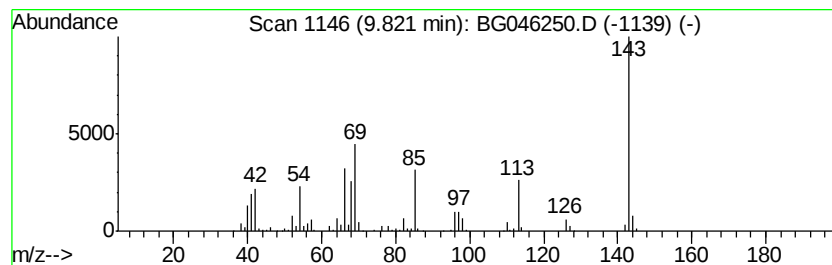
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Quant Title : SVOA CALIBRATION

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	15.32 ng/ul	543279	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		1-(1-Butoxypropan-2-yloxy)propan...	274	C15H30O4	1000367-10-9	12
4		4(1H)-Pyrimidinone, 6-amino-2,3-...	143	C4H5N3O5	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\BG081320\
Data File : BG046250.D
Acq On : 13 Aug 2020 15:39
Operator : CG/JU
Sample : PB130901BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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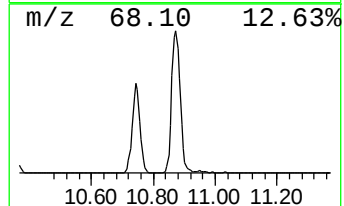
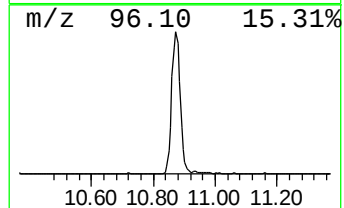
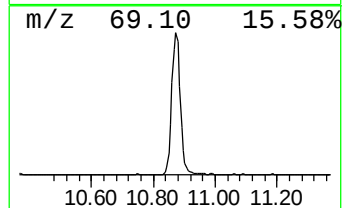
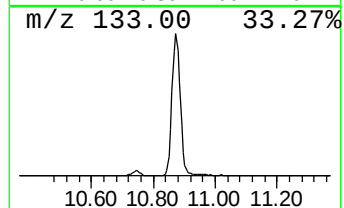
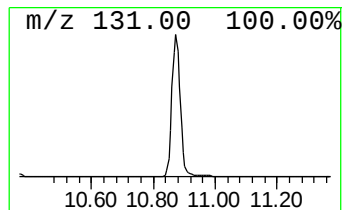
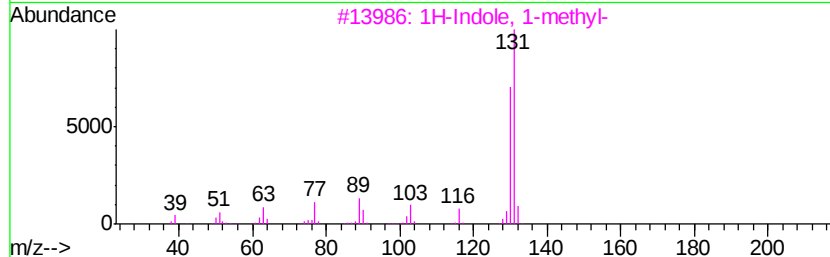
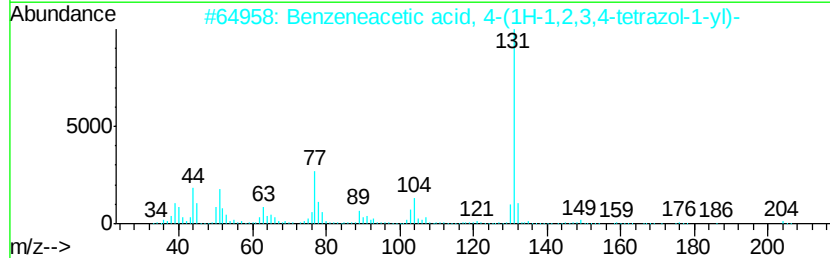
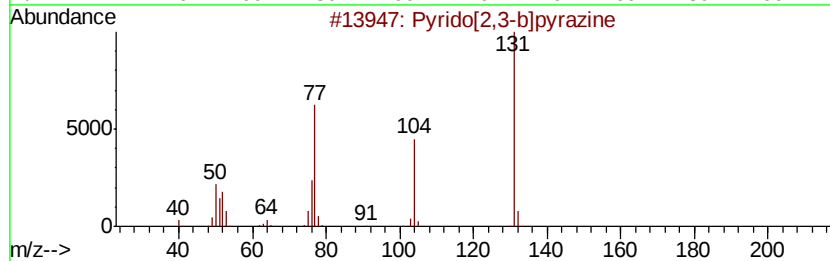
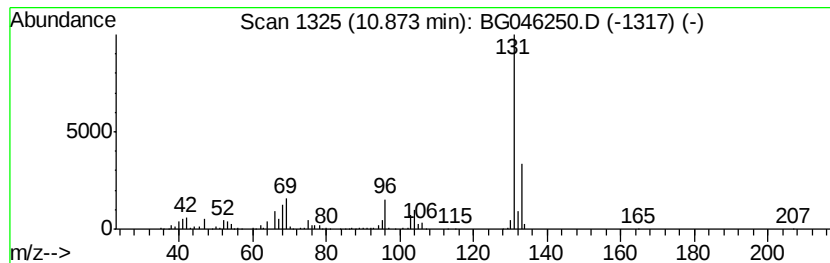
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-05 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.87	24.93 ng/ul	884201	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		1H-Indole, 1-methyl-	131	C9H9N	000603-76-9	28
4		Pyrido[2,3-b]pyrazine	131	C7H5N3	000322-46-3	9
5		1H-Indole, 6-methyl-	131	C9H9N	003420-02-8	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\
Data File : BG046250.D
Acq On : 13 Aug 2020 15:39
Operator : CG/JU
Sample : PB130901BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

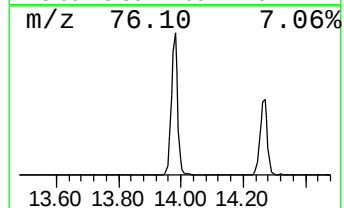
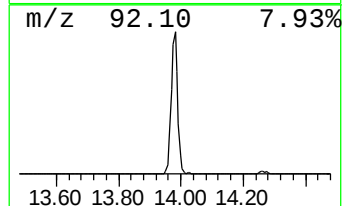
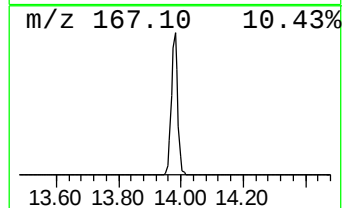
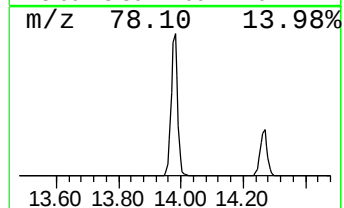
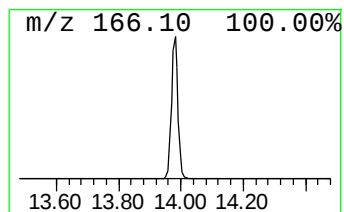
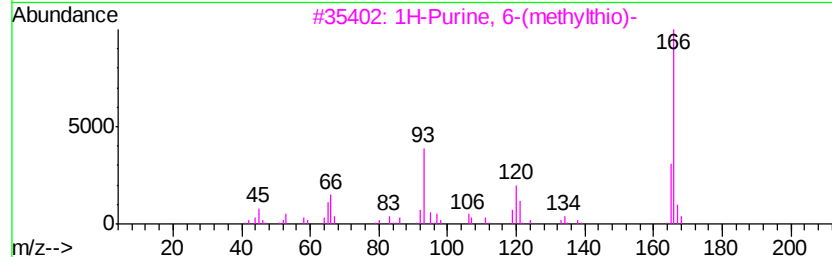
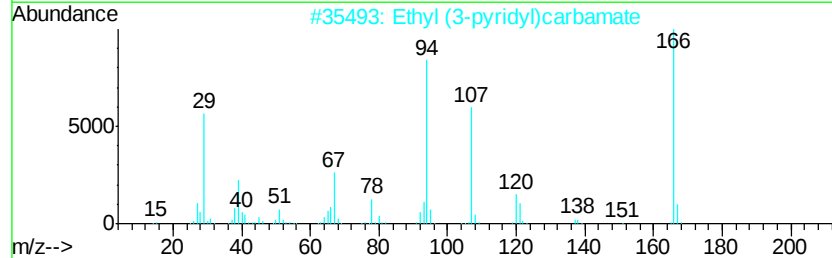
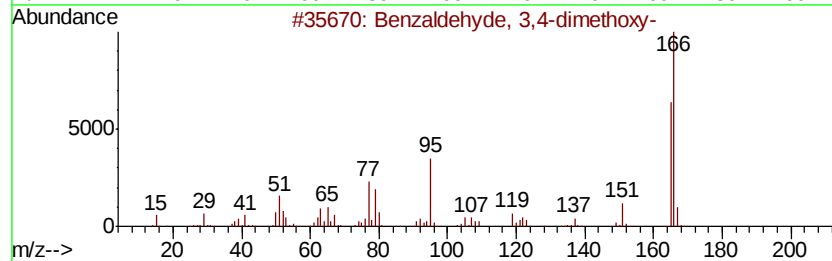
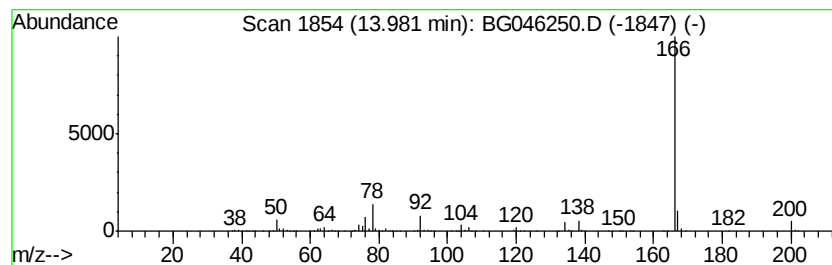
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BNA_G
ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-06 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.98	27.42 ng/ul	1448640	Acenaphthene-d10	14.57
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzaldehyde, 3,4-dimethoxy-	166 C9H10O3	000120-14-9 45
2		Ethyl (3-pyridyl)carbamate	166 C8H10N2O2	006276-11-5 42
3		1H-Purine, 6-(methylthio)-	166 C6H6N4S	000050-66-8 38
4		Phenol, 5-(dimethylamino)-2-nitr...	166 C8H10N2O2	016761-04-9 38
5		Benzenamine, N,N-dimethyl-4-nitro-	166 C8H10N2O2	000100-23-2 38



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\
Data File : BG046250.D
Acq On : 13 Aug 2020 15:39
Operator : CG/JU
Sample : PB130901BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK01

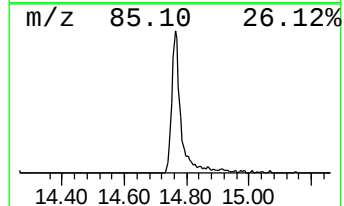
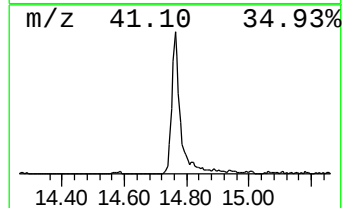
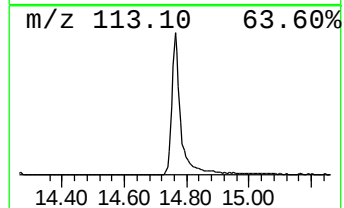
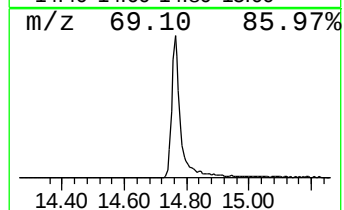
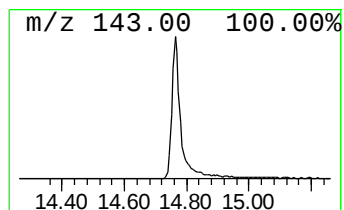
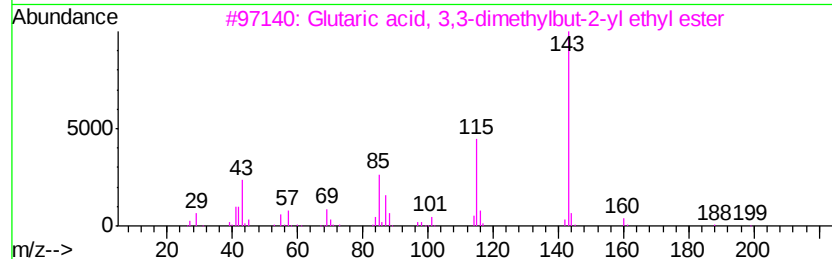
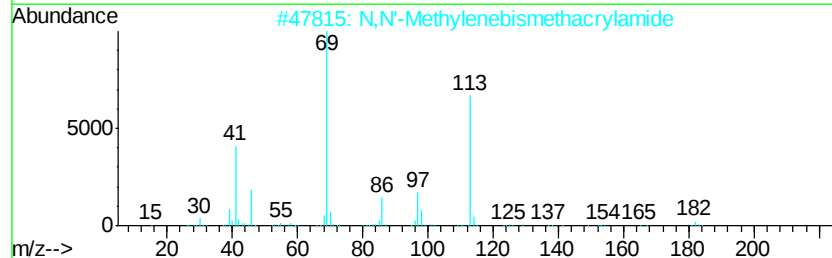
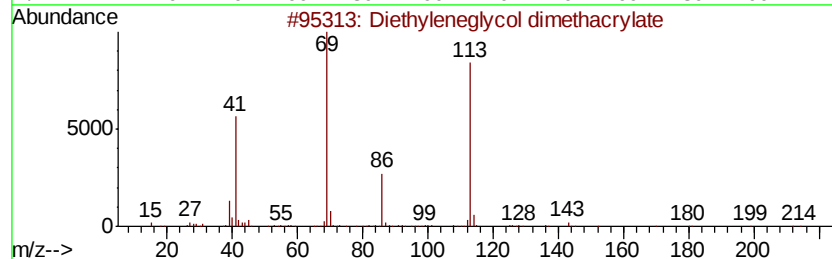
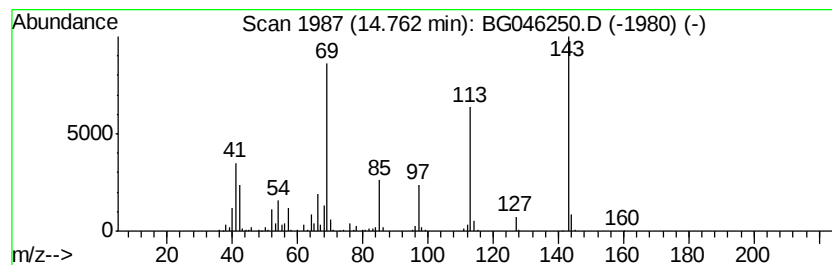
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-07 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.76	11.39 ng/ul	601869	Acenaphthene-d10	14.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diethyleneglycol dimethacrylate	242	C12H18O5	002358-84-1	12
2		N,N'-Methylenebismethacrylamide	182	C9H14N2O2	002359-15-1	12
3		Glutaric acid, 3,3-dimethylbut-2...	244	C13H24O4	1000359-74-7	12
4		1,2-Benzenedicarbonitrile, 4-amino-	143	C8H5N3	056765-79-8	10
5		Cyclopentanecarboxylic acid, eth...	140	C8H12O2	016523-06-1	10



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\
Data File : BG046250.D
Acq On : 13 Aug 2020 15:39
Operator : CG/JU
Sample : PB130901BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK01

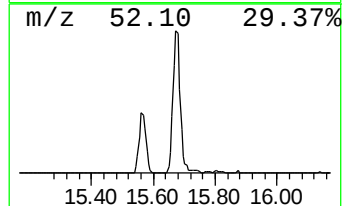
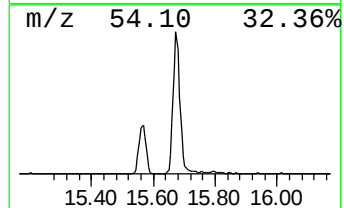
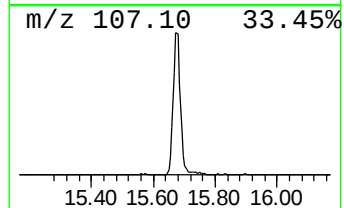
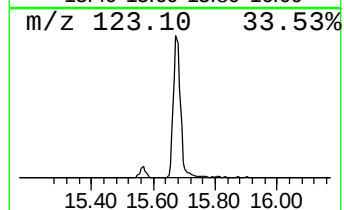
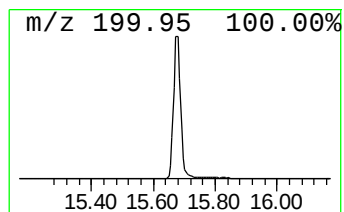
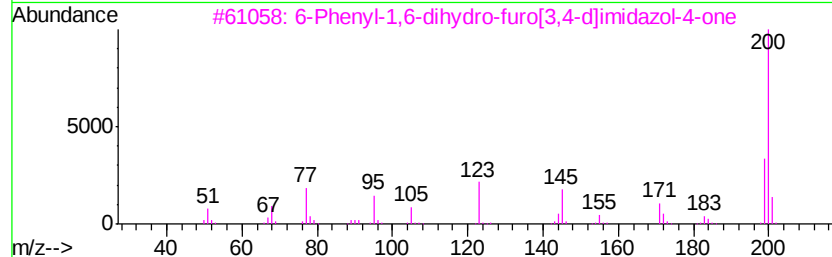
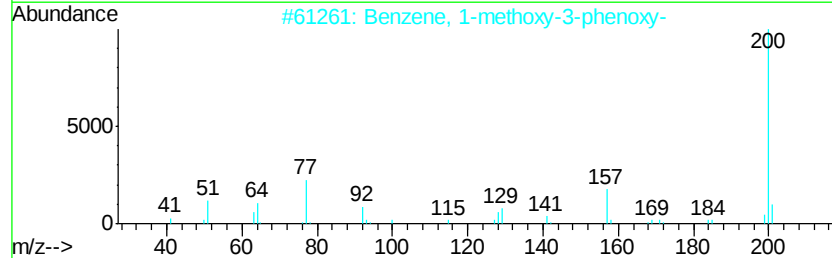
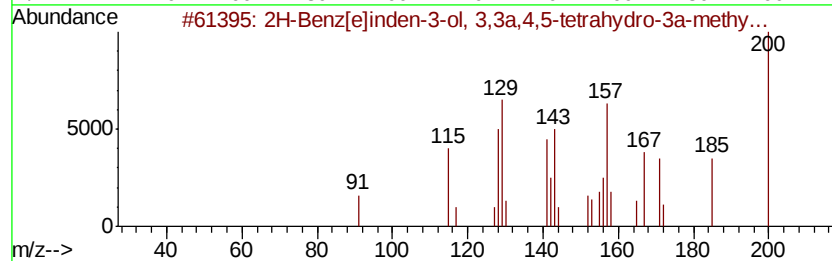
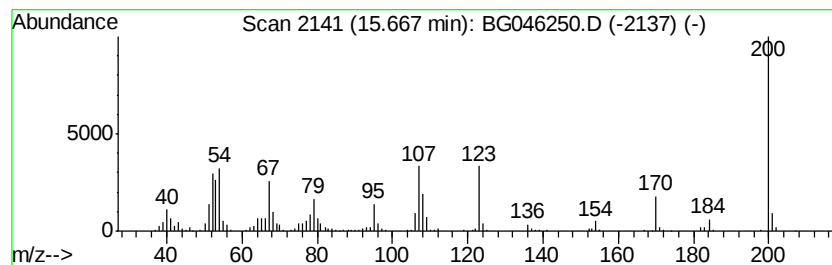
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-08 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.67	11.55 ng/ul	610328	Acenaphthene-d10	14.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2H-Benz[e]inden-3-ol, 3,3a,4,5-t...	200	C14H16O	071805-91-9	35
2		Benzene, 1-methoxy-3-phenoxy-	200	C13H12O2	001655-68-1	22
3		6-Phenyl-1,6-dihydro-furo[3,4-d]...	200	C11H8N2O2	313263-12-6	17
4		N-[2-Methyl-4-pyridyl]-4-aminoph...	200	C12H12N2O	1000252-23-9	16
5		1,6-Dioxaspiro[4.4]non-3-ene, 2-...	200	C13H12O2	016863-61-9	16



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG081320\
Data File : BG046250.D
Acq On : 13 Aug 2020 15:39
Operator : CG/JU
Sample : PB130901BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK01

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...	3.16	3.2	ng/ul	77085	1	7.95	485664	20.0
2-Furancarboxylic...	7.11	26.8	ng/ul	651168	1	7.95	485664	20.0
unknown-01	7.28	21.8	ng/ul	529864	1	7.95	485664	20.0
unknown-02	7.48	27.0	ng/ul	655538	1	7.95	485664	20.0
unknown-03	8.65	33.5	ng/ul	812697	1	7.95	485664	20.0
1H-Imidazole, 4-m...	9.10	26.4	ng/ul	640753	1	7.95	485664	20.0
unknown-04	9.82	15.3	ng/ul	543279	2	10.74	709427	20.0
unknown-05	10.87	24.9	ng/ul	884201	2	10.74	709427	20.0
unknown-06	13.98	27.4	ng/ul	1448640	3	14.57	1056760	20.0
unknown-07	14.76	11.4	ng/ul	601869	3	14.57	1056760	20.0
unknown-08	15.67	11.6	ng/ul	610328	3	14.57	1056760	20.0