

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
 Data File : BG046372.D
 Acq On : 20 Aug 2020 10:33
 Operator : CG/JU
 Sample : PB130998BL
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK98

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

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

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

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.141	5	9	17	rVB2	38177	64385	2.59%	0.285%
2	3.235	22	25	34	rVB2	8256	15342	0.62%	0.068%
3	3.394	48	52	64	rVB	19585	33267	1.34%	0.147%
4	4.140	176	179	182	rBV5	2370	3371	0.14%	0.015%
5	7.101	677	683	695	rBV	289523	531572	21.42%	2.352%
6	7.266	705	711	721	rBV	228132	395530	15.94%	1.750%
7	7.471	738	746	759	rBV	316142	544406	21.94%	2.409%
8	7.935	818	825	834	rBV	242193	412167	16.61%	1.824%
9	8.640	938	945	958	rBV	368613	660426	26.61%	2.922%
10	9.093	1013	1022	1038	rBV	268261	497346	20.04%	2.201%
11	9.816	1137	1145	1161	rBV	239459	430097	17.33%	1.903%
12	10.356	1230	1237	1254	rBV	371882	708190	28.54%	3.134%
13	10.732	1290	1301	1314	rBV	341347	623090	25.11%	2.757%
14	10.861	1314	1323	1336	rBV	358075	677131	27.29%	2.996%
15	12.324	1567	1572	1579	rBV3	5559	10026	0.40%	0.044%
16	13.970	1846	1852	1866	rBV	718058	1086929	43.80%	4.810%
17	14.257	1894	1901	1917	rBV	854706	1371590	55.27%	6.069%
18	14.569	1947	1954	1965	rBV	572916	904541	36.45%	4.002%
19	14.763	1980	1987	2006	rBV	261792	480921	19.38%	2.128%
20	15.562	2115	2123	2135	rBV	1133015	1747418	70.42%	7.732%
21	15.673	2137	2142	2153	rBV	265092	414205	16.69%	1.833%
22	15.803	2158	2164	2169	rVB3	12119	20594	0.83%	0.091%
23	16.343	2249	2256	2261	rBV4	4443	6617	0.27%	0.029%
24	17.089	2380	2383	2385	rBV2	2759	2725	0.11%	0.012%
25	17.307	2414	2420	2430	rBV	746869	1156966	46.62%	5.119%
26	17.407	2431	2437	2449	rVB	1275799	1958600	78.93%	8.667%
27	17.589	2466	2468	2475	rVB7	1753	2907	0.12%	0.013%
28	19.146	2730	2733	2737	rVB6	2178	3550	0.14%	0.016%
29	19.257	2749	2752	2756	rVV5	4755	6376	0.26%	0.028%
30	19.328	2758	2764	2769	rVV	16862	26008	1.05%	0.115%
31	19.581	2800	2807	2812	rBV	15897	27538	1.11%	0.122%
32	19.639	2815	2817	2819	rBV2	4279	2935	0.12%	0.013%
33	19.692	2820	2826	2836	rBV	1655336	2397711	96.62%	10.610%
34	19.869	2854	2856	2857	rBV2	3972	2819	0.11%	0.012%

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35	20.239	2917	2919	2923	rBV5	3953	5234	0.21%	0.023%
36	21.555	3137	3143	3154	rVB	839361	1366108	55.05%	6.045%
37	23.805	3522	3526	3533	rVB2	32915	63688	2.57%	0.282%
38	24.451	3626	3636	3651	rBV	898551	2481528	100.00%	10.980%
39	24.663	3662	3672	3691	rVB2	473183	1455594	58.66%	6.441%

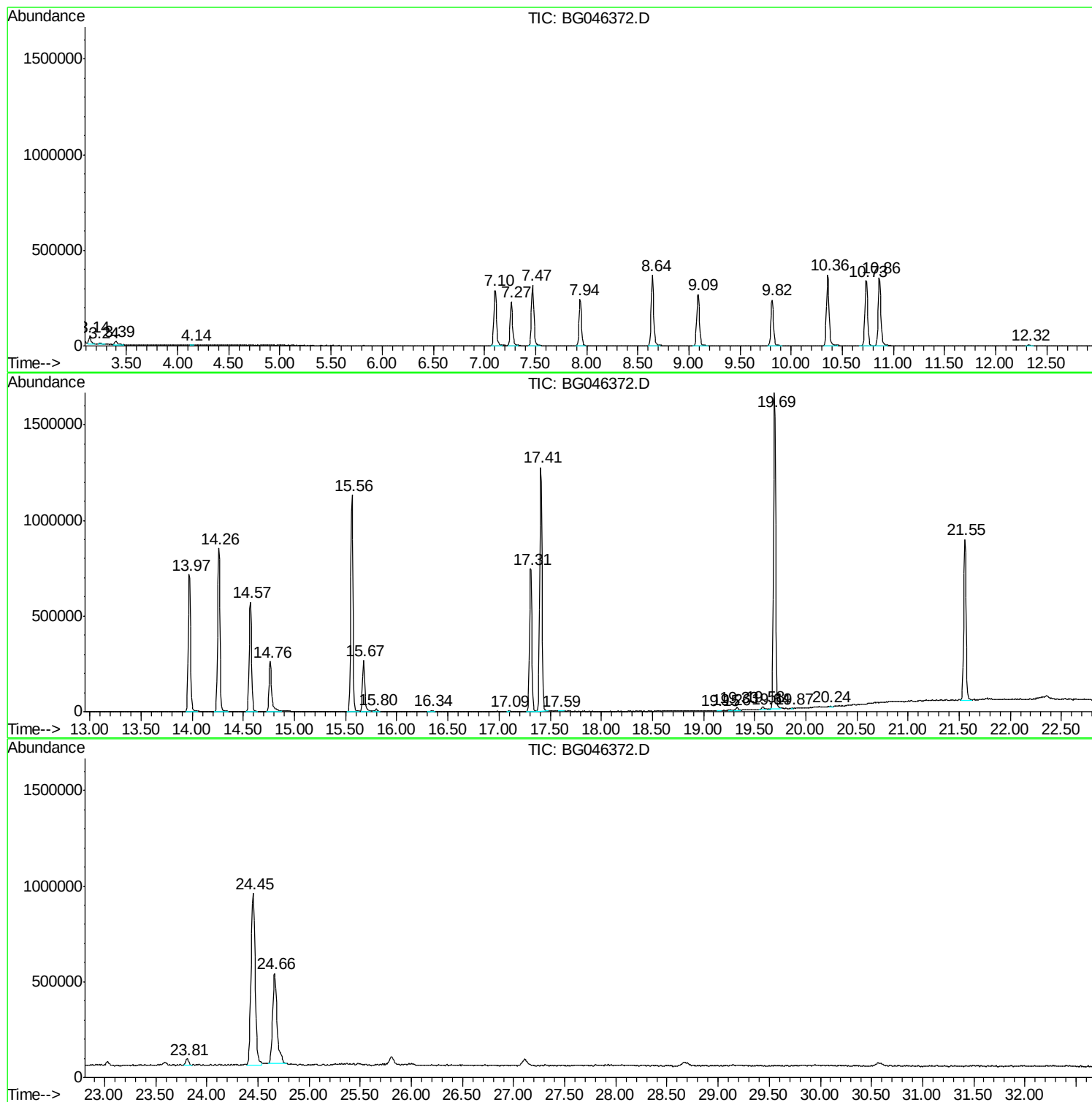
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG081320PAH.M
Quant Title : SVOA CALIBRATION

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



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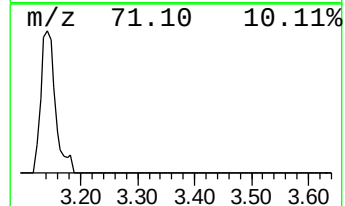
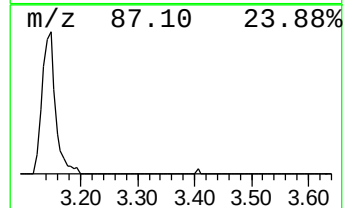
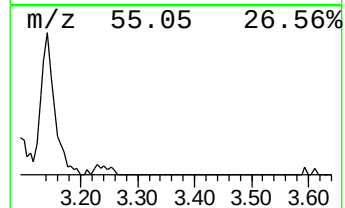
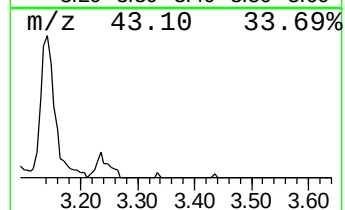
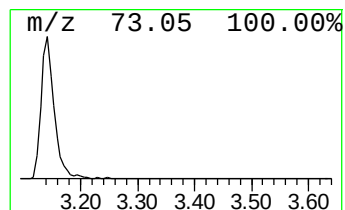
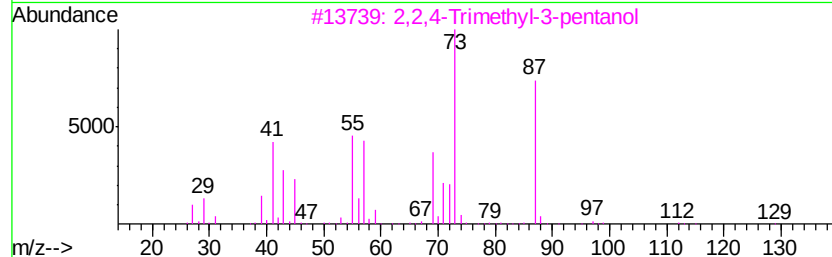
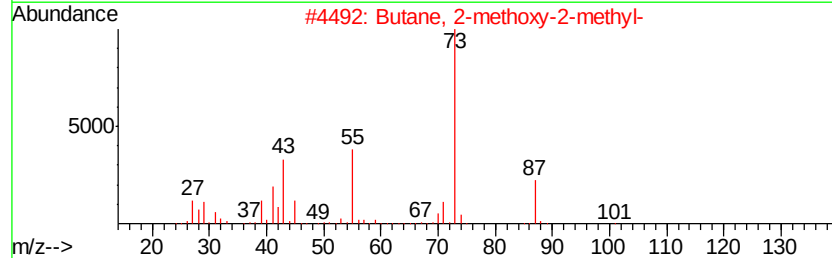
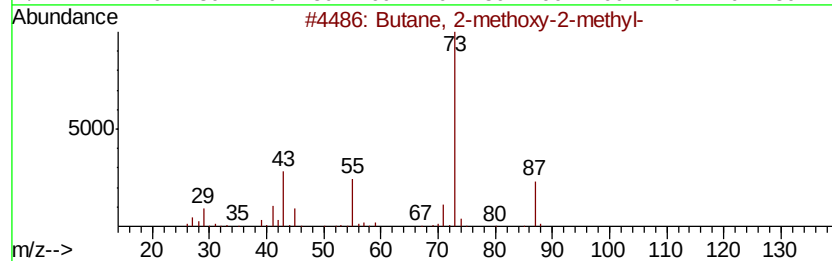
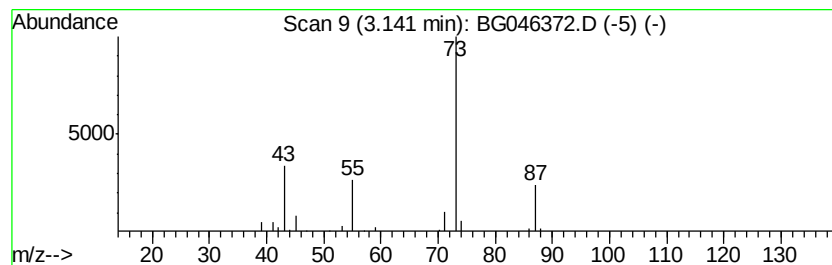
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Quant Title : SVOA CALIBRATION

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	3.12 ng/ul	64385	1,4-Dichlorobenzene-d4	7.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5			Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	1000186-02-3	9



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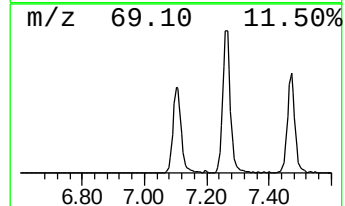
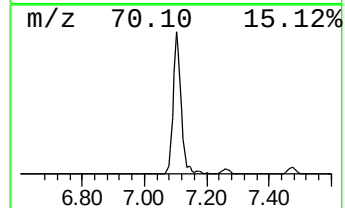
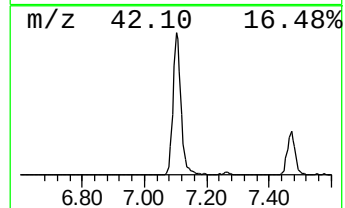
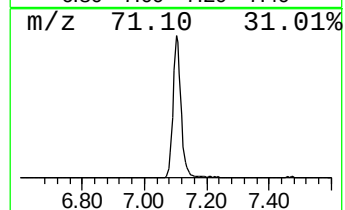
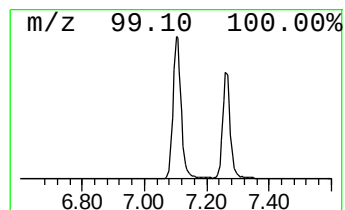
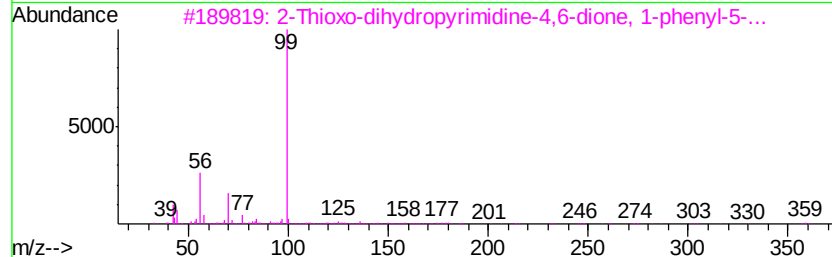
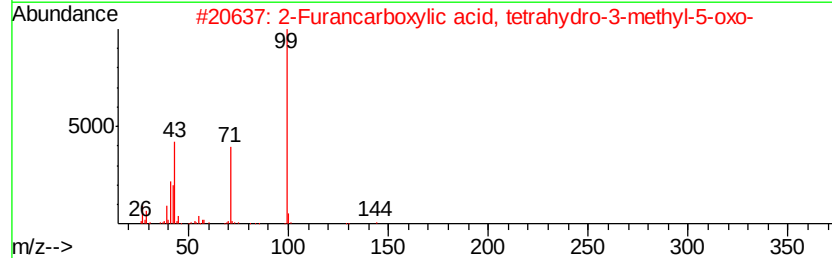
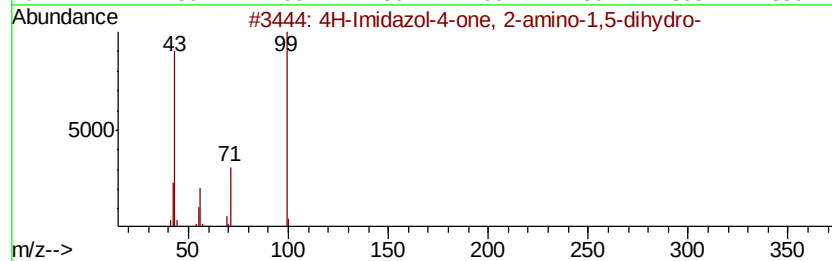
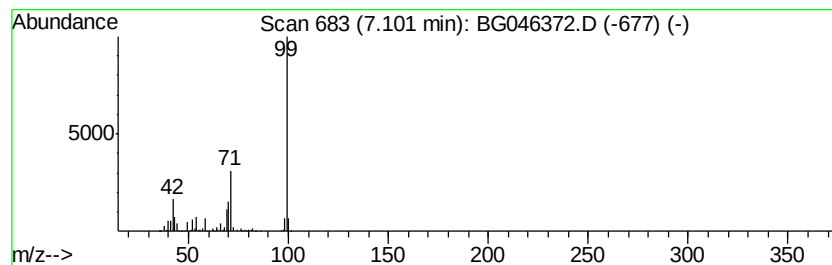
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG081320PAH.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 4H-Imidazol-4-one, 2-amino-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.10	25.79 ng/ul	531572	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Imidazol-4-one, 2-amino-1,5-d...	99	C3H5N3O	000503-86-6	64
2		2-Furancarboxylic acid, tetrahyd...	144	C6H8O4	022073-04-7	59
3		2-Thioxo-dihydropyrimidine-4,6-d...	359	C17H21N5O2S	1000304-28-6	50
4		Hexanoic acid, anhydride	214	C12H22O3	002051-49-2	45
5		Phenol-d6-	100	C6D6O	013127-88-3	43



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ClientSampled :
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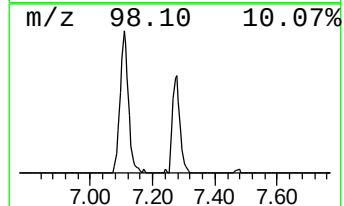
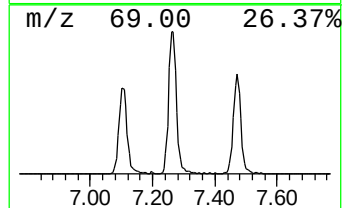
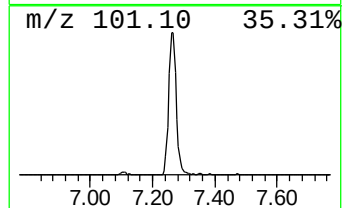
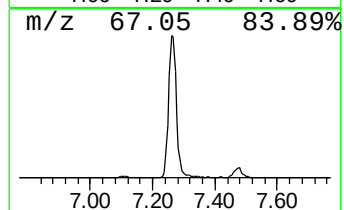
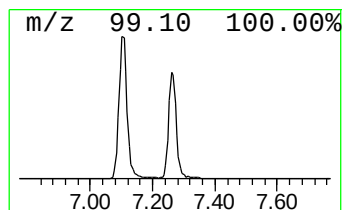
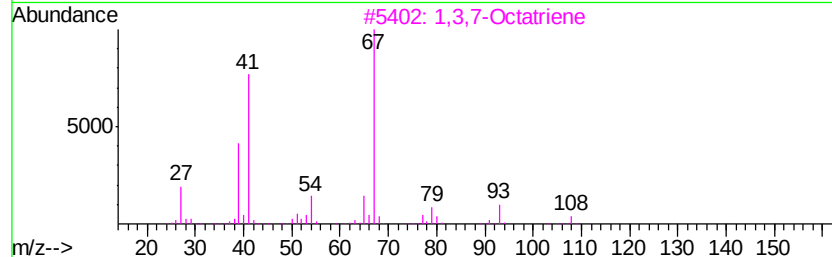
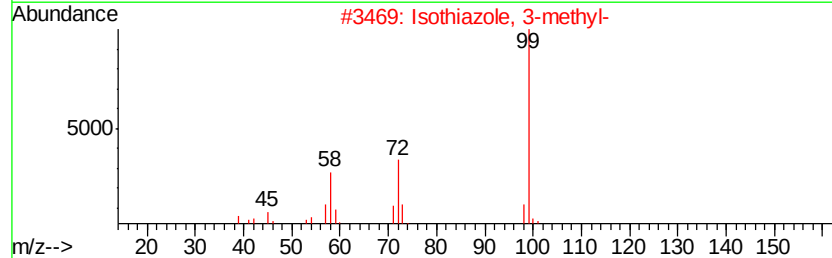
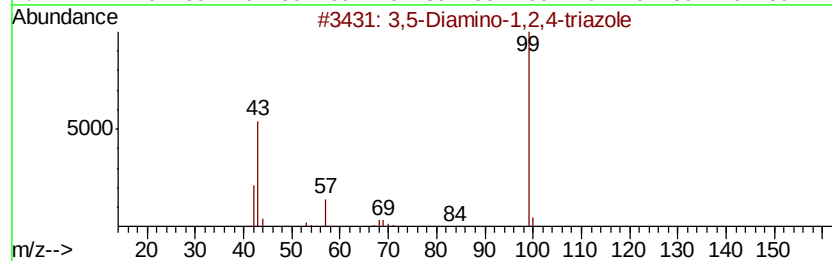
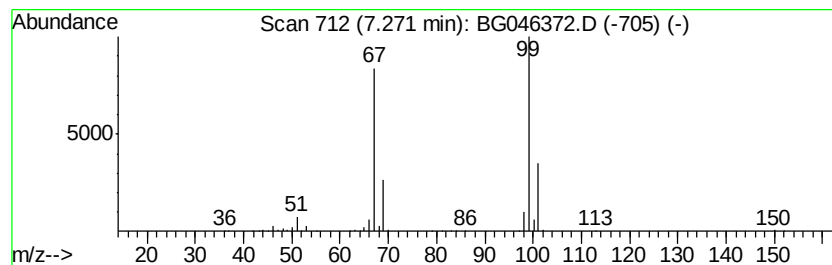
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Quant Title : SVOA CALIBRATION

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	19.19 ng/ul	395530	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3,5-Diamino-1,2,4-triazole	99	C2H5N5	001455-77-2	9
2		Isothiazole, 3-methyl-	99	C4H5NS	000693-92-5	9
3		1,3,7-Octatriene	108	C8H12	001002-35-3	9
4		1-Pentanamine, 5-(methylthio)-	133	C6H15NS	055021-78-8	9
5		Acetamide, N-2-thienyl-	141	C6H7NOS	013053-81-1	9



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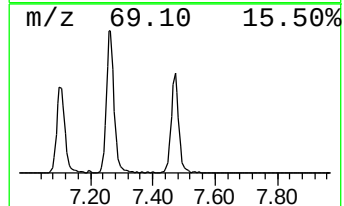
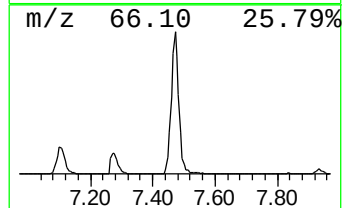
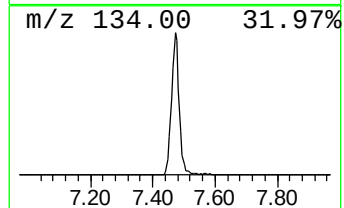
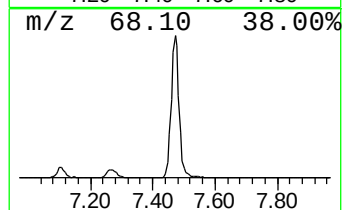
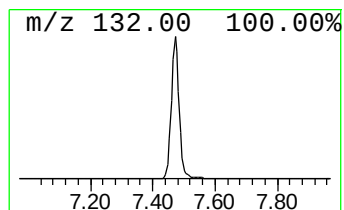
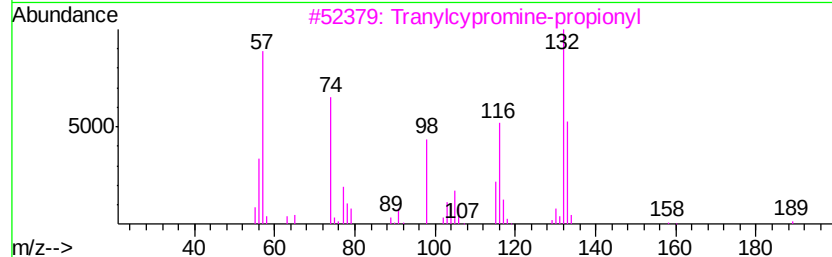
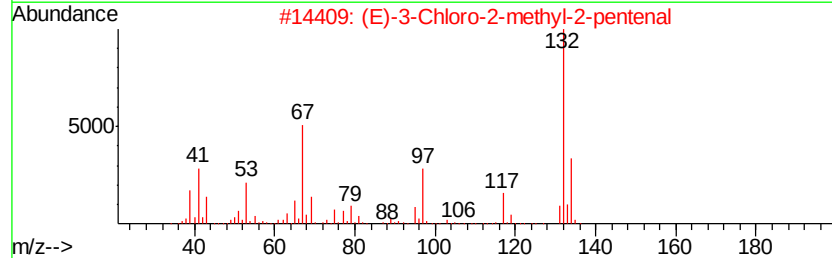
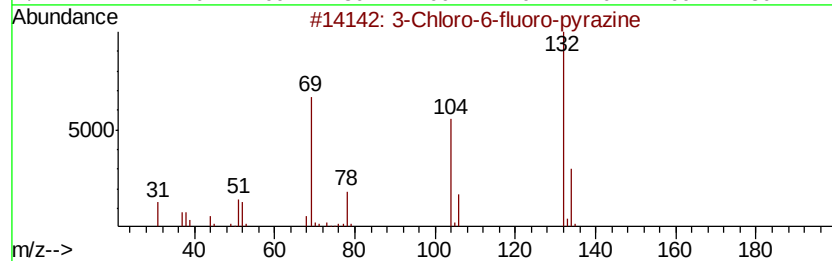
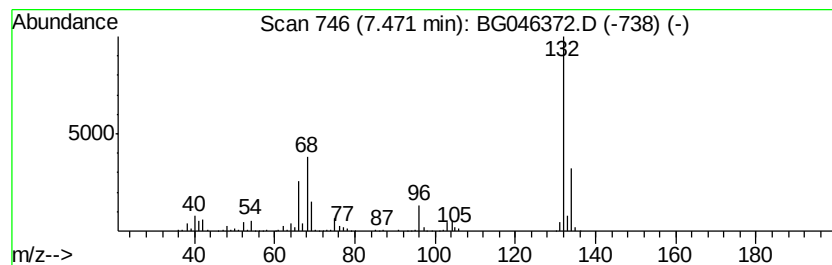
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown-02 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.47	26.42 ng/ul	544406	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	16
4		5-Aminoindole	132	C8H8N2	005192-03-0	16
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	9



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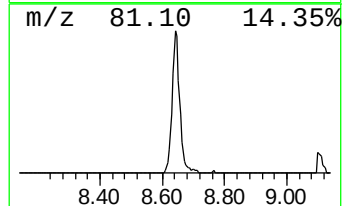
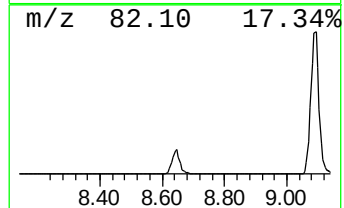
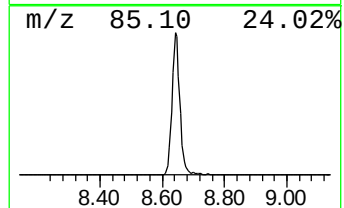
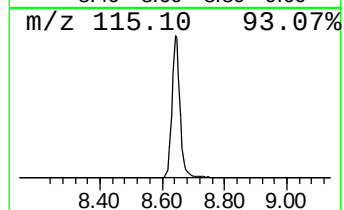
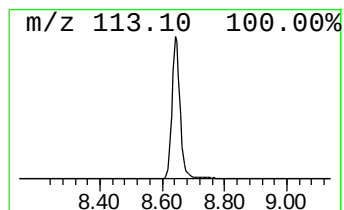
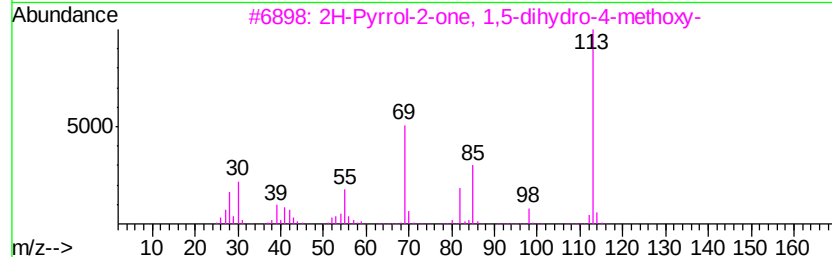
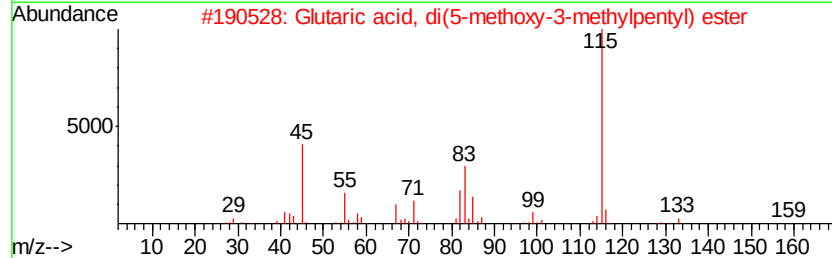
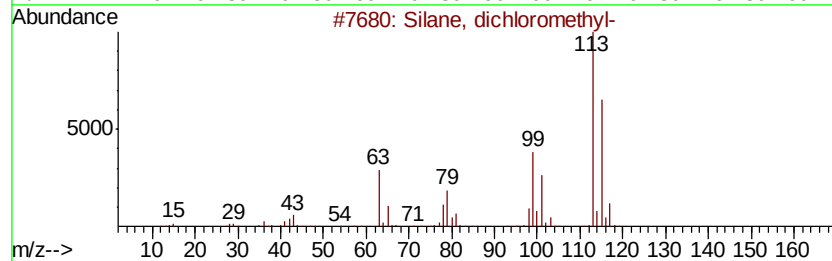
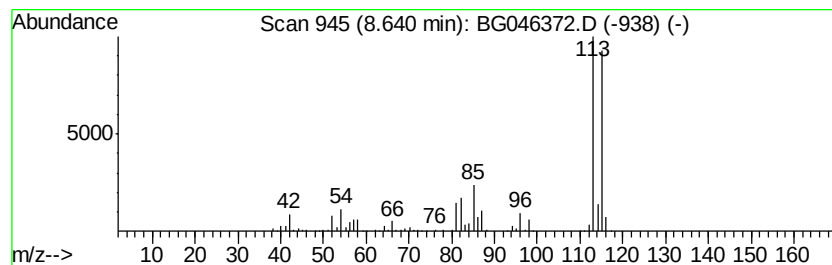
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Quant Title : SVOA CALIBRATION

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TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown-03 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.64	32.05 ng/ul	660426	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Silane, dichloromethyl-	114	CH4Cl2Si	000075-54-7	38
2		Glutaric acid, di(5-methoxy-3-me...	360	C19H36O6	1000360-08-3	35
3		2H-Pyrrol-2-one, 1,5-dihydro-4-m...	113	C5H7NO2	069778-83-2	32
4		Succinic acid, di(5-methoxy-3-me...	346	C18H34O6	1000330-26-5	32
5		Heptanoic acid, anhydride	242	C14H26O3	000626-27-7	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046372.D
Acq On : 20 Aug 2020 10:33
Operator : CG/JU
Sample : PB130998BL
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK98

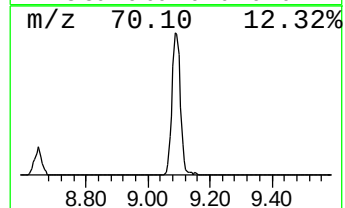
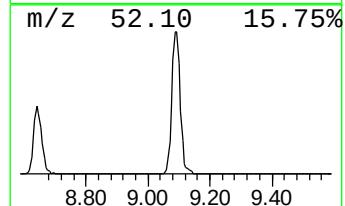
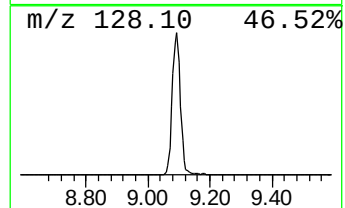
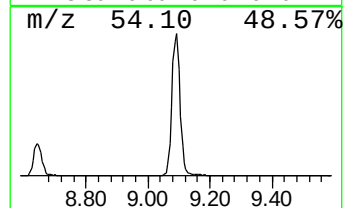
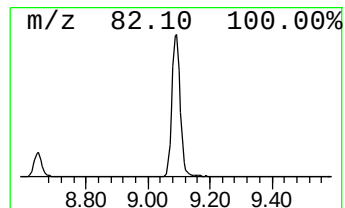
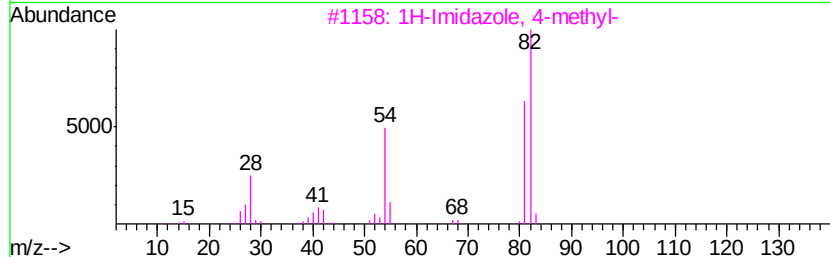
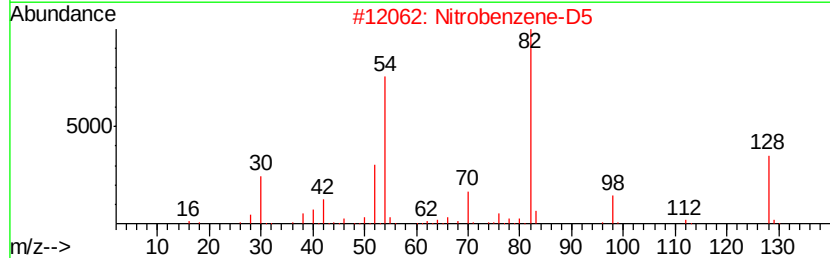
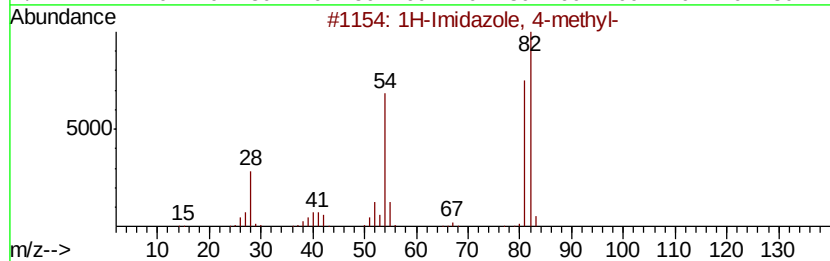
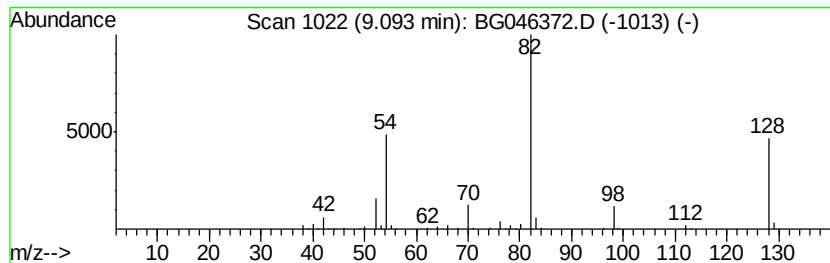
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Quant Title : SVOA CALIBRATION

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	24.13 ng/ul	497346	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		Nitrobenzene-D5	128	C6D5N02	004165-60-0	38
3		1H-Imidazole, 4-methyl-	82	C4H6N2	000822-36-6	38
4		Nitrobenzene-D5	128	C6D5N02	004165-60-0	10
5		1H-Imidazole, 2-methyl-	82	C4H6N2	000693-98-1	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
 Data File : BG046372.D
 Acq On : 20 Aug 2020 10:33
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 Sample : PB130998BL
 Misc :
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 ClientSampleId :
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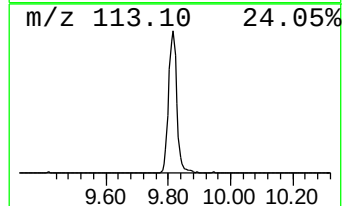
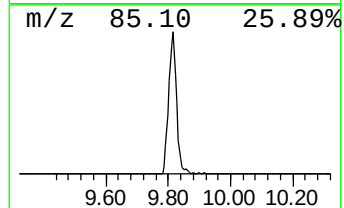
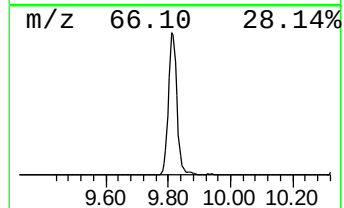
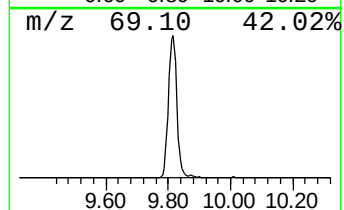
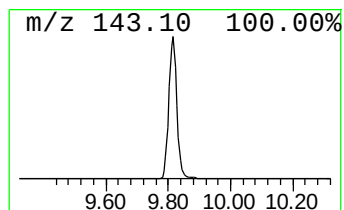
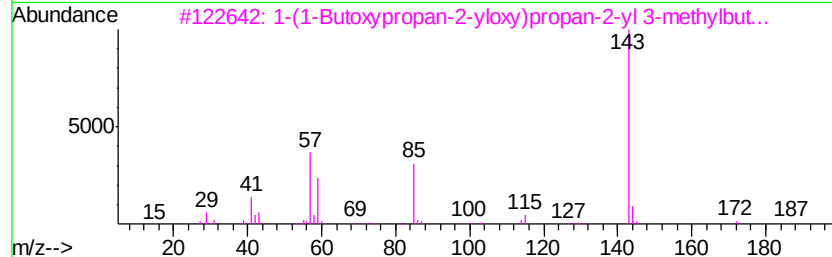
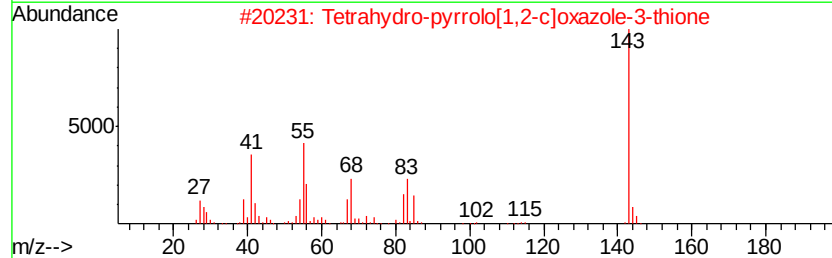
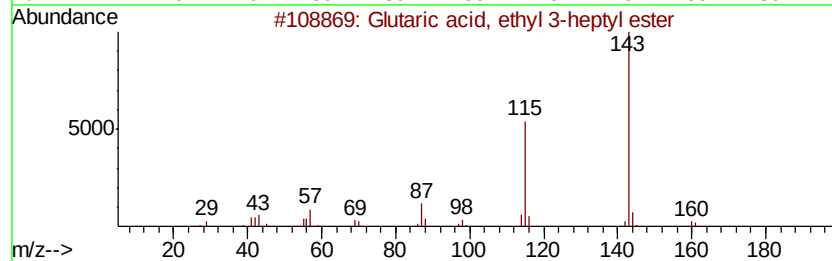
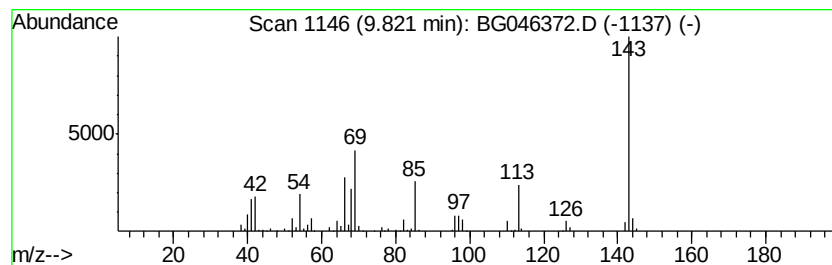
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 Quant Title : SVOA CALIBRATION

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

 Peak Number 7 unknown-04 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.82	13.81 ng/ul	430097	Napthalene-d8	10.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Glutaric acid, ethyl 3-heptyl ester	258	C14H26O4	1000359-73-0	22
2		Tetrahydro-pyrrolo[1,2-c]oxazole...	143	C6H9NOS	1000190-53-6	17
3		1-(1-Butoxypropan-2-yloxy)propan...	274	C15H30O4	1000367-12-9	16
4		1-(1-Butoxypropan-2-yloxy)propan...	274	C15H30O4	1000367-13-0	16
5		6-Chloro-N-methyl-3-pyridazinamine	143	C5H6ClN3	014959-32-1	12



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046372.D
Acq On : 20 Aug 2020 10:33
Operator : CG/JU
Sample : PB130998BL
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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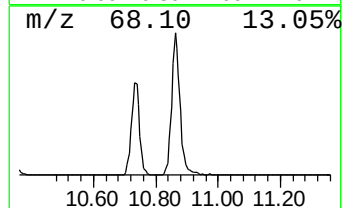
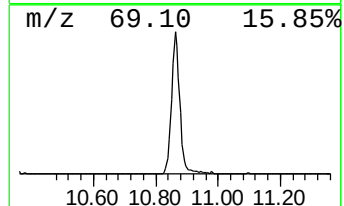
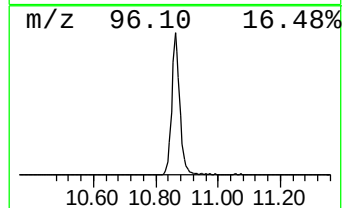
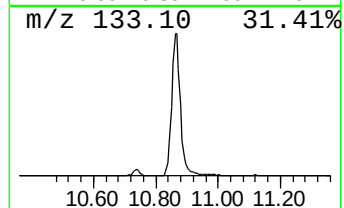
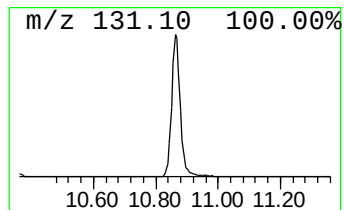
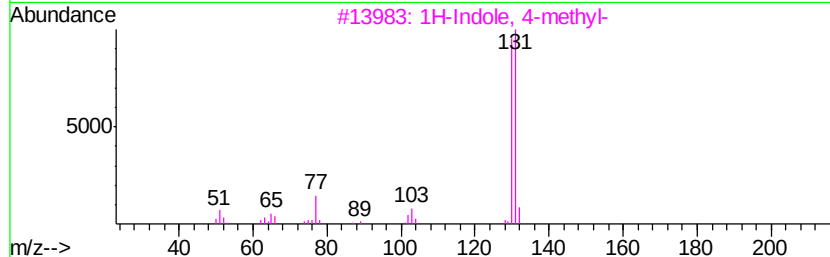
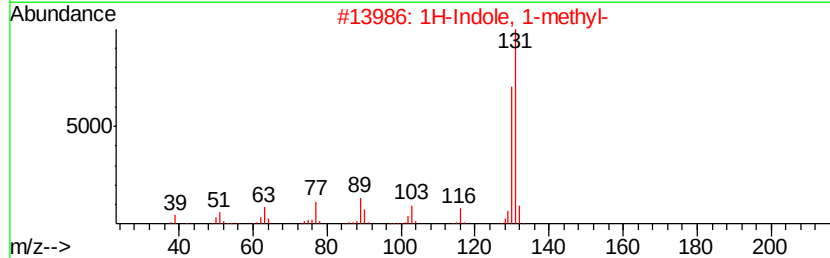
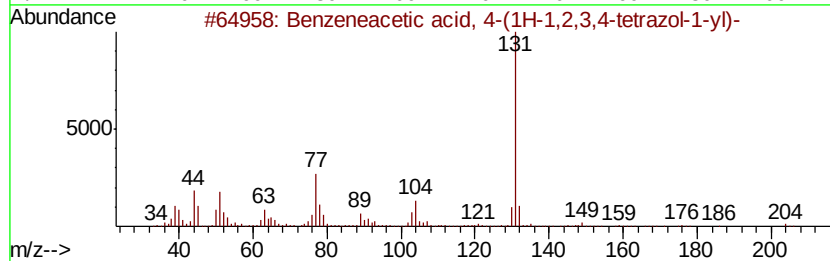
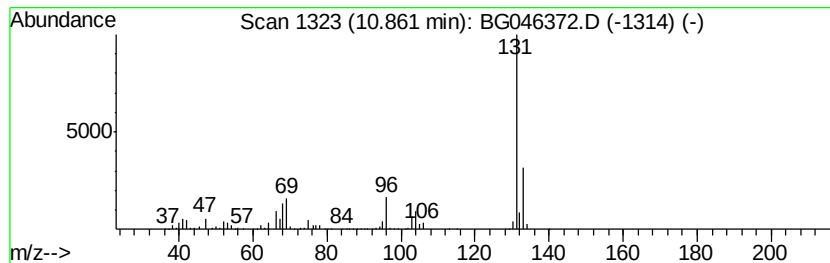
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Quant Title : SVOA CALIBRATION

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	21.73 ng/ul	677131	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		1H-Indole, 4-methyl-	131	C9H9N	016096-32-5	9
4		1H-Indole, 5-methyl-	131	C9H9N	000614-96-0	9
5		1H-Indole, 7-methyl-	131	C9H9N	000933-67-5	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046372.D
Acq On : 20 Aug 2020 10:33
Operator : CG/JU
Sample : PB130998BL
Misc :
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Instrument :
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ClientSampleId :
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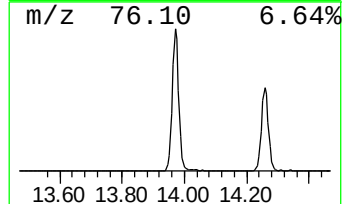
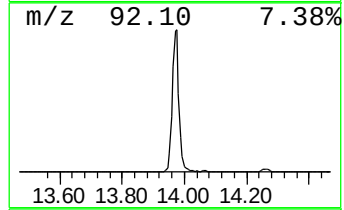
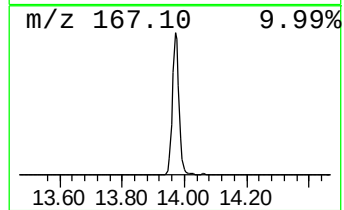
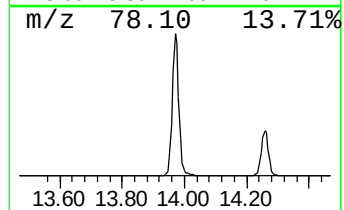
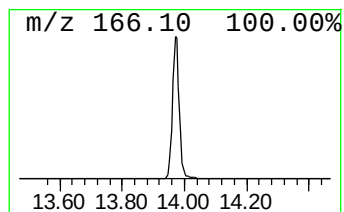
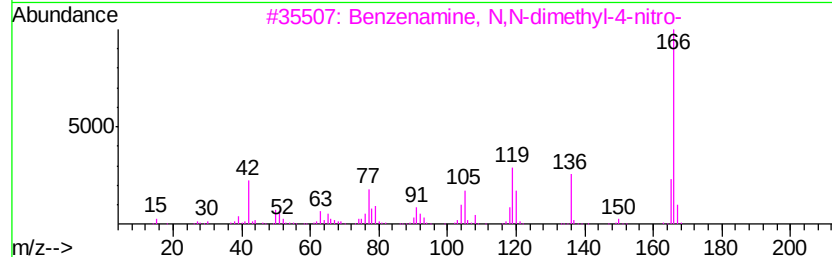
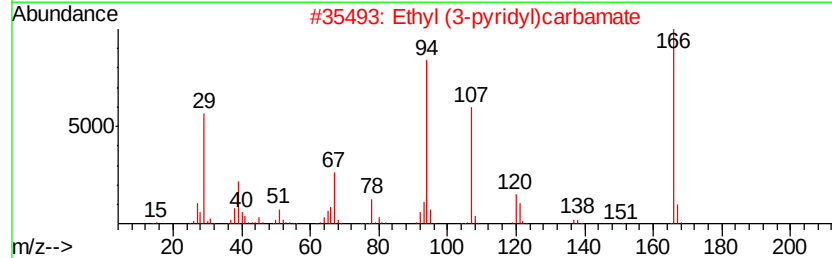
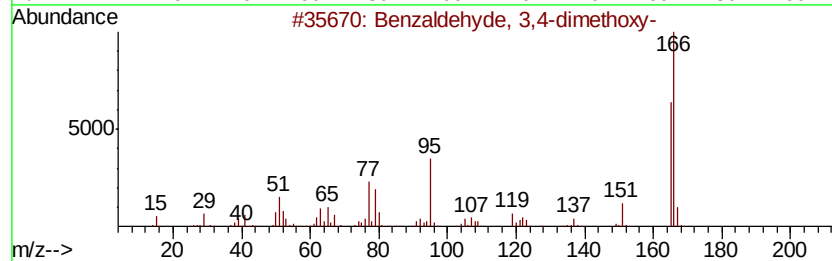
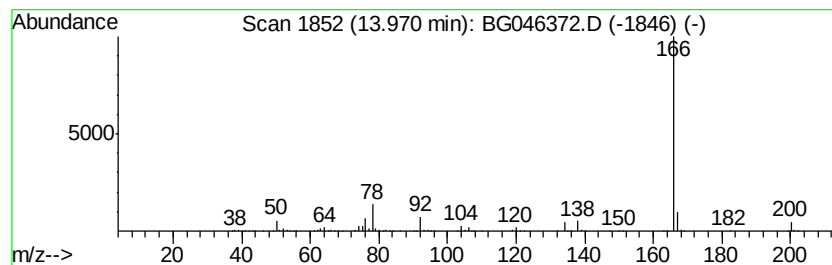
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-06 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	24.03 ng/ul	1086930	Acenaphthene-d10	14.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzaldehyde, 3,4-dimethoxy-	166	C9H10O3	000120-14-9	45
2		Ethyl (3-pyridyl)carbamate	166	C8H10N2O2	006276-11-5	39
3		Benzenamine, N,N-dimethyl-4-nitro-	166	C8H10N2O2	000100-23-2	38
4		3,5-Diamino-2-methylbenzoic acid	166	C8H10N2O2	090007-30-0	36
5		6H-Purine-6-thione, 1,7-dihydro-...	166	C6H6N4S	001126-23-4	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046372.D
Acq On : 20 Aug 2020 10:33
Operator : CG/JU
Sample : PB130998BL
Misc :
ALS Vial : 36 Sample Multiplier: 1

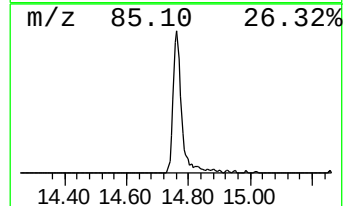
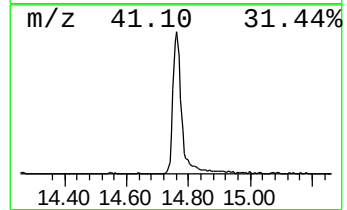
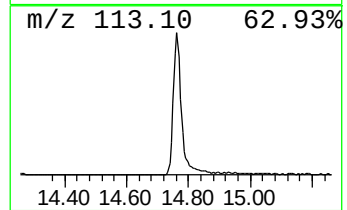
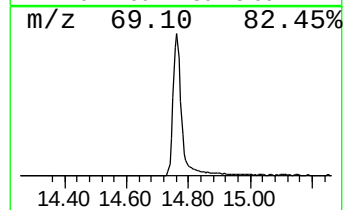
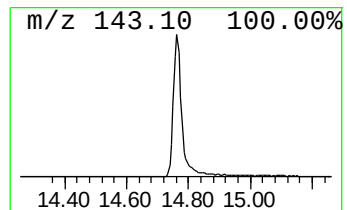
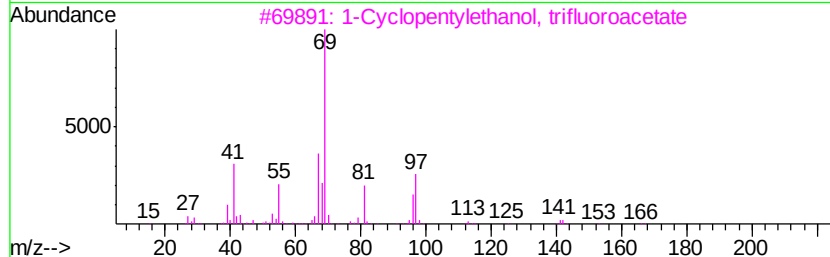
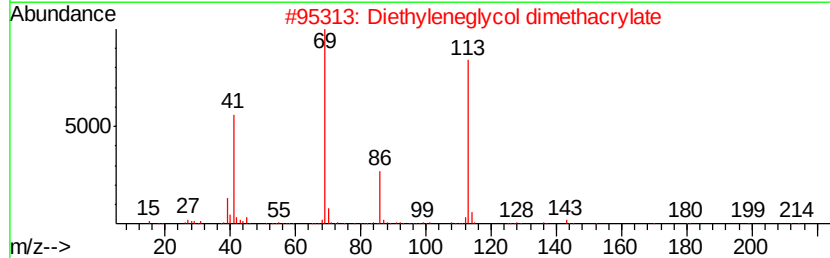
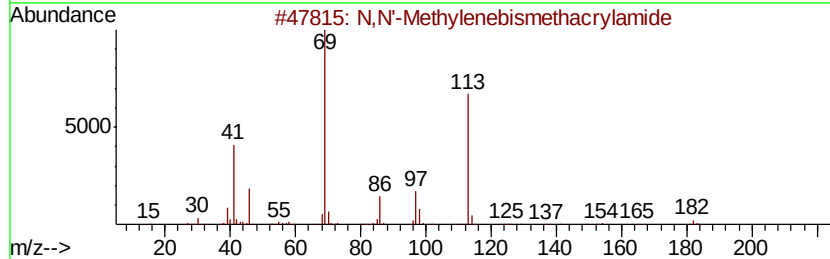
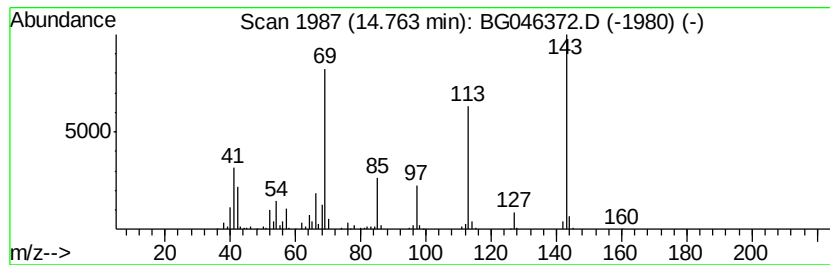
Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG081320PAH.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-07 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.76	10.63 ng/ul	480921	Acenaphthene-d10	14.57
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		N,N'-Methylenebismethacrylamide	182 C9H14N2O2	002359-15-1 25
2		Diethyleneglycol dimethacrylate	242 C12H18O5	002358-84-1 17
3		1-Cyclopentylethanol, trifluoroa...	210 C9H13F3O2	1000364-11-9 16
4		6-Aza-2-thiothymine	143 C4H5N3OS	000615-76-9 12
5		Propanenitrile, 2,2'-azobis[2-me...	164 C8H12N4	000078-67-1 12



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081920\
Data File : BG046372.D
Acq On : 20 Aug 2020 10:33
Operator : CG/JU
Sample : PB130998BL
Misc :
ALS Vial : 36 Sample Multiplier: 1

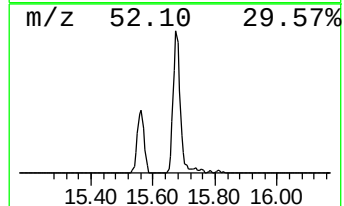
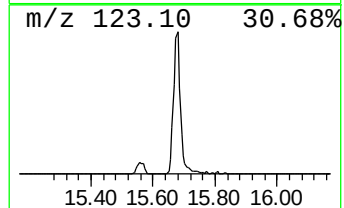
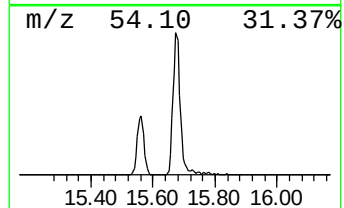
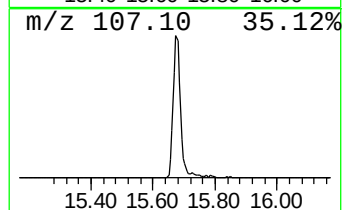
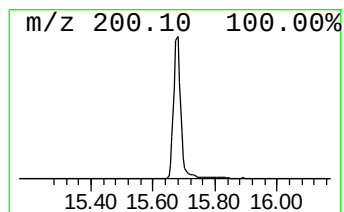
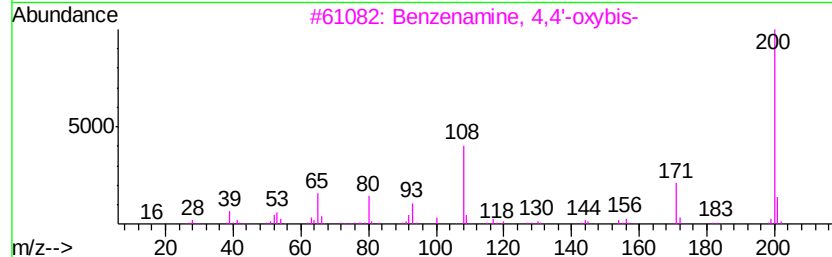
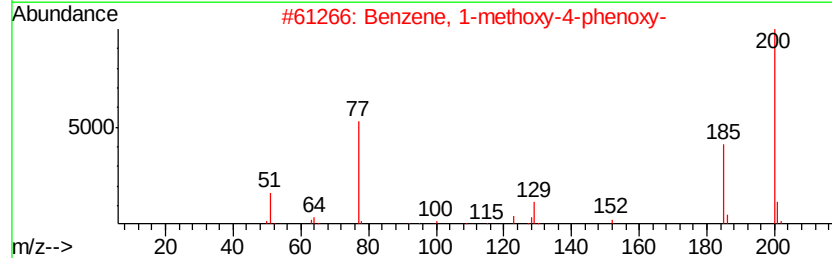
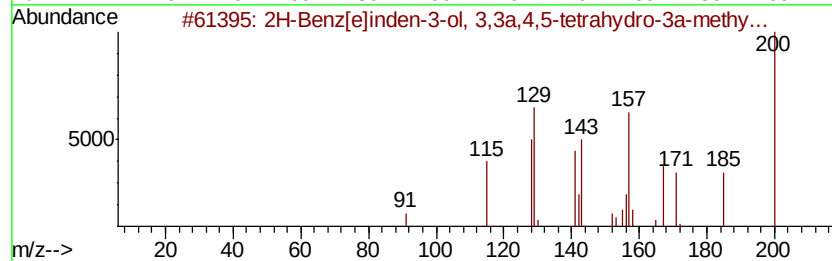
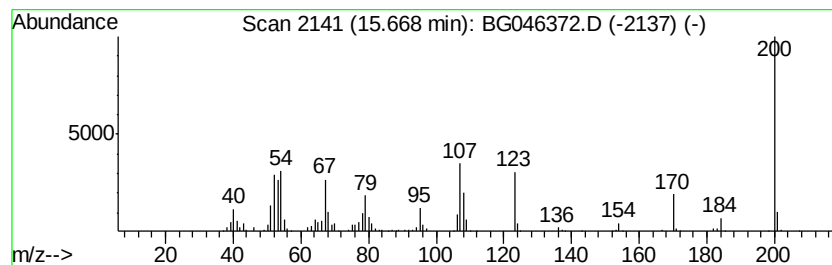
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-08 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.67	9.16 ng/ul	414205	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	43
2	Benzene, 1-methoxy-4-phenoxy-	200 C13H12O2	001655-69-2	14
3	Benzenamine, 4,4'-oxybis-	200 C12H12N2O	000101-80-4	14
4	2,2'-Bipyridyl-5-carboxylic acid	200 C11H8N2O2	001970-80-5	10
5	2-Aminocaprylic acid, N-propoxyc...	469 C28H55N04	1000325-43-2	10



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG081920\
Data File : BG046372.D
Acq On : 20 Aug 2020 10:33
Operator : CG/JU
Sample : PB130998BL
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK98

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	3.1	ng/ul	64385	1	7.94	412167	20.0
4H-Imidazol-4-one...	7.10	25.8	ng/ul	531572	1	7.94	412167	20.0
unknown-01	7.27	19.2	ng/ul	395530	1	7.94	412167	20.0
unknown-02	7.47	26.4	ng/ul	544406	1	7.94	412167	20.0
unknown-03	8.64	32.0	ng/ul	660426	1	7.94	412167	20.0
1H-Imidazole, 4-m...	9.09	24.1	ng/ul	497346	1	7.94	412167	20.0
unknown-04	9.82	13.8	ng/ul	430097	2	10.73	623090	20.0
unknown-05	10.86	21.7	ng/ul	677131	2	10.73	623090	20.0
unknown-06	13.97	24.0	ng/ul	1086930	3	14.57	904541	20.0
unknown-07	14.76	10.6	ng/ul	480921	3	14.57	904541	20.0
unknown-08	15.67	9.2	ng/ul	414205	3	14.57	904541	20.0