

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
 Data File : BG018908.D
 Acq On : 26 Sep 2015 12:07
 Operator : UM/NP
 Sample : G3797-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 D9G27

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	2.873	4	6	13	rVB5	7151	10922	0.56%	0.054%
2	3.096	38	44	52	rBV	75453	128489	6.61%	0.639%
3	3.167	52	56	66	rVB	181477	237055	12.19%	1.179%
4	6.063	541	549	556	rBV	72249	110297	5.67%	0.549%
5	6.527	624	628	631	rVV2	7534	12500	0.64%	0.062%
6	6.568	631	635	641	rVV	12946	20882	1.07%	0.104%
7	6.633	641	646	652	rVV	22567	37362	1.92%	0.186%
8	7.168	729	737	746	rBV2	42228	112459	5.78%	0.559%
9	7.315	756	762	772	rBV	152051	242720	12.48%	1.207%
10	7.414	774	779	784	rVB4	8704	14686	0.76%	0.073%
11	7.526	790	798	803	rBV2	203326	338699	17.42%	1.685%
12	7.573	803	806	813	rVV	53166	78777	4.05%	0.392%
13	7.761	832	838	846	rVB	122466	190630	9.80%	0.948%
14	7.949	865	870	871	rVV4	7042	12145	0.62%	0.060%
15	7.990	871	877	883	rVV	154789	266864	13.72%	1.327%
16	8.049	883	887	894	rVV3	14589	26673	1.37%	0.133%
17	8.143	897	903	910	rVV7	5671	10633	0.55%	0.053%
18	8.225	910	917	925	rVV	132165	234208	12.04%	1.165%
19	8.296	925	929	934	rVV4	9565	17104	0.88%	0.085%
20	8.513	963	966	971	rBV2	7029	11500	0.59%	0.057%
21	8.601	976	981	984	rBV4	6773	11148	0.57%	0.055%
22	8.701	991	998	1005	rBV	135364	226992	11.67%	1.129%
23	8.954	1036	1041	1047	rBV5	6502	14770	0.76%	0.073%
24	9.083	1057	1063	1068	rBV2	21998	39620	2.04%	0.197%
25	9.148	1068	1074	1086	rVB	164827	293703	15.10%	1.461%
26	9.306	1093	1101	1107	rBV3	24215	55745	2.87%	0.277%
27	9.530	1132	1139	1145	rBV	36012	61416	3.16%	0.305%
28	9.712	1162	1170	1176	rBV3	5720	11171	0.57%	0.056%
29	9.870	1189	1197	1200	rBV	136463	250203	12.87%	1.245%
30	9.906	1200	1203	1208	rVV	104422	153728	7.90%	0.765%
31	9.964	1208	1213	1218	rVV	119169	192526	9.90%	0.958%
32	10.029	1218	1224	1230	rVV5	55882	130215	6.70%	0.648%
33	10.088	1230	1234	1242	rVB2	30676	55755	2.87%	0.277%
34	10.264	1260	1264	1268	rBV2	22697	33228	1.71%	0.165%

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35	10.417	1284	1290	1302	rVB	242992	428825	22.05%	2.133%
36	10.552	1307	1313	1320	rBV2	78635	137858	7.09%	0.686%
37	10.793	1347	1354	1360	rVV	214771	400880	20.61%	1.994%
38	11.075	1398	1402	1408	rVB7	8994	18543	0.95%	0.092%
39	11.151	1409	1415	1420	rBV2	23362	43978	2.26%	0.219%
40	11.316	1432	1443	1448	rBV	120493	199666	10.27%	0.993%
41	11.375	1448	1453	1459	rVV	148049	235062	12.09%	1.169%
42	11.439	1459	1464	1472	rVB2	31572	62236	3.20%	0.310%
43	11.527	1472	1479	1485	rBV	60525	115647	5.95%	0.575%
44	11.586	1485	1489	1494	rVB8	7161	13418	0.69%	0.067%
45	11.680	1501	1505	1508	rVB	11955	17538	0.90%	0.087%
46	11.833	1526	1531	1539	rVB4	6942	16092	0.83%	0.080%
47	11.956	1542	1552	1556	rBV	91884	157448	8.10%	0.783%
48	12.003	1556	1560	1562	rVV2	63903	104271	5.36%	0.519%
49	12.027	1562	1564	1567	rVV	69710	79583	4.09%	0.396%
50	12.068	1567	1571	1576	rVB	50951	69523	3.57%	0.346%
51	12.244	1594	1601	1602	rVV3	18021	29377	1.51%	0.146%
52	12.291	1606	1609	1617	rVV2	22948	40560	2.09%	0.202%
53	12.509	1641	1646	1653	rVB5	5342	11763	0.60%	0.059%
54	12.843	1700	1703	1708	rVV2	21476	29888	1.54%	0.149%
55	12.973	1718	1725	1733	rBV	984002	1514707	77.89%	7.534%
56	13.026	1733	1734	1737	rVV	12148	13327	0.69%	0.066%
57	13.061	1737	1740	1746	rVB	22226	36108	1.86%	0.180%
58	13.225	1761	1768	1775	rBV	1354755	1944753	100.00%	9.674%
59	13.360	1787	1791	1798	rVB5	10167	18804	0.97%	0.094%
60	13.531	1811	1820	1825	rBV3	44285	94408	4.85%	0.470%
61	13.660	1838	1842	1847	rBV7	8173	13860	0.71%	0.069%
62	13.725	1849	1853	1860	rBV2	27447	51248	2.64%	0.255%
63	13.819	1865	1869	1873	rVB5	8456	13452	0.69%	0.067%
64	13.919	1882	1886	1891	rVB5	6977	11707	0.60%	0.058%
65	14.019	1891	1903	1907	rBV	447706	682470	35.09%	3.395%
66	14.060	1907	1910	1917	rVV	78941	114389	5.88%	0.569%
67	14.136	1917	1923	1928	rVV6	18998	32892	1.69%	0.164%
68	14.254	1936	1943	1946	rBV5	7738	14168	0.73%	0.070%
69	14.306	1946	1952	1967	rVV	469136	774951	39.85%	3.855%
70	14.465	1974	1979	1989	rVB	41291	72125	3.71%	0.359%
71	14.612	1992	2004	2012	rBV2	372030	607961	31.26%	3.024%

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72	14.823	2033	2040	2048	rVV2	54386	107223	5.51%	0.533%
73	15.199	2100	2104	2108	rVV4	14531	21934	1.13%	0.109%
74	15.293	2111	2120	2125	rBV5	15550	38685	1.99%	0.192%
75	15.346	2125	2129	2134	rVV	60300	80873	4.16%	0.402%
76	15.423	2134	2142	2148	rVV	1087483	1479561	76.08%	7.360%
77	15.605	2166	2173	2185	rVV	407181	638646	32.84%	3.177%
78	15.722	2185	2193	2204	rVV	162921	245262	12.61%	1.220%
79	15.811	2204	2208	2211	rVV2	9137	12651	0.65%	0.063%
80	15.957	2229	2233	2241	rVB3	39998	68098	3.50%	0.339%
81	16.527	2326	2330	2336	rVV2	31116	45421	2.34%	0.226%
82	16.598	2338	2342	2347	rVB2	7968	11516	0.59%	0.057%
83	16.974	2403	2406	2410	rVV2	13359	22586	1.16%	0.112%
84	17.003	2410	2411	2418	rVV5	10157	11662	0.60%	0.058%
85	17.097	2425	2427	2431	rVV4	8321	11373	0.58%	0.057%
86	17.303	2456	2462	2465	rBV3	18347	29400	1.51%	0.146%
87	17.356	2465	2471	2478	rVV2	546876	823218	42.33%	4.095%
88	17.450	2481	2487	2495	rVB	167407	264093	13.58%	1.314%
89	17.626	2512	2517	2522	rVB2	31478	40201	2.07%	0.200%
90	17.955	2568	2573	2578	rVV7	7233	14573	0.75%	0.072%
91	18.014	2578	2583	2590	rVV	339102	446126	22.94%	2.219%
92	18.184	2608	2612	2614	rBV3	10883	14299	0.74%	0.071%
93	18.308	2626	2633	2638	rVB2	44719	72006	3.70%	0.358%
94	19.436	2820	2825	2830	rBV9	7500	15042	0.77%	0.075%
95	19.735	2870	2876	2885	rVB	536836	745097	38.31%	3.706%
96	21.269	3134	3137	3141	rVB4	8052	10545	0.54%	0.052%
97	21.610	3188	3195	3204	rBV	742536	1203592	61.89%	5.987%
98	23.848	3571	3576	3580	rBV8	8763	17365	0.89%	0.086%
99	24.559	3687	3697	3709	rVB	188404	519665	26.72%	2.585%
100	24.776	3724	3734	3748	rVB	446110	1224721	62.98%	6.092%

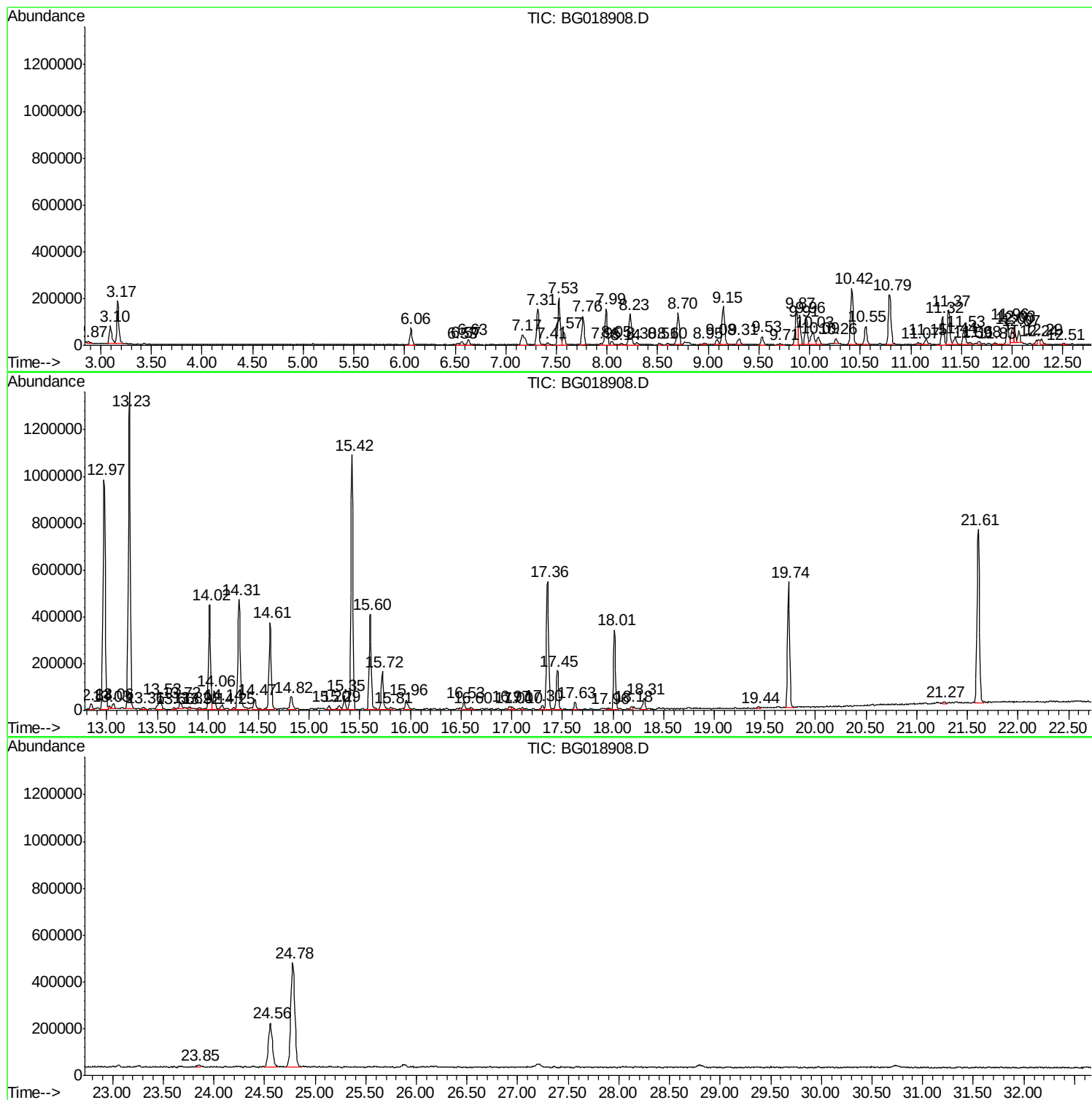
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG091015.M
Quant Title : SVOA CALIBRATION

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



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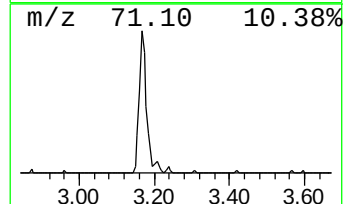
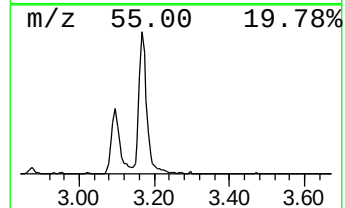
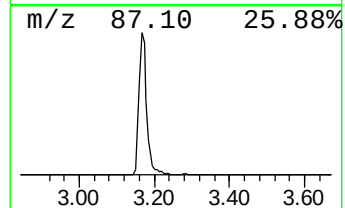
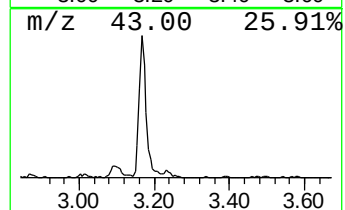
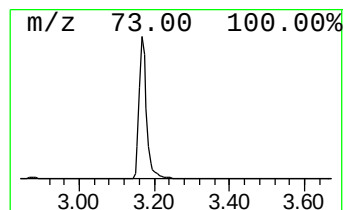
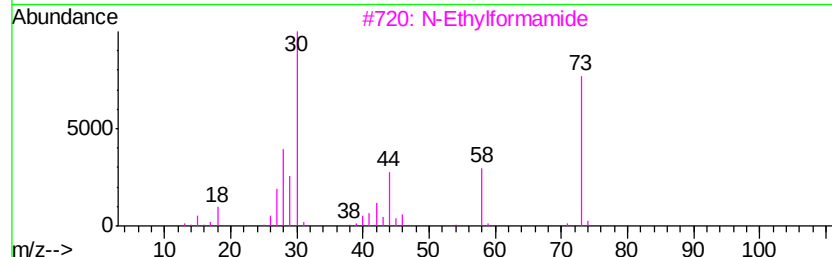
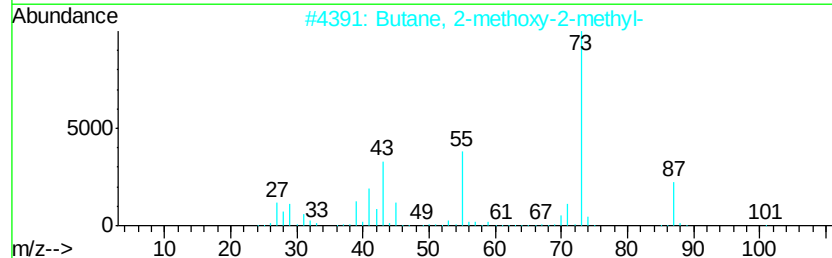
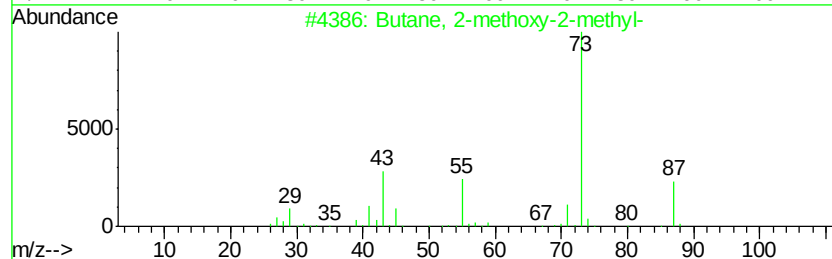
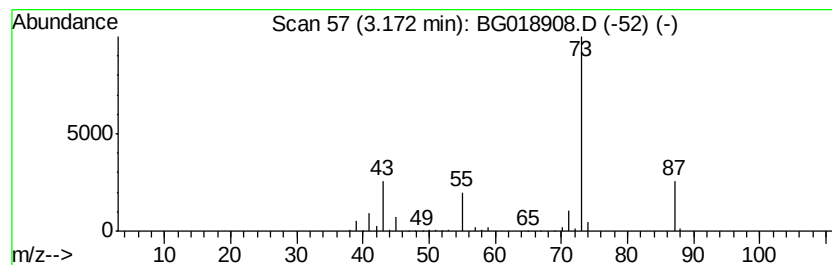
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG091015.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.17	17.77 ng/ul	237055	1,4-Dichlorobenzene-d4	7.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	90
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3		N-Ethylformamide	73	C3H7NO	000627-45-2	9
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	9



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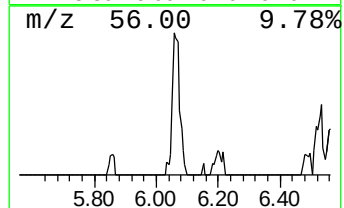
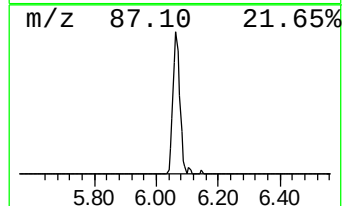
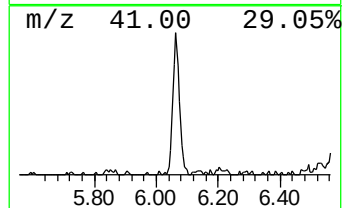
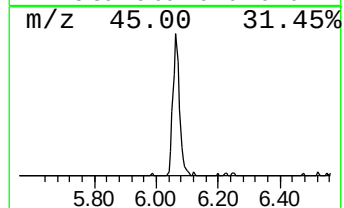
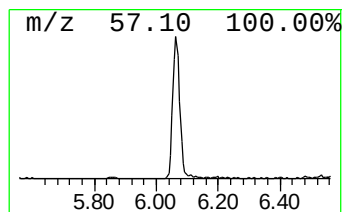
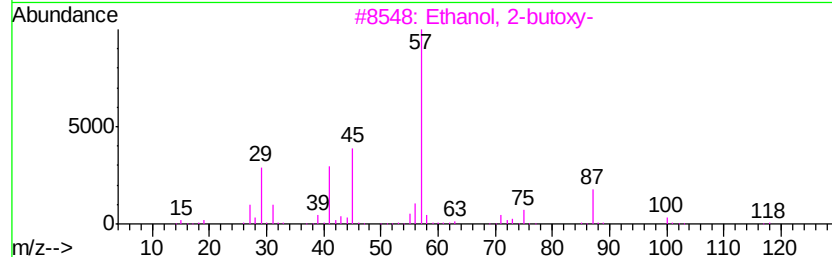
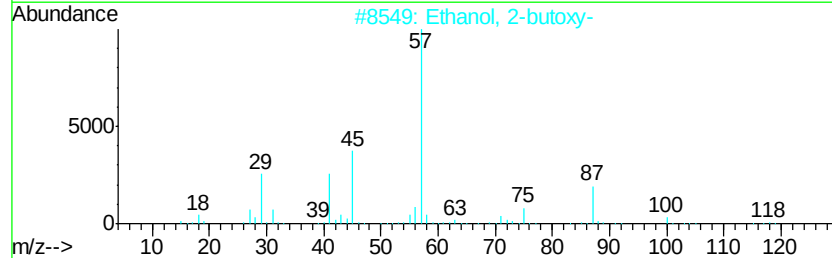
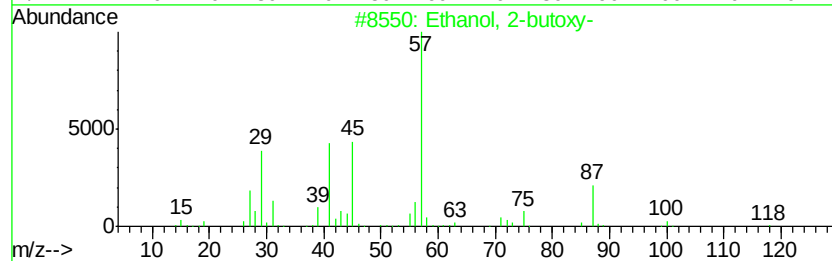
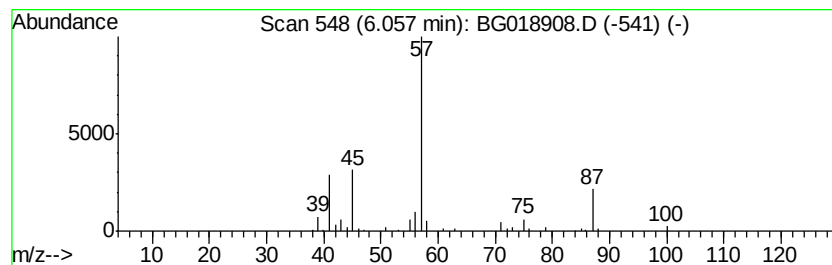
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG091015.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Ethanol, 2-butoxy- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.06	8.27 ng/ul	110297	1,4-Dichlorobenzene-d4	7.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	83
2		Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	64
3		Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	64
4		Propanoic acid, 2-methylpropyl e...	130	C7H14O2	000540-42-1	37
5		3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	37



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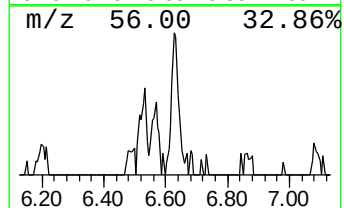
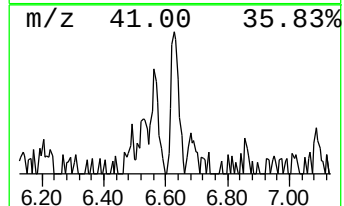
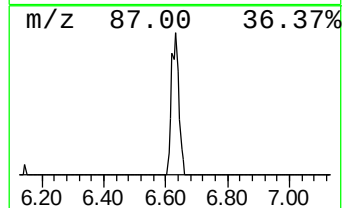
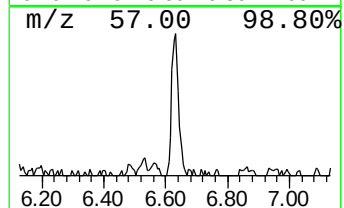
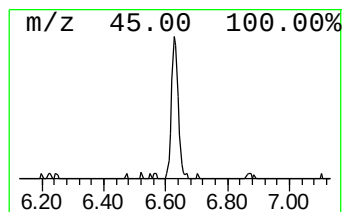
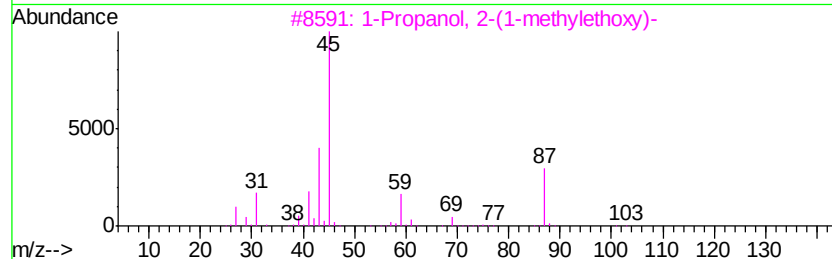
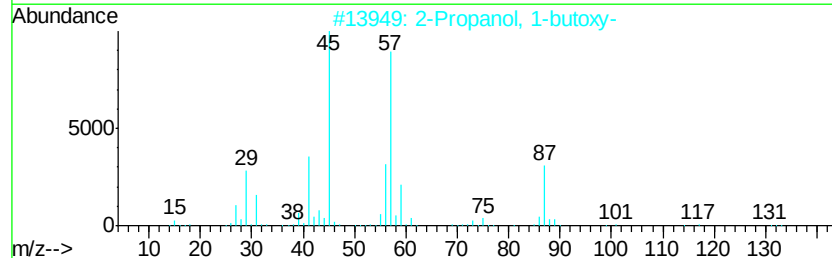
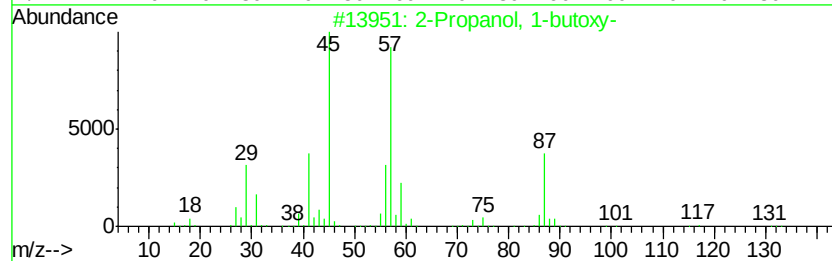
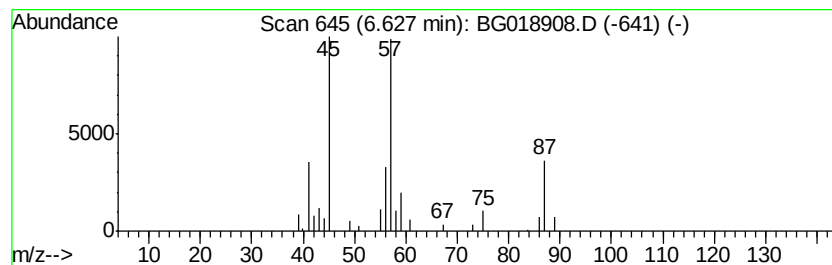
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG091015.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 2-Propanol, 1-butoxy- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	2.80 ng/ul	37362	1,4-Dichlorobenzene-d4	7.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propanol, 1-butoxy-	132	C7H16O2	005131-66-8	78
2		2-Propanol, 1-butoxy-	132	C7H16O2	005131-66-8	50
3		1-Propanol, 2-(1-methylethoxy)-	118	C6H14O2	003944-37-4	42
4		3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	42
5		Propane, 2,2'-[ethylidenebis(oxy...]	146	C8H18O2	004285-59-0	38



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
Operator : UM/NP
Sample : G3797-02
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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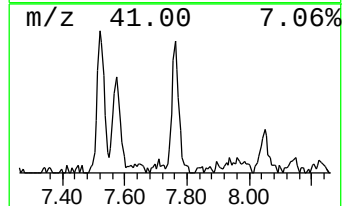
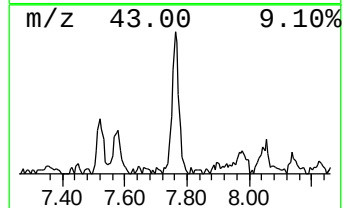
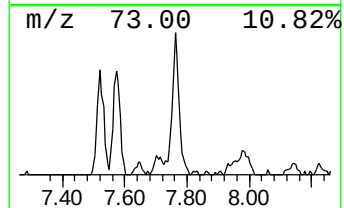
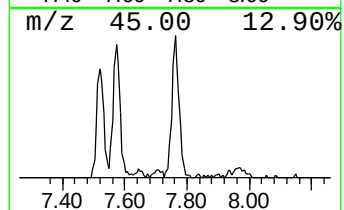
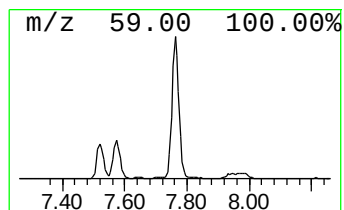
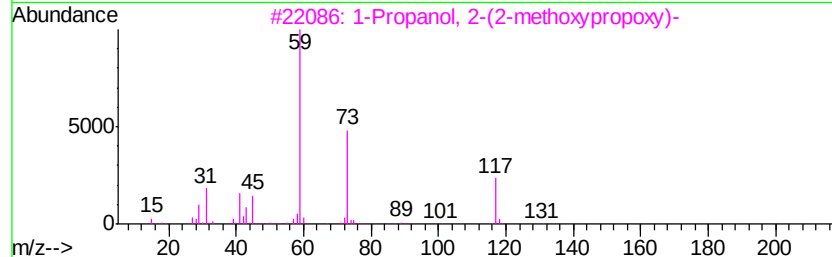
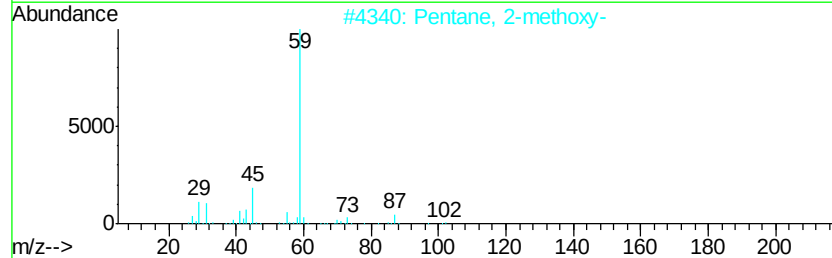
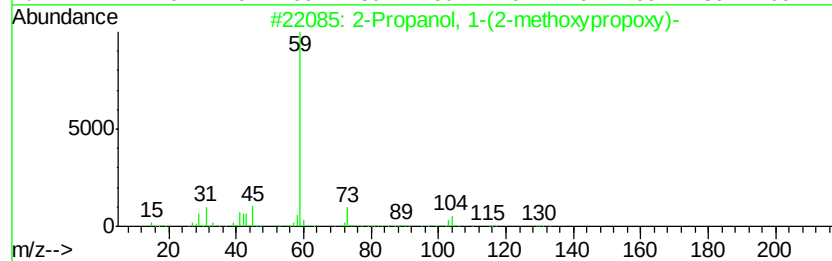
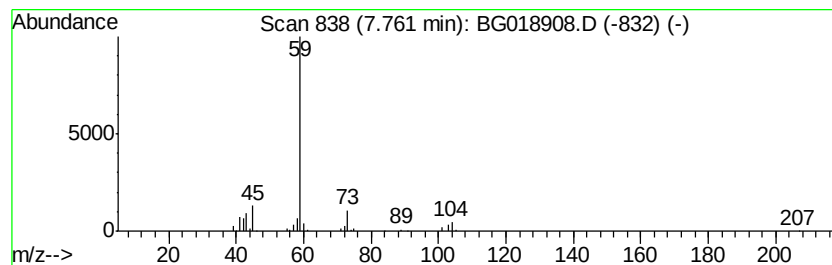
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 2-Propanol, 1-(2-methoxypro... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.76	14.29 ng/ul	190630	1,4-Dichlorobenzene-d4	7.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	83
2		Pentane, 2-methoxy-	102	C6H14O	006795-88-6	64
3		1-Propanol, 2-(2-methoxypropoxy)-	148	C7H16O3	013588-28-8	64
4		2-Propanol, 1,1'-[(1-methyl-1,2-...]	192	C9H20O4	001638-16-0	64
5		2-Propanol, 1-[1-methyl-2-(2-pro...]	174	C9H18O3	055956-25-7	56



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
Operator : UM/NP
Sample : G3797-02
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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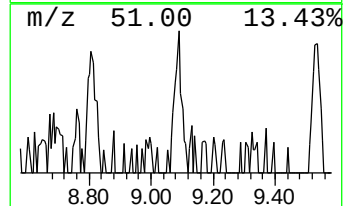
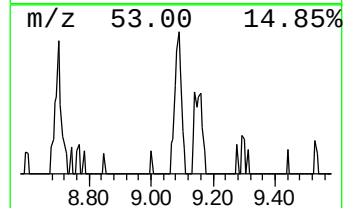
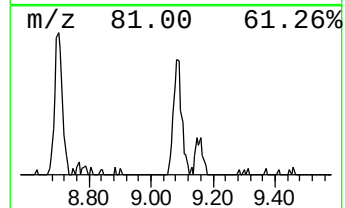
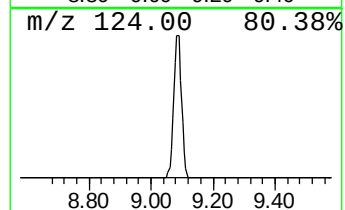
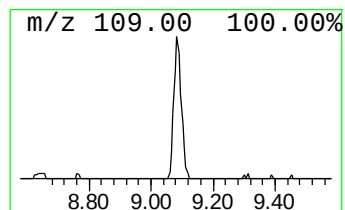
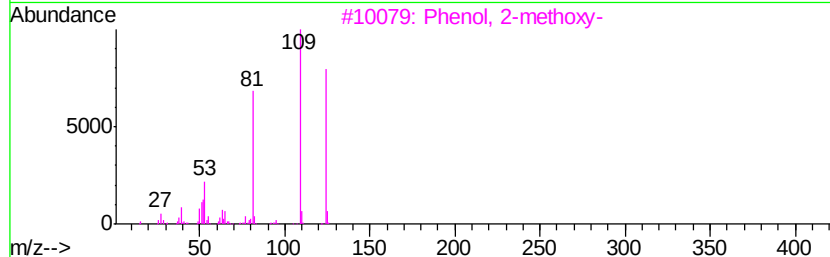
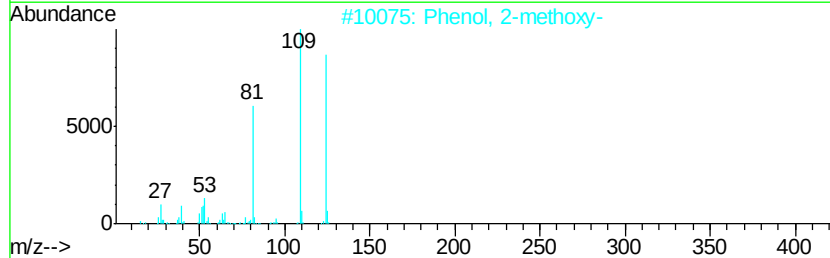
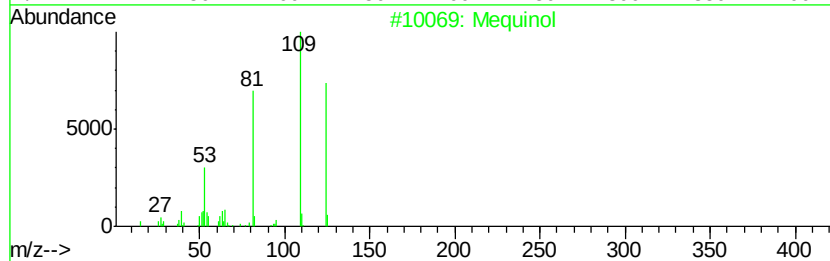
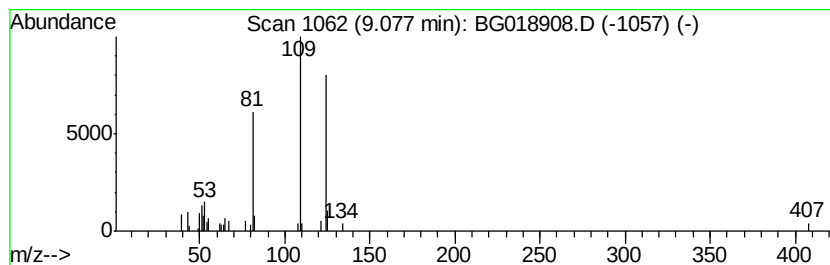
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Mequinol Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.08	2.97 ng/ul	39620	1,4-Dichlorobenzene-d4	7.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Mequinol	124	C7H8O2	000150-76-5	87
2		Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	86
3		Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	86
4		2-Cyclopenten-1-one, 2,3,4-trime...	124	C8H12O	028790-86-5	78
5		Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	78



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
Operator : UM/NP
Sample : G3797-02
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
D9G27

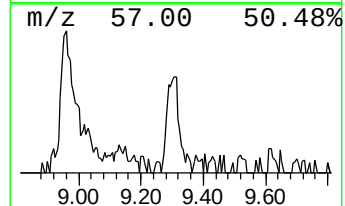
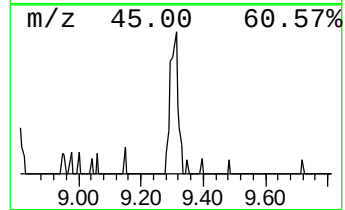
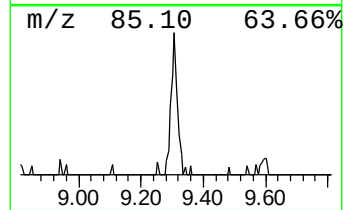
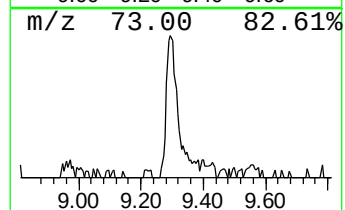
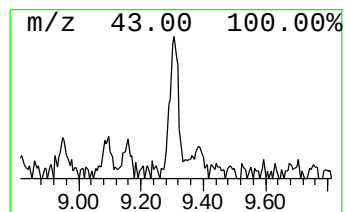
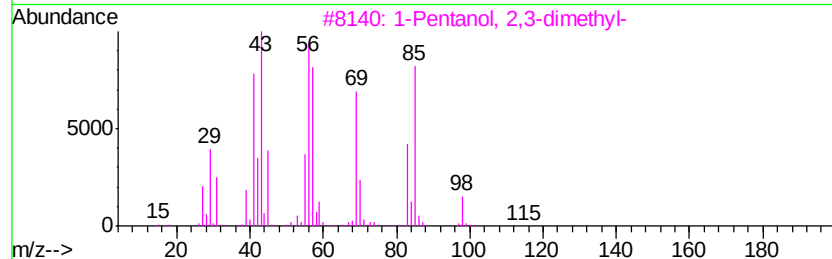
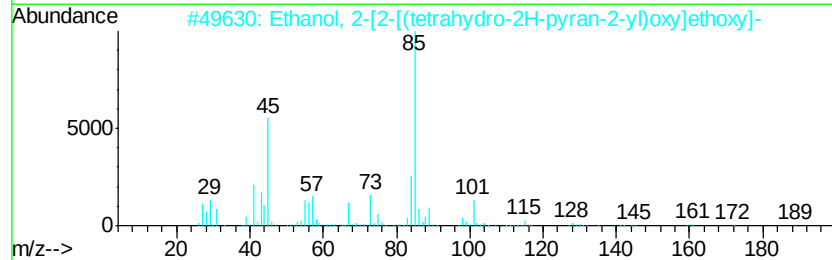
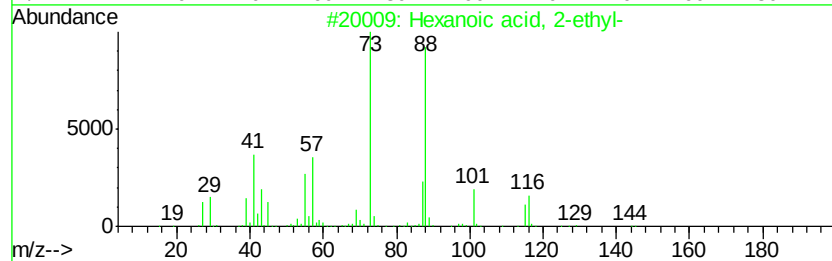
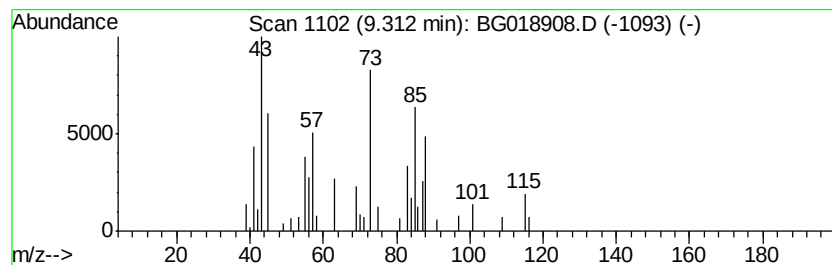
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-01 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.31	4.18 ng/ul	55745	1,4-Dichlorobenzene-d4	7.99

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	27
2			Ethanol, 2-[2-[(tetrahydro-2H-py...	190	C9H18O4	002163-11-3	25
3			1-Pentanol, 2,3-dimethyl-	116	C7H16O	010143-23-4	25
4			1-Pentanol, 2,3-dimethyl-	116	C7H16O	010143-23-4	23
5			Butanoic acid, ethyl ester	116	C6H12O2	000105-54-4	23



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
Operator : UM/NP
Sample : G3797-02
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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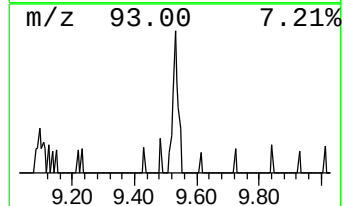
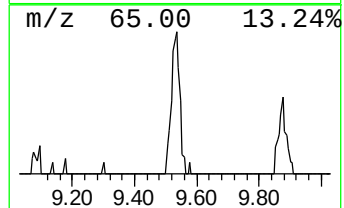
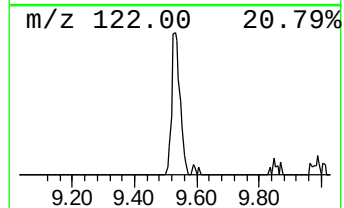
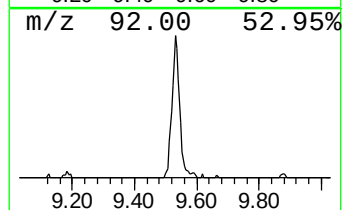
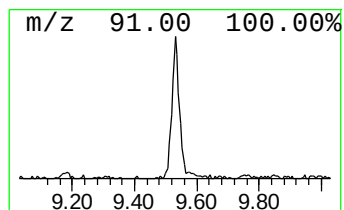
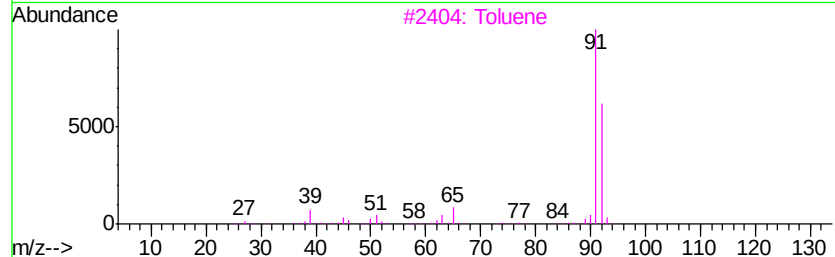
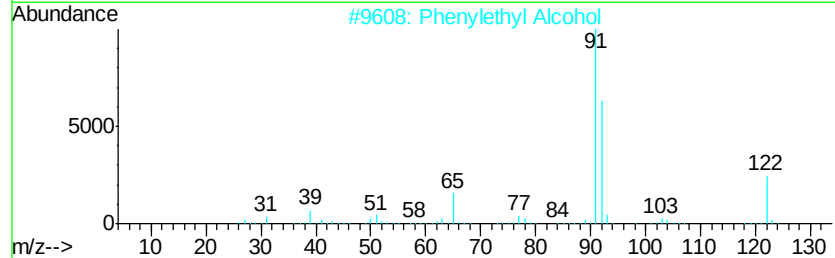
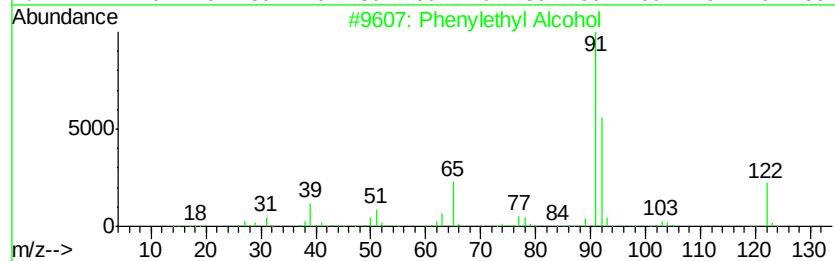
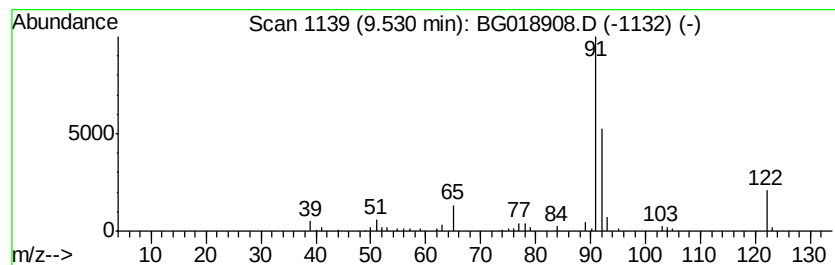
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Phenylethyl Alcohol Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.53	3.06 ng/ul	61416	Napthalene-d8	10.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenylethyl Alcohol	122	C8H10O	000060-12-8	91
2		Phenylethyl Alcohol	122	C8H10O	000060-12-8	91
3		Toluene	92	C7H8	000108-88-3	53
4		Toluene	92	C7H8	000108-88-3	53
5		Phenylethyl Alcohol	122	C8H10O	000060-12-8	52



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
Operator : UM/NP
Sample : G3797-02
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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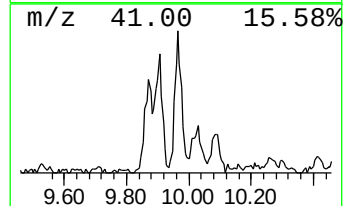
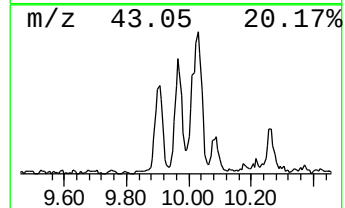
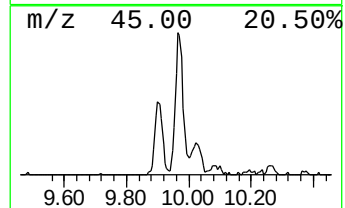
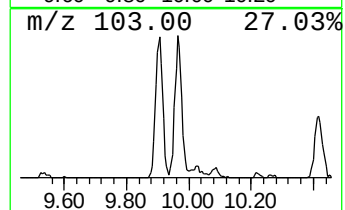
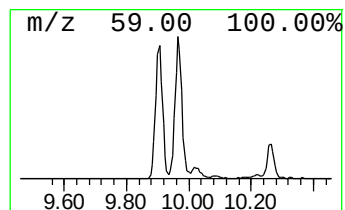
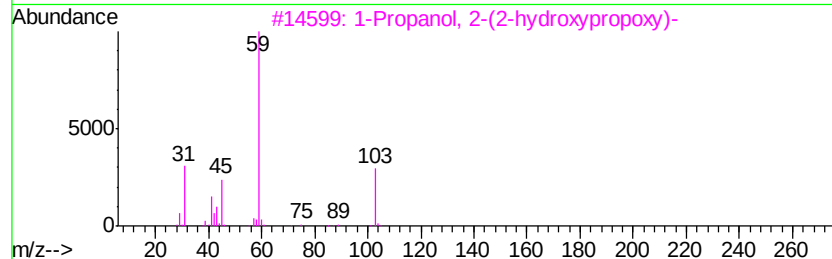
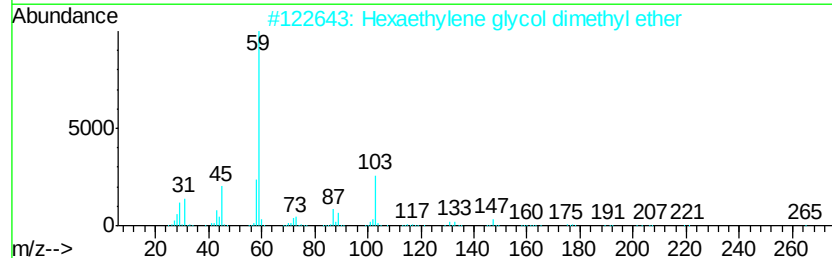
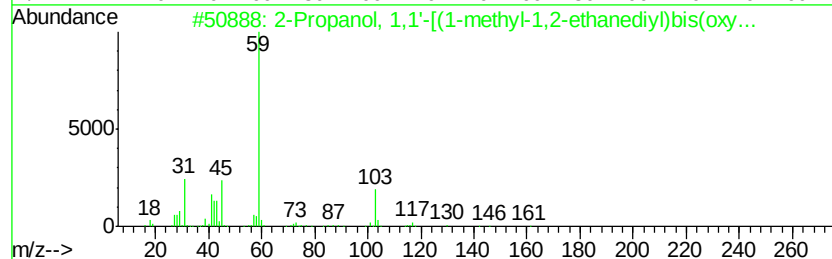
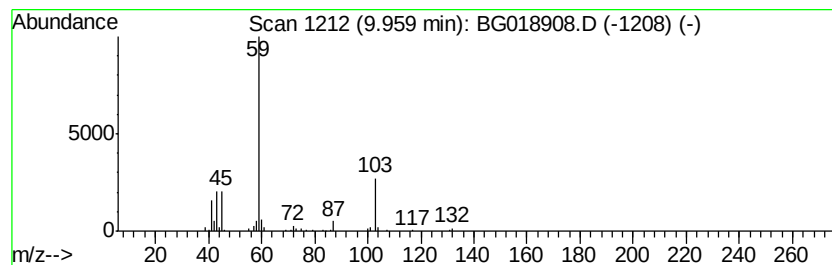
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 2-Propanol, 1,1'-[(1-methyl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.96	9.61 ng/ul	192526	Napthalene-d8	10.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propanol, 1,1'-[(1-methyl-1,2-...	192	C9H20O4	001638-16-0	78
2			Hexaethylene glycol dimethyl ether	310	C14H30O7	001072-40-8	74
3			1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72
4			2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	72
5			Methane, diethoxy-	104	C5H12O2	000462-95-3	72



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
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Sample : G3797-02
Misc :
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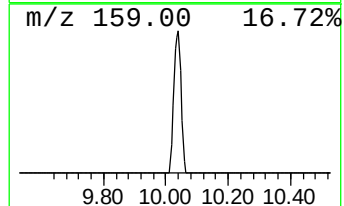
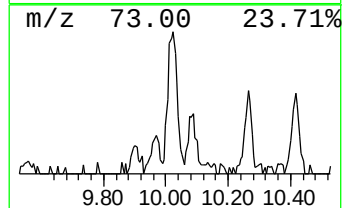
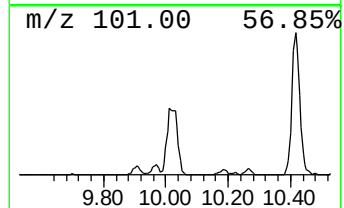
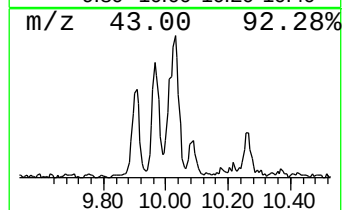
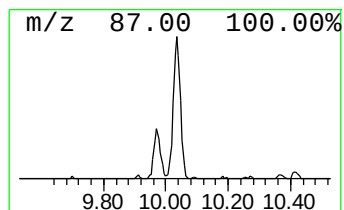
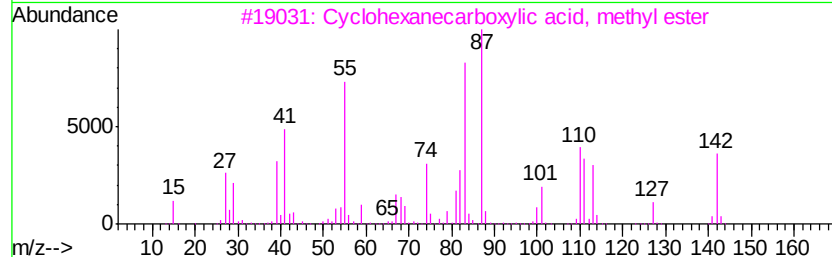
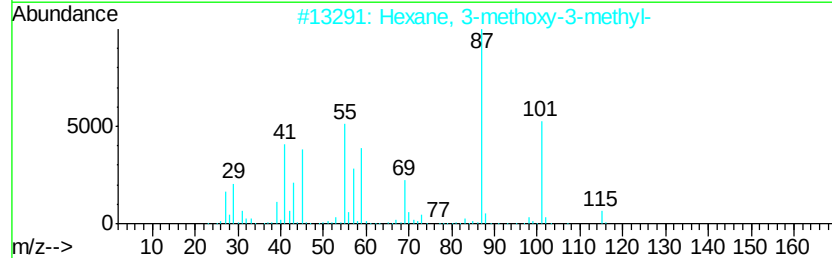
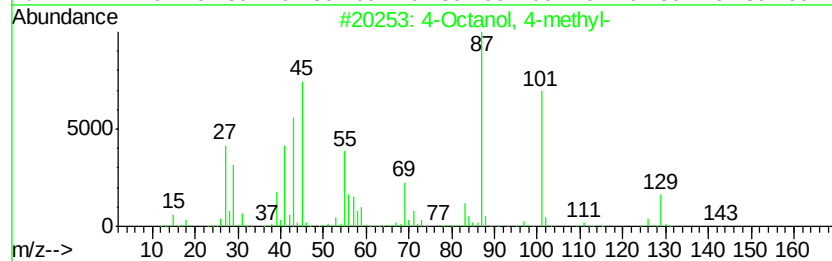
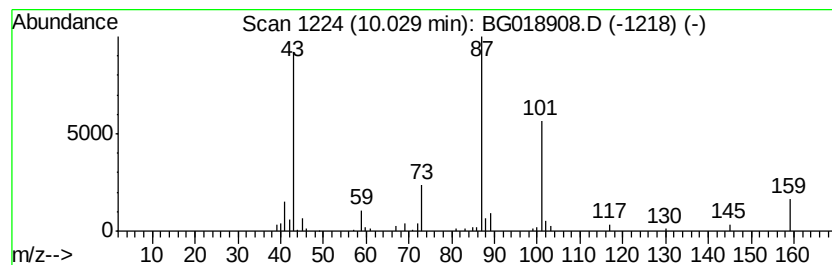
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-02 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.03	6.50 ng/ul	130215	Naphthalene-d8	10.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Octanol, 4-methyl-	144	C9H20O	023418-37-3	45
2			Hexane, 3-methoxy-3-methyl-	130	C8H18O	074630-91-4	45
3			Cyclohexanecarboxylic acid, meth...	142	C8H14O2	004630-82-4	43
4			Hydrazine, 1,1-diethyl-2-pentyl-	158	C9H22N2	067398-41-8	38
5			Ethyl acetoacetate ethylene acetal	174	C8H14O4	006413-10-1	38



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
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Sample : G3797-02
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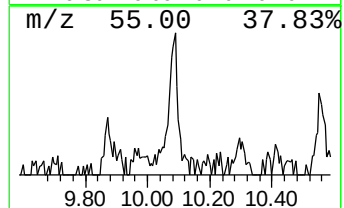
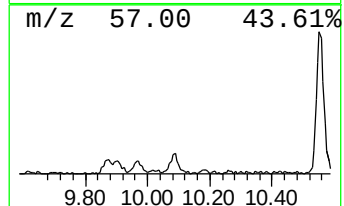
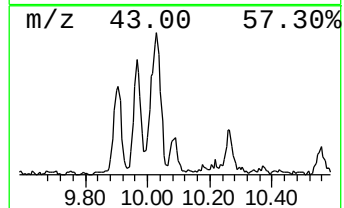
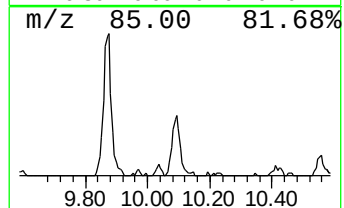
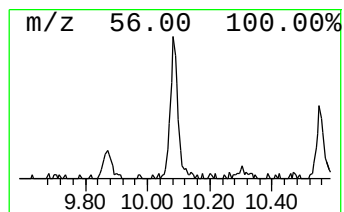
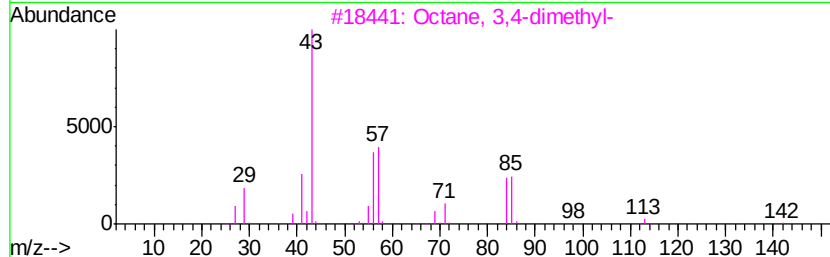
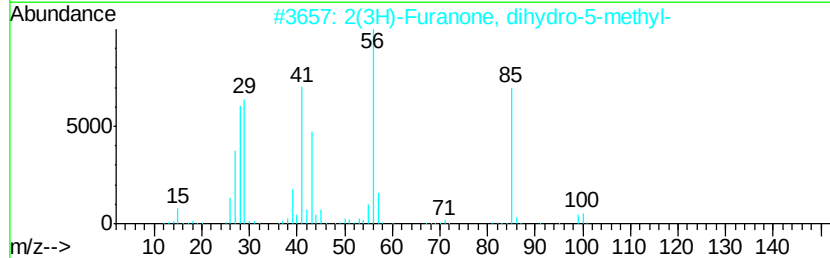
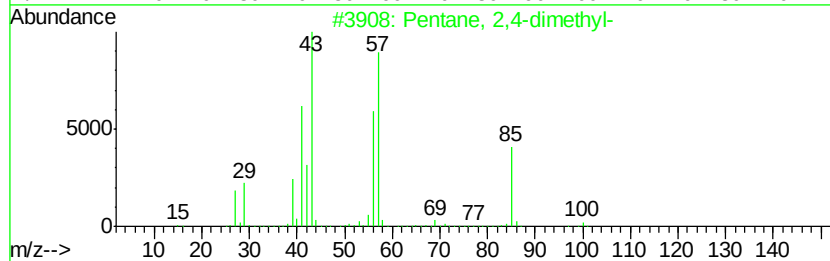
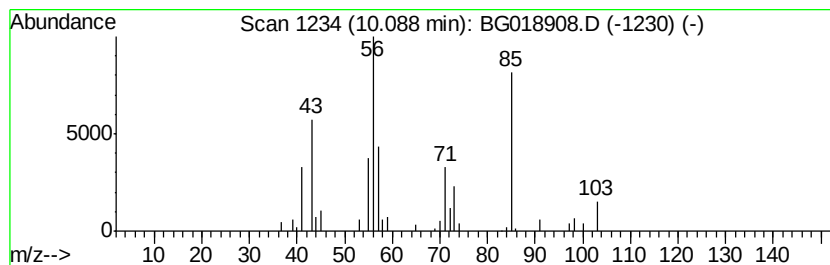
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.09	2.78 ng/ul	55755	Napthalene-d8	10.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	53
2		2(3H)-Furanone, dihydro-5-methyl-	100	C5H8O2	000108-29-2	40
3		Octane, 3,4-dimethyl-	142	C10H22	015869-92-8	38
4		2-Methoxy-2-methylbut-3-ene	100	C6H12O	040426-44-6	35
5		2(3H)-Furanone, dihydro-5-methyl-	100	C5H8O2	000108-29-2	33



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
Operator : UM/NP
Sample : G3797-02
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
D9G27

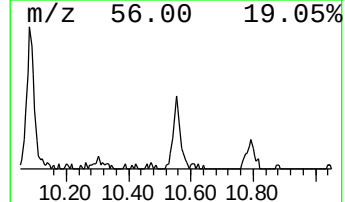
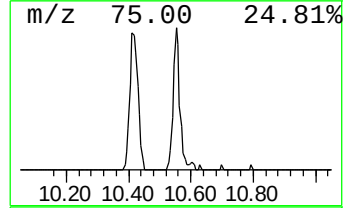
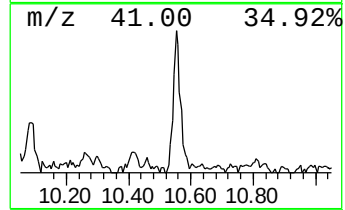
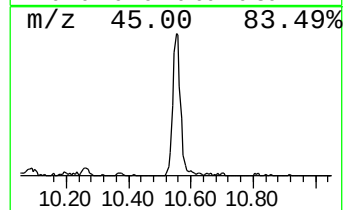
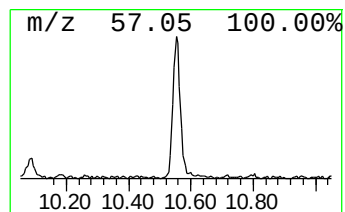
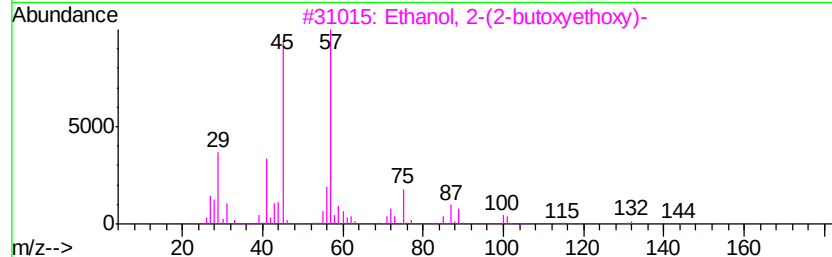
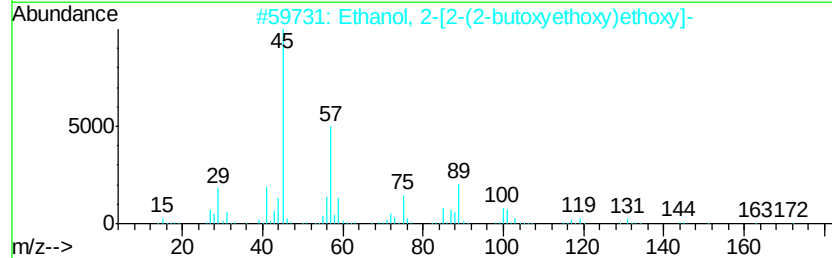
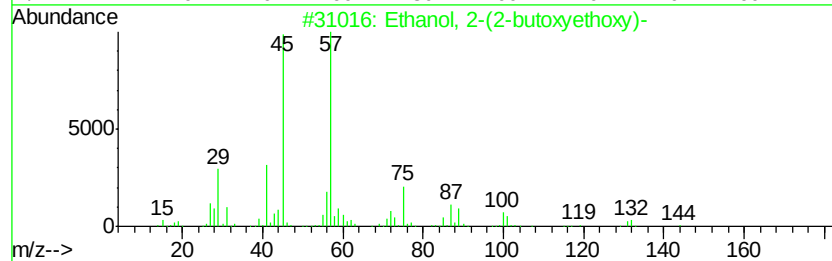
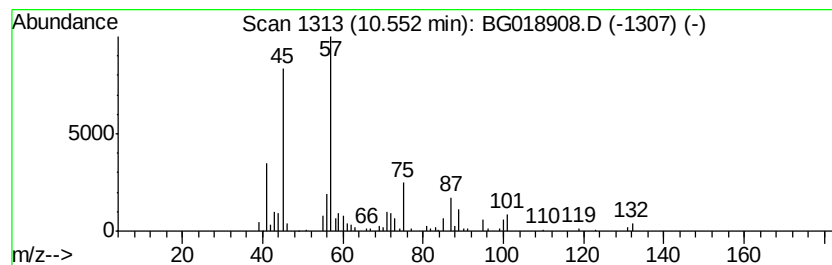
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.55	6.88 ng/ul	137858	Naphthalene-d8	10.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	80
2		Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	72
3		Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	64
4		Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	64
5		Propanoic acid, 2-hydroxy-, 2-me...	146	C7H14O3	000585-24-0	59



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
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Misc :
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Instrument :
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ClientSampleId :
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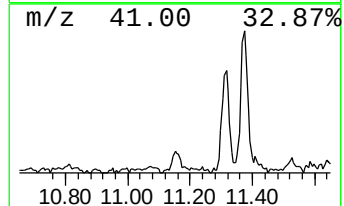
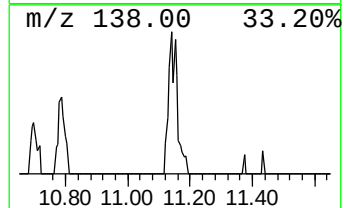
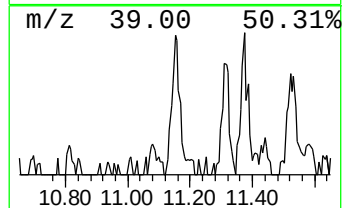
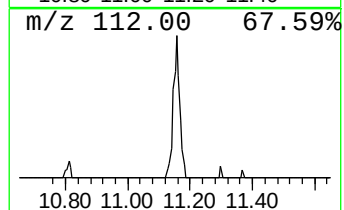
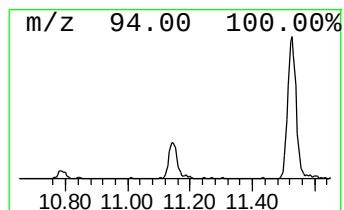
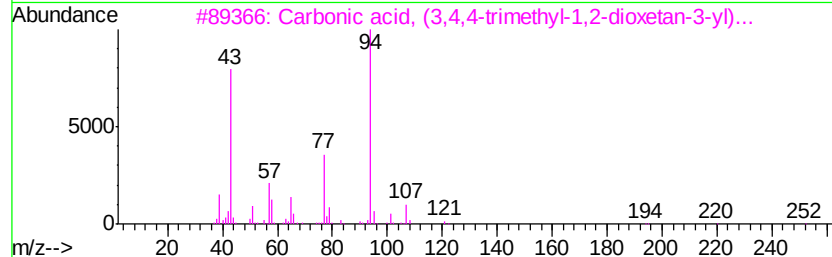
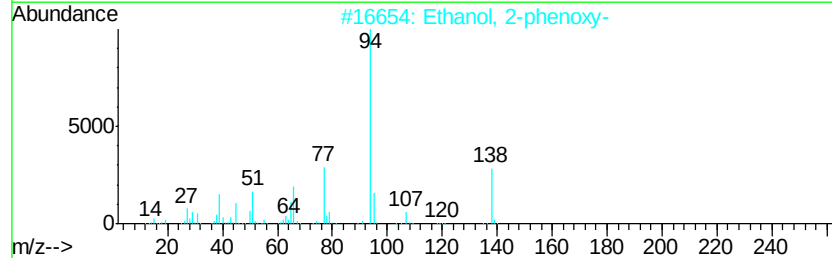
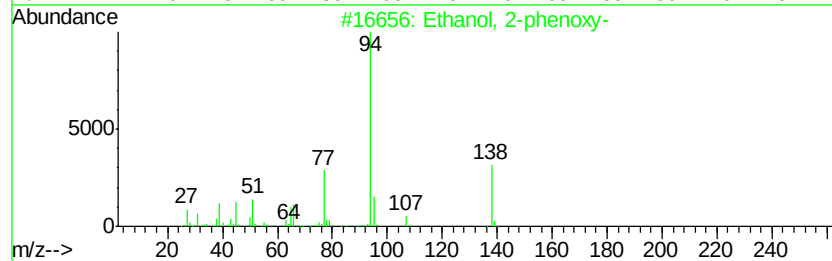
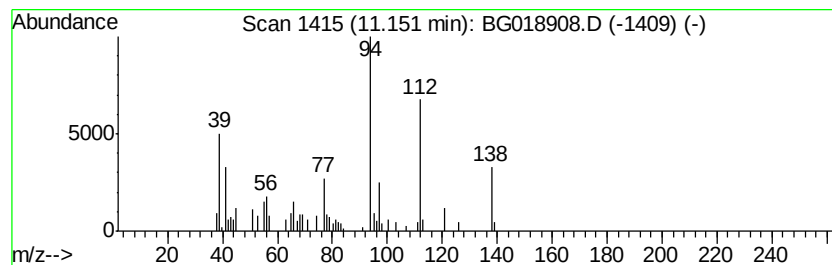
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 unknown-03 Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.15	2.19 ng/ul	43978	Napthalene-d8	10.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	43
2		Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	43
3		Carbonic acid, (3,4,4-trimethyl-...	252	C13H16O5	109123-69-5	35
4		3-Methylpyridazine	94	C5H6N2	001632-76-4	32
5		2-Aminopyridine	94	C5H6N2	000504-29-0	30



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
Operator : UM/NP
Sample : G3797-02
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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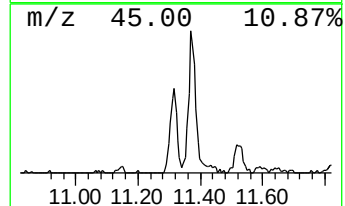
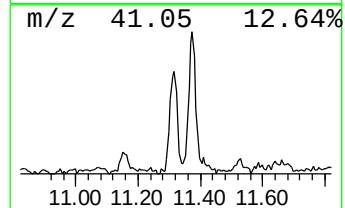
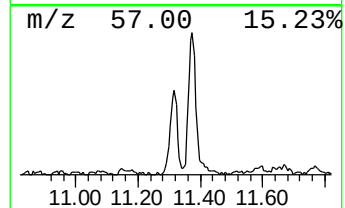
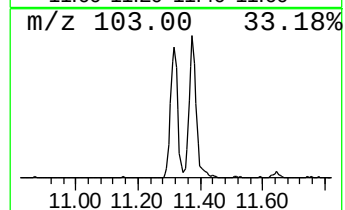
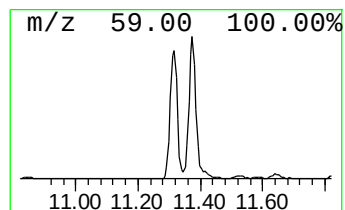
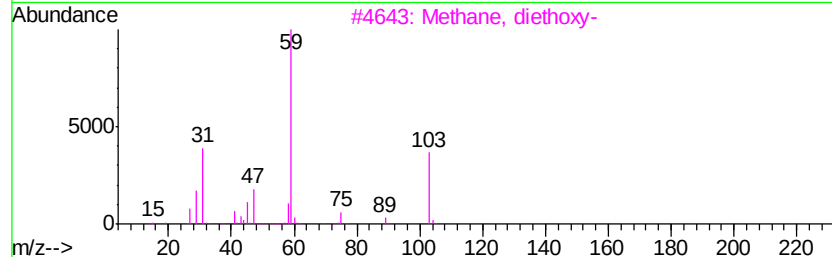
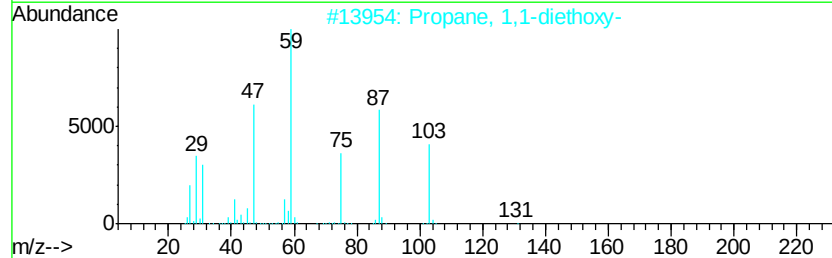
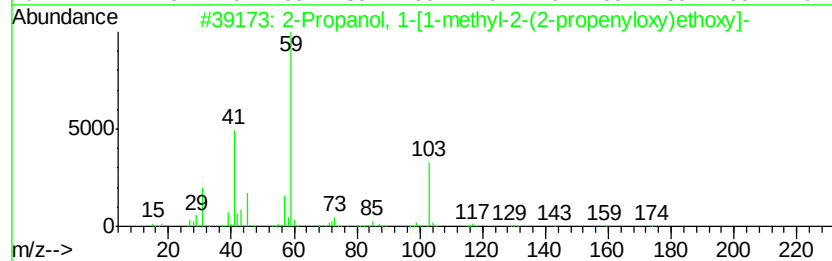
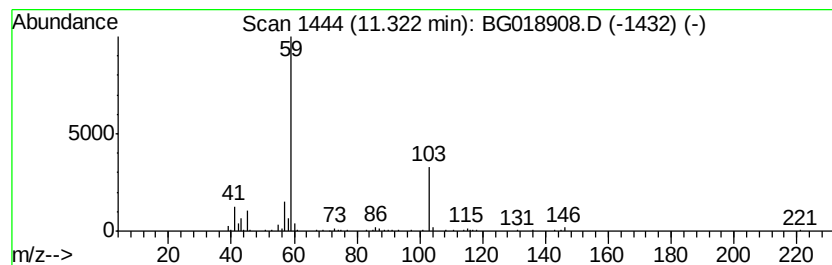
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 2-Propanol, 1-[1-methyl-2-(... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.32	9.96 ng/ul	199666	Napthalene-d8	10.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	83
2		Propane, 1,1-diethoxy-	132	C7H16O2	004744-08-5	83
3		Methane, diethoxy-	104	C5H12O2	000462-95-3	72
4		1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72
5		1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
Operator : UM/NP
Sample : G3797-02
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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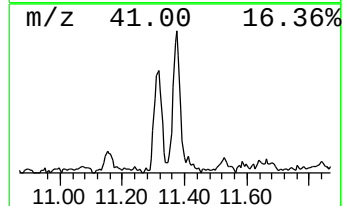
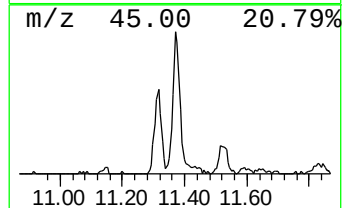
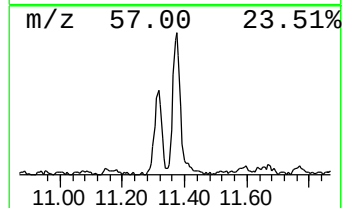
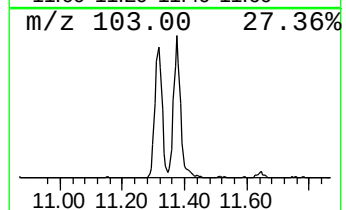
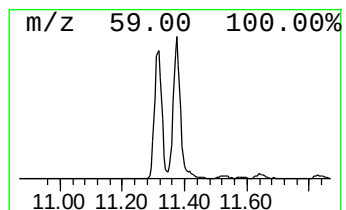
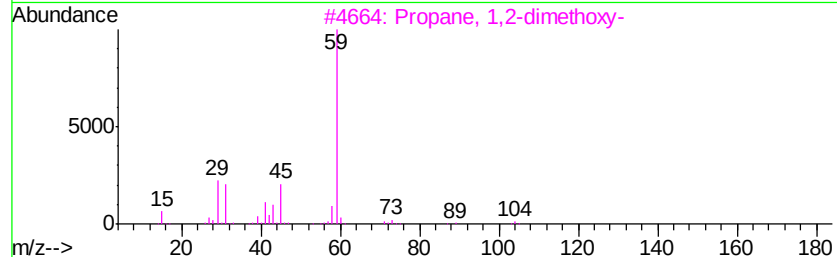
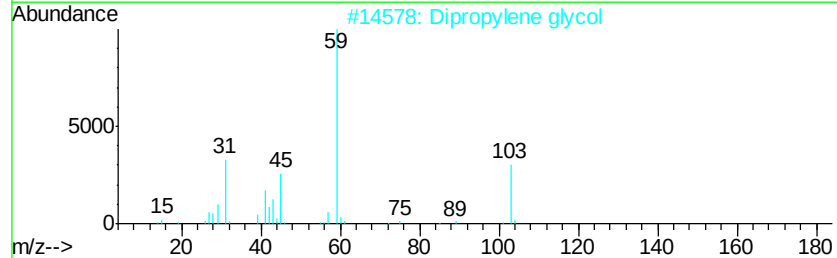
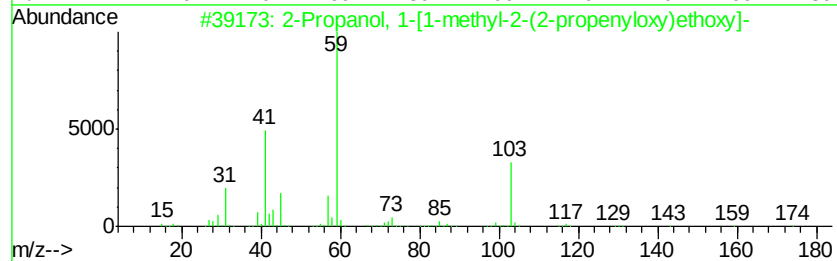
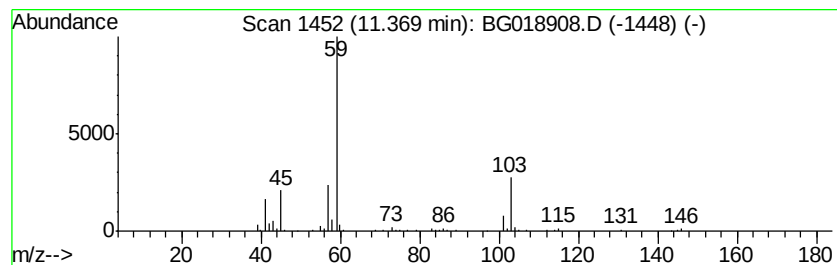
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Dipropylene glycol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.37	11.73 ng/ul	235062	Naphthalene-d8	10.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	72
2		Dipropylene glycol	134	C6H14O3	025265-71-8	72
3		Propane, 1,2-dimethoxy-	104	C5H12O2	007778-85-0	58
4		2-Propanol, 1,1'-[(1-methyl-1,2-...	192	C9H20O4	001638-16-0	56
5		2-Propanol, 1,1'-[(1-methyl-1,2-...	192	C9H20O4	001638-16-0	56



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
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Sample : G3797-02
Misc :
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Instrument :
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ClientSampleId :
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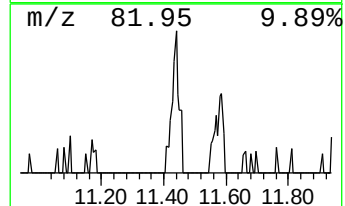
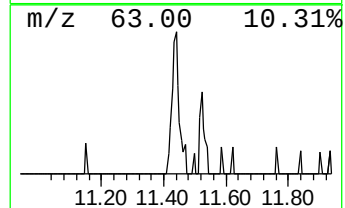
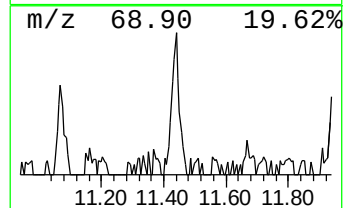
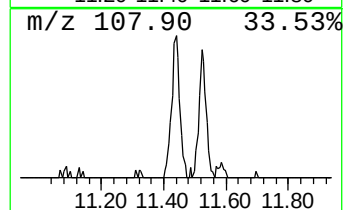
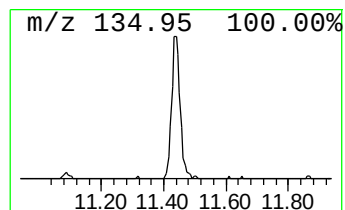
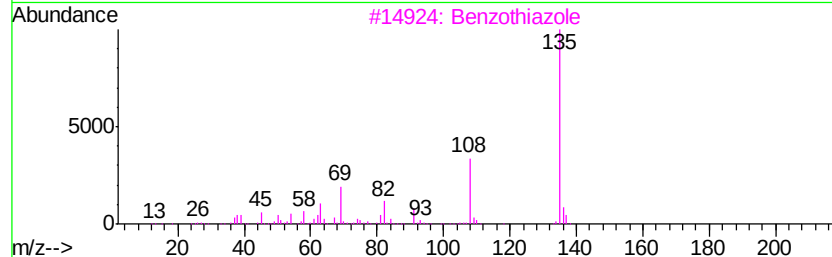
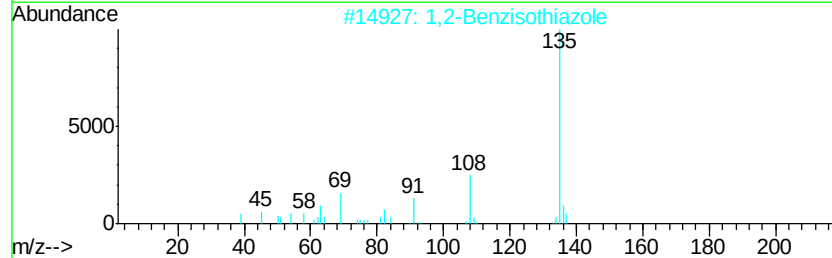
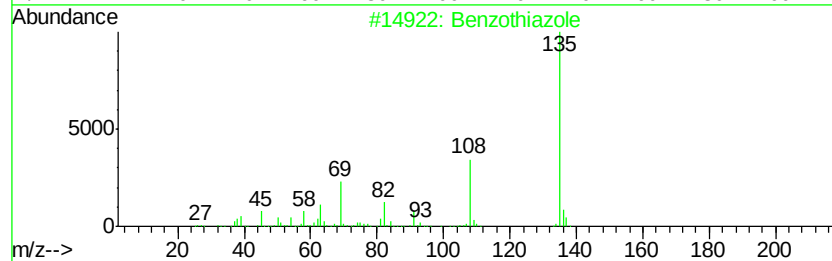
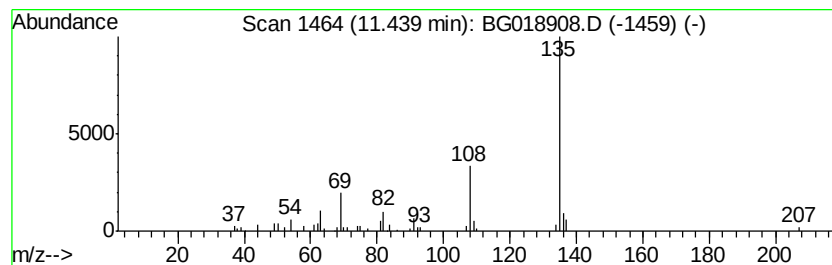
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Benzothiazole Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.44	3.10 ng/ul	62236	Napthalene-d8	10.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzothiazole	135	C7H5NS	000095-16-9	94
2			1,2-Benzisothiazole	135	C7H5NS	000272-16-2	91
3			Benzothiazole	135	C7H5NS	000095-16-9	91
4			Benzothiazole	135	C7H5NS	000095-16-9	90
5			Thieno[2,3-c]pyridine	135	C7H5NS	000272-12-8	72



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
Operator : UM/NP
Sample : G3797-02
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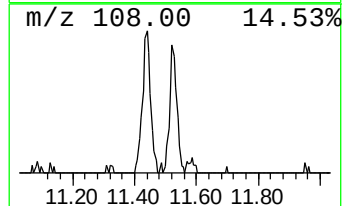
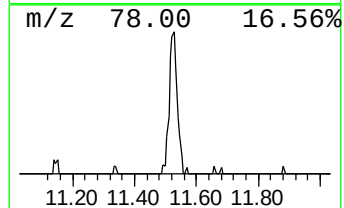
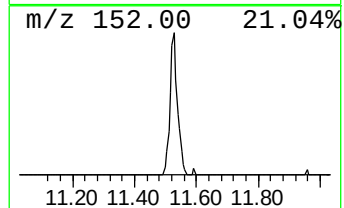
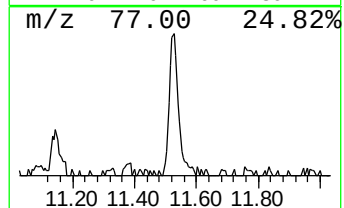
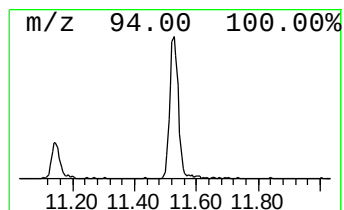
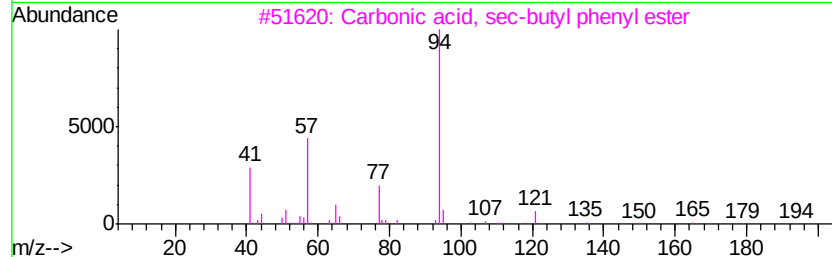
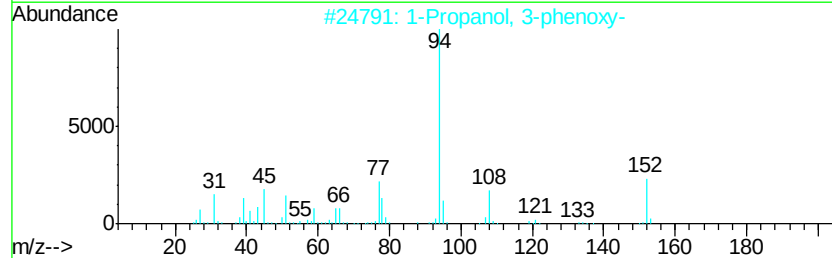
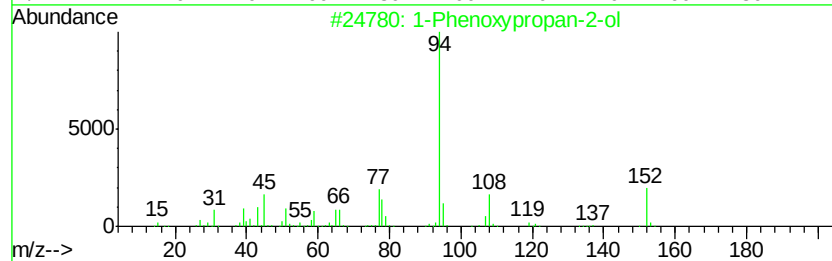
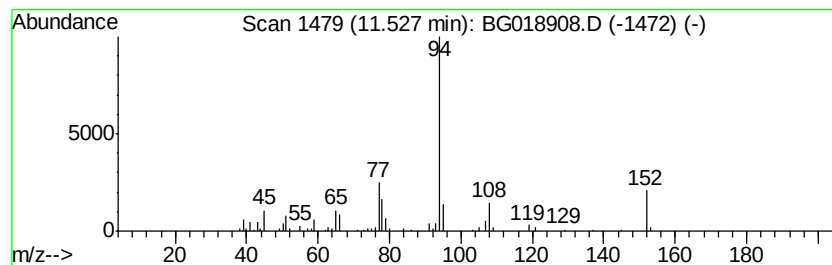
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 1-Phenoxypropan-2-ol Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.53	5.77 ng/ul	115647	Napthalene-d8	10.79
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1-Phenoxypropan-2-ol	152 C9H12O2	000770-35-4 93
2		1-Propanol, 3-phenoxy-	152 C9H12O2	006180-61-6 87
3		Carbonic acid, sec-butyl phenyl ...	194 C11H14O3	013183-17-0 59
4		1-Propanol, 2-phenoxy-	152 C9H12O2	004169-04-4 59
5		Carbonic acid, neopentyl phenyl ...	208 C12H16O3	013183-19-2 59



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
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Instrument :
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ClientSampleId :
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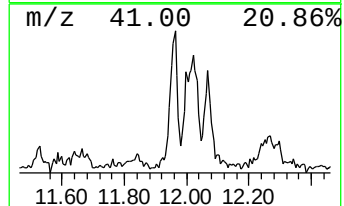
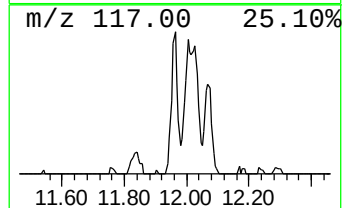
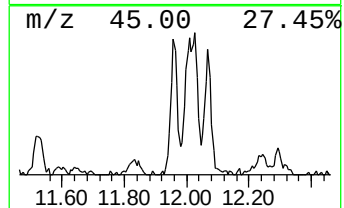
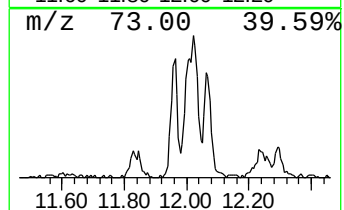
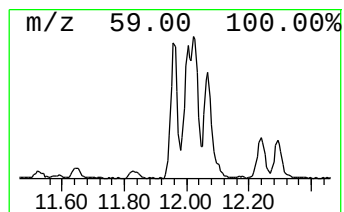
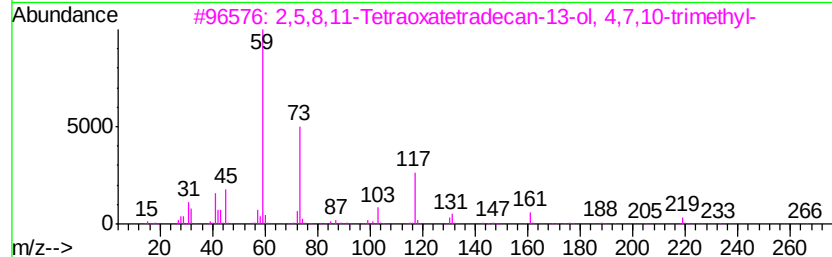
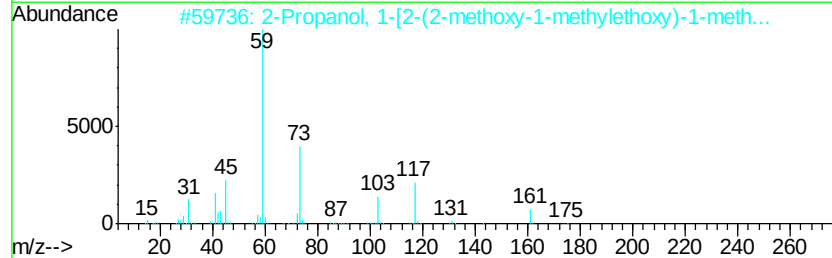
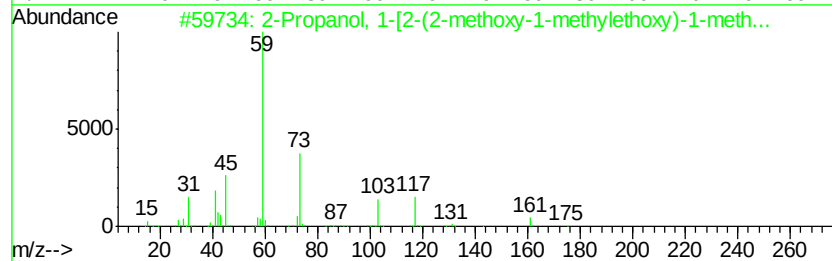
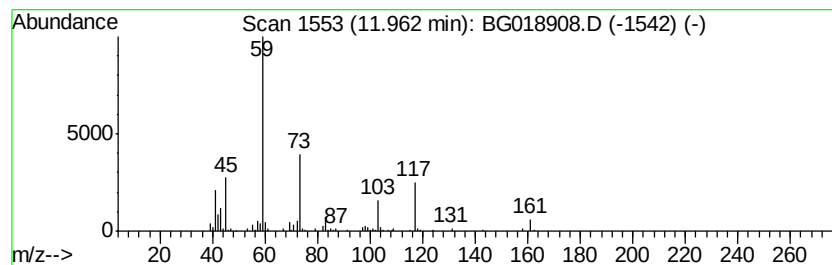
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 2-Propanol, 1-[2-(2-methoxy... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.96	7.86 ng/ul	157448	Naphthalene-d8	10.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	80
2			2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	72
3			2,5,8,11-Tetraoxatetradecan-13-o...	264	C13H28O5	020324-34-9	72
4			2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	72
5			Dipropylene glycol monomethyl ether	148	C7H16O3	034590-94-8	59



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
Operator : UM/NP
Sample : G3797-02
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
D9G27

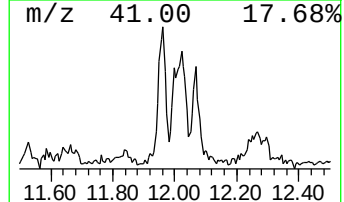
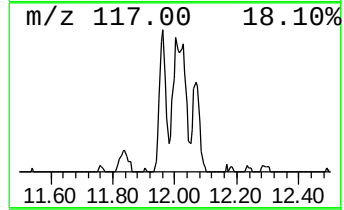
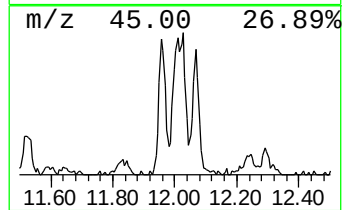
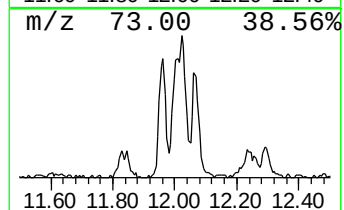
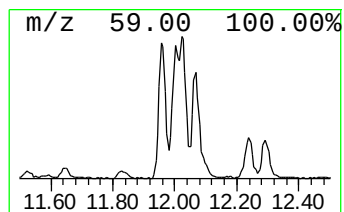
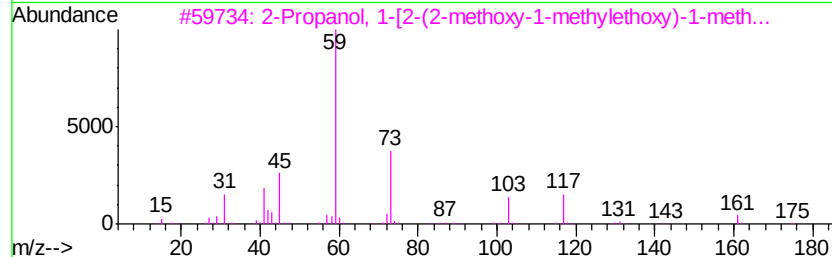
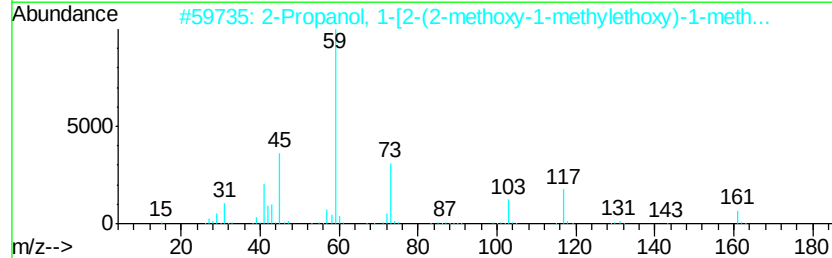
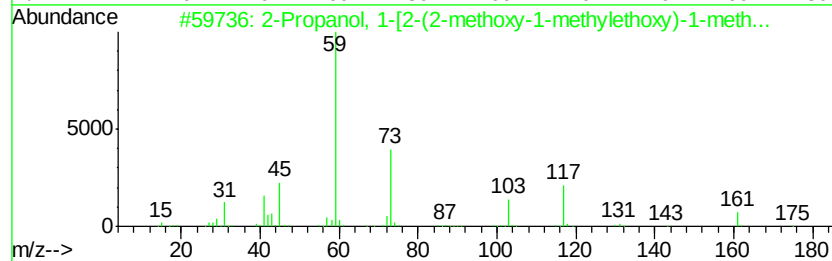
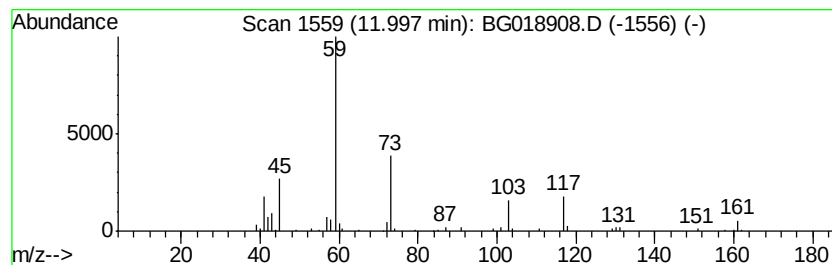
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 1-Propanol, 2-(2-methoxypro... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.00	5.20 ng/ul	104271	Napthalene-d8	10.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	90
2		2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	83
3		2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	78
4		1-Propanol, 2-(2-methoxypropoxy)-	148	C7H16O3	013588-28-8	64
5		Tri(1,2-propyleneglycol), monome...	206	C10H22O4	1000262-26-6	56



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
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Sample : G3797-02
Misc :
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Instrument :
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ClientSampleId :
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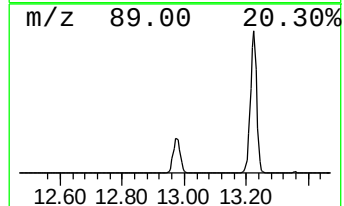
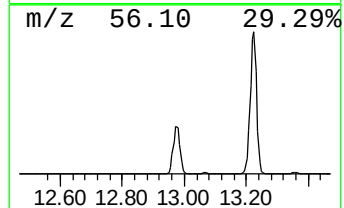
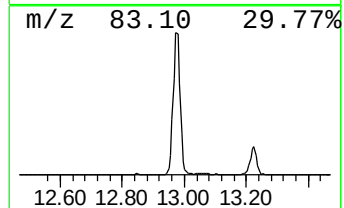
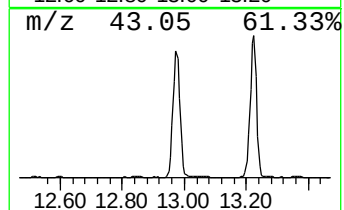
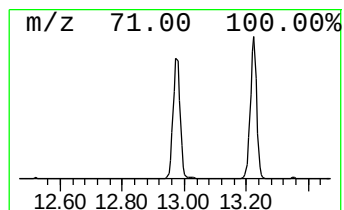
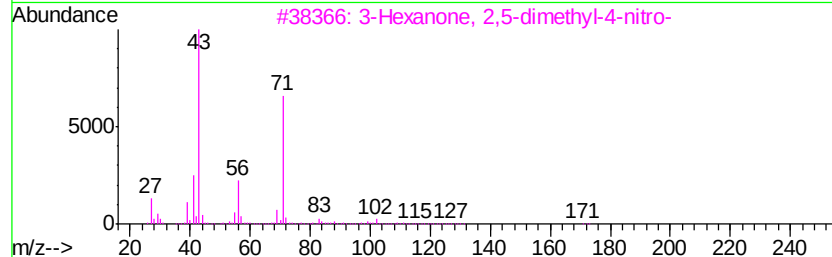
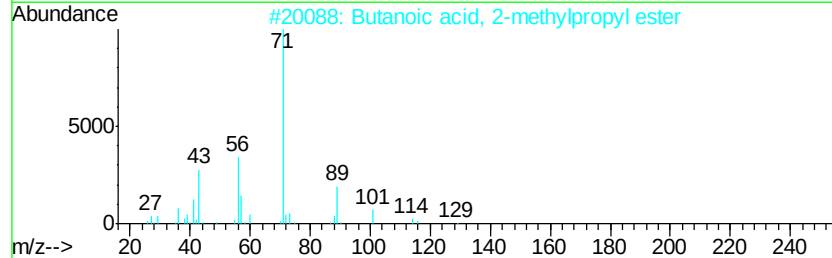
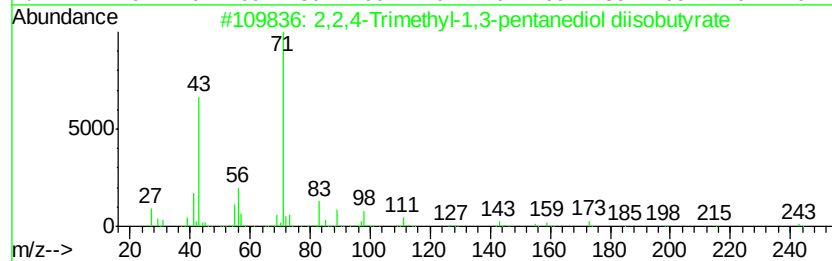
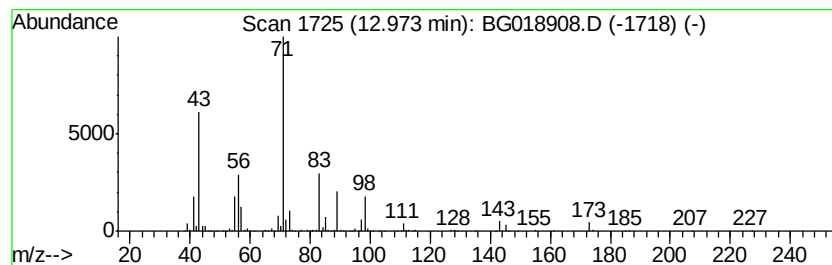
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 2,2,4-Trimethyl-1,3-pentane... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	49.83 ng/ul	1514710	Acenaphthene-d10	14.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2,4-Trimethyl-1,3-pentanediol ...	286	C16H30O4	006846-50-0	53
2		Butanoic acid, 2-methylpropyl ester	144	C8H16O2	000539-90-2	42
3		3-Hexanone, 2,5-dimethyl-4-nitro-	173	C8H15NO3	059906-54-6	38
4		Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	38
5		Propanoic acid, 2-methyl-, 2,2-d...	216	C12H24O3	074367-33-2	37



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
Operator : UM/NP
Sample : G3797-02
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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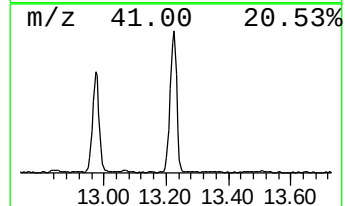
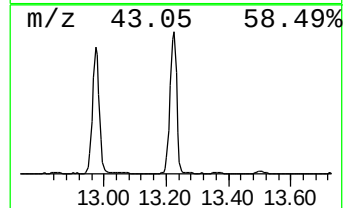
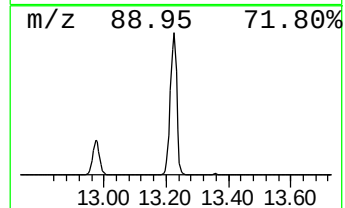
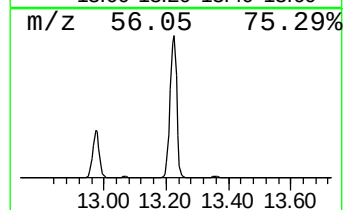
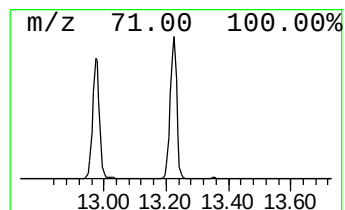
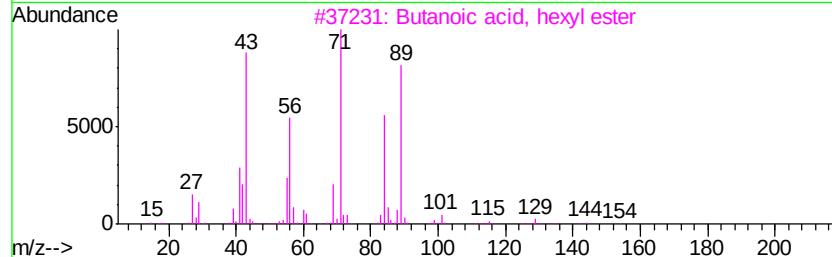
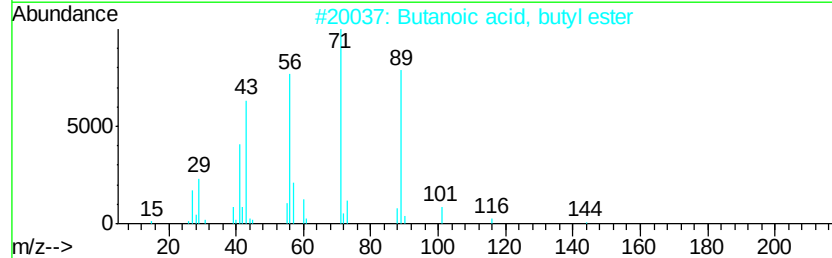
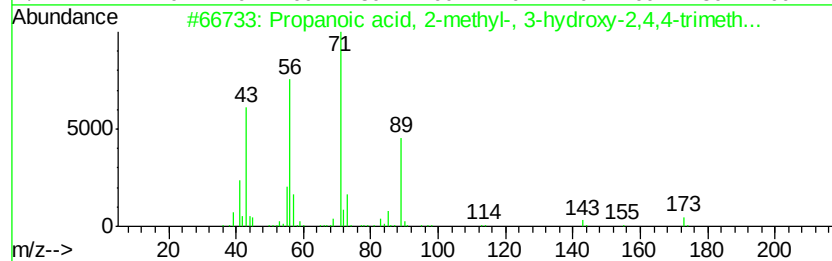
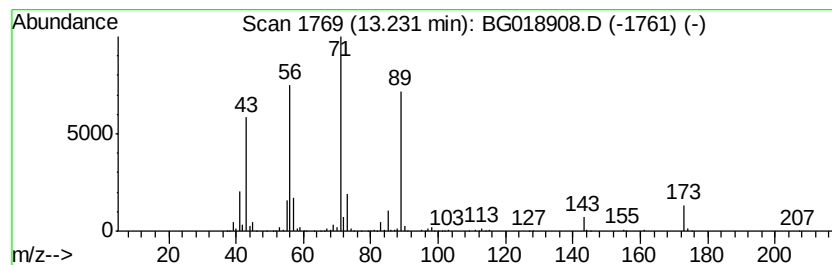
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 Propanoic acid, 2-methyl-, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.23	63.98 ng/ul	1944750	Acenaphthene-d10	14.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propanoic acid, 2-methyl-, 3-hyd...	216	C12H24O3	074367-34-3	86
2		Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	83
3		Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	64
4		Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	50
5		Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	50



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
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Sample : G3797-02
Misc :
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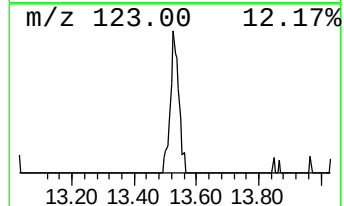
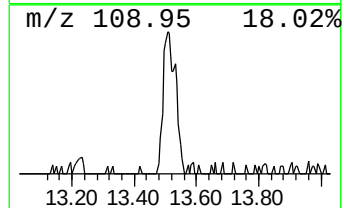
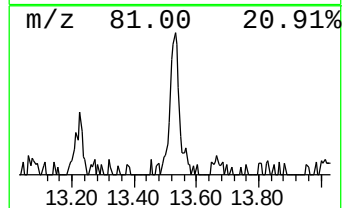
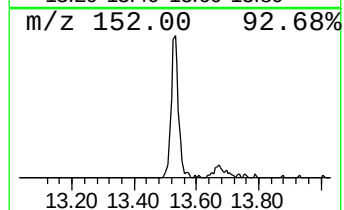
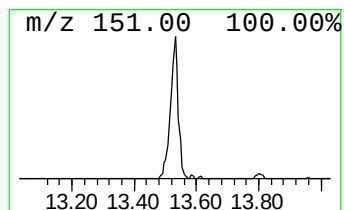
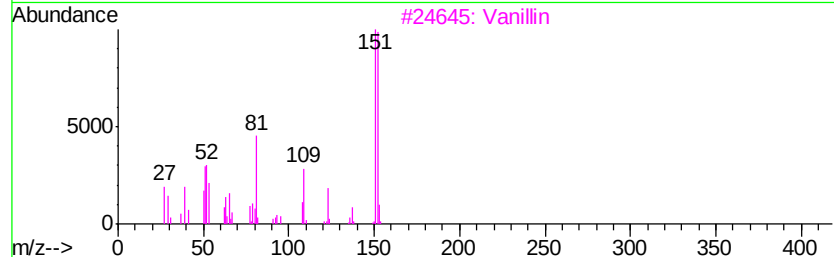
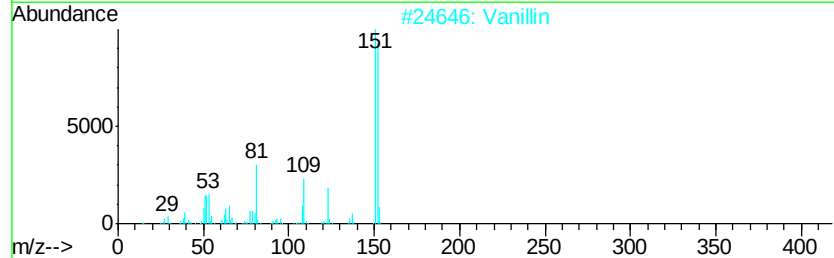
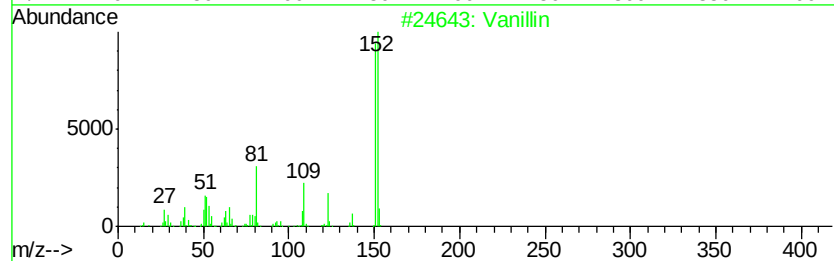
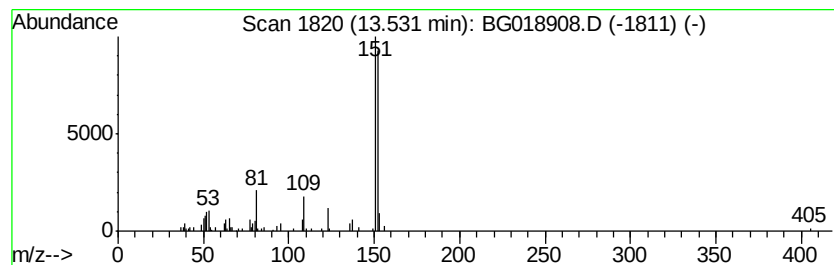
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 Vanillin Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.53	3.11 ng/ul	94408	Acenaphthene-d10	14.61

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Vanillin			152	C8H8O3	000121-33-5	94
2	Vanillin			152	C8H8O3	000121-33-5	94
3	Vanillin			152	C8H8O3	000121-33-5	93
4	Vanillin			152	C8H8O3	000121-33-5	93
5	Benzaldehyde, 3-hydroxy-4-methoxy-			152	C8H8O3	000621-59-0	93



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
Operator : UM/NP
Sample : G3797-02
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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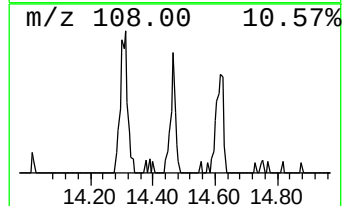
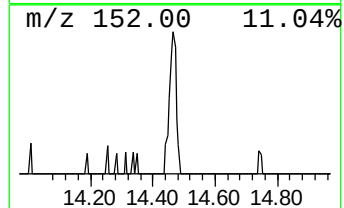
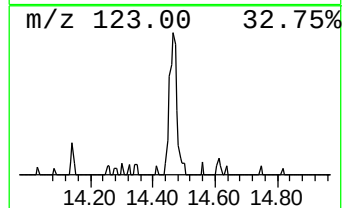
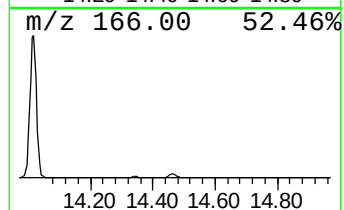
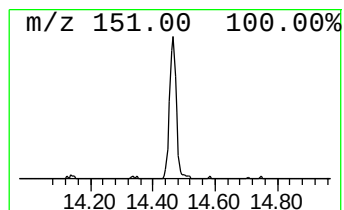
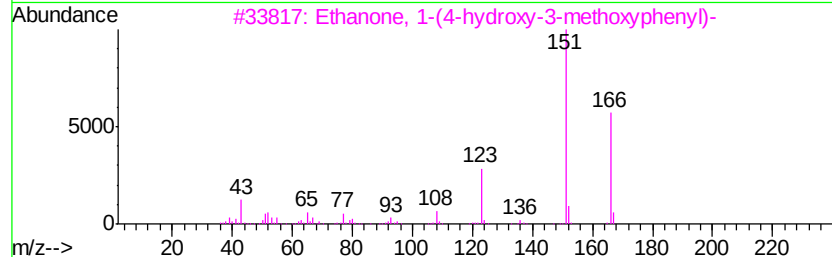
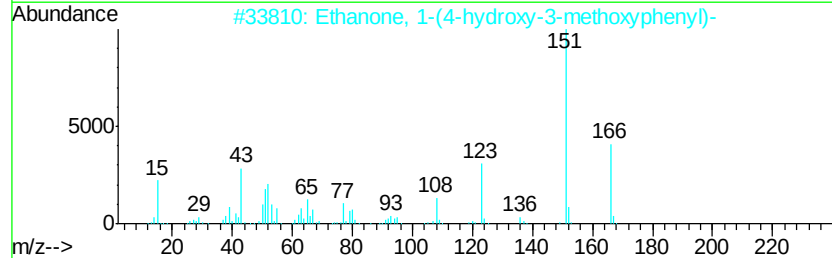
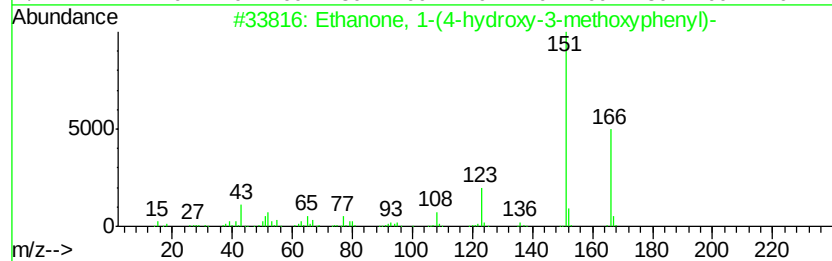
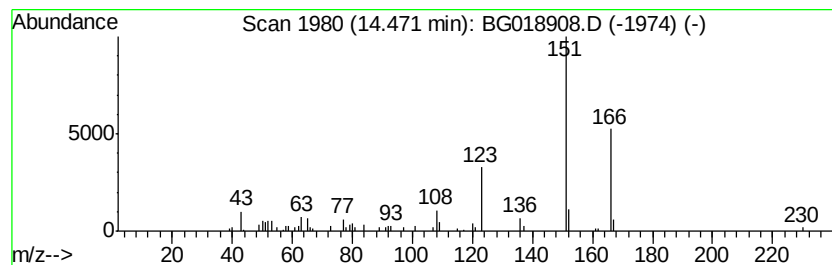
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 Ethanone, 1-(4-hydroxy-3-me... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.47	2.37 ng/ul	72125	Acenaphthene-d10	14.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanone, 1-(4-hydroxy-3-methoxy...	166	C9H10O3	000498-02-2	91
2		Ethanone, 1-(4-hydroxy-3-methoxy...	166	C9H10O3	000498-02-2	91
3		Ethanone, 1-(4-hydroxy-3-methoxy...	166	C9H10O3	000498-02-2	90
4		Ethanone, 1-(3-hydroxy-4-methoxy...	166	C9H10O3	006100-74-9	90
5		Ethanone, 1-(4-hydroxy-3-methoxy...	166	C9H10O3	000498-02-2	90



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
Operator : UM/NP
Sample : G3797-02
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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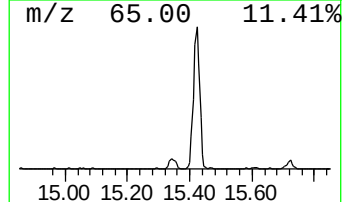
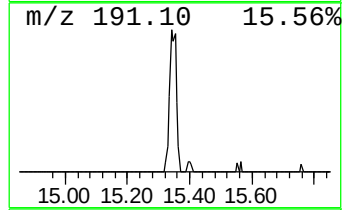
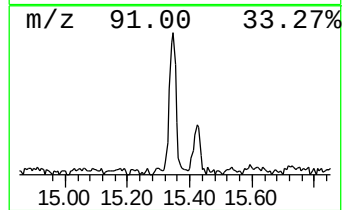
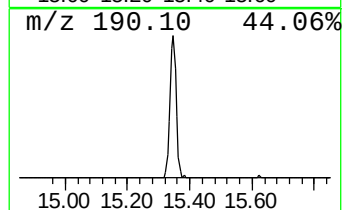
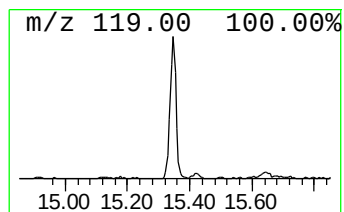
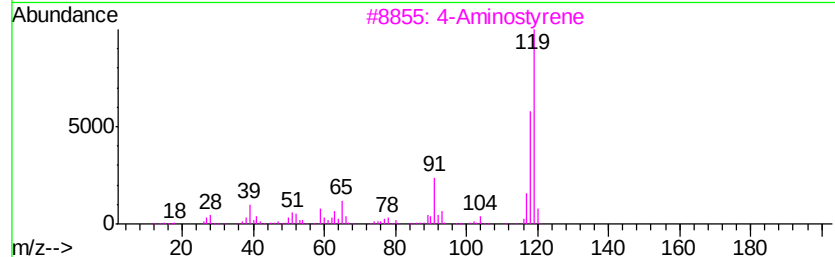
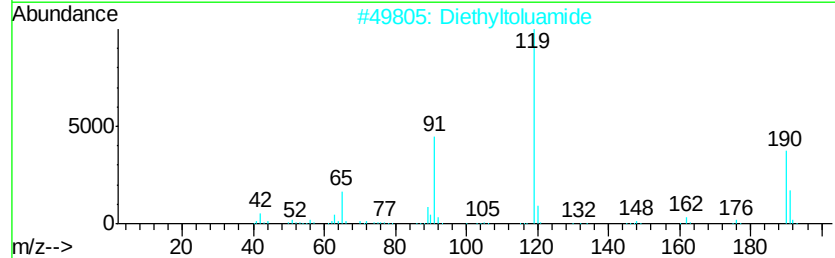
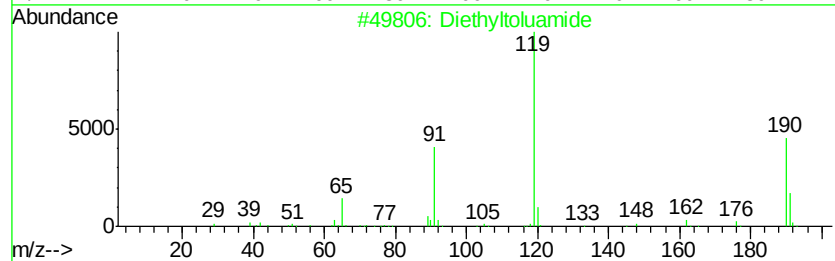
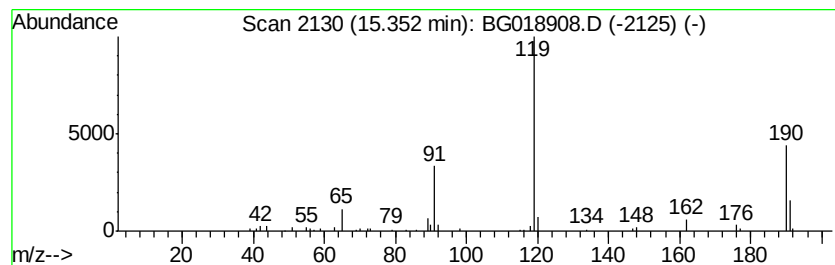
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 Diethyltoluamide Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.35	2.66 ng/ul	80873	Acenaphthene-d10	14.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diethyltoluamide	191	C12H17NO	000134-62-3	93
2		Diethyltoluamide	191	C12H17NO	000134-62-3	72
3		4-Aminostyrene	119	C8H9N	001520-21-4	52
4		Benzoxazol, 2,3-dihydro-2-thioxo...	269	C15H11N02S	037442-06-1	50
5		3-Buten-1-one, 1-(4-methylphenyl)-	160	C11H12O	130649-66-0	50



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
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Sample : G3797-02
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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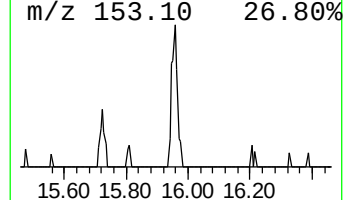
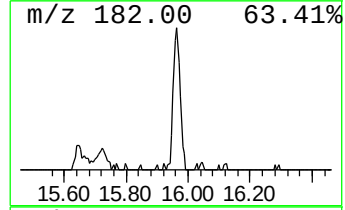
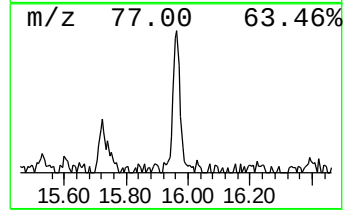
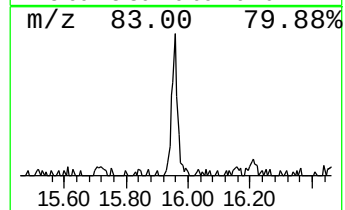
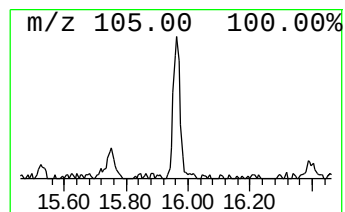
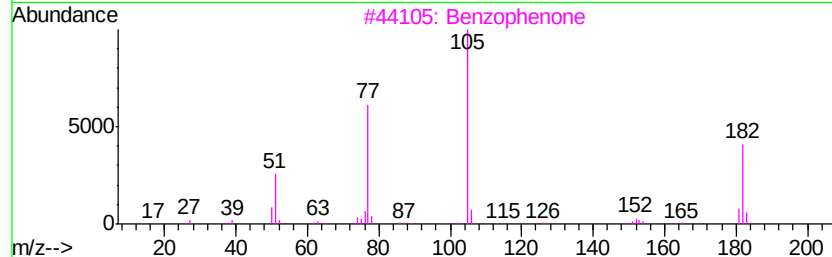
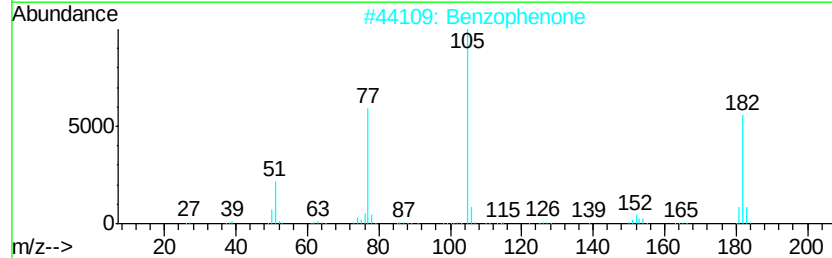
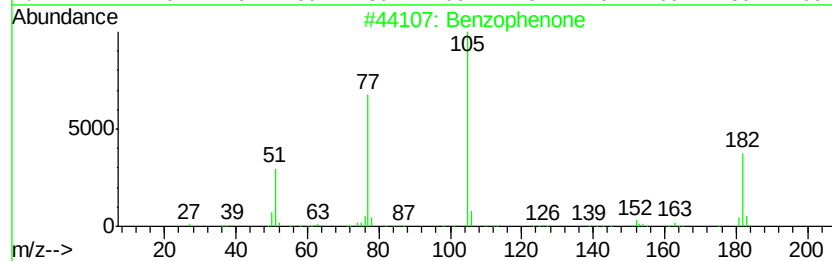
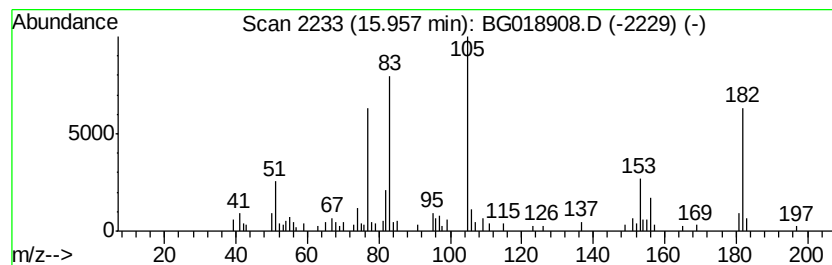
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Quant Title : SVOA CALIBRATION

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

Peak Number 29 Benzophenone Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.96	2.24 ng/ul	68098	Acenaphthene-d10	14.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzophenone	182	C13H10O	000119-61-9	50
2		Benzophenone	182	C13H10O	000119-61-9	43
3		Benzophenone	182	C13H10O	000119-61-9	43
4		Benzophenone	182	C13H10O	000119-61-9	43
5		Benzophenone	182	C13H10O	000119-61-9	38



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
Data File : BG018908.D
Acq On : 26 Sep 2015 12:07
Operator : UM/NP
Sample : G3797-02
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
D9G27

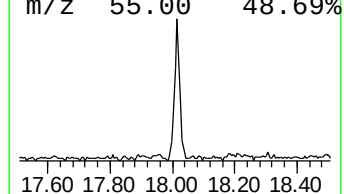
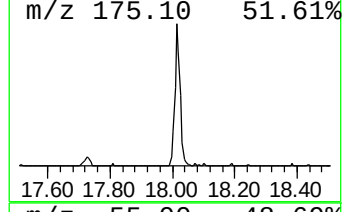
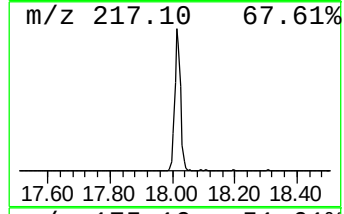
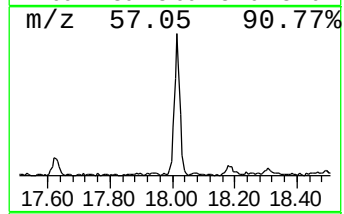
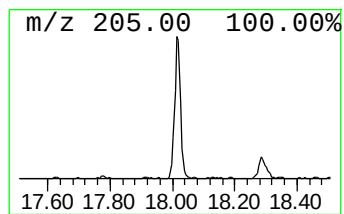
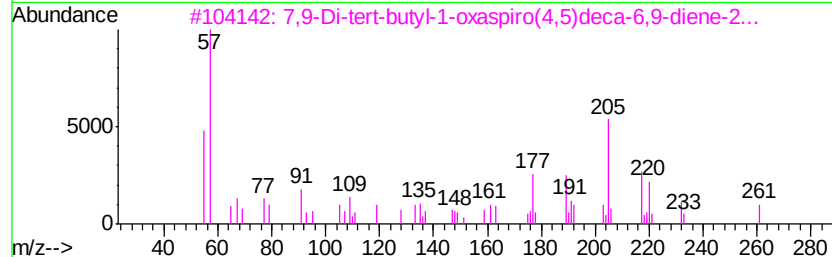
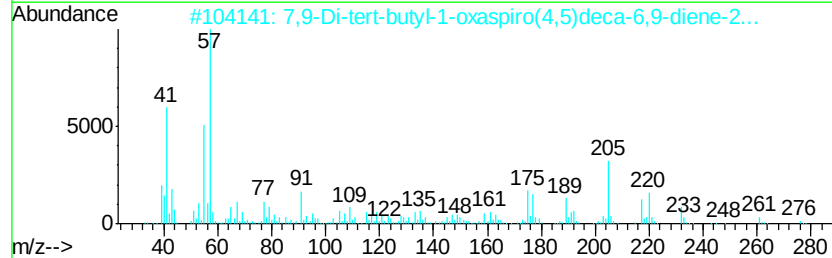
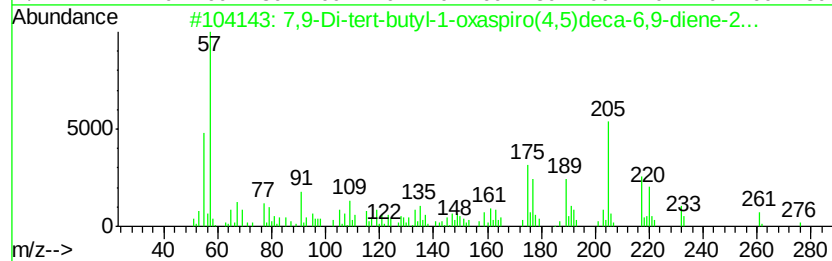
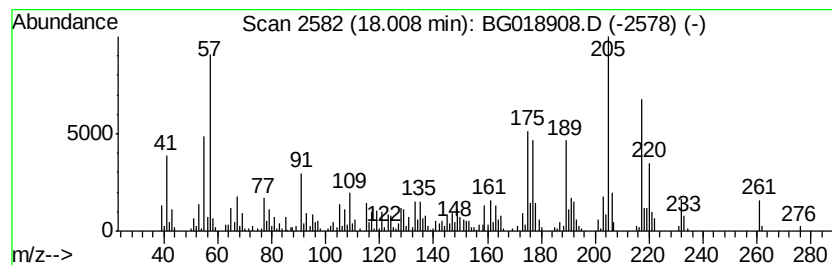
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG091015.M
Quant Title : SVOA CALIBRATION

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

Peak Number 30 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.01	10.84 ng/ul	446126	Phenanthrene-d10	17.35

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	98
2			7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	83
3			7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	76
4			2,5-di-tert-Butyl-1,4-benzoquinone	220	C14H2O2	002460-77-7	53
5			Benzenamine, N,N-diethyl-4-(2-ni...	220	C12H16N2O2	064462-51-7	30



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092515\
 Data File : BG018908.D
 Acq On : 26 Sep 2015 12:07
 Operator : UM/NP
 Sample : G3797-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 D9G27

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...	3.17	17.8	ng/ul	237055	1	7.99	266864	20.0
Ethanol, 2-butoxy-	6.06	8.3	ng/ul	110297	1	7.99	266864	20.0
2-Propanol, 1-but...	6.63	2.8	ng/ul	37362	1	7.99	266864	20.0
2-Propanol, 1-(2-...	7.76	14.3	ng/ul	190630	1	7.99	266864	20.0
Mequinol	9.08	3.0	ng/ul	39620	1	7.99	266864	20.0
unknown-01	9.31	4.2	ng/ul	55745	1	7.99	266864	20.0
Phenylethyl Alcohol	9.53	3.1	ng/ul	61416	2	10.79	400880	20.0
2-Propanol, 1,1'-...	9.96	9.6	ng/ul	192526	2	10.79	400880	20.0
unknown-02	10.03	6.5	ng/ul	130215	2	10.79	400880	20.0
(DEL) Alkane: Str...	10.09	2.8	ng/ul	55755	2	10.79	400880	20.0
Ethanol, 2-(2-but...	10.55	6.9	ng/ul	137858	2	10.79	400880	20.0
unknown-03	11.15	2.2	ng/ul	43978	2	10.79	400880	20.0
2-Propanol, 1-[1-...	11.32	10.0	ng/ul	199666	2	10.79	400880	20.0
Dipropylene glycol	11.37	11.7	ng/ul	235062	2	10.79	400880	20.0
Benzothiazole	11.44	3.1	ng/ul	62236	2	10.79	400880	20.0
1-Phenoxypropan-2-ol	11.53	5.8	ng/ul	115647	2	10.79	400880	20.0
2-Propanol, 1-[2-...	11.96	7.9	ng/ul	157448	2	10.79	400880	20.0
1-Propanol, 2-(2-...	12.00	5.2	ng/ul	104271	2	10.79	400880	20.0
Tri(1,2-propylene...	12.03	4.0	ng/ul	79583	2	10.79	400880	20.0
2,2,4-Trimethyl-1...	12.97	49.8	ng/ul	1514710	3	14.61	607961	20.0
Propanoic acid, 2...	13.23	64.0	ng/ul	1944750	3	14.61	607961	20.0
Vanillin	13.53	3.1	ng/ul	94408	3	14.61	607961	20.0
Ethanone, 1-(4-hy...	14.47	2.4	ng/ul	72125	3	14.61	607961	20.0
Diethyltoluamide	15.35	2.7	ng/ul	80873	3	14.61	607961	20.0
Benzophenone	15.96	2.2	ng/ul	68098	3	14.61	607961	20.0
7,9-Di-tert-butyl...	18.01	10.8	ng/ul	446126	4	17.35	823218	20.0