

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021115\
 Data File : BG015981.D
 Acq On : 11 Feb 2015 11:22
 Operator : TP/IZ
 Sample : G1227-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 X1H33

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\SOM01.2-EPA-BG021015.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	2.819	19	22	29	rVB	25165	30764	0.79%	0.043%
2	2.960	41	46	54	rBV	331912	519476	13.28%	0.723%
3	3.037	54	59	69	rVB	433926	558793	14.29%	0.778%
4	3.430	124	126	133	rVB6	3928	5213	0.13%	0.007%
5	3.853	193	198	205	rVB	34720	46607	1.19%	0.065%
6	4.035	226	229	232	rBV5	3641	4774	0.12%	0.007%
7	4.740	345	349	354	rBV4	6617	8274	0.21%	0.012%
8	4.940	378	383	387	rBV5	5423	8345	0.21%	0.012%
9	5.016	390	396	401	rVV5	5221	7705	0.20%	0.011%
10	5.128	412	415	421	rVB4	7495	9982	0.26%	0.014%
11	6.979	721	730	733	rBV	438158	777280	19.87%	1.082%
12	7.155	754	760	769	rBV	359100	528891	13.52%	0.736%
13	7.349	786	793	804	rBV	392755	612388	15.66%	0.852%
14	7.448	804	810	816	rVB2	17777	26787	0.68%	0.037%
15	7.819	865	873	881	rBV	265478	412540	10.55%	0.574%
16	8.042	905	911	917	rBV3	13916	22698	0.58%	0.032%
17	8.512	982	991	996	rBV	540239	870911	22.27%	1.212%
18	8.570	996	1001	1008	rVB	747657	1149959	29.40%	1.600%
19	8.970	1062	1069	1073	rBV	370970	624266	15.96%	0.869%
20	9.017	1073	1077	1086	rVV	231443	368014	9.41%	0.512%
21	9.087	1086	1089	1093	rVB5	3475	5170	0.13%	0.007%
22	9.692	1184	1192	1199	rBV	293926	482039	12.32%	0.671%
23	9.781	1199	1207	1220	rVB	779523	1281010	32.75%	1.782%
24	10.033	1245	1250	1256	rVB5	3363	5789	0.15%	0.008%
25	10.221	1275	1282	1292	rBV	491723	800877	20.48%	1.114%
26	10.609	1341	1348	1351	rBV	418311	683145	17.47%	0.951%
27	10.656	1351	1356	1364	rVV	1300939	2182568	55.80%	3.037%
28	10.738	1364	1370	1379	rVB2	538278	933419	23.86%	1.299%
29	11.637	1514	1523	1534	rBV	1179765	2025994	51.80%	2.819%
30	11.737	1534	1540	1550	rVB2	21404	54350	1.39%	0.076%
31	12.136	1605	1608	1612	rBV3	3181	5396	0.14%	0.008%
32	12.195	1612	1618	1625	rVV2	10739	19057	0.49%	0.027%
33	12.377	1643	1649	1660	rBV2	24403	55387	1.42%	0.077%
34	12.547	1671	1678	1684	rBV2	6139	10418	0.27%	0.014%

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35	12.859	1721	1731	1734	rBV	408869	786697	20.11%	1.095%
36	12.923	1735	1742	1755	rVB2	450721	1524028	38.96%	2.121%
37	13.282	1798	1803	1807	rVB4	6053	9085	0.23%	0.013%
38	13.857	1895	1901	1906	rBV	988319	1310555	33.51%	1.824%
39	13.904	1906	1909	1913	rVB	13238	16557	0.42%	0.023%
40	14.139	1942	1949	1956	rBV	1064550	1618019	41.37%	2.251%
41	14.445	1992	2001	2008	rBV2	650687	972667	24.87%	1.353%
42	14.639	2027	2034	2042	rBV	462040	734260	18.77%	1.022%
43	14.809	2056	2063	2083	rBV	169424	329561	8.43%	0.459%
44	15.438	2163	2170	2177	rBV	1402279	2043600	52.25%	2.844%
45	15.543	2182	2188	2198	rBV	387115	530280	13.56%	0.738%
46	15.702	2206	2215	2222	rBV	1335550	1847016	47.22%	2.570%
47	15.831	2231	2237	2243	rBV	1106639	1349644	34.51%	1.878%
48	16.084	2275	2280	2285	rBV2	18008	24967	0.64%	0.035%
49	16.342	2319	2324	2325	rBV2	9507	12099	0.31%	0.017%
50	16.383	2325	2331	2337	rVV	1796949	2440459	62.39%	3.396%
51	16.466	2339	2345	2351	rBV	117841	162775	4.16%	0.226%
52	16.595	2363	2367	2371	rBV6	3051	5386	0.14%	0.007%
53	16.654	2371	2377	2382	rVV	1736419	2269550	58.02%	3.158%
54	16.689	2382	2383	2393	rVB6	14363	14916	0.38%	0.021%
55	16.830	2401	2407	2419	rBV	1232733	1767410	45.19%	2.459%
56	17.006	2429	2437	2446	rVB	2450590	3542724	90.58%	4.929%
57	17.147	2455	2461	2464	rBV	1408401	1713093	43.80%	2.384%
58	17.182	2464	2467	2477	rVV2	829271	1153356	29.49%	1.605%
59	17.282	2477	2484	2497	rVB2	1506836	2239302	57.25%	3.116%
60	17.558	2527	2531	2536	rBV6	3116	5343	0.14%	0.007%
61	17.746	2560	2563	2566	rBV5	6453	7836	0.20%	0.011%
62	17.811	2566	2574	2581	rVV	1742943	2279559	58.28%	3.172%
63	18.157	2627	2633	2639	rBV	1720235	2246251	57.43%	3.126%
64	18.410	2672	2676	2683	rVB2	48727	71811	1.84%	0.100%
65	18.563	2698	2702	2705	rVB3	3940	5613	0.14%	0.008%
66	18.616	2707	2711	2715	rVB4	3471	5217	0.13%	0.007%
67	18.768	2732	2737	2741	rBV	22034	27570	0.70%	0.038%
68	19.209	2806	2812	2816	rBV2	22263	30039	0.77%	0.042%
69	19.274	2819	2823	2824	rBV4	3237	4308	0.11%	0.006%
70	19.315	2824	2830	2836	rVV	80315	114774	2.93%	0.160%
71	19.374	2836	2840	2844	rVB5	3454	6077	0.16%	0.008%

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72	19.421	2844	2848	2851	rBV4	4918	7335	0.19%	0.010%
73	19.462	2851	2855	2859	rVB	14492	17925	0.46%	0.025%
74	19.573	2867	2874	2881	rBV	1945761	2561963	65.50%	3.565%
75	19.714	2895	2898	2901	rVB5	4536	4877	0.12%	0.007%
76	19.808	2910	2914	2919	rBV2	26410	37448	0.96%	0.052%
77	19.937	2931	2936	2938	rBV5	2712	4835	0.12%	0.007%
78	20.149	2961	2972	2981	rVB	3476174	3911368	100.00%	5.442%
79	20.507	3028	3033	3039	rBV	2269434	2473393	63.24%	3.442%
80	20.918	3098	3103	3107	rBV7	3632	6752	0.17%	0.009%
81	21.024	3119	3121	3126	rVB5	3020	4174	0.11%	0.006%
82	21.077	3126	3130	3135	rBV5	6386	9466	0.24%	0.013%
83	21.153	3140	3143	3147	rVV3	11032	14083	0.36%	0.020%
84	21.283	3160	3165	3172	rBV	3058030	3349869	85.64%	4.661%
85	21.353	3172	3177	3185	rVB	1079888	1367106	34.95%	1.902%
86	22.181	3312	3318	3336	rVB	2002263	2688351	68.73%	3.741%
87	22.446	3361	3363	3368	rVB4	9086	11822	0.30%	0.016%
88	22.898	3433	3440	3448	rVB7	15588	44250	1.13%	0.062%
89	22.980	3451	3454	3459	rVB4	9078	14129	0.36%	0.020%
90	23.139	3477	3481	3485	rBV6	11998	18738	0.48%	0.026%
91	23.233	3492	3497	3504	rBV3	45610	70067	1.79%	0.097%
92	23.515	3537	3545	3554	rBV2	1323240	2497030	63.84%	3.474%
93	23.662	3562	3570	3584	rVB	677558	1301643	33.28%	1.811%
94	24.549	3717	3721	3723	rBV5	7725	10706	0.27%	0.015%
95	26.752	4082	4096	4116	rBV	946704	3080498	78.76%	4.286%

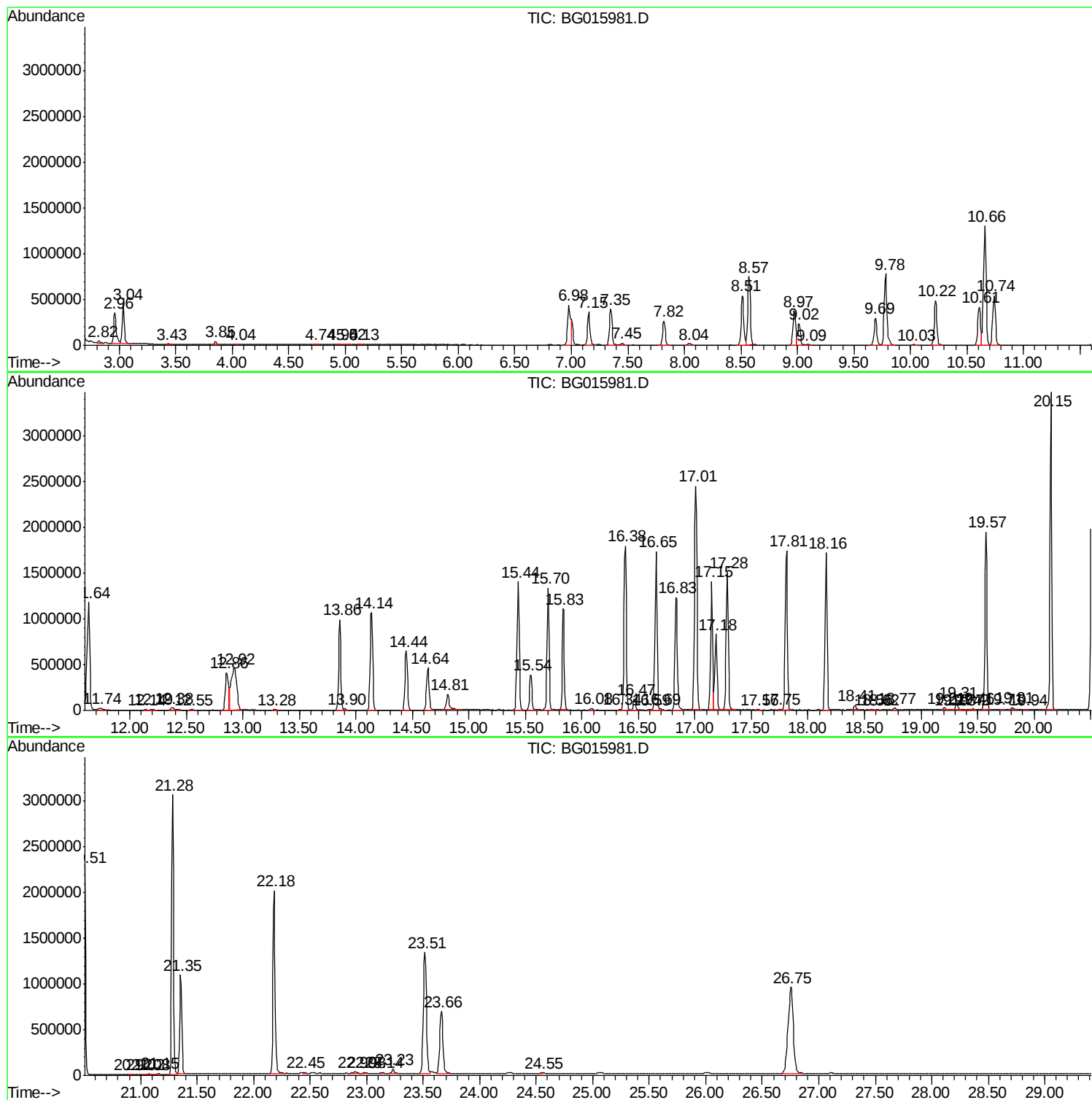
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG021015.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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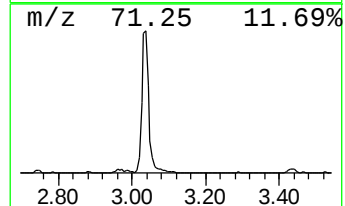
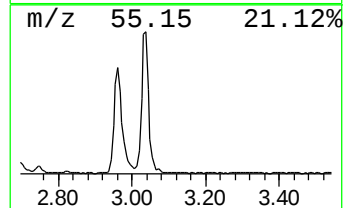
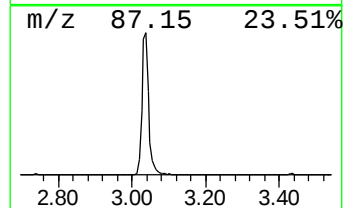
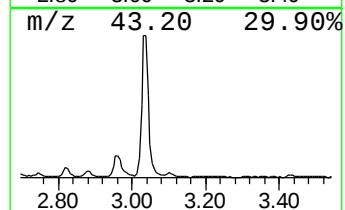
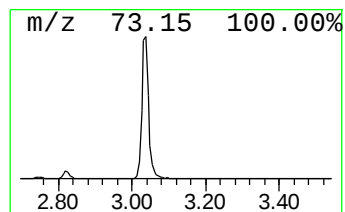
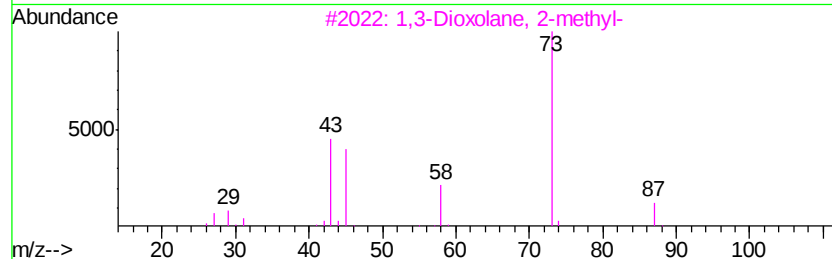
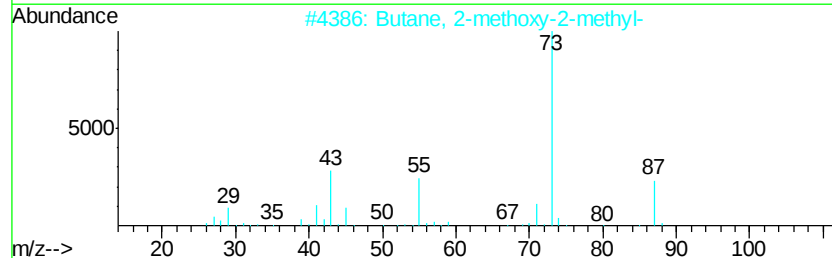
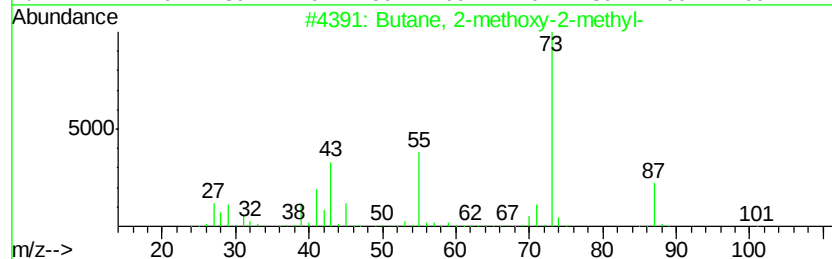
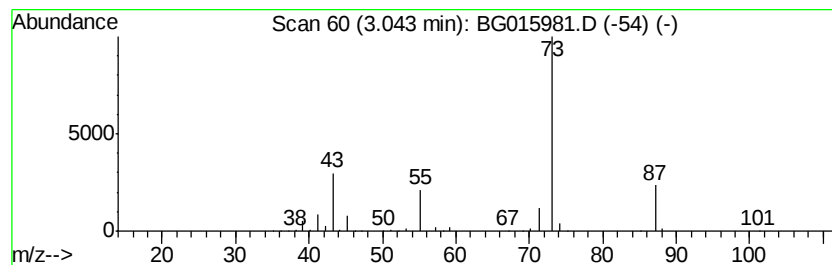
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG021015.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.04	27.09 ng/ul	558793	1,4-Dichlorobenzene-d4	7.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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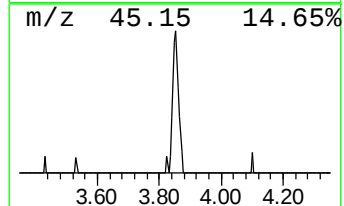
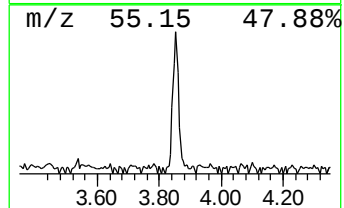
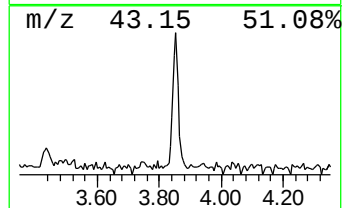
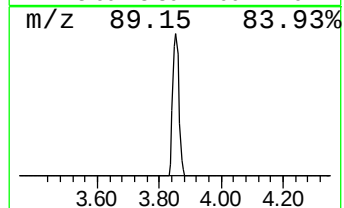
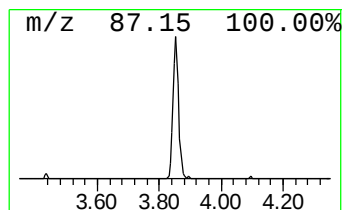
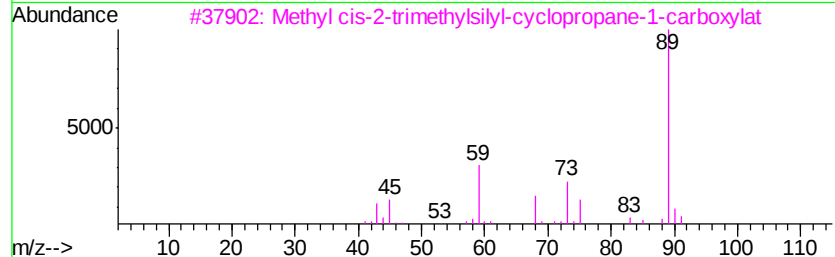
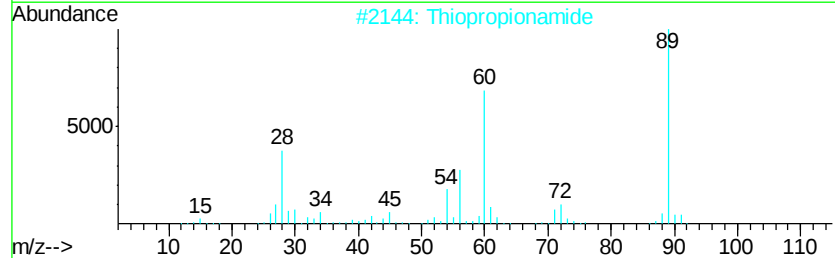
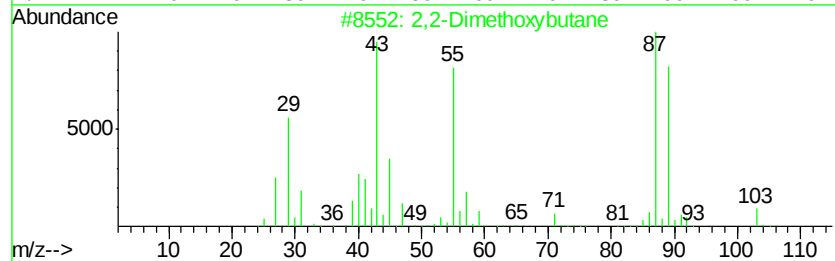
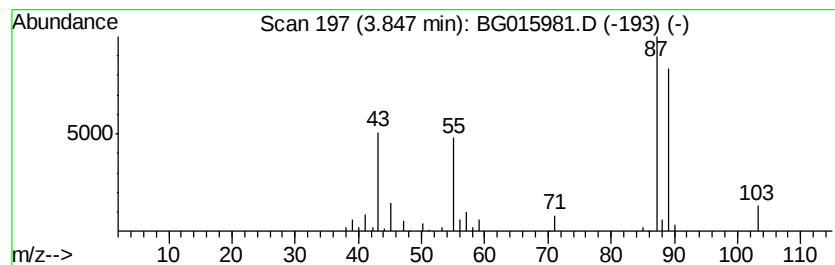
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 2,2-Dimethoxybutane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.85	2.26 ng/ul	46607	1,4-Dichlorobenzene-d4	7.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,2-Dimethoxybutane	118	C6H14O2	003453-99-4	74
2			Thiopropionamide	89	C3H7NS	000631-58-3	53
3			Methyl cis-2-trimethylsilyl-cycl...	172	C8H16O2Si	1000144-79-0	45
4			Thiazole, tetrahydro-	89	C3H7NS	000504-78-9	9
5			1,3-Dioxane	88	C4H8O2	000505-22-6	9



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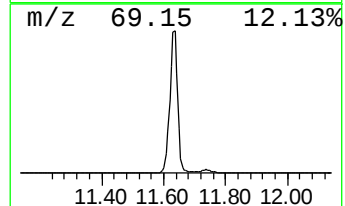
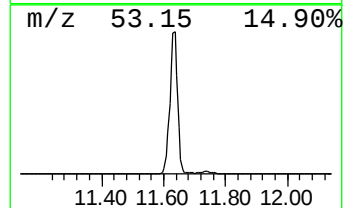
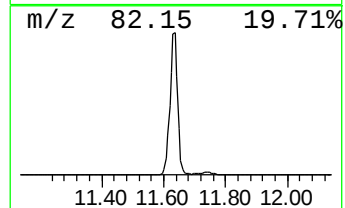
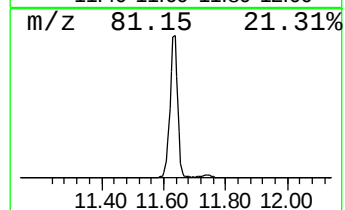
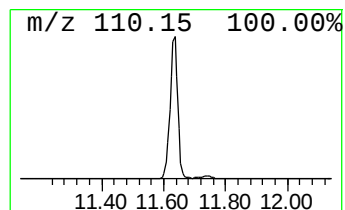
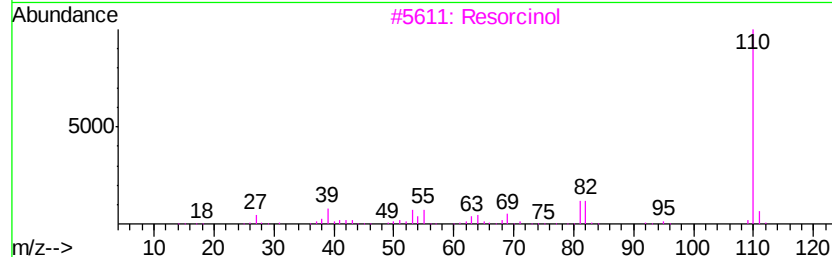
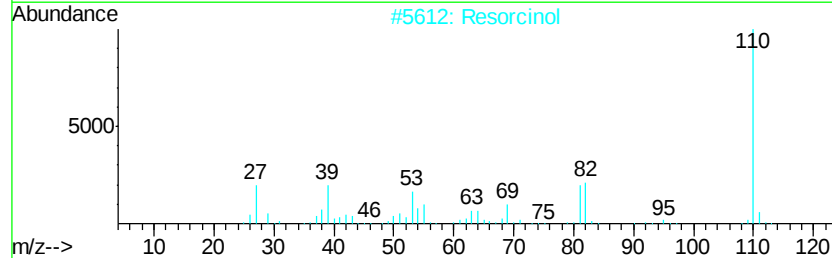
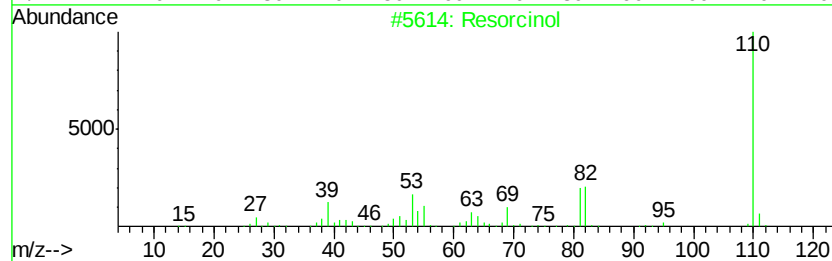
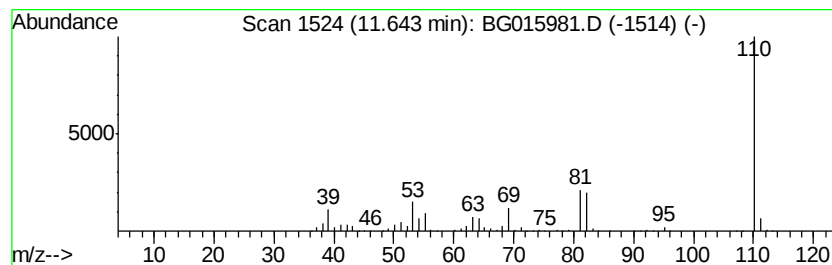
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Peak Number 4 Resorcinol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.64	59.31 ng/ul	2025990	Napthalene-d8	10.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Resorcinol	110	C6H6O2	000108-46-3	97
2		Resorcinol	110	C6H6O2	000108-46-3	94
3		Resorcinol	110	C6H6O2	000108-46-3	91
4		Resorcinol	110	C6H6O2	000108-46-3	91
5		Hydroquinone	110	C6H6O2	000123-31-9	64



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ClientSampleId :
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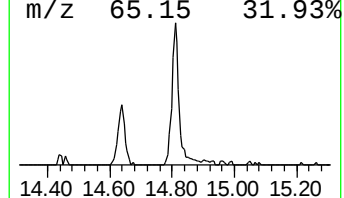
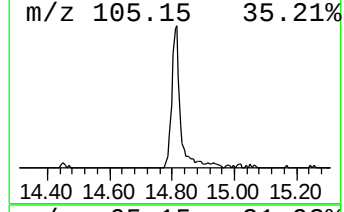
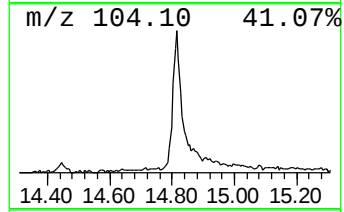
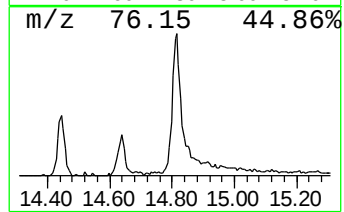
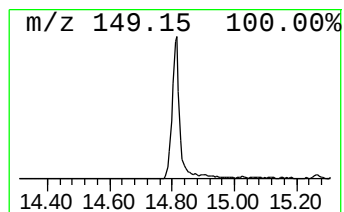
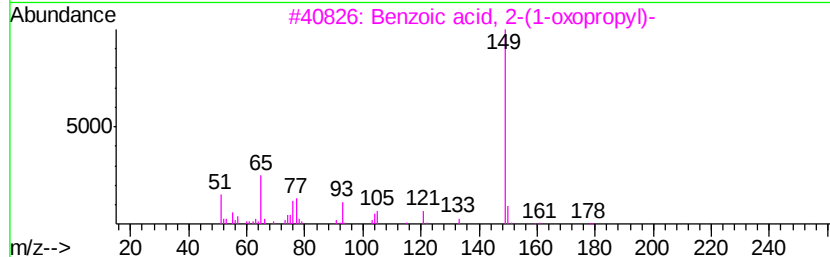
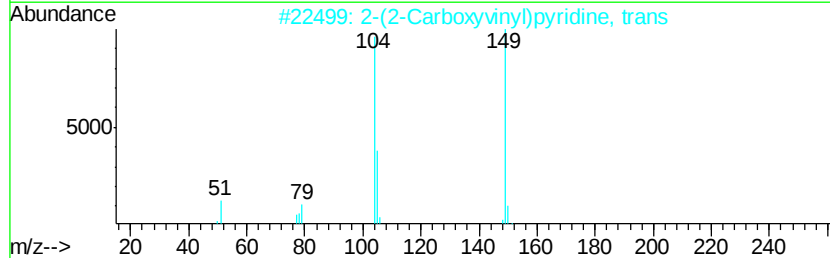
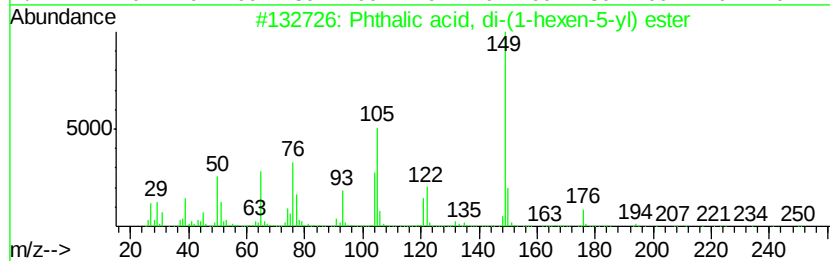
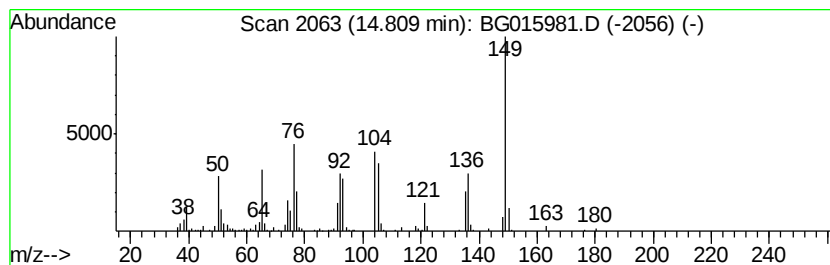
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Phthalic acid, di-(1-hexen-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.81	6.78 ng/ul	329561	Acenaphthene-d10	14.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phthalic acid, di-(1-hexen-5-yl)...	330	C20H26O4	1000160-80-2	50
2		2-(2-Carboxyvinyl)pyridine, trans	149	C8H7NO2	007340-22-9	38
3		Benzoic acid, 2-(1-oxopropyl)-	178	C10H10O3	002360-45-4	38
4		Benzoic acid, 2-(1-oxopropyl)-	178	C10H10O3	002360-45-4	37
5		Benzofuran-2-one, 3-methyl-3-aza...	149	C8H7NO2	1000289-12-2	37



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021115\
Data File : BG015981.D
Acq On : 11 Feb 2015 11:22
Operator : TP/IZ
Sample : G1227-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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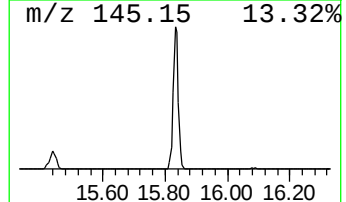
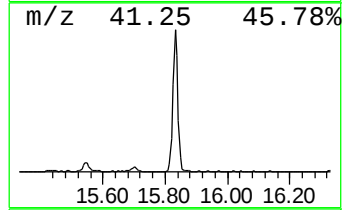
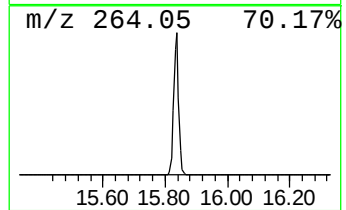
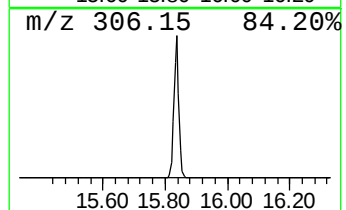
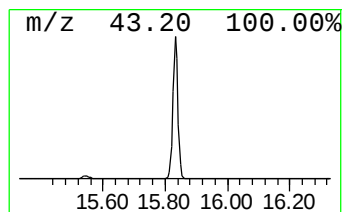
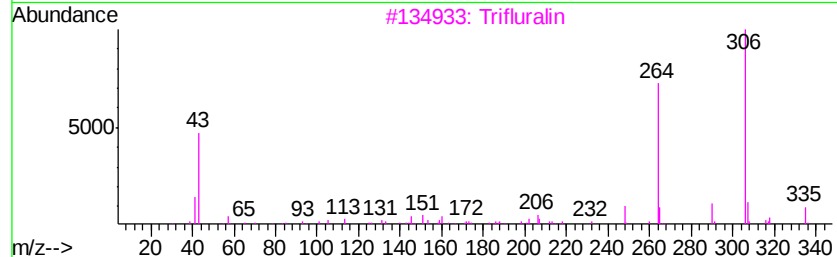
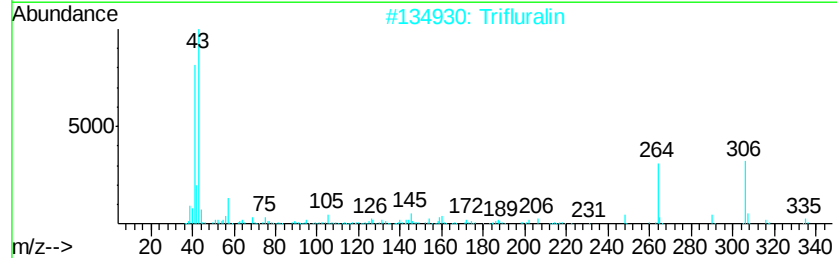
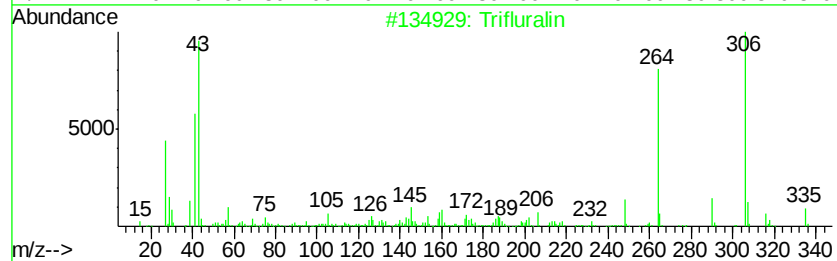
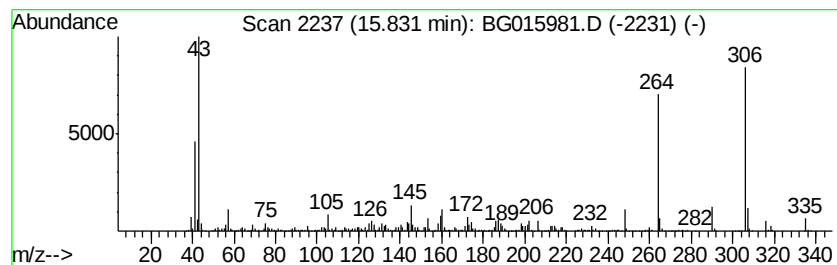
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Trifluralin Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.83	23.40 ng/ul	1349640	Phenanthrene-d10	17.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trifluralin	335	C13H16F3N3O4	001582-09-8	97
2		Trifluralin	335	C13H16F3N3O4	001582-09-8	95
3		Trifluralin	335	C13H16F3N3O4	001582-09-8	95
4		Trifluralin	335	C13H16F3N3O4	001582-09-8	70
5		Azuleno[6,5-b]furan-2(3H)-one, 6...	366	C19H26O7	051292-55-8	59



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021115\
Data File : BG015981.D
Acq On : 11 Feb 2015 11:22
Operator : TP/IZ
Sample : G1227-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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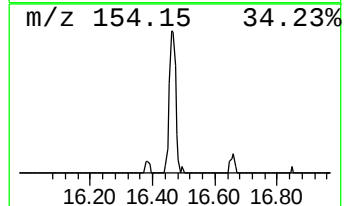
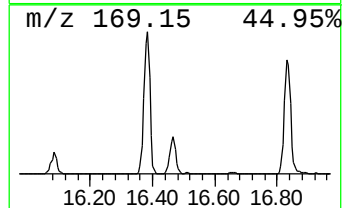
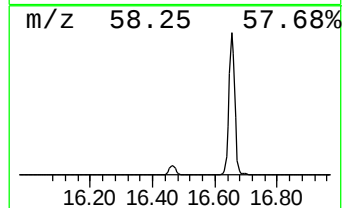
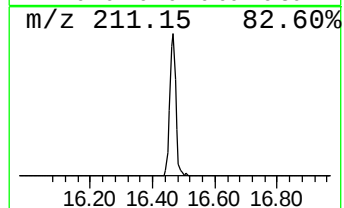
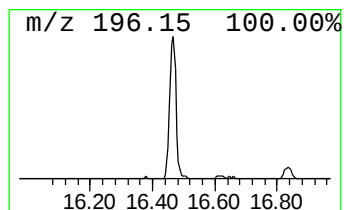
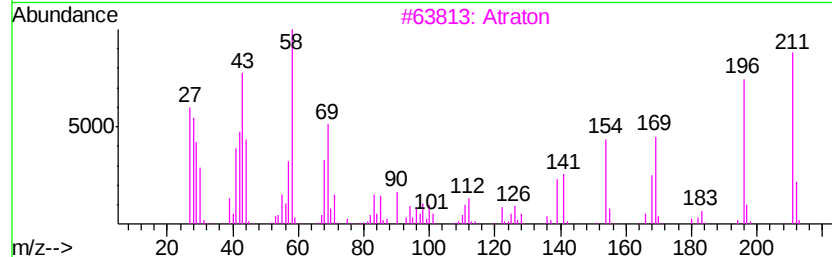
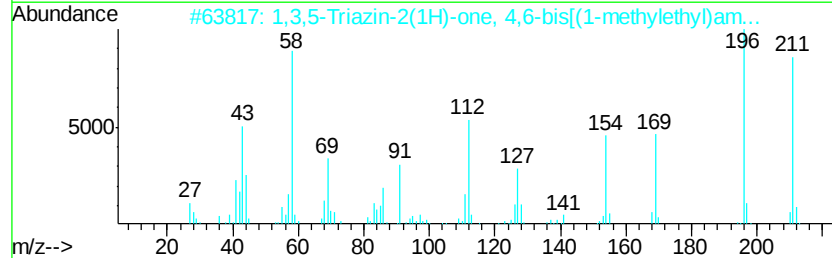
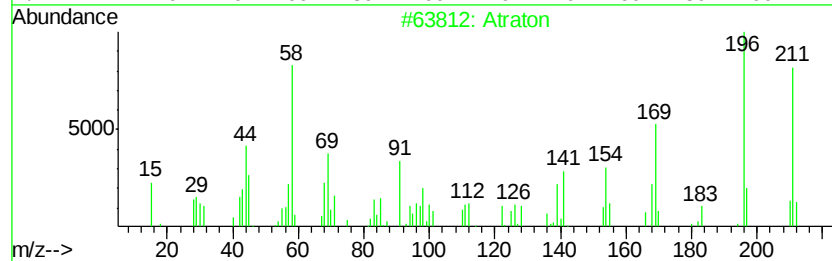
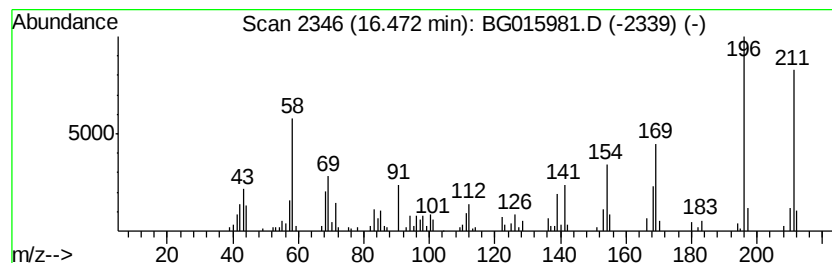
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Atraton Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.47	2.82 ng/ul	162775	Phenanthrene-d10	17.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Atraton	211	C9H17N5O	001610-17-9	91
2		1,3,5-Triazin-2(1H)-one, 4,6-bis...	211	C9H17N5O	007374-53-0	89
3		Atraton	211	C9H17N5O	001610-17-9	87
4		3,6-Diamino-9-methylcarbazole	211	C13H13N3	046498-17-3	55
5		2-Naphthaleneacetonitrile, 6-met...	211	C14H13NO	086603-94-3	50



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021115\
Data File : BG015981.D
Acq On : 11 Feb 2015 11:22
Operator : TP/IZ
Sample : G1227-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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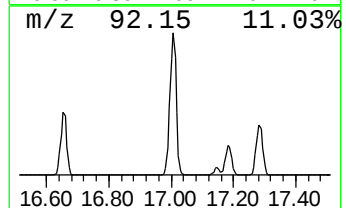
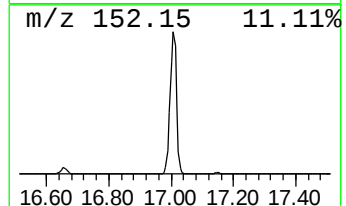
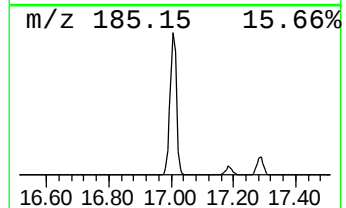
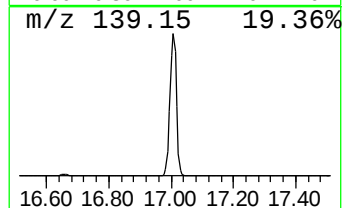
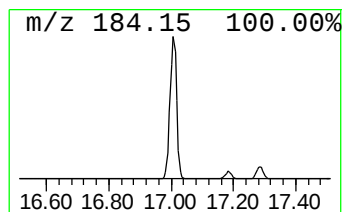
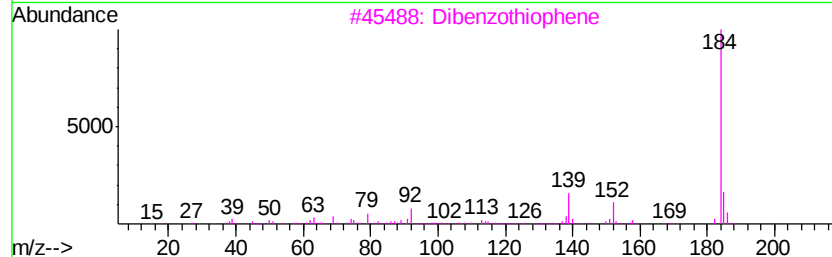
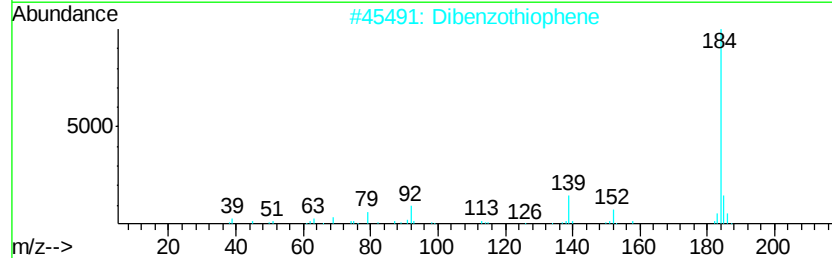
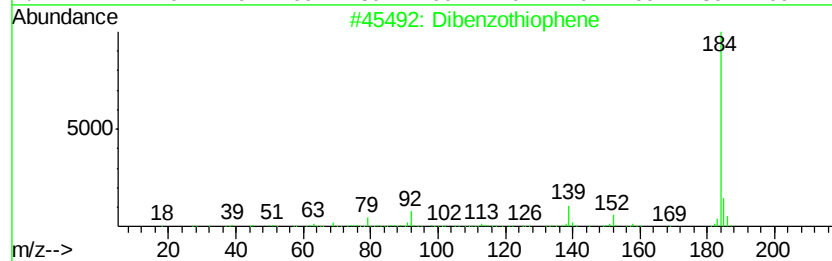
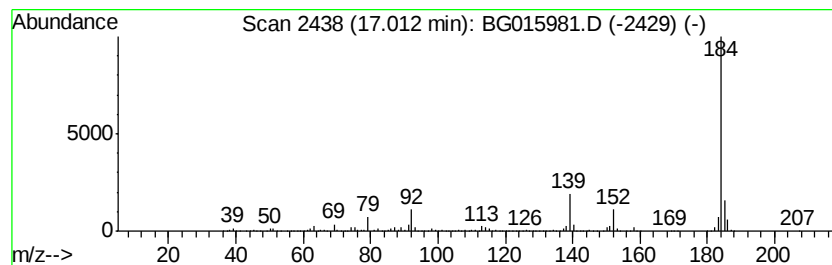
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Dibenzothiophene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.01	61.43 ng/ul	3542720	Phenanthrene-d10	17.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dibenzothiophene	184	C12H8S	000132-65-0	97
2		Dibenzothiophene	184	C12H8S	000132-65-0	96
3		Dibenzothiophene	184	C12H8S	000132-65-0	96
4		Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	95
5		Dibenzothiophene	184	C12H8S	000132-65-0	95



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021115\
Data File : BG015981.D
Acq On : 11 Feb 2015 11:22
Operator : TP/IZ
Sample : G1227-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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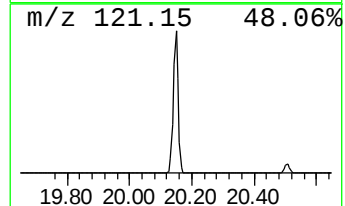
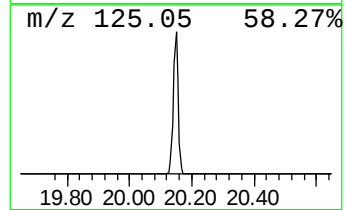
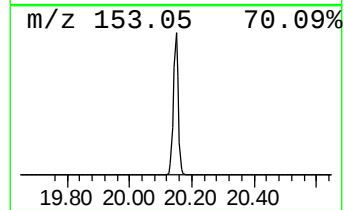
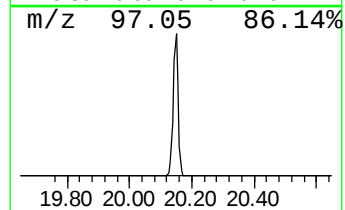
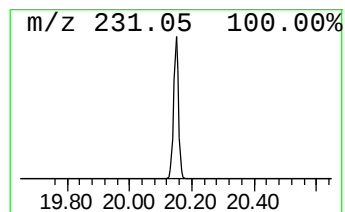
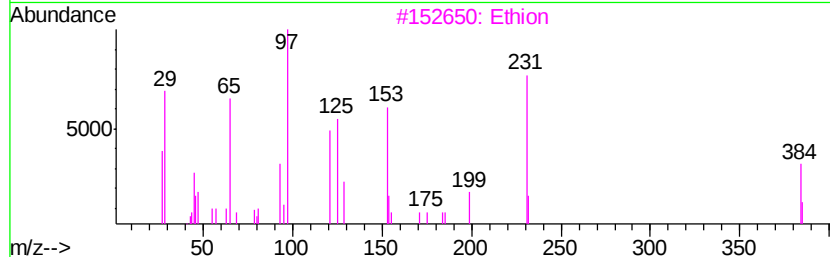
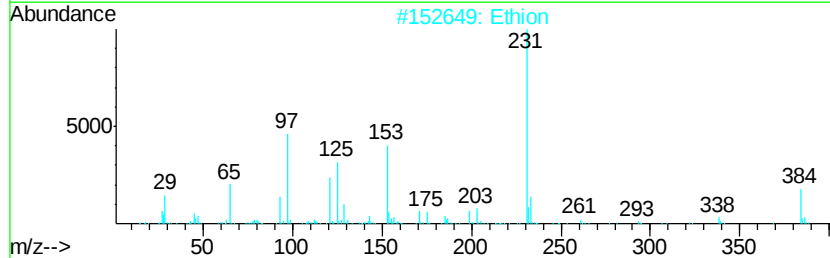
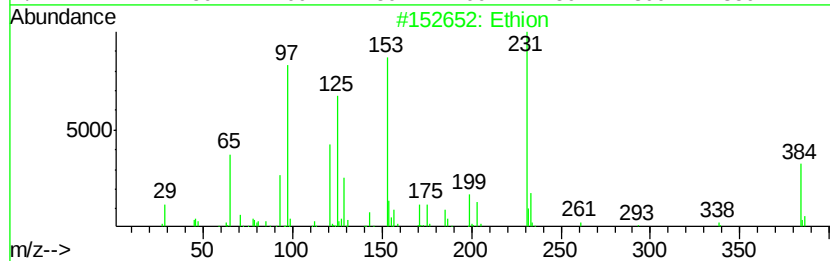
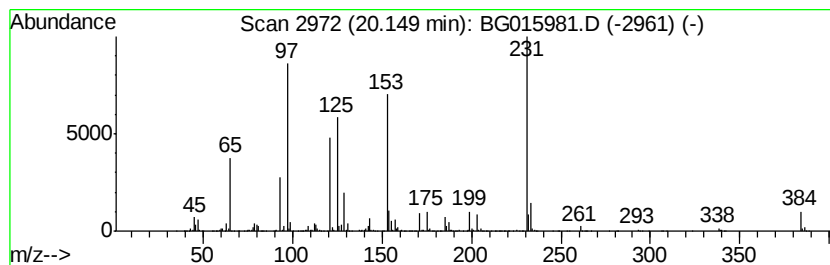
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Ethion Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.15	57.22 ng/ul	3911370	Chrysene-d12	21.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethion	384	C9H22O4P2S4	000563-12-2	98
2		Ethion	384	C9H22O4P2S4	000563-12-2	96
3		Ethion	384	C9H22O4P2S4	000563-12-2	87
4		Ethion	384	C9H22O4P2S4	000563-12-2	58
5		Methoxycarbonyl dimethoxythiopho...	276	C6H13O4PS3	1000251-32-8	43



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021115\
Data File : BG015981.D
Acq On : 11 Feb 2015 11:22
Operator : TP/IZ
Sample : G1227-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG021015.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...	3.04	27.1	ng/ul	558793	1	7.82	412540	20.0
2,2-Dimethoxybutane	3.85	2.3	ng/ul	46607	1	7.82	412540	20.0
Resorcinol	11.64	59.3	ng/ul	2025990	2	10.61	683145	20.0
Phthalic acid, di...	14.81	6.8	ng/ul	329561	3	14.44	972667	20.0
Trifluralin	15.83	23.4	ng/ul	1349640	4	17.18	1153360	20.0
Atraton	16.47	2.8	ng/ul	162775	4	17.18	1153360	20.0
Dibenzothiophene	17.01	61.4	ng/ul	3542720	4	17.18	1153360	20.0
Ethion	20.15	57.2	ng/ul	3911370	5	21.35	1367110	20.0