

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG091515\
 Data File : BG018699.D
 Acq On : 16 Sep 2015 1:23
 Operator : UM/IZ
 Sample : G3641-06DL 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 B0AM4DL

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:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG091015.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	2.884	6	8	12	rVB2	12558	15120	0.34%	0.065%
2	3.107	41	46	53	rBV2	49902	85727	1.92%	0.368%
3	3.178	53	58	69	rVB	706479	1000961	22.39%	4.298%
4	3.477	105	109	117	rBV	20124	32484	0.73%	0.139%
5	3.777	156	160	162	rVB2	4189	5520	0.12%	0.024%
6	4.112	212	217	221	rVB8	4714	7406	0.17%	0.032%
7	5.140	390	392	399	rVB5	5437	8141	0.18%	0.035%
8	5.446	438	444	450	rVV	78509	118910	2.66%	0.511%
9	5.933	520	527	535	rVB	290007	483299	10.81%	2.075%
10	6.292	580	588	597	rVB7	7489	17848	0.40%	0.077%
11	6.450	612	615	619	rBV5	4474	6473	0.14%	0.028%
12	6.527	624	628	632	rVB3	7071	10738	0.24%	0.046%
13	7.167	731	737	746	rBV3	37249	72188	1.62%	0.310%
14	7.326	756	764	773	rBV	124106	199512	4.46%	0.857%
15	7.420	773	780	787	rBV	815394	1246752	27.89%	5.353%
16	7.537	793	800	809	rVB2	126474	222844	4.99%	0.957%
17	8.001	873	879	889	rVB2	198464	339595	7.60%	1.458%
18	8.425	944	951	957	rBV	300660	488711	10.93%	2.098%
19	8.472	957	959	965	rVB4	7339	10720	0.24%	0.046%
20	8.542	965	971	973	rBV4	4883	6041	0.14%	0.026%
21	8.618	976	984	992	rBV	180077	300323	6.72%	1.290%
22	8.701	992	998	1006	rVB	106950	187576	4.20%	0.805%
23	9.159	1070	1076	1085	rVB	136373	237631	5.32%	1.020%
24	9.382	1108	1114	1117	rVB5	4777	7083	0.16%	0.030%
25	9.882	1188	1199	1206	rBV2	113398	209309	4.68%	0.899%
26	10.422	1282	1291	1299	rBV	198138	355213	7.95%	1.525%
27	10.610	1317	1323	1328	rVB5	10918	20855	0.47%	0.090%
28	10.804	1346	1356	1371	rBV	297013	534877	11.97%	2.297%
29	10.939	1373	1379	1383	rVB3	9212	14839	0.33%	0.064%
30	11.127	1404	1411	1418	rBV	189882	333120	7.45%	1.430%
31	11.932	1537	1548	1556	rBV	2667100	4469646	100.00%	19.192%
32	12.009	1556	1561	1568	rVB4	27728	47408	1.06%	0.204%
33	12.167	1578	1588	1595	rBV3	38039	81425	1.82%	0.350%
34	12.878	1707	1709	1712	rVV3	6719	7145	0.16%	0.031%

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Integration Parameters: LSCINT.P

Integrator: RTE
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 Sampling : 1
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Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG091015.M
 Title : SVOA CALIBRATION

35	12.984	1723	1727	1731	rVB4	3288	5057	0.11%	0.022%
36	13.195	1760	1763	1767	rBV2	5964	9588	0.21%	0.041%
37	13.472	1808	1810	1816	rBV4	4670	5602	0.13%	0.024%
38	13.624	1830	1836	1843	rBV7	5489	11605	0.26%	0.050%
39	13.942	1883	1890	1895	rBV3	18356	33272	0.74%	0.143%
40	13.994	1895	1899	1900	rVV3	11883	17203	0.38%	0.074%
41	14.030	1900	1905	1912	rVB	396475	575689	12.88%	2.472%
42	14.323	1946	1955	1961	rBV	461583	734375	16.43%	3.153%
43	14.629	1998	2007	2015	rBV3	517809	890617	19.93%	3.824%
44	14.770	2028	2031	2035	rVB4	5390	7000	0.16%	0.030%
45	14.829	2035	2041	2045	rVV2	29538	54117	1.21%	0.232%
46	14.858	2045	2046	2050	rVV3	5624	5832	0.13%	0.025%
47	14.929	2054	2058	2063	rVB6	6782	10547	0.24%	0.045%
48	15.029	2069	2075	2080	rBV6	4779	8710	0.19%	0.037%
49	15.099	2082	2087	2093	rBV	44870	63675	1.42%	0.273%
50	15.434	2139	2144	2148	rBV3	9283	13090	0.29%	0.056%
51	15.616	2166	2175	2183	rBV	667837	1000150	22.38%	4.294%
52	15.728	2189	2194	2200	rBV	111965	166514	3.73%	0.715%
53	15.857	2211	2216	2220	rVB6	7693	14911	0.33%	0.064%
54	16.121	2254	2261	2266	rBV3	10763	20225	0.45%	0.087%
55	16.198	2268	2274	2278	rBV6	7842	11175	0.25%	0.048%
56	16.503	2322	2326	2330	rVB5	7473	10922	0.24%	0.047%
57	16.568	2335	2337	2341	rVV4	4747	6356	0.14%	0.027%
58	16.627	2341	2347	2354	rVV6	9619	17223	0.39%	0.074%
59	16.920	2392	2397	2403	rVB3	40752	62592	1.40%	0.269%
60	17.050	2415	2419	2425	rVB5	7897	11425	0.26%	0.049%
61	17.155	2433	2437	2444	rVB5	10032	17253	0.39%	0.074%
62	17.367	2467	2473	2480	rVV2	779656	1145160	25.62%	4.917%
63	17.426	2480	2483	2484	rVV	12128	11614	0.26%	0.050%
64	17.467	2484	2490	2496	rVV2	781326	1137624	25.45%	4.885%
65	18.013	2581	2583	2587	rVV4	5024	6125	0.14%	0.026%
66	18.066	2587	2592	2598	rVV2	86874	119425	2.67%	0.513%
67	18.143	2600	2605	2607	rVB3	4902	8258	0.18%	0.035%
68	18.195	2610	2614	2617	rVB6	2880	4886	0.11%	0.021%
69	18.260	2620	2625	2629	rBV6	5379	9128	0.20%	0.039%
70	18.325	2633	2636	2639	rVV	11071	12584	0.28%	0.054%
71	18.883	2726	2731	2733	rBV4	5031	6840	0.15%	0.029%

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Integration Parameters: LSCINT.P

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72	18.994	2746	2750	2754	rVV7	5852	10146	0.23%	0.044%
73	19.059	2758	2761	2762	rBV2	6971	6600	0.15%	0.028%
74	19.294	2796	2801	2805	rBV3	23415	31741	0.71%	0.136%
75	19.388	2814	2817	2823	rBV7	8443	12882	0.29%	0.055%
76	19.647	2855	2861	2866	rVB2	9619	17421	0.39%	0.075%
77	19.752	2870	2879	2885	rBV	937245	1338532	29.95%	5.747%
78	19.929	2903	2909	2911	rBV6	4040	7881	0.18%	0.034%
79	20.058	2930	2931	2934	rBV3	8046	7507	0.17%	0.032%
80	20.405	2986	2990	2991	rBV4	4110	5146	0.12%	0.022%
81	20.422	2991	2993	2997	rVB5	5590	6491	0.15%	0.028%
82	20.575	3017	3019	3022	rVB4	7366	7421	0.17%	0.032%
83	20.675	3032	3036	3040	rVB7	14481	21692	0.49%	0.093%
84	20.892	3070	3073	3076	rBV5	10754	13634	0.31%	0.059%
85	21.627	3192	3198	3205	rBV	860120	1372577	30.71%	5.894%
86	21.938	3246	3251	3255	rBV2	66180	115410	2.58%	0.496%
87	22.155	3286	3288	3289	rBV2	6691	4928	0.11%	0.021%
88	23.107	3445	3450	3453	rBV7	8970	16337	0.37%	0.070%
89	23.536	3521	3523	3524	rBV	7831	4985	0.11%	0.021%
90	23.883	3580	3582	3583	rBV2	7364	6452	0.14%	0.028%
91	24.588	3692	3702	3715	rBV	494378	1336961	29.91%	5.741%
92	24.811	3728	3740	3751	rBV2	487119	1405514	31.45%	6.035%
93	25.945	3925	3933	3940	rBV2	15987	48594	1.09%	0.209%
94	28.160	4308	4310	4311	rBV2	8288	5368	0.12%	0.023%
95	28.601	4381	4385	4388	rBV6	6591	9036	0.20%	0.039%
96	28.877	4428	4432	4435	rBV6	11331	18838	0.42%	0.081%
97	29.294	4501	4503	4505	rBV3	6064	4756	0.11%	0.020%
98	29.788	4585	4587	4590	rVB3	8325	6660	0.15%	0.029%
99	32.144	4986	4988	4990	rBV3	6983	6872	0.15%	0.030%
100	32.473	5042	5044	5046	rBV3	7155	7203	0.16%	0.031%

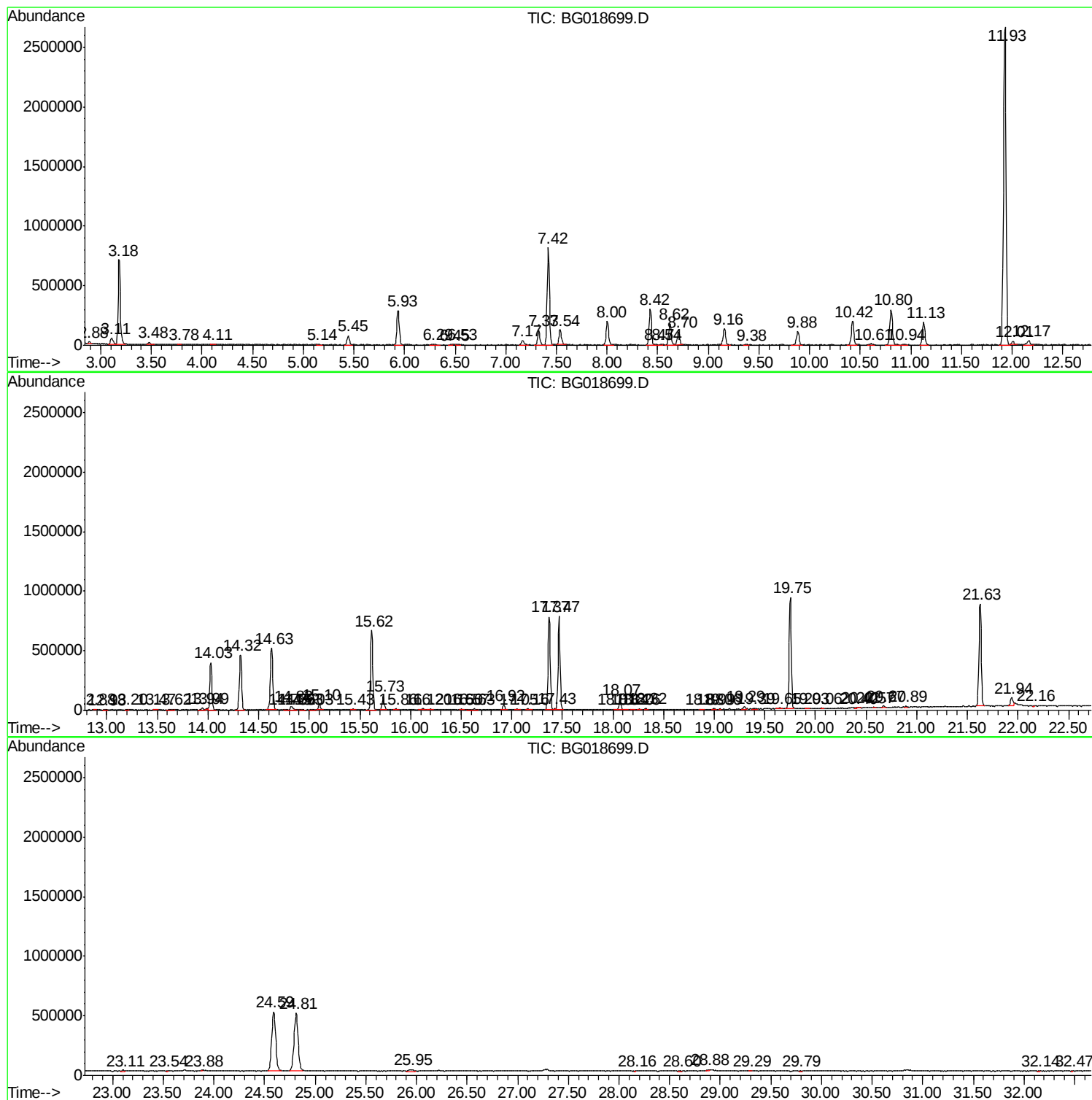
Sum of corrected areas: 23289394

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Instrument :
BNA_G
ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG091015.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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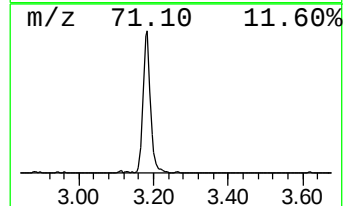
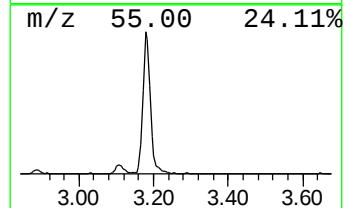
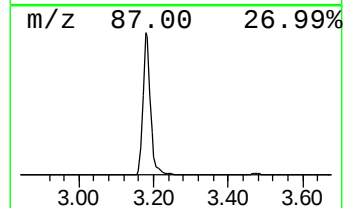
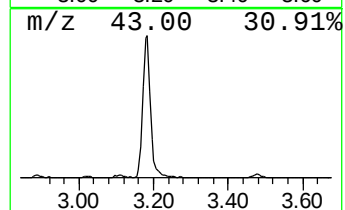
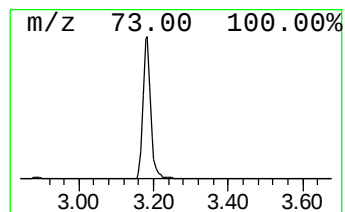
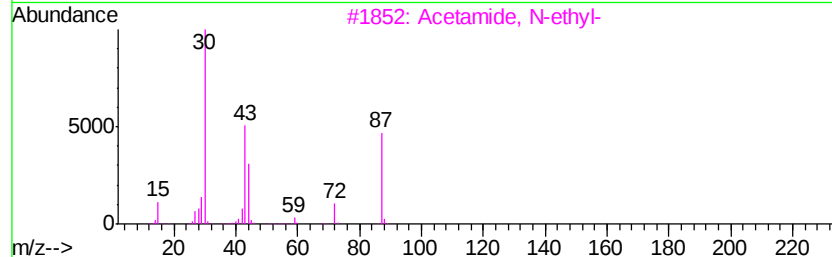
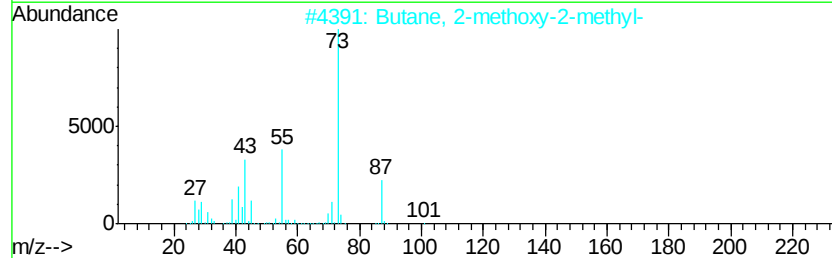
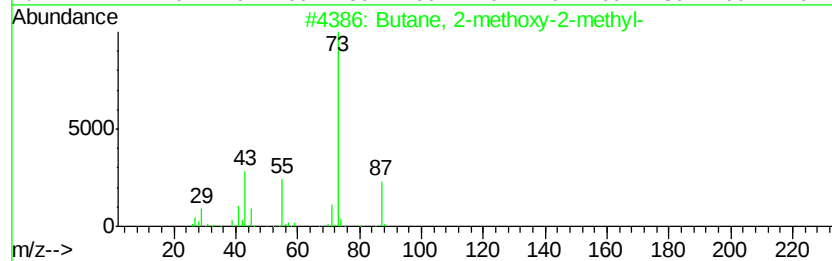
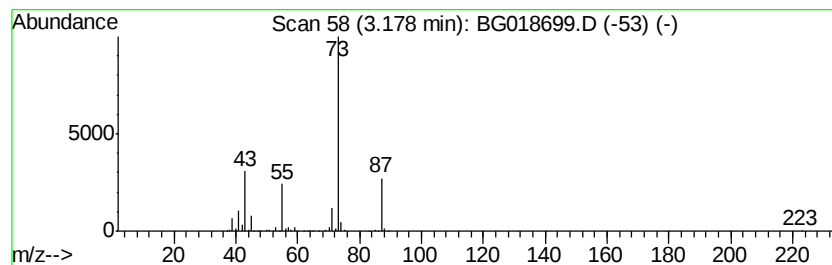
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG091015.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.18	58.95 ng/ul	1000960	1,4-Dichlorobenzene-d4	8.00

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83	
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78	
3	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9	
4	N-Ethylformamide	73	C3H7NO	000627-45-2	9	
5	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9	



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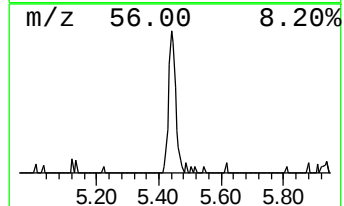
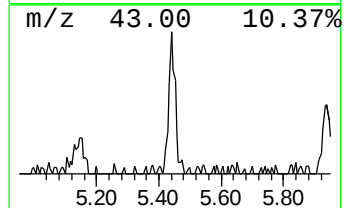
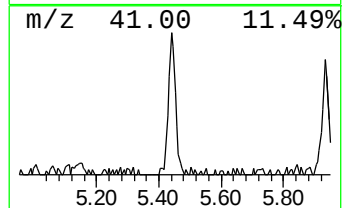
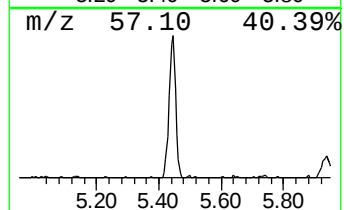
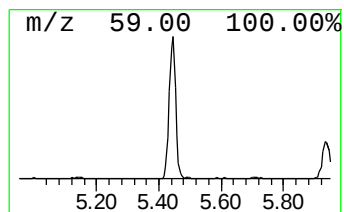
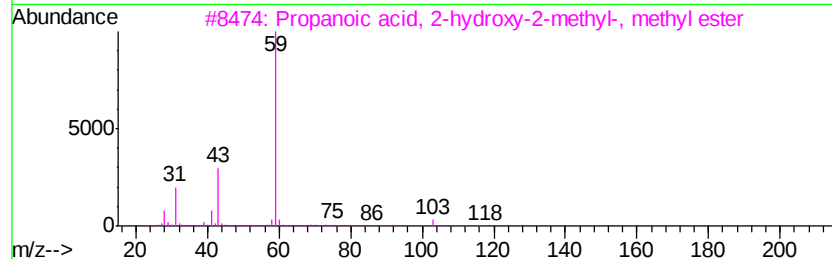
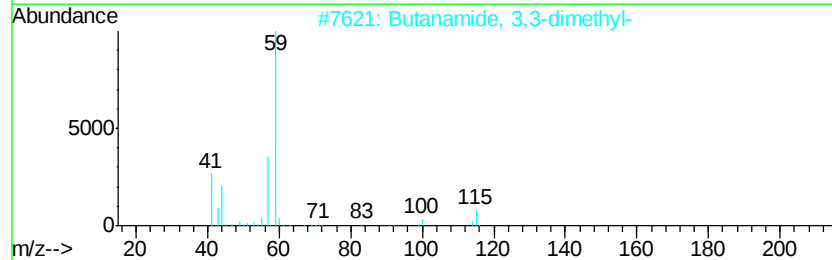
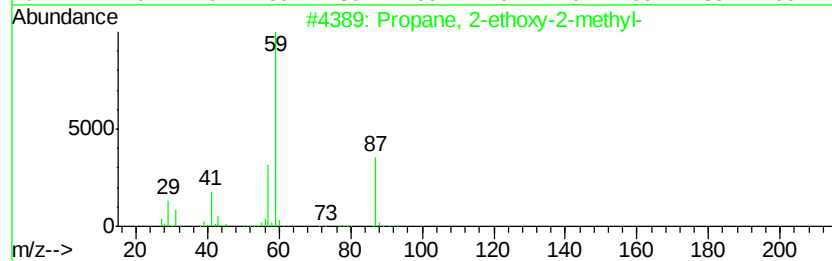
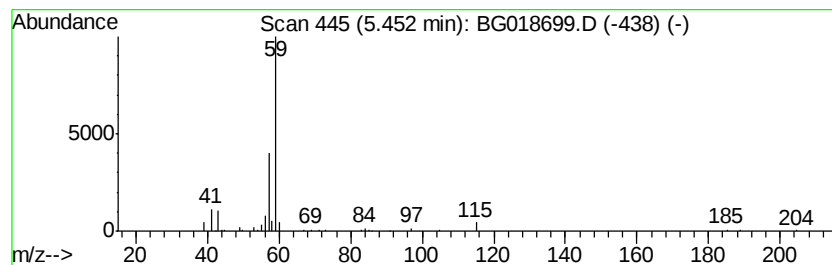
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG091015.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Propane, 2-ethoxy-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.45	7.00 ng/ul	118910	1,4-Dichlorobenzene-d4	8.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	56
2		Butanamide, 3,3-dimethyl-	115	C6H13NO	000926-04-5	56
3		Propanoic acid, 2-hydroxy-2-methyl-	118	C5H10O3	002110-78-3	38
4		2-Hydroxy-2,5-dimethyl-hept-6-en-	156	C9H16O2	1000192-55-8	9
5		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	9



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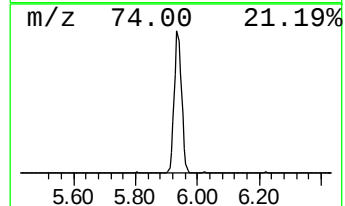
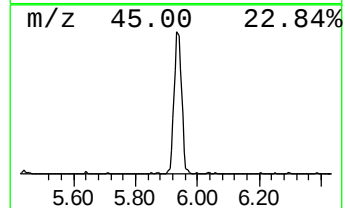
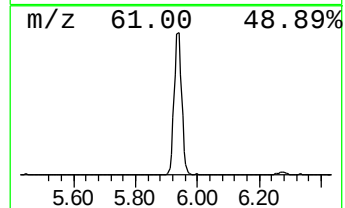
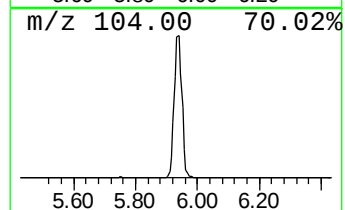
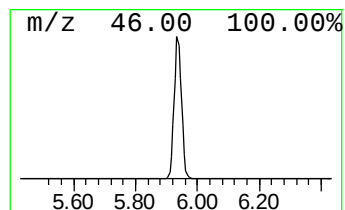
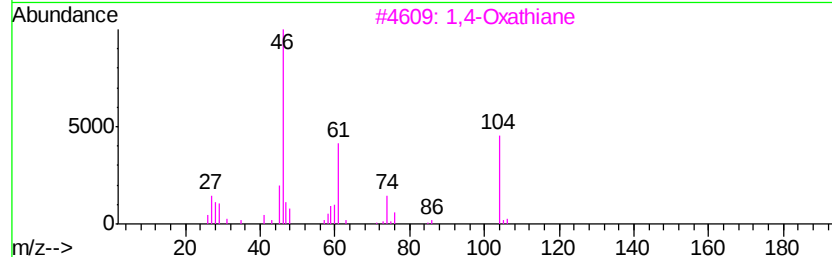
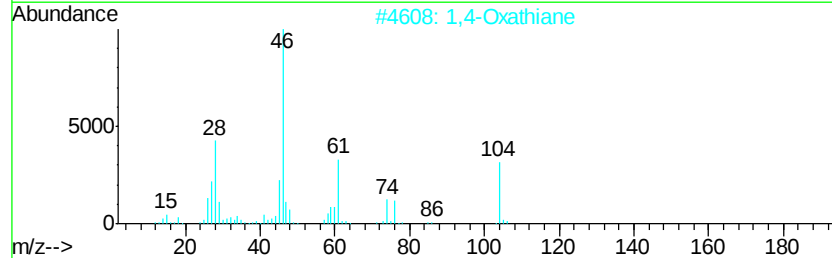
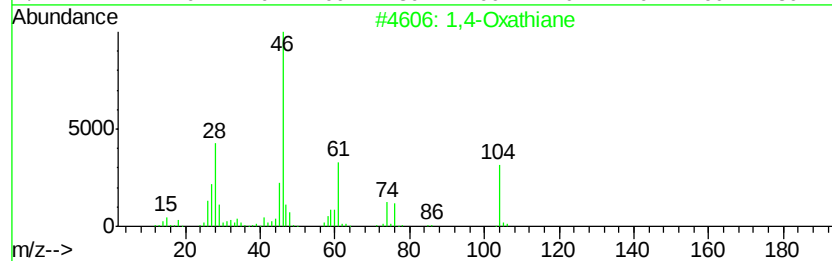
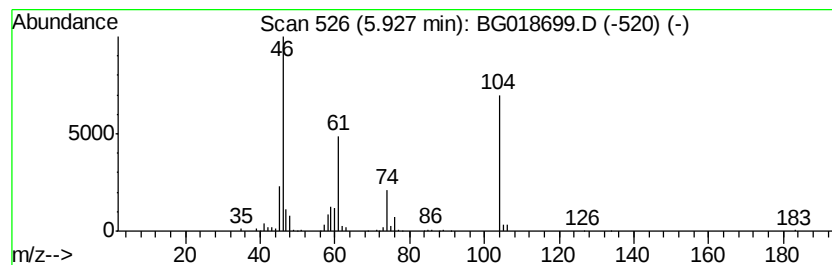
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1,4-Oxathiane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.93	28.46 ng/ul	483299	1,4-Dichlorobenzene-d4	8.00
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1,4-Oxathiane	104 C4H8OS	015980-15-1	91
2	1,4-Oxathiane	104 C4H8OS	015980-15-1	91
3	1,4-Oxathiane	104 C4H8OS	015980-15-1	91
4	Thietane	74 C3H6S	000287-27-4	43
5	Dihydro-3-(2H)-thiophenone	102 C4H6OS	001003-04-9	43



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG091515\
Data File : BG018699.D
Acq On : 16 Sep 2015 1:23
Operator : UM/IZ
Sample : G3641-06DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B0AM4DL

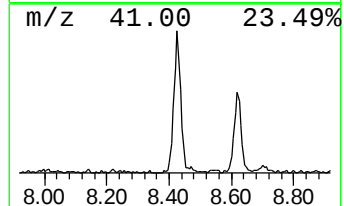
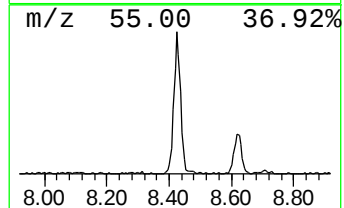
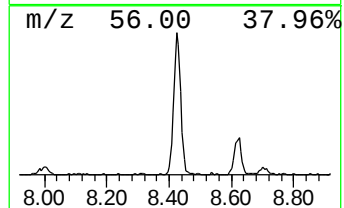
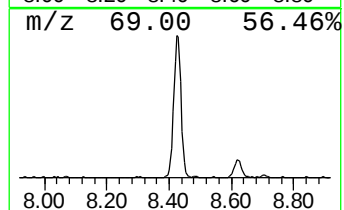
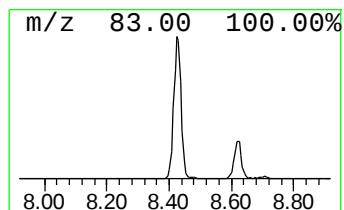
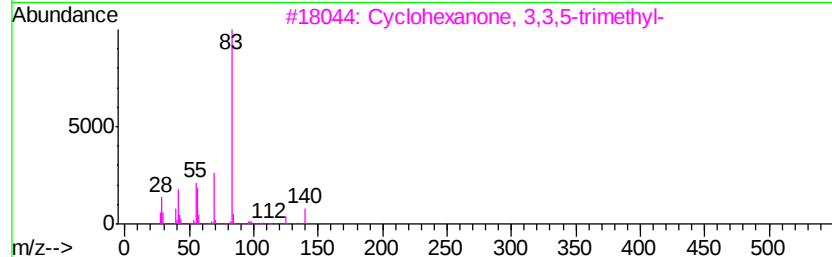
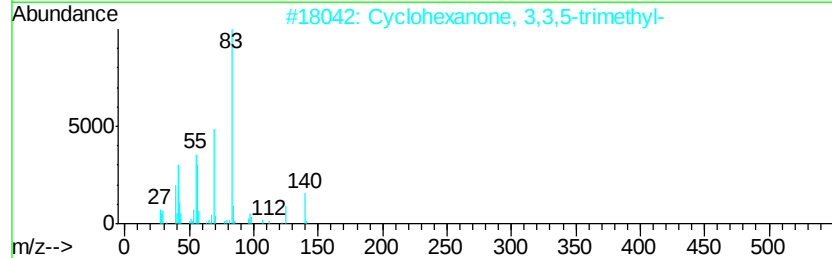
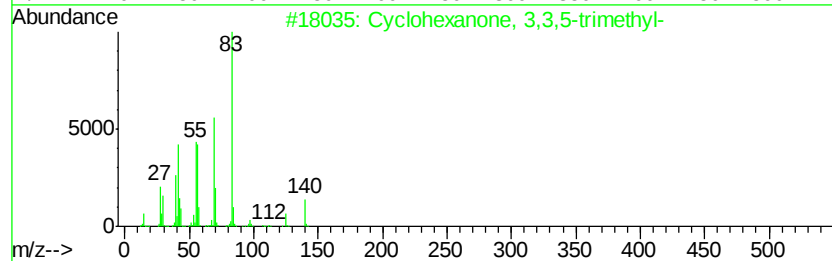
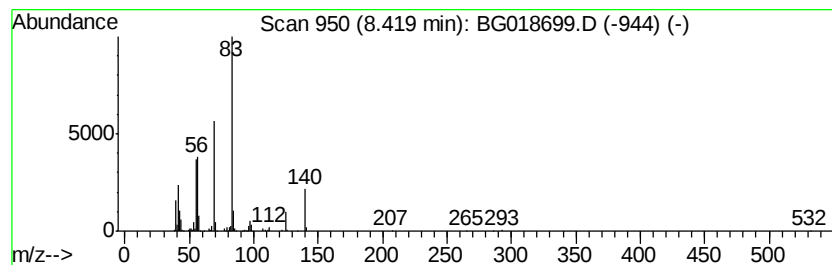
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Cyclohexanone, 3,3,5-trimet... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.42	28.78 ng/ul	488711	1,4-Dichlorobenzene-d4	8.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	91
2		Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	87
3		Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	64
4		Cyclohexane, butyl-	140	C10H20	001678-93-9	59
5		Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	58



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG091515\
Data File : BG018699.D
Acq On : 16 Sep 2015 1:23
Operator : UM/IZ
Sample : G3641-06DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B0AM4DL

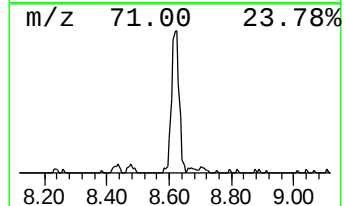
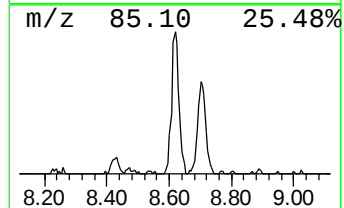
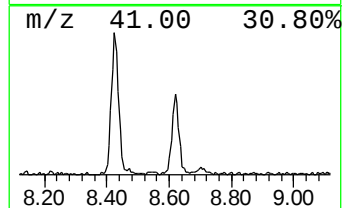
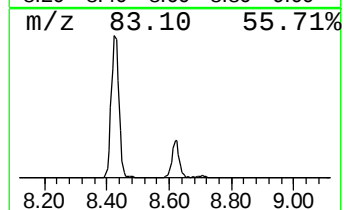
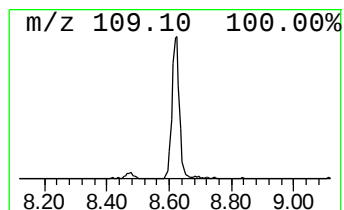
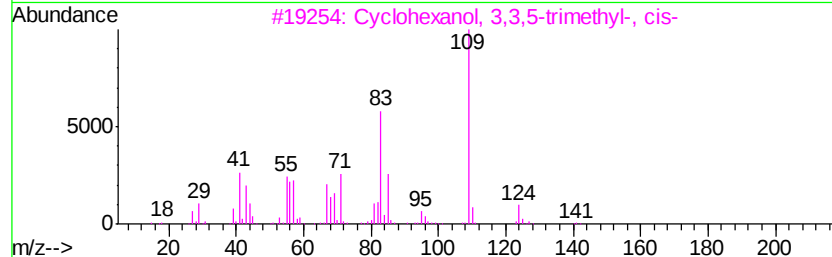
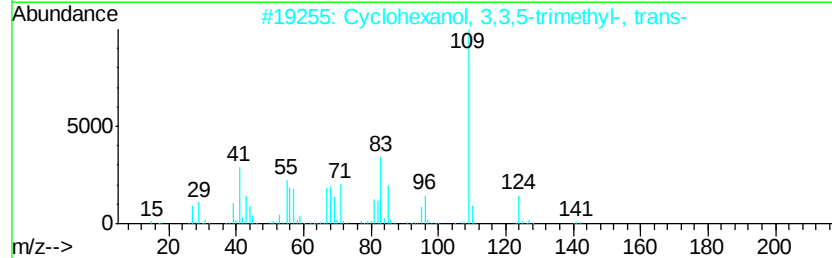
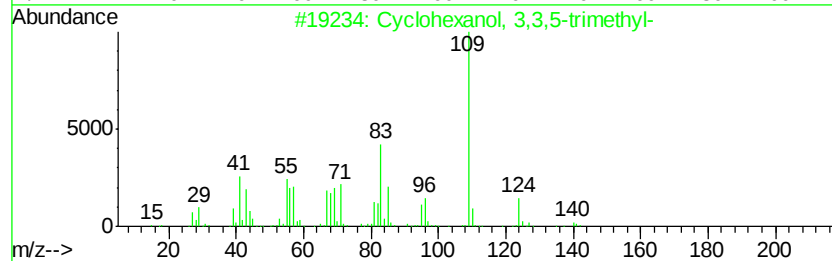
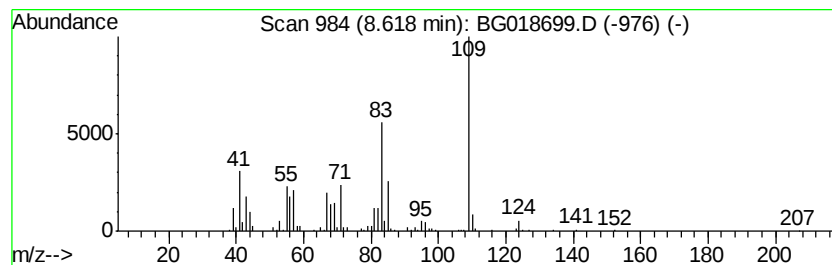
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Cyclohexanol, 3,3,5-trimeth... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.62	17.69 ng/ul	300323	1,4-Dichlorobenzene-d4	8.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanol, 3,3,5-trimethyl-	142	C9H18O	000116-02-9	64
2		Cyclohexanol, 3,3,5-trimethyl-, ...	142	C9H18O	000767-54-4	53
3		Cyclohexanol, 3,3,5-trimethyl-, ...	142	C9H18O	000933-48-2	43
4		Cyclohexanol, 3,3,5-trimethyl-, ...	184	C11H20O2	024691-16-5	40
5		Cyclohexanol, 3,3,5-trimethyl-	142	C9H18O	000116-02-9	38



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG091515\
Data File : BG018699.D
Acq On : 16 Sep 2015 1:23
Operator : UM/IZ
Sample : G3641-06DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B0AM4DL

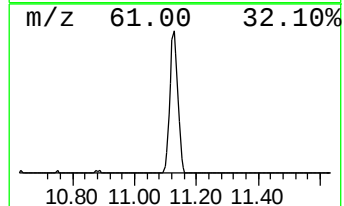
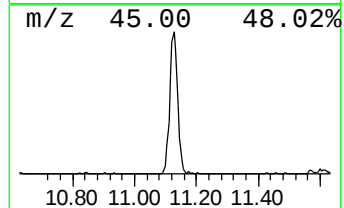
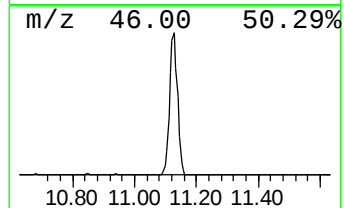
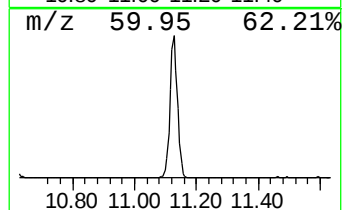
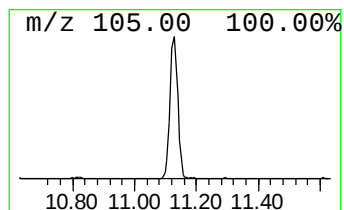
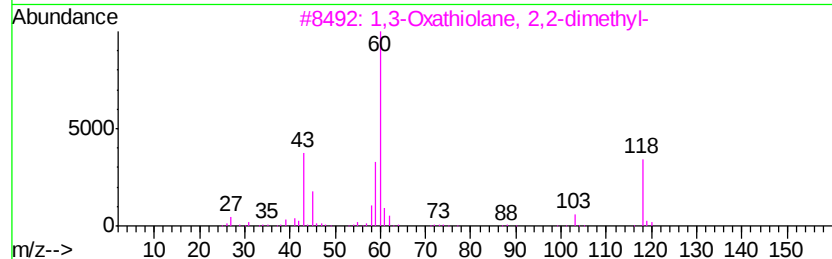
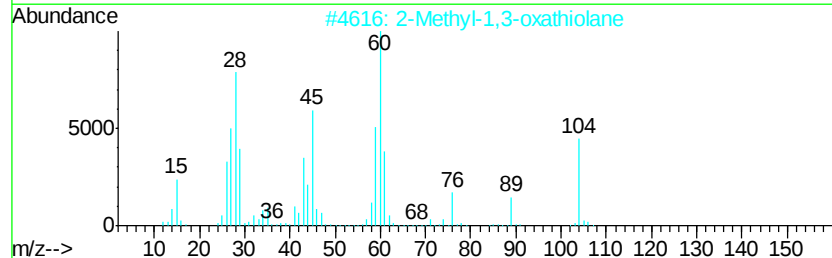
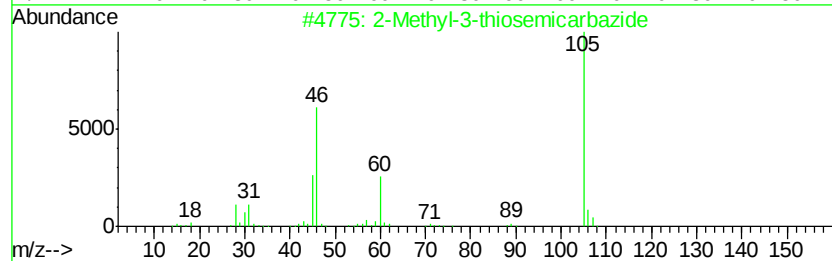
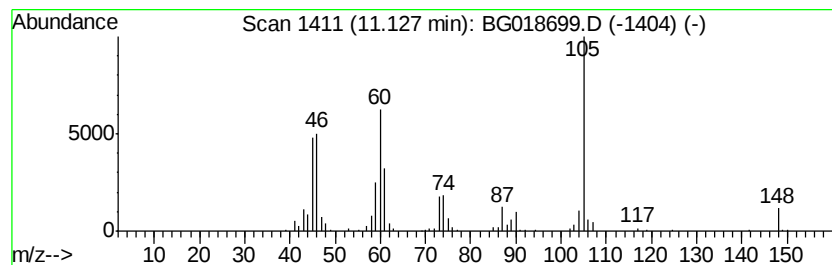
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.13	12.46 ng/ul	333120	Naphthalene-d8	10.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Methyl-3-thiosemicarbazide	105	C2H7N3S	021185-13-7	49
2		2-Methyl-1,3-oxathiolane	104	C4H8OS	017642-74-9	38
3		1,3-Oxathiolane, 2,2-dimethyl-	118	C5H10OS	005684-31-1	38
4		1,3-Dithiane-4,6-dione	148	C4H4O2S2	052133-76-3	27
5		1,2,4-Thiadiazole, 5-amino-3-(me...	147	C3H5N3S2	006913-13-9	17



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG091515\
Data File : BG018699.D
Acq On : 16 Sep 2015 1:23
Operator : UM/IZ
Sample : G3641-06DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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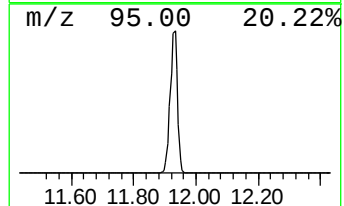
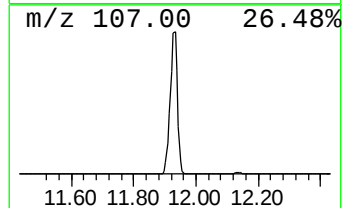
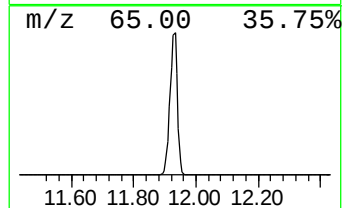
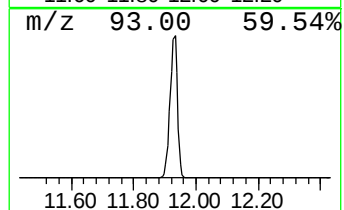
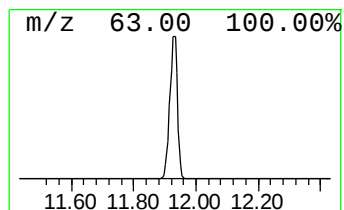
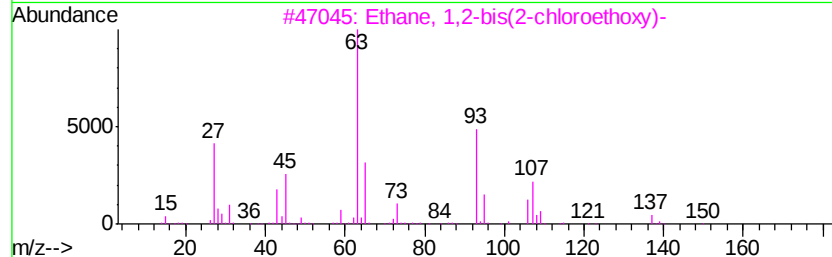
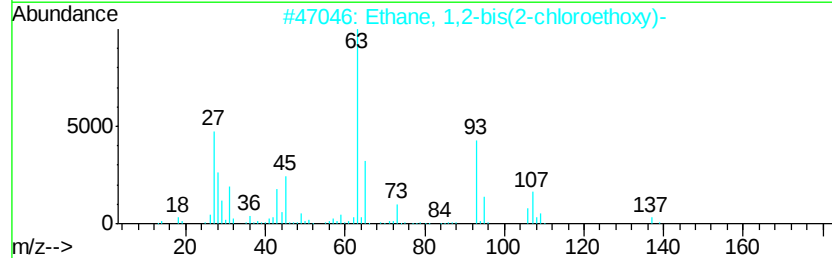
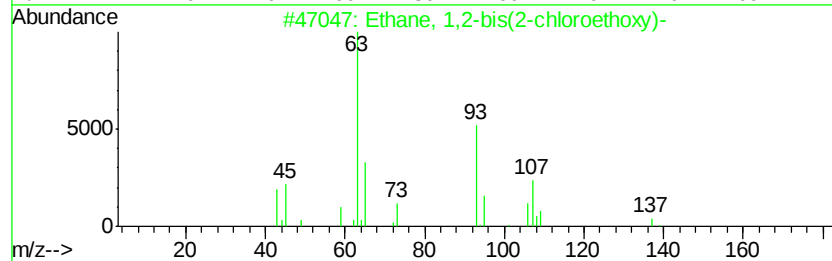
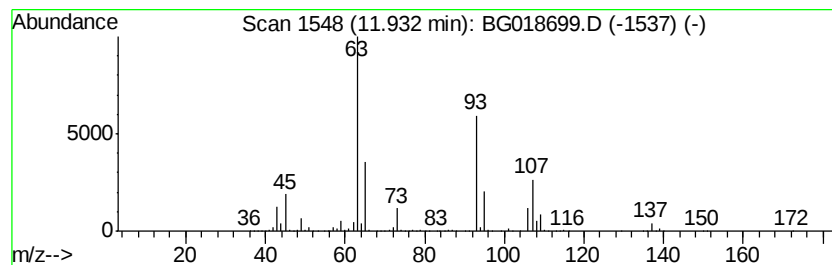
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Ethane, 1,2-bis(2-chloroeth... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	167.13 ng/ul	4469650	Naphthalene-d8	10.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethane, 1,2-bis(2-chloroethoxy)-	186	C6H12Cl2O2	000112-26-5	91
2		Ethane, 1,2-bis(2-chloroethoxy)-	186	C6H12Cl2O2	000112-26-5	90
3		Ethane, 1,2-bis(2-chloroethoxy)-	186	C6H12Cl2O2	000112-26-5	90
4		Bis(2-chloroethyl) ether	142	C4H8Cl2O	000111-44-4	72
5		.alpha.-Chloroethyl chloroformate	142	C3H4Cl2O2	050893-53-3	64



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG091515\
Data File : BG018699.D
Acq On : 16 Sep 2015 1:23
Operator : UM/IZ
Sample : G3641-06DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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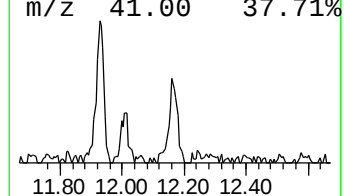
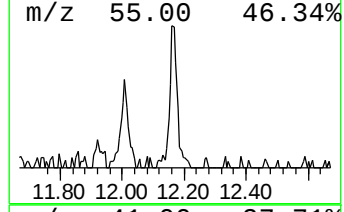
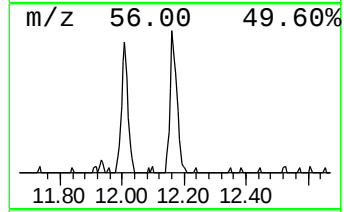
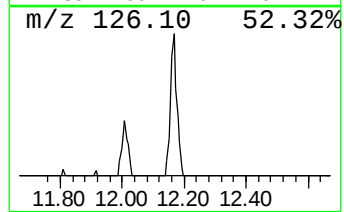
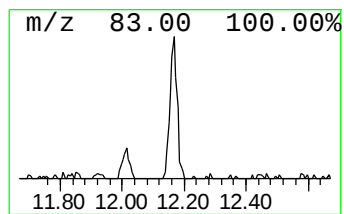
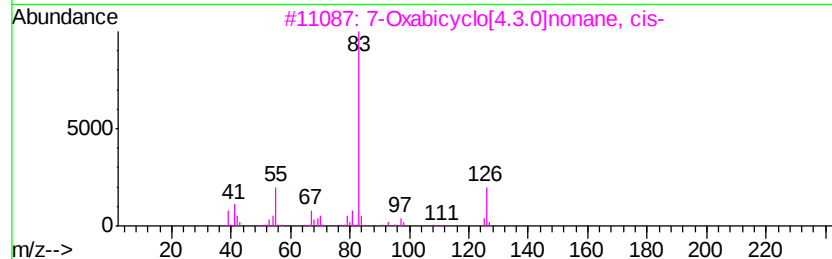
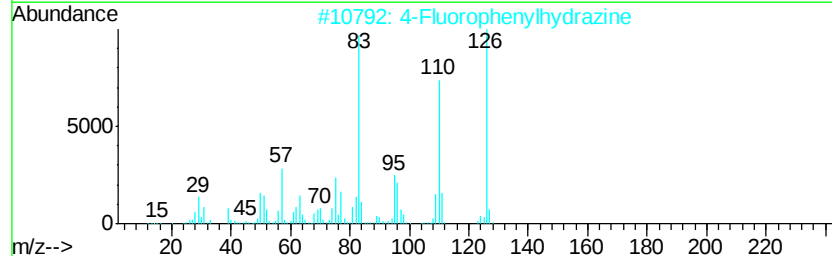
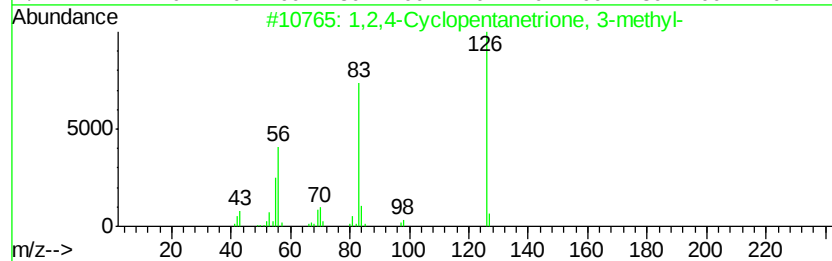
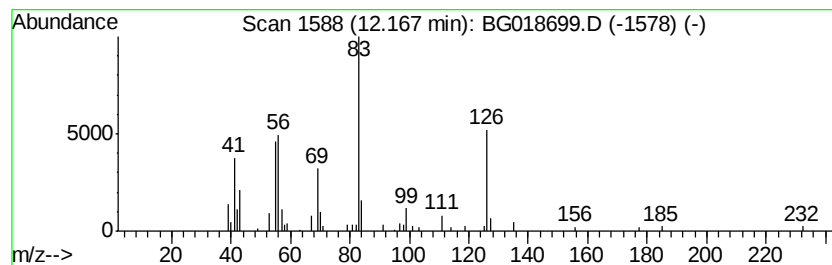
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 1,2,4-Cyclopentanetrione, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.17	3.04 ng/ul	81425	Naphthalene-d8	10.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2,4-Cyclopentanetrione, 3-methyl-	126	C6H6O3	004505-54-8	72
2		4-Fluorophenylhydrazine	126	C6H7FN2	000823-85-8	52
3		7-Oxabicyclo[4.3.0]nonane, cis-	126	C8H14O	013149-01-4	50
4		2-Fluorophenylhydrazine	126	C6H7FN2	002368-80-1	50
5		Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	40



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG091515\
Data File : BG018699.D
Acq On : 16 Sep 2015 1:23
Operator : UM/IZ
Sample : G3641-06DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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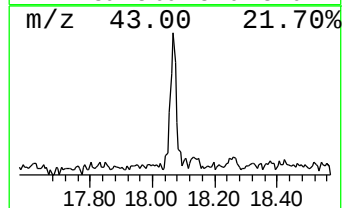
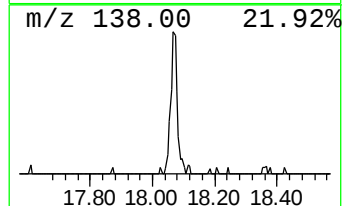
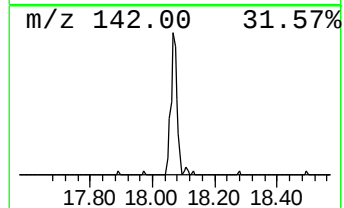
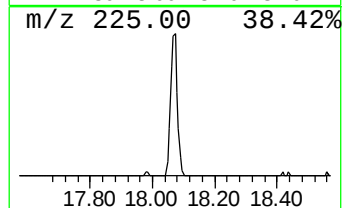
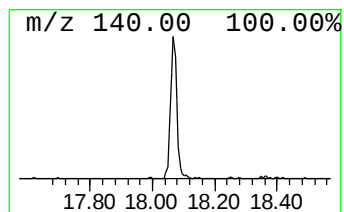
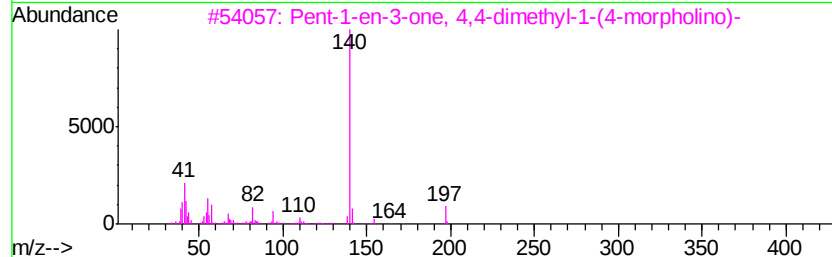
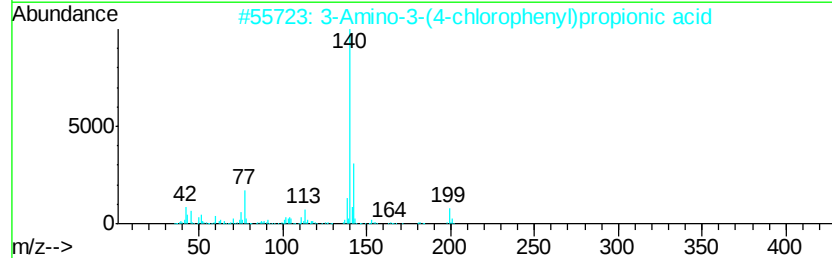
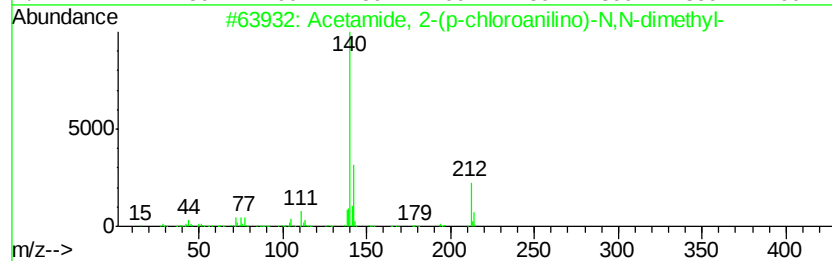
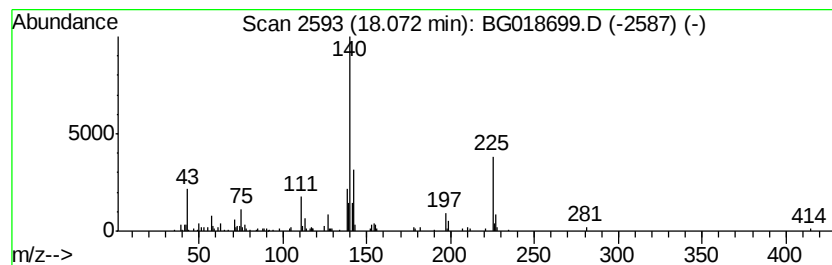
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Acetamide, 2-(p-chloroanili... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.07	2.09 ng/ul	119425	Phenanthrene-d10	17.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Acetamide, 2-(p-chloroanilino)-N...	212	C10H13ClN2O	090875-42-6	53
2		3-Amino-3-(4-chlorophenyl)propio...	199	C9H10ClN02	019947-39-8	52
3		Pent-1-en-3-one, 4,4-dimethyl-1-...	197	C11H19N02	1000259-86-0	43
4		p-Chloro-N-ethylaniline	155	C8H10ClN	013519-75-0	38
5		Ethanol, 2-(4-chlorophenyl)amino-	171	C8H10ClNO	002933-81-5	38



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG091515\
Data File : BG018699.D
Acq On : 16 Sep 2015 1:23
Operator : UM/IZ
Sample : G3641-06DL 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B0AM4DL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG091015.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.18	59.0	ng/ul	1000960	1	8.00	339595	20.0
Propane, 2-ethoxy...	5.45	7.0	ng/ul	118910	1	8.00	339595	20.0
1,4-Oxathiane	5.93	28.5	ng/ul	483299	1	8.00	339595	20.0
Cyclohexanone, 3,...	8.42	28.8	ng/ul	488711	1	8.00	339595	20.0
Cyclohexanol, 3,3...	8.62	17.7	ng/ul	300323	1	8.00	339595	20.0
unknown-01	11.13	12.5	ng/ul	333120	2	10.80	534877	20.0
Ethane, 1,2-bis(2...	11.93	167.1	ng/ul	4469650	2	10.80	534877	20.0
1,2,4-Cyclopentan...	12.17	3.0	ng/ul	81425	2	10.80	534877	20.0
Acetamide, 2-(p-c...	18.07	2.1	ng/ul	119425	4	17.37	1145160	20.0