

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
 Data File : BG046342.D
 Acq On : 19 Aug 2020 11:21
 Operator : CG/JU
 Sample : PB130973BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK73

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	19	rVB	34690	53933	2.38%	0.253%
2	3.395	49	52	61	rVB	17337	27893	1.23%	0.131%
3	4.141	176	179	186	rVB4	4493	7689	0.34%	0.036%
4	5.069	335	337	343	rVB4	1473	2277	0.10%	0.011%
5	7.108	677	684	695	rBV	261078	464870	20.51%	2.177%
6	7.266	704	711	721	rVV	210217	354710	15.65%	1.661%
7	7.472	739	746	755	rBV	283787	481184	21.23%	2.253%
8	7.936	819	825	835	rVB	239730	411512	18.16%	1.927%
9	8.641	938	945	958	rBV	338347	594705	26.24%	2.784%
10	9.088	1014	1021	1032	rBV	248962	448970	19.81%	2.102%
11	9.816	1137	1145	1153	rBV	212298	381165	16.82%	1.785%
12	10.357	1230	1237	1257	rBV	350352	632440	27.90%	2.961%
13	10.733	1294	1301	1310	rVV	355089	623184	27.50%	2.918%
14	10.862	1316	1323	1338	rBV	319971	603574	26.63%	2.826%
15	12.325	1566	1572	1578	rBV3	4315	8107	0.36%	0.038%
16	13.970	1845	1852	1865	rBV	673694	970969	42.84%	4.546%
17	14.258	1894	1901	1913	rBV	784566	1229678	54.26%	5.757%
18	14.570	1946	1954	1966	rVB	567422	933246	41.18%	4.369%
19	14.764	1978	1987	2004	rBV	216645	424063	18.71%	1.985%
20	15.557	2115	2122	2131	rBV	1052954	1593680	70.32%	7.462%
21	15.674	2135	2142	2154	rVV	239237	375908	16.59%	1.760%
22	15.798	2158	2163	2168	rVB2	11830	17662	0.78%	0.083%
23	16.344	2252	2256	2259	rVB4	3820	5328	0.24%	0.025%
24	17.096	2379	2384	2388	rVB4	3116	4379	0.19%	0.021%
25	17.308	2413	2420	2429	rBV	810089	1204872	53.16%	5.641%
26	17.407	2430	2437	2445	rVV	1189948	1779740	78.52%	8.333%
27	19.123	2726	2729	2732	rBV4	2309	2466	0.11%	0.012%
28	19.258	2749	2752	2756	rBV3	1710	2614	0.12%	0.012%
29	19.329	2759	2764	2769	rVB2	15321	20942	0.92%	0.098%
30	19.576	2802	2806	2812	rVB	12431	18189	0.80%	0.085%
31	19.628	2812	2815	2819	rBV6	3555	4768	0.21%	0.022%
32	19.693	2819	2826	2839	rBV	1500277	2209488	97.49%	10.345%
33	19.946	2865	2869	2871	rBV5	2290	3537	0.16%	0.017%
34	20.128	2898	2900	2902	rBV3	3024	2416	0.11%	0.011%

LSC Area Percent Report

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35	20.257	2921	2922	2925	rBV3	3012	3743	0.17%	0.018%
36	20.827	3017	3019	3020	rBV2	4826	2775	0.12%	0.013%
37	21.556	3136	3143	3151	rBV	872467	1417638	62.55%	6.637%
38	23.019	3390	3392	3400	rVB4	17975	29023	1.28%	0.136%
39	24.446	3624	3635	3648	rBV	859607	2266476	100.00%	10.612%
40	24.658	3660	3671	3692	rVB2	511773	1552249	68.49%	7.268%
41	25.809	3860	3867	3875	rVB2	34802	104683	4.62%	0.490%
42	27.108	4081	4088	4097	rVB4	28580	81829	3.61%	0.383%

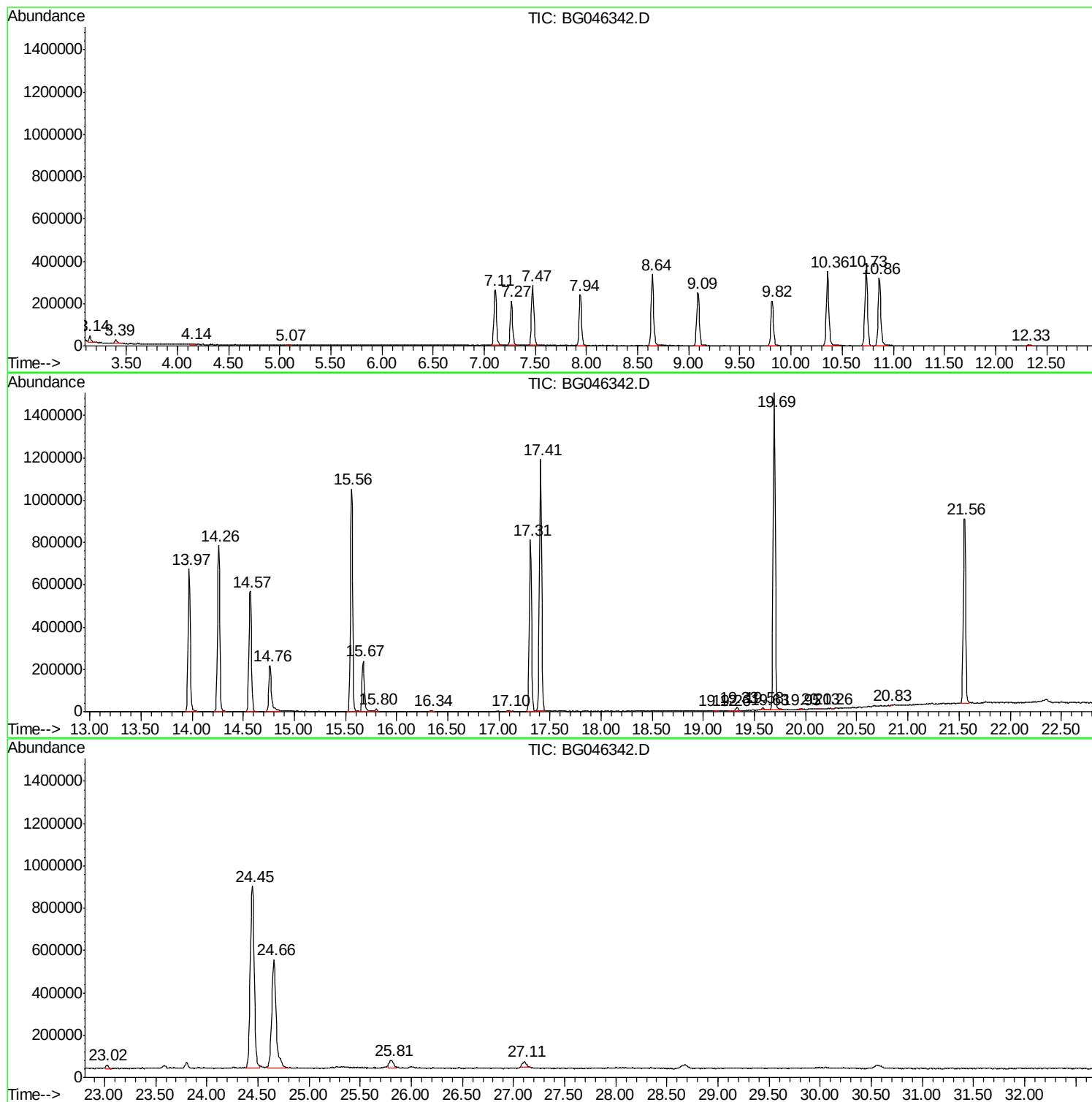
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG081320PAH.M
Quant Title : SVOA CALIBRATION

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



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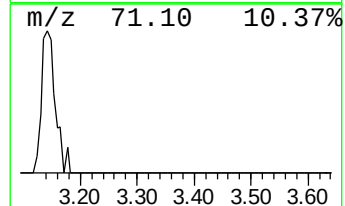
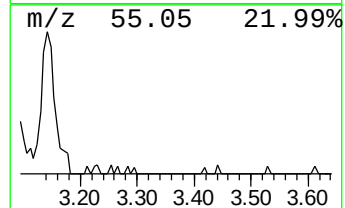
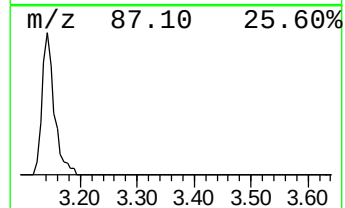
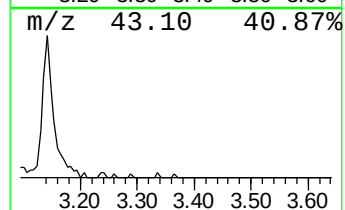
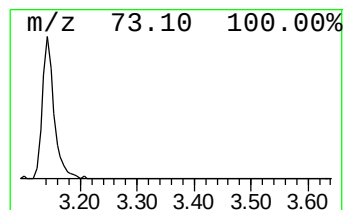
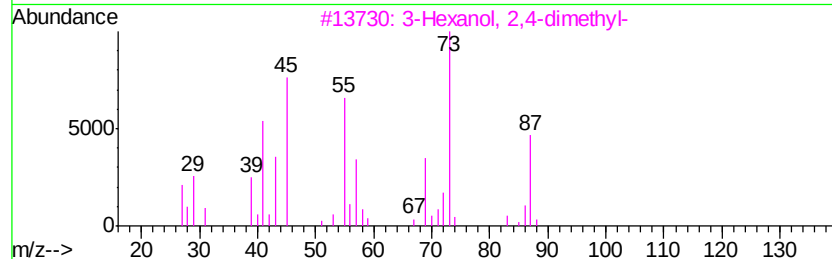
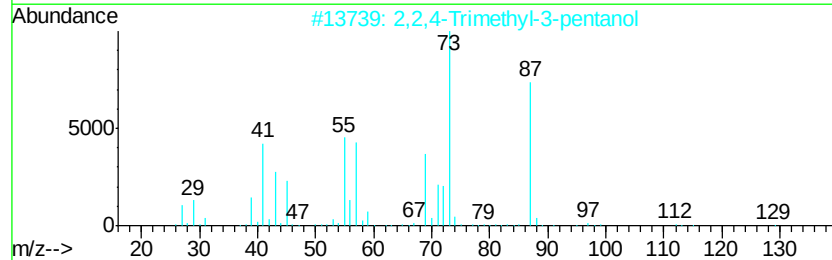
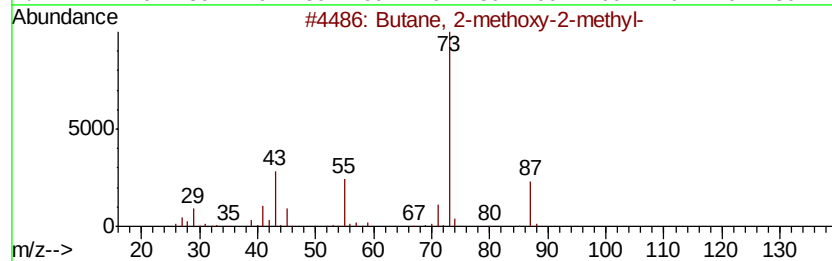
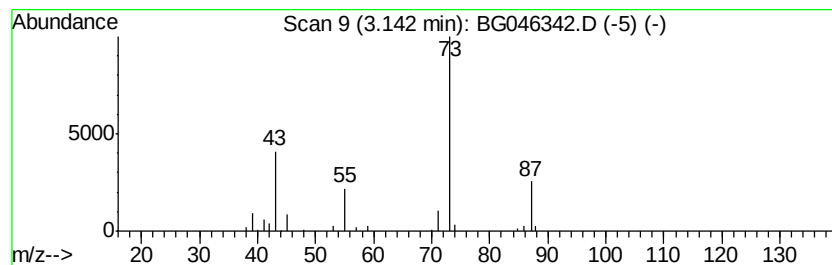
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Quant Title : SVOA CALIBRATION

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.
3.14	2.62 ng/ul	53933	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	64
3		3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	42
4		3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	25
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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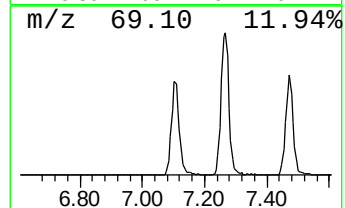
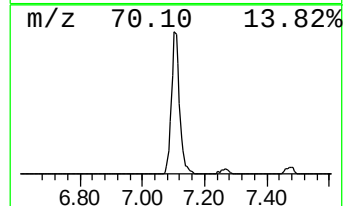
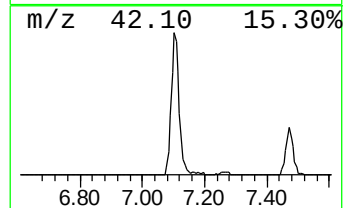
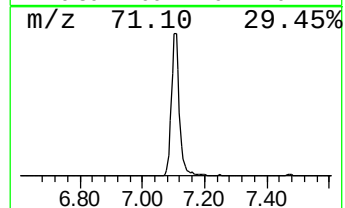
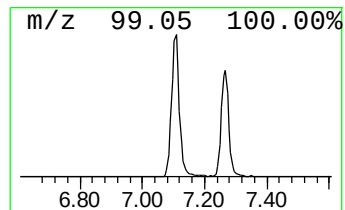
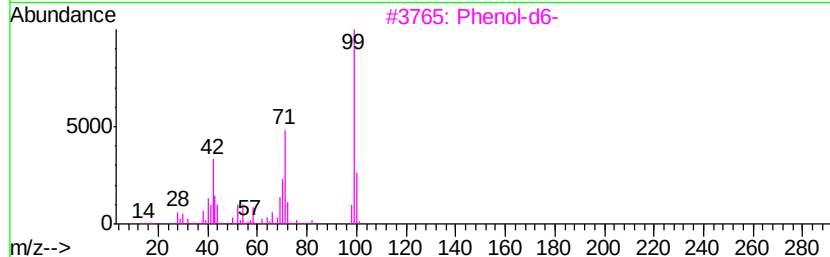
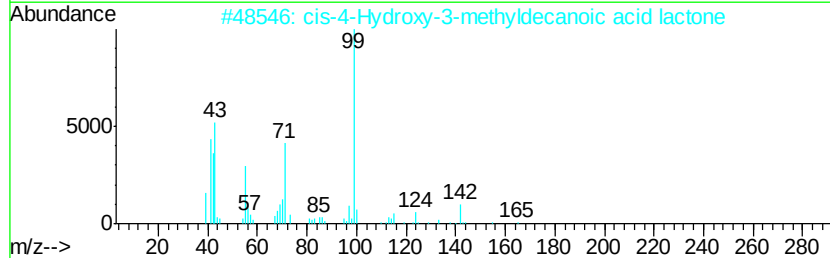
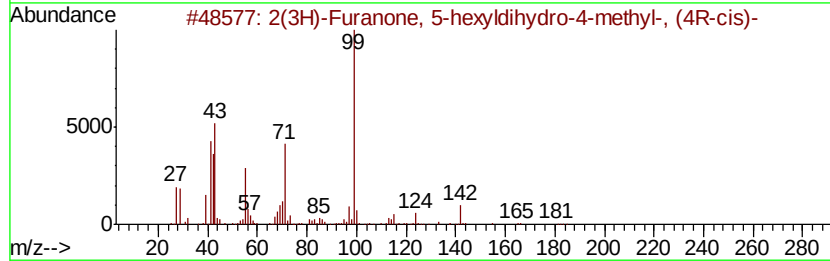
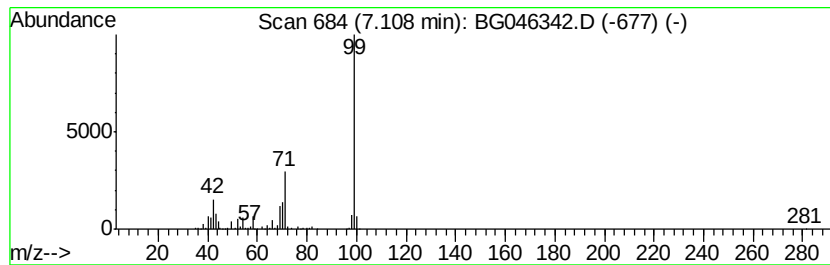
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 2(3H)-Furanone, 5-hexyldihy... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.11	22.59 ng/ul	464870	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2(3H)-Furanone, 5-hexyldihydro-4...	184	C11H20O2	121644-12-0	64
2		cis-4-Hydroxy-3-methyldecanoic a...	184	C11H20O2	147254-32-8	64
3		Phenol-d6-	100	C6D6O	013127-88-3	56
4		Hexanoic acid, 4-tridecyl ester	298	C19H38O2	1000279-29-3	56
5		Hexanoic acid, 3-tridecyl ester	298	C19H38O2	1000279-29-2	56



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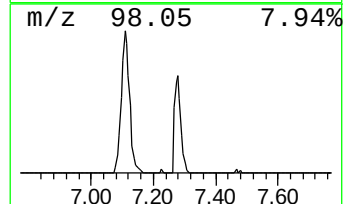
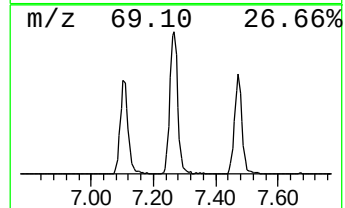
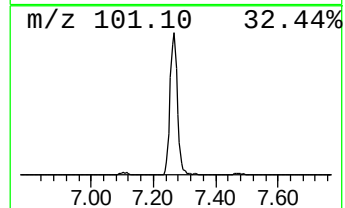
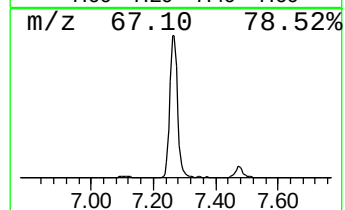
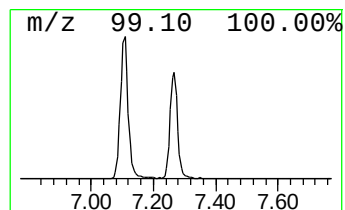
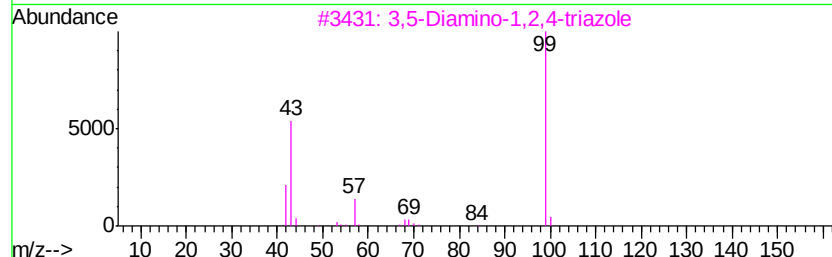
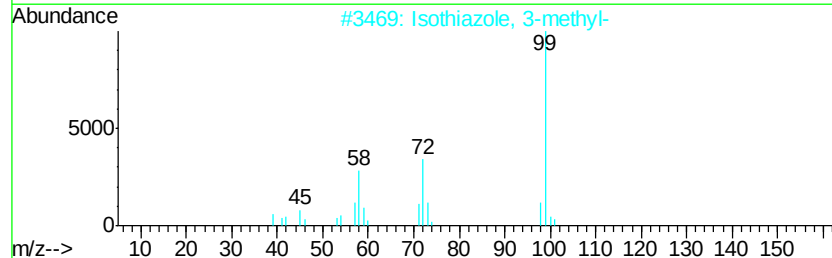
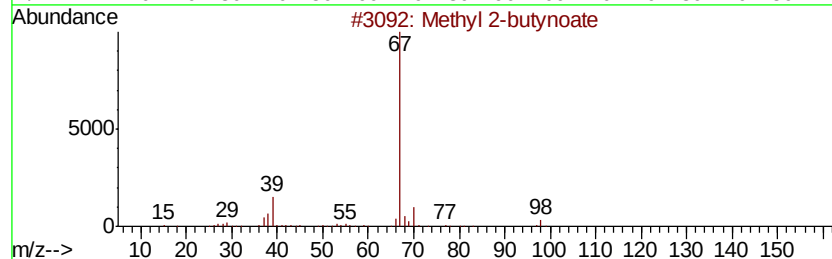
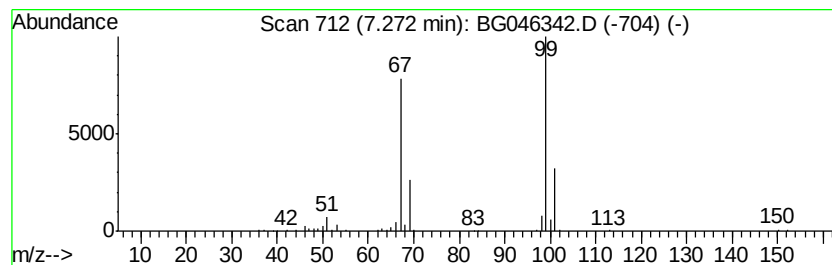
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Quant Title : SVOA CALIBRATION

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	17.24 ng/ul	354710	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyl 2-butyrate	98	C5H6O2	023326-27-4	9
2		Isothiazole, 3-methyl-	99	C4H5NS	000693-92-5	9
3		3,5-Diamino-1,2,4-triazole	99	C2H5N5	001455-77-2	9
4		4-Methylthiazole	99	C4H5NS	000693-95-8	9
5		1-Butyne, 3,3-dimethyl-	82	C6H10	000917-92-0	9



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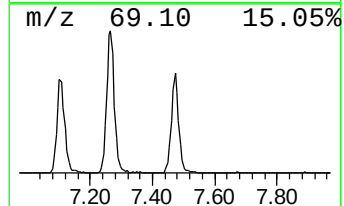
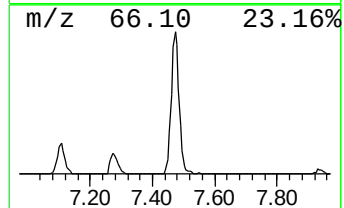
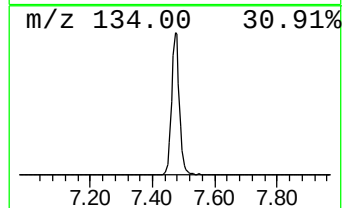
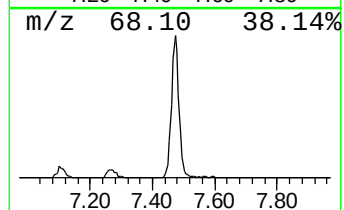
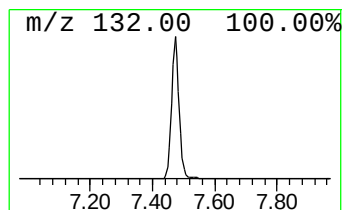
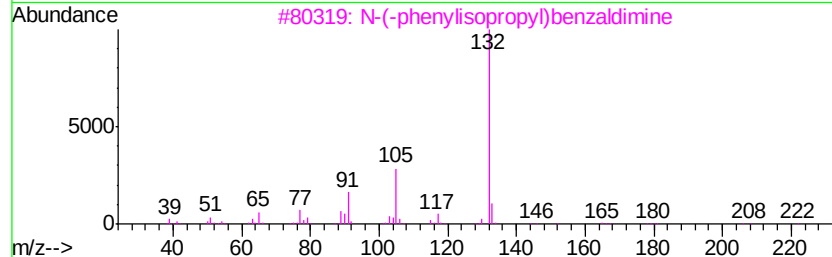
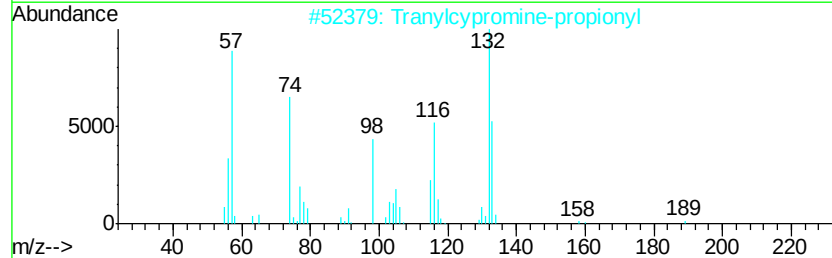
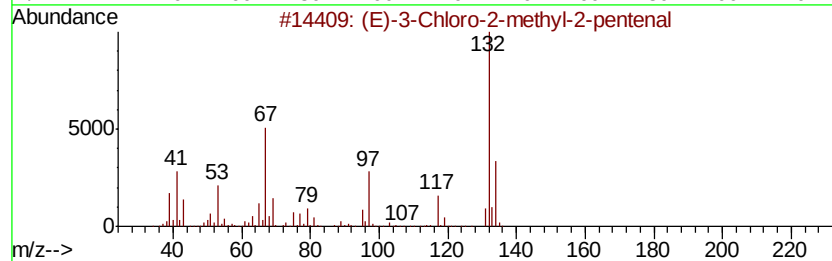
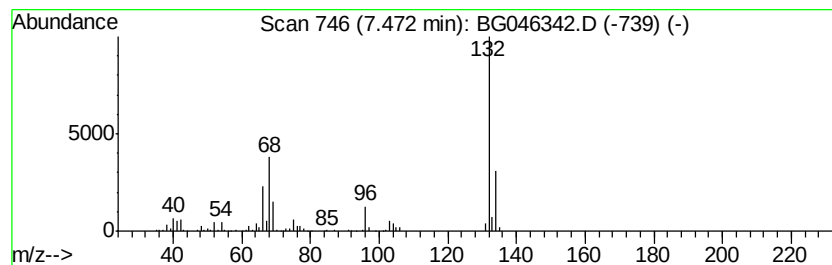
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Peak Number 4 unknown-02 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.47	23.39 ng/ul	481184	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		N-(phenylisopropyl)benzalimine	223	C16H17N	1000379-01-3	10
4		5-Aminoindole	132	C8H8N2	005192-03-0	10
5		Xenon	132	Xe	007440-63-3	9



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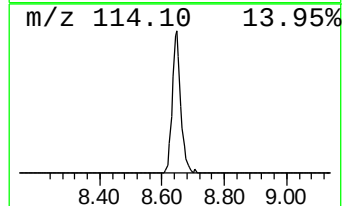
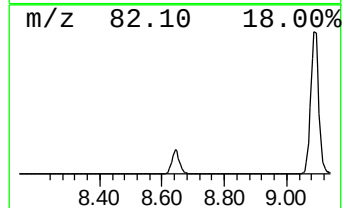
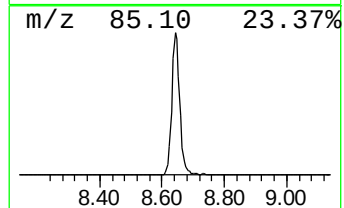
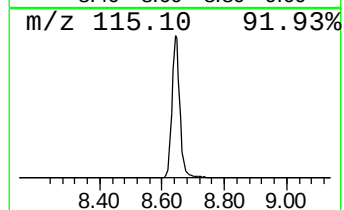
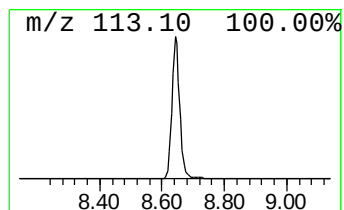
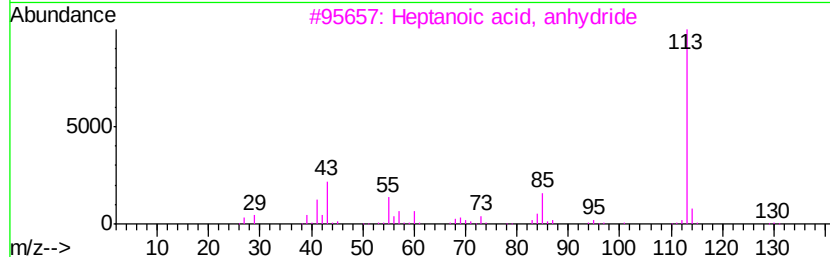
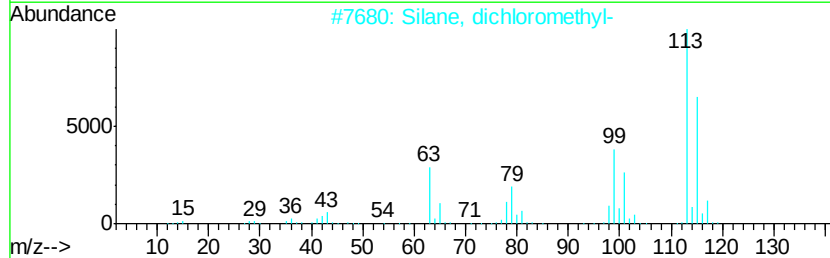
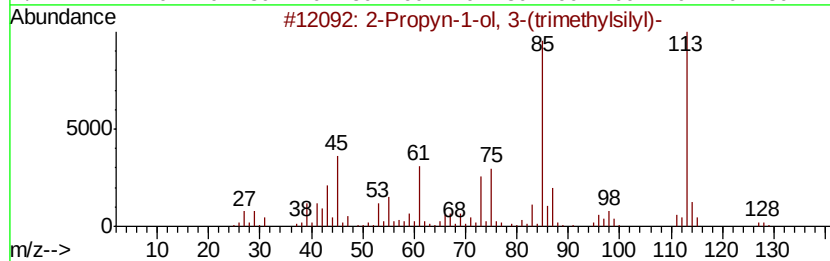
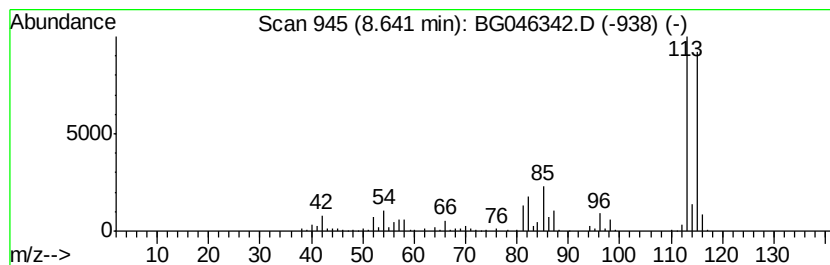
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Quant Title : SVOA CALIBRATION

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.
8.64	28.90 ng/ul	594705	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propyn-1-ol, 3-(trimethylsilyl)-	128	C6H12OSi	005272-36-6	43
2		Silane, dichloromethyl-	114	CH4Cl2Si	000075-54-7	38
3		Heptanoic acid, anhydride	242	C14H26O3	000626-27-7	32
4		2H-Pyrrol-2-one, 1,5-dihydro-4-m...	113	C5H7NO2	069778-83-2	32
5		Piperidine-2,5-dione	113	C5H7NO2	052065-78-8	17



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046342.D
Acq On : 19 Aug 2020 11:21
Operator : CG/JU
Sample : PB130973BL
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK73

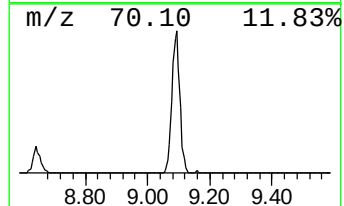
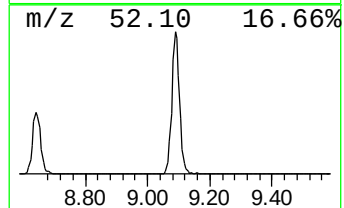
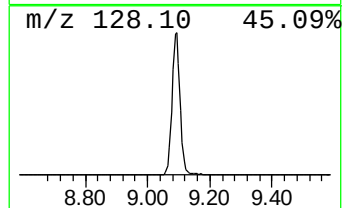
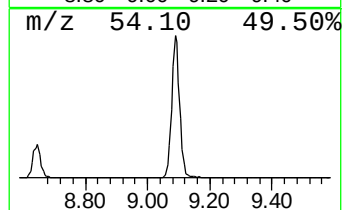
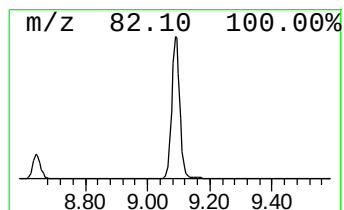
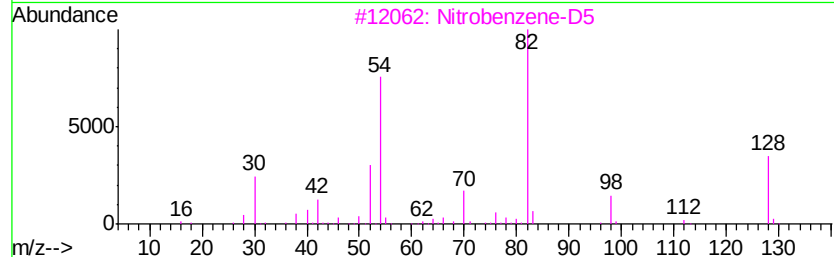
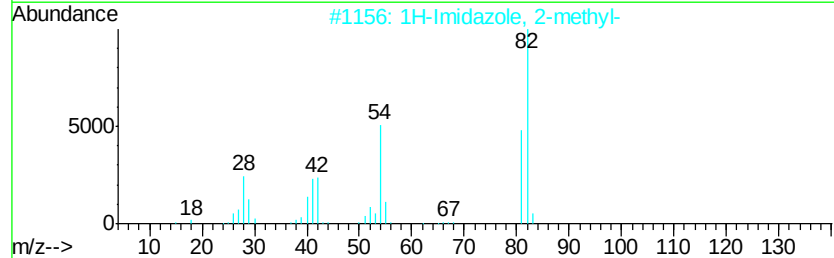
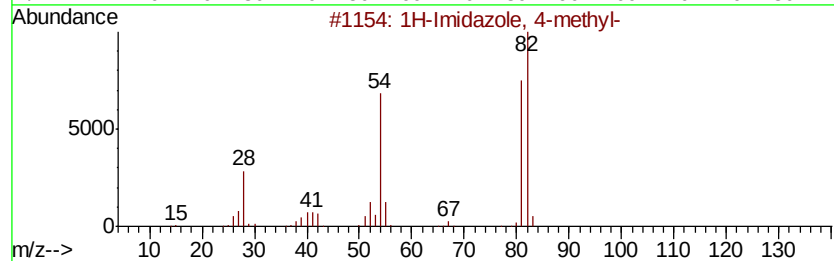
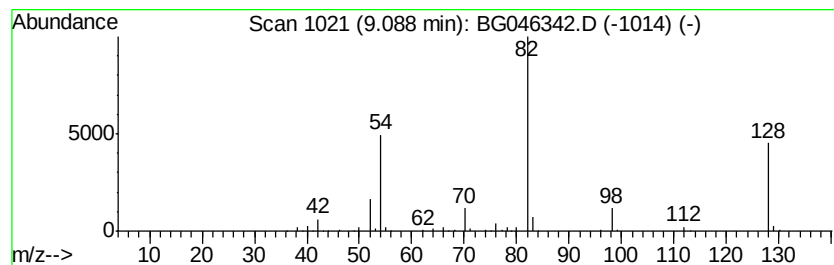
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Quant Title : SVOA CALIBRATION

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	21.82 ng/ul	448970	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Imidazole, 4-methyl-	82	C4H6N2	000822-36-6	50
2		1H-Imidazole, 2-methyl-	82	C4H6N2	000693-98-1	40
3		Nitrobenzene-D5	128	C6D5NO2	004165-60-0	38
4		1H-Imidazole, 2-methyl-	82	C4H6N2	000693-98-1	38
5		Nitrobenzene-D5	128	C6D5NO2	004165-60-0	16



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046342.D
Acq On : 19 Aug 2020 11:21
Operator : CG/JU
Sample : PB130973BL
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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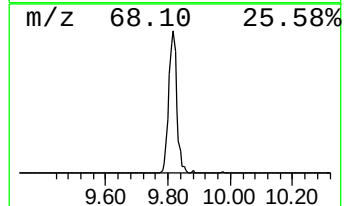
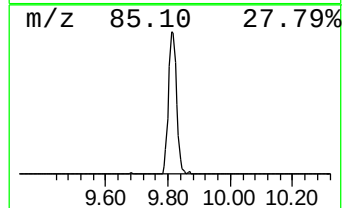
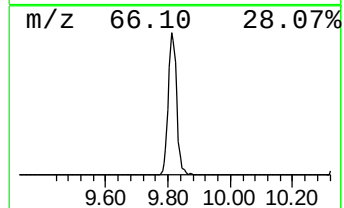
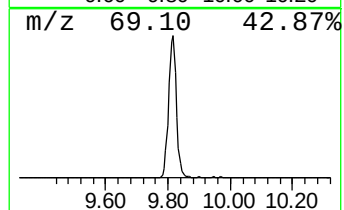
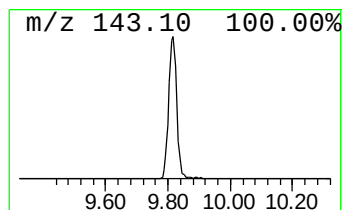
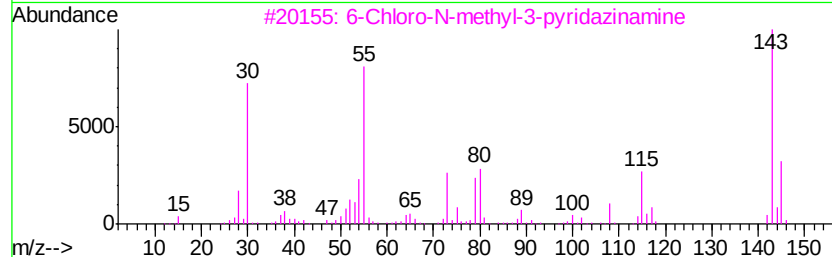
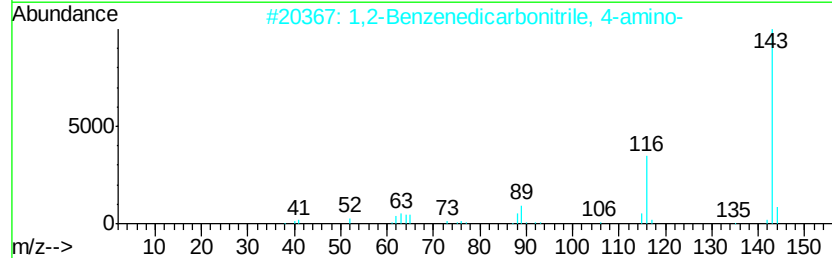
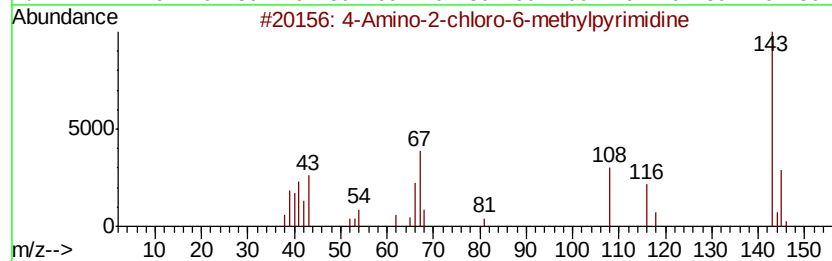
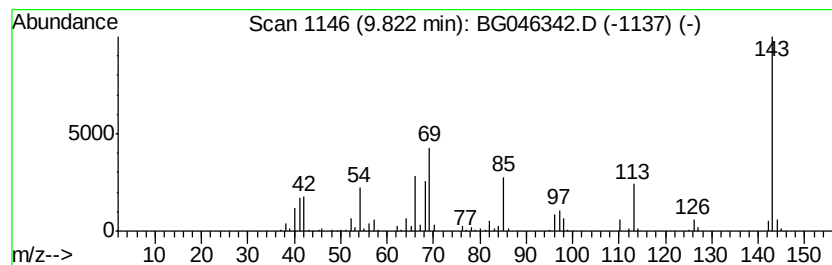
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-04 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.82	12.23 ng/ul	381165	Napthalene-d8	10.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Amino-2-chloro-6-methylpyrimidine	143	C5H6ClN3	014394-60-6	33
2		1,2-Benzenedicarbonitrile, 4-amino-	143	C8H5N3	056765-79-8	18
3		6-Chloro-N-methyl-3-pyridazinamine	143	C5H6ClN3	014959-32-1	12
4		1-(1-Butoxypropan-2-yloxy)propan...	274	C15H30O4	1000367-12-9	10
5		1,2-Benzenedicarbonitrile, 4-amino-	143	C8H5N3	056765-79-8	10



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046342.D
Acq On : 19 Aug 2020 11:21
Operator : CG/JU
Sample : PB130973BL
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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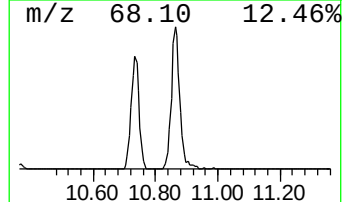
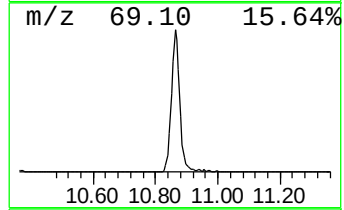
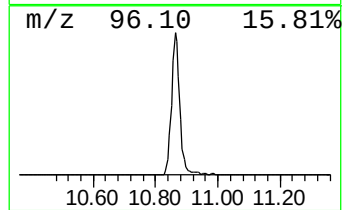
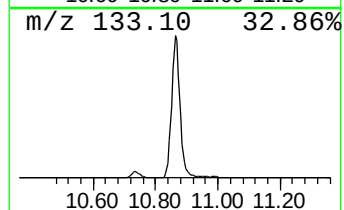
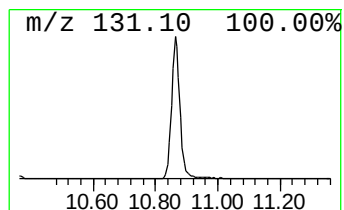
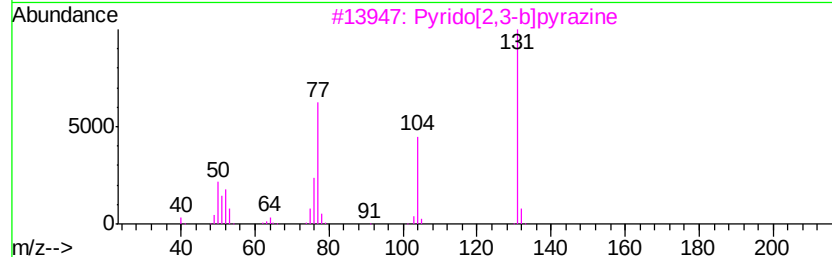
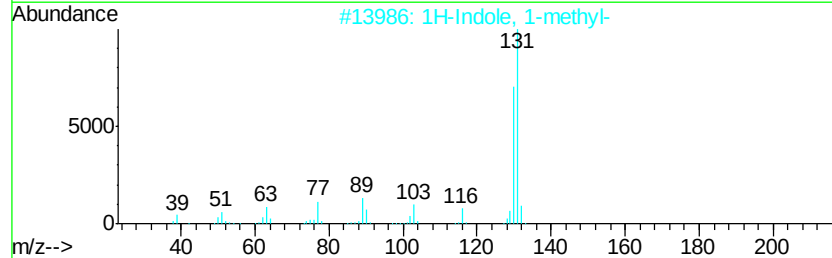
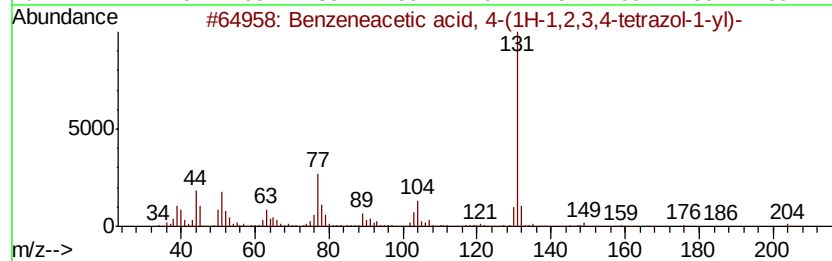
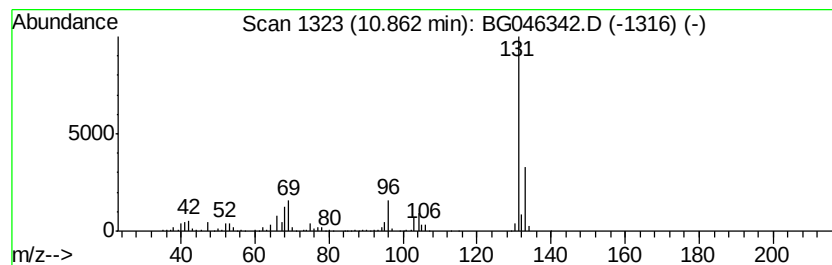
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Quant Title : SVOA CALIBRATION

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.
10.86	19.37 ng/ul	603574	Naphthalene-d8	10.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzeneacetic acid, 4-(1H-1,2,3,...	204	C9H8N4O2	1000351-56-5	33
2		1H-Indole, 1-methyl-	131	C9H9N	000603-76-9	28
3		Pyrido[2,3-b]pyrazine	131	C7H5N3	000322-46-3	25
4		Pyrido[2,3-b]pyrazine	131	C7H5N3	000322-46-3	9
5		4-Pyrimidinamine, 2,6-difluoro-	131	C4H3F2N3	000675-12-7	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046342.D
Acq On : 19 Aug 2020 11:21
Operator : CG/JU
Sample : PB130973BL
Misc :
ALS Vial : 5 Sample Multiplier: 1

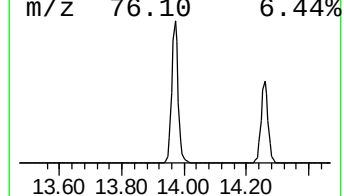
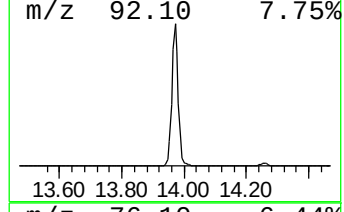
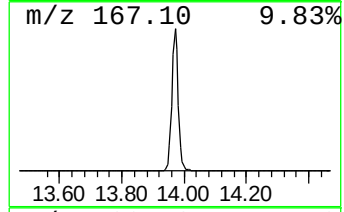
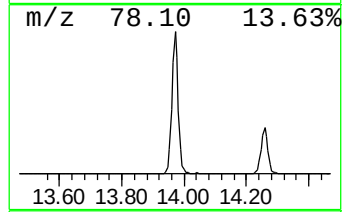
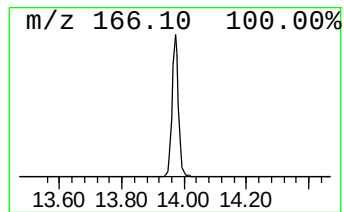
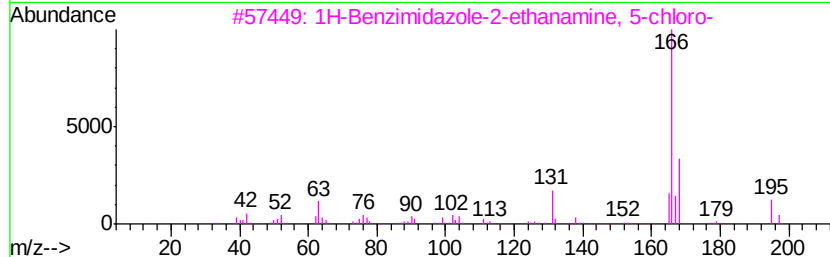
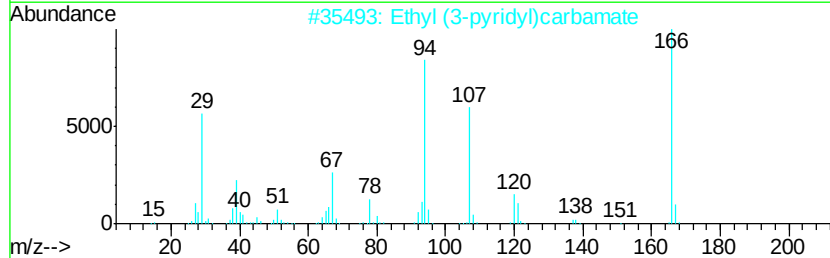
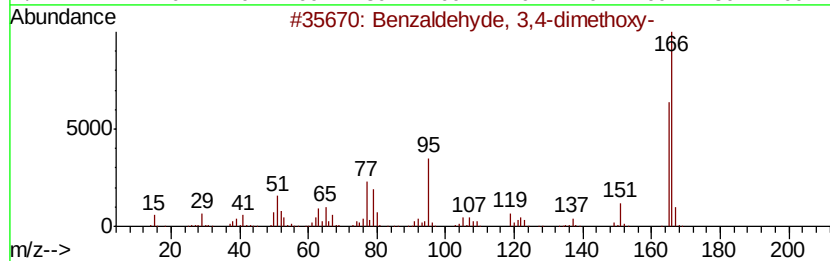
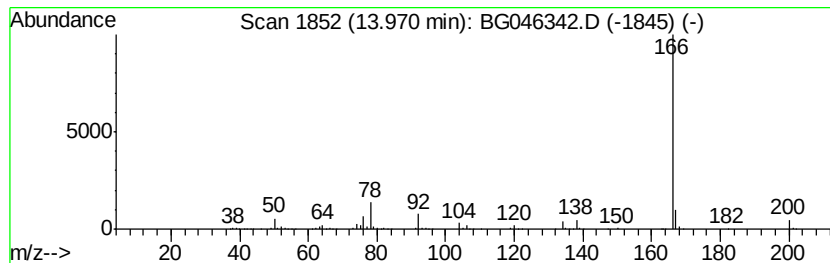
Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG081320PAH.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-06 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.97	20.81 ng/ul	970969	Acenaphthene-d10	14.56	
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	39
3	1H-Benzimidazole-2-ethanamine, 5...	195	C9H10ClN3	135875-16-0	38
4	1H-Purine, 6-(methylthio)-	166	C6H6N4S	000050-66-8	38
5	Benzenamine, N,N-dimethyl-4-nitro-	166	C8H10N2O2	000100-23-2	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046342.D
Acq On : 19 Aug 2020 11:21
Operator : CG/JU
Sample : PB130973BL
Misc :
ALS Vial : 5 Sample Multiplier: 1

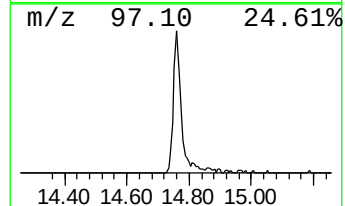
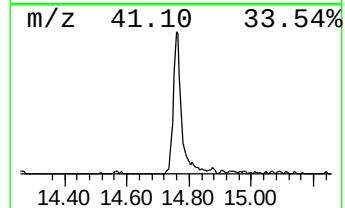
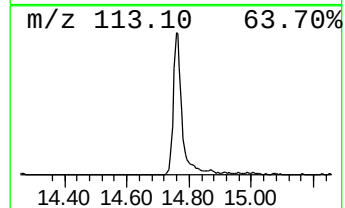
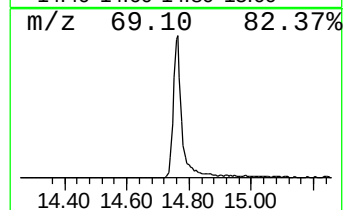
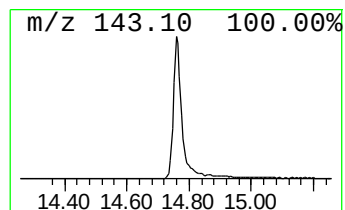
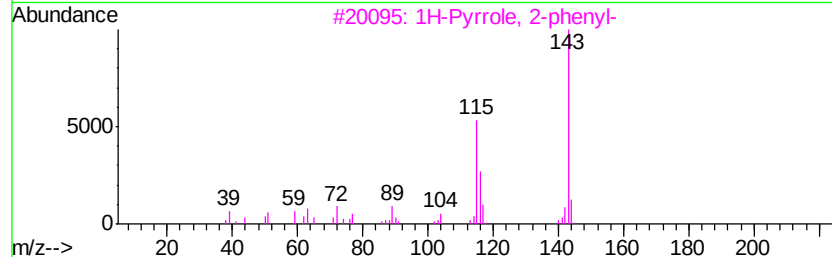
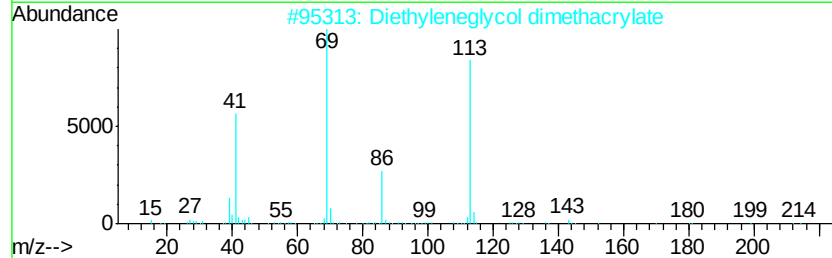
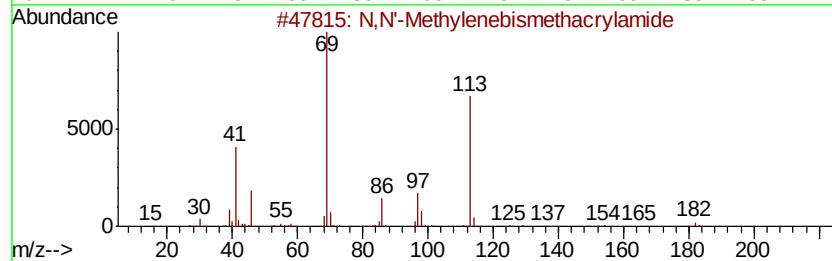
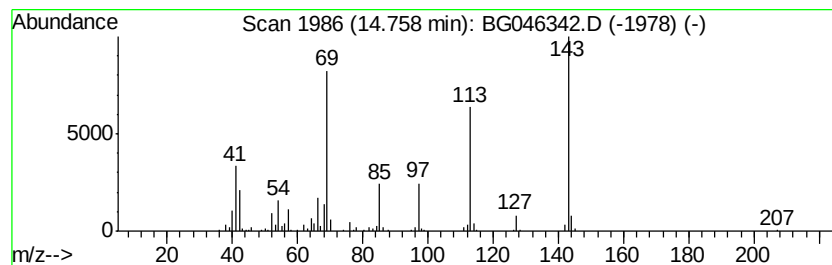
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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	9.09 ng/ul	424063	Acenaphthene-d10	14.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		N,N'-Methylenebismethacrylamide	182 C9H14N2O2	002359-15-1 38
2		Diethyleneglycol dimethacrylate	242 C12H18O5	002358-84-1 35
3		1H-Pyrrole, 2-phenyl-	143 C10H9N	003042-22-6 22
4		Glutaric acid, ethyl 4-methylhep...	272 C15H28O4	1000377-14-9 16
5		6-Chloro-N-methyl-3-pyridazinamine	143 C5H6ClN3	014959-32-1 14



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046342.D
Acq On : 19 Aug 2020 11:21
Operator : CG/JU
Sample : PB130973BL
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK73

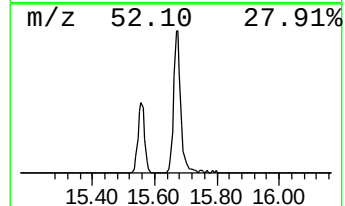
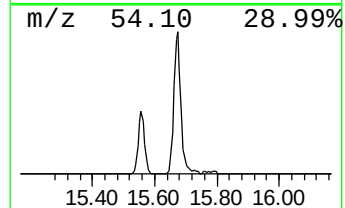
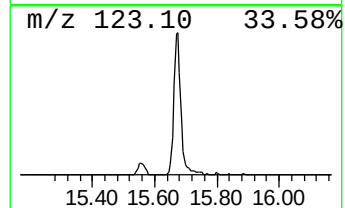
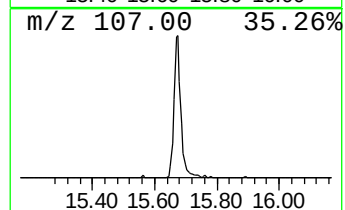
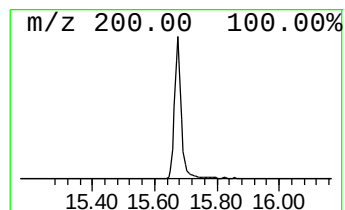
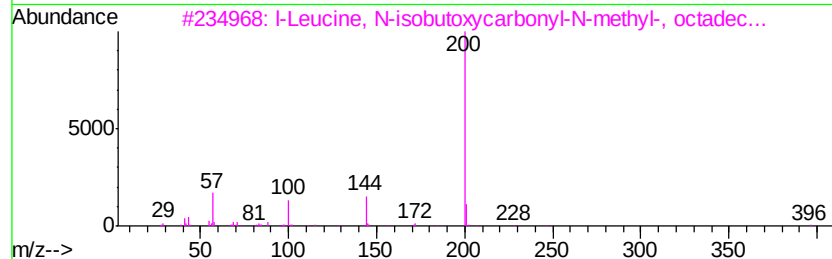
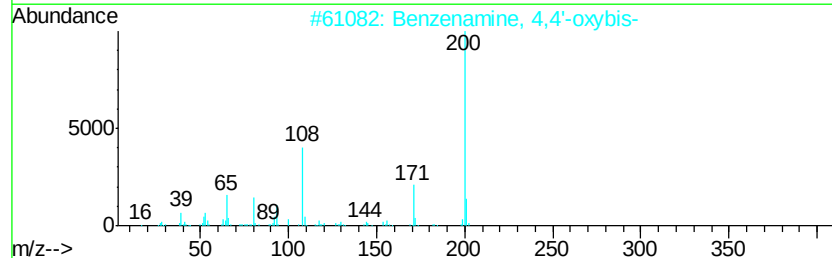
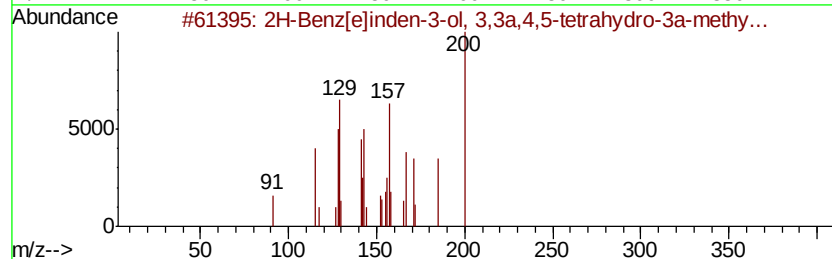
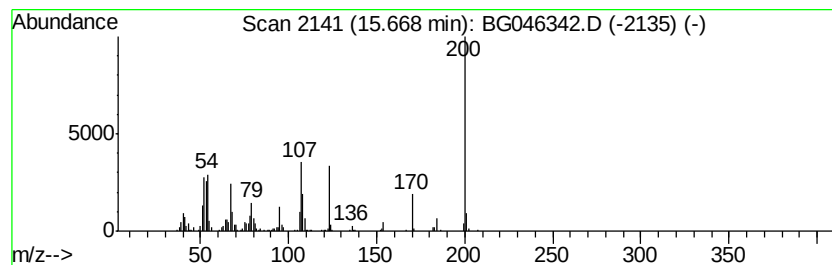
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Quant Title : SVOA CALIBRATION

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	8.06 ng/ul	375908	Acenaphthene-d10	14.56

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2H-Benz[e]indene-3-ol, 3,3a,4,5-t...	200	C14H16O	071805-91-9	38
2			Benzenamine, 4,4'-oxybis-	200	C12H12N2O	000101-80-4	14
3			l-Leucine, N-isobutoxycarbonyl-N...	497	C30H59NO4	1000321-88-1	10
4			Benzenamine, 4,4'-oxybis-	200	C12H12N2O	000101-80-4	9
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 : BG046342.D
Acq On : 19 Aug 2020 11:21
Operator : CG/JU
Sample : PB130973BL
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK73

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	2.6	ng/ul	53933	1	7.94	411512	20.0
2(3H)-Furanone, 5...	7.11	22.6	ng/ul	464870	1	7.94	411512	20.0
unknown-01	7.27	17.2	ng/ul	354710	1	7.94	411512	20.0
unknown-02	7.47	23.4	ng/ul	481184	1	7.94	411512	20.0
unknown-03	8.64	28.9	ng/ul	594705	1	7.94	411512	20.0
1H-Imidazole, 4-m...	9.09	21.8	ng/ul	448970	1	7.94	411512	20.0
unknown-04	9.82	12.2	ng/ul	381165	2	10.73	623184	20.0
unknown-05	10.86	19.4	ng/ul	603574	2	10.73	623184	20.0
unknown-06	13.97	20.8	ng/ul	970969	3	14.56	933246	20.0
unknown-07	14.76	9.1	ng/ul	424063	3	14.56	933246	20.0
unknown-08	15.67	8.1	ng/ul	375908	3	14.56	933246	20.0