

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
 Data File : BG023673.D
 Acq On : 13 Aug 2016 2:00
 Operator : UM/SJ
 Sample : H4385-18
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 P003-SS001-0206-01

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-BG080416.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	4.460	272	276	284	rBV	124899	187136	2.89%	0.158%
2	7.268	747	754	764	rBV2	372016	691520	10.68%	0.583%
3	7.427	775	781	791	rVB	287237	490427	7.57%	0.413%
4	7.644	811	818	824	rBV	367925	634771	9.80%	0.535%
5	8.114	890	898	907	rVB	402264	697935	10.78%	0.588%
6	8.831	1012	1020	1028	rBV	446831	819127	12.65%	0.690%
7	9.289	1091	1098	1110	rBV	376110	702202	10.84%	0.592%
8	10.023	1214	1223	1232	rVV	340252	631639	9.75%	0.532%
9	10.570	1307	1316	1336	rVB	481073	959283	14.81%	0.808%
10	10.951	1374	1381	1387	rBV	614461	1105513	17.07%	0.932%
11	11.092	1396	1405	1421	rBV	395373	806866	12.46%	0.680%
12	12.608	1656	1663	1671	rBV	189537	316693	4.89%	0.267%
13	12.831	1691	1701	1707	rBV	226195	403581	6.23%	0.340%
14	13.959	1881	1893	1900	rBV3	207121	567567	8.76%	0.478%
15	14.106	1911	1918	1923	rVV	408030	772914	11.93%	0.651%
16	14.182	1923	1931	1935	rVV	1060281	1906473	29.44%	1.607%
17	14.229	1935	1939	1948	rVB	709118	1028886	15.89%	0.867%
18	14.341	1952	1958	1962	rVV	199652	338444	5.23%	0.285%
19	14.482	1975	1982	1995	rBV2	1063394	2224503	34.35%	1.875%
20	14.787	2029	2034	2040	rVV2	1036850	1686827	26.05%	1.421%
21	14.852	2040	2045	2052	rVV4	195623	344786	5.32%	0.291%
22	15.011	2063	2072	2079	rBV3	378143	862429	13.32%	0.727%
23	15.181	2089	2101	2108	rVV3	307338	622648	9.61%	0.525%
24	15.257	2108	2114	2120	rVB	284201	428929	6.62%	0.361%
25	15.404	2131	2139	2143	rBV3	220029	437976	6.76%	0.369%
26	15.457	2143	2148	2152	rVB	201222	284200	4.39%	0.239%
27	15.575	2160	2168	2171	rBV5	157793	415974	6.42%	0.351%
28	15.604	2171	2173	2179	rVB	194519	256710	3.96%	0.216%
29	15.786	2199	2204	2209	rVV	1463207	2340860	36.14%	1.973%
30	15.839	2209	2213	2222	rVV	417925	833340	12.87%	0.702%
31	15.921	2222	2227	2232	rVV	387252	689262	10.64%	0.581%
32	16.133	2255	2263	2271	rVB3	135925	353587	5.46%	0.298%
33	16.279	2284	2288	2296	rVB	152626	297750	4.60%	0.251%
34	16.509	2324	2327	2337	rVB3	187371	355450	5.49%	0.300%

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35	16.849	2376	2385	2389	rVV3	287514	695729	10.74%	0.586%
36	16.896	2389	2393	2399	rVB2	229219	345323	5.33%	0.291%
37	17.002	2406	2411	2416	rVB2	155875	256778	3.96%	0.216%
38	17.078	2416	2424	2428	rBV5	120640	311561	4.81%	0.263%
39	17.208	2441	2446	2452	rBV2	176484	314621	4.86%	0.265%
40	17.272	2453	2457	2464	rVV6	172080	269262	4.16%	0.227%
41	17.372	2469	2474	2479	rVV	399478	597773	9.23%	0.504%
42	17.554	2499	2505	2508	rBV2	1284249	2020391	31.20%	1.703%
43	17.601	2508	2513	2518	rVV	3656218	5937481	91.68%	5.003%
44	17.654	2518	2522	2526	rVV	1535758	2415195	37.29%	2.035%
45	17.689	2526	2528	2532	rVV	554951	615802	9.51%	0.519%
46	17.748	2532	2538	2542	rVB3	183776	335977	5.19%	0.283%
47	17.965	2570	2575	2581	rVV3	176402	388119	5.99%	0.327%
48	18.071	2589	2593	2596	rVV	139431	187679	2.90%	0.158%
49	18.136	2600	2604	2612	rVB	482876	796850	12.30%	0.671%
50	18.283	2624	2629	2637	rVB	325947	647087	9.99%	0.545%
51	18.359	2637	2642	2647	rBV4	151051	290115	4.48%	0.244%
52	18.435	2649	2655	2659	rVV	1558476	2590389	40.00%	2.183%
53	18.482	2659	2663	2669	rVB	1841983	2674945	41.30%	2.254%
54	18.565	2672	2677	2682	rBV2	293881	484074	7.47%	0.408%
55	18.623	2682	2687	2691	rVV2	1365713	2664884	41.15%	2.246%
56	18.664	2691	2694	2699	rVB	1048224	1505705	23.25%	1.269%
57	18.829	2718	2722	2726	rVB2	226221	318821	4.92%	0.269%
58	18.929	2734	2739	2743	rVV2	656090	981386	15.15%	0.827%
59	18.982	2743	2748	2756	rVB2	615090	1115582	17.23%	0.940%
60	19.170	2772	2780	2784	rBV2	520133	1027776	15.87%	0.866%
61	19.223	2785	2789	2793	rVV	714097	912583	14.09%	0.769%
62	19.264	2793	2796	2801	rVB2	477294	638284	9.86%	0.538%
63	19.346	2804	2810	2814	rBV	1452994	2428075	37.49%	2.046%
64	19.399	2814	2819	2823	rVV3	471318	954195	14.73%	0.804%
65	19.440	2823	2826	2829	rVB	433129	528039	8.15%	0.445%
66	19.487	2829	2834	2839	rBV3	512981	1131470	17.47%	0.953%
67	19.616	2850	2856	2861	rVB	3962213	6073800	93.78%	5.118%
68	19.669	2861	2865	2868	rBV2	309177	440613	6.80%	0.371%
69	19.704	2868	2871	2876	rVB2	372479	520644	8.04%	0.439%
70	19.763	2877	2881	2885	rBV	227963	334565	5.17%	0.282%
71	19.851	2892	2896	2900	rBV2	268360	448088	6.92%	0.378%

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72	19.951	2907	2913	2915	rBV2	2051566	3499744	54.04%	2.949%
73	19.980	2915	2918	2924	rVB	4644061	6476505	100.00%	5.458%
74	20.045	2924	2929	2934	rVB	736046	1116432	17.24%	0.941%
75	20.121	2934	2942	2943	rBV5	239205	507018	7.83%	0.427%
76	20.209	2953	2957	2963	rVB4	306561	550610	8.50%	0.464%
77	20.303	2965	2973	2979	rBV6	641327	1723157	26.61%	1.452%
78	20.468	2991	3001	3008	rVB	841814	2305790	35.60%	1.943%
79	20.562	3014	3017	3022	rVB	459686	604001	9.33%	0.509%
80	20.627	3024	3028	3032	rVB	908300	1134156	17.51%	0.956%
81	20.767	3048	3052	3056	rBV	842187	1187742	18.34%	1.001%
82	20.814	3056	3060	3072	rVB2	794226	1504582	23.23%	1.268%
83	21.249	3130	3134	3141	rVB2	529321	949940	14.67%	0.801%
84	21.349	3148	3151	3156	rVB	210817	304092	4.70%	0.256%
85	21.443	3158	3167	3170	rVV3	624396	1434420	22.15%	1.209%
86	21.484	3171	3174	3179	rVV	450314	639987	9.88%	0.539%
87	21.537	3179	3183	3187	rVV3	364449	639889	9.88%	0.539%
88	21.602	3191	3194	3201	rVB	370166	641242	9.90%	0.540%
89	21.866	3232	3239	3246	rBV2	2165199	5907588	91.22%	4.978%
90	21.931	3246	3250	3262	rVB	1795944	3339521	51.56%	2.814%
91	22.089	3274	3277	3282	rVB2	201581	283645	4.38%	0.239%
92	22.166	3286	3290	3294	rBV3	223925	390268	6.03%	0.329%
93	22.647	3364	3372	3378	rBV	582375	1249686	19.30%	1.053%
94	22.724	3381	3385	3392	rVB	322641	571610	8.83%	0.482%
95	22.935	3417	3421	3425	rBV3	246311	412715	6.37%	0.348%
96	24.204	3630	3637	3645	rBV2	889803	3085169	47.64%	2.600%
97	24.974	3761	3768	3775	rBV	551095	1604757	24.78%	1.352%
98	25.062	3776	3783	3789	rVV2	761413	2193832	33.87%	1.849%
99	25.138	3789	3796	3808	rVB	962510	2710436	41.85%	2.284%
100	25.291	3813	3822	3830	rBV2	684193	2003728	30.94%	1.689%

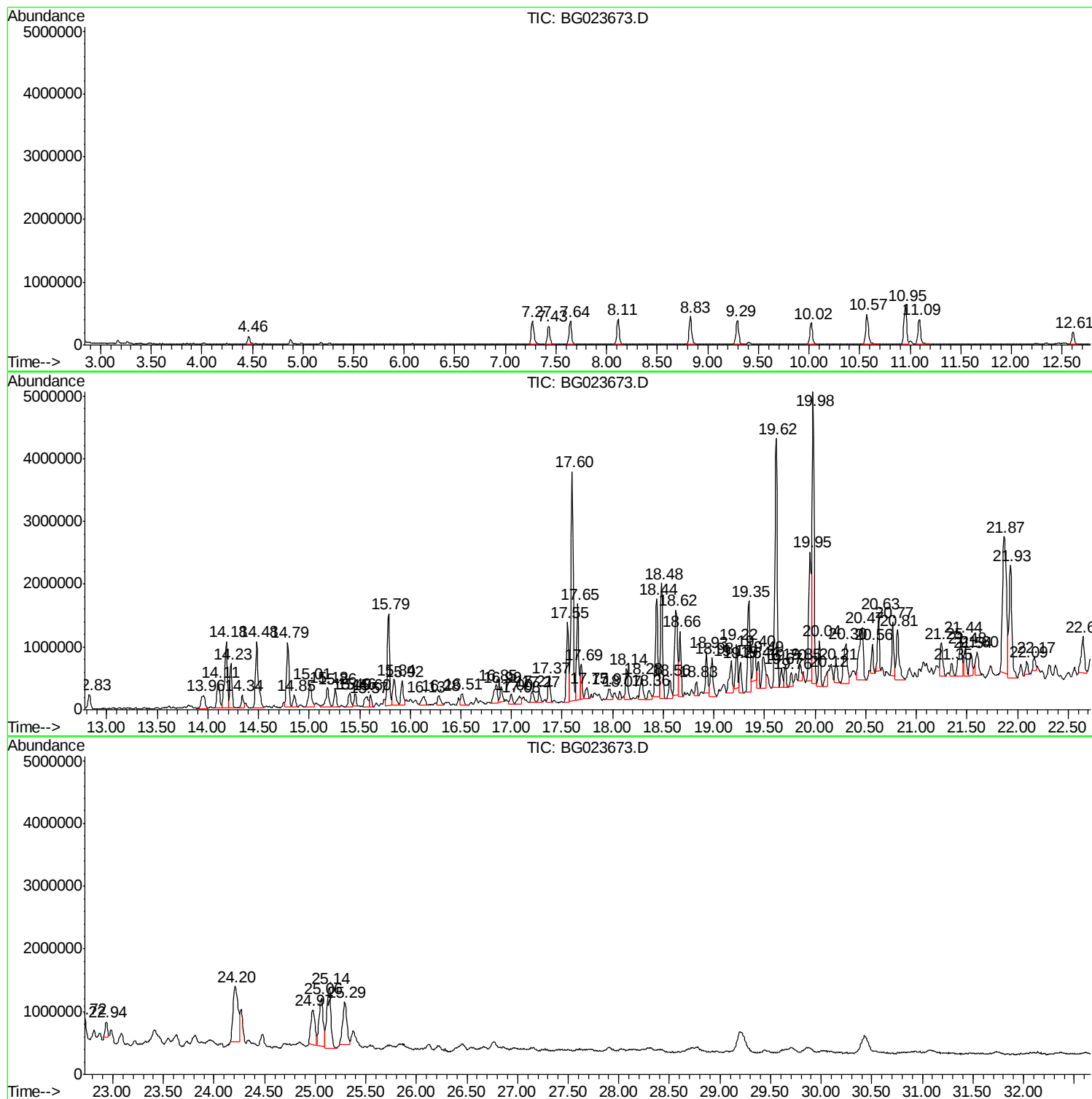
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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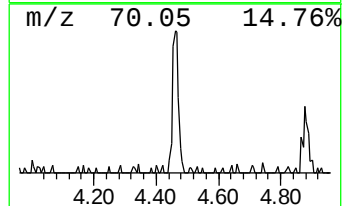
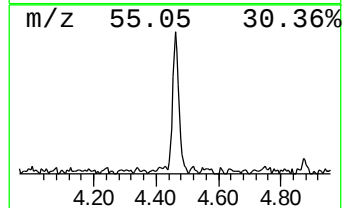
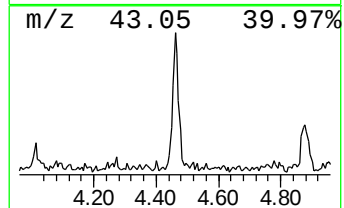
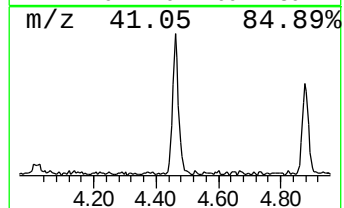
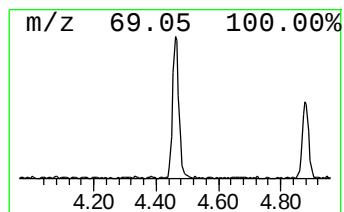
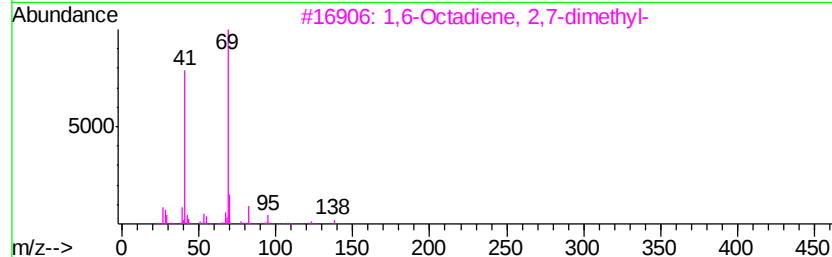
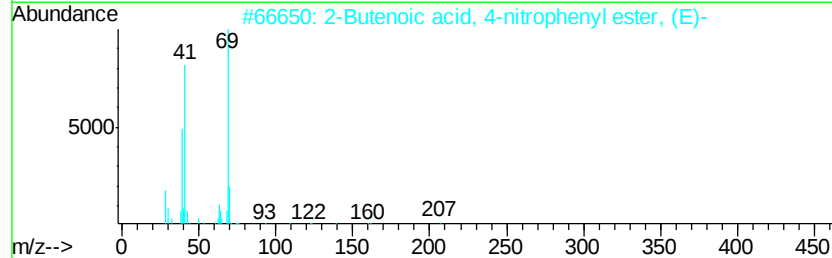
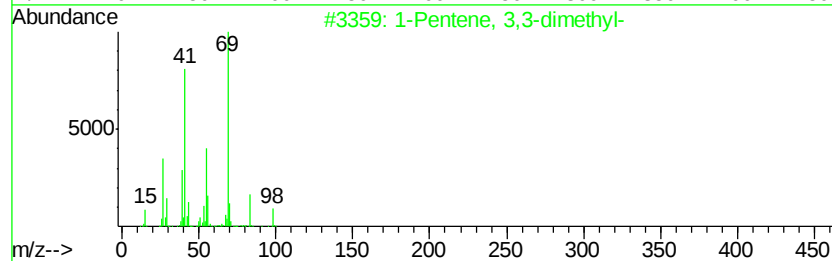
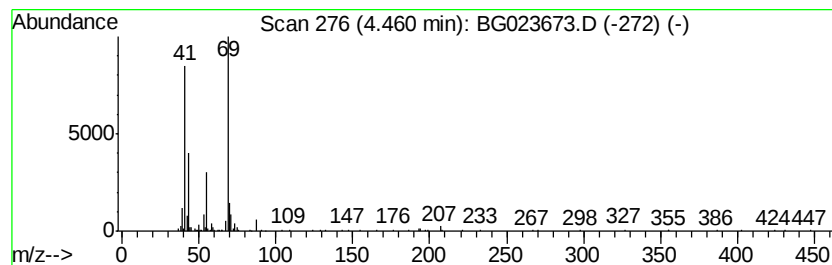
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Cyclic4.46 Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.46	5.36 ng/ul	187136	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Pentene, 3,3-dimethyl-	98	C7H14	003404-73-7	59
2		2-Butenoic acid, 4-nitrophenyl e...	207	C10H9NO4	014617-88-0	53
3		1,6-Octadiene, 2,7-dimethyl-	138	C10H18	040195-09-3	38
4		1,5-Heptadiene, 2,6-dimethyl-	124	C9H16	006709-39-3	9
5		6,11-Dimethyl-2,6,10-dodecatrien...	208	C14H24O	1000196-53-3	9



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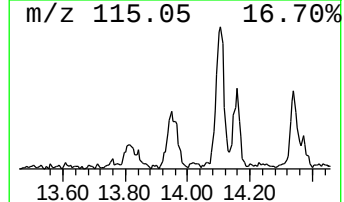
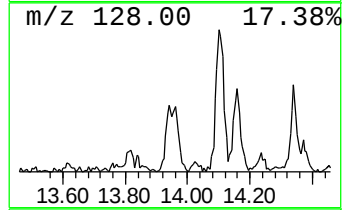
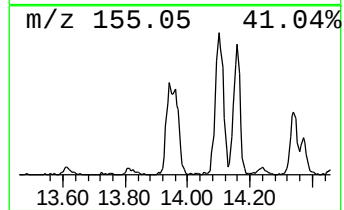
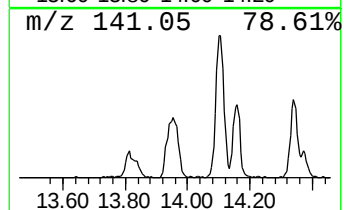
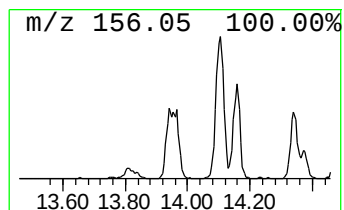
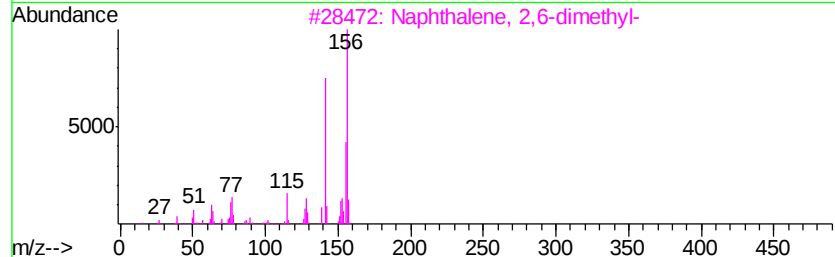
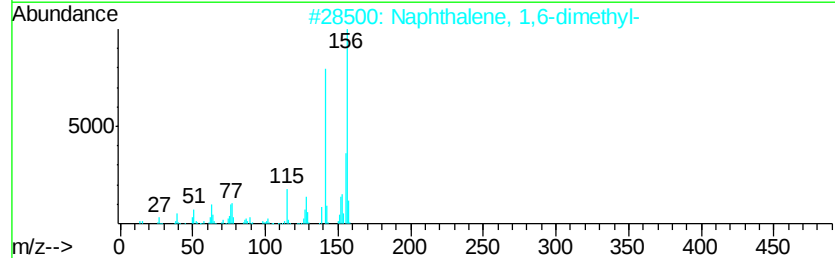
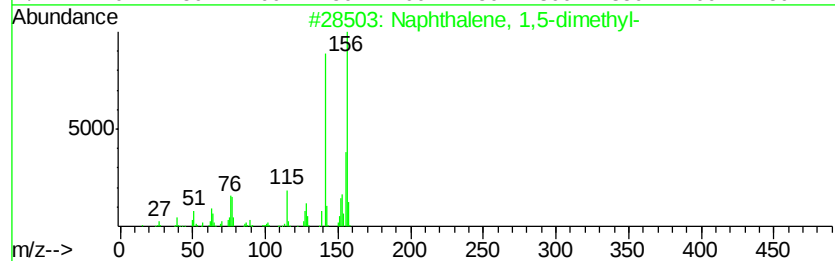
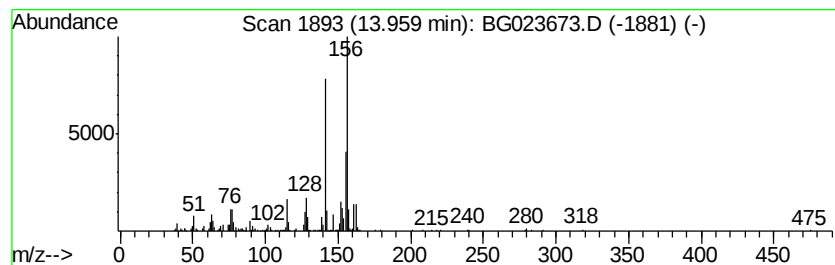
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Peak Number 3 Naphthalene, 1,5-dimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.96	6.73 ng/ul	567567	Acenaphthene-d10	14.79
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,5-dimethyl-	156 C12H12	000571-61-9 98
2		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 98
3		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 97
4		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 96
5		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 96



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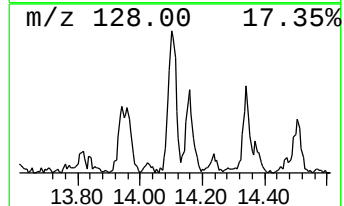
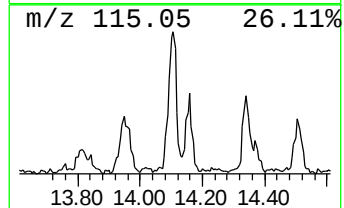
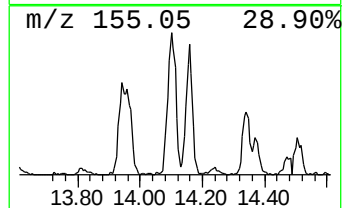
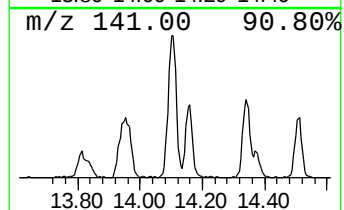
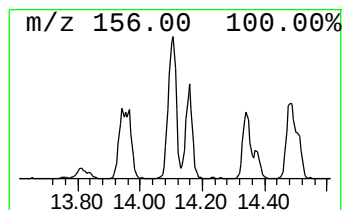
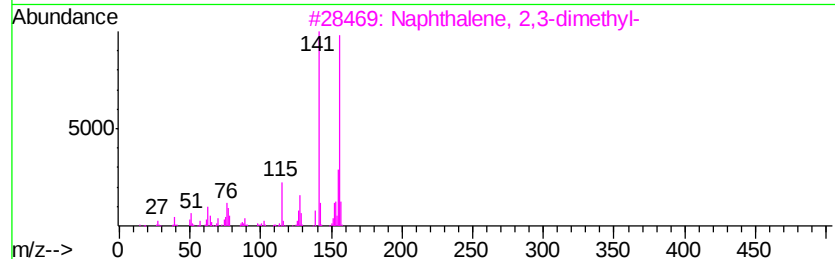
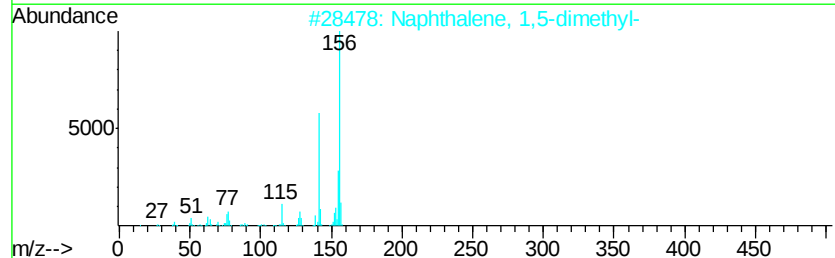
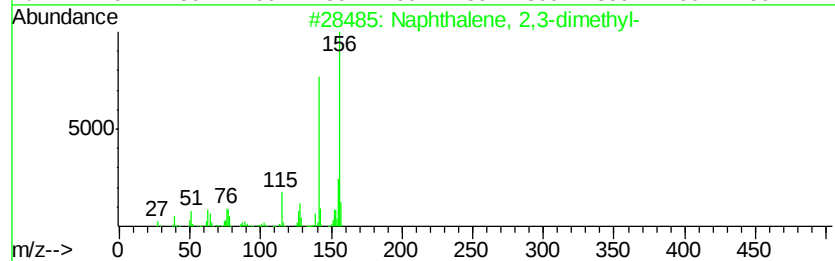
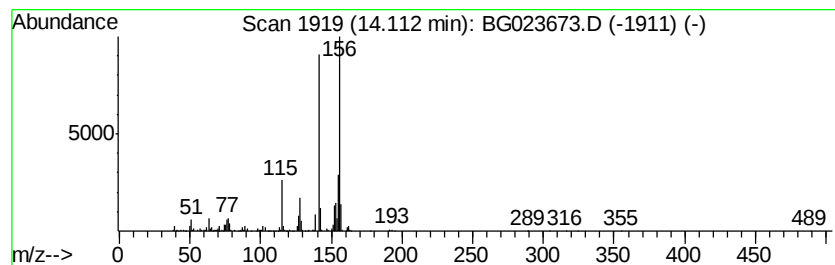
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Naphthalene, 2,3-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.11	9.16 ng/ul	772914	Acenaphthene-d10	14.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
3		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
Data File : BG023673.D
Acq On : 13 Aug 2016 2:00
Operator : UM/SJ
Sample : H4385-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
P003-SS001-0206-01

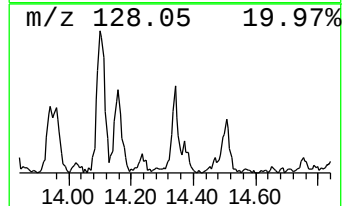
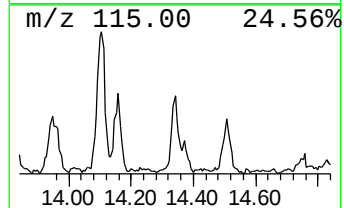
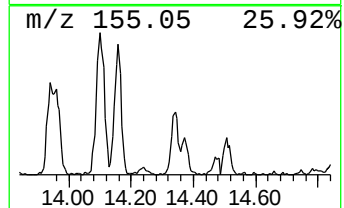
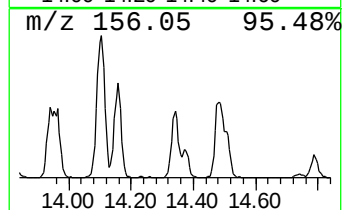
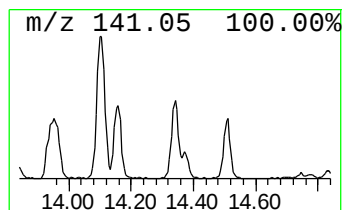
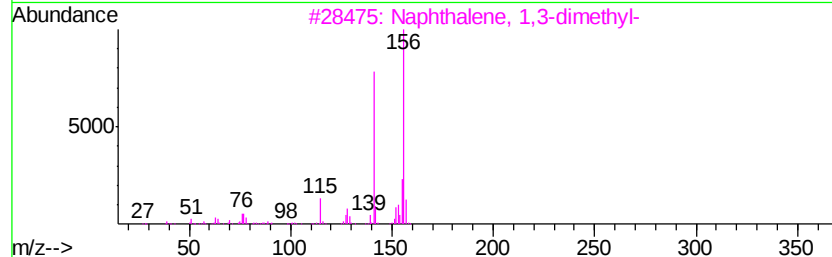
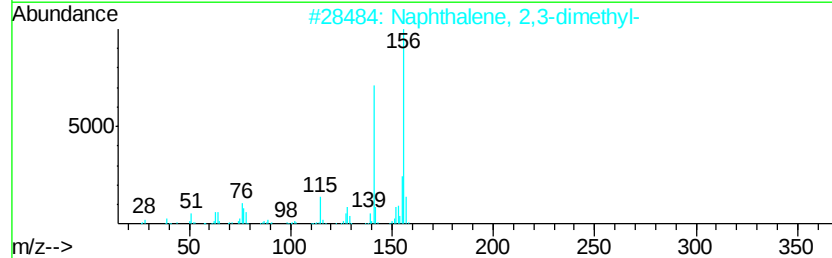
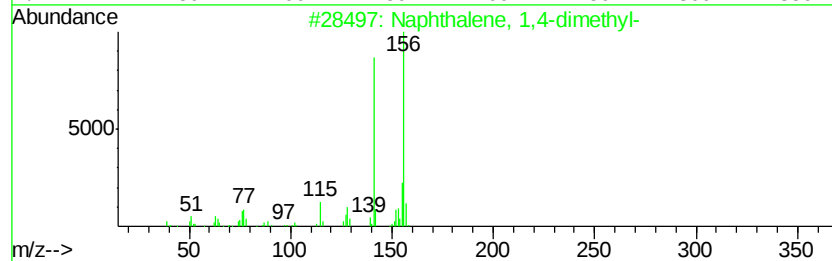
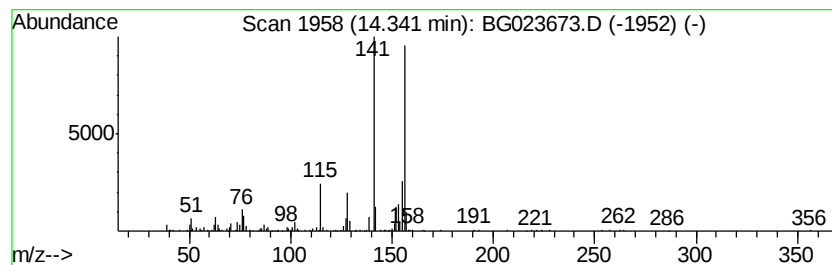
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Naphthalene, 1,4-dimethyl- Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.34	4.01 ng/ul	338444	Acenaphthene-d10	14.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
2		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	94
3		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	94
4		Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	93
5		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	93



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
Data File : BG023673.D
Acq On : 13 Aug 2016 2:00
Operator : UM/SJ
Sample : H4385-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
P003-SS001-0206-01

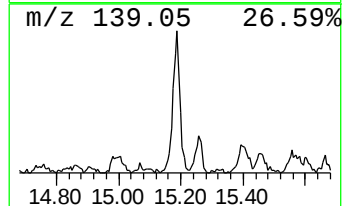
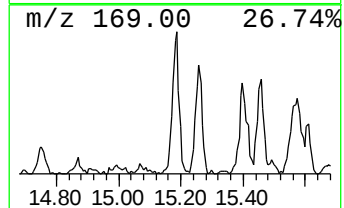
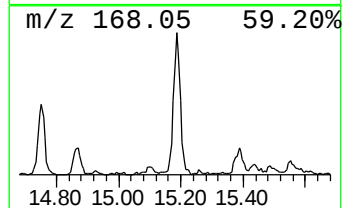
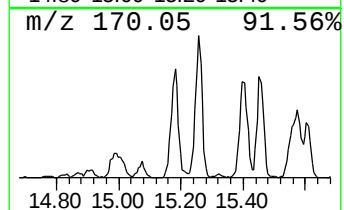
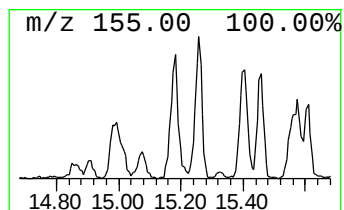
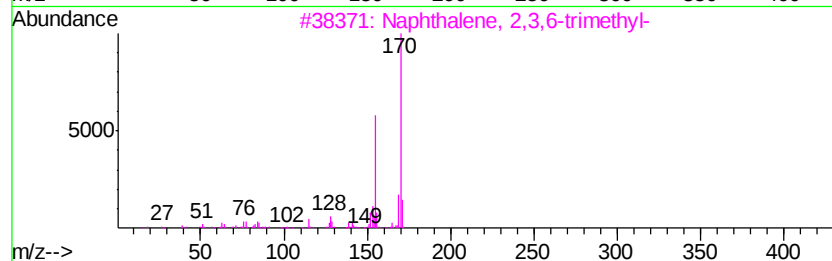
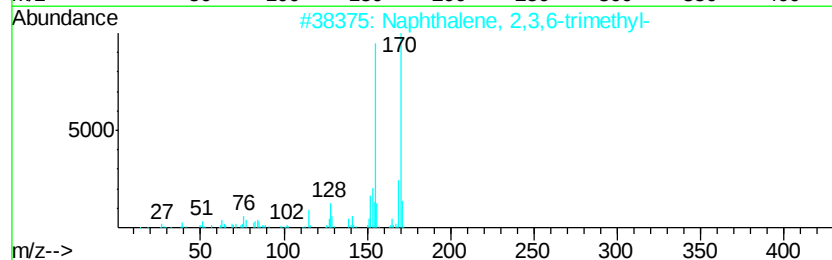
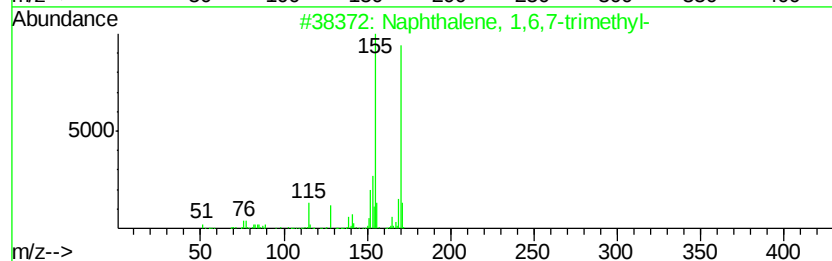
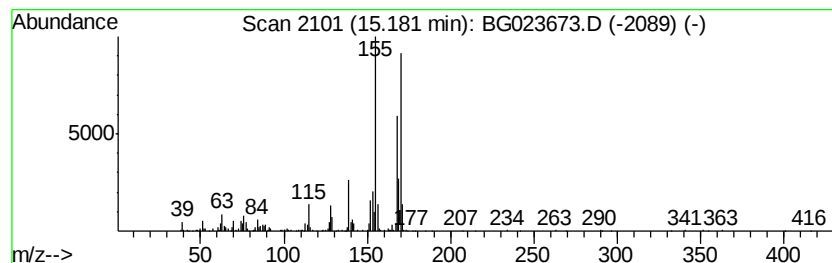
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Naphthalene, 1,6,7-trimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.18	7.38 ng/ul	622648	Acenaphthene-d10	14.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
2		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	95
3		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	95
4		Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	95
5		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	95



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
Data File : BG023673.D
Acq On : 13 Aug 2016 2:00
Operator : UM/SJ
Sample : H4385-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
P003-SS001-0206-01

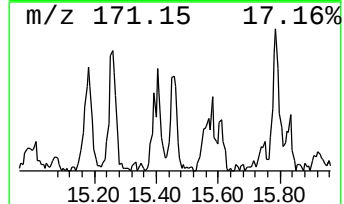
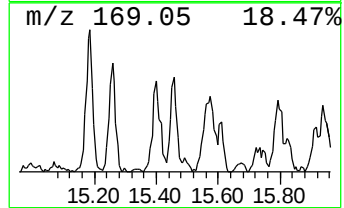
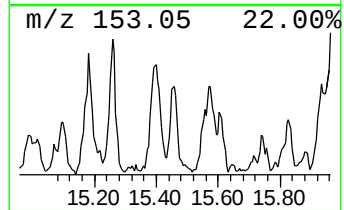
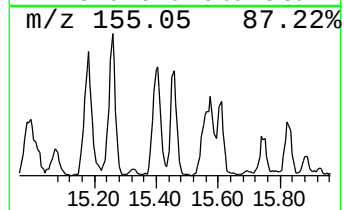
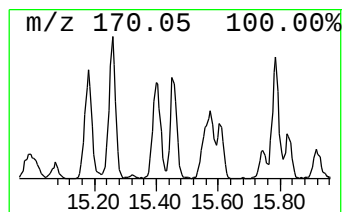
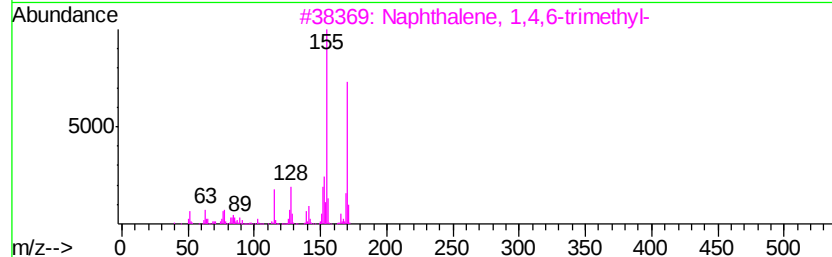
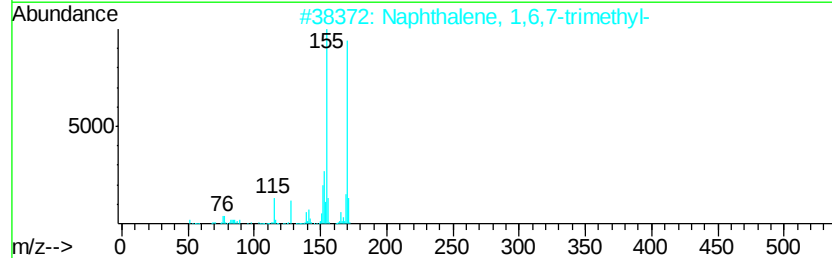
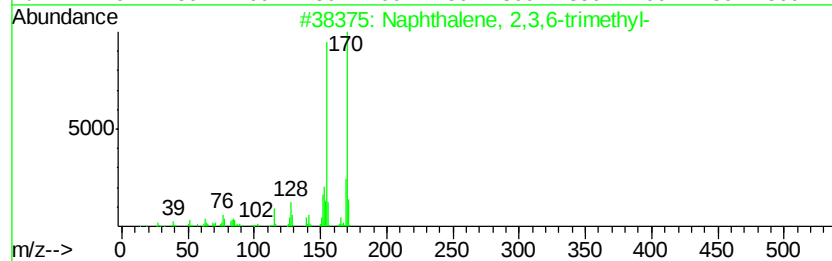
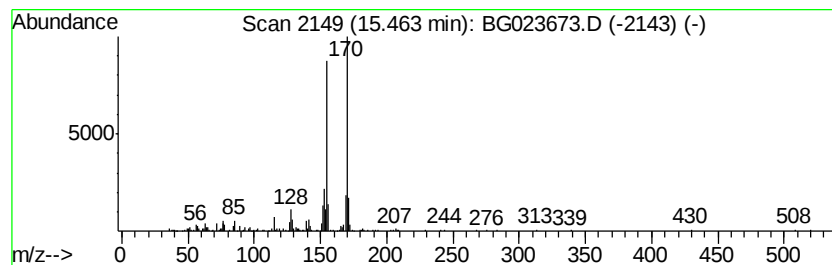
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Naphthalene, 2,3,6-trimethyl- Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.46	3.37 ng/ul	284200	Acenaphthene-d10	14.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
2		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	97
3		Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	96
4		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
5		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	95



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
Data File : BG023673.D
Acq On : 13 Aug 2016 2:00
Operator : UM/SJ
Sample : H4385-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
P003-SS001-0206-01

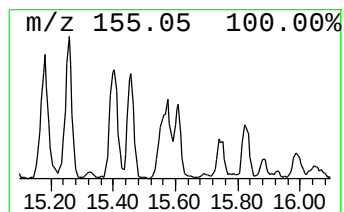
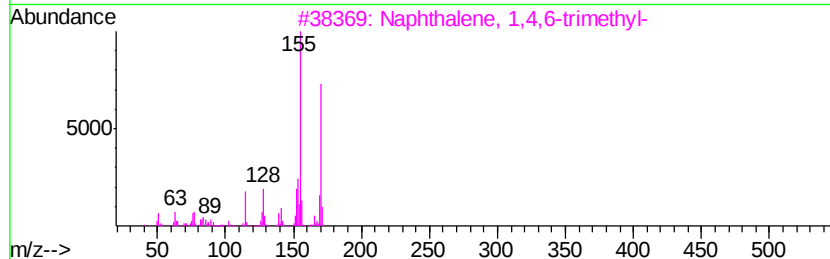
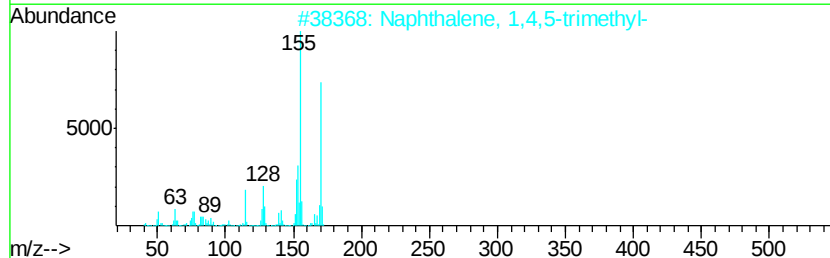
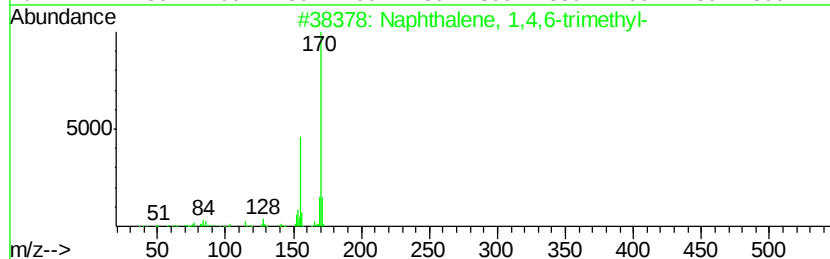
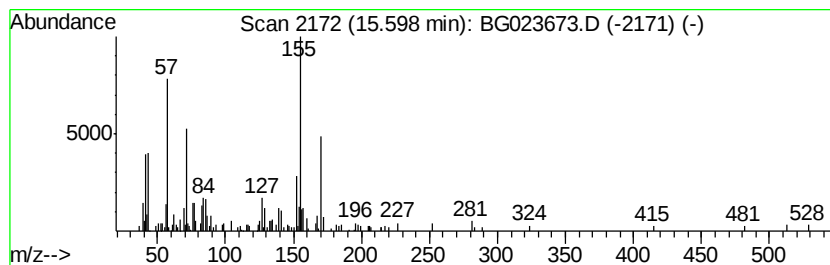
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Naphthalene, 1,4,6-trimethyl- Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.60	3.04 ng/ul	256710	Acenaphthene-d10	14.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	64
2			Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	60
3			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	50
4			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	47
5			trans, cis-2-Methyl-1-thiadecalin	170	C10H18S	042900-29-8	43



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
Data File : BG023673.D
Acq On : 13 Aug 2016 2:00
Operator : UM/SJ
Sample : H4385-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
P003-SS001-0206-01

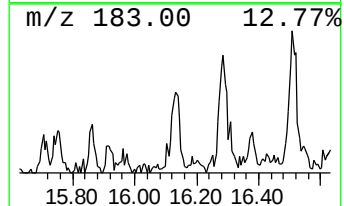
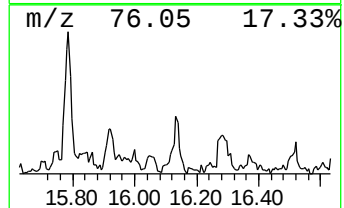
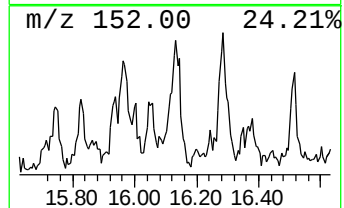
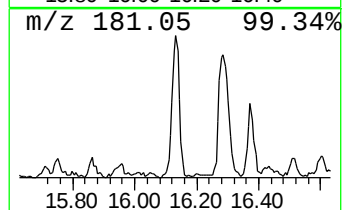
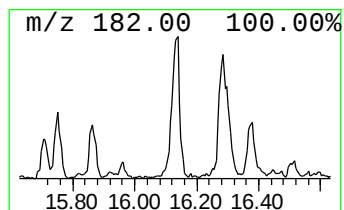
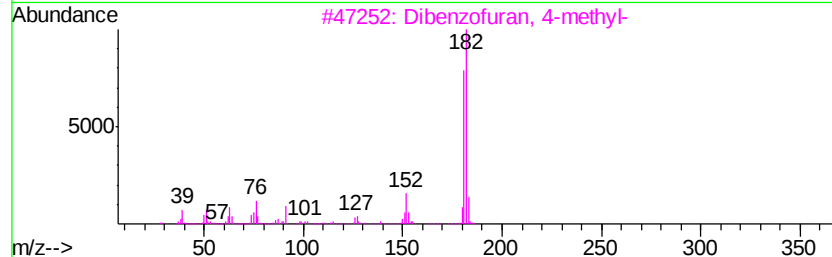
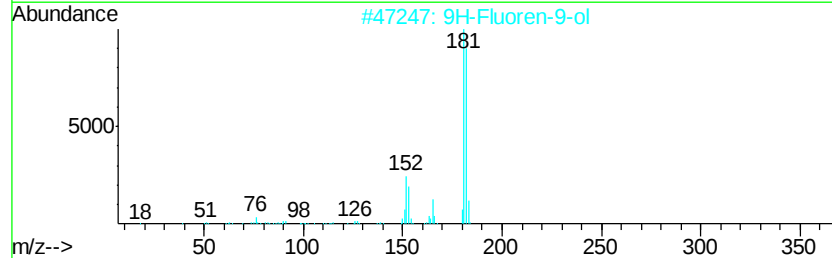
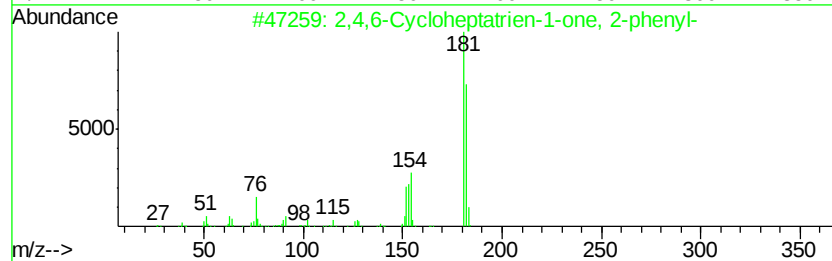
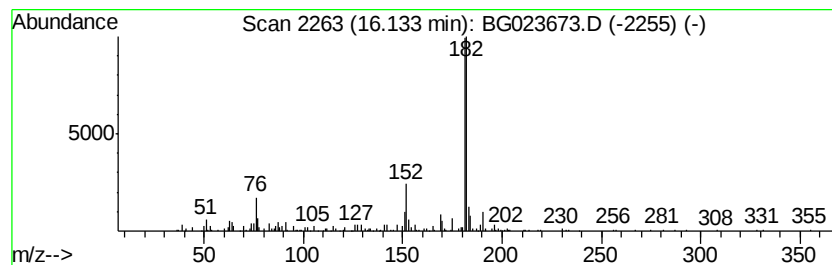
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 2,4,6-Cycloheptatrien-1-one... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.13	4.19 ng/ul	353587	Acenaphthene-d10	14.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	90
2		9H-Fluoren-9-ol	182	C13H10O	001689-64-1	87
3		Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	80
4		9H-Fluoren-9-ol	182	C13H10O	001689-64-1	64
5		Pyridine, 4,4'-(1,2-ethenediyl)bis-	182	C12H10N2	001135-32-6	59



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
Data File : BG023673.D
Acq On : 13 Aug 2016 2:00
Operator : UM/SJ
Sample : H4385-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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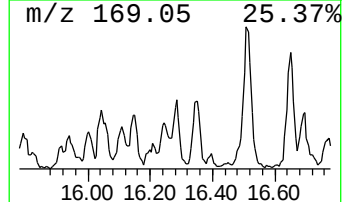
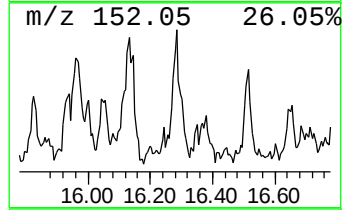
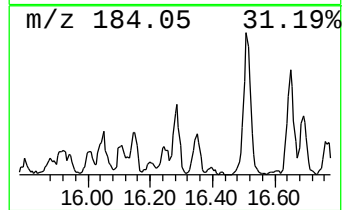
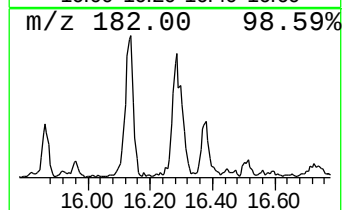
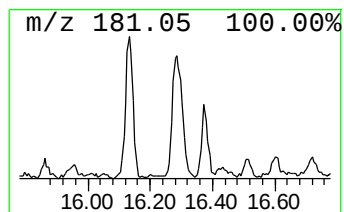
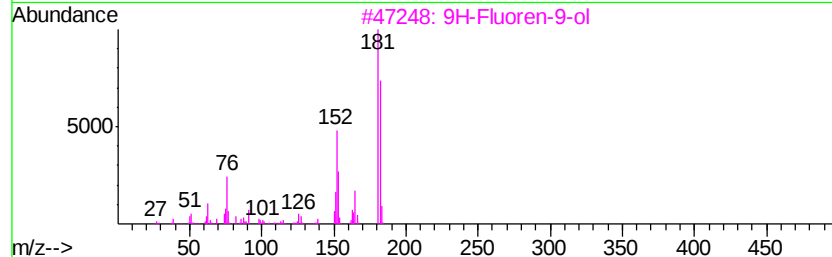
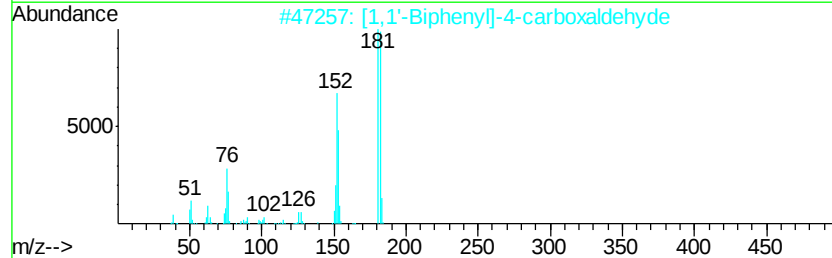
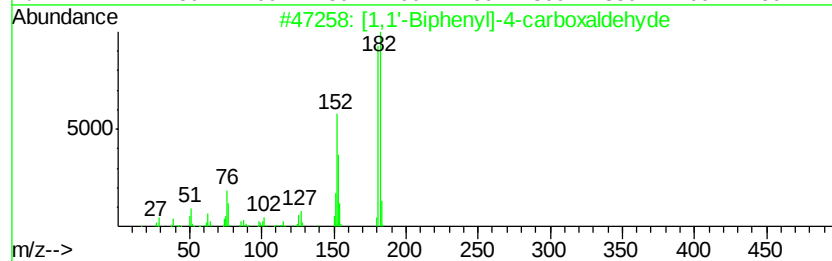
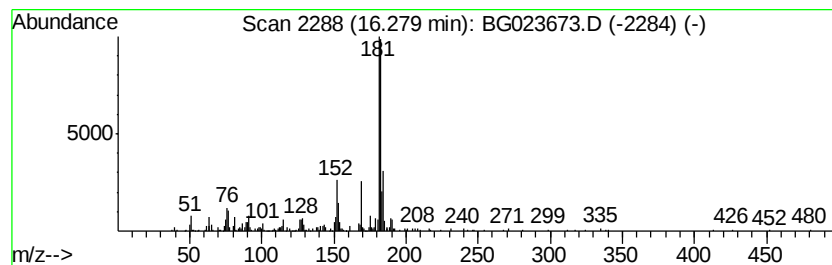
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 [1,1'-Biphenyl]-4-carboxald... Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.28	2.95 ng/ul	297750	Phenanthrene-d10	17.55

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	72
2			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	64
3			9H-Fluoren-9-ol	182	C13H10O	001689-64-1	64
4			9H-Fluoren-9-ol	182	C13H10O	001689-64-1	64
5			2-(4-Chlorophenyl)-1,3,2-dioxabo...	182	C8H8BClO2	069519-09-1	58



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
Data File : BG023673.D
Acq On : 13 Aug 2016 2:00
Operator : UM/SJ
Sample : H4385-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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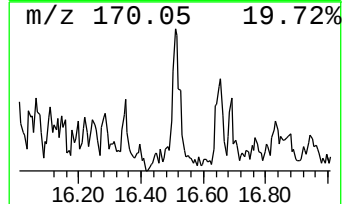
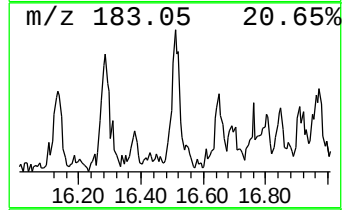
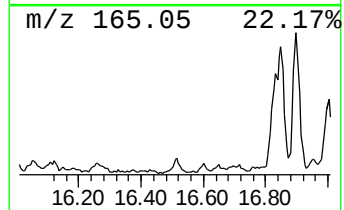
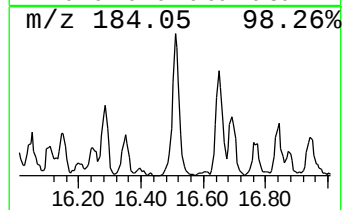
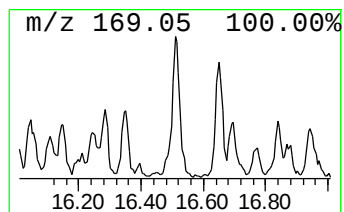
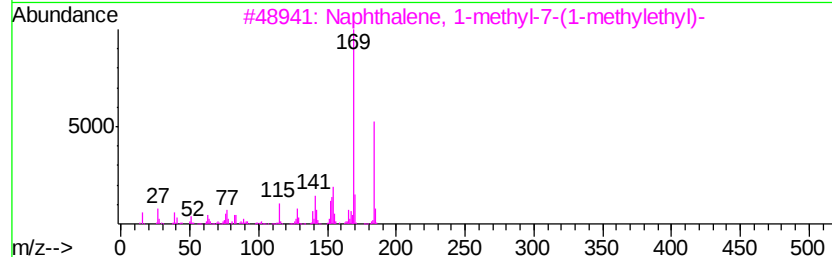
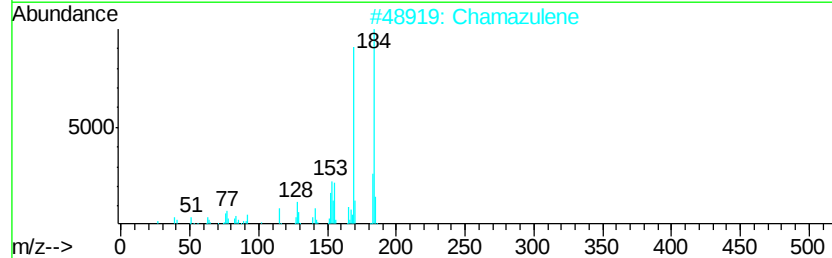
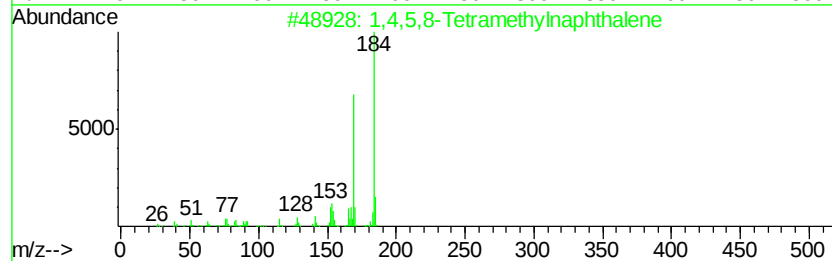
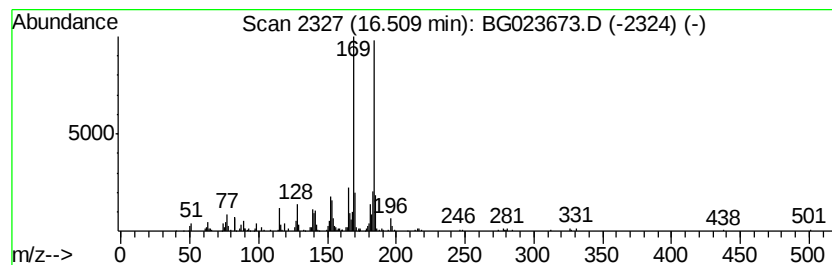
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 1,4,5,8-Tetramethylnaphthalene Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.51	3.52 ng/ul	355450	Phenanthrene-d10	17.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,4,5,8-Tetramethylnaphthalene	184	C14H16	002717-39-7	87
2		Chamazulene	184	C14H16	000529-05-5	76
3		Naphthalene, 1-methyl-7-(1-methy...	184	C14H16	000490-65-3	76
4		Naphthalene, 1,2,3,4-tetramethyl-	184	C14H16	003031-15-0	76
5		Benzene, (4,5,5-trimethyl-1,3-cy...	184	C14H16	033930-85-7	64



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
Data File : BG023673.D
Acq On : 13 Aug 2016 2:00
Operator : UM/SJ
Sample : H4385-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
P003-SS001-0206-01

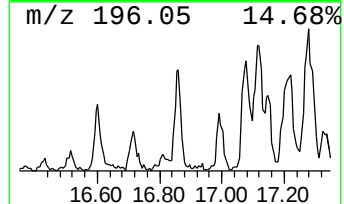
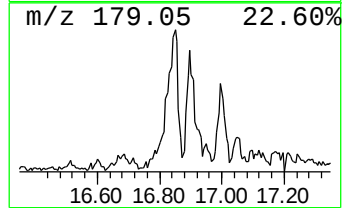
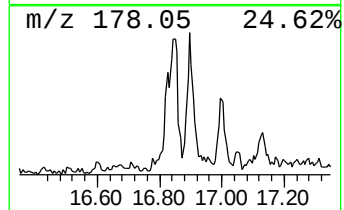
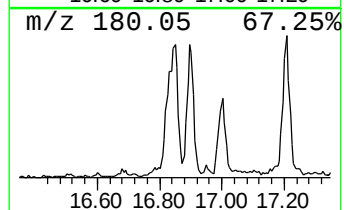
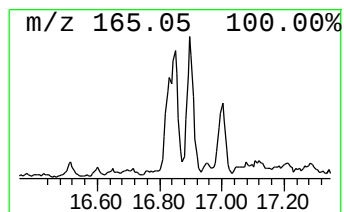
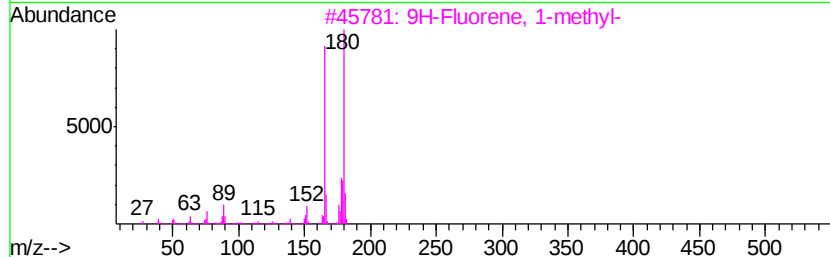
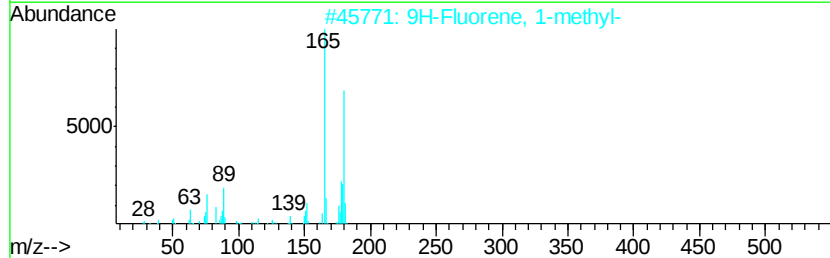
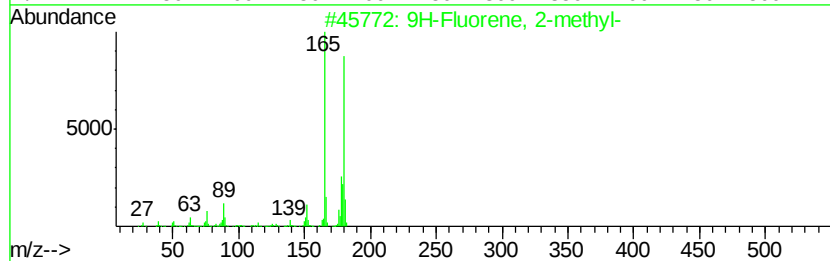
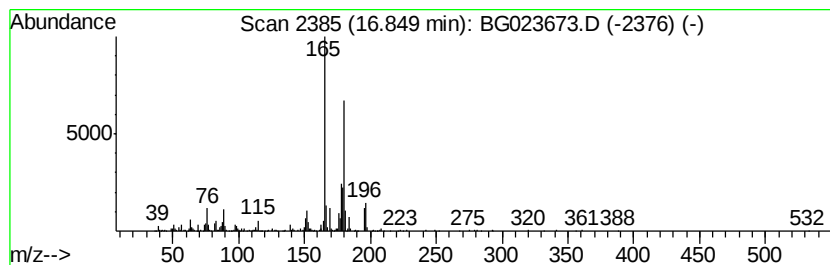
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 9H-Fluorene, 2-methyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.85	6.89 ng/ul	695729	Phenanthrene-d10	17.55

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	86
2			9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	86
3			9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	83
4			(3H)Benzo[c]pyrrole, 3-methyl-3-...	208	C14H12N2	059341-20-7	76
5			9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	70



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
Data File : BG023673.D
Acq On : 13 Aug 2016 2:00
Operator : UM/SJ
Sample : H4385-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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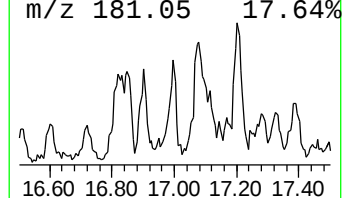
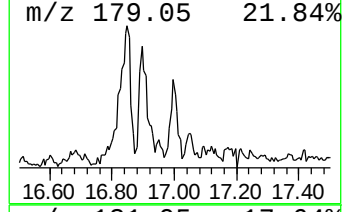
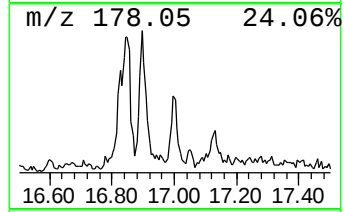
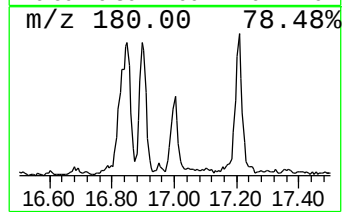
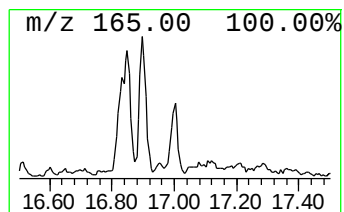
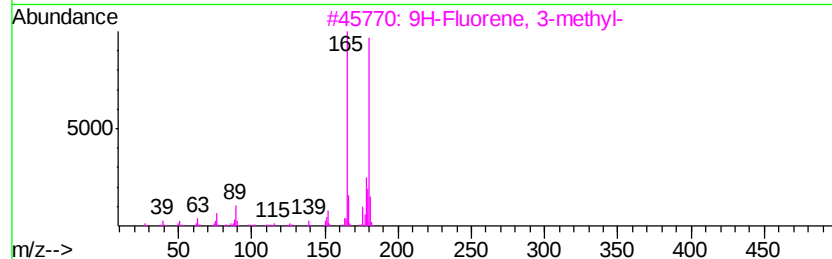
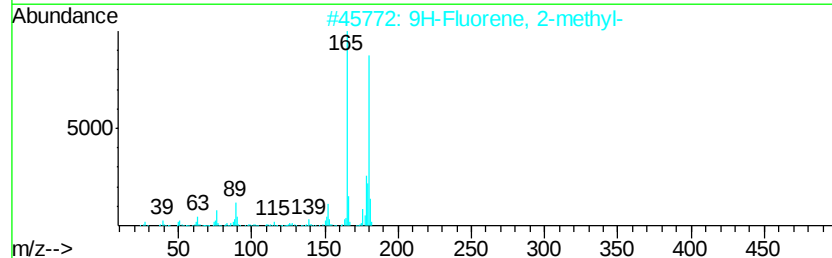
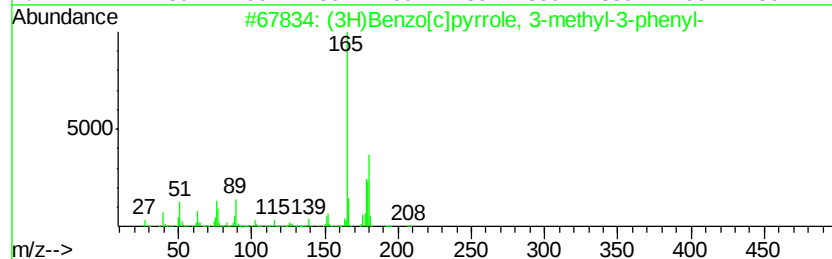
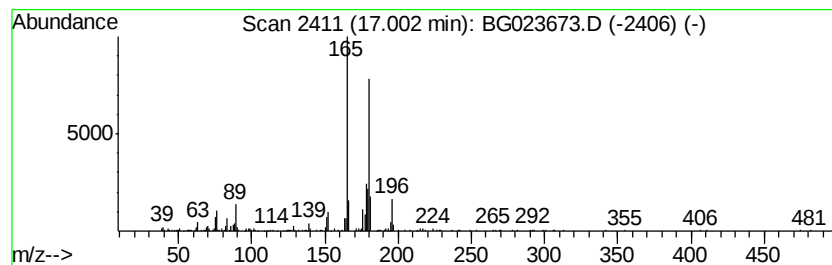
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 (3H)Benzo[c]pyrrole, 3-meth... Concentration Rank 53

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.00	2.54 ng/ul	256778	Phenanthrene-d10	17.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(3H)Benzo[c]pyrrole, 3-methyl-3-...	208	C14H12N2	059341-20-7	95
2		9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	93
3		9H-Fluorene, 3-methyl-	180	C14H12	002523-39-9	93
4		9H-Fluorene, 9-methyl-	180	C14H12	002523-37-7	91
5		9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	91



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
Data File : BG023673.D
Acq On : 13 Aug 2016 2:00
Operator : UM/SJ
Sample : H4385-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
P003-SS001-0206-01

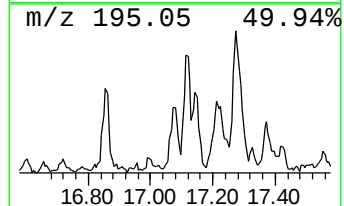
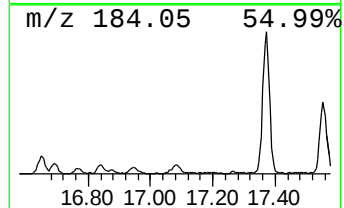
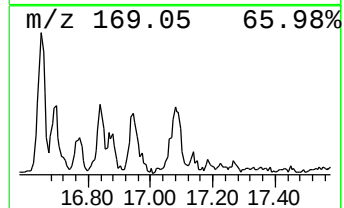
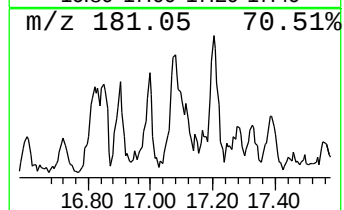
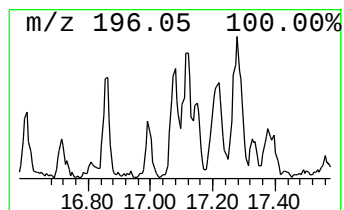
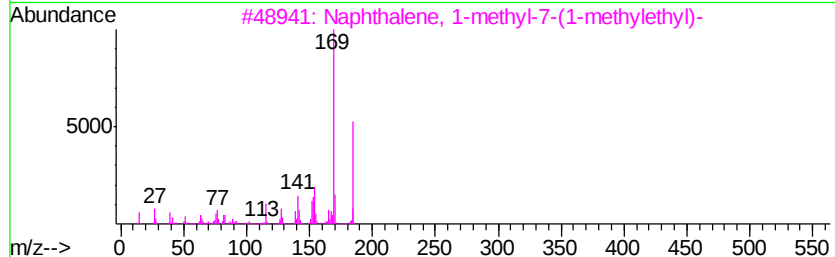
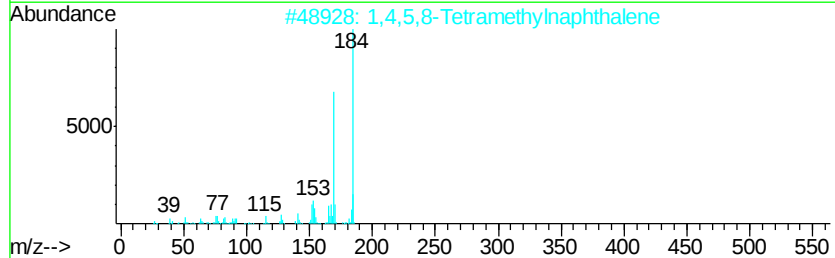
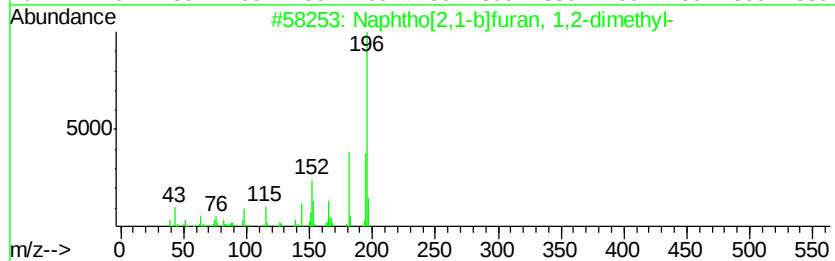
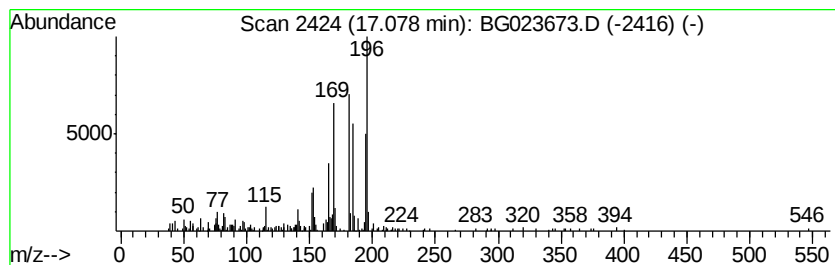
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 unknown-01 Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.08	3.08 ng/ul	311561	Phenanthrene-d10	17.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphtho[2,1-b]furan, 1,2-dimethyl-	196	C14H12O	129812-23-3	43
2		1,4,5,8-Tetramethylnaphthalene	184	C14H16	002717-39-7	38
3		Naphthalene, 1-methyl-7-(1-methy...	184	C14H16	000490-65-3	35
4		Naphthalene, 1,2,3,4-tetramethyl-	184	C14H16	003031-15-0	35
5		Naphthalene, 1-methyl-7-(1-methy...	184	C14H16	000490-65-3	35



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
Data File : BG023673.D
Acq On : 13 Aug 2016 2:00
Operator : UM/SJ
Sample : H4385-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
P003-SS001-0206-01

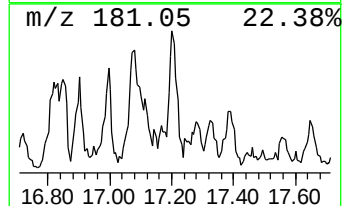
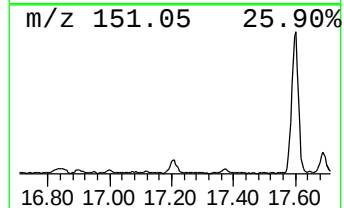
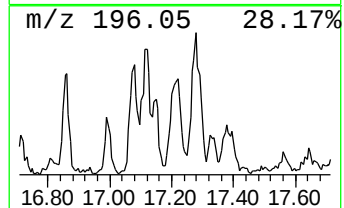
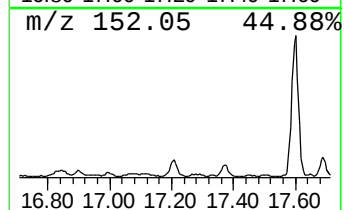
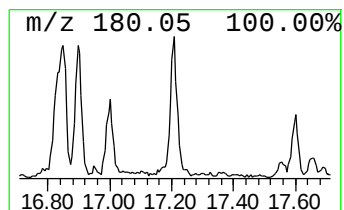
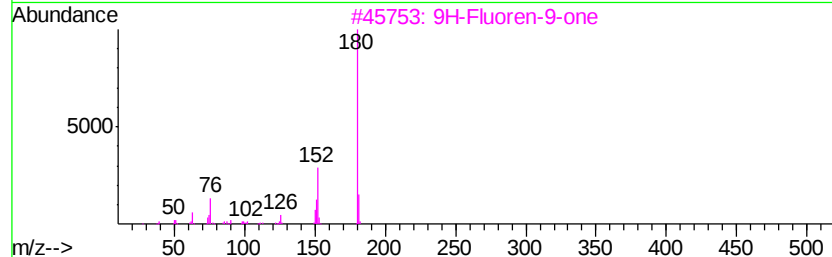
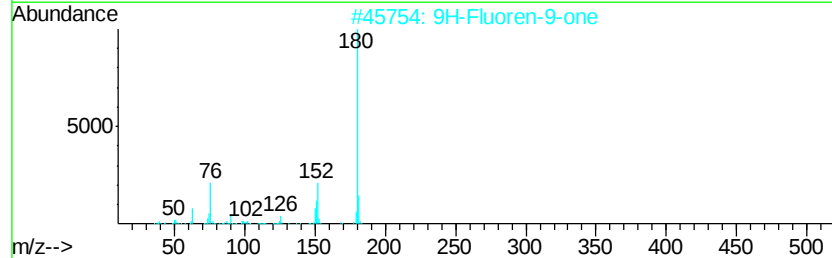
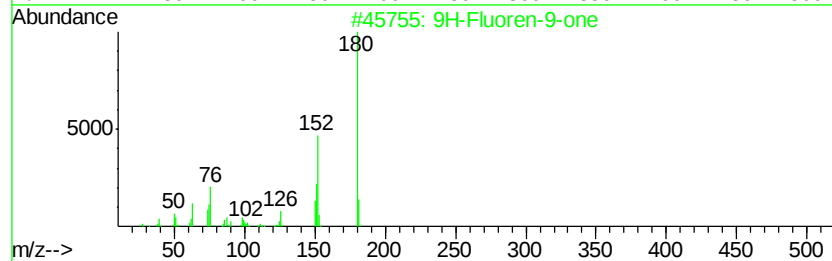
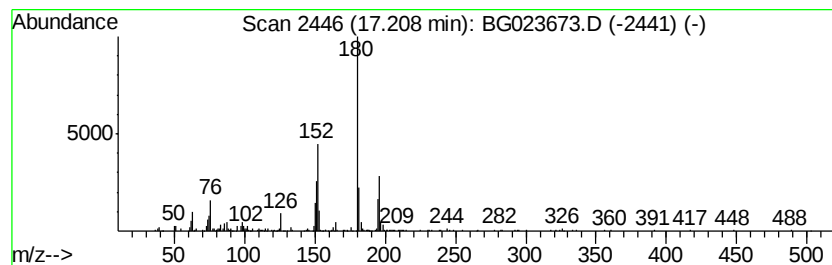
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 9H-Fluoren-9-one Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.21	3.11 ng/ul	314621	Phenanthrene-d10	17.55

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9H-Fluoren-9-one	180	C13H8O	000486-25-9	93
2			9H-Fluoren-9-one	180	C13H8O	000486-25-9	81
3			9H-Fluoren-9-one	180	C13H8O	000486-25-9	81
4			Diphenic anhydride	224	C14H8O3	006050-13-1	74
5			9H-Fluoren-9-one	180	C13H8O	000486-25-9	64



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
Data File : BG023673.D
Acq On : 13 Aug 2016 2:00
Operator : UM/SJ
Sample : H4385-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
P003-SS001-0206-01

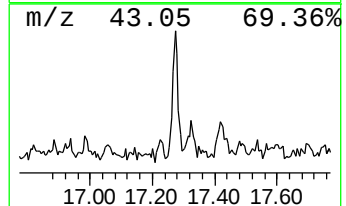
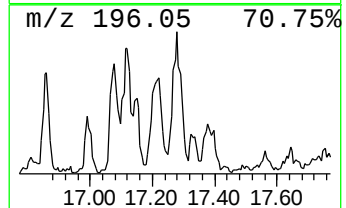
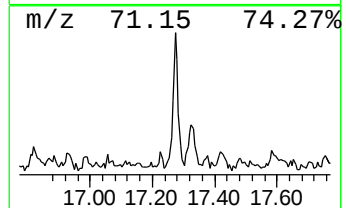
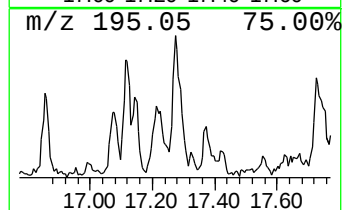
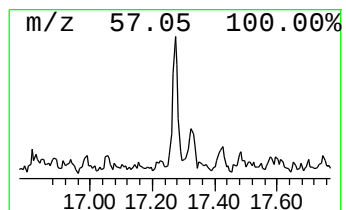
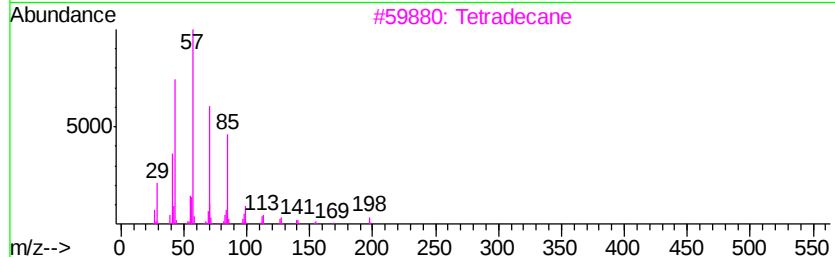
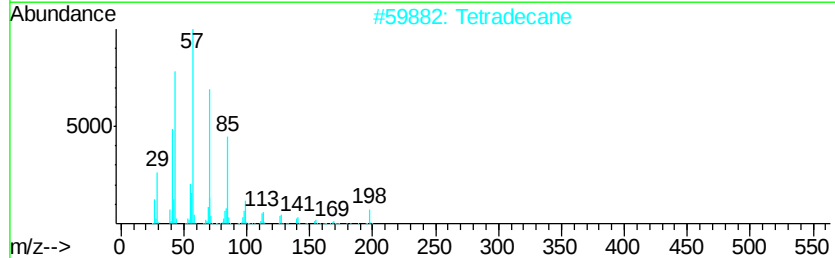
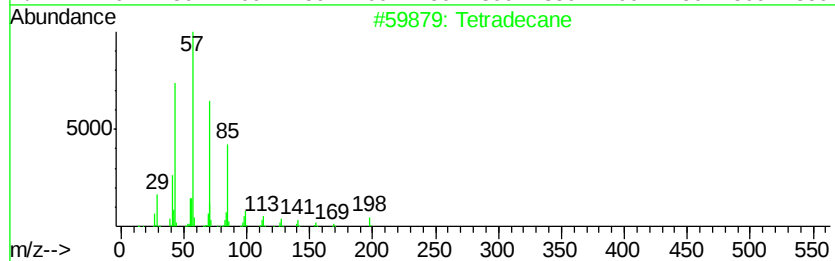
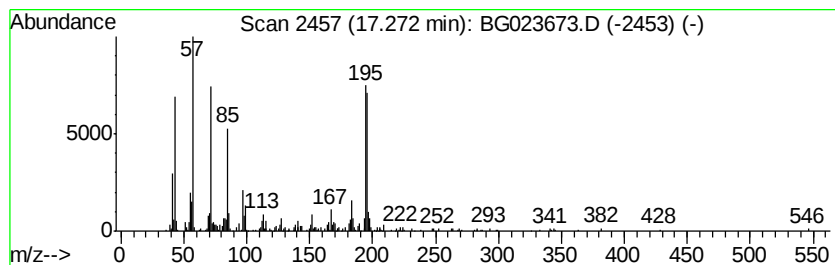
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.27	2.67 ng/ul	269262	Phenanthrene-d10	17.55

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	92
2			Tetradecane	198	C14H30	000629-59-4	59
3			Tetradecane	198	C14H30	000629-59-4	50
4			Tritetracontane	605	C43H88	007098-21-7	43
5			Tetradecane	198	C14H30	000629-59-4	42



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
Data File : BG023673.D
Acq On : 13 Aug 2016 2:00
Operator : UM/SJ
Sample : H4385-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
P003-SS001-0206-01

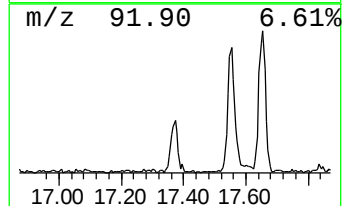
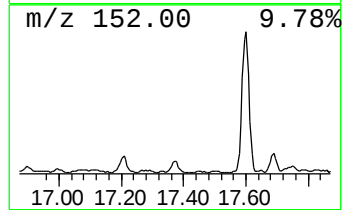
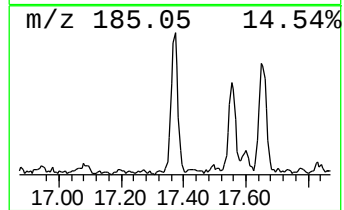
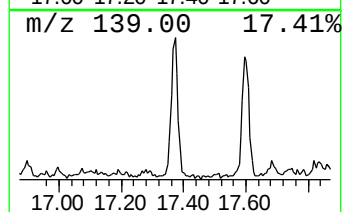
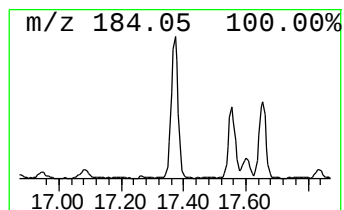
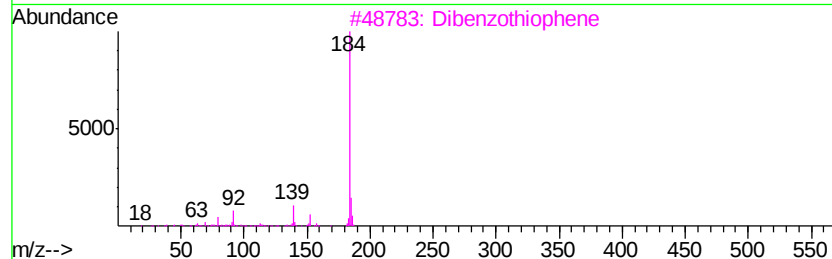
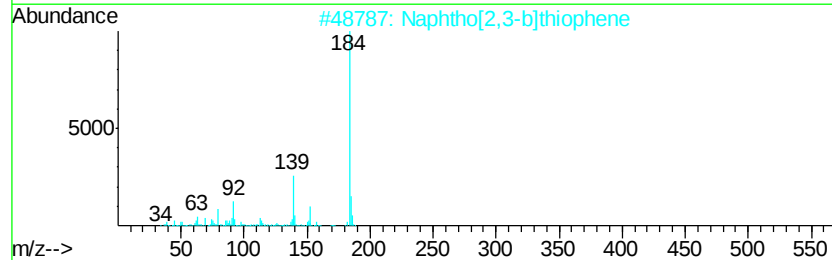
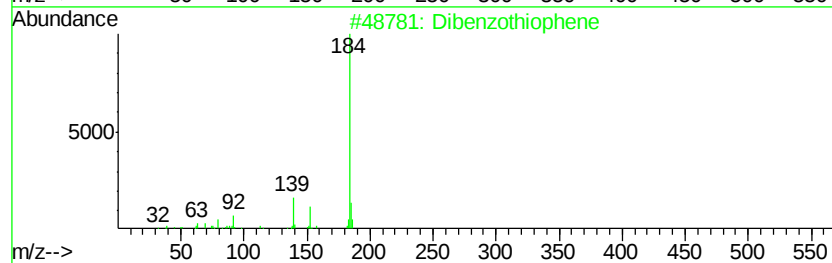
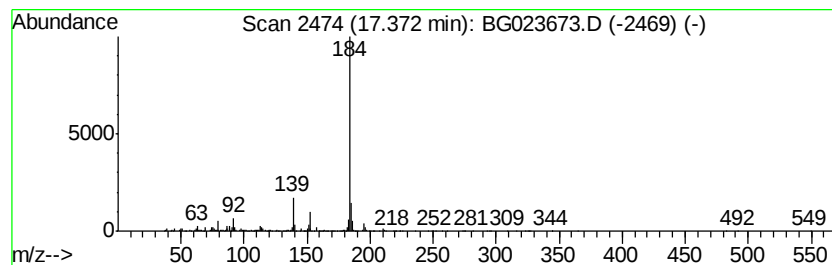
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 22 Dibenzothiophene Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.37	5.92 ng/ul	597773	Phenanthrene-d10	17.55

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzothiophene	184	C12H8S	000132-65-0	96
2			Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	95
3			Dibenzothiophene	184	C12H8S	000132-65-0	93
4			Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	93
5			Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	93



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
Data File : BG023673.D
Acq On : 13 Aug 2016 2:00
Operator : UM/SJ
Sample : H4385-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
P003-SS001-0206-01

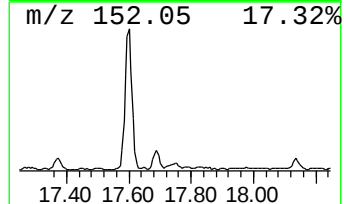
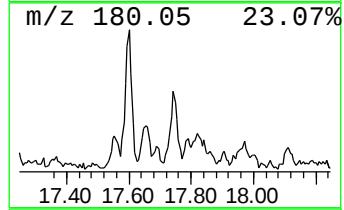
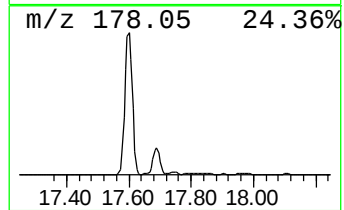
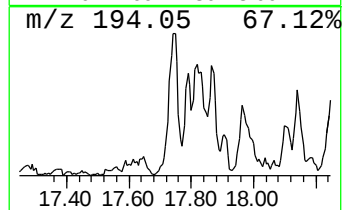
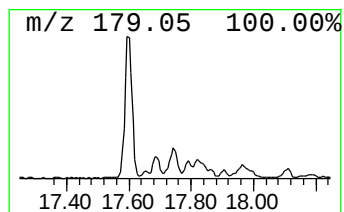
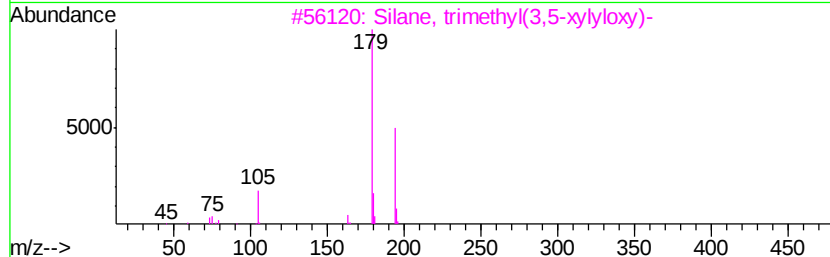
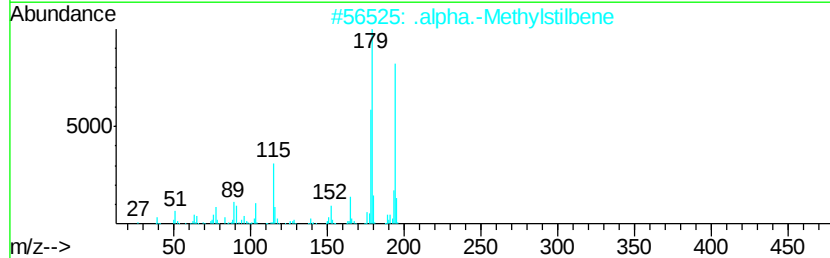
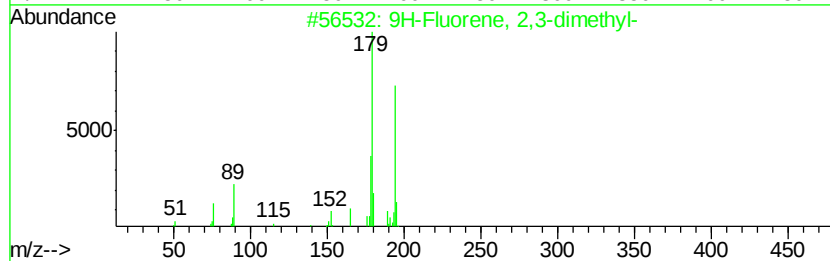
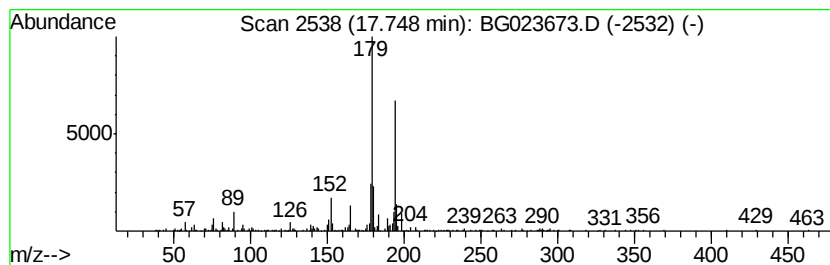
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 23 9H-Fluorene, 2,3-dimethyl- Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.75	3.33 ng/ul	335977	Phenanthrene-d10	17.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9H-Fluorene, 2,3-dimethyl-	194	C15H14	004612-63-9	87
2		.alpha.-Methylstilbene	194	C15H14	000779-51-1	59
3		Silane, trimethyl(3,5-xylyloxy)-	194	C11H18OSi	017994-05-7	58
4		.alpha.-Methylstilbene	194	C15H14	000779-51-1	53
5		Silane, trimethyl(3,5-xylyloxy)-	194	C11H18OSi	017994-05-7	50



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
Data File : BG023673.D
Acq On : 13 Aug 2016 2:00
Operator : UM/SJ
Sample : H4385-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
P003-SS001-0206-01

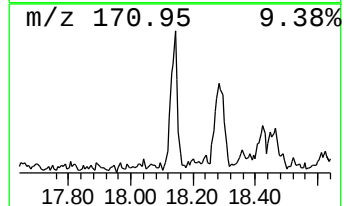
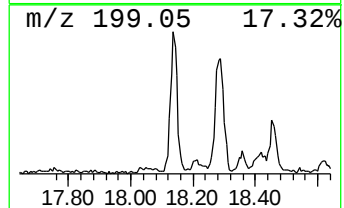
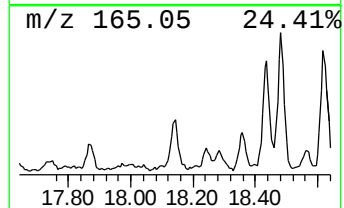
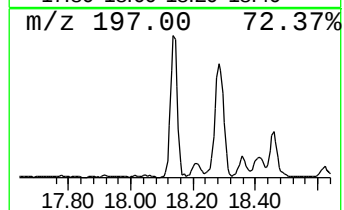
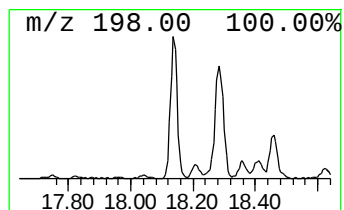
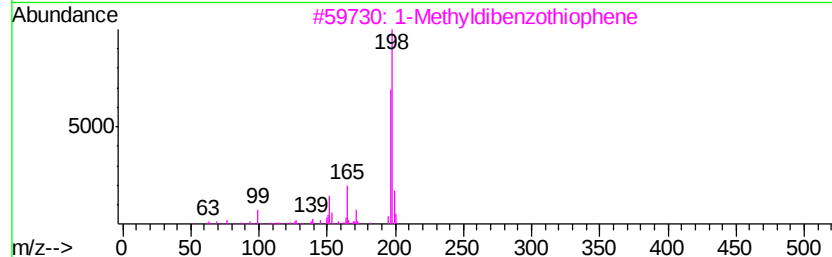
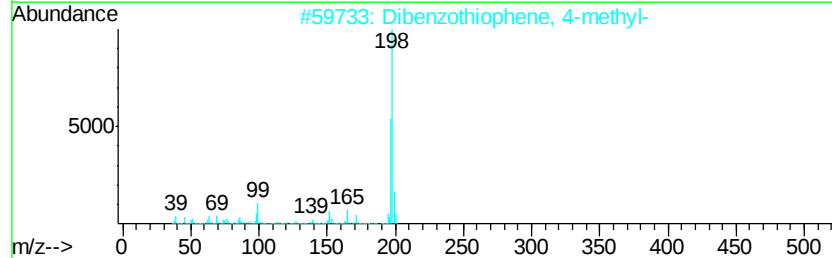
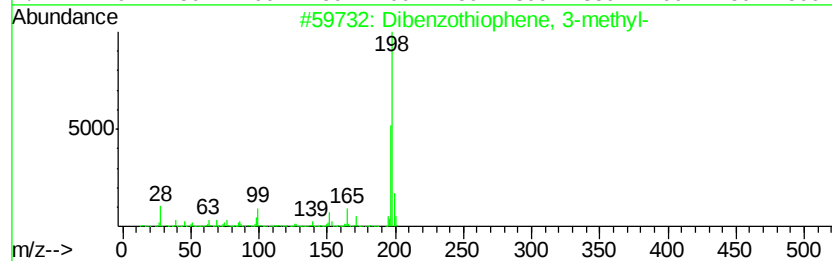
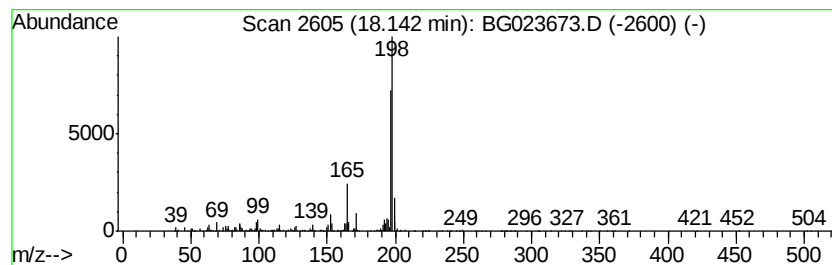
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 24 Dibenzothiophene, 3-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.14	7.89 ng/ul	796850	Phenanthrene-d10	17.55

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzothiophene, 3-methyl-	198	C13H10S	016587-52-3	91
2			Dibenzothiophene, 4-methyl-	198	C13H10S	007372-88-5	87
3			1-Methyldibenzothiophene	198	C13H10S	031317-07-4	87
4			Benzenamine, 4,4'-methylenebis-	198	C13H14N2	000101-77-9	64
5			Tacrine	198	C13H14N2	000321-64-2	64



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
Data File : BG023673.D
Acq On : 13 Aug 2016 2:00
Operator : UM/SJ
Sample : H4385-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
P003-SS001-0206-01

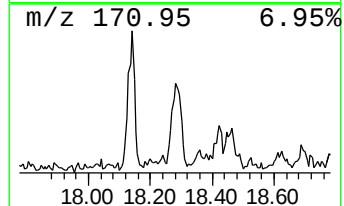
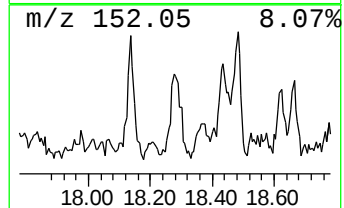
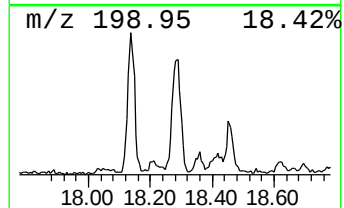
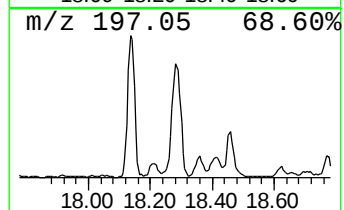
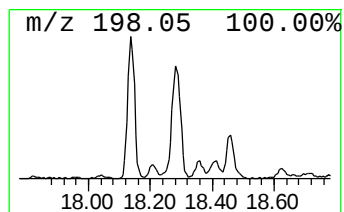
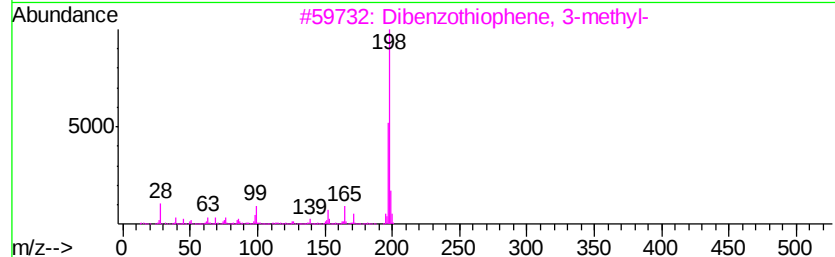
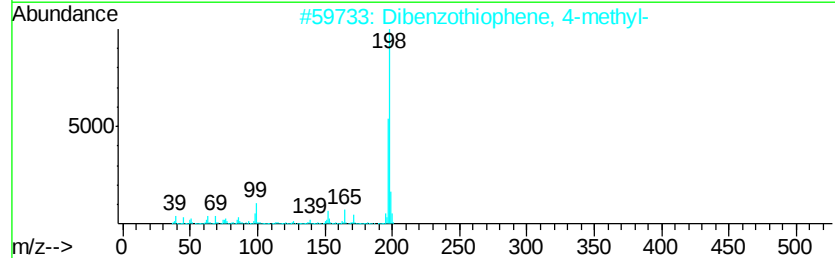
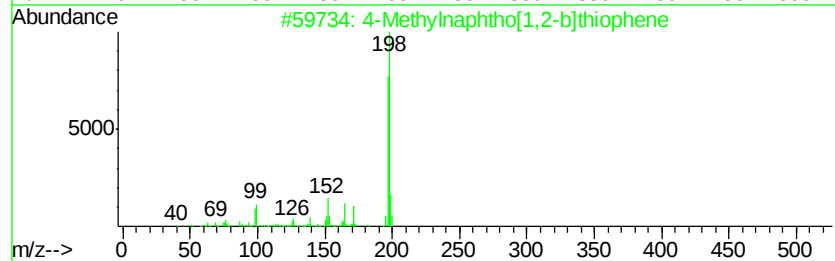
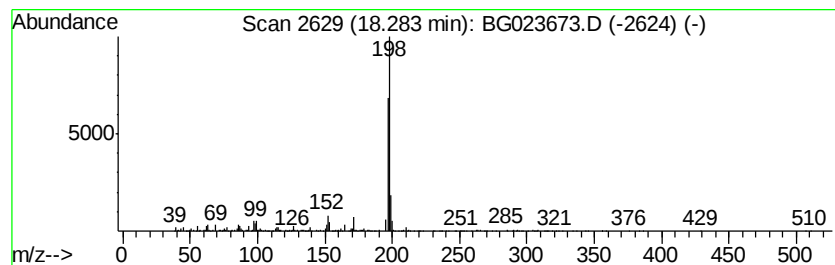
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 25 4-Methylnaphtho[1,2-b]thiop... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.28	6.41 ng/ul	647087	Phenanthrene-d10	17.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Methylnaphtho[1,2-b]thiophene	198	C13H10S	067388-11-8	94
2		Dibenzothiophene, 4-methyl-	198	C13H10S	007372-88-5	91
3		Dibenzothiophene, 3-methyl-	198	C13H10S	016587-52-3	91
4		1-Methyldibenzothiophene	198	C13H10S	031317-07-4	78
5		Benzenamine, 4,4'-methylenebis-	198	C13H14N2	000101-77-9	72



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
Data File : BG023673.D
Acq On : 13 Aug 2016 2:00
Operator : UM/SJ
Sample : H4385-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
P003-SS001-0206-01

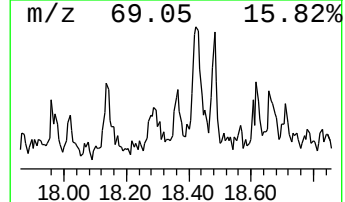
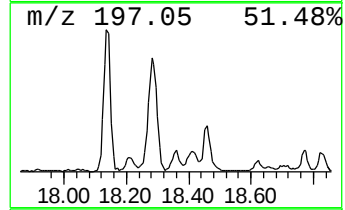
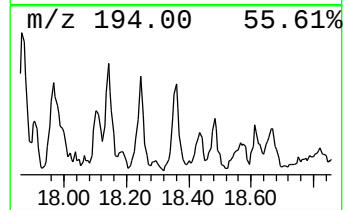
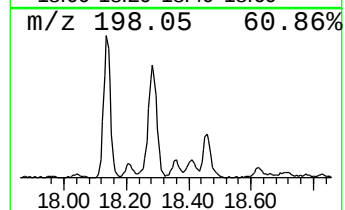
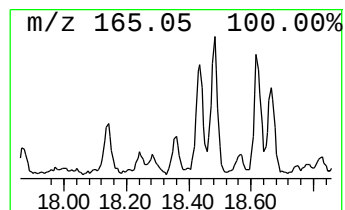
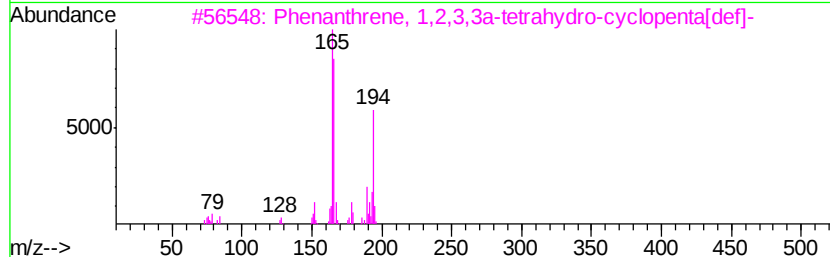
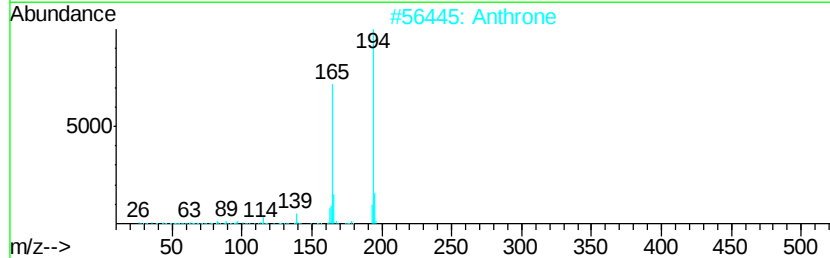
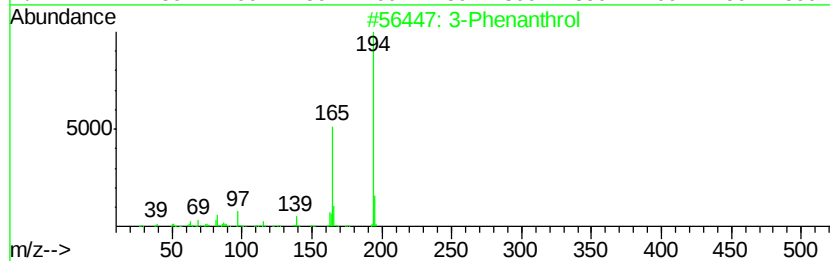
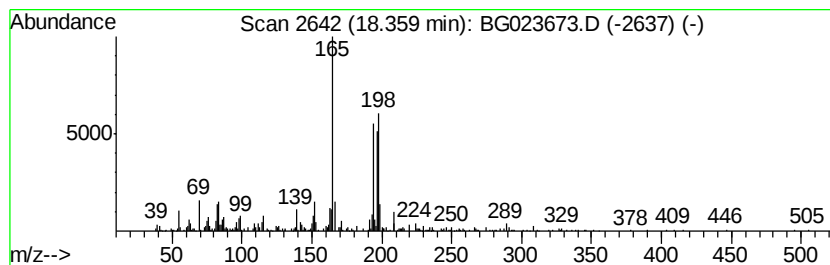
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 26 3-Phenanthrol Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.36	2.87 ng/ul	290115	Phenanthrene-d10	17.55

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Phenanthrol	194	C14H10O	000605-87-8	50
2			Anthrone	194	C14H10O	000090-44-8	50
3			Phenanthrene, 1,2,3,3a-tetrahydr...	194	C15H14	091077-10-0	45
4			2-Fluorencarboxaldehyde	194	C14H10O	030084-90-3	43
5			Anthrone	194	C14H10O	000090-44-8	43



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
Data File : BG023673.D
Acq On : 13 Aug 2016 2:00
Operator : UM/SJ
Sample : H4385-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
P003-SS001-0206-01

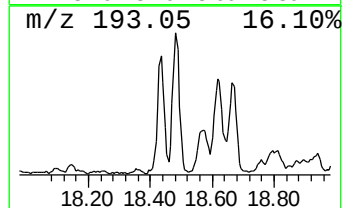
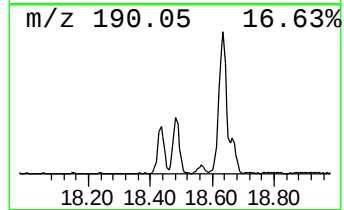
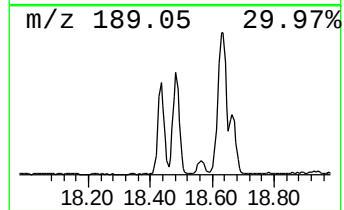
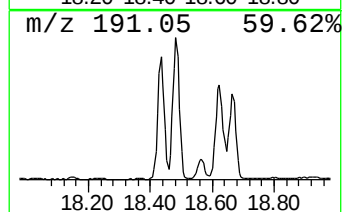
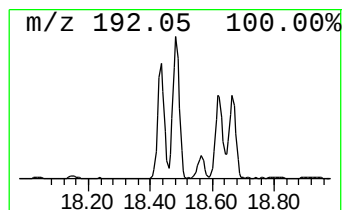
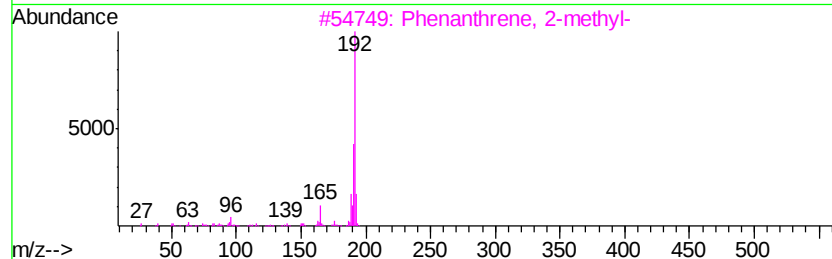
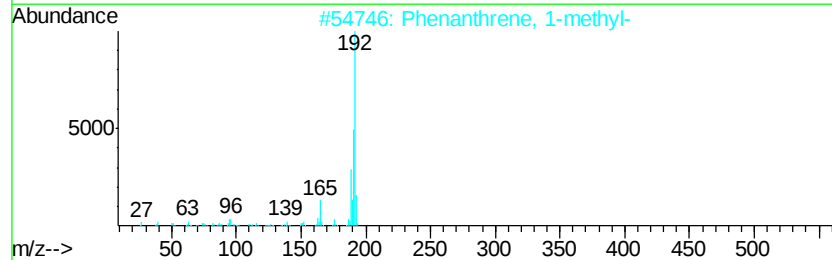
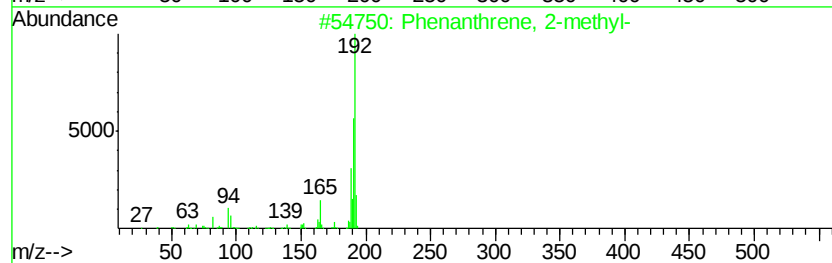
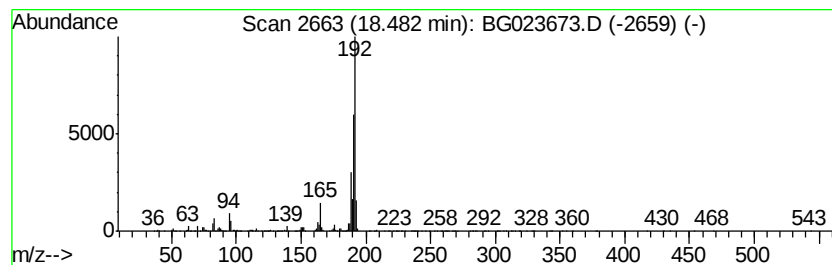
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 28 Phenanthrene, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.48	26.48 ng/ul	2674950	Phenanthrene-d10	17.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4		1H-Cyclopropa[1]phenanthrene, 1a, ...	192	C15H12	000949-41-7	96
5		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	95



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
Data File : BG023673.D
Acq On : 13 Aug 2016 2:00
Operator : UM/SJ
Sample : H4385-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
P003-SS001-0206-01

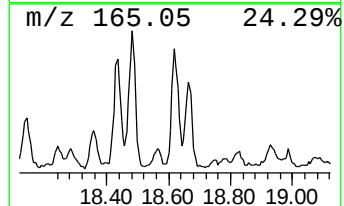
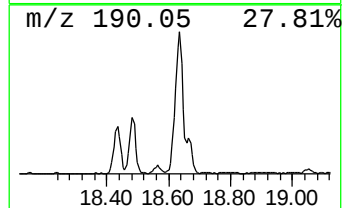
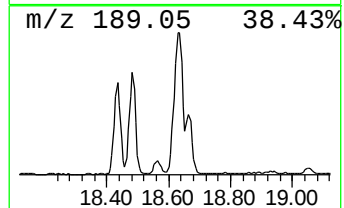
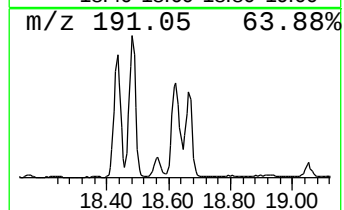
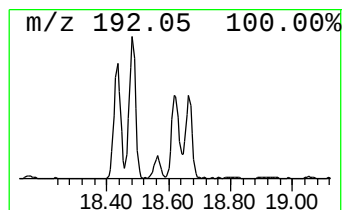
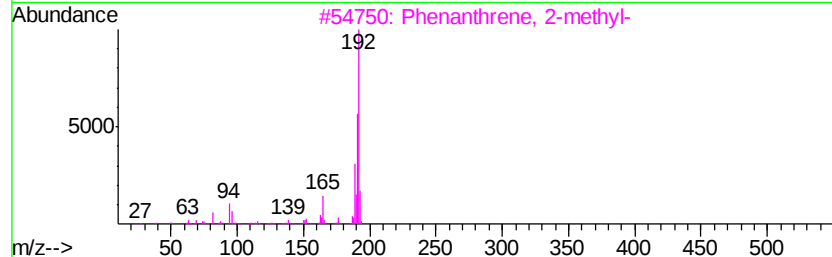
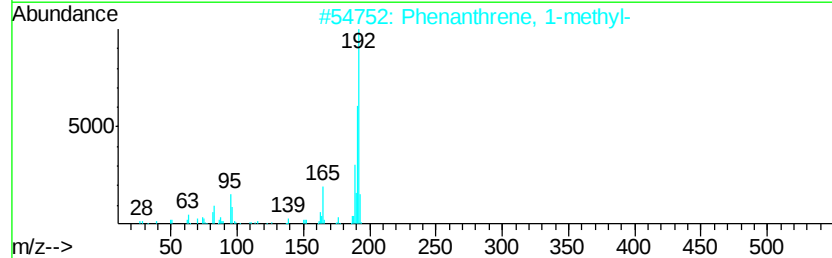
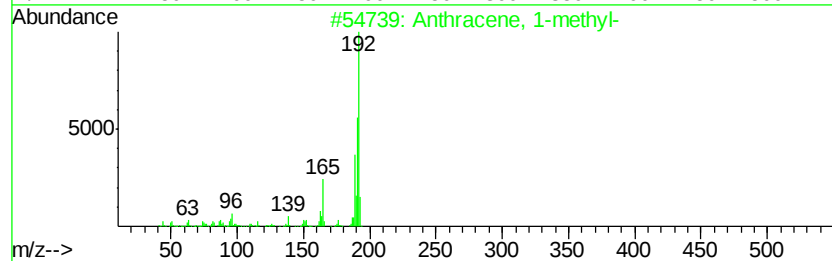
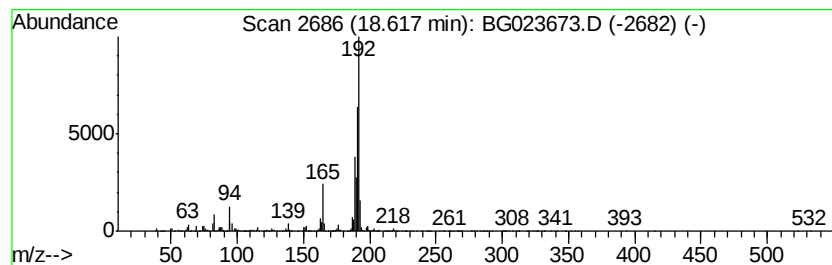
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 30 Anthracene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.62	26.38 ng/ul	2664880	Phenanthrene-d10	17.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	78
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	64
4		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	60
5		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	60



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081216\
 Data File : BG023673.D
 Acq On : 13 Aug 2016 2:00
 Operator : UM/SJ
 Sample : H4385-18
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 P003-SS001-0206-01

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Cyc...	4.46	5.4	ng/ul	187136	1	8.11	697935	20.0
Naphthalene, 1,5-...	13.96	6.7	ng/ul	567567	3	14.79	1686830	20.0
Naphthalene, 2,3-...	14.11	9.2	ng/ul	772914	3	14.79	1686830	20.0
Naphthalene, 1,4-...	14.34	4.0	ng/ul	338444	3	14.79	1686830	20.0
Naphthalene, 1,6,...	15.18	7.4	ng/ul	622648	3	14.79	1686830	20.0
Naphthalene, 2,3,...	15.46	3.4	ng/ul	284200	3	14.79	1686830	20.0
Naphthalene, 1,4,...	15.60	3.0	ng/ul	256710	3	14.79	1686830	20.0
2,4,6-Cycloheptat...	16.13	4.2	ng/ul	353587	3	14.79	1686830	20.0
[1,1'-Biphenyl]-4...	16.28	3.0	ng/ul	297750	4	17.55	2020390	20.0
1,4,5,8-Tetrameth...	16.51	3.5	ng/ul	355450	4	17.55	2020390	20.0
9H-Fluorene, 2-me...	16.85	6.9	ng/ul	695729	4	17.55	2020390	20.0
(3H)Benzo[c]pyrro...	17.00	2.5	ng/ul	256778	4	17.55	2020390	20.0
unknown-01	17.08	3.1	ng/ul	311561	4	17.55	2020390	20.0
9H-Fluoren-9-one	17.21	3.1	ng/ul	314621	4	17.55	2020390	20.0
(DEL) Alkane: Str...	17.27	2.7	ng/ul	269262	4	17.55	2020390	20.0
Dibenzothiophene	17.37	5.9	ng/ul	597773	4	17.55	2020390	20.0
9H-Fluorene, 2,3-...	17.75	3.3	ng/ul	335977	4	17.55	2020390	20.0
Dibenzothiophene,...	18.14	7.9	ng/ul	796850	4	17.55	2020390	20.0
4-Methylnaphtho[1...	18.28	6.4	ng/ul	647087	4	17.55	2020390	20.0
3-Phenanthrol	18.36	2.9	ng/ul	290115	4	17.55	2020390	20.0
Phenanthrene, 2-m...	18.48	26.5	ng/ul	2674950	4	17.55	2020390	20.0
Anthracene, 1-met...	18.62	26.4	ng/ul	2664880	4	17.55	2020390	20.0