

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG070715\
 Data File : BG017792.D
 Acq On : 7 Jul 2015 12:19
 Operator : TP/UM
 Sample : G2860-05
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 G93J0

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-BG070615.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	2.813	17	21	25	rBV	92030	140575	5.73%	0.532%
2	2.884	29	33	46	rVB	1931390	2452929	100.00%	9.287%
3	3.648	159	163	168	rVV4	22367	34641	1.41%	0.131%
4	4.746	346	350	360	rVB3	21186	41569	1.69%	0.157%
5	6.779	687	696	709	rVV	148403	280743	11.45%	1.063%
6	6.950	718	725	731	rBV	101121	160124	6.53%	0.606%
7	7.138	751	757	766	rVB2	151830	244686	9.98%	0.926%
8	7.243	770	775	785	rVB2	4558	13290	0.54%	0.050%
9	7.602	828	836	844	rBV	237154	392304	15.99%	1.485%
10	8.307	947	956	962	rVV	171999	292336	11.92%	1.107%
11	8.424	970	976	984	rVV	109298	191678	7.81%	0.726%
12	8.548	994	997	1002	rVV5	7590	12467	0.51%	0.047%
13	8.753	1026	1032	1041	rVB	136167	219760	8.96%	0.832%
14	9.470	1148	1154	1161	rVV2	111157	194505	7.93%	0.736%
15	9.641	1178	1183	1187	rBV5	7117	14602	0.60%	0.055%
16	9.829	1211	1215	1221	rVV4	11321	17826	0.73%	0.067%
17	10.005	1238	1245	1254	rVB	185654	307530	12.54%	1.164%
18	10.387	1301	1310	1316	rBV	377090	653314	26.63%	2.473%
19	10.434	1316	1318	1324	rVV	17554	21785	0.89%	0.082%
20	10.528	1327	1334	1347	rVB	100474	190700	7.77%	0.722%
21	10.927	1396	1402	1406	rVV6	10908	21378	0.87%	0.081%
22	11.174	1439	1444	1448	rVB	10118	14527	0.59%	0.055%
23	12.055	1588	1594	1599	rVB6	11496	18911	0.77%	0.072%
24	12.279	1628	1632	1637	rBV4	10451	15128	0.62%	0.057%
25	12.860	1728	1731	1735	rVB3	11043	17941	0.73%	0.068%
26	13.571	1847	1852	1858	rVB4	14672	26122	1.06%	0.099%
27	13.671	1864	1869	1874	rVV	280364	394384	16.08%	1.493%
28	13.718	1874	1877	1882	rVB	96008	123909	5.05%	0.469%
29	13.812	1888	1893	1896	rBV5	10497	16863	0.69%	0.064%
30	13.947	1907	1916	1929	rBV	366720	592010	24.13%	2.241%
31	14.176	1950	1955	1959	rVB2	17786	23813	0.97%	0.090%
32	14.259	1959	1969	1976	rBV2	612675	942891	38.44%	3.570%
33	14.323	1976	1980	1988	rVB7	12652	23854	0.97%	0.090%
34	14.459	1997	2003	2016	rBV	109232	193207	7.88%	0.731%

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35	14.623	2025	2031	2033	rBV7	10836	17631	0.72%	0.067%
36	14.658	2033	2037	2045	rVB7	15988	38274	1.56%	0.145%
37	14.793	2057	2060	2065	rBV6	9014	14440	0.59%	0.055%
38	14.882	2071	2075	2080	rVV7	10260	18087	0.74%	0.068%
39	15.046	2100	2103	2107	rVV5	9369	15095	0.62%	0.057%
40	15.134	2113	2118	2122	rVV3	29572	38477	1.57%	0.146%
41	15.252	2132	2138	2144	rVV	457575	647667	26.40%	2.452%
42	15.310	2144	2148	2154	rVV4	28399	45675	1.86%	0.173%
43	15.375	2154	2159	2166	rVB2	59955	84911	3.46%	0.321%
44	15.534	2182	2186	2191	rVB6	22326	38404	1.57%	0.145%
45	15.757	2220	2224	2226	rBV5	13627	20279	0.83%	0.077%
46	15.998	2261	2265	2267	rVV3	43035	65229	2.66%	0.247%
47	16.021	2267	2269	2274	rVB	39085	44392	1.81%	0.168%
48	16.521	2350	2354	2358	rBV	67507	85925	3.50%	0.325%
49	16.662	2374	2378	2382	rVB2	27640	37854	1.54%	0.143%
50	16.738	2387	2391	2394	rBV5	12892	18580	0.76%	0.070%
51	16.791	2396	2400	2403	rVV4	32937	46524	1.90%	0.176%
52	16.832	2403	2407	2412	rVB3	40802	71579	2.92%	0.271%
53	17.008	2431	2437	2441	rBV2	730072	1051971	42.89%	3.983%
54	17.050	2441	2444	2449	rVB	407413	533761	21.76%	2.021%
55	17.108	2449	2454	2458	rBV	465640	646001	26.34%	2.446%
56	17.384	2498	2501	2503	rBV4	13864	14098	0.57%	0.053%
57	17.420	2503	2507	2510	rVV2	23327	29722	1.21%	0.113%
58	17.531	2522	2526	2531	rVB6	41609	61663	2.51%	0.233%
59	17.590	2533	2536	2540	rVV3	23276	30120	1.23%	0.114%
60	17.790	2567	2570	2571	rBV2	11806	14463	0.59%	0.055%
61	17.890	2583	2587	2590	rBV	68449	96289	3.93%	0.365%
62	17.937	2590	2595	2600	rVB	103796	163807	6.68%	0.620%
63	18.013	2605	2608	2613	rVB	30138	35792	1.46%	0.136%
64	18.078	2614	2619	2623	rVV3	106395	186491	7.60%	0.706%
65	18.119	2624	2626	2630	rVB	61442	65554	2.67%	0.248%
66	18.184	2635	2637	2641	rBV5	10170	17111	0.70%	0.065%
67	18.225	2641	2644	2651	rVV9	17904	31917	1.30%	0.121%
68	18.395	2669	2673	2676	rBV	59321	74523	3.04%	0.282%
69	18.430	2676	2679	2685	rVB6	51217	82742	3.37%	0.313%
70	18.495	2686	2690	2696	rVV	39658	58516	2.39%	0.222%
71	18.724	2726	2729	2734	rVB3	33496	47537	1.94%	0.180%

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72	18.812	2740	2744	2749	rBV	70573	121078	4.94%	0.458%
73	18.865	2750	2753	2758	rVV6	34731	58248	2.37%	0.221%
74	18.953	2764	2768	2775	rBV3	64856	121181	4.94%	0.459%
75	19.077	2784	2789	2795	rVB	762340	1027958	41.91%	3.892%
76	19.412	2841	2846	2849	rBV2	599853	926972	37.79%	3.509%
77	19.441	2849	2851	2861	rVB	872229	1061927	43.29%	4.020%
78	19.623	2879	2882	2886	rBV2	38062	59212	2.41%	0.224%
79	19.923	2927	2933	2934	rBV6	55301	107839	4.40%	0.408%
80	20.040	2951	2953	2958	rVB3	62817	67798	2.76%	0.257%
81	20.099	2959	2963	2967	rVV2	128295	177089	7.22%	0.670%
82	20.240	2983	2987	2991	rBV2	120253	153868	6.27%	0.583%
83	20.287	2992	2995	2999	rVV	105803	139697	5.70%	0.529%
84	20.322	2999	3001	3009	rVB6	62250	91719	3.74%	0.347%
85	20.686	3060	3063	3069	rVB	132850	197278	8.04%	0.747%
86	20.781	3076	3079	3084	rVB	89012	114716	4.68%	0.434%
87	20.833	3085	3088	3091	rBV2	87936	128810	5.25%	0.488%
88	20.945	3102	3107	3111	rBV3	139776	251475	10.25%	0.952%
89	21.204	3145	3151	3156	rBV2	1089989	2135764	87.07%	8.086%
90	21.245	3156	3158	3165	rVB	521739	641135	26.14%	2.427%
91	21.321	3168	3171	3174	rBV	117649	145117	5.92%	0.549%
92	21.697	3233	3235	3238	rBV4	66111	90384	3.68%	0.342%
93	21.756	3239	3245	3249	rVB4	119972	263174	10.73%	0.996%
94	22.267	3329	3332	3340	rVB6	170574	304830	12.43%	1.154%
95	22.766	3412	3417	3422	rBV	468054	1062784	43.33%	4.024%
96	23.242	3493	3498	3502	rBV	351271	630878	25.72%	2.388%
97	23.295	3502	3507	3510	rVV2	335393	633792	25.84%	2.400%
98	23.336	3510	3514	3523	rVB	560880	1047710	42.71%	3.967%
99	23.436	3525	3531	3536	rBV	556686	1045838	42.64%	3.960%
100	25.692	3908	3915	3925	rVB3	247182	711590	29.01%	2.694%

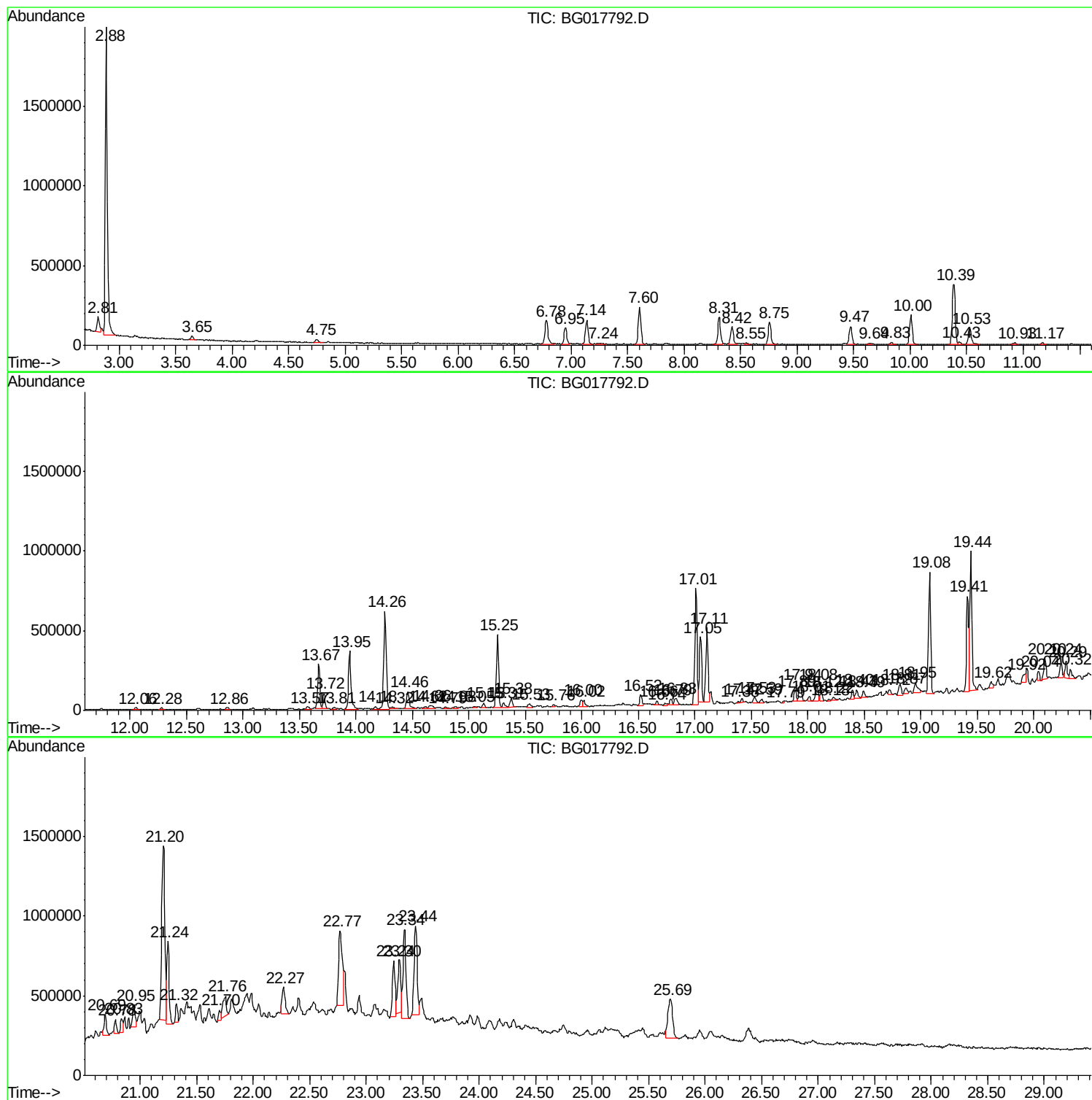
Sum of corrected areas: 26413264

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG070615.M
Quant Title : SVOA CALIBRATION

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



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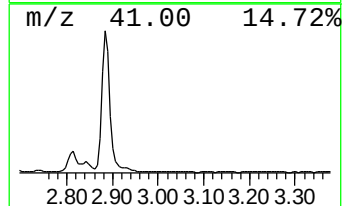
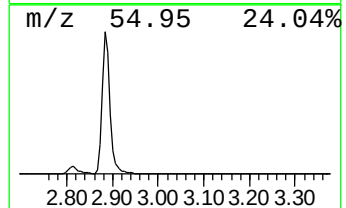
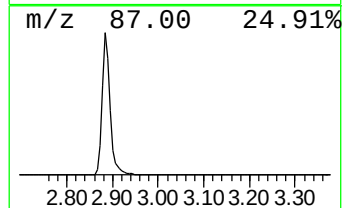
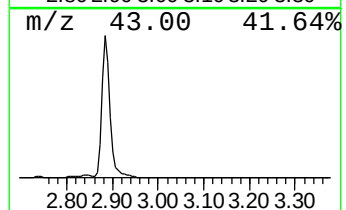
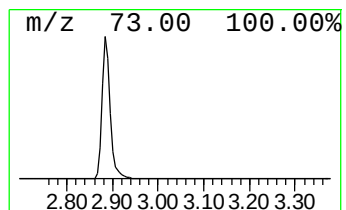
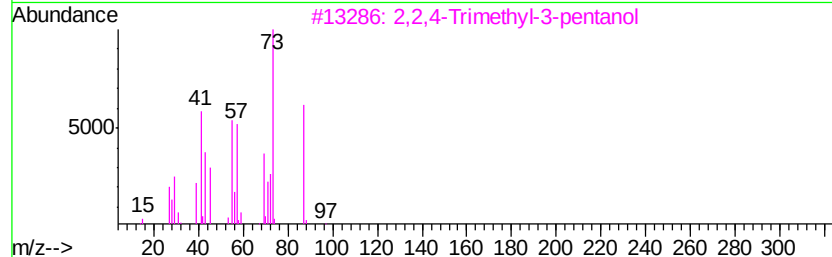
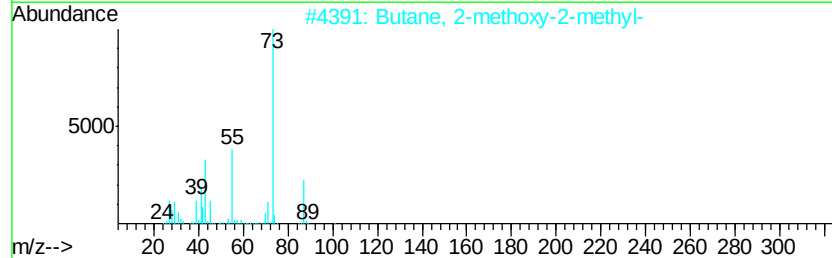
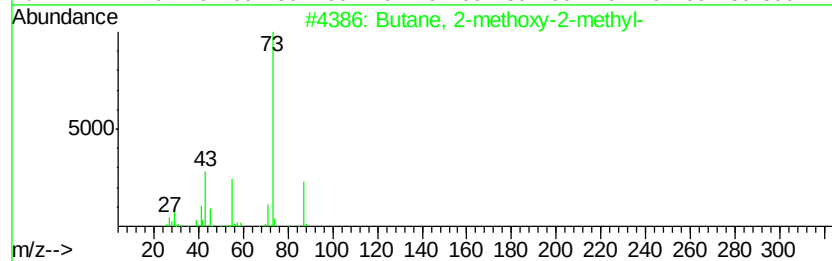
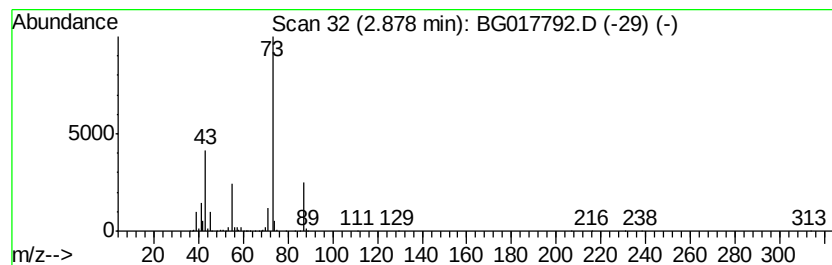
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG070615.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.88	125.05 ng/ul	2452930	1,4-Dichlorobenzene-d4	7.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	50
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	35
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17



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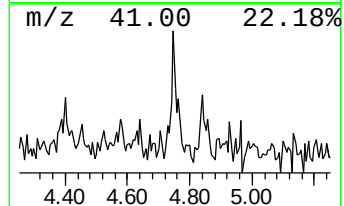
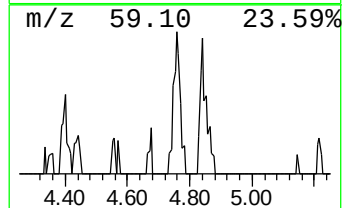
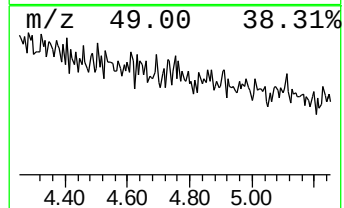
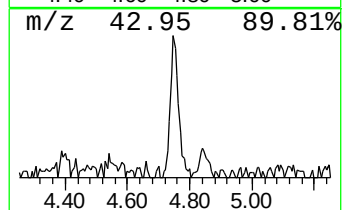
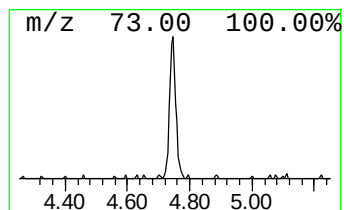
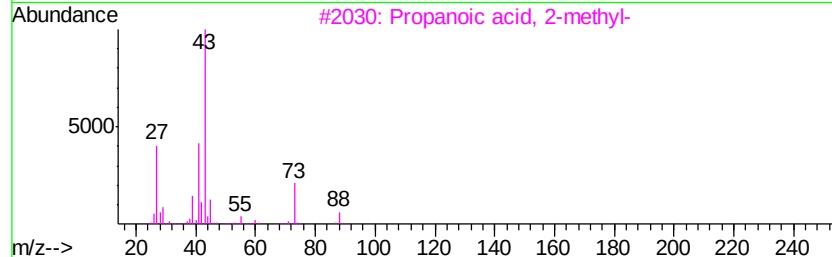
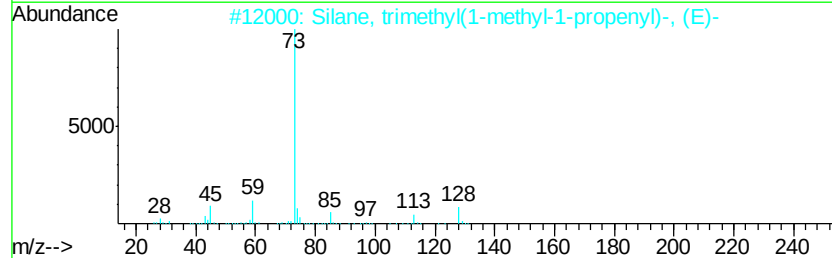
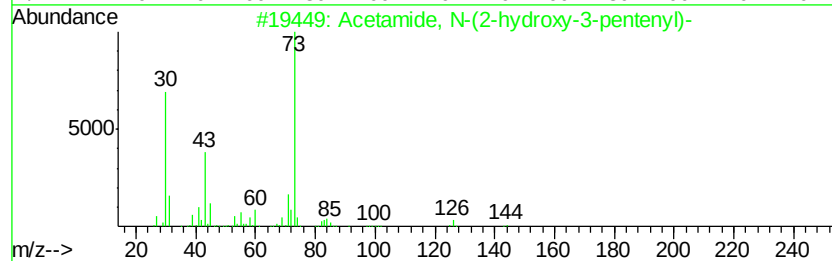
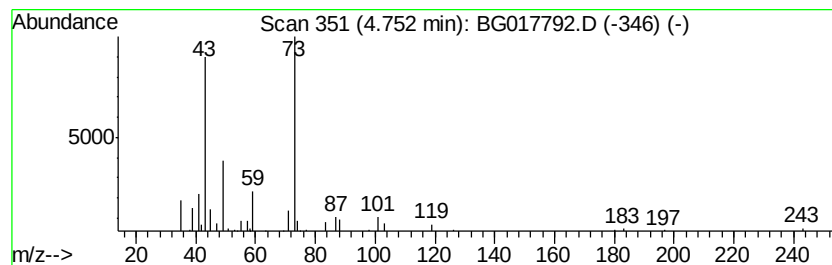
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.75	2.12 ng/ul	41569	1,4-Dichlorobenzene-d4	7.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Acetamide, N-(2-hydroxy-3-penten...	143	C7H13NO2	093393-95-4	38
2		Silane, trimethyl(1-methyl-1-pro...	128	C7H16Si	010111-13-4	25
3		Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	9
4		1,4-Butanediamine	88	C4H12N2	000110-60-1	9
5		Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	9



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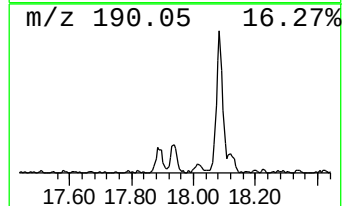
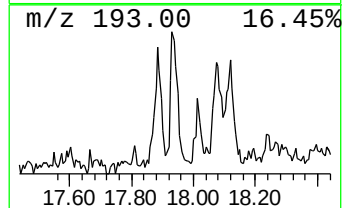
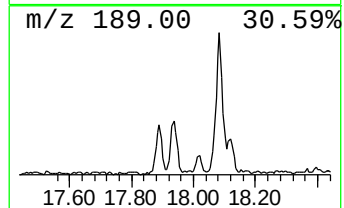
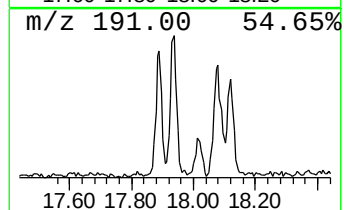
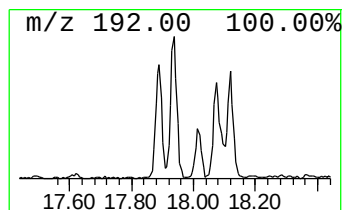
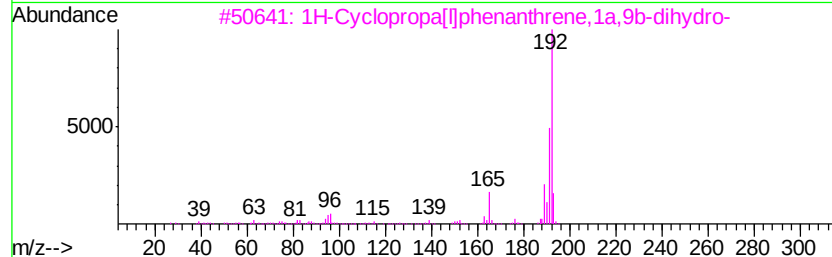
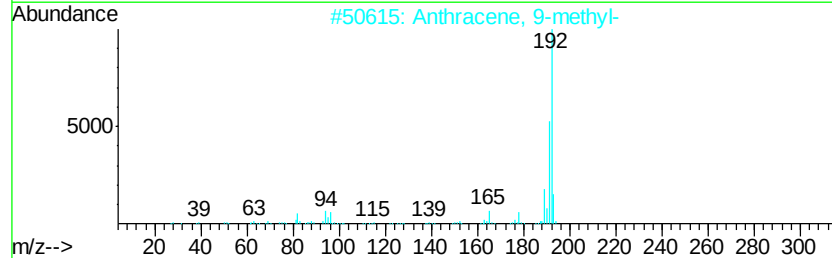
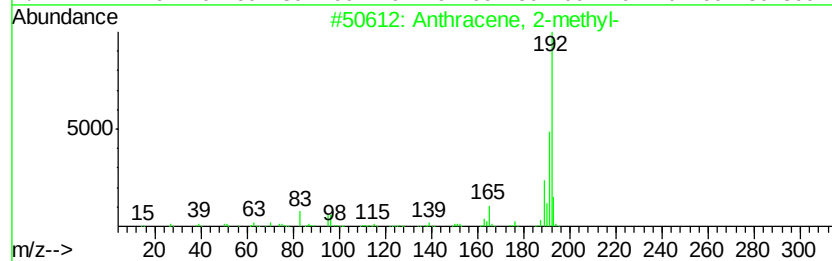
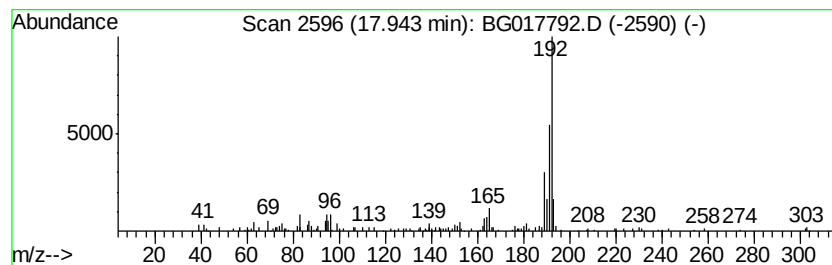
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Peak Number 4 Anthracene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.94	3.11 ng/ul	163807	Phenanthrene-d10	17.01	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
2	Anthracene, 9-methyl-	192	C15H12	000779-02-2	95
3	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	94
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG070715\
Data File : BG017792.D
Acq On : 7 Jul 2015 12:19
Operator : TP/UM
Sample : G2860-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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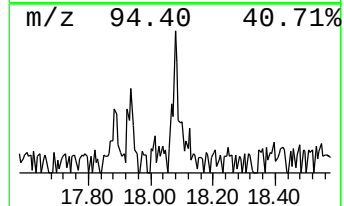
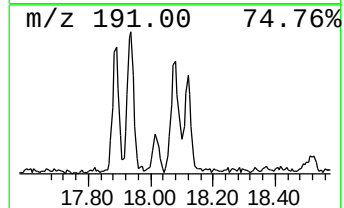
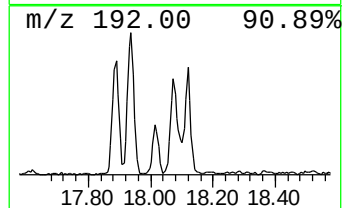
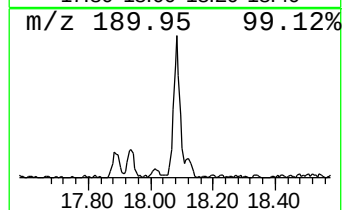
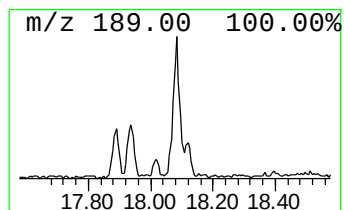
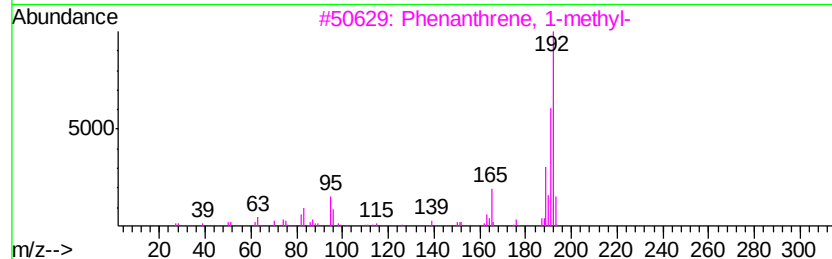
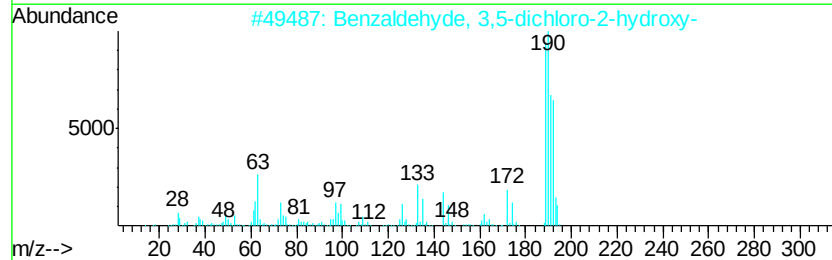
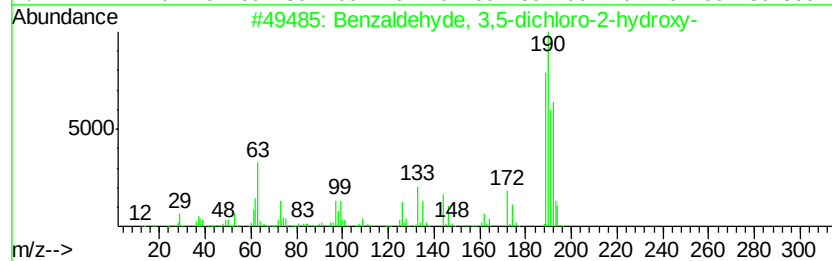
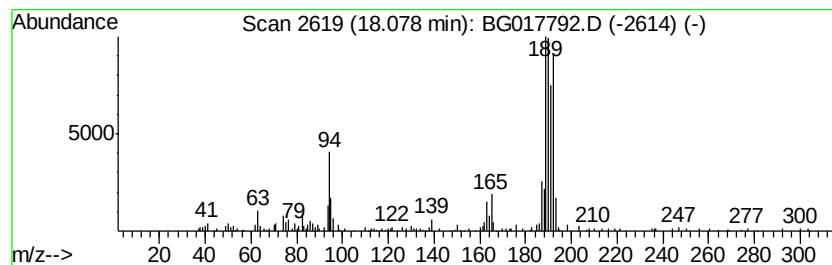
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzaldehyde, 3,5-dichloro-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.08	3.55 ng/ul	186491	Phenanthrene-d10	17.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	64
2		Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	64
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	46
4		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	41
5		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	41



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG070715\
Data File : BG017792.D
Acq On : 7 Jul 2015 12:19
Operator : TP/UM
Sample : G2860-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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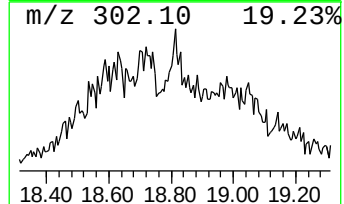
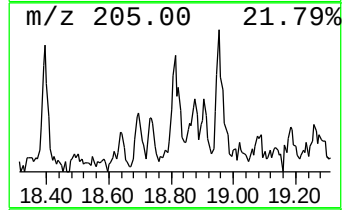
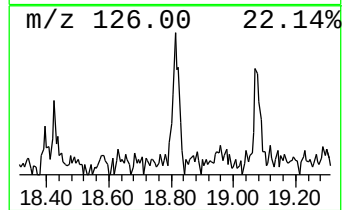
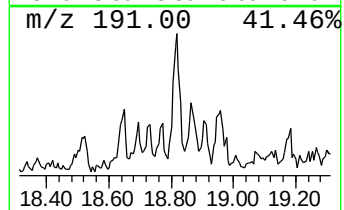
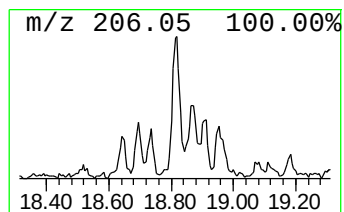
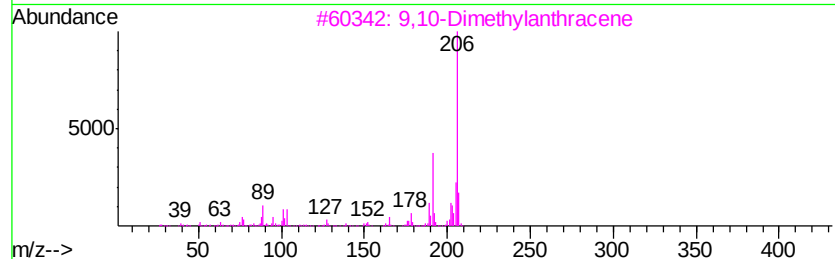
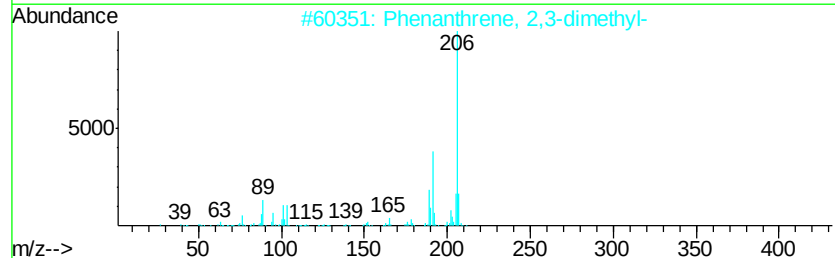
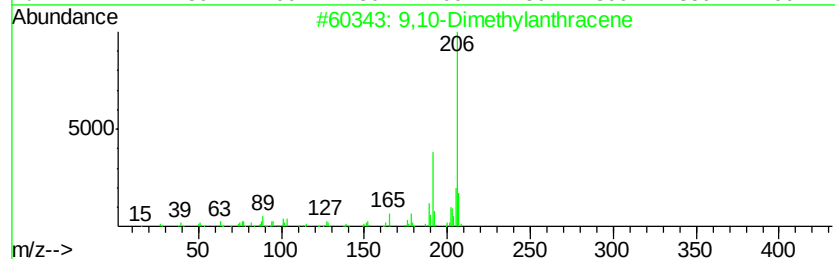
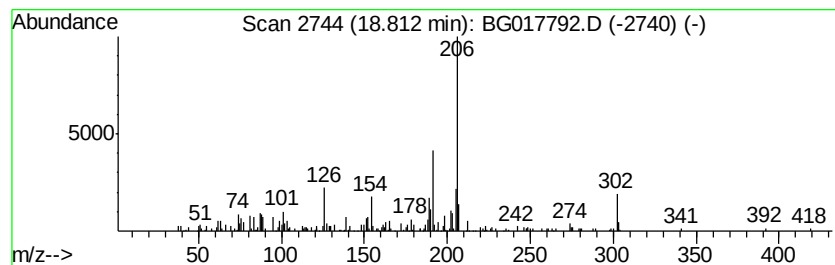
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 9,10-Dimethylantracene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.81	2.30 ng/ul	121078	Phenanthrene-d10	17.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Dimethylantracene	206	C16H14	000781-43-1	70		
2	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	64		
3	9,10-Dimethylantracene	206	C16H14	000781-43-1	62		
4	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	62		
5	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	62		



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG070715\
Data File : BG017792.D
Acq On : 7 Jul 2015 12:19
Operator : TP/UM
Sample : G2860-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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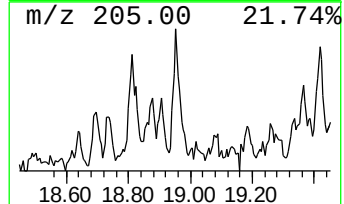
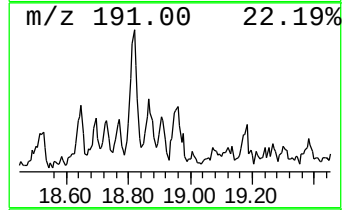
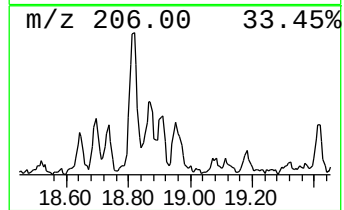
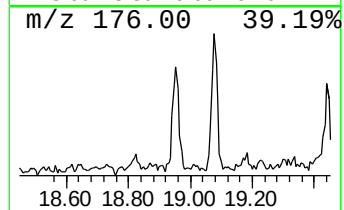
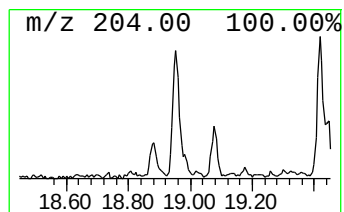
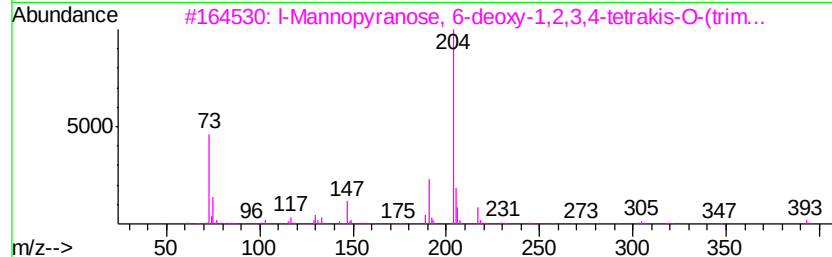
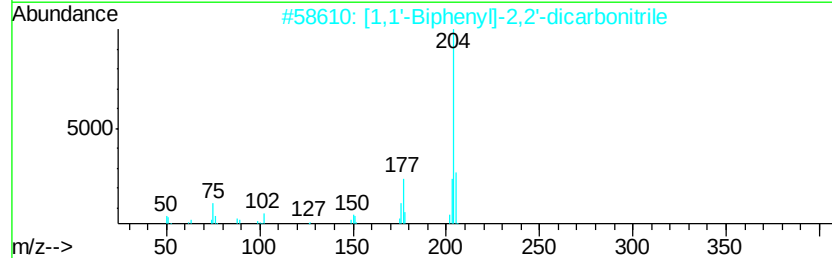
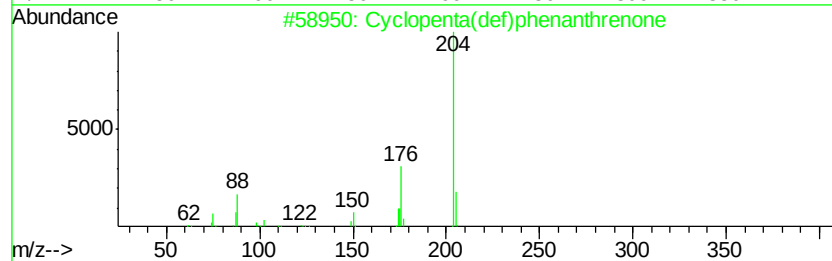
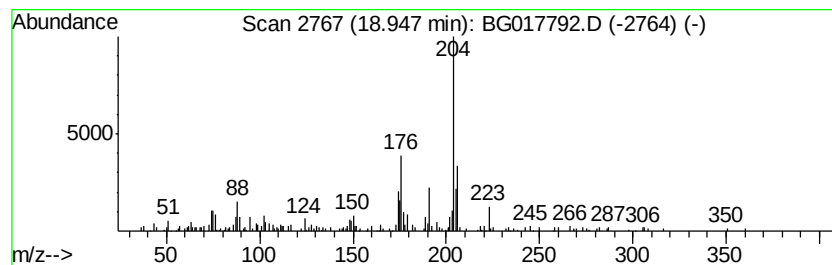
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Cyclopenta(def)phenanthrenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.95	2.30 ng/ul	121181	Phenanthrene-d10	17.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	58
2		[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	46
3		l-Mannopyranose, 6-deoxy-1,2,3,4...	452	C18H44O5Si4	108392-01-4	43
4		[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	43
5		1-Ethyl-1,2,3,4-tetrahydro-6,7-m...	219	C12H13N03	132767-88-5	38



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG070715\
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Acq On : 7 Jul 2015 12:19
Operator : TP/UM
Sample : G2860-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

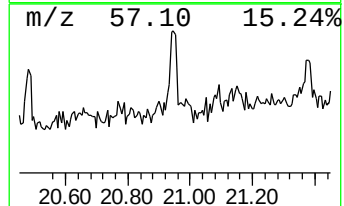
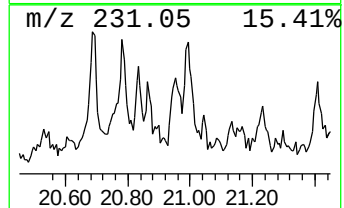
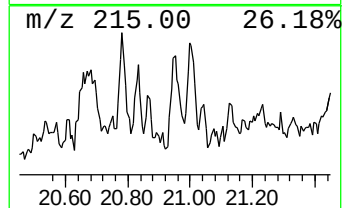
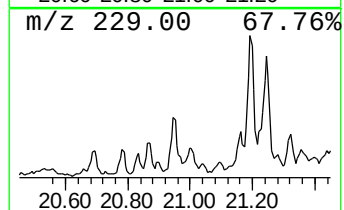
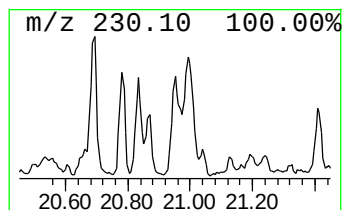
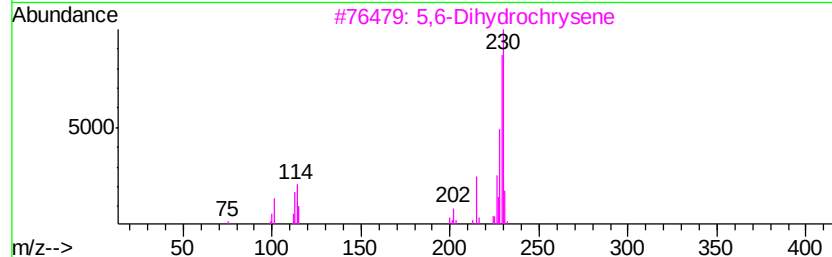
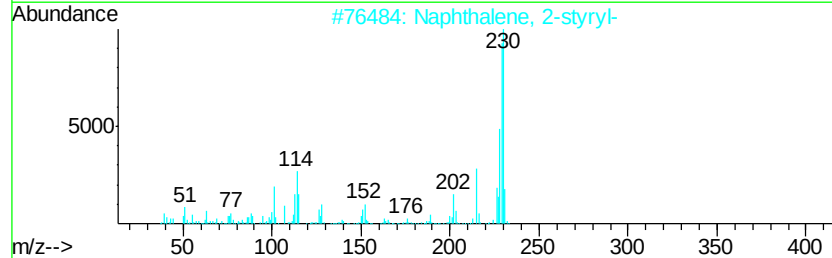
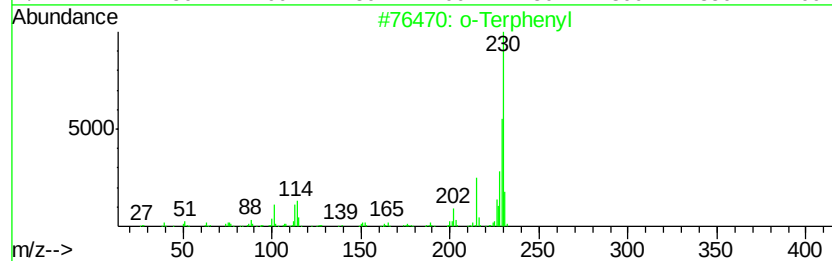
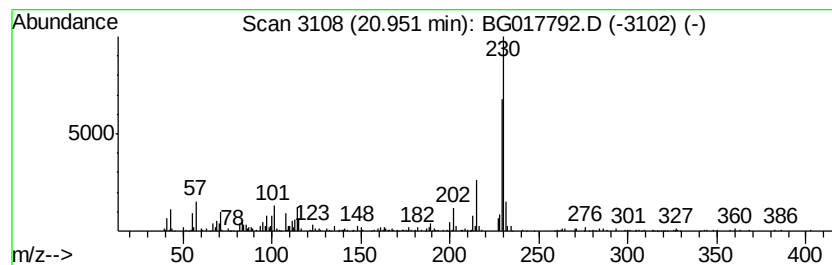
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 o-Terphenyl Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.95	2.35 ng/ul	251475	Chrysene-d12	21.21
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		o-Terphenyl	230 C18H14	000084-15-1 95
2		Naphthalene, 2-styryl-	230 C18H14	002039-70-5 93
3		5,6-Dihydrochrysene	230 C18H14	002091-92-1 90
4		6,6-Diphenylfulvene	230 C18H14	002175-90-8 81
5		Naphthacene, 5,12-dihydro-	230 C18H14	000959-02-4 80



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG070715\
Data File : BG017792.D
Acq On : 7 Jul 2015 12:19
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Sample : G2860-05
Misc :
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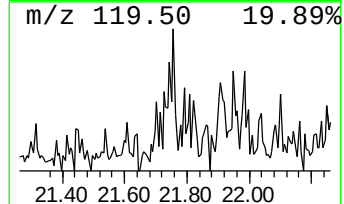
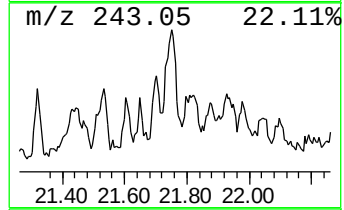
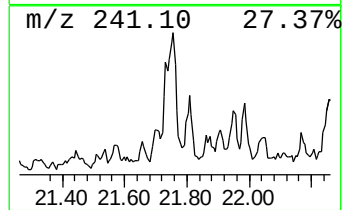
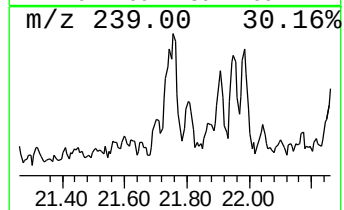
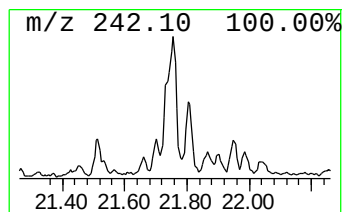
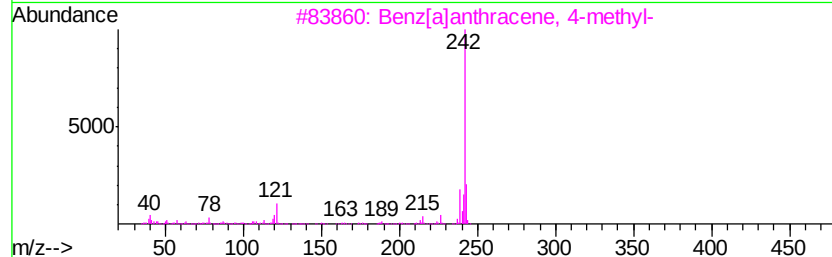
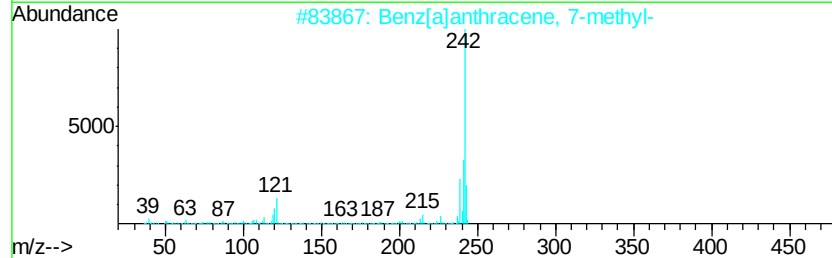
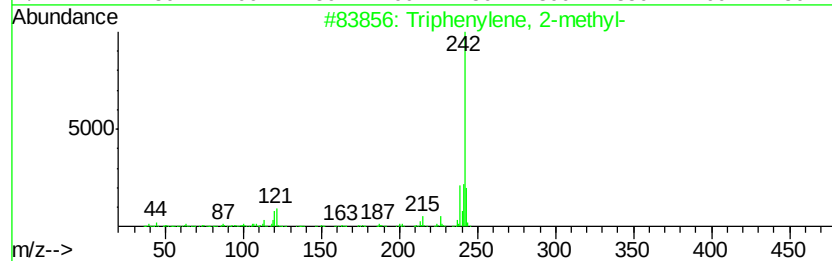
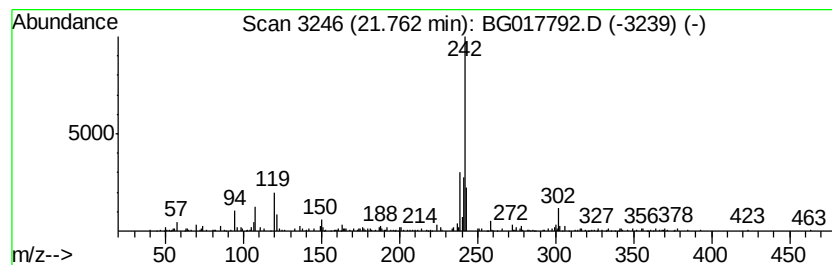
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Triphenylene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.76	2.46 ng/ul	263174	Chrysene-d12	21.21
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Triphenylene, 2-methyl-	242 C19H14	001705-84-6 86
2		Benz[a]anthracene, 7-methyl-	242 C19H14	002541-69-7 83
3		Benz[a]anthracene, 4-methyl-	242 C19H14	000316-49-4 64
4		Benz[a]anthracene, 5-methyl-	242 C19H14	002319-96-2 64
5		Benz[a]anthracene, 11-methyl-	242 C19H14	006111-78-0 64



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG070715\
Data File : BG017792.D
Acq On : 7 Jul 2015 12:19
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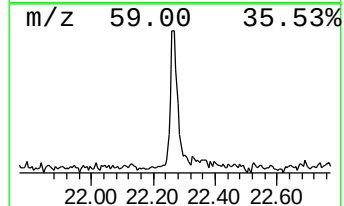
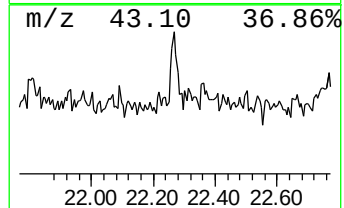
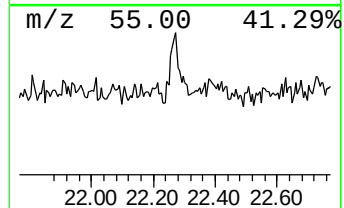
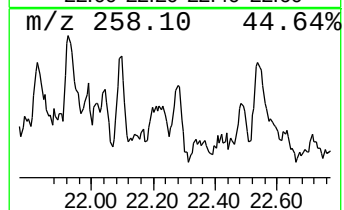
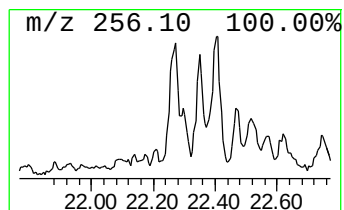
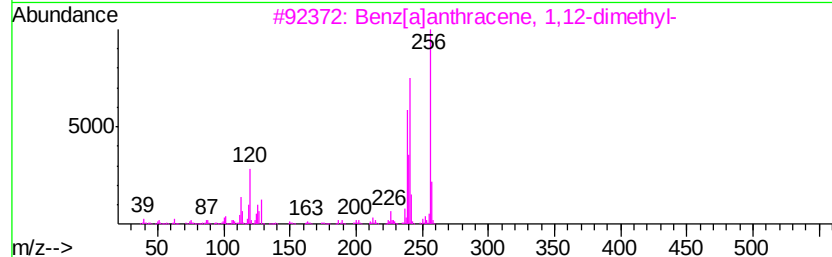
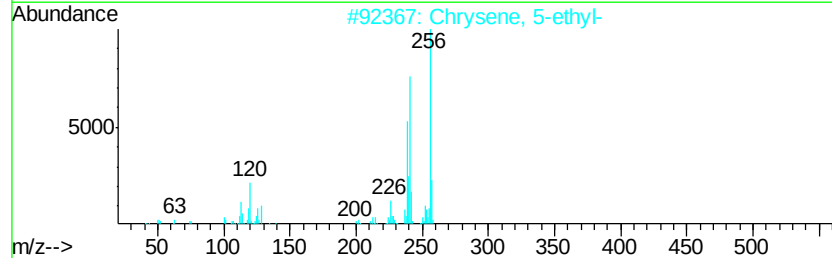
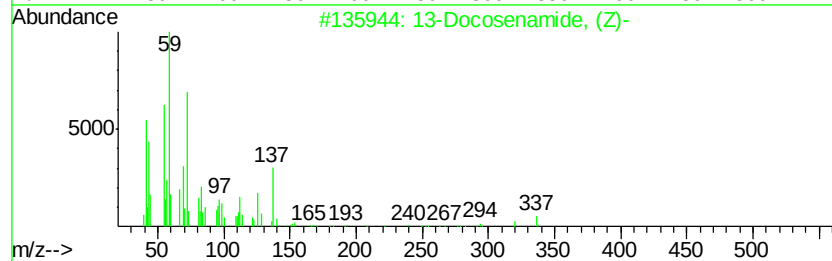
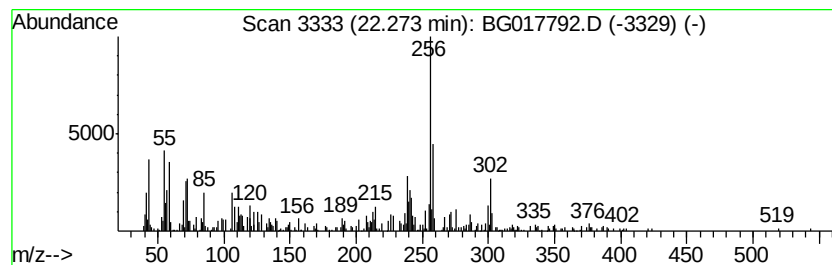
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 13-Docosenamide, (Z)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.27	2.85 ng/ul	304830	Chrysene-d12	21.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	53
2		Chrysene, 5-ethyl-	256	C20H16	054986-62-8	50
3		Benz[a]anthracene, 1,12-dimethyl-	256	C20H16	000313-74-6	41
4		Benz[a]anthracene, 1,12-dimethyl-	256	C20H16	000313-74-6	38
5		Benz[a]anthracene, 7,12-dimethyl-	256	C20H16	000057-97-6	38



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG070715\
Data File : BG017792.D
Acq On : 7 Jul 2015 12:19
Operator : TP/UM
Sample : G2860-05
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG070615.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
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unknown-01	4.75	2.1	ng/ul	41569	1	7.60	392304	20.0
Anthracene, 2-met...	17.94	3.1	ng/ul	163807	4	17.01	1051970	20.0
Benzaldehyde, 3,5...	18.08	3.5	ng/ul	186491	4	17.01	1051970	20.0
9,10-Dimethylanthe...	18.81	2.3	ng/ul	121078	4	17.01	1051970	20.0
Cyclopenta(def)ph...	18.95	2.3	ng/ul	121181	4	17.01	1051970	20.0
o-Terphenyl	20.95	2.4	ng/ul	251475	5	21.21	2135760	20.0
Triphenylene, 2-m...	21.76	2.5	ng/ul	263174	5	21.21	2135760	20.0
13-Docosenamide, ...	22.27	2.9	ng/ul	304830	5	21.21	2135760	20.0