

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092115\
 Data File : BG018791.D
 Acq On : 21 Sep 2015 20:33
 Operator : TP
 Sample : G3696-06
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AQ8

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	2.886	4	8	16	rVB3	12369	17388	0.96%	0.099%
2	3.103	41	45	52	rBV	62627	101344	5.59%	0.574%
3	3.180	52	58	73	rVB	978264	1406638	77.60%	7.969%
4	3.438	98	102	104	rBV4	2388	3381	0.19%	0.019%
5	4.085	205	212	215	rBV6	2108	3392	0.19%	0.019%
6	5.113	381	387	390	rVB4	1957	3368	0.19%	0.019%
7	6.012	535	540	544	rBV4	1484	3037	0.17%	0.017%
8	7.163	731	736	742	rBV	34536	55650	3.07%	0.315%
9	7.322	756	763	774	rVB	140203	232471	12.82%	1.317%
10	7.534	792	799	807	rVB	130048	221594	12.22%	1.255%
11	7.998	872	878	887	rVV	140322	232689	12.84%	1.318%
12	8.703	992	998	1007	rVV	103402	180254	9.94%	1.021%
13	9.108	1063	1067	1069	rBV4	3923	5133	0.28%	0.029%
14	9.155	1069	1075	1083	rVV	153333	271189	14.96%	1.536%
15	9.525	1133	1138	1142	rVB3	4587	7244	0.40%	0.041%
16	9.608	1144	1152	1160	rVB2	197046	334384	18.45%	1.894%
17	9.884	1192	1199	1205	rBV	125197	226310	12.48%	1.282%
18	9.937	1205	1208	1218	rVB3	26026	42946	2.37%	0.243%
19	10.031	1220	1224	1229	rBV	10398	14075	0.78%	0.080%
20	10.424	1284	1291	1301	rBV	198361	357516	19.72%	2.026%
21	10.724	1336	1342	1348	rVB	21503	36803	2.03%	0.209%
22	10.800	1348	1355	1364	rBV	201112	359860	19.85%	2.039%
23	10.935	1373	1378	1387	rBV	15113	28678	1.58%	0.162%
24	11.353	1441	1449	1455	rBV2	86891	162364	8.96%	0.920%
25	11.452	1457	1466	1472	rVV2	39666	64813	3.58%	0.367%
26	11.611	1489	1493	1498	rBV4	8117	12277	0.68%	0.070%
27	11.840	1525	1532	1540	rBV8	5746	11995	0.66%	0.068%
28	11.958	1548	1552	1558	rVB2	10026	15552	0.86%	0.088%
29	12.052	1563	1568	1572	rVV4	3347	5915	0.33%	0.034%
30	12.193	1586	1592	1596	rBV4	1989	3764	0.21%	0.021%
31	12.246	1596	1601	1606	rBV5	7411	11278	0.62%	0.064%
32	12.469	1636	1639	1642	rBV3	2859	3685	0.20%	0.021%
33	12.516	1642	1647	1652	rVV5	8201	15722	0.87%	0.089%
34	12.604	1657	1662	1663	rVV2	5648	9031	0.50%	0.051%

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 Title : SVOA CALIBRATION

35	12.639	1663	1668	1676	rVV4	9341	17851	0.98%	0.101%
36	12.733	1680	1684	1687	rVV2	4921	6677	0.37%	0.038%
37	12.892	1704	1711	1717	rBV5	4404	9439	0.52%	0.053%
38	12.986	1720	1727	1732	rVB5	5061	10012	0.55%	0.057%
39	13.438	1796	1804	1812	rVB	367459	589441	32.52%	3.339%
40	13.568	1819	1826	1830	rVV3	10983	16819	0.93%	0.095%
41	13.632	1830	1837	1844	rVB	205899	333930	18.42%	1.892%
42	13.809	1860	1867	1873	rVB2	6639	11293	0.62%	0.064%
43	13.938	1884	1889	1891	rBV3	2822	3524	0.19%	0.020%
44	14.026	1898	1904	1910	rBV	497025	726785	40.09%	4.118%
45	14.073	1910	1912	1916	rVB3	9066	10727	0.59%	0.061%
46	14.320	1947	1954	1960	rVV	527209	839126	46.29%	4.754%
47	14.378	1960	1964	1969	rVB2	55704	80950	4.47%	0.459%
48	14.437	1969	1974	1979	rBV5	1875	3834	0.21%	0.022%
49	14.519	1982	1988	1993	rVB	17438	30260	1.67%	0.171%
50	14.625	1993	2006	2013	rBV3	378272	654339	36.10%	3.707%
51	14.772	2027	2031	2035	rVB5	2273	3443	0.19%	0.020%
52	14.831	2035	2041	2051	rBV2	34133	70546	3.89%	0.400%
53	15.542	2157	2162	2164	rBV2	2086	3126	0.17%	0.018%
54	15.618	2168	2175	2185	rBV	727768	1136826	62.71%	6.441%
55	15.730	2188	2194	2202	rBV	194214	285958	15.78%	1.620%
56	15.853	2210	2215	2221	rVB3	7519	11866	0.65%	0.067%
57	16.576	2331	2338	2340	rBV4	2514	4365	0.24%	0.025%
58	17.152	2431	2436	2442	rVB	3861	5373	0.30%	0.030%
59	17.369	2466	2473	2483	rBV2	583549	863664	47.64%	4.893%
60	17.469	2483	2490	2497	rBV	917872	1354615	74.73%	7.675%
61	18.333	2632	2637	2641	rBV4	5190	8579	0.47%	0.049%
62	18.750	2703	2708	2710	rBV5	2164	3310	0.18%	0.019%
63	18.967	2742	2745	2748	rBV5	1938	3144	0.17%	0.018%
64	19.085	2762	2765	2767	rBV2	3151	2995	0.17%	0.017%
65	19.314	2800	2804	2807	rBV4	2963	4979	0.27%	0.028%
66	19.637	2854	2859	2865	rBV2	11898	19280	1.06%	0.109%
67	19.749	2872	2878	2888	rBV	1255855	1812721	100.00%	10.270%
68	19.978	2911	2917	2921	rVB8	5360	7572	0.42%	0.043%
69	20.266	2965	2966	2969	rBV3	3050	2609	0.14%	0.015%
70	20.295	2969	2971	2973	rVV3	2684	2654	0.15%	0.015%
71	20.336	2976	2978	2982	rVV5	3030	4299	0.24%	0.024%

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72	21.347	3148	3150	3155	rVB5	4548	6621	0.37%	0.038%
73	21.623	3190	3197	3210	rBV	707846	1166024	64.32%	6.606%
74	21.858	3235	3237	3239	rVB3	3547	3061	0.17%	0.017%
75	21.876	3239	3240	3244	rBV4	5467	5432	0.30%	0.031%
76	22.052	3269	3270	3272	rVB2	6826	3674	0.20%	0.021%
77	22.105	3277	3279	3280	rBV2	3721	2604	0.14%	0.015%
78	22.522	3348	3350	3351	rBV2	3739	2643	0.15%	0.015%
79	23.092	3444	3447	3449	rBV4	5531	6349	0.35%	0.036%
80	23.262	3474	3476	3479	rBV3	5671	5470	0.30%	0.031%
81	23.697	3547	3550	3561	rVB3	10324	27638	1.52%	0.157%
82	23.867	3578	3579	3580	rBV	5333	3545	0.20%	0.020%
83	24.314	3653	3655	3657	rBV3	3338	2554	0.14%	0.014%
84	24.596	3692	3703	3715	rBV	655044	1774203	97.88%	10.052%
85	24.813	3729	3740	3753	rBV2	408853	1160236	64.01%	6.573%
86	24.995	3769	3771	3775	rBV5	3504	5115	0.28%	0.029%
87	25.783	3903	3905	3908	rBV4	3846	4420	0.24%	0.025%
88	25.924	3927	3929	3935	rVB7	5876	9312	0.51%	0.053%
89	26.676	4055	4057	4059	rBV3	3849	2877	0.16%	0.016%
90	26.928	4098	4100	4103	rBV4	4520	3364	0.19%	0.019%
91	27.099	4127	4129	4133	rBV4	3939	4395	0.24%	0.025%
92	27.216	4147	4149	4150	rBV2	4587	2929	0.16%	0.017%
93	28.039	4287	4289	4290	rBV2	3314	2671	0.15%	0.015%
94	28.092	4296	4298	4300	rBV3	3068	2627	0.14%	0.015%
95	29.126	4473	4474	4477	rBV3	4403	3603	0.20%	0.020%
96	29.514	4537	4540	4543	rVB5	4830	6330	0.35%	0.036%
97	29.901	4604	4606	4607	rBV2	3728	2562	0.14%	0.015%
98	30.371	4684	4686	4689	rVB4	5345	4275	0.24%	0.024%
99	30.418	4692	4694	4697	rVB4	4227	4551	0.25%	0.026%
100	31.394	4858	4860	4861	rBV	3730	2925	0.16%	0.017%

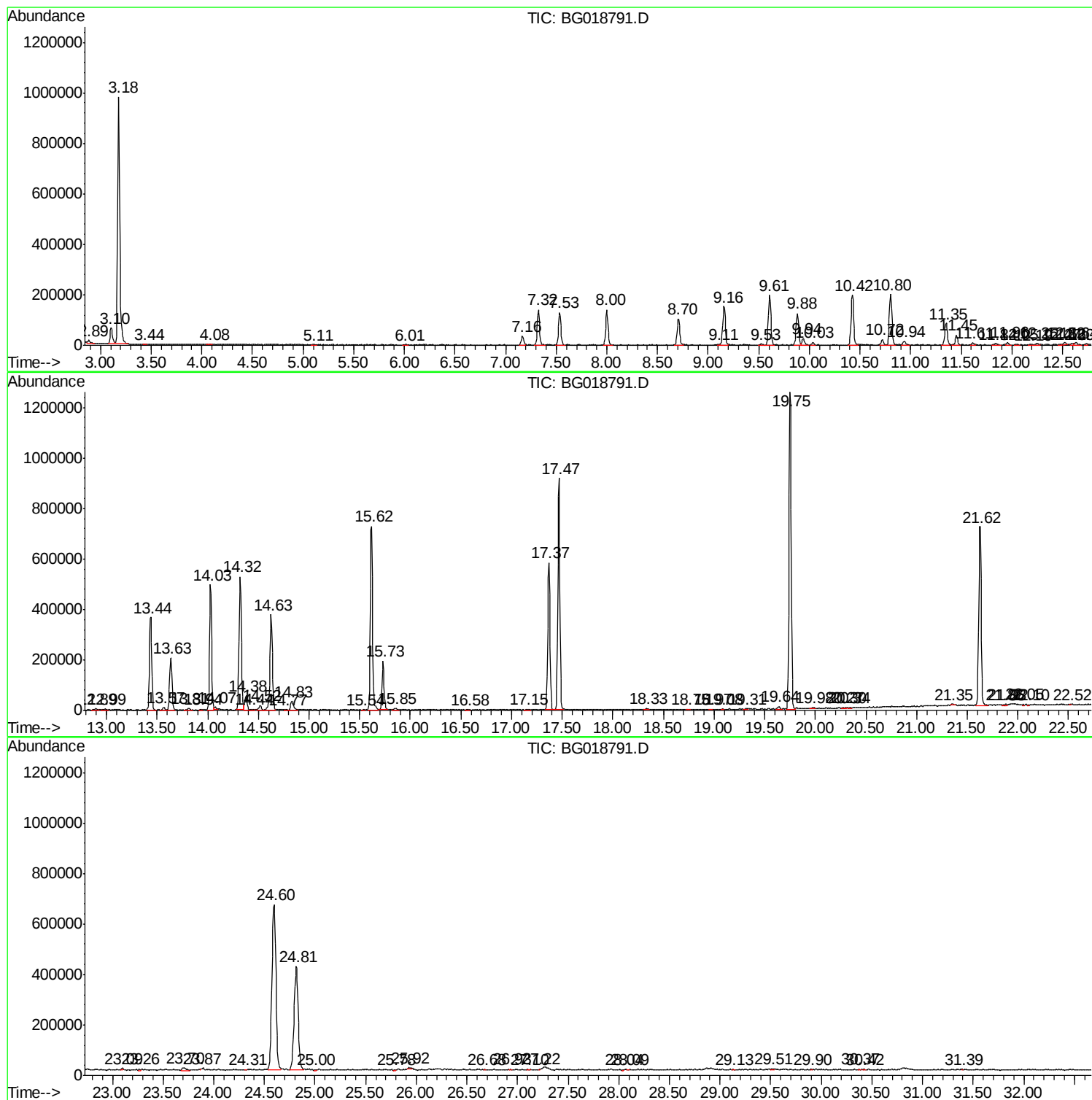
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M
Quant Title : SVOA CALIBRATION

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



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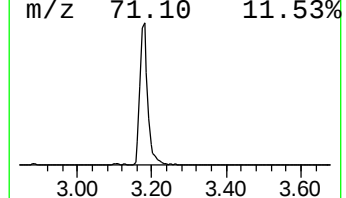
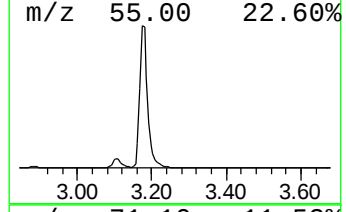
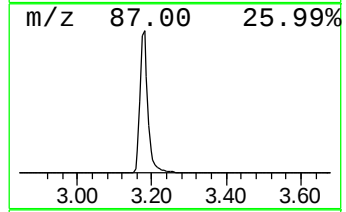
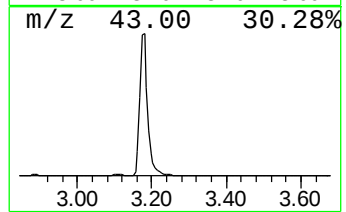
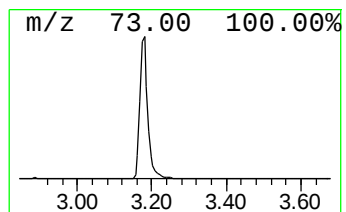
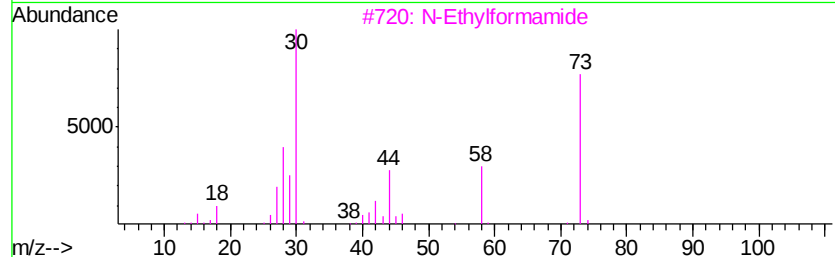
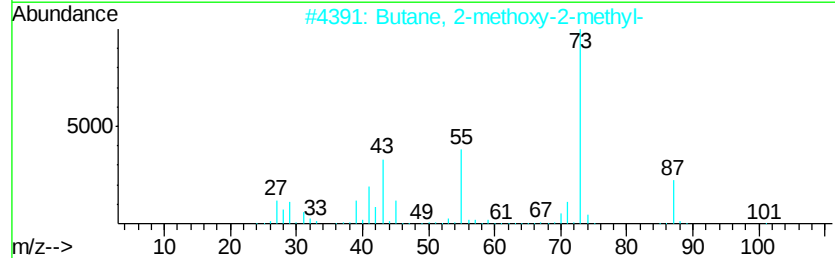
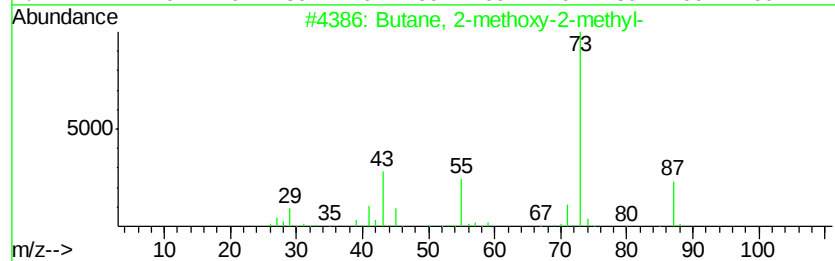
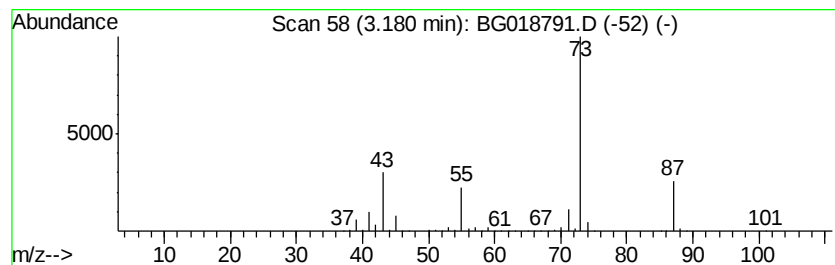
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.18	120.90 ng/ul	1406640	1,4-Dichlorobenzene-d4	8.00

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	78
3		N-Ethylformamide	73	C3H7NO	000627-45-2	9
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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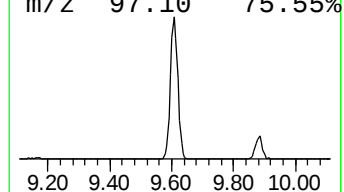
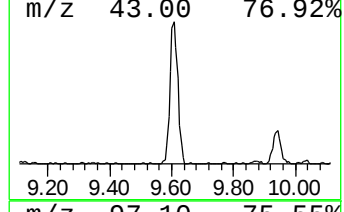
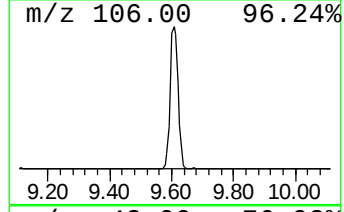
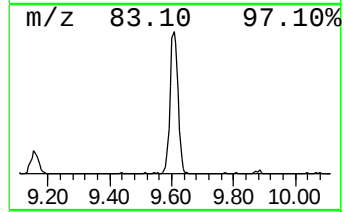
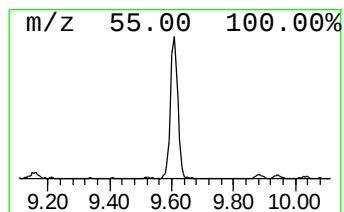
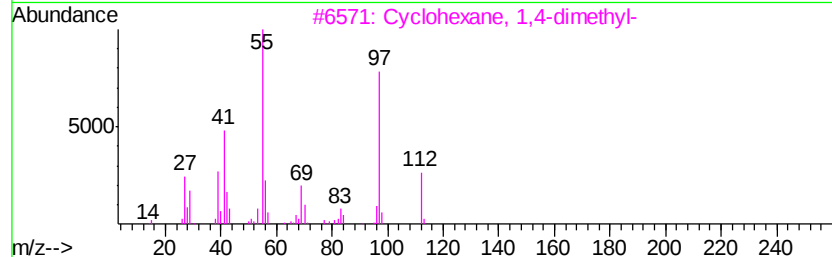
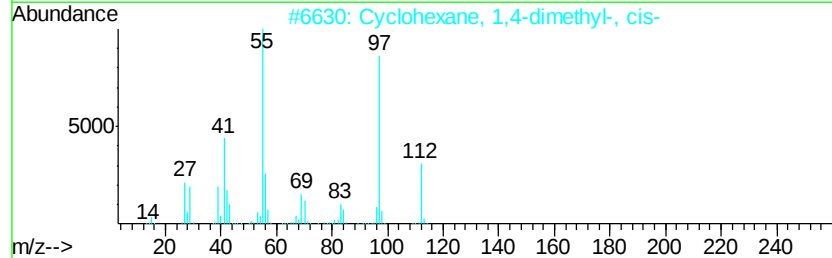
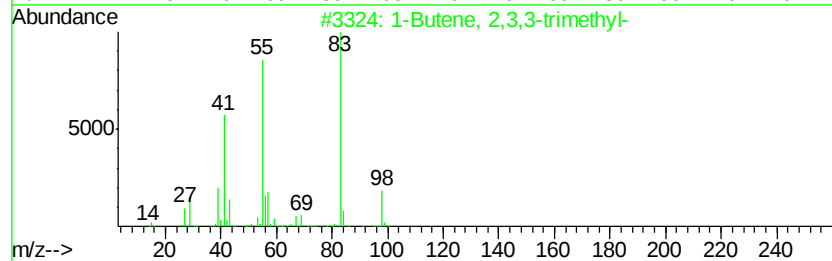
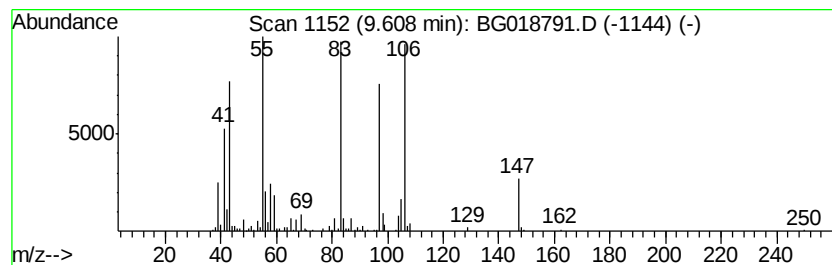
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.61	18.58 ng/ul	334384	Naphthalene-d8	10.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Butene, 2,3,3-trimethyl-	98	C7H14	000594-56-9	27
2		Cyclohexane, 1,4-dimethyl-, cis-	112	C8H16	000624-29-3	27
3		Cyclohexane, 1,4-dimethyl-	112	C8H16	000589-90-2	27
4		Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	27
5		Cyclohexane, 1,4-dimethyl-, cis-	112	C8H16	000624-29-3	27



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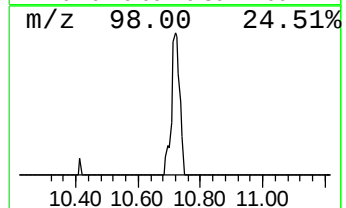
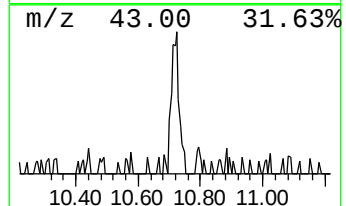
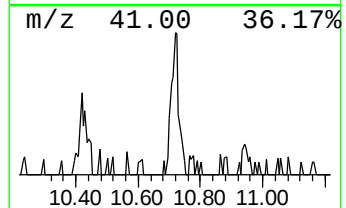
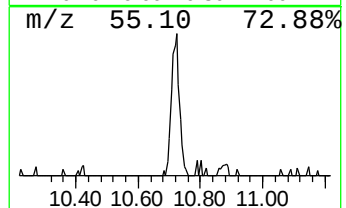
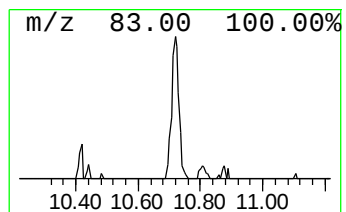
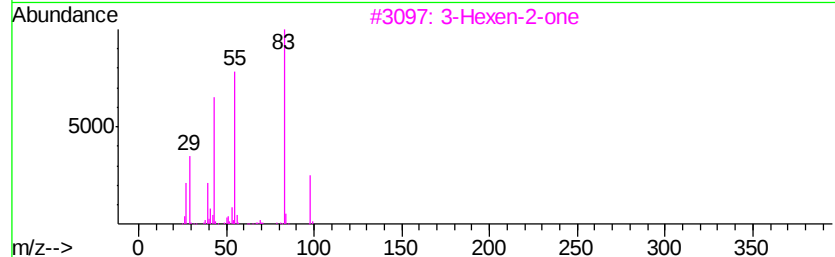
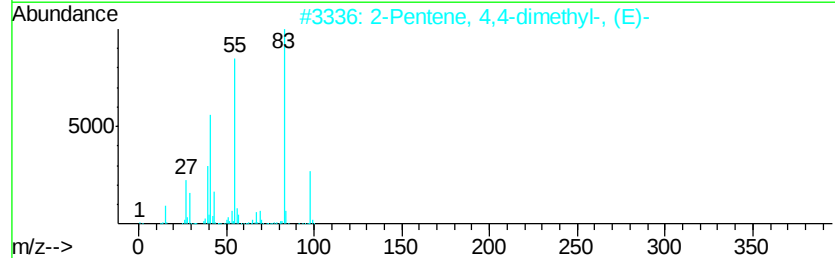
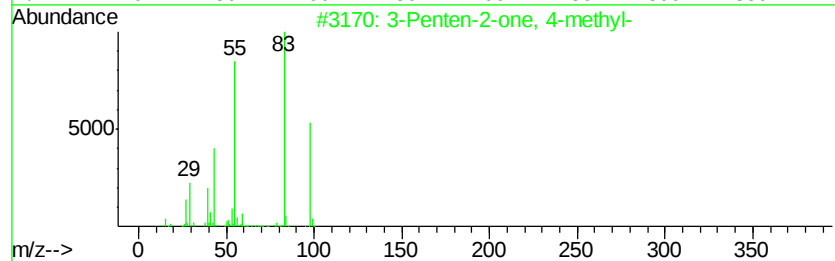
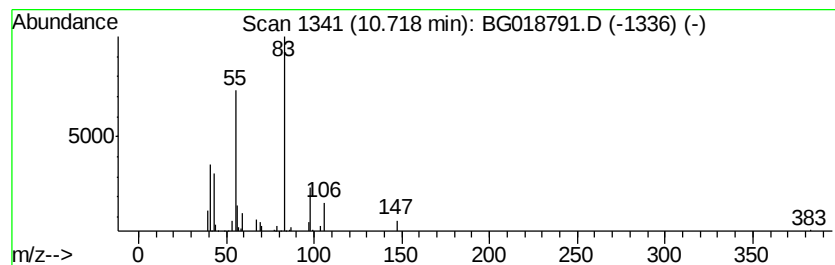
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 3-Penten-2-one, 4-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.72	2.05 ng/ul	36803	Naphthalene-d8	10.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	59
2		2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	58
3		3-Hexen-2-one	98	C6H10O	000763-93-9	53
4		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	53
5		2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	53



Data Path : Z:\HPCHEM1\BNA G\DATA\BG092115\
Data File : BG018791.D
Acq On : 21 Sep 2015 20:33
Operator : TP
Sample : G3696-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ8

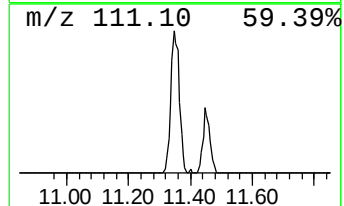
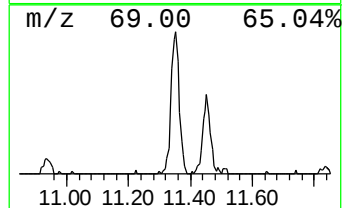
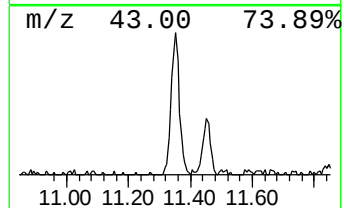
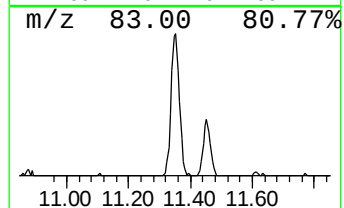
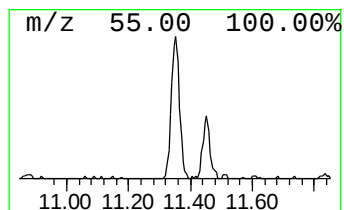
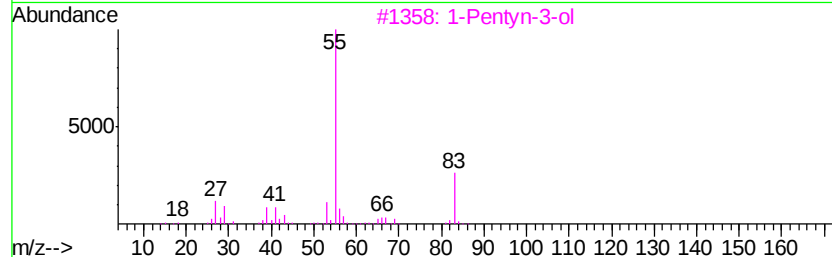
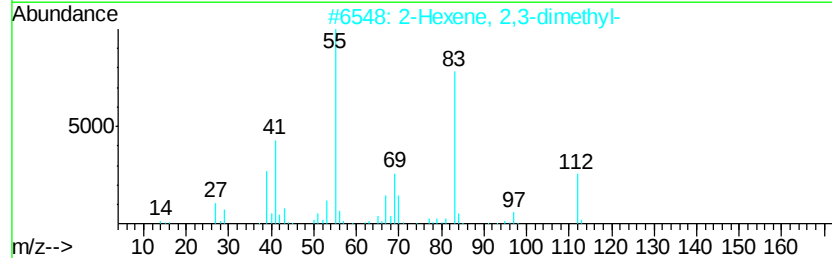
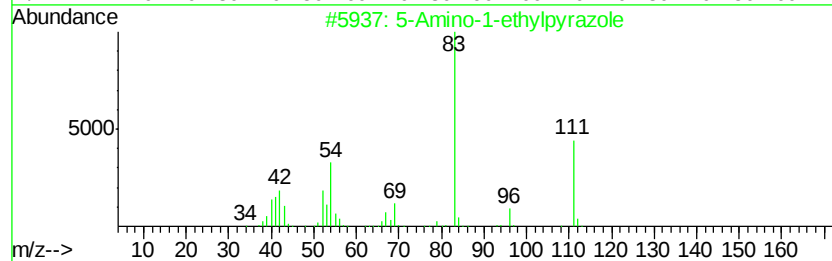
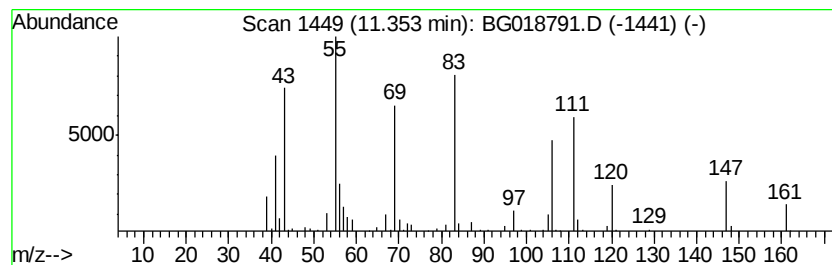
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.35	9.02 ng/ul	162364	Naphthalene-d8	10.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Amino-1-ethylpyrazole	111	C5H9N3	003528-58-3	35
2		2-Hexene, 2,3-dimethyl-	112	C8H16	007145-20-2	30
3		1-Pentyn-3-ol	84	C5H8O	004187-86-4	30
4		3-Ethyl-4-methyl-2-pentene	112	C8H16	019780-68-8	30
5		1-Pentene, 2,3,3-trimethyl-	112	C8H16	000560-23-6	30



Data Path : Z:\HPCHEM1\BNA G\DATA\BG092115\
Data File : BG018791.D
Acq On : 21 Sep 2015 20:33
Operator : TP
Sample : G3696-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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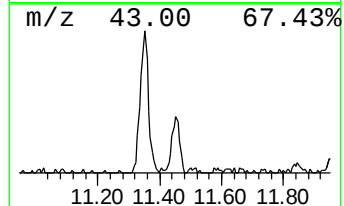
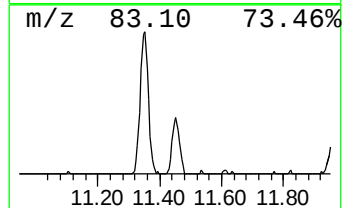
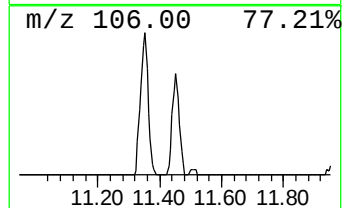
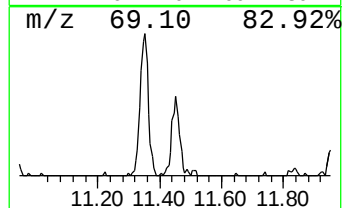
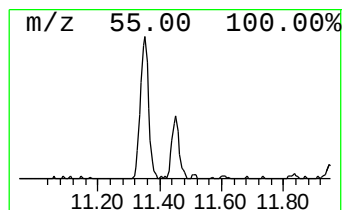
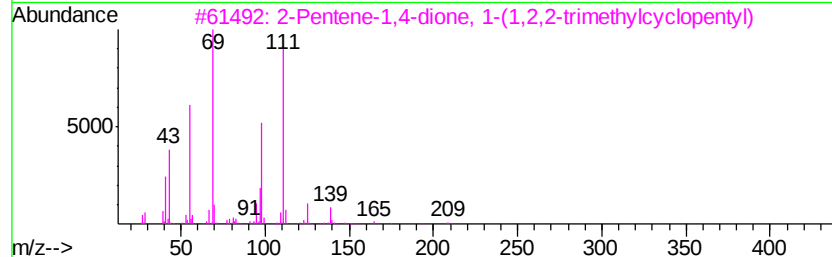
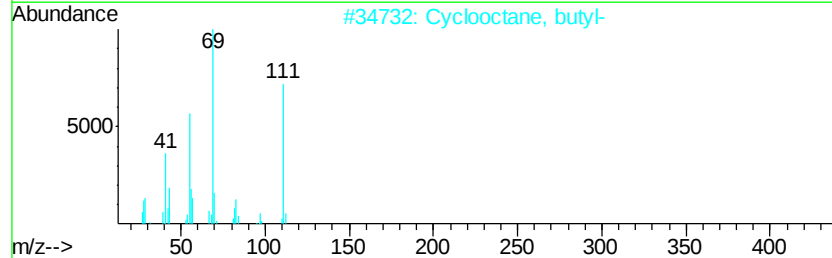
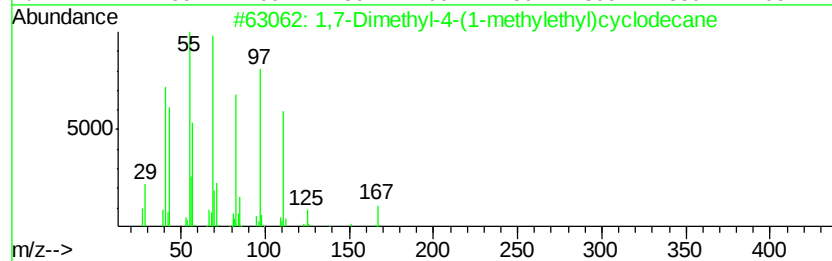
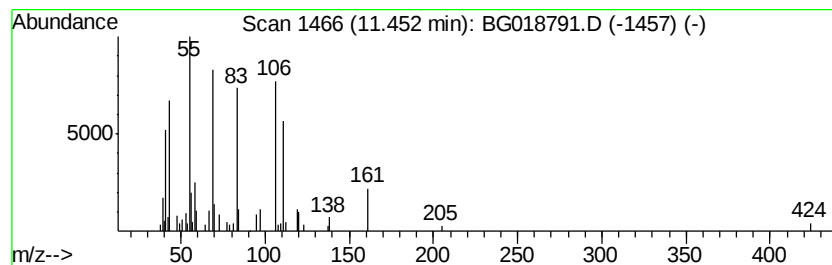
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-03 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.45	3.60 ng/ul	64813	Naphthalene-d8	10.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,7-Dimethyl-4-(1-methylethyl)cv...	210	C15H30	000645-10-3	43
2		Cyclooctane, butyl-	168	C12H24	016538-93-5	22
3		2-Pentene-1,4-dione, 1-(1,2,2-tr...	208	C13H20O2	1000196-77-9	22
4		Cyclooctane, (1-methylpropyl)-	168	C12H24	016538-89-9	22
5		2,4-Dimethyl-3-hexene(c,t)	112	C8H16	1000118-16-4	18



Data Path : Z:\HPCHEM1\BNA G\DATA\BG092115\
Data File : BG018791.D
Acq On : 21 Sep 2015 20:33
Operator : TP
Sample : G3696-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

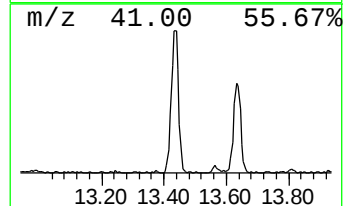
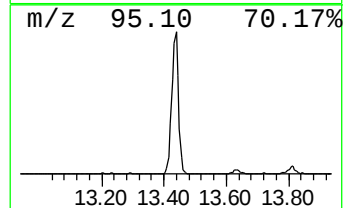
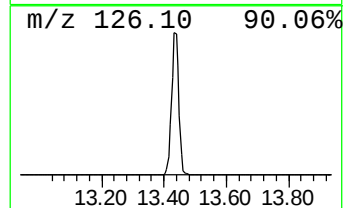
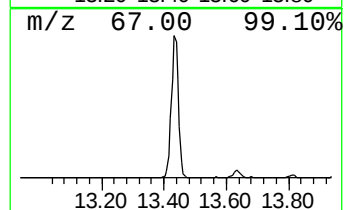
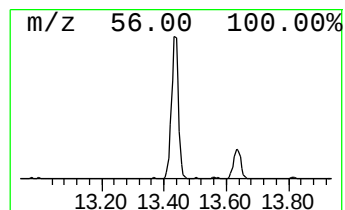
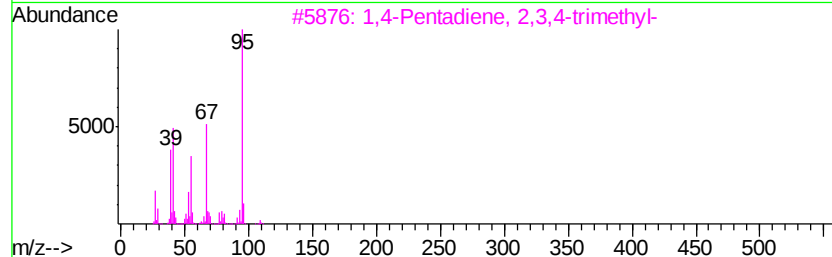
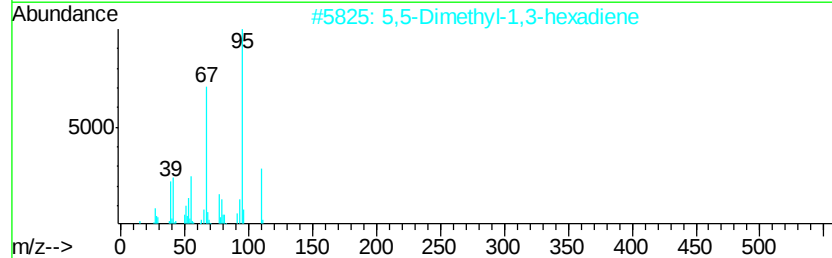
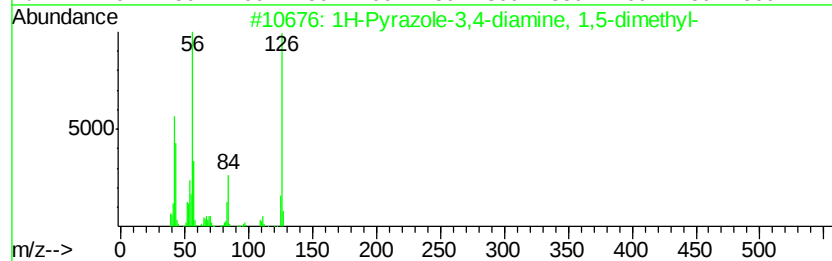
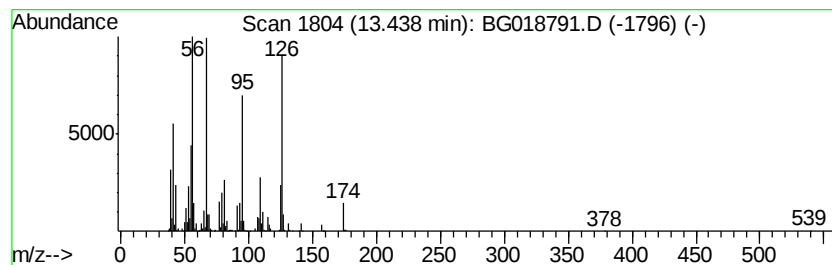
Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-04 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.44	18.02 ng/ul	589441	Acenaphthene-d10	14.63
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-Pyrazole-3,4-diamine, 1,5-dim...	126 C5H10N4	076368-87-1 38
2		5,5-Dimethyl-1,3-hexadiene	110 C8H14	001515-79-3 30
3		1,4-Pentadiene, 2,3,4-trimethyl-	110 C8H14	072014-90-5 22
4		Cyclopentene, 1,2,3-trimethyl-	110 C8H14	000473-91-6 22
5		1,4-Pentadiene, 2,3,3-trimethyl-	110 C8H14	000756-02-5 22



Data Path : Z:\HPCHEM1\BNA G\DATA\BG092115\
Data File : BG018791.D
Acq On : 21 Sep 2015 20:33
Operator : TP
Sample : G3696-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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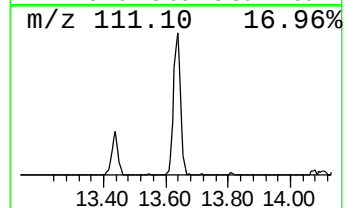
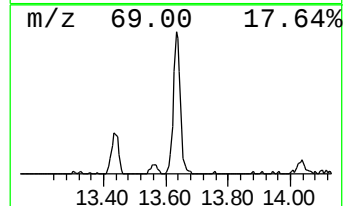
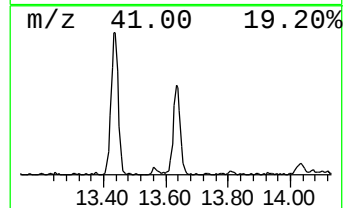
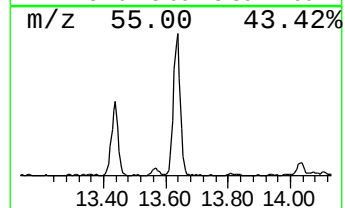
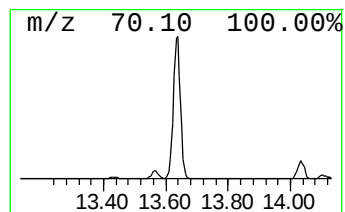
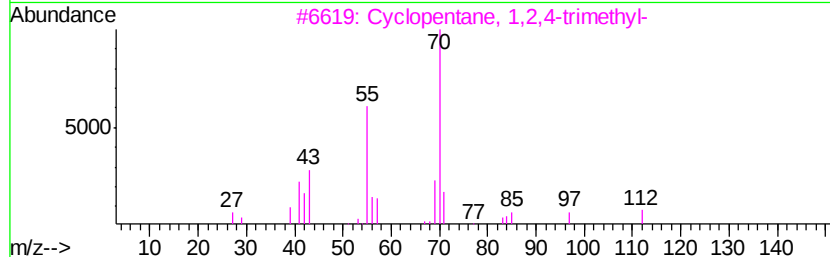
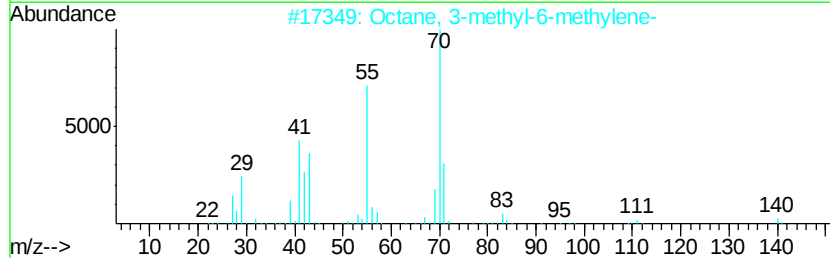
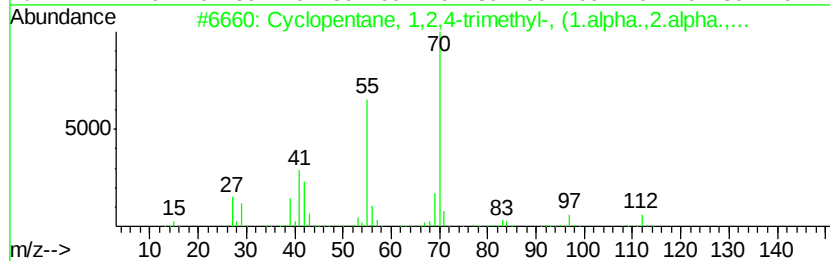
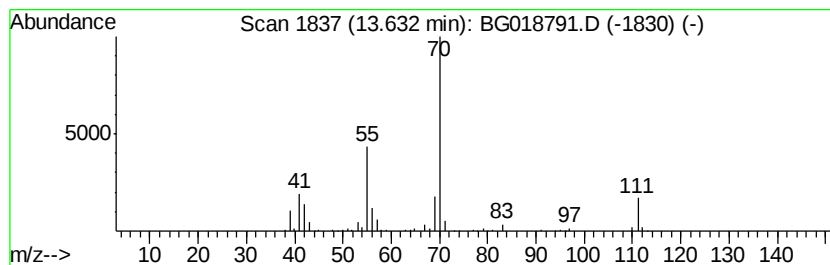
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Cyclic13.63 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.63	10.21 ng/ul	333930	Acenaphthene-d10	14.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, 1,2,4-trimethyl-, ...	112	C8H16	004850-28-6	72
2		Octane, 3-methyl-6-methylene-	140	C10H20	074630-07-2	56
3		Cyclopentane, 1,2,4-trimethyl-	112	C8H16	002815-58-9	56
4		Cyclopentane, 1,2,4-trimethyl-	112	C8H16	002815-58-9	56
5		2,3-Dimethyl-1-hexene	112	C8H16	016746-86-4	45



Data Path : Z:\HPCHEM1\BNA G\DATA\BG092115\
Data File : BG018791.D
Acq On : 21 Sep 2015 20:33
Operator : TP
Sample : G3696-06
Misc :
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Instrument :
BNA_G
ClientSampleId :
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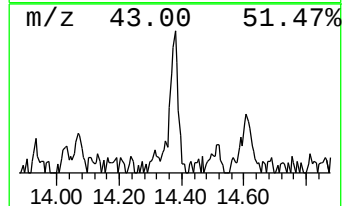
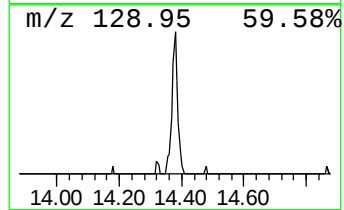
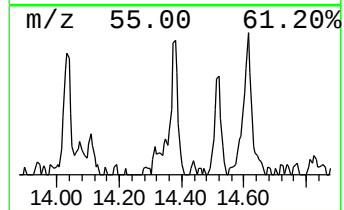
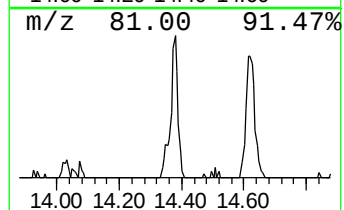
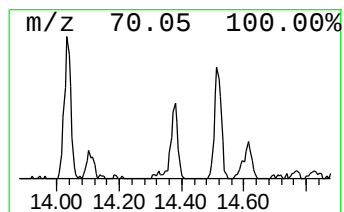
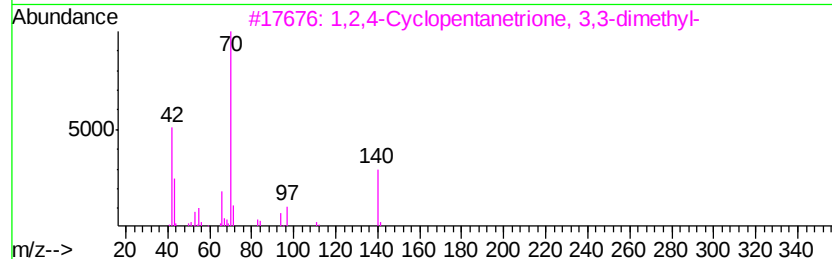
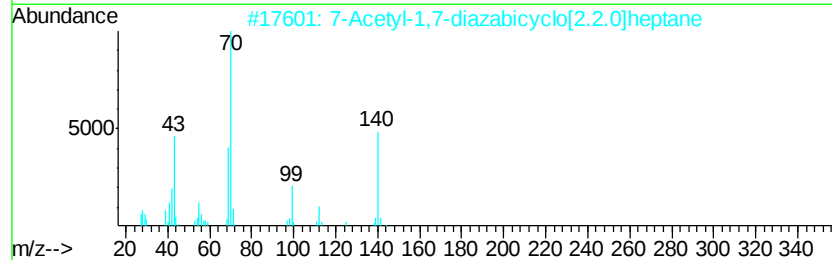
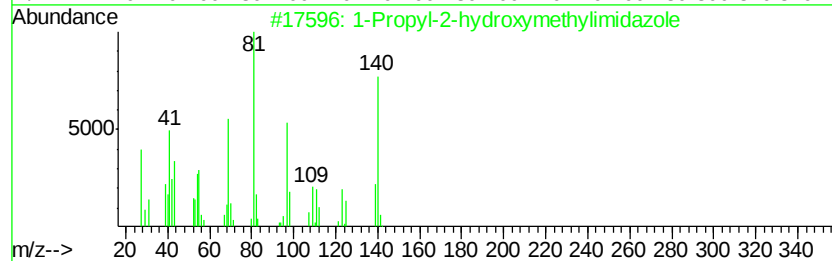
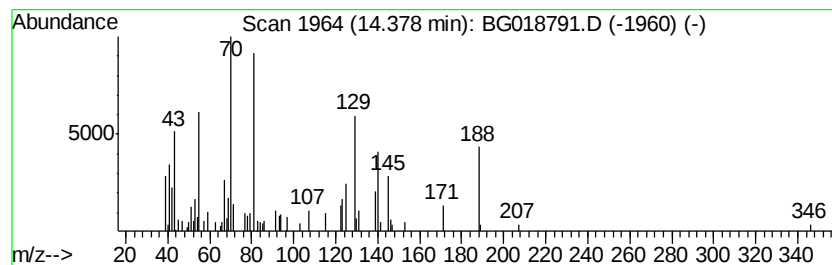
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-05 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.38	2.47 ng/ul	80950	Acenaphthene-d10	14.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Propyl-2-hydroxymethylimidazole	140	C7H12N2O	143886-50-4	27
2		7-Acetyl-1,7-diazabicyclo[2.2.0]...	140	C7H12N2O	036553-03-4	22
3		1,2,4-Cyclopentanetrione, 3,3-di...	140	C7H8O3	017530-56-2	18
4		2-Nonene, 3-methyl-, (E)-	140	C10H20	017003-99-5	14
5		1-Norbornanecarboxylic acid, 7-o...	168	C9H12O3	090673-79-3	12



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092115\
Data File : BG018791.D
Acq On : 21 Sep 2015 20:33
Operator : TP
Sample : G3696-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ8

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	120.9	ng/ul	1406640	1	8.00	232689	20.0
unknown-01	9.61	18.6	ng/ul	334384	2	10.80	359860	20.0
3-Penten-2-one, 4...	10.72	2.0	ng/ul	36803	2	10.80	359860	20.0
unknown-02	11.35	9.0	ng/ul	162364	2	10.80	359860	20.0
unknown-03	11.45	3.6	ng/ul	64813	2	10.80	359860	20.0
unknown-04	13.44	18.0	ng/ul	589441	3	14.63	654339	20.0
(DEL) Alkane: Cyc...	13.63	10.2	ng/ul	333930	3	14.63	654339	20.0
unknown-05	14.38	2.5	ng/ul	80950	3	14.63	654339	20.0