

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122015\
 Data File : BG020341.D
 Acq On : 20 Dec 2015 14:13
 Operator : UM/SJ
 Sample : G4803-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C04D9

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-BG121415.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.114	41	47	56	rBV	45878	87593	12.02%	1.227%
2	3.191	56	60	73	rVB	34842	53312	7.32%	0.747%
3	5.171	392	397	402	rBB	3361	4895	0.67%	0.069%
4	5.705	483	488	492	rBV2	1768	2984	0.41%	0.042%
5	6.081	546	552	557	rBB	3058	4563	0.63%	0.064%
6	7.245	744	750	755	rBV	23054	38303	5.26%	0.537%
7	7.304	757	760	764	rVB3	3715	4944	0.68%	0.069%
8	7.409	771	778	786	rVB	71302	113183	15.54%	1.586%
9	7.621	807	814	823	rBV	78119	132872	18.24%	1.862%
10	8.091	887	894	901	rBV	66585	110221	15.13%	1.544%
11	8.796	1005	1014	1021	rBV	70536	118103	16.21%	1.655%
12	9.260	1086	1093	1100	rVV	95557	168173	23.08%	2.357%
13	9.336	1103	1106	1109	rVV2	1158	2019	0.28%	0.028%
14	9.912	1196	1204	1207	rBV3	1477	2457	0.34%	0.034%
15	9.983	1210	1216	1225	rVB	84856	150923	20.72%	2.115%
16	10.271	1258	1265	1267	rBV5	3674	7137	0.98%	0.100%
17	10.312	1270	1272	1278	rVB3	3099	4800	0.66%	0.067%
18	10.523	1301	1308	1318	rBV	132772	228755	31.40%	3.205%
19	10.606	1318	1322	1326	rVB3	1936	2995	0.41%	0.042%
20	10.658	1326	1331	1337	rBV5	1711	3925	0.54%	0.055%
21	10.723	1337	1342	1343	rBV2	2256	2766	0.38%	0.039%
22	10.788	1343	1353	1361	rVB2	26065	70946	9.74%	0.994%
23	10.911	1366	1374	1386	rVB	118227	245329	33.67%	3.438%
24	11.035	1388	1395	1402	rBV3	7274	14065	1.93%	0.197%
25	11.140	1409	1413	1418	rBV3	1910	3095	0.42%	0.043%
26	11.334	1440	1446	1452	rVB5	6803	13206	1.81%	0.185%
27	11.387	1452	1455	1460	rVB3	2025	3202	0.44%	0.045%
28	11.446	1460	1465	1466	rBV2	2205	3186	0.44%	0.045%
29	11.493	1470	1473	1478	rVB2	2302	2870	0.39%	0.040%
30	11.546	1478	1482	1487	rVB2	3229	4638	0.64%	0.065%
31	11.610	1487	1493	1501	rVB2	17041	30304	4.16%	0.425%
32	11.745	1513	1516	1517	rBV	1868	1965	0.27%	0.028%
33	11.810	1522	1527	1531	rVB5	2199	3754	0.52%	0.053%
34	11.881	1534	1539	1541	rBV3	4329	6721	0.92%	0.094%

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35	11.933	1545	1548	1551	rBV2	1426	2034	0.28%	0.029%
36	12.086	1570	1574	1580	rBV4	2496	5037	0.69%	0.071%
37	12.151	1580	1585	1590	rVB4	3085	5081	0.70%	0.071%
38	12.233	1590	1599	1604	rBV	12187	24805	3.40%	0.348%
39	12.268	1604	1605	1609	rVB3	4054	3753	0.52%	0.053%
40	12.327	1609	1615	1616	rBV3	5976	8621	1.18%	0.121%
41	12.351	1616	1619	1624	rVV5	6227	10775	1.48%	0.151%
42	12.427	1628	1632	1635	rBV4	2706	3411	0.47%	0.048%
43	12.498	1641	1644	1649	rVB6	3144	4627	0.64%	0.065%
44	12.556	1649	1654	1659	rBV5	4612	8272	1.14%	0.116%
45	12.650	1667	1670	1676	rVB4	9195	15212	2.09%	0.213%
46	12.768	1684	1690	1691	rBV4	1707	3142	0.43%	0.044%
47	12.827	1697	1700	1705	rVB6	2112	3837	0.53%	0.054%
48	13.003	1721	1730	1732	rBV7	4203	9500	1.30%	0.133%
49	13.103	1743	1747	1753	rVB7	3328	5371	0.74%	0.075%
50	13.203	1761	1764	1766	rVV3	1874	2109	0.29%	0.030%
51	13.326	1781	1785	1790	rVB4	2098	2979	0.41%	0.042%
52	13.385	1790	1795	1800	rBV7	3247	6089	0.84%	0.085%
53	13.449	1800	1806	1807	rBV4	1854	2620	0.36%	0.037%
54	13.508	1811	1816	1821	rVV6	7410	12945	1.78%	0.181%
55	13.655	1835	1841	1844	rBV6	4367	8797	1.21%	0.123%
56	13.767	1856	1860	1866	rVB4	6065	10214	1.40%	0.143%
57	13.819	1866	1869	1872	rBV4	1382	2315	0.32%	0.032%
58	13.855	1872	1875	1879	rVB4	1351	2071	0.28%	0.029%
59	13.943	1885	1890	1893	rBV6	2848	4864	0.67%	0.068%
60	13.978	1893	1896	1899	rVB4	4155	4780	0.66%	0.067%
61	14.043	1902	1907	1913	rBV6	4926	11104	1.52%	0.156%
62	14.119	1913	1920	1928	rVB	245075	359260	49.31%	5.034%
63	14.331	1953	1956	1960	rVV7	3591	5178	0.71%	0.073%
64	14.419	1964	1971	1980	rVB	267740	421823	57.90%	5.911%
65	14.536	1984	1991	1996	rBV8	3281	7447	1.02%	0.104%
66	14.630	2004	2007	2012	rBV5	3268	5269	0.72%	0.074%
67	14.724	2012	2023	2029	rBV2	160386	258457	35.48%	3.622%
68	14.789	2029	2034	2038	rVB4	5207	8240	1.13%	0.115%
69	14.918	2049	2056	2062	rBV2	23466	41269	5.66%	0.578%
70	15.071	2079	2082	2083	rBV3	2511	2568	0.35%	0.036%
71	15.089	2083	2085	2088	rVV4	4062	5656	0.78%	0.079%

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72	15.112	2088	2089	2095	rVB5	4870	5318	0.73%	0.075%
73	15.188	2097	2102	2107	rVV8	5387	10444	1.43%	0.146%
74	15.253	2111	2113	2117	rVB4	4100	4806	0.66%	0.067%
75	15.447	2143	2146	2147	rBV3	3540	3950	0.54%	0.055%
76	15.512	2152	2157	2162	rVV2	23629	40276	5.53%	0.564%
77	15.553	2162	2164	2167	rVV4	2728	3382	0.46%	0.047%
78	15.600	2170	2172	2175	rVB4	3338	3484	0.48%	0.049%
79	15.659	2180	2182	2185	rBV4	2451	2487	0.34%	0.035%
80	15.711	2185	2191	2201	rVB	369157	548768	75.32%	7.690%
81	15.823	2204	2210	2217	rBV	117535	173035	23.75%	2.425%
82	15.946	2228	2231	2236	rBV5	5906	9909	1.36%	0.139%
83	16.228	2277	2279	2282	rBV4	2353	2424	0.33%	0.034%
84	16.481	2320	2322	2325	rVB3	2729	2743	0.38%	0.038%
85	16.792	2373	2375	2380	rVB5	4436	5908	0.81%	0.083%
86	17.039	2412	2417	2423	rBV3	26497	45205	6.20%	0.633%
87	17.468	2484	2490	2496	rBV2	199985	309053	42.42%	4.331%
88	17.568	2500	2507	2515	rVB	380347	583008	80.02%	8.169%
89	17.727	2531	2534	2539	rBV6	4858	9568	1.31%	0.134%
90	18.338	2635	2638	2642	rVB4	13153	15105	2.07%	0.212%
91	18.631	2684	2688	2695	rBV	40956	54900	7.54%	0.769%
92	19.072	2760	2763	2768	rVB2	20118	22869	3.14%	0.320%
93	19.607	2851	2854	2862	rVB9	19459	36456	5.00%	0.511%
94	19.848	2889	2895	2900	rBV	480064	696331	95.58%	9.757%
95	21.740	3212	3217	3225	rVB	231915	391241	53.70%	5.482%
96	23.226	3466	3470	3477	rVB3	31915	65736	9.02%	0.921%
97	24.795	3727	3737	3748	rVB	255503	728547	100.00%	10.209%
98	25.018	3765	3775	3785	rVB	139835	408356	56.05%	5.722%
99	28.015	4283	4285	4286	rBV2	3589	2001	0.27%	0.028%
100	31.416	4862	4864	4865	rBV	3663	3467	0.48%	0.049%

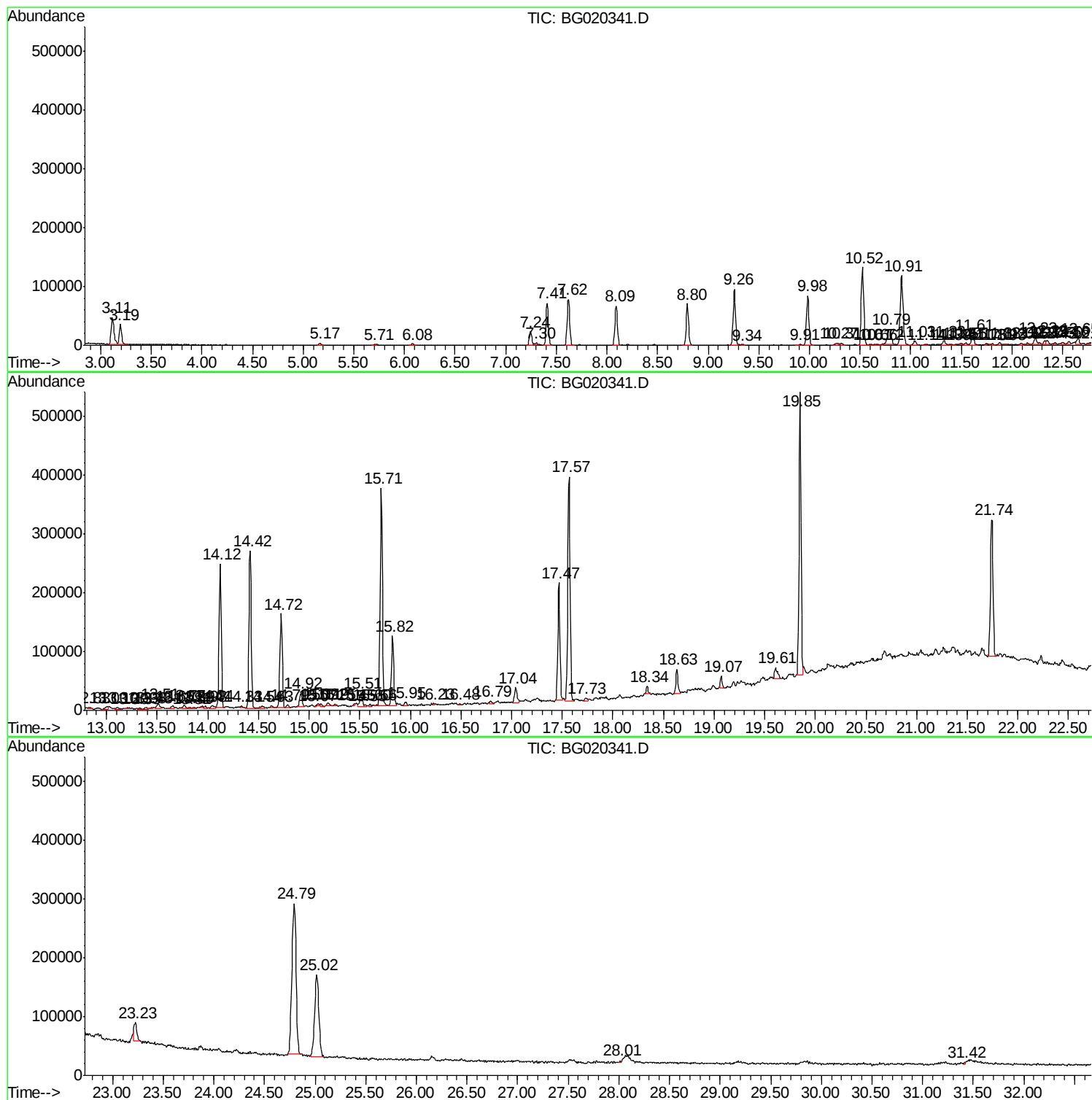
Sum of corrected areas: 7136513

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

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



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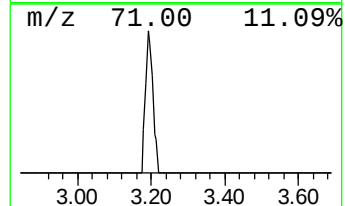
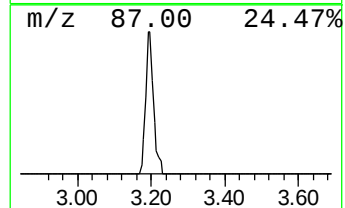
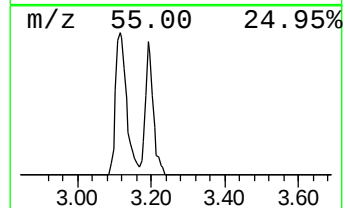
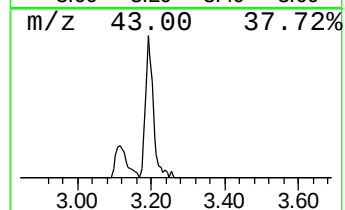
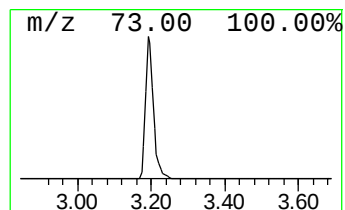
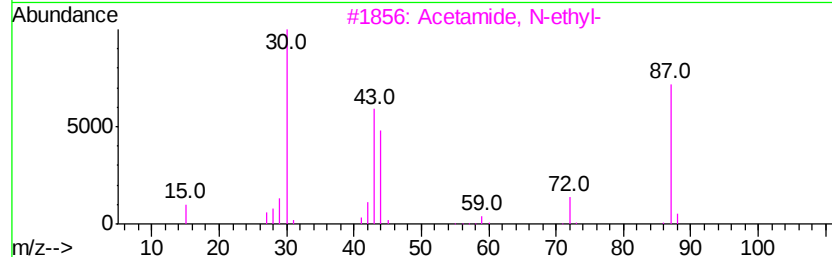
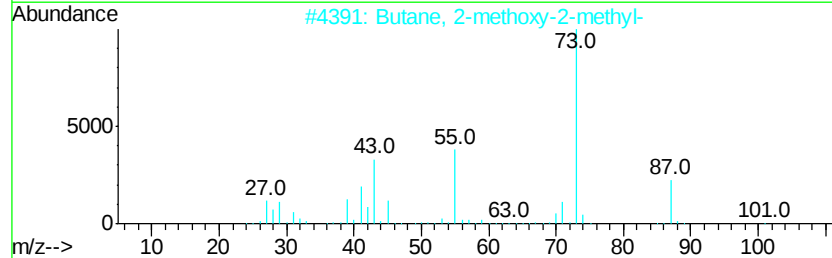
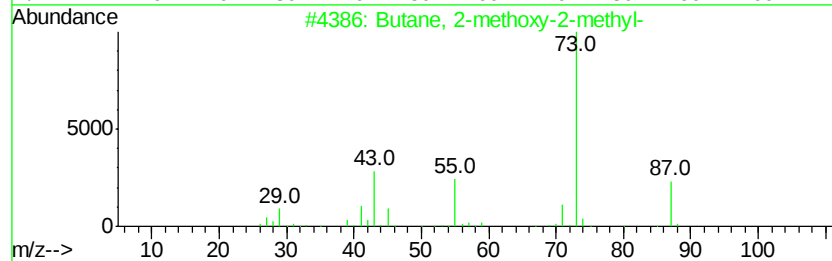
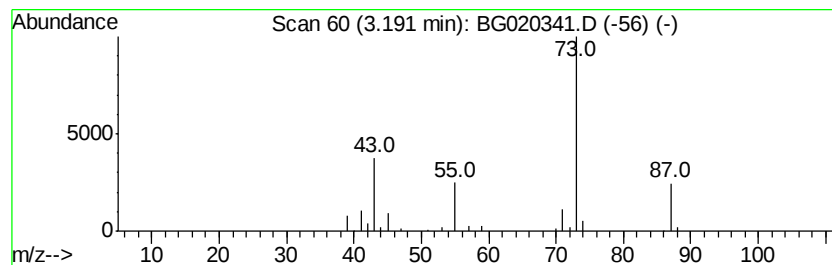
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.19	9.67 ng/ul	53312	1,4-Dichlorobenzene-d4	8.09

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



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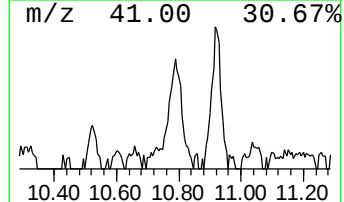
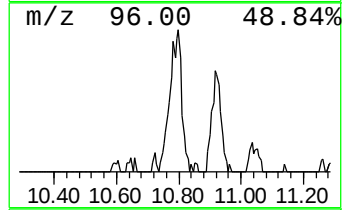
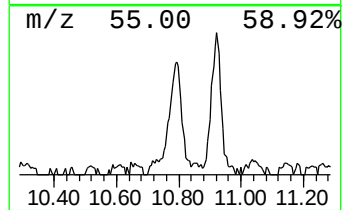
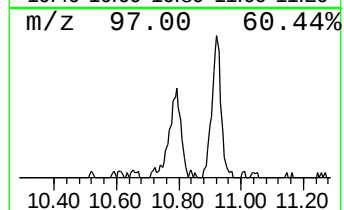
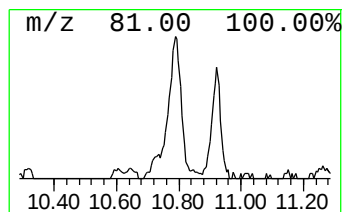
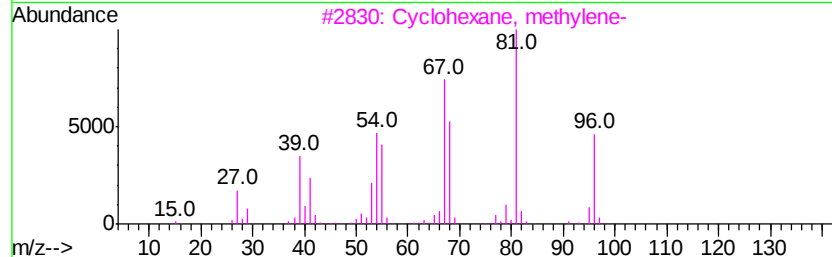
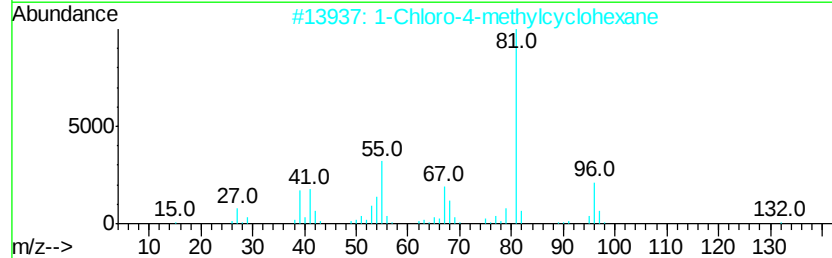
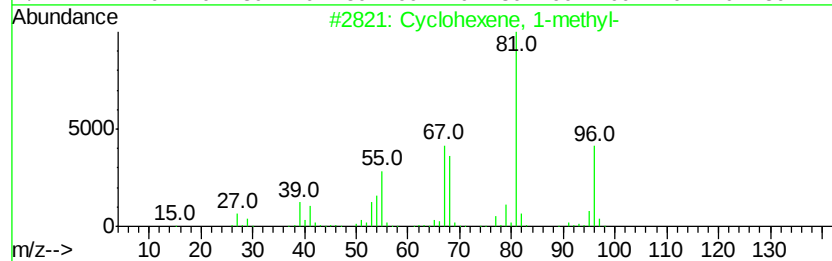
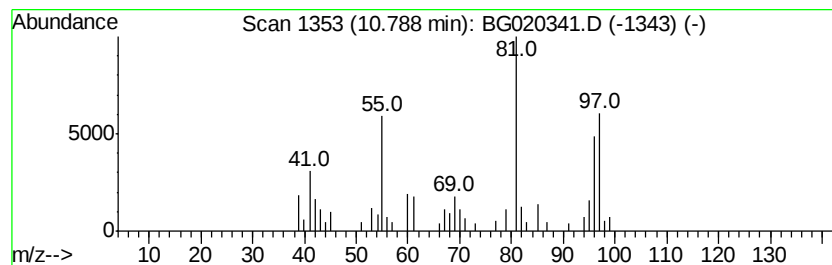
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Branched10.79 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.79	5.78 ng/ul	70946	Napthalene-d8	10.91
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexene, 1-methyl-	96 C7H12	000591-49-1 49
2		1-Chloro-4-methylcyclohexane	132 C7H13Cl	000931-68-0 47
3		Cyclohexane, methylene-	96 C7H12	001192-37-6 47
4		Cyclobutane, (1-methylethylidene)-	96 C7H12	001528-22-9 43
5		1-Methyl-2-trifluoroacetoxycyclo...	210 C9H13F3O2	1000216-63-9 40



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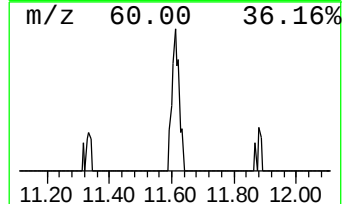
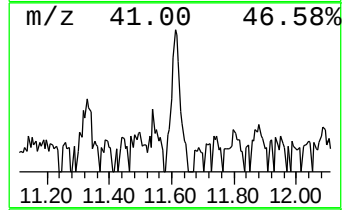
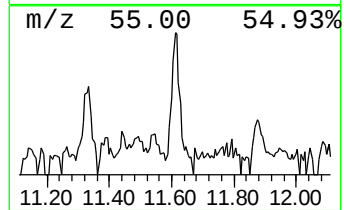
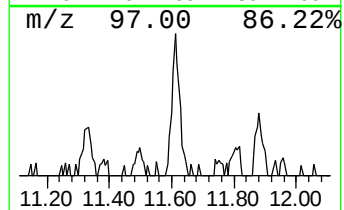
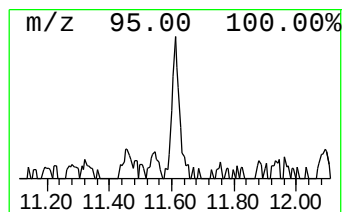
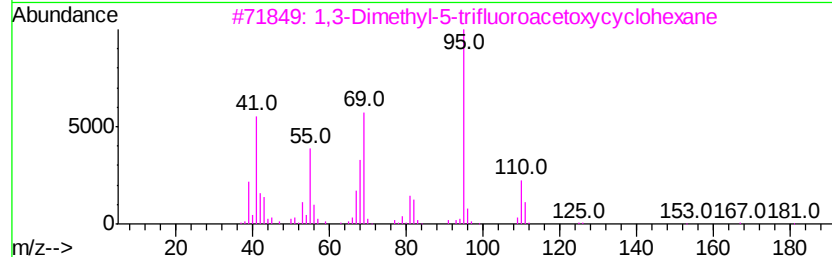
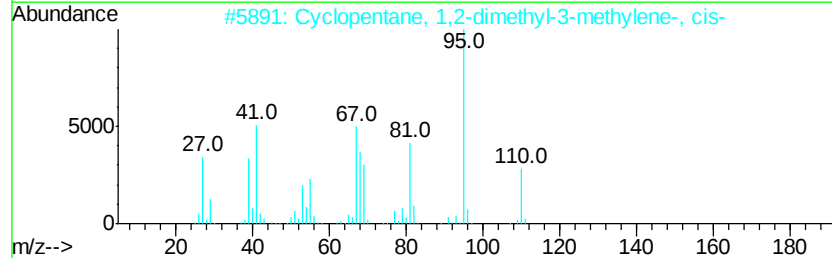
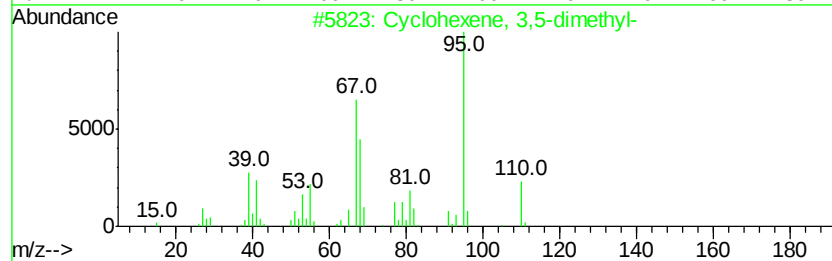
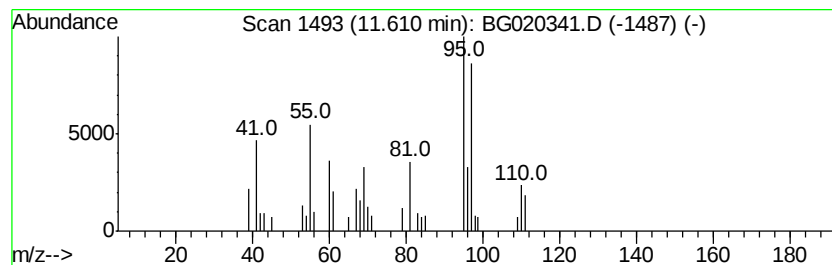
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TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Branched11.61 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.61	2.47 ng/ul	30304	Napthalene-d8	10.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexene, 3,5-dimethyl-	110	C8H14	000823-17-6	35
2		Cyclopentane, 1,2-dimethyl-3-met...	110	C8H14	091884-67-2	35
3		1,3-Dimethyl-5-trifluoroacetoxyc...	224	C10H15F3O2	1000216-63-7	32
4		5-Amino-3-methylpyrazole	97	C4H7N3	031230-17-8	25
5		5,5-Dimethyl-1,3-hexadiene	110	C8H14	001515-79-3	25



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122015\
Data File : BG020341.D
Acq On : 20 Dec 2015 14:13
Operator : UM/SJ
Sample : G4803-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C04D9

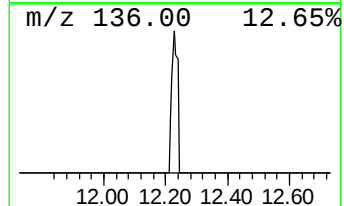
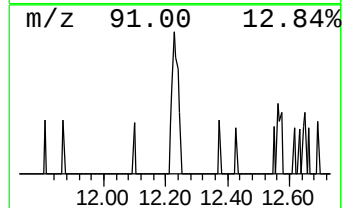
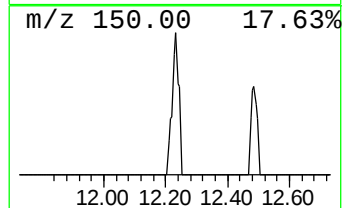
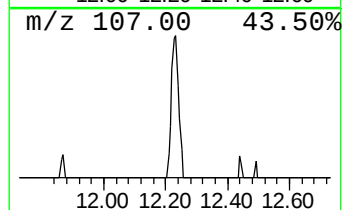
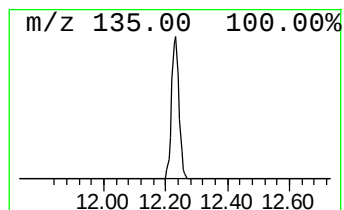
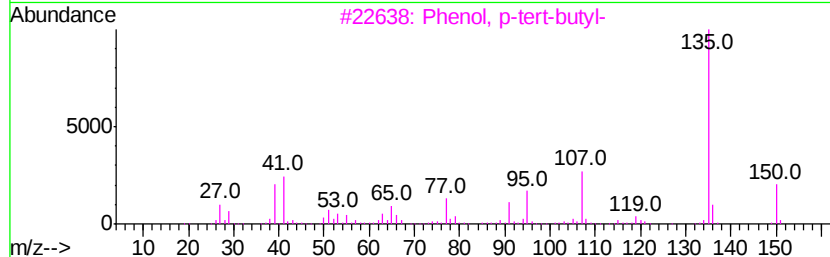
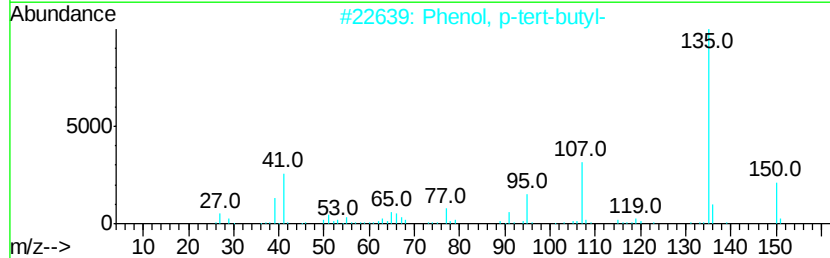
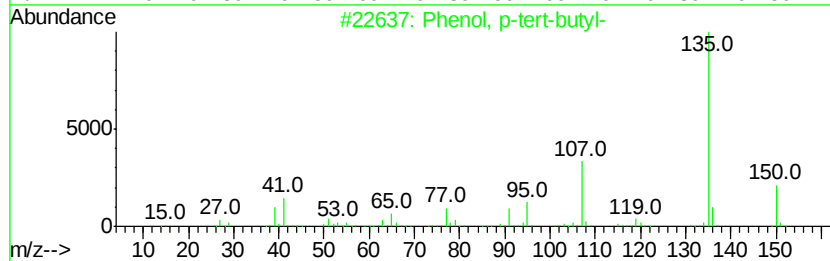
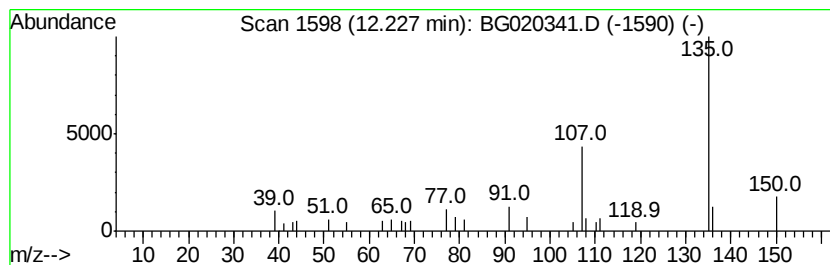
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Phenol, p-tert-butyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.23	2.02 ng/ul	24805	Naphthalene-d8	10.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	87
2		Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	86
3		Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	74
4		Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	72
5		4-Hydroxy-2-methylacetophenone	150	C9H10O2	000875-59-2	58



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122015\
Data File : BG020341.D
Acq On : 20 Dec 2015 14:13
Operator : UM/SJ
Sample : G4803-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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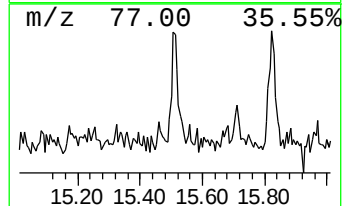
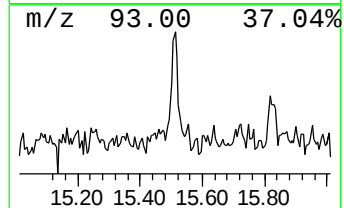
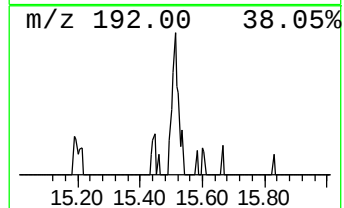
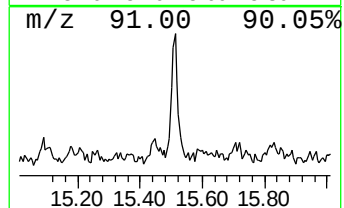
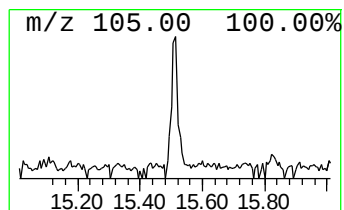
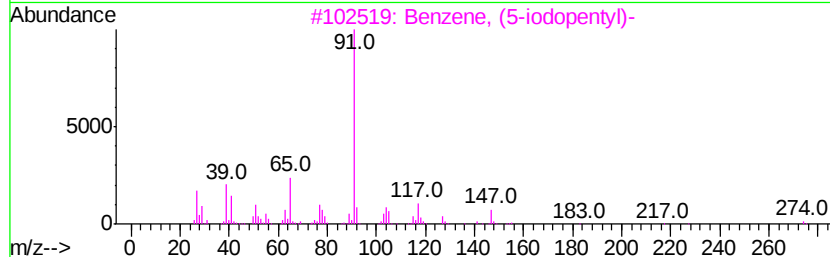
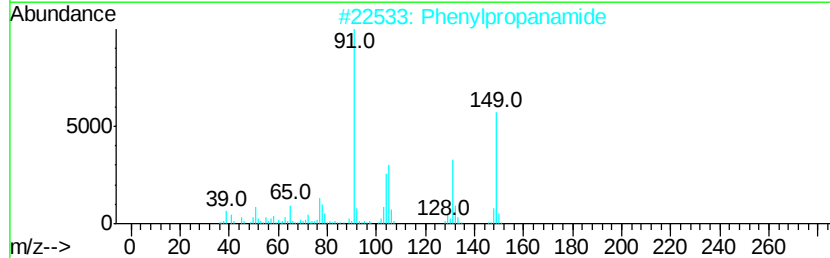
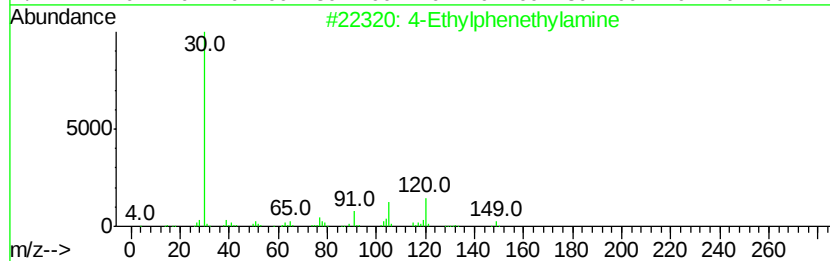
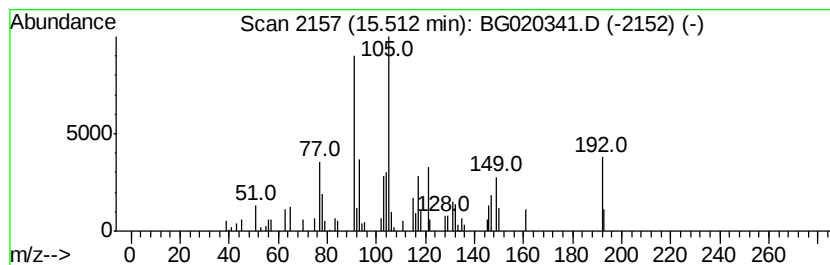
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.51	3.12 ng/ul	40276	Acenaphthene-d10	14.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Ethylphenethylamine	149	C10H15N	064353-29-3	30
2		Phenylpropanamide	149	C9H11NO	000102-93-2	14
3		Benzene, (5-iodopentyl)-	274	C11H15I	099858-37-4	10
4		Benzenemethanamine, N-butyl-N-ni...	192	C11H16N2O	040055-48-9	10
5		Benzenepropanoic acid, octyl ester	262	C17H26O2	037826-57-6	10



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122015\
Data File : BG020341.D
Acq On : 20 Dec 2015 14:13
Operator : UM/SJ
Sample : G4803-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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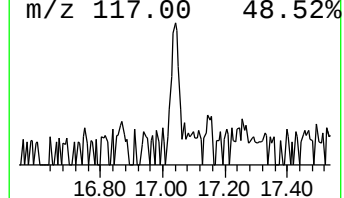
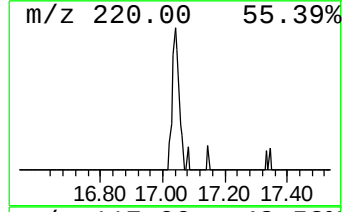
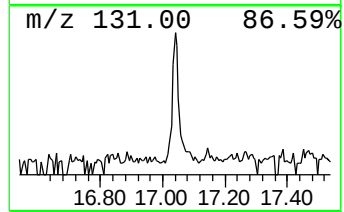
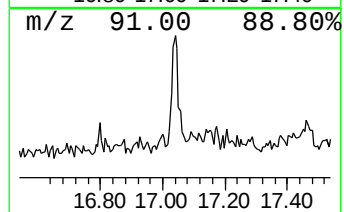
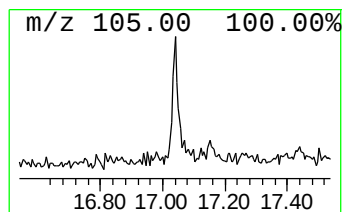
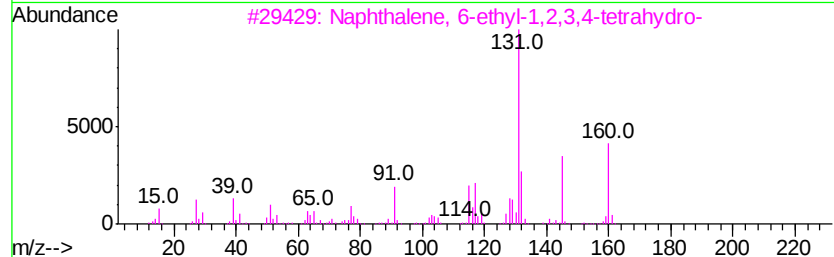
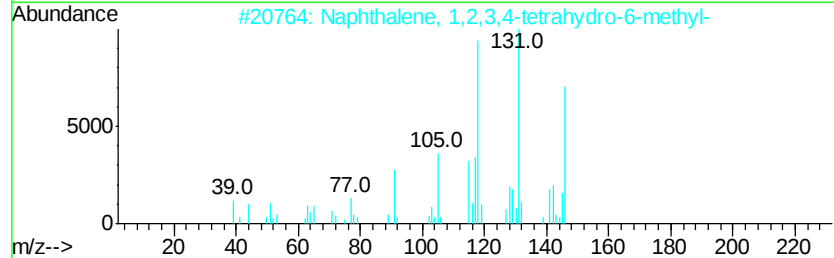
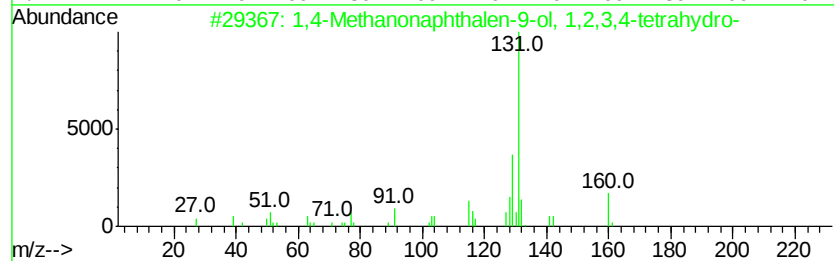
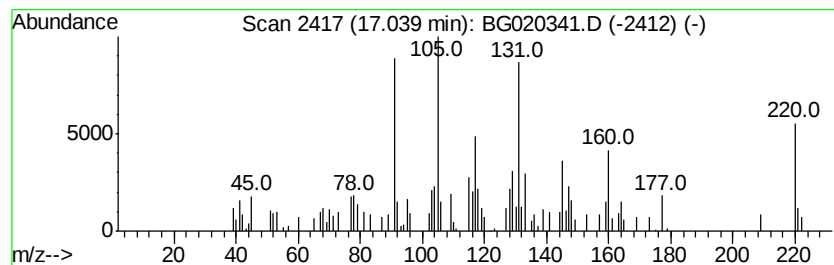
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1,4-Methanonaphthalen-9-ol,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.04	2.93 ng/ul	45205	Phenanthrene-d10	17.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,4-Methanonaphthalen-9-ol, 1,2,...	160	C11H12O	055255-94-2	52
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	38
3		Naphthalene, 6-ethyl-1,2,3,4-tet...	160	C12H16	022531-20-0	38
4		Naphthalene, 1-ethyl-1,2,3,4-tet...	160	C12H16	013556-58-6	30
5		Thiophene, tetrahydro-3-phenyl-,...	180	C10H12OS	093134-22-6	27



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122015\
Data File : BG020341.D
Acq On : 20 Dec 2015 14:13
Operator : UM/SJ
Sample : G4803-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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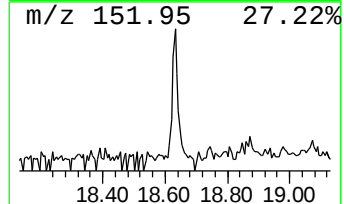
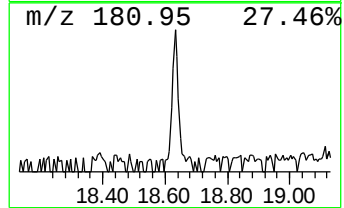
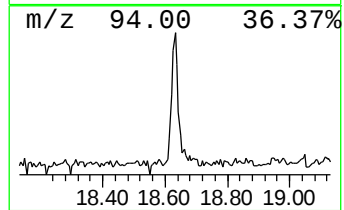
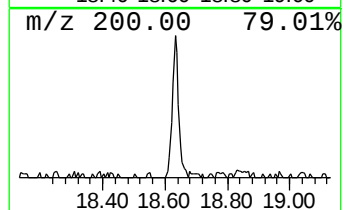
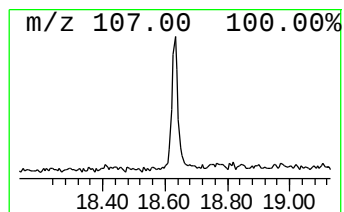
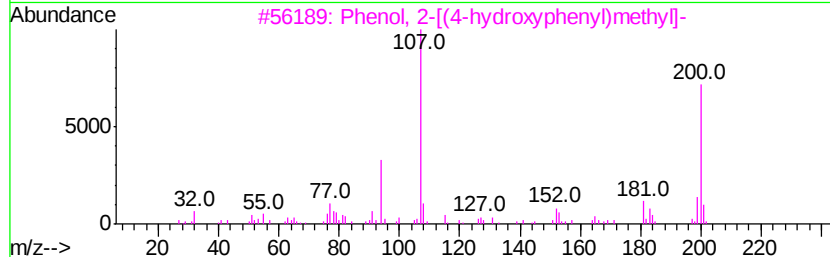
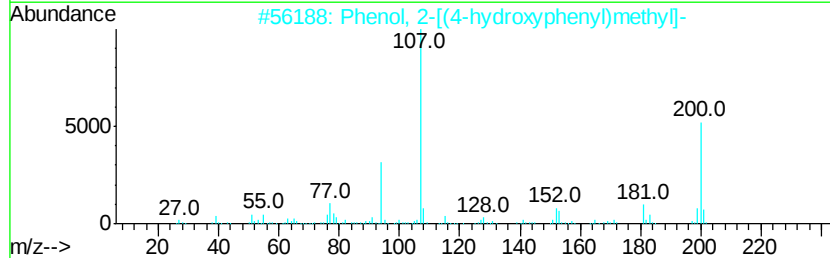
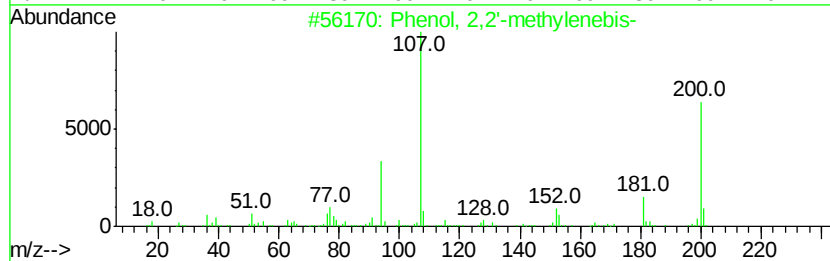
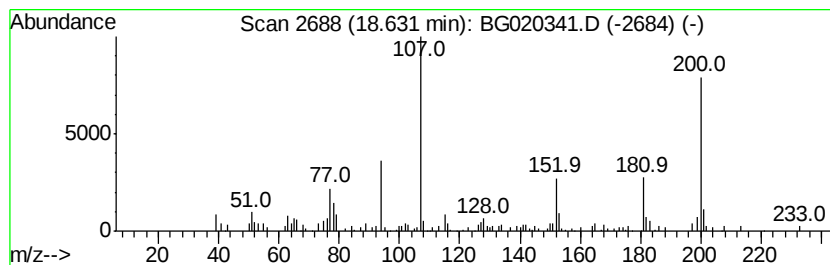
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Phenol, 2,2'-methylenebis- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.63	3.55 ng/ul	54900	Phenanthrene-d10	17.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	81
2		Phenol, 2-[(4-hydroxyphenyl)meth...	200	C13H12O2	002467-03-0	80
3		Phenol, 2-[(4-hydroxyphenyl)meth...	200	C13H12O2	002467-03-0	80
4		Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	72
5		Phenol, 4,4'-methylenebis-	200	C13H12O2	000620-92-8	59



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122015\
Data File : BG020341.D
Acq On : 20 Dec 2015 14:13
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Sample : G4803-04
Misc :
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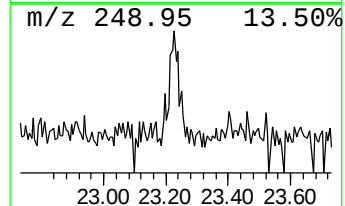
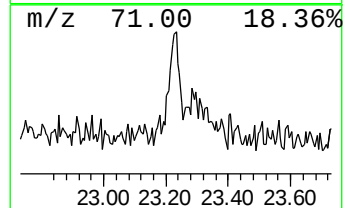
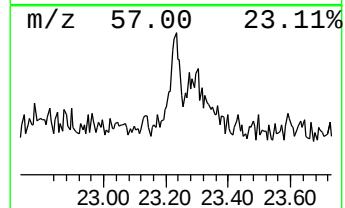
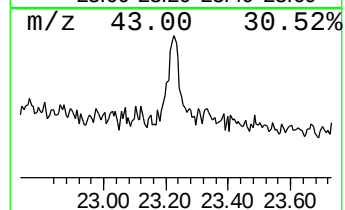
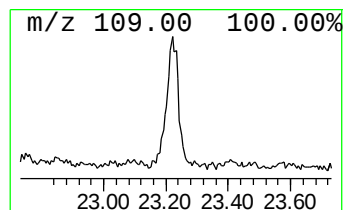
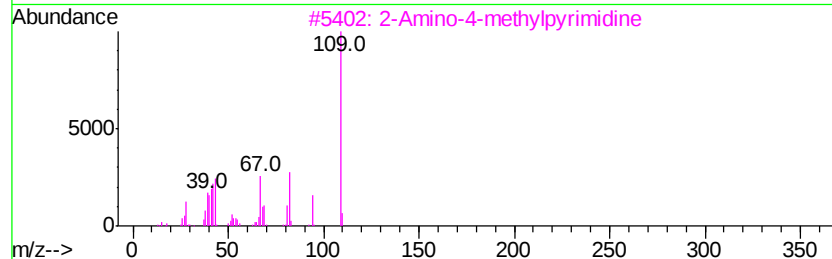
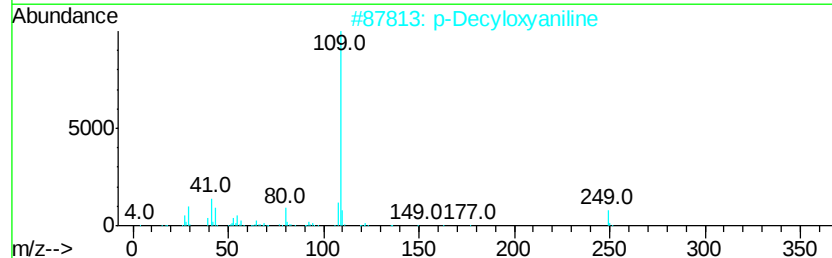
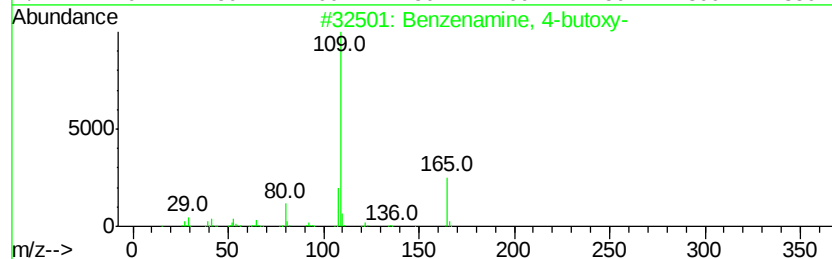
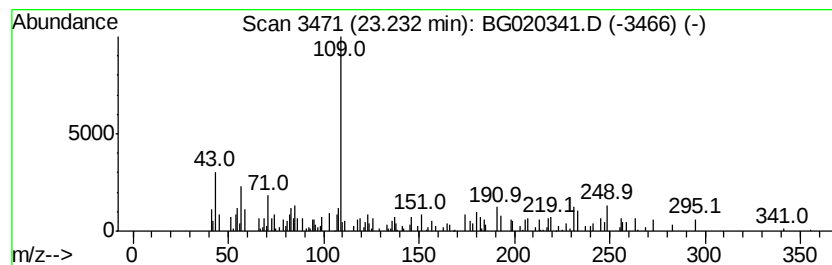
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.23	3.36 ng/ul	65736	Chrysene-d12	21.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenamine, 4-butoxy-	165	C10H15NO	004344-55-2	43
2		p-Decyloxvaniline	249	C16H27NO	039905-47-0	40
3		2-Amino-4-methylpyrimidine	109	C5H7N3	000108-52-1	38
4		Bicyclo[2.2.1]heptan-2-one, 5-br...	230	C10H15BrO	010293-00-2	38
5		1,2-Diazaspiro[4.4]nonen-3-carbo...	252	C14H24N2O2	1000164-25-9	38



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG122015\
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ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.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...	3.19	9.7	ng/ul	53312	1	8.09	110221	20.0
(DEL) Alkane: Bra...	10.79	5.8	ng/ul	70946	2	10.91	245329	20.0
(DEL) Alkane: Bra...	11.61	2.5	ng/ul	30304	2	10.91	245329	20.0
Phenol, p-tert-bu...	12.23	2.0	ng/ul	24805	2	10.91	245329	20.0
unknown-01	15.51	3.1	ng/ul	40276	3	14.72	258457	20.0
1,4-Methanonaphth...	17.04	2.9	ng/ul	45205	4	17.47	309053	20.0
Phenol, 2,2'-meth...	18.63	3.5	ng/ul	54900	4	17.47	309053	20.0
unknown-02	23.23	3.4	ng/ul	65736	5	21.74	391241	20.0