

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032015\
 Data File : BG016088.D
 Acq On : 21 Mar 2015 5:11
 Operator : TP/IZ
 Sample : G1565-11
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AF1

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG031715.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	2.732	3	7	13	rVB2	31155	46894	0.75%	0.086%
2	2.944	38	43	51	rBV	296230	537475	8.62%	0.989%
3	3.020	51	56	72	rVB	624594	896949	14.39%	1.651%
4	4.923	375	380	386	rVB2	27961	46634	0.75%	0.086%
5	6.797	690	699	707	rBV5	7141	20805	0.33%	0.038%
6	6.962	720	727	740	rBV	307624	512675	8.22%	0.944%
7	7.132	749	756	764	rBV	244617	370063	5.94%	0.681%
8	7.332	780	790	800	rBV	307913	474223	7.61%	0.873%
9	7.426	800	806	813	rVB3	12248	23364	0.37%	0.043%
10	7.802	862	870	877	rBV	362624	566110	9.08%	1.042%
11	8.495	980	988	996	rBV	393767	609954	9.79%	1.123%
12	8.624	1003	1010	1014	rBV2	12230	23192	0.37%	0.043%
13	8.953	1060	1066	1073	rVV	266603	426926	6.85%	0.786%
14	9.670	1179	1188	1195	rBV	220361	369702	5.93%	0.681%
15	9.770	1199	1205	1209	rBV7	8557	17080	0.27%	0.031%
16	9.817	1209	1213	1219	rVB2	8414	17190	0.28%	0.032%
17	10.204	1270	1279	1289	rBV	360846	594055	9.53%	1.094%
18	10.357	1299	1305	1316	rBV	269563	411197	6.60%	0.757%
19	10.586	1333	1344	1349	rBV	520879	894388	14.35%	1.646%
20	10.633	1349	1352	1358	rVV	57000	93383	1.50%	0.172%
21	10.721	1361	1367	1381	rVV	110316	215118	3.45%	0.396%
22	11.491	1493	1498	1506	rVB7	10084	24155	0.39%	0.044%
23	11.773	1538	1546	1550	rBV7	8653	16269	0.26%	0.030%
24	12.243	1621	1626	1635	rVB	42124	70406	1.13%	0.130%
25	12.466	1658	1664	1673	rVB	44152	79955	1.28%	0.147%
26	12.777	1713	1717	1722	rVB	17286	24612	0.39%	0.045%
27	13.030	1755	1760	1767	rVB	28025	39860	0.64%	0.073%
28	13.253	1792	1798	1804	rVB2	33388	62922	1.01%	0.116%
29	13.588	1849	1855	1856	rBV2	36910	56523	0.91%	0.104%
30	13.741	1875	1881	1886	rVV	63190	113809	1.83%	0.209%
31	13.800	1886	1891	1892	rVV	47328	63037	1.01%	0.116%
32	13.835	1892	1897	1902	rVV	563280	804344	12.90%	1.481%
33	13.876	1902	1904	1908	rVB	37314	44773	0.72%	0.082%
34	13.952	1913	1917	1919	rVV5	12339	19851	0.32%	0.037%

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35	13.982	1919	1922	1925	rVV	31258	41998	0.67%	0.077%
36	14.011	1925	1927	1933	rVB5	16346	21749	0.35%	0.040%
37	14.117	1939	1945	1955	rVB	737448	1092517	17.53%	2.011%
38	14.258	1965	1969	1975	rVB8	14432	27312	0.44%	0.050%
39	14.322	1975	1980	1985	rBV	37453	52954	0.85%	0.097%
40	14.422	1988	1997	2003	rVV2	724314	1196227	19.19%	2.202%
41	14.487	2003	2008	2016	rVV	347435	517385	8.30%	0.952%
42	14.552	2016	2019	2023	rVV2	22600	34653	0.56%	0.064%
43	14.610	2023	2029	2041	rVV	231915	408297	6.55%	0.752%
44	14.734	2043	2050	2052	rVV3	23139	55083	0.88%	0.101%
45	14.763	2052	2055	2059	rVV3	35764	65071	1.04%	0.120%
46	14.822	2059	2065	2072	rVV	296898	470330	7.55%	0.866%
47	14.898	2074	2078	2082	rVV	41843	59540	0.96%	0.110%
48	14.945	2082	2086	2091	rVB6	13367	24793	0.40%	0.046%
49	15.039	2098	2102	2106	rBV	32857	50896	0.82%	0.094%
50	15.092	2106	2111	2115	rVB	32407	48962	0.79%	0.090%
51	15.215	2125	2132	2134	rBV6	31419	57519	0.92%	0.106%
52	15.239	2134	2136	2139	rVV3	25819	35534	0.57%	0.065%
53	15.274	2139	2142	2146	rVB	38110	45937	0.74%	0.085%
54	15.415	2157	2166	2170	rBV	896207	1278230	20.51%	2.353%
55	15.468	2170	2175	2180	rVV	548650	777816	12.48%	1.432%
56	15.527	2180	2185	2190	rVV	199373	275245	4.42%	0.507%
57	15.591	2193	2196	2199	rVV3	39993	47530	0.76%	0.087%
58	15.632	2199	2203	2207	rVV3	48372	70868	1.14%	0.130%
59	15.679	2208	2211	2214	rVB2	27909	31215	0.50%	0.057%
60	15.762	2221	2225	2233	rVB	124914	208037	3.34%	0.383%
61	15.838	2234	2238	2241	rBV6	17176	28299	0.45%	0.052%
62	15.908	2245	2250	2257	rVB	114011	246425	3.95%	0.454%
63	15.979	2258	2262	2263	rBV4	21778	28609	0.46%	0.053%
64	16.132	2284	2288	2297	rVB4	85443	173407	2.78%	0.319%
65	16.273	2309	2312	2316	rBV2	18556	25660	0.41%	0.047%
66	16.472	2342	2346	2350	rVB2	91288	131512	2.11%	0.242%
67	16.519	2350	2354	2359	rVB3	49912	72453	1.16%	0.133%
68	16.619	2368	2371	2376	rVB	51394	67590	1.08%	0.124%
69	16.696	2382	2384	2387	rBV2	30985	29874	0.48%	0.055%
70	16.737	2388	2391	2395	rVB2	43576	47449	0.76%	0.087%
71	16.819	2401	2405	2411	rVB2	92796	137688	2.21%	0.253%

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72	16.895	2414	2418	2420	rBV	56751	87906	1.41%	0.162%
73	16.983	2428	2433	2442	rVB	214631	356605	5.72%	0.656%
74	17.166	2458	2464	2467	rBV	771684	1191408	19.11%	2.193%
75	17.207	2467	2471	2476	rVV	3002356	4256416	68.29%	7.835%
76	17.260	2476	2480	2483	rVV2	910238	1350538	21.67%	2.486%
77	17.295	2483	2486	2491	rVV	932004	1244003	19.96%	2.290%
78	17.354	2491	2496	2501	rVV2	108055	182230	2.92%	0.335%
79	17.565	2527	2532	2538	rVB	349352	497390	7.98%	0.916%
80	17.659	2545	2548	2550	rBV2	44797	54465	0.87%	0.100%
81	18.035	2608	2612	2614	rBV2	302522	429261	6.89%	0.790%
82	18.082	2617	2620	2627	rVB	457608	683008	10.96%	1.257%
83	18.164	2630	2634	2639	rBV	172144	216102	3.47%	0.398%
84	18.235	2640	2646	2650	rBV2	761982	1261821	20.24%	2.323%
85	18.481	2685	2688	2692	rBV5	90536	102554	1.65%	0.189%
86	18.534	2694	2697	2701	rVV	245954	337983	5.42%	0.622%
87	18.575	2701	2704	2710	rVB2	135577	225045	3.61%	0.414%
88	18.951	2765	2768	2771	rBV	173347	248407	3.99%	0.457%
89	19.098	2790	2793	2799	rBV2	140952	217749	3.49%	0.401%
90	19.222	2808	2814	2820	rBV	4181282	6233146	100.00%	11.474%
91	19.556	2866	2871	2873	rBV3	1625392	2680002	43.00%	4.933%
92	19.586	2873	2876	2884	rVV	3574018	4878106	78.26%	8.979%
93	19.650	2884	2887	2893	rVB2	176121	293632	4.71%	0.541%
94	19.821	2913	2916	2920	rVB	265062	344402	5.53%	0.634%
95	19.897	2926	2929	2936	rBV4	204605	446059	7.16%	0.821%
96	20.073	2955	2959	2965	rVB	802285	1190957	19.11%	2.192%
97	20.173	2973	2976	2980	rVB	728522	810165	13.00%	1.491%
98	21.325	3168	3172	3178	rBV3	2499953	4741682	76.07%	8.728%
99	21.378	3178	3181	3189	rVB	2084687	2640781	42.37%	4.861%
100	21.489	3197	3200	3204	rBV2	473976	550722	8.84%	1.014%

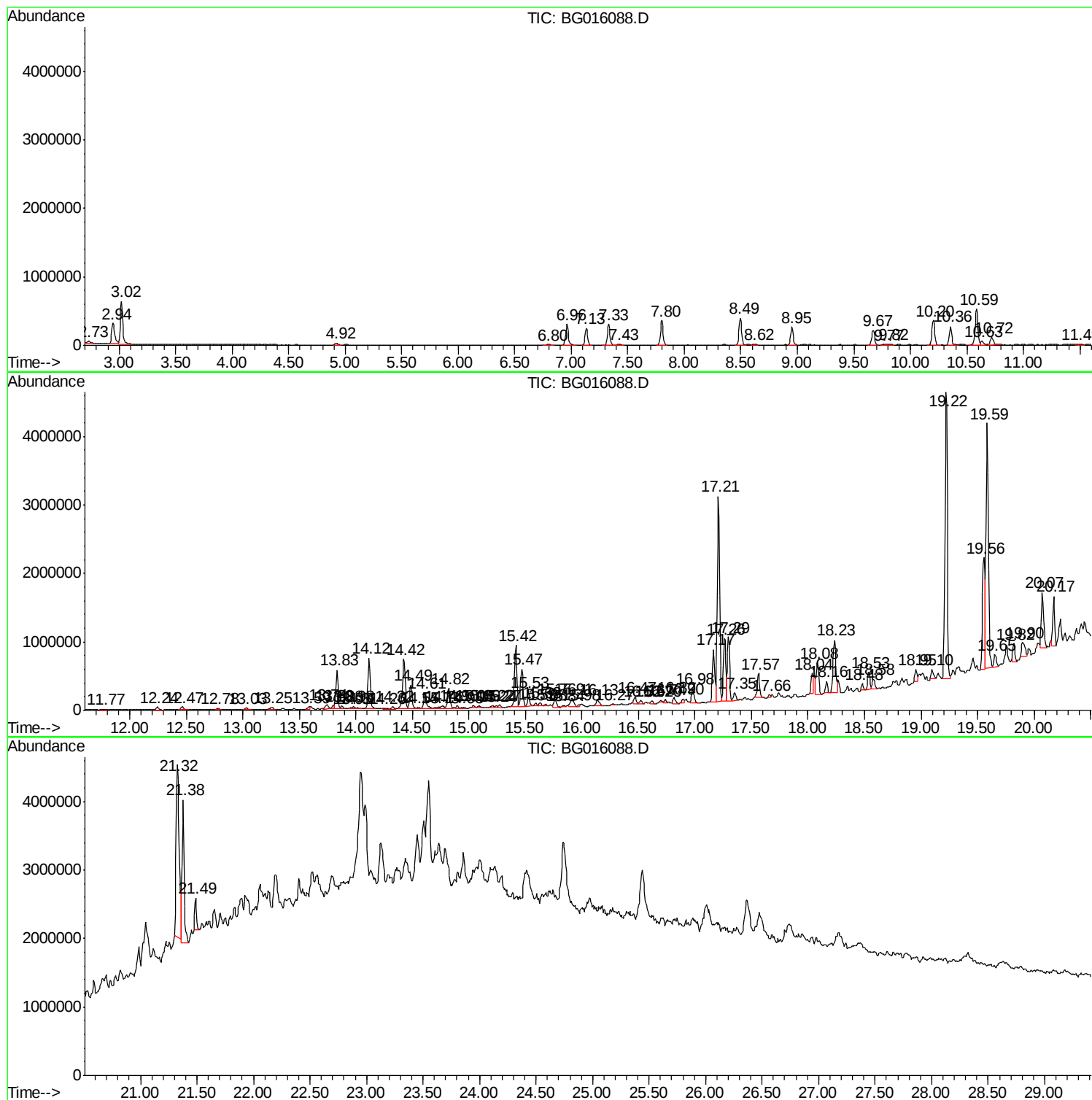
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG031715.M
Quant Title : SVOA CALIBRATION

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



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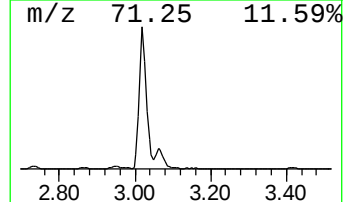
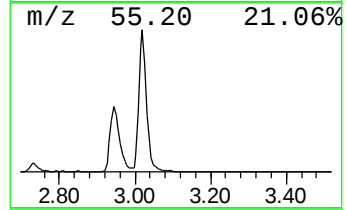
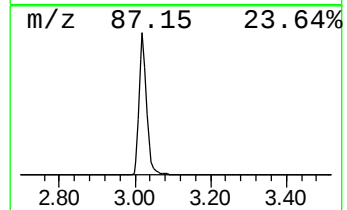
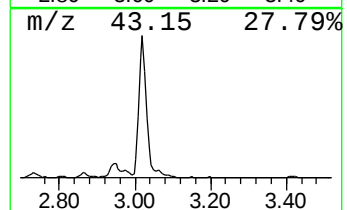
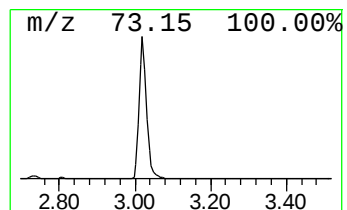
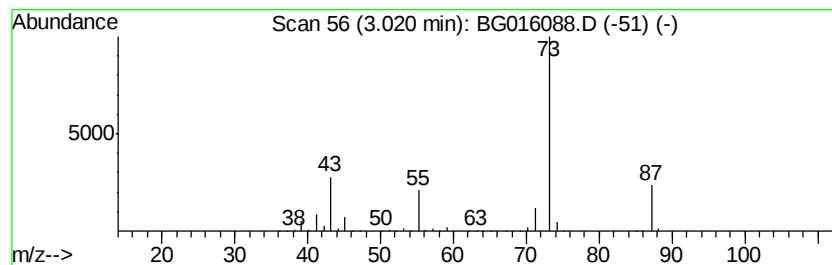
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG031715.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.02	31.69 ng/ul	896949	1,4-Dichlorobenzene-d4	7.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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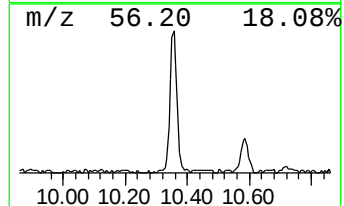
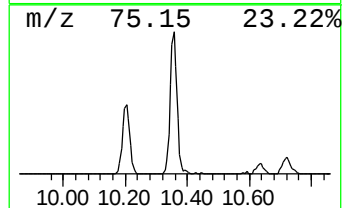
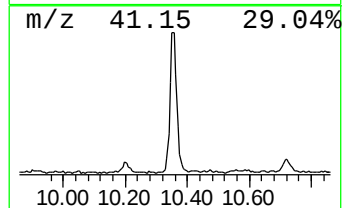
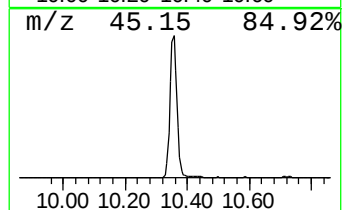
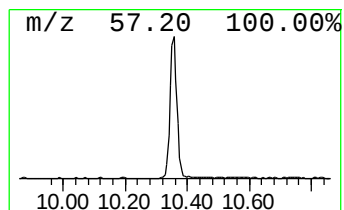
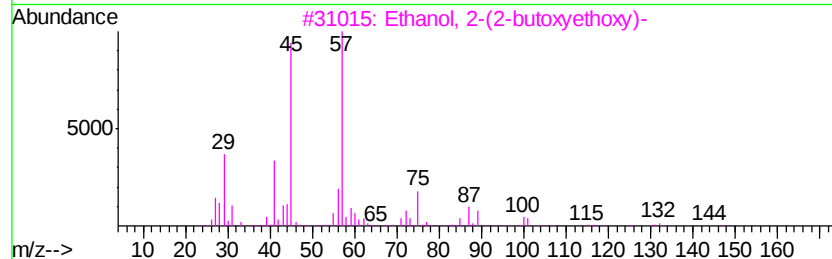
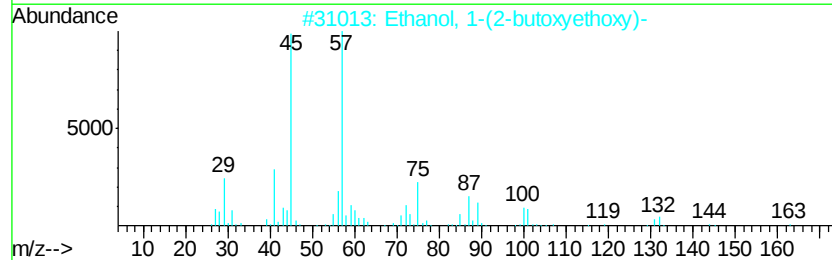
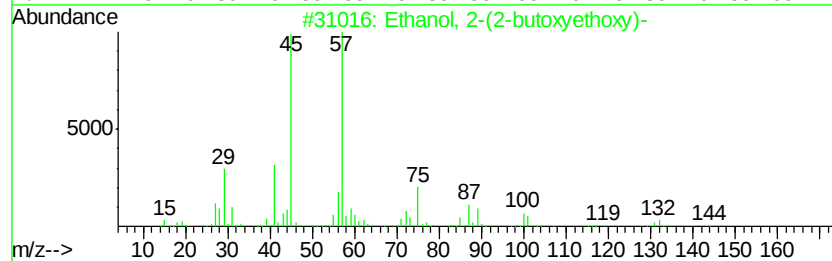
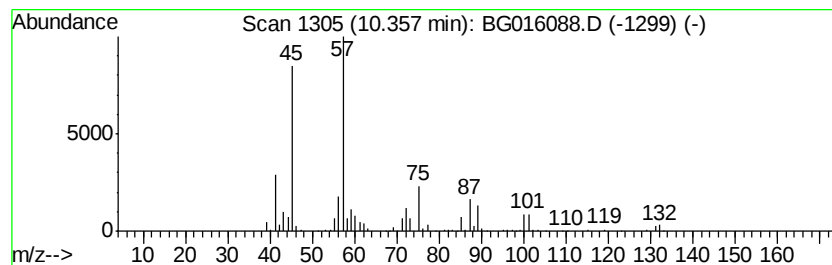
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG031715.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.36	9.20 ng/ul	411197	Naphthalene-d8	10.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2		Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
3		Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4		Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
5		Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83



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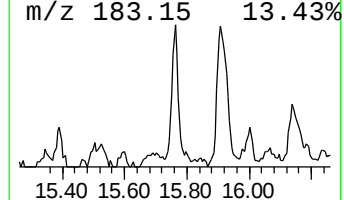
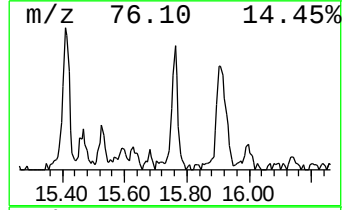
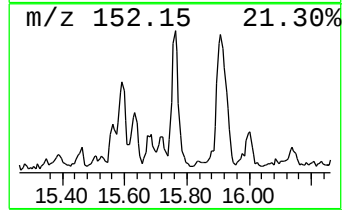
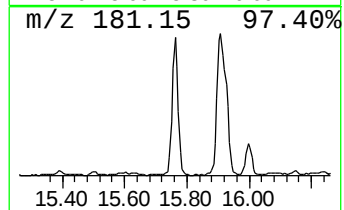
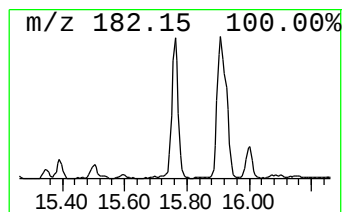
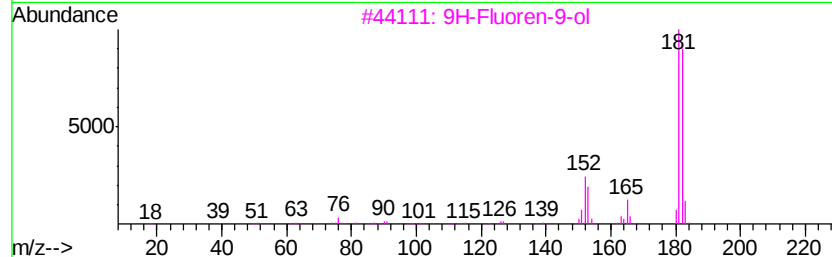
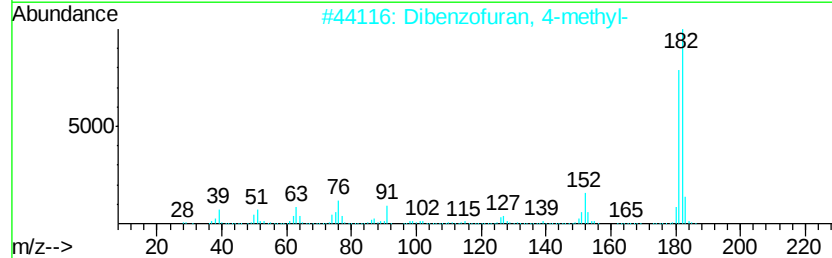
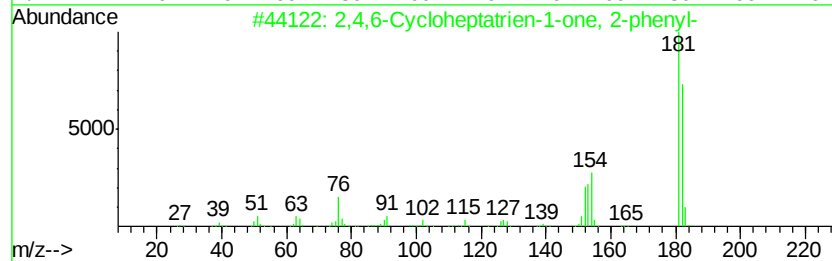
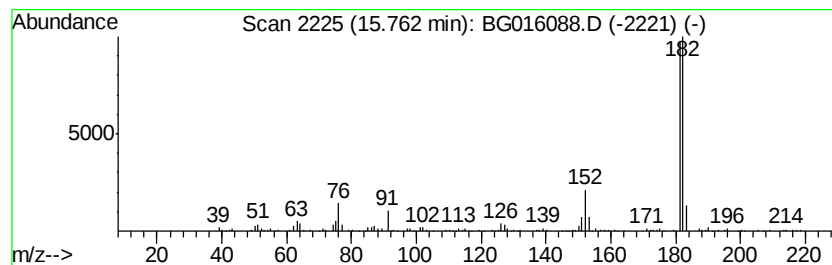
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 2,4,6-Cycloheptatrien-1-one... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.76	3.48 ng/ul	208037	Acenaphthene-d10	14.42

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	91
2			Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	91
3			9H-Fluoren-9-ol	182	C13H10O	001689-64-1	78
4			9H-Fluoren-9-ol	182	C13H10O	001689-64-1	72
5			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	64



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ClientSampleId :
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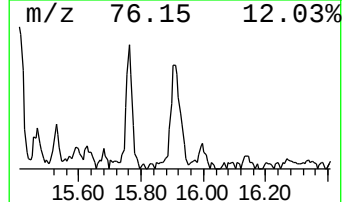
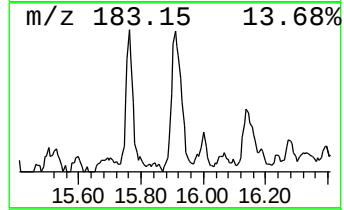
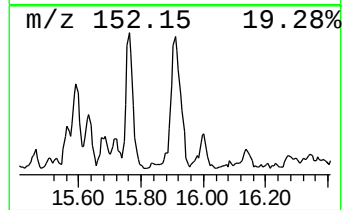
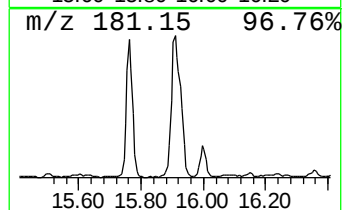
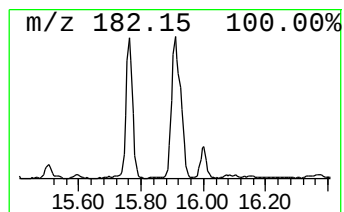
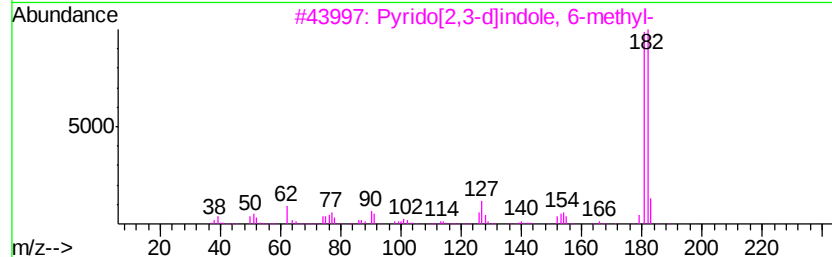
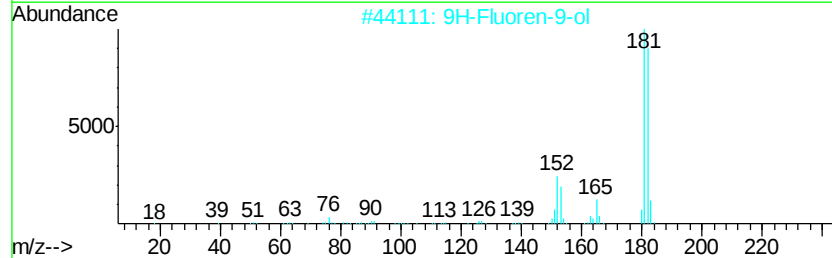
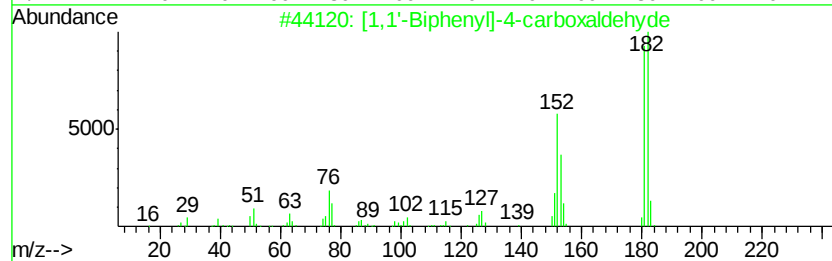
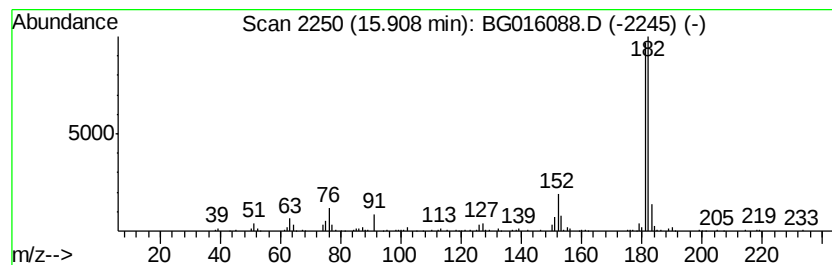
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 [1,1'-Biphenyl]-4-carboxald... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.91	4.14 ng/ul	246425	Phenanthrene-d10	17.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	80
2		9H-Fluoren-9-ol	182	C13H10O	001689-64-1	78
3		Pyrido[2,3-d]indole, 6-methyl-	182	C12H10N2	108349-67-3	72
4		Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	70
5		9H-Xanthene	182	C13H10O	000092-83-1	60



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032015\
Data File : BG016088.D
Acq On : 21 Mar 2015 5:11
Operator : TP/IZ
Sample : G1565-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleID :
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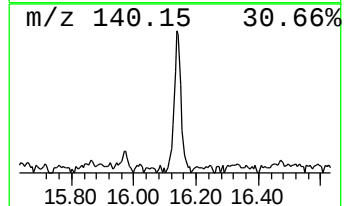
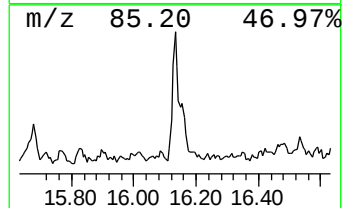
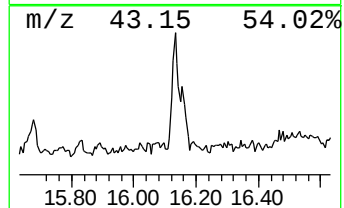
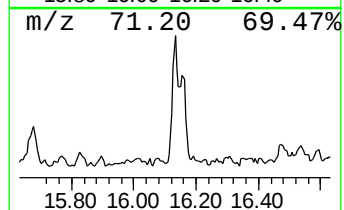
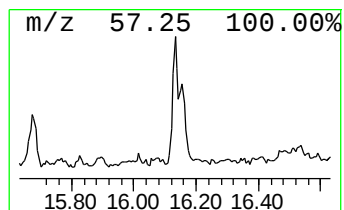
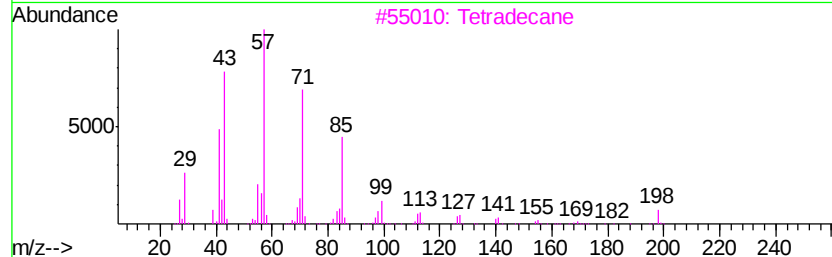
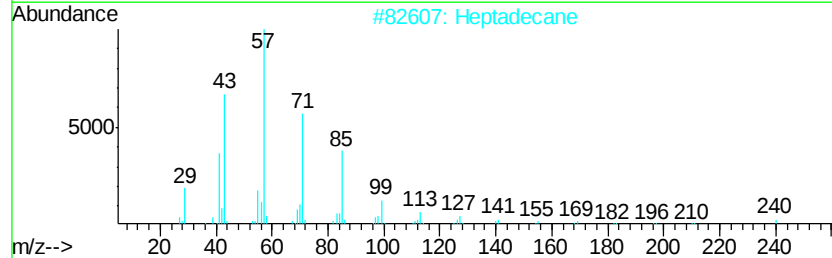
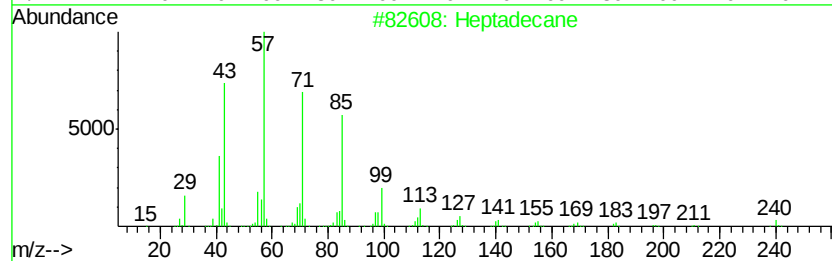
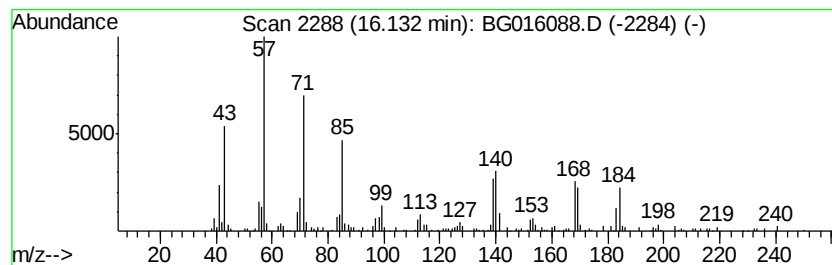
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Heptadecane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.13	2.91 ng/ul	173407	Phenanthrene-d10	17.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	86
2	Heptadecane		240	C17H36	000629-78-7	83
3	Tetradecane		198	C14H30	000629-59-4	81
4	Dodecane		170	C12H26	000112-40-3	81
5	Tetradecane		198	C14H30	000629-59-4	50



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032015\
Data File : BG016088.D
Acq On : 21 Mar 2015 5:11
Operator : TP/IZ
Sample : G1565-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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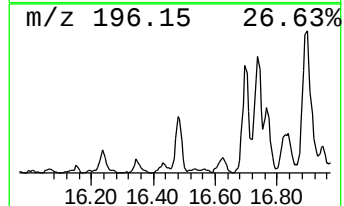
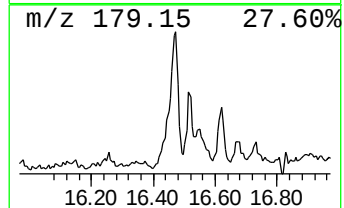
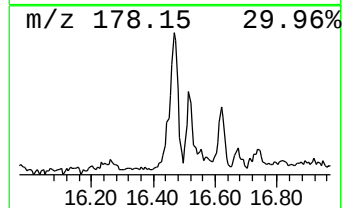
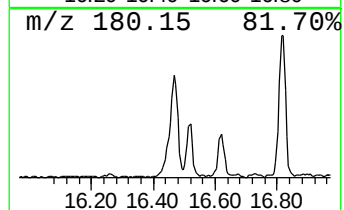
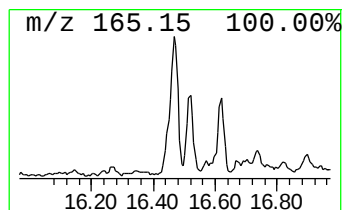
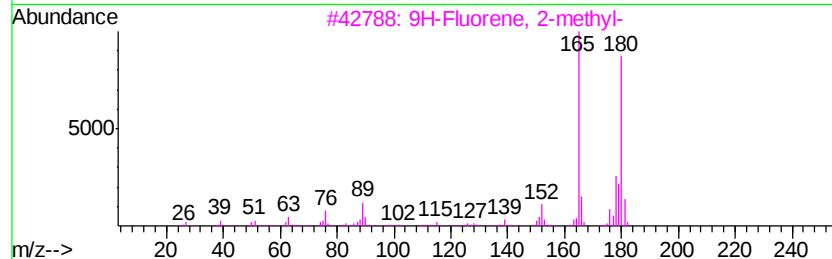
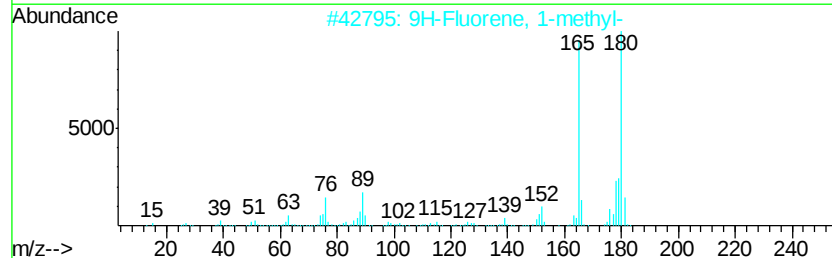
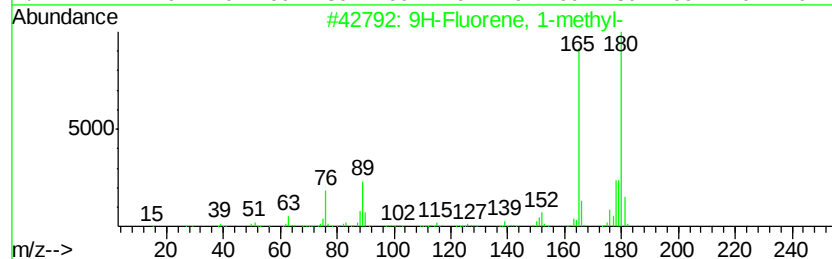
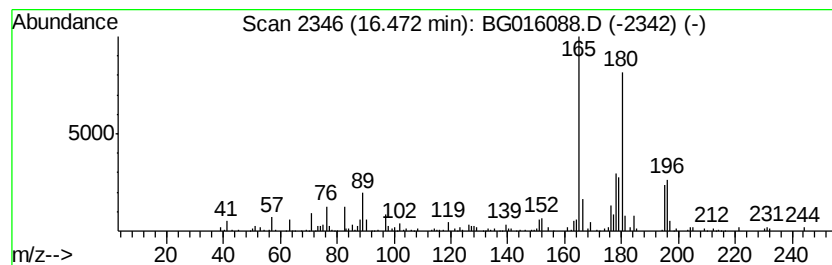
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9H-Fluorene, 1-methyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.47	2.21 ng/ul	131512	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	86
2		9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	86
3		9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	86
4		9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	86
5		9H-Fluorene, 9-methyl-	180	C14H12	002523-37-7	86



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032015\
Data File : BG016088.D
Acq On : 21 Mar 2015 5:11
Operator : TP/IZ
Sample : G1565-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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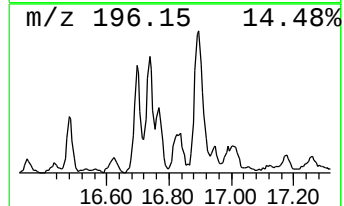
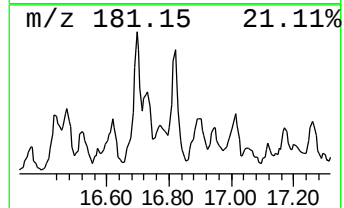
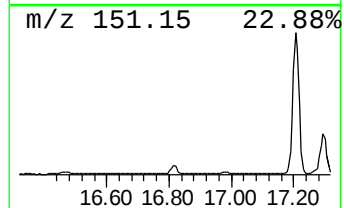
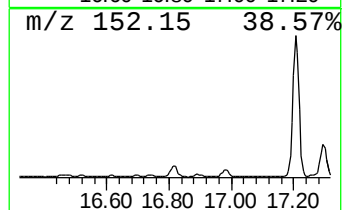
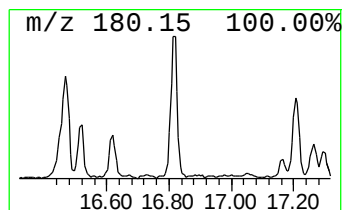
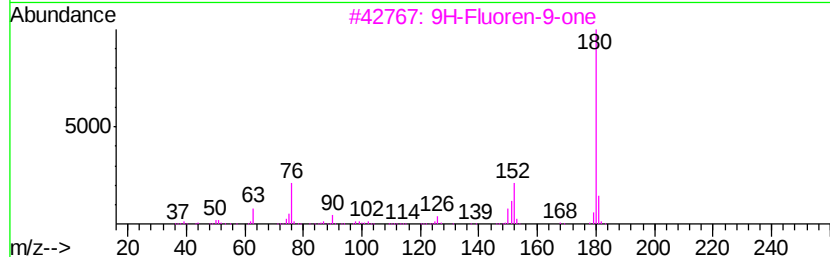
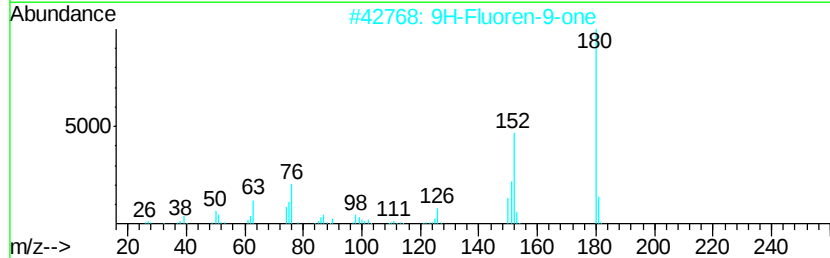
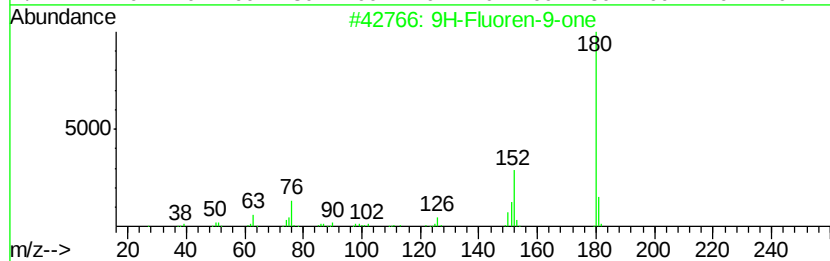
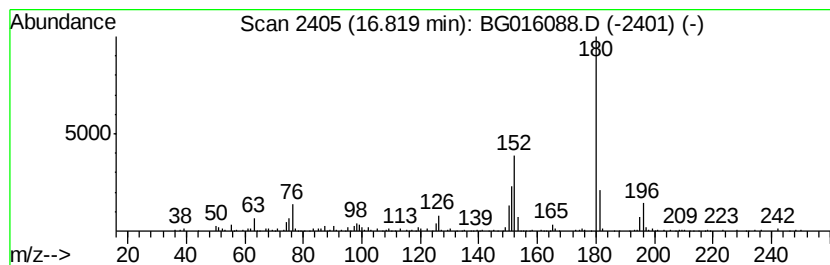
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 9H-Fluoren-9-one Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.82	2.31 ng/ul	137688	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9H-Fluoren-9-one	180	C13H8O	000486-25-9	94
2		9H-Fluoren-9-one	180	C13H8O	000486-25-9	91
3		9H-Fluoren-9-one	180	C13H8O	000486-25-9	87
4		9,10-Phenanthrenedione	208	C14H8O2	000084-11-7	80
5		Benzo[h]cinnoline	180	C12H8N2	000230-31-9	59



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032015\
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Sample : G1565-11
Misc :
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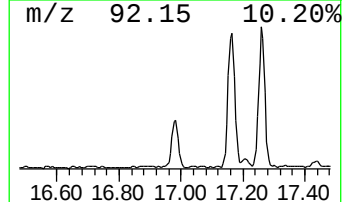
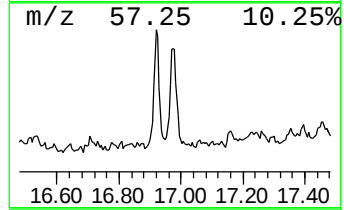
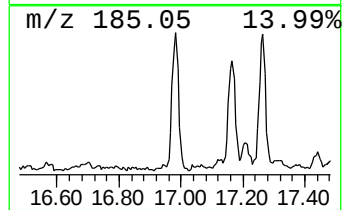
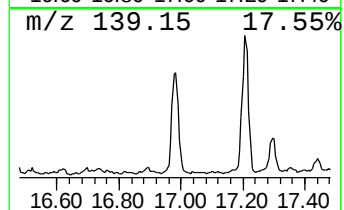
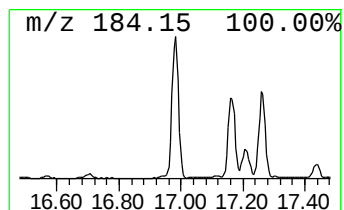
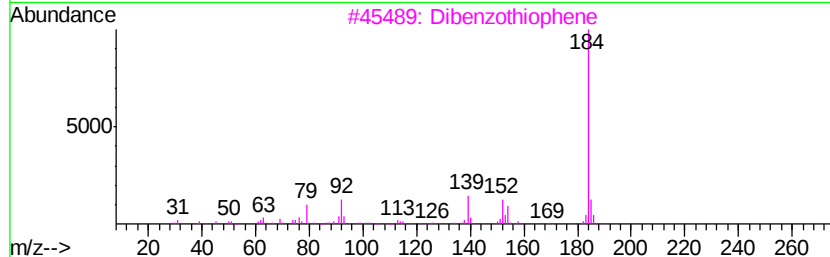
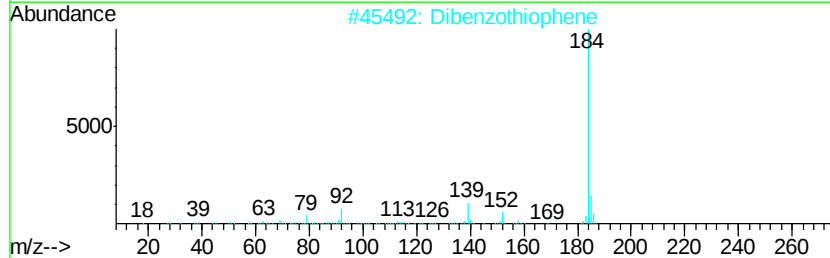
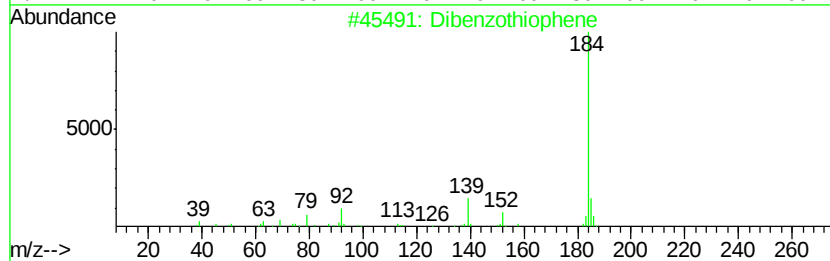
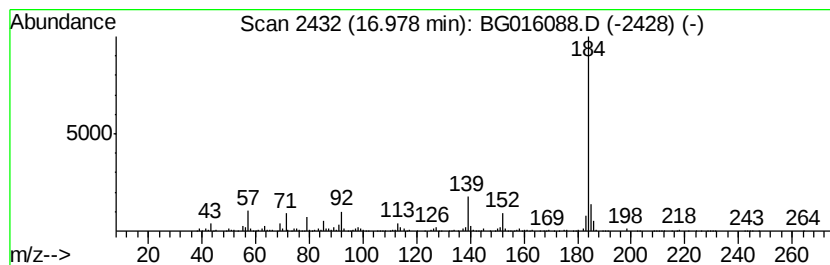
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG031715.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 Dibenzothiophene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.98	5.99 ng/ul	356605	Phenanthrene-d10	17.17	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzothiophene	184	C12H8S	000132-65-0	96
2	Dibenzothiophene	184	C12H8S	000132-65-0	95
3	Dibenzothiophene	184	C12H8S	000132-65-0	94
4	Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	94
5	Dibenzothiophene	184	C12H8S	000132-65-0	93



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032015\
Data File : BG016088.D
Acq On : 21 Mar 2015 5:11
Operator : TP/IZ
Sample : G1565-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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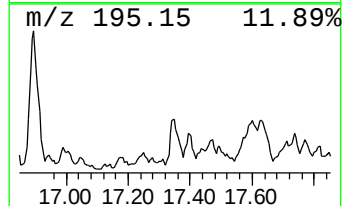
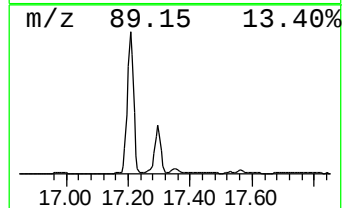
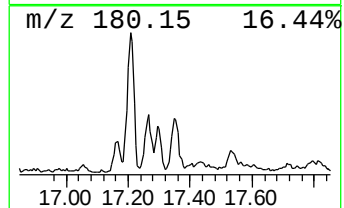
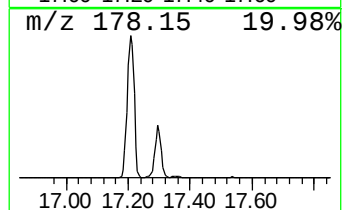
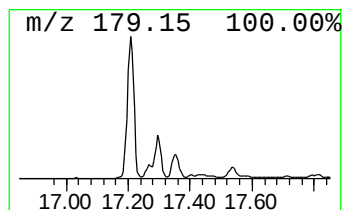
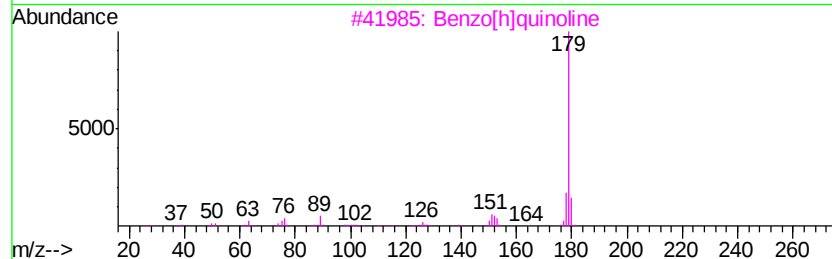
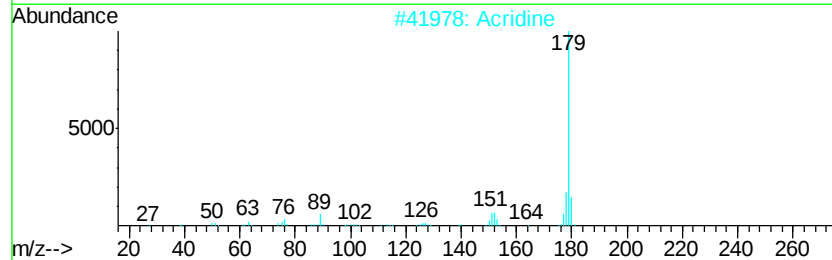
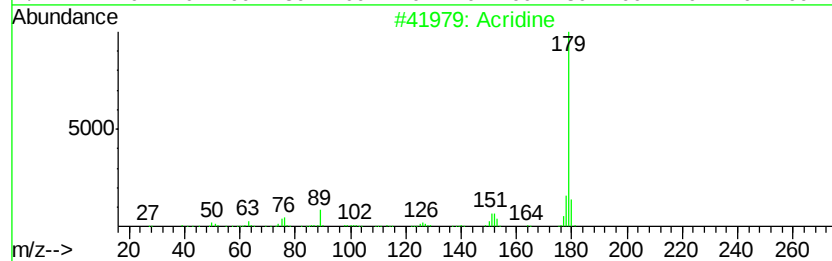
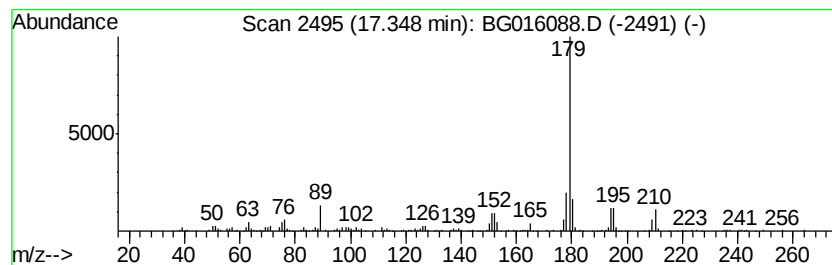
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Acridine Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.35	3.06 ng/ul	182230	Phenanthrene-d10	17.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acridine	179	C13H9N	000260-94-6	96
2			Acridine	179	C13H9N	000260-94-6	93
3			Benzo[h]quinoline	179	C13H9N	000230-27-3	93
4			Benzo[g]quinoline	179	C13H9N	000260-36-6	89
5			p-Phenylbenzonitrile	179	C13H9N	002920-38-9	83



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032015\
Data File : BG016088.D
Acq On : 21 Mar 2015 5:11
Operator : TP/IZ
Sample : G1565-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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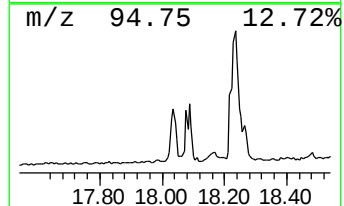
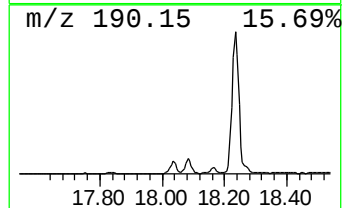
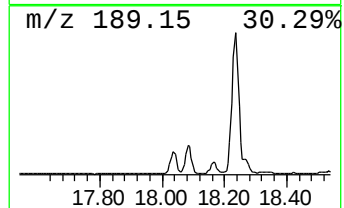
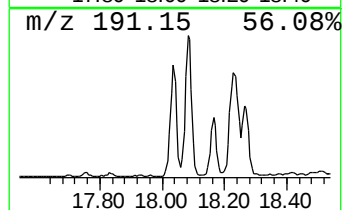
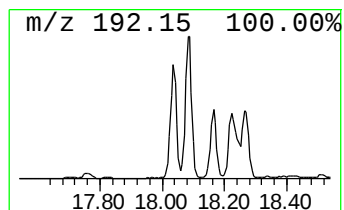
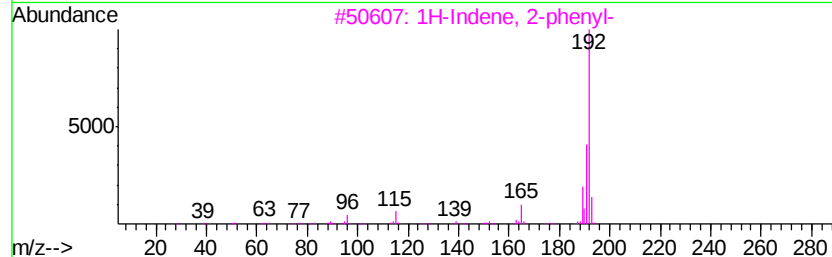
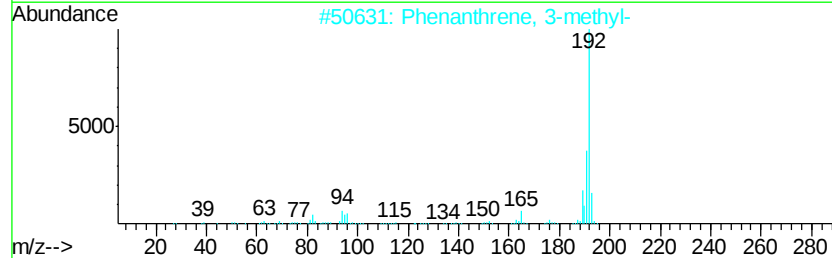
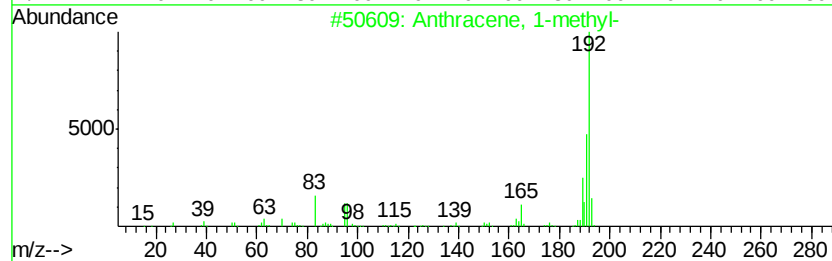
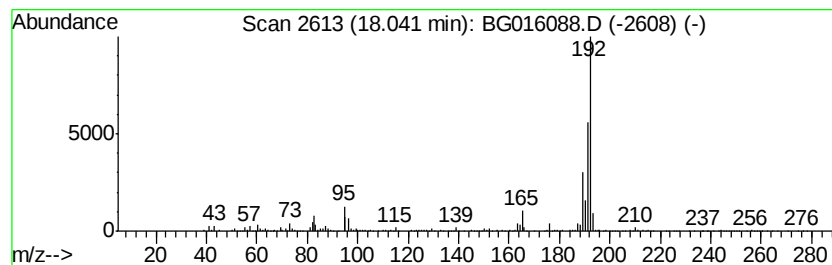
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Anthracene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.04	7.21 ng/ul	429261	Phenanthrene-d10	17.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 1-methyl-	192	C15H12	000610-48-0	72
2			Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	70
3			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	68
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	68
5			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	64



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Data File : BG016088.D
Acq On : 21 Mar 2015 5:11
Operator : TP/IZ
Sample : G1565-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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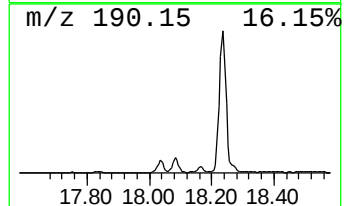
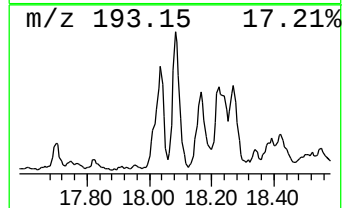
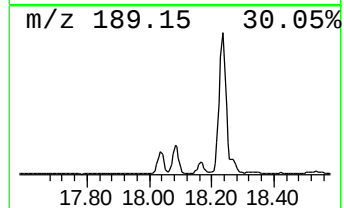
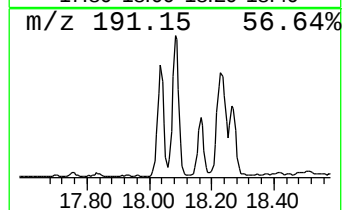
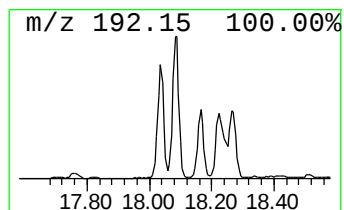
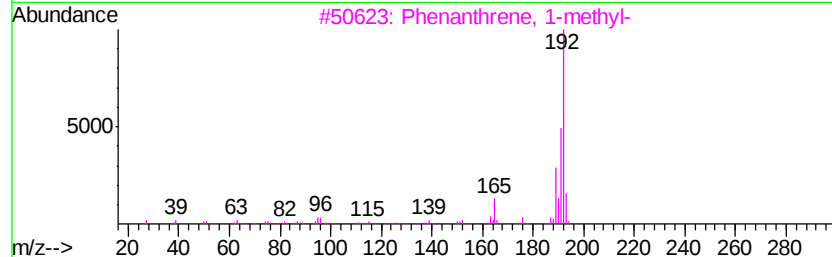
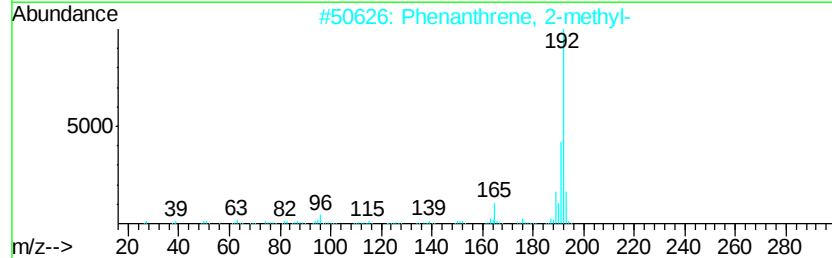
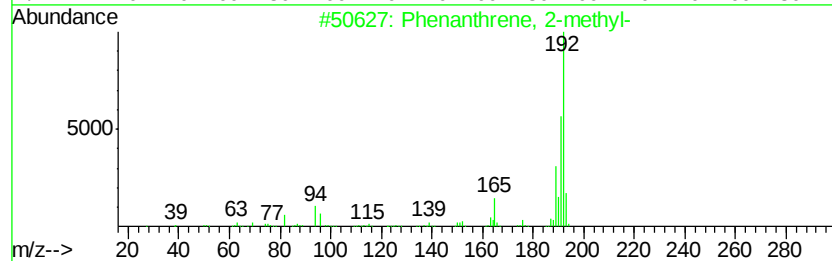
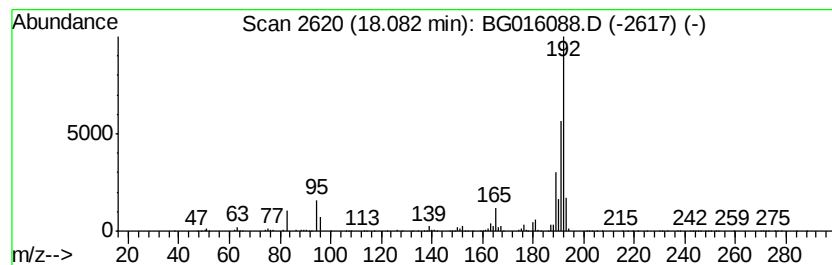
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Phenanthrene, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.08	11.47 ng/ul	683008	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
5		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032015\
Data File : BG016088.D
Acq On : 21 Mar 2015 5:11
Operator : TP/IZ
Sample : G1565-11
Misc :
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Instrument :
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ClientSampleId :
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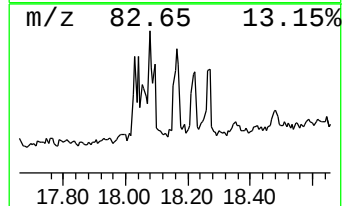
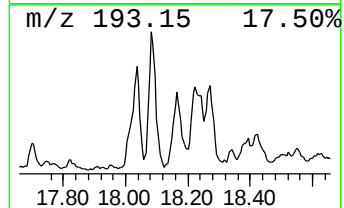
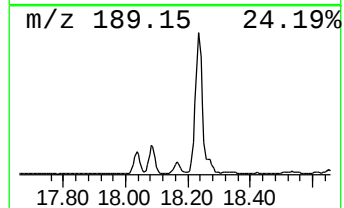
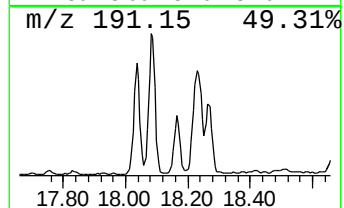
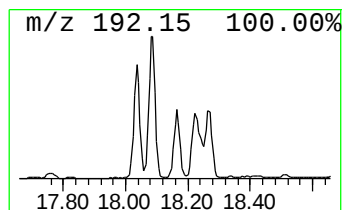
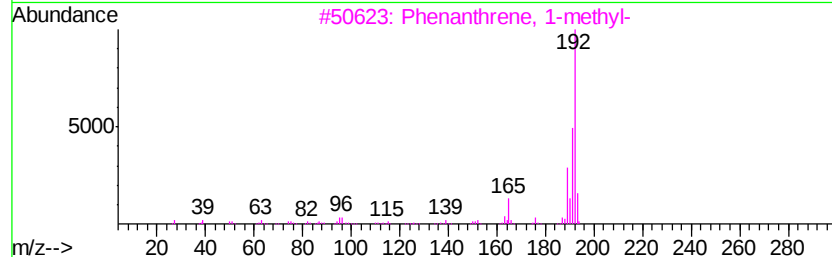
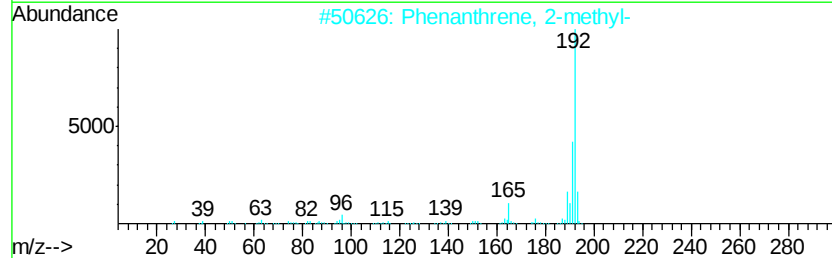
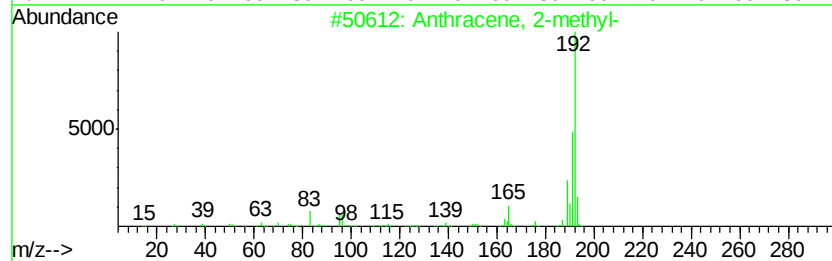
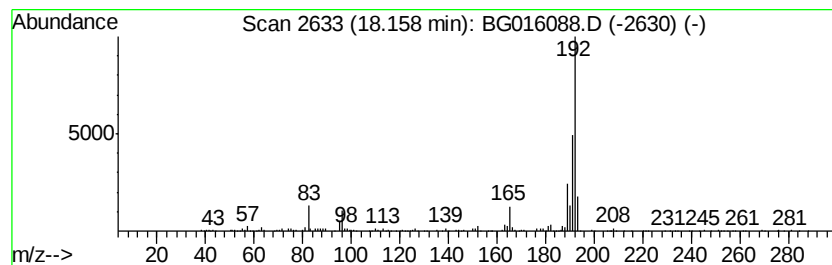
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Anthracene, 2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.16	3.63 ng/ul	216102	Phenanthrene-d10	17.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	97
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
4			1H-Cyclopropa[1]phenanthrene, 1a, ...	192	C15H12	000949-41-7	96
5			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032015\
Data File : BG016088.D
Acq On : 21 Mar 2015 5:11
Operator : TP/IZ
Sample : G1565-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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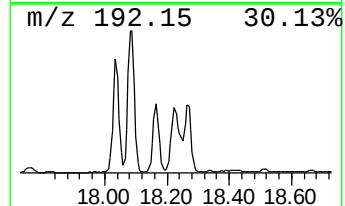
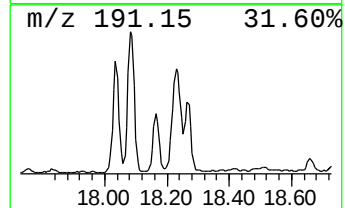
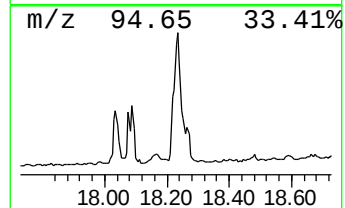
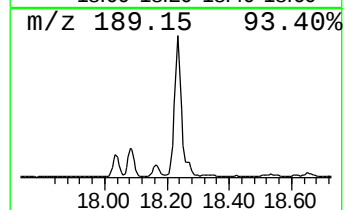
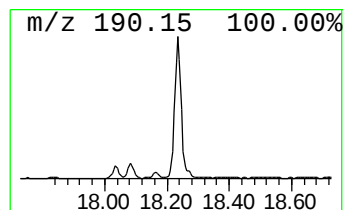
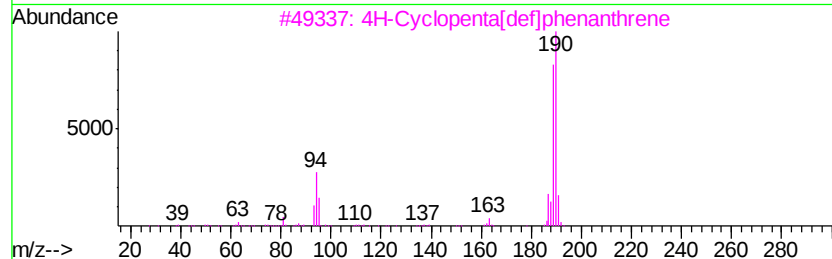
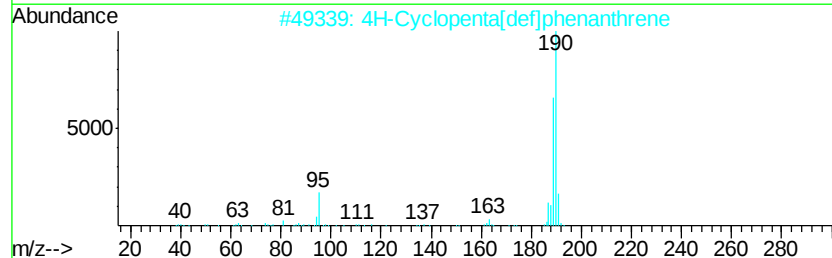
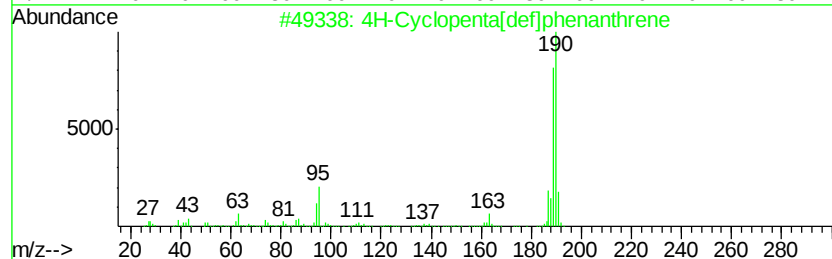
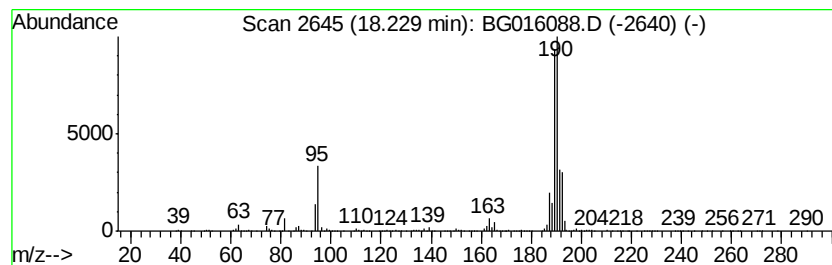
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.23	21.18 ng/ul	1261820	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	91
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	81
3		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	81
4		2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	47
5		Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	38



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032015\
Data File : BG016088.D
Acq On : 21 Mar 2015 5:11
Operator : TP/IZ
Sample : G1565-11
Misc :
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Instrument :
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ClientSampleId :
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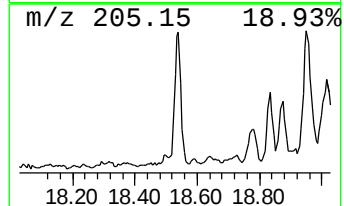
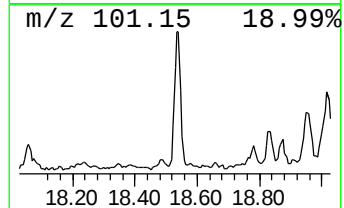
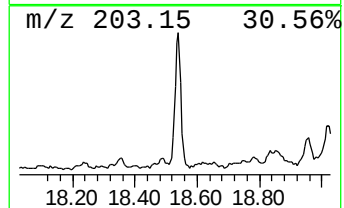
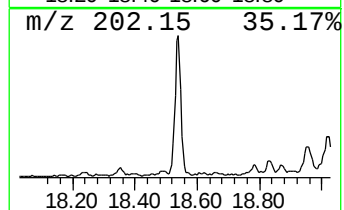
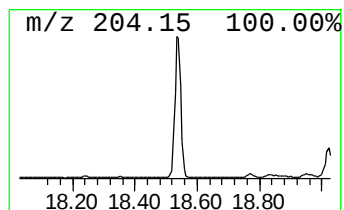
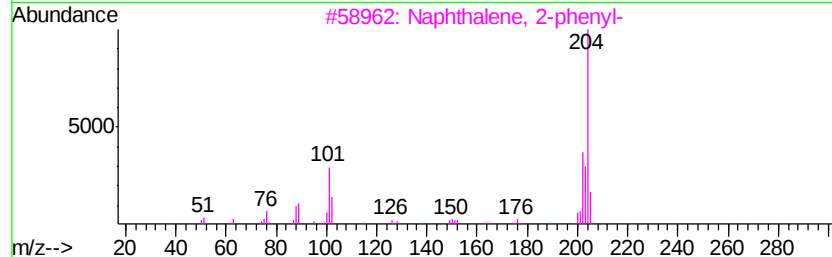
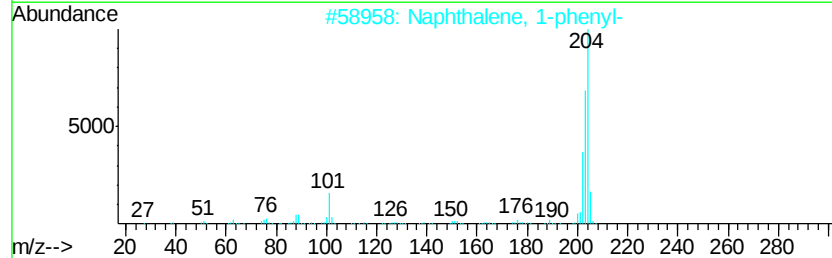
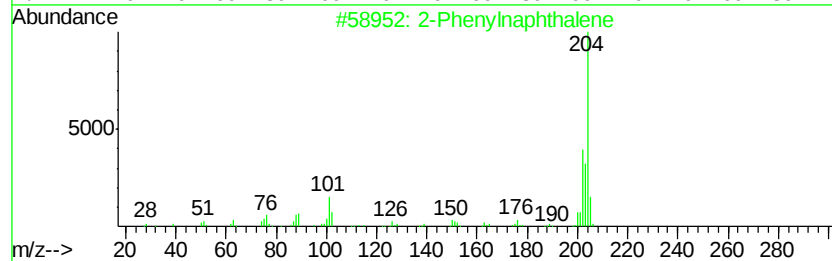
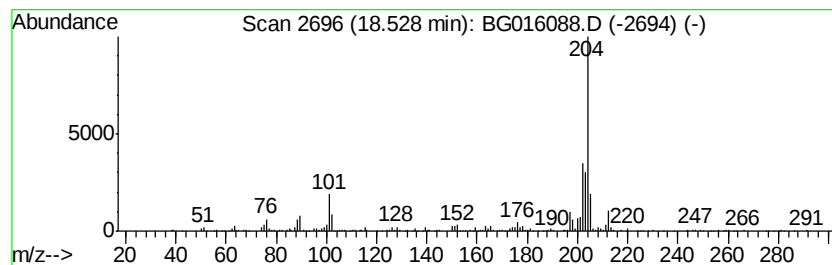
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 2-Phenylnaphthalene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.53	5.67 ng/ul	337983	Phenanthrene-d10	17.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Phenylnaphthalene	204	C16H12	035465-71-5	96
2			Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	95
3			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	94
4			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	94
5			9,10-di(Chloromethyl)anthracene	274	C16H12Cl2	010387-13-0	91



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032015\
Data File : BG016088.D
Acq On : 21 Mar 2015 5:11
Operator : TP/IZ
Sample : G1565-11
Misc :
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Instrument :
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ClientSampleId :
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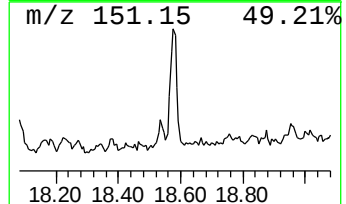
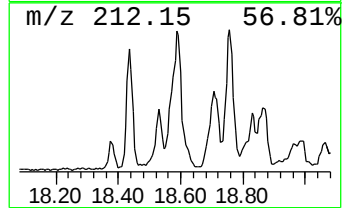
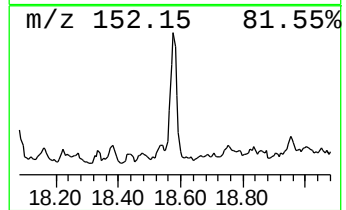
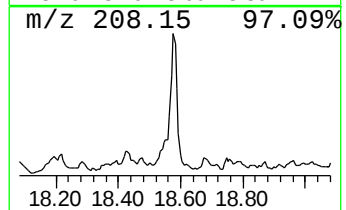
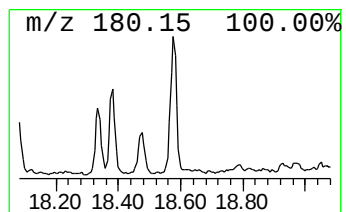
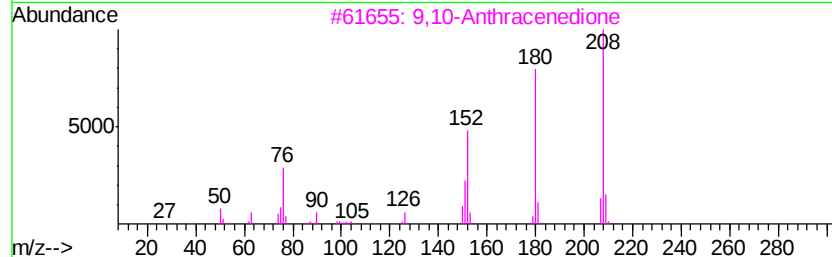
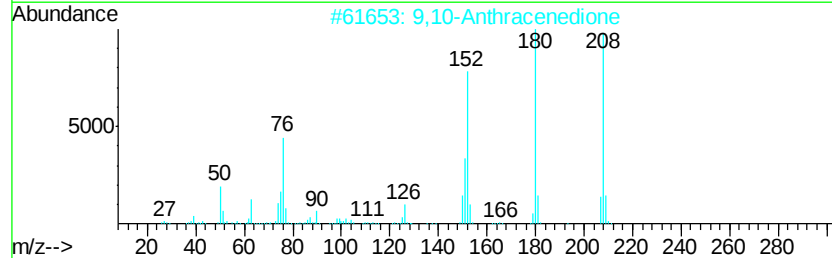
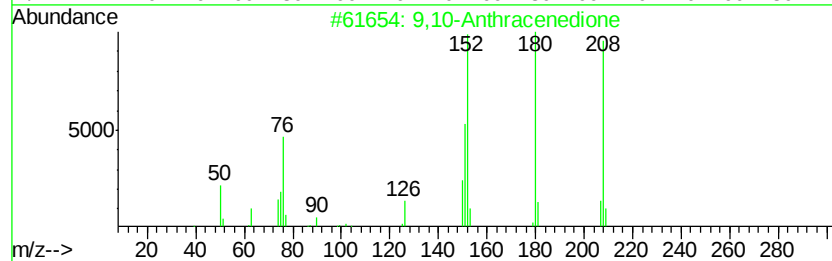
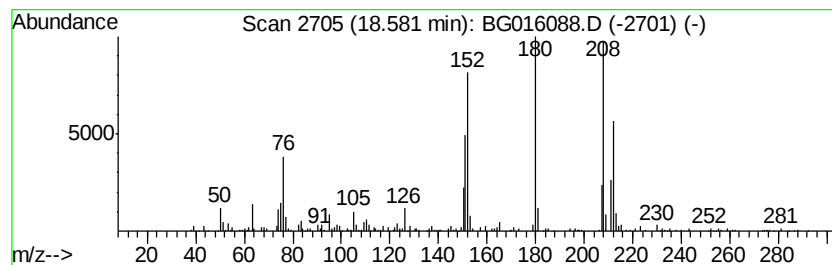
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 9,10-Anthracenedione Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.58	3.78 ng/ul	225045	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
2		9,10-Anthracenedione	208	C14H8O2	000084-65-1	93
3		9,10-Anthracenedione	208	C14H8O2	000084-65-1	90
4		9H-Fluoren-9-one	180	C13H8O	000486-25-9	87
5		9H-Fluoren-9-one	180	C13H8O	000486-25-9	60



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032015\
Data File : BG016088.D
Acq On : 21 Mar 2015 5:11
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Misc :
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Instrument :
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ClientSampled :
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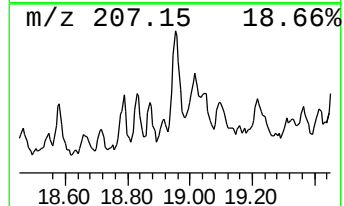
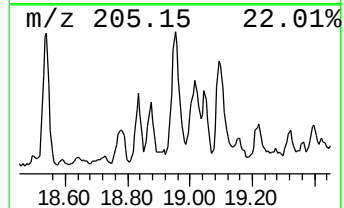
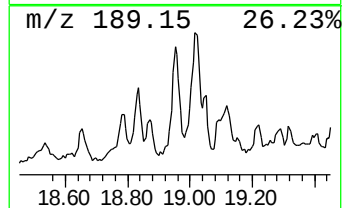
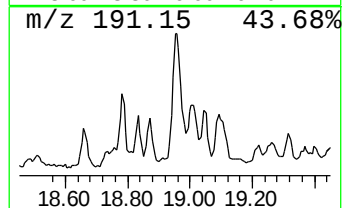
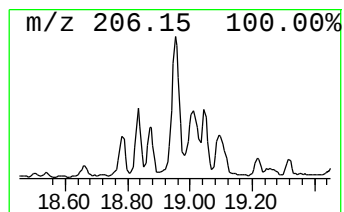
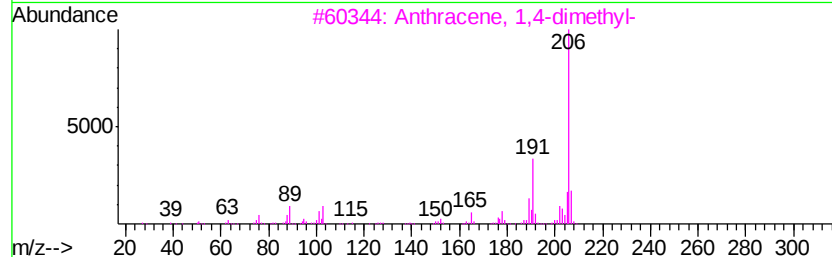
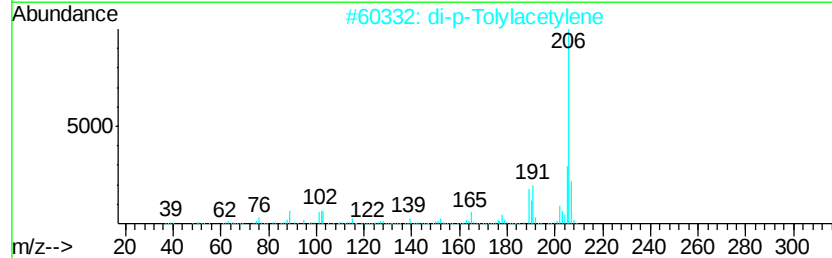
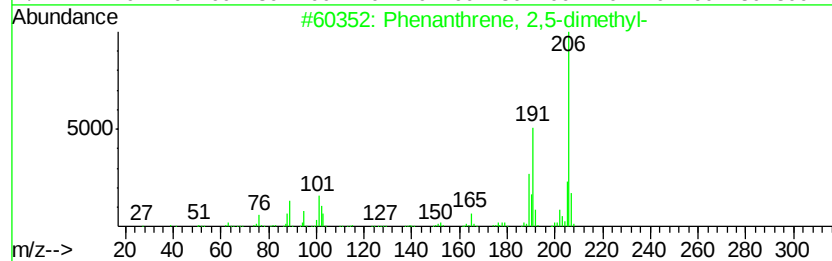
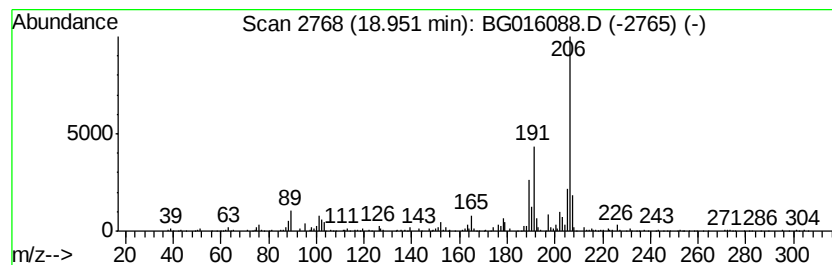
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Phenanthrene, 2,5-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.95	4.17 ng/ul	248407	Phenanthrene-d10	17.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	97
2			di-p-Tolylacetylene	206	C16H14	002789-88-0	93
3			Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	93
4			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	93
5			9,10-Dimethylantracene	206	C16H14	000781-43-1	91



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032015\
Data File : BG016088.D
Acq On : 21 Mar 2015 5:11
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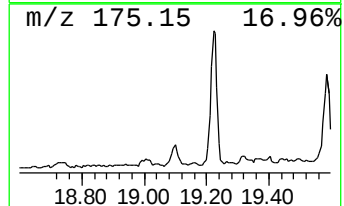
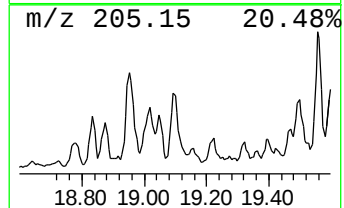
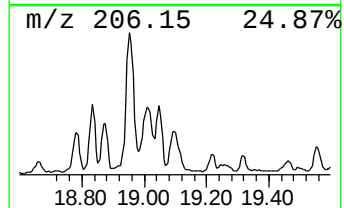
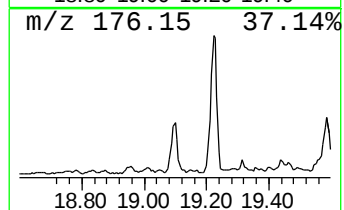
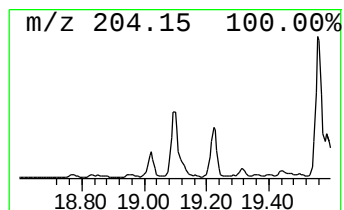
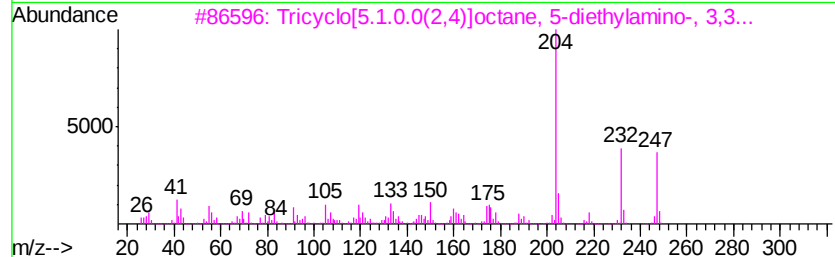
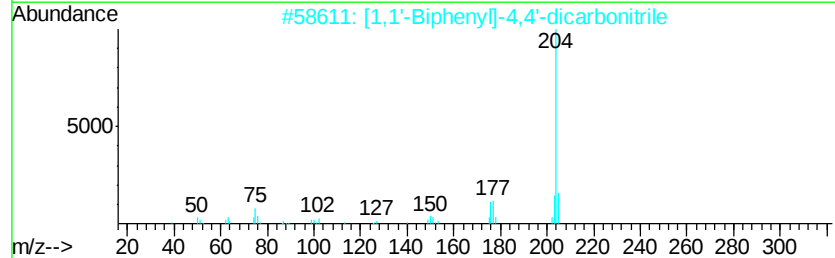
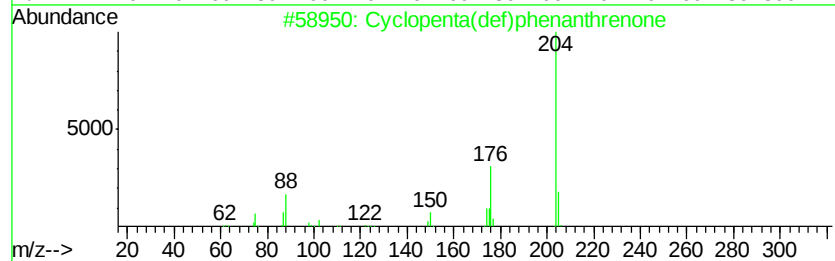
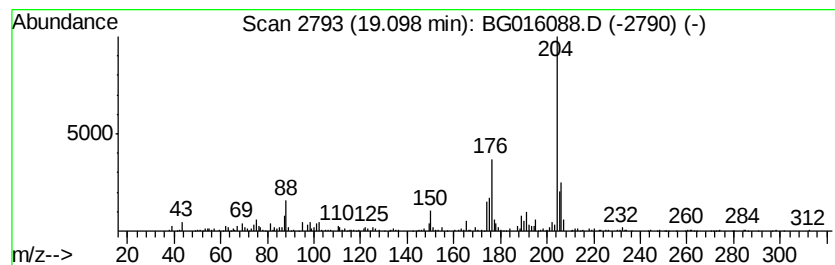
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 Cyclopenta(def)phenanthrenone Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.10	3.66 ng/ul	217749	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	96
2		[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	53
3		Tricyclo[5.1.0.0(2,4)]octane, 5-...	247	C17H29N	1000063-59-3	50
4		[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	49
5		[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	47



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032015\
Data File : BG016088.D
Acq On : 21 Mar 2015 5:11
Operator : TP/IZ
Sample : G1565-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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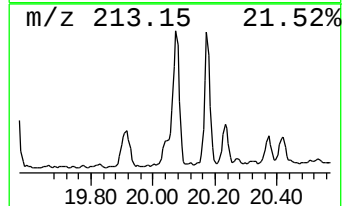
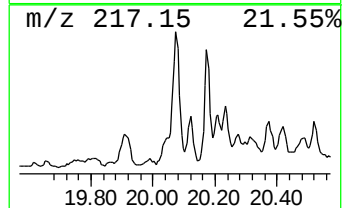
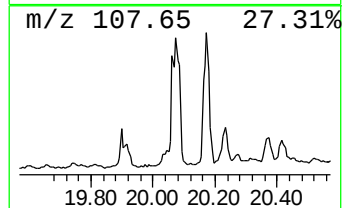
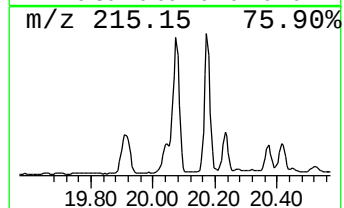
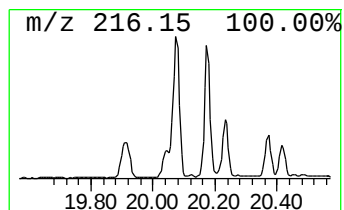
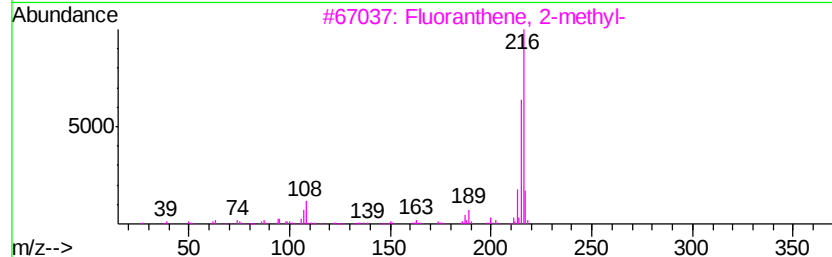
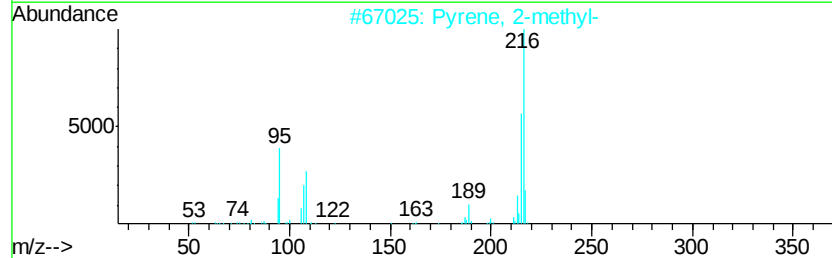
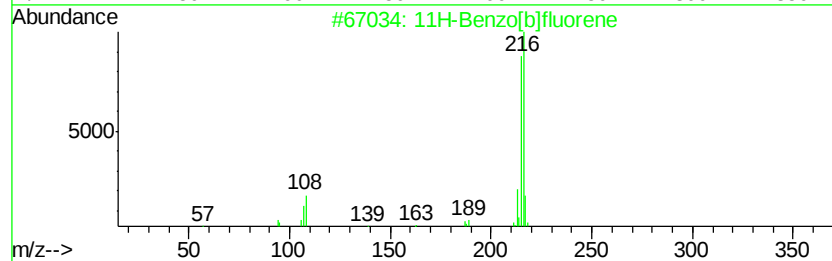
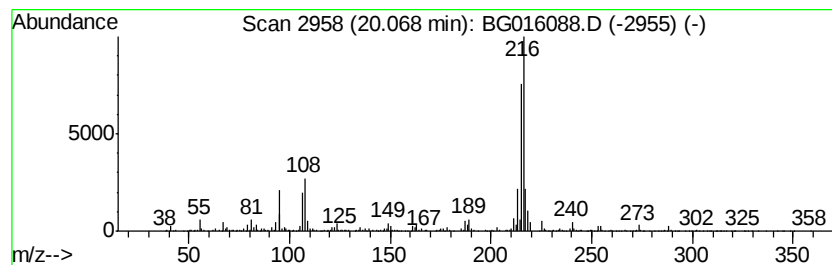
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 11H-Benzo[b]fluorene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.07	5.02 ng/ul	1190960	Chrysene-d12	21.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
2		Pyrene, 2-methyl-	216	C17H12	003442-78-2	91
3		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	90
4		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
5		7H-Benzo[c]fluorene	216	C17H12	000205-12-9	87



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032015\
Data File : BG016088.D
Acq On : 21 Mar 2015 5:11
Operator : TP/IZ
Sample : G1565-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

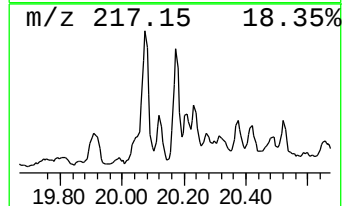
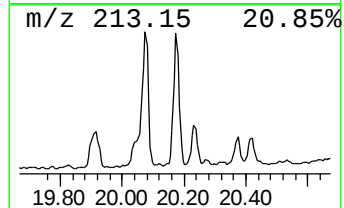
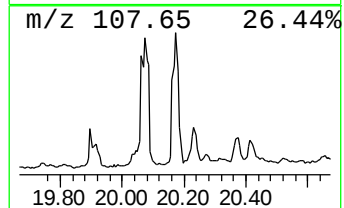
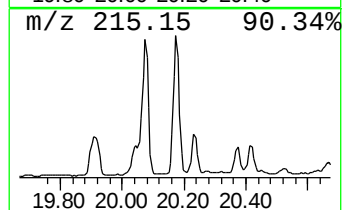
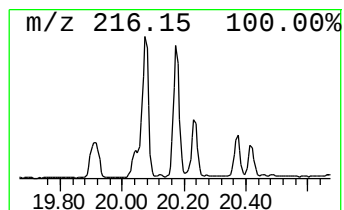
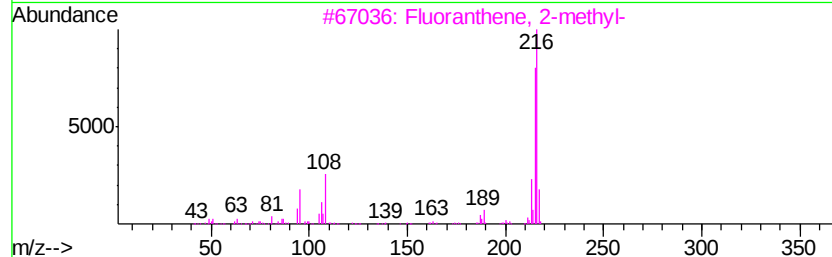
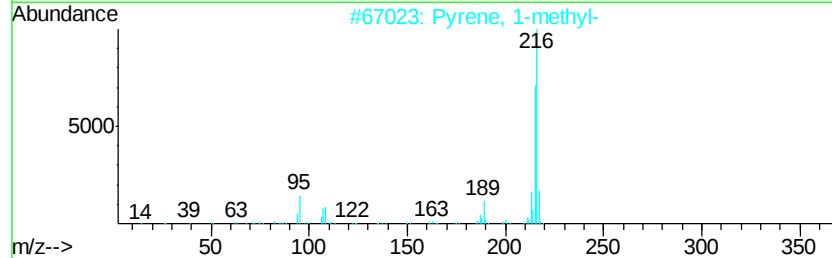
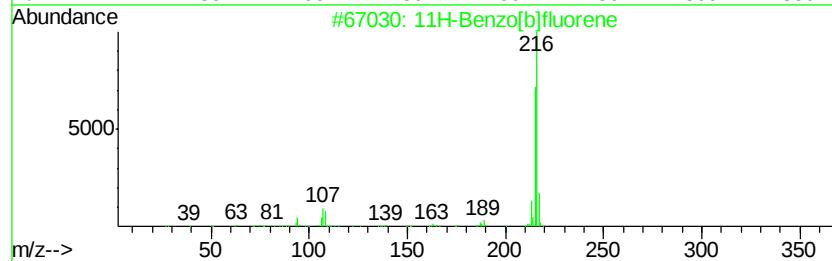
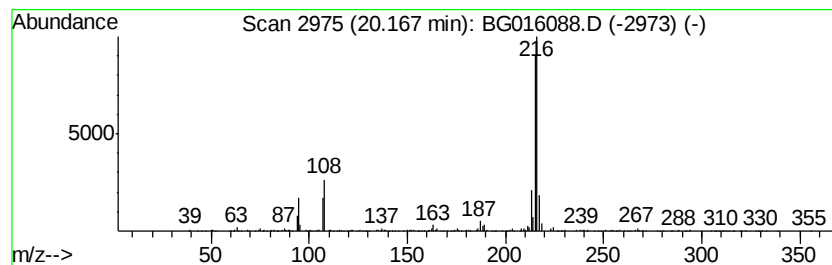
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG031715.M
Quant Title : SVOA CALIBRATION

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

Peak Number 20 11H-Benzo[b]fluorene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.17	3.42 ng/ul	810165	Chrysene-d12	21.34
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		11H-Benzo[b]fluorene	216 C17H12	000243-17-4 93
2		Pyrene, 1-methyl-	216 C17H12	002381-21-7 93
3		Fluoranthene, 2-methyl-	216 C17H12	033543-31-6 92
4		11H-Benzo[b]fluorene	216 C17H12	000243-17-4 91
5		11H-Benzo[a]fluorene	216 C17H12	000238-84-6 90



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032015\
Data File : BG016088.D
Acq On : 21 Mar 2015 5:11
Operator : TP/IZ
Sample : G1565-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AF1

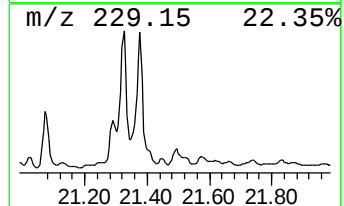
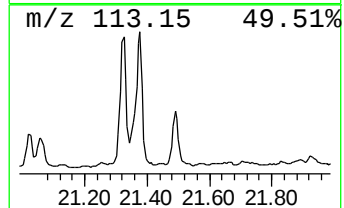
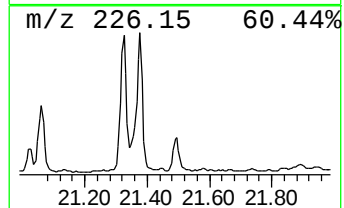
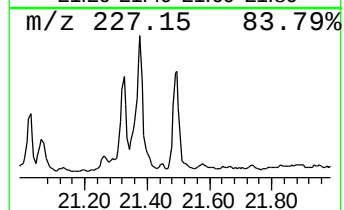
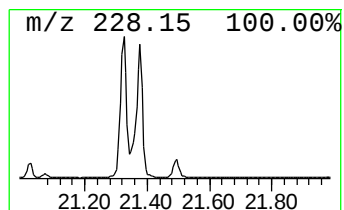
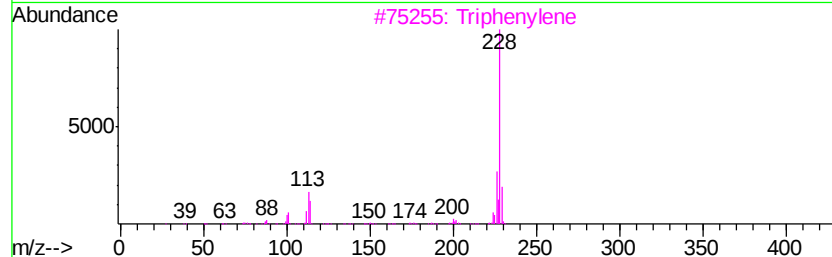
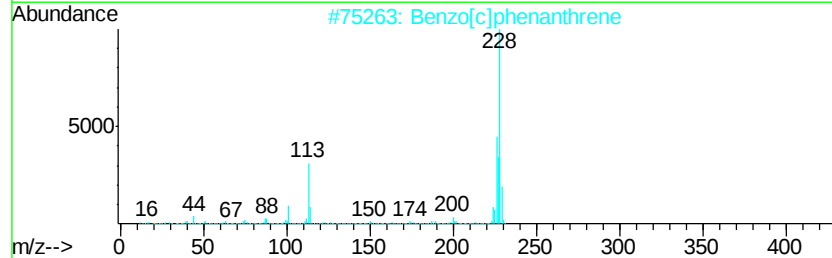
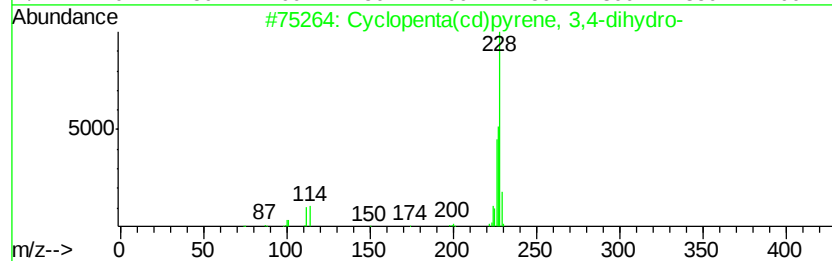
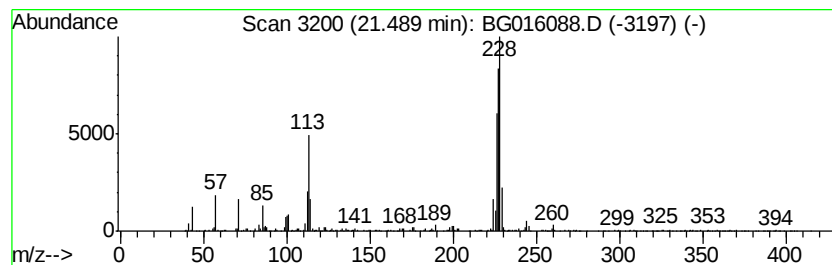
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 unknown21.49 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.49	2.32 ng/ul	550722	Chrysene-d12	21.34

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(cd)pyrene, 3,4-dihydro-	228	C18H12	025732-74-5	45
2			Benzo[c]phenanthrene	228	C18H12	000195-19-7	43
3			Triphenylene	228	C18H12	000217-59-4	38
4			Triphenylene	228	C18H12	000217-59-4	38
5			Chrysene	228	C18H12	000218-01-9	38



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032015\
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 Acq On : 21 Mar 2015 5:11
 Operator : TP/IZ
 Sample : G1565-11
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG031715.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...	3.02	31.7	ng/ul	896949	1	7.80	566110	20.0
Ethanol, 2-(2-but...	10.36	9.2	ng/ul	411197	2	10.59	894388	20.0
2,4,6-Cycloheptat...	15.76	3.5	ng/ul	208037	3	14.42	1196230	20.0
[1,1'-Biphenyl]-4...	15.91	4.1	ng/ul	246425	4	17.17	1191410	20.0
Heptadecane	16.13	2.9	ng/ul	173407	4	17.17	1191410	20.0
9H-Fluorene, 1-me...	16.47	2.2	ng/ul	131512	4	17.17	1191410	20.0
9H-Fluoren-9-one	16.82	2.3	ng/ul	137688	4	17.17	1191410	20.0
Dibenzothiophene	16.98	6.0	ng/ul	356605	4	17.17	1191410	20.0
Acridine	17.35	3.1	ng/ul	182230	4	17.17	1191410	20.0
Anthracene, 1-met...	18.04	7.2	ng/ul	429261	4	17.17	1191410	20.0
Phenanthrene, 2-m...	18.08	11.5	ng/ul	683008	4	17.17	1191410	20.0
Anthracene, 2-met...	18.16	3.6	ng/ul	216102	4	17.17	1191410	20.0
4H-Cyclopenta[def...	18.23	21.2	ng/ul	1261820	4	17.17	1191410	20.0
2-Phenylnaphthalene	18.53	5.7	ng/ul	337983	4	17.17	1191410	20.0
9,10-Anthracenedione	18.58	3.8	ng/ul	225045	4	17.17	1191410	20.0
Phenanthrene, 2,5...	18.95	4.2	ng/ul	248407	4	17.17	1191410	20.0
Cyclopenta(def)ph...	19.10	3.7	ng/ul	217749	4	17.17	1191410	20.0
11H-Benzo[b]fluorene	20.07	5.0	ng/ul	1190960	5	21.34	4741680	20.0
11H-Benzo[b]fluorene	20.17	3.4	ng/ul	810165	5	21.34	4741680	20.0
unknown21.49	21.49	2.3	ng/ul	550722	5	21.34	4741680	20.0