

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
 Data File : BG018082.D
 Acq On : 24 Jul 2015 1:33
 Operator : TP/UM
 Sample : G3035-16
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C01K4

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-BG072315.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	2.763	11	13	17	rVB2	12790	12493	0.59%	0.036%
2	6.694	678	682	693	rVB	97977	161984	7.63%	0.464%
3	6.859	705	710	717	rBV	81055	121509	5.72%	0.348%
4	7.047	736	742	749	rVB	113733	174944	8.24%	0.501%
5	7.511	815	821	830	rBV	209856	325366	15.33%	0.932%
6	8.227	937	943	951	rVV2	122760	190177	8.96%	0.545%
7	8.333	955	961	971	rVB	35844	57859	2.73%	0.166%
8	8.668	1011	1018	1027	rVB	105317	174911	8.24%	0.501%
9	9.385	1133	1140	1145	rBV	91705	147267	6.94%	0.422%
10	9.920	1225	1231	1239	rBV	149956	247352	11.65%	0.709%
11	10.161	1266	1272	1278	rBV2	82993	129585	6.10%	0.371%
12	10.296	1287	1295	1301	rBV	332475	560983	26.43%	1.607%
13	10.437	1311	1319	1331	rBV	113451	190942	9.00%	0.547%
14	10.678	1354	1360	1366	rBV	33959	52879	2.49%	0.152%
15	11.330	1467	1471	1478	rBV6	5947	13683	0.64%	0.039%
16	11.747	1537	1542	1547	rBV3	10513	16343	0.77%	0.047%
17	11.964	1575	1579	1587	rVB3	10487	20857	0.98%	0.060%
18	12.346	1638	1644	1649	rBV7	8736	22194	1.05%	0.064%
19	12.605	1681	1688	1691	rVV8	6297	12592	0.59%	0.036%
20	12.716	1703	1707	1711	rVB6	9983	13335	0.63%	0.038%
21	12.781	1711	1718	1721	rBV6	10275	17149	0.81%	0.049%
22	12.951	1741	1747	1753	rBV5	21073	48834	2.30%	0.140%
23	13.010	1753	1757	1762	rVB3	18861	24970	1.18%	0.072%
24	13.104	1769	1773	1776	rBV5	7271	11379	0.54%	0.033%
25	13.351	1809	1815	1819	rBV8	6346	15471	0.73%	0.044%
26	13.498	1836	1840	1844	rVB5	14162	19152	0.90%	0.055%
27	13.598	1851	1857	1861	rBV	304057	411387	19.38%	1.179%
28	13.645	1861	1865	1869	rVB	81641	102129	4.81%	0.293%
29	13.721	1875	1878	1884	rVB7	14271	22577	1.06%	0.065%
30	13.868	1897	1903	1908	rVV	320471	460056	21.67%	1.318%
31	13.915	1908	1911	1916	rVV6	16249	27360	1.29%	0.078%
32	14.009	1923	1927	1933	rVB	41240	57271	2.70%	0.164%
33	14.097	1936	1942	1946	rBV3	31777	54347	2.56%	0.156%
34	14.179	1946	1956	1964	rBV2	501223	849327	40.01%	2.433%

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Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG072315.M
 Title : SVOA CALIBRATION

35	14.244	1964	1967	1970	rBV	25574	36900	1.74%	0.106%
36	14.397	1988	1993	2005	rVB	108998	170601	8.04%	0.489%
37	14.726	2045	2049	2055	rVB8	16621	26161	1.23%	0.075%
38	14.902	2075	2079	2085	rBV8	21164	40763	1.92%	0.117%
39	14.973	2087	2091	2095	rBV2	40485	57003	2.69%	0.163%
40	15.055	2102	2105	2110	rBV2	44694	60602	2.85%	0.174%
41	15.178	2121	2126	2132	rVB	401852	548620	25.85%	1.572%
42	15.237	2132	2136	2144	rBV3	33765	71740	3.38%	0.206%
43	15.307	2145	2148	2155	rVB	53751	71156	3.35%	0.204%
44	15.924	2249	2253	2254	rBV2	59362	72795	3.43%	0.209%
45	16.718	2385	2388	2391	rBV2	45319	56177	2.65%	0.161%
46	16.935	2420	2425	2429	rBV	623269	902295	42.51%	2.585%
47	16.976	2429	2432	2437	rVB	292720	361075	17.01%	1.035%
48	17.035	2437	2442	2446	rBV	367194	507322	23.90%	1.454%
49	17.857	2578	2582	2586	rBV2	77638	123115	5.80%	0.353%
50	18.016	2605	2609	2613	rVB3	183317	240762	11.34%	0.690%
51	18.075	2616	2619	2623	rVB2	66548	70477	3.32%	0.202%
52	18.304	2655	2658	2664	rBV2	81209	121249	5.71%	0.347%
53	18.792	2739	2741	2746	rVB6	57283	64326	3.03%	0.184%
54	18.874	2751	2755	2764	rVB	576240	777491	36.63%	2.228%
55	19.009	2773	2778	2782	rVB	566773	731872	34.48%	2.097%
56	19.085	2788	2791	2798	rVB4	93766	153745	7.24%	0.441%
57	19.162	2799	2804	2809	rVB	734958	905034	42.64%	2.593%
58	19.220	2812	2814	2818	rVV2	50510	57367	2.70%	0.164%
59	19.344	2830	2835	2837	rBV2	499249	703466	33.14%	2.016%
60	19.373	2837	2840	2844	rVB	478014	610293	28.75%	1.749%
61	19.503	2858	2862	2867	rBV	126023	148992	7.02%	0.427%
62	19.591	2873	2877	2878	rBV	255929	313294	14.76%	0.898%
63	19.608	2878	2880	2886	rVB2	414664	455228	21.45%	1.304%
64	19.732	2897	2901	2905	rBV	433062	505229	23.80%	1.448%
65	19.779	2906	2909	2916	rVB2	197751	321476	15.14%	0.921%
66	19.873	2920	2925	2930	rVB	1495041	1832119	86.31%	5.249%
67	19.926	2930	2934	2938	rBV	184004	218673	10.30%	0.627%
68	20.078	2956	2960	2964	rVB	212490	259755	12.24%	0.744%
69	20.155	2968	2973	2978	rBV	1736402	2027777	95.53%	5.810%
70	20.208	2978	2982	2994	rVB3	492612	678224	31.95%	1.943%
71	20.307	2995	2999	3008	rVB2	516513	782548	36.87%	2.242%

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72	20.401	3012	3015	3019	rBV3	119917	140326	6.61%	0.402%
73	20.472	3019	3027	3033	rBV	1170910	2091624	98.54%	5.993%
74	20.619	3048	3052	3058	rVB	683943	816183	38.45%	2.339%
75	20.683	3059	3063	3066	rBV2	333686	399603	18.83%	1.145%
76	20.777	3075	3079	3083	rBV3	182056	251020	11.83%	0.719%
77	20.889	3093	3098	3102	rVB2	579633	693347	32.66%	1.987%
78	20.942	3103	3107	3113	rVB	355533	444687	20.95%	1.274%
79	21.001	3113	3117	3119	rBV	172551	201101	9.47%	0.576%
80	21.024	3119	3121	3126	rVB2	82286	93656	4.41%	0.268%
81	21.071	3126	3129	3132	rBV	350727	379624	17.88%	1.088%
82	21.142	3135	3141	3143	rVV2	932521	1469040	69.21%	4.209%
83	21.177	3143	3147	3153	rVB	1776794	2122705	100.00%	6.082%
84	21.312	3164	3170	3174	rVB2	72733	138872	6.54%	0.398%
85	21.483	3194	3199	3204	rVB	600869	819953	38.63%	2.349%
86	21.535	3205	3208	3214	rVB2	219805	266670	12.56%	0.764%
87	21.600	3216	3219	3225	rBV2	222360	302474	14.25%	0.867%
88	21.729	3239	3241	3246	rVB2	85349	95828	4.51%	0.275%
89	21.794	3248	3252	3256	rBV3	105196	161427	7.60%	0.463%
90	21.941	3274	3277	3281	rBV4	121667	171036	8.06%	0.490%
91	22.152	3310	3313	3317	rVB3	152039	170856	8.05%	0.490%
92	22.675	3396	3402	3407	rBV2	314854	700462	33.00%	2.007%
93	22.840	3426	3430	3438	rVB	56955	106815	5.03%	0.306%
94	23.139	3476	3481	3484	rBV	156575	264186	12.45%	0.757%
95	23.186	3485	3489	3493	rVV	245884	419806	19.78%	1.203%
96	23.233	3493	3497	3504	rVB	322454	547625	25.80%	1.569%
97	23.327	3507	3513	3519	rBV2	554452	999356	47.08%	2.863%
98	23.868	3601	3605	3613	rVB3	87192	165767	7.81%	0.475%
99	25.531	3882	3888	3898	rVB2	132050	345837	16.29%	0.991%
100	26.201	3997	4002	4016	rVB2	90308	267442	12.60%	0.766%

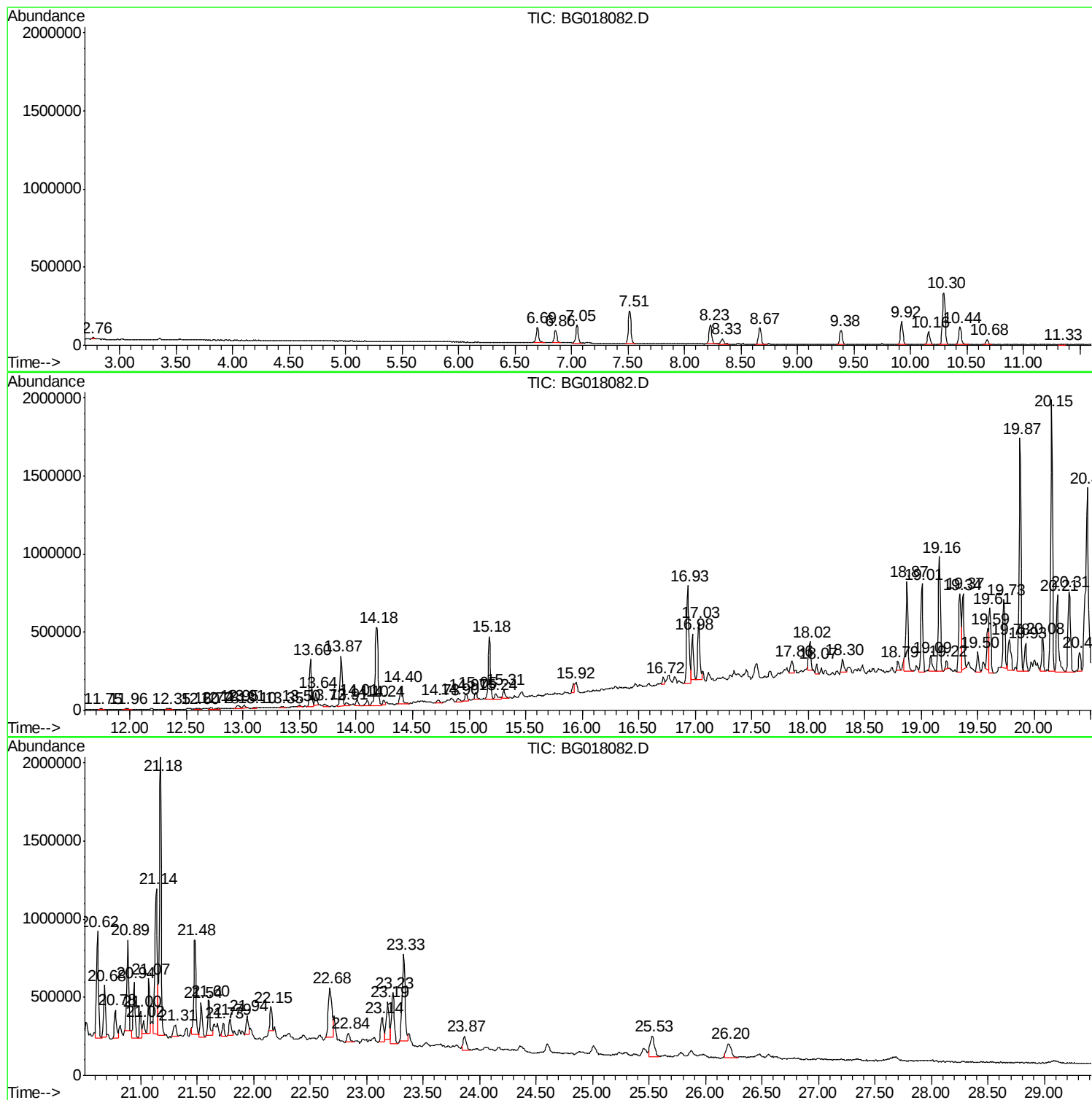
Sum of corrected areas: 34901714

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG072315.M
Quant Title : SVOA CALIBRATION

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



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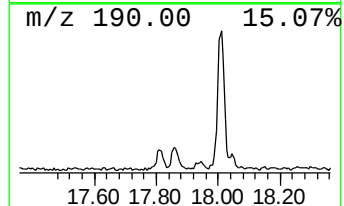
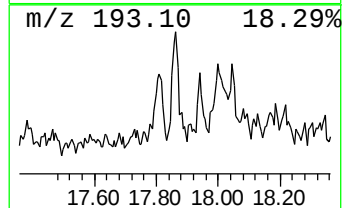
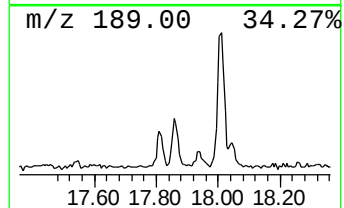
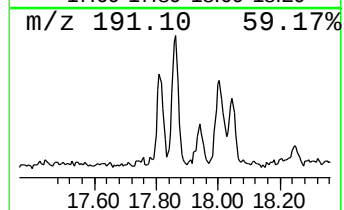
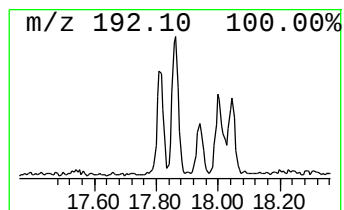
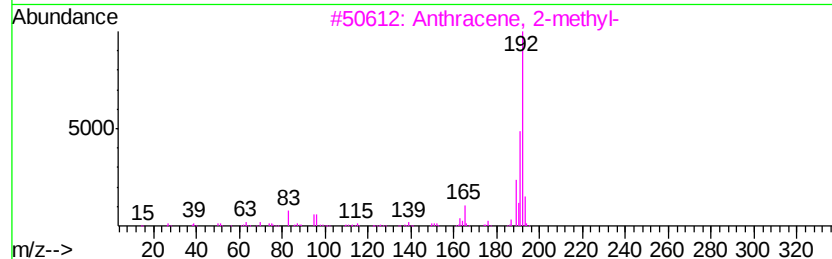
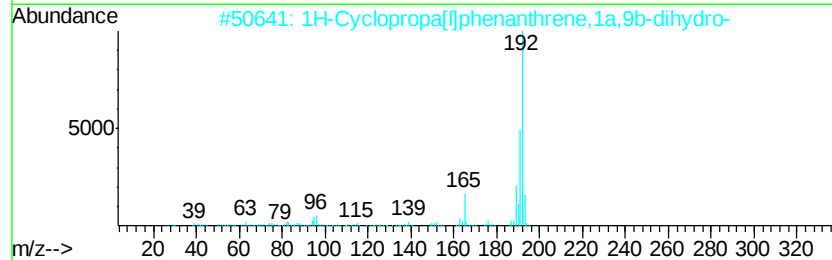
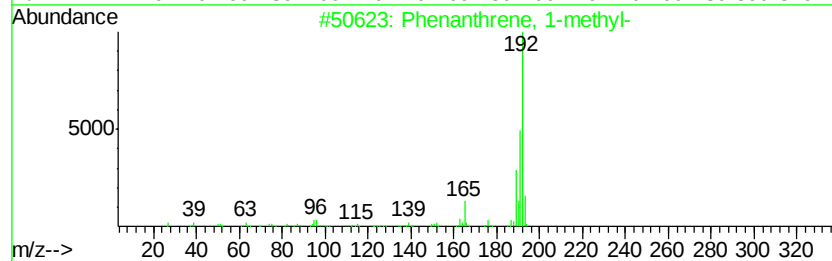
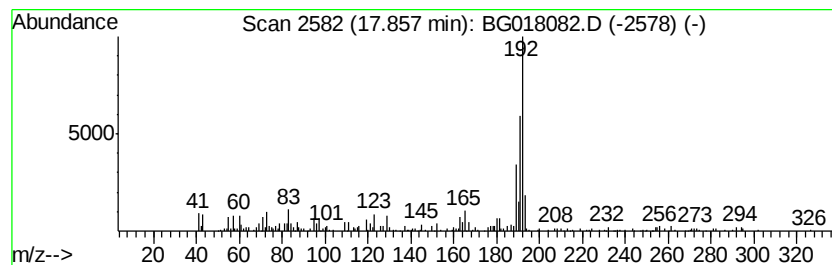
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG072315.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Phenanthrene, 1-methyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.86	2.73 ng/ul	123115	Phenanthrene-d10		16.93	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	89
2		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	89
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	81
4		Anthracene, 1-methyl-	192	C15H12	000610-48-0	81
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	81



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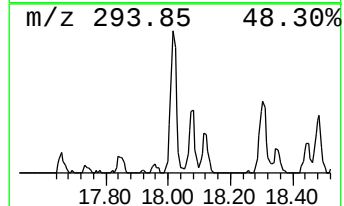
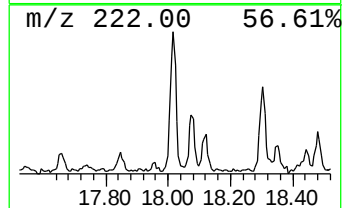
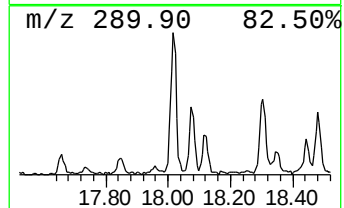
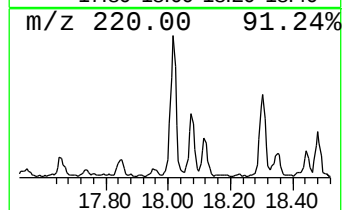
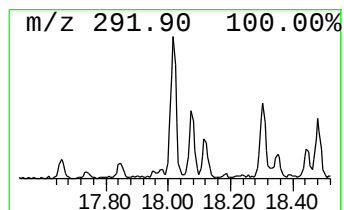
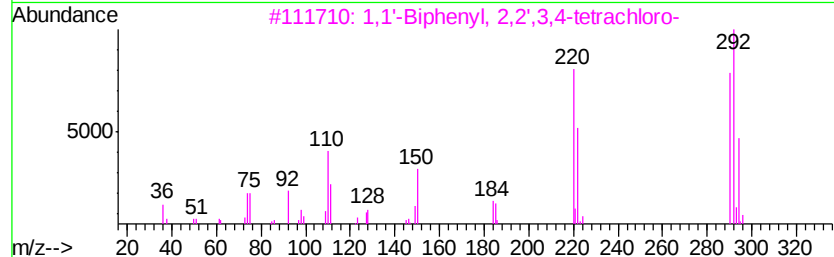
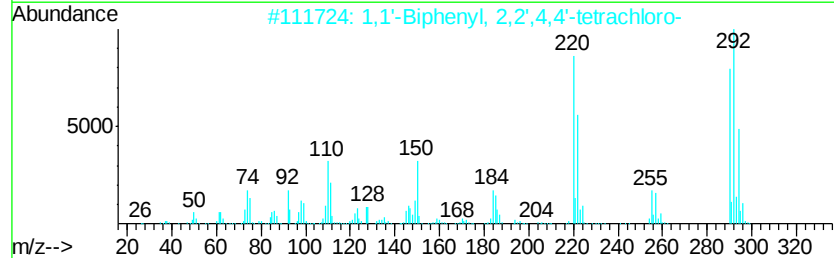
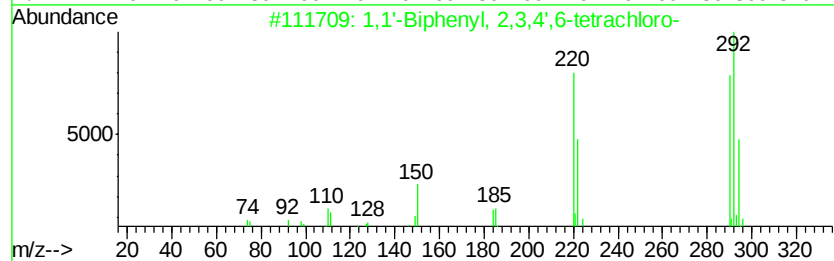
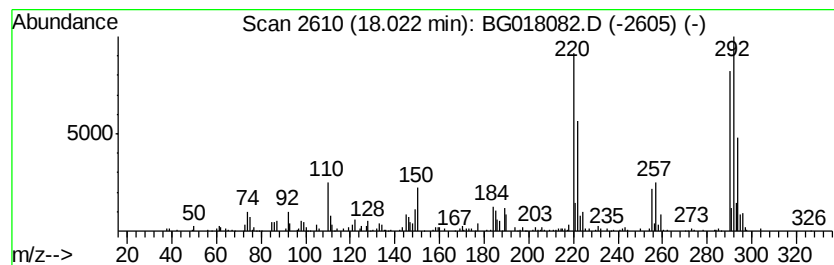
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 1,1'-Biphenyl, 2,3,4',6-tet... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.02	5.34 ng/ul	240762	Phenanthrene-d10	16.93

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,4',6-tetrachl...	290	C12H6Cl4	052663-58-8	99		
2	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99		
3	1,1'-Biphenyl, 2,2',3,4-tetrachl...	290	C12H6Cl4	052663-59-9	99		
4	1,1'-Biphenyl, 2,4,4',6-tetrachl...	290	C12H6Cl4	032598-12-2	98		
5	1,1'-Biphenyl, 2,3',5,5'-tetrach...	290	C12H6Cl4	041464-42-0	95		



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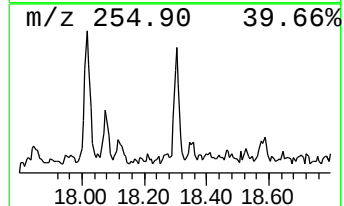
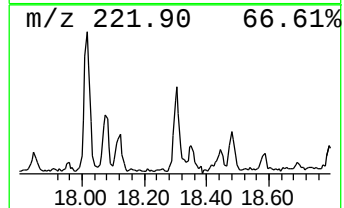
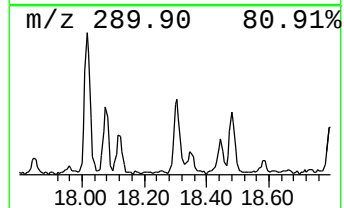
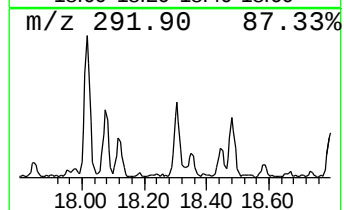
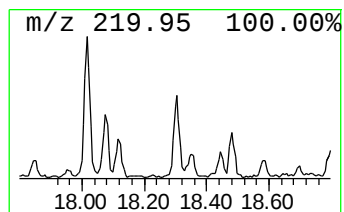
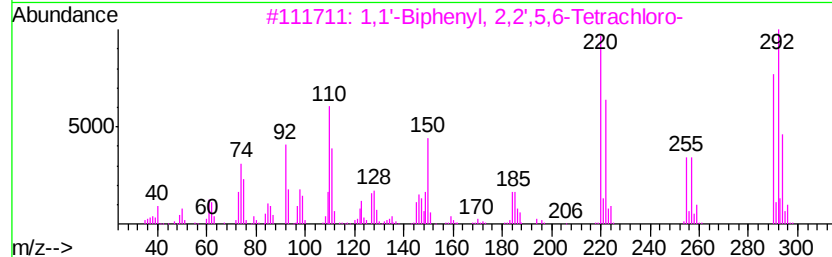
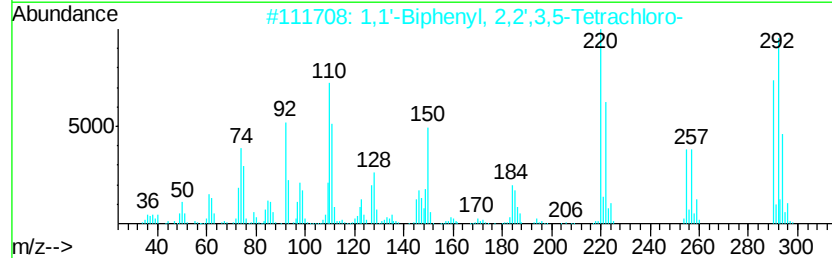
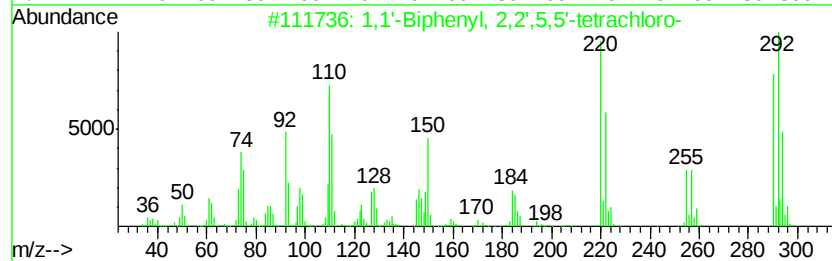
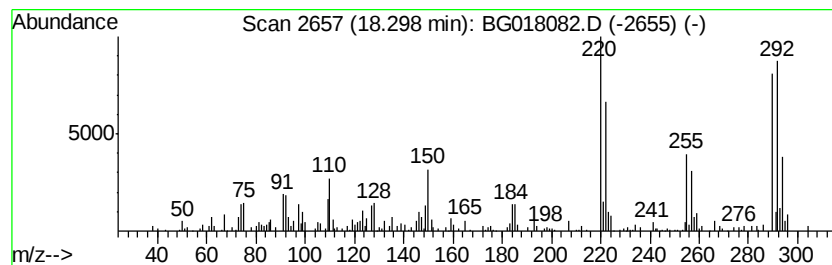
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1,1'-Biphenyl, 2,2',5,5'-te... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.30	2.69 ng/ul	121249	Phenanthrene-d10	16.93		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5,5'-tetrachl...	290	C12H6Cl4	035693-99-3	96	
2	1,1'-Biphenyl, 2,2',3,5-Tetrachl...	290	C12H6Cl4	070362-46-8	96	
3	1,1'-Biphenyl, 2,2',5,6-Tetrachl...	290	C12H6Cl4	041464-41-9	96	
4	1,1'-Biphenyl, 2,2',5,6-Tetrachl...	290	C12H6Cl4	041464-41-9	95	
5	1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	95	



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
Data File : BG018082.D
Acq On : 24 Jul 2015 1:33
Operator : TP/UM
Sample : G3035-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C01K4

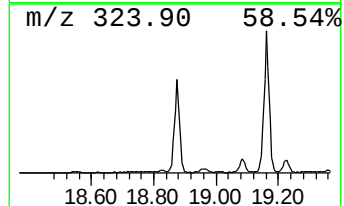
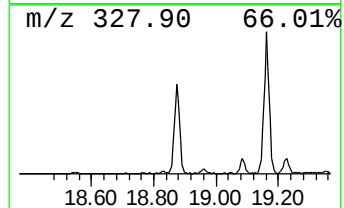
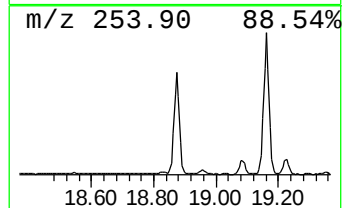
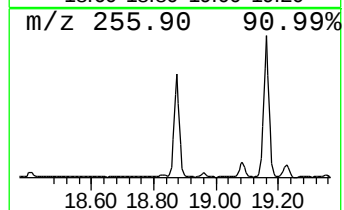
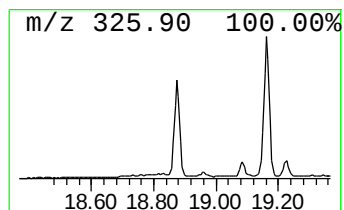
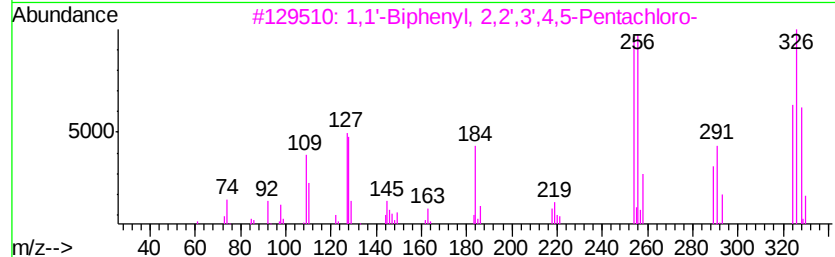
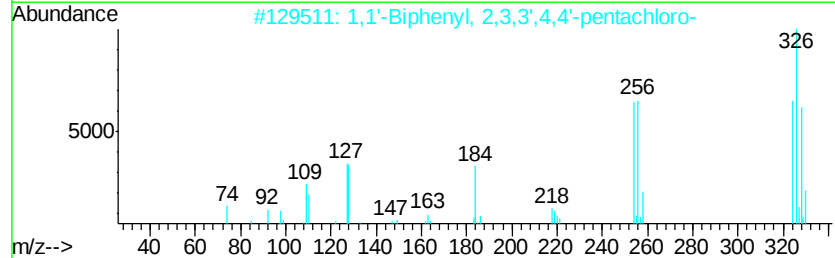
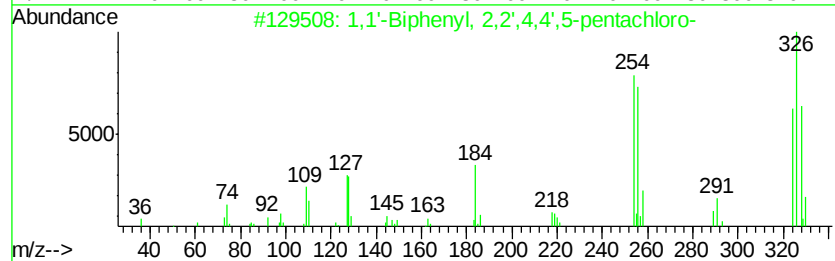
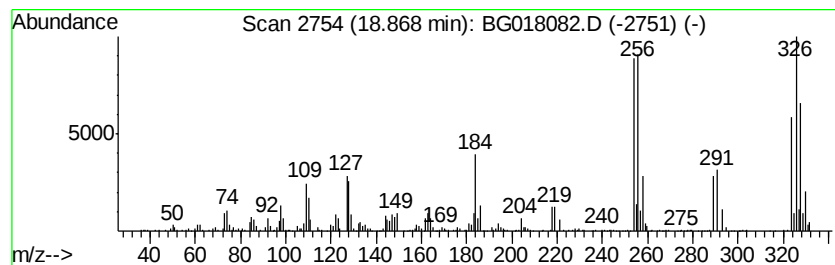
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,1'-Biphenyl, 2,2',4,4',5-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.87	17.23 ng/ul	777491	Phenanthrene-d10	16.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',4,4',5-penta...	324	C12H5Cl5	038380-01-7	99
2		1,1'-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	99
3		1,1'-Biphenyl, 2,2',3',4,5-Penta...	324	C12H5Cl5	041464-51-1	99
4		1,1'-Biphenyl, 2,2',4,5,5'-penta...	324	C12H5Cl5	037680-73-2	99
5		1,1'-Biphenyl, 2,2',4,5',6-Penta...	324	C12H5Cl5	060145-21-3	99



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
Data File : BG018082.D
Acq On : 24 Jul 2015 1:33
Operator : TP/UM
Sample : G3035-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C01K4

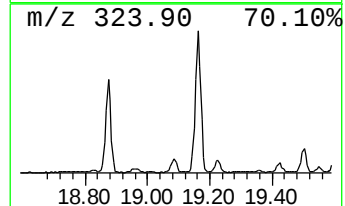
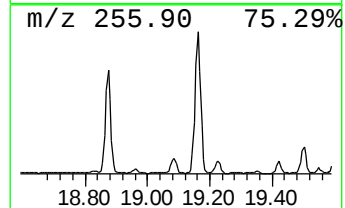
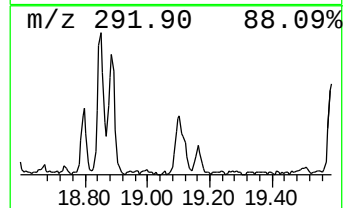
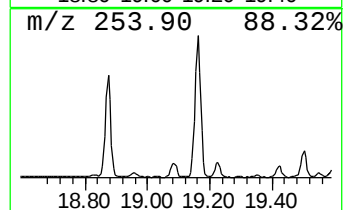
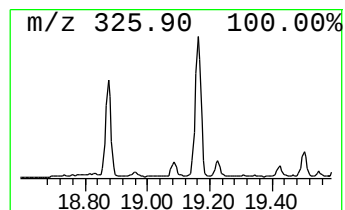
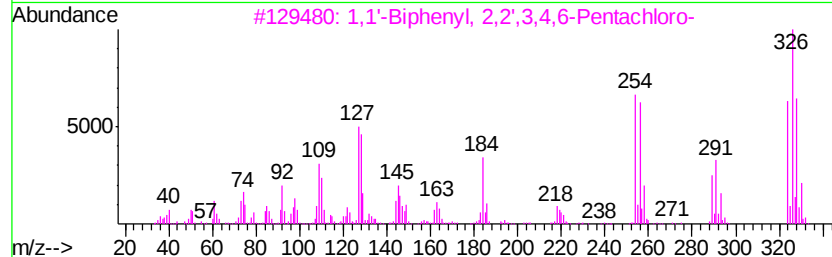
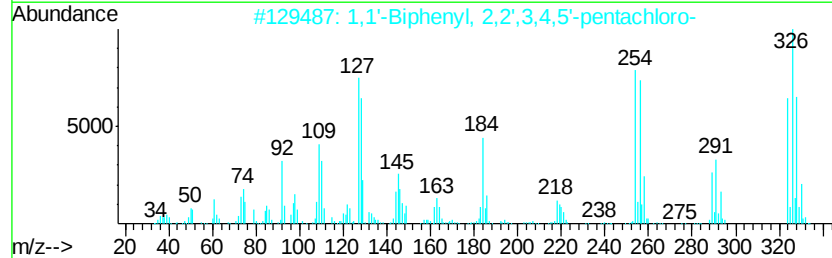
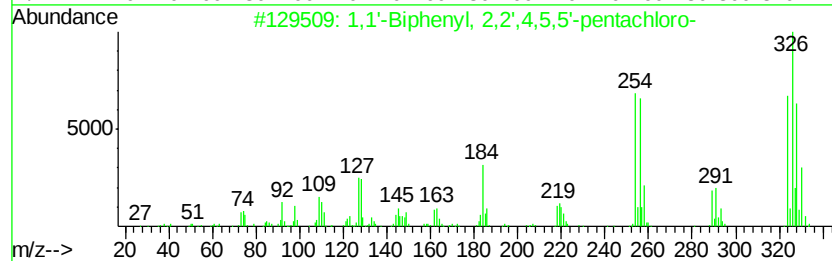
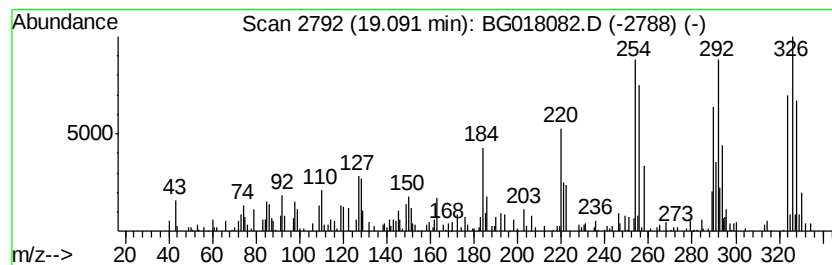
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1,1'-Biphenyl, 2,2',4,5,5'-... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.09	2.09 ng/ul	153745	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',4,5,5'-penta...	324	C12H5Cl5	037680-73-2	99
2		1,1'-Biphenyl, 2,2',3,4,5'-penta...	324	C12H5Cl5	038380-02-8	99
3		1,1'-Biphenyl, 2,2',3,4,6-Pentac...	324	C12H5Cl5	055215-17-3	99
4		1,1'-Biphenyl, 2,2',3,5,6-Pentac...	324	C12H5Cl5	073575-56-1	99
5		1,1'-Biphenyl, 2,3,4',5,6-Pentac...	324	C12H5Cl5	068194-11-6	98



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
Data File : BG018082.D
Acq On : 24 Jul 2015 1:33
Operator : TP/UM
Sample : G3035-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampled :
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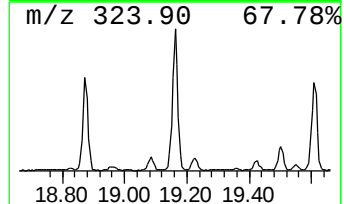
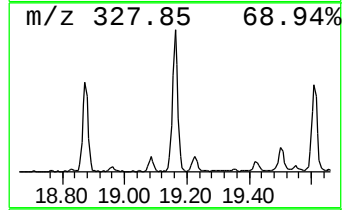
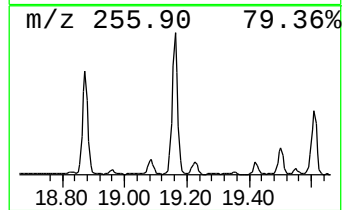
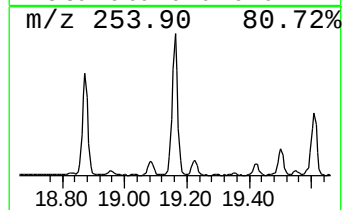
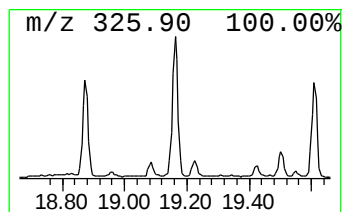
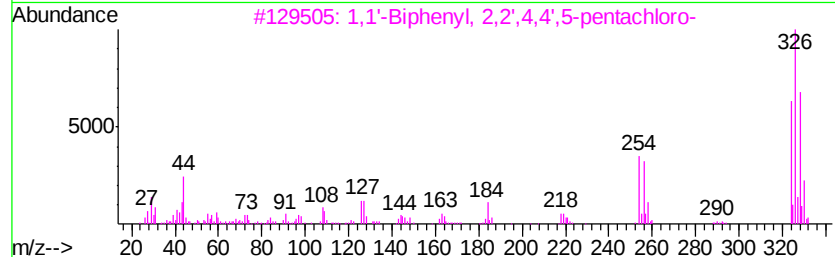
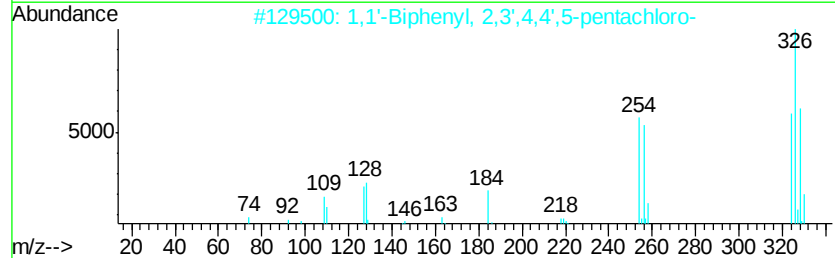
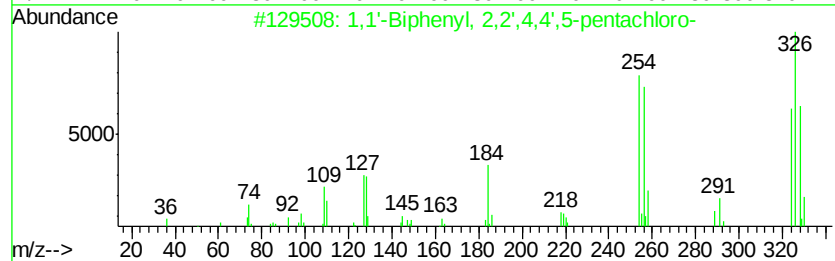
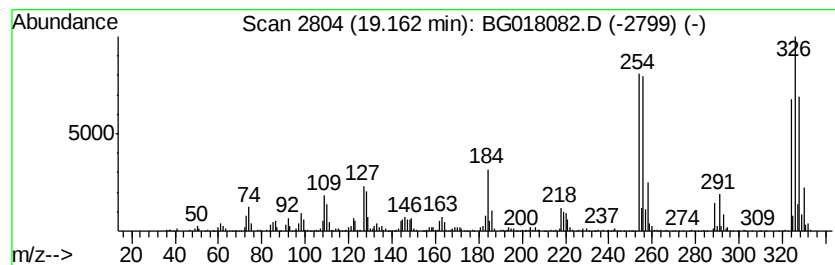
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1,1'-Biphenyl, 2,3,3',4,4'-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.16	12.32 ng/ul	905034	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',4,4',5-penta...	324	C12H5Cl5	038380-01-7	99
2		1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	99
3		1,1'-Biphenyl, 2,2',4,4',5-penta...	324	C12H5Cl5	038380-01-7	99
4		1,1'-Biphenyl, 2,2',4,5,5'-penta...	324	C12H5Cl5	037680-73-2	99
5		1,1'-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	99



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
Data File : BG018082.D
Acq On : 24 Jul 2015 1:33
Operator : TP/UM
Sample : G3035-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampled :
C01K4

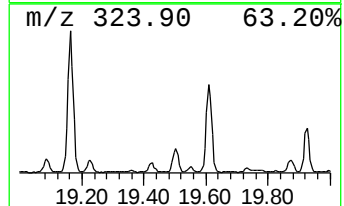
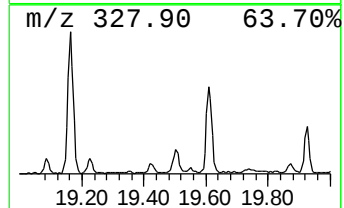
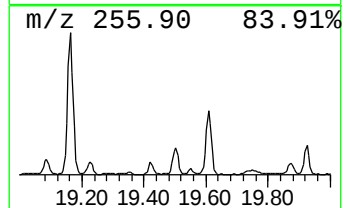
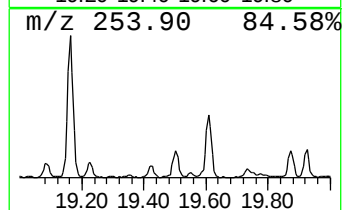
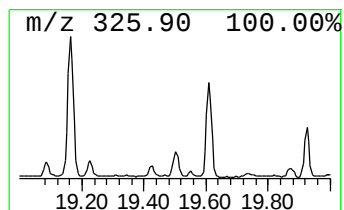
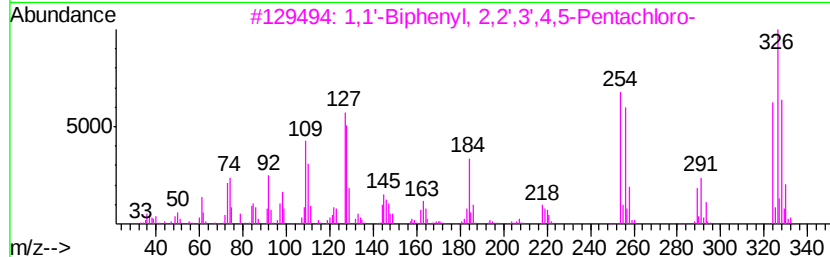
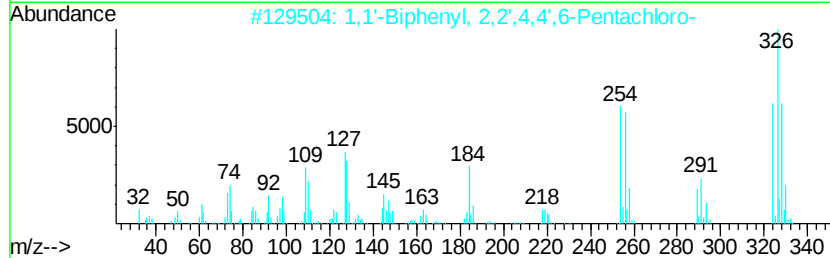
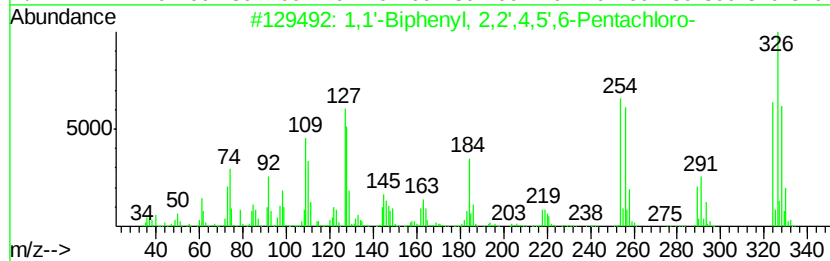
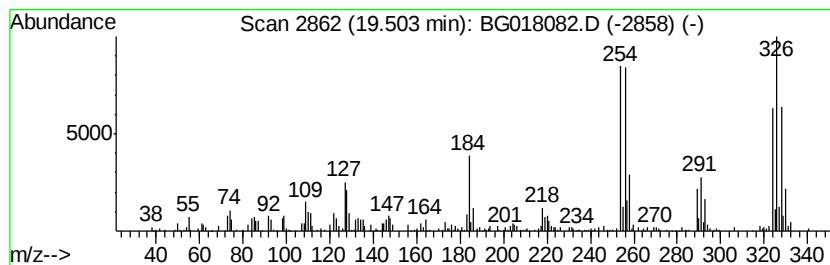
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1,1'-Biphenyl, 2,2',4,5',6-... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.50	2.03 ng/ul	148992	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',4,5',6-Penta...	324	C12H5Cl5	060145-21-3	99
2		1,1'-Biphenyl, 2,2',4,4',6-Penta...	324	C12H5Cl5	039485-83-1	99
3		1,1'-Biphenyl, 2,2',3',4,5-Penta...	324	C12H5Cl5	041464-51-1	99
4		1,1'-Biphenyl, 2,2',4,4',5-penta...	324	C12H5Cl5	038380-01-7	99
5		1,1'-Biphenyl, 2,2',3,5,5'-penta...	324	C12H5Cl5	052663-61-3	98



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
Data File : BG018082.D
Acq On : 24 Jul 2015 1:33
Operator : TP/UM
Sample : G3035-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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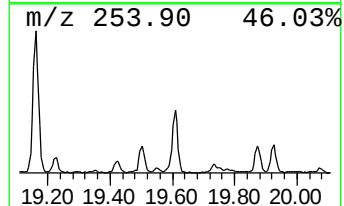
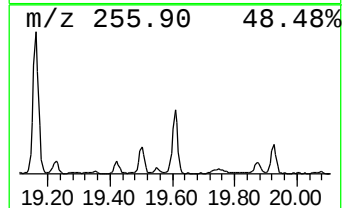
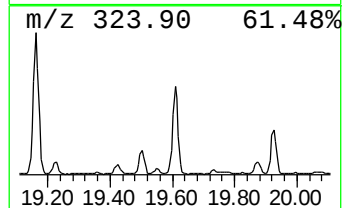
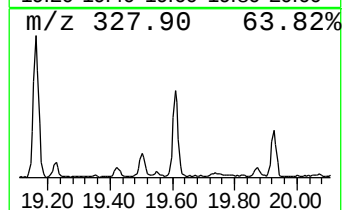
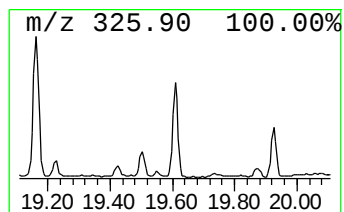
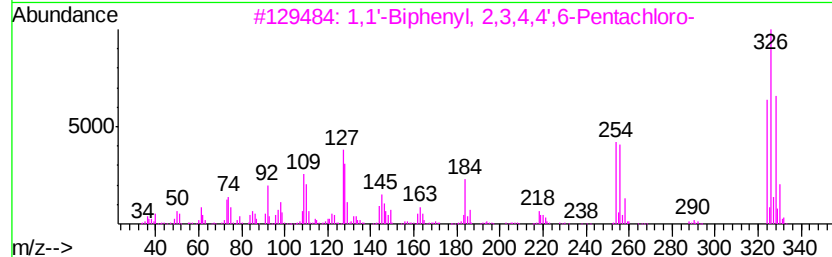
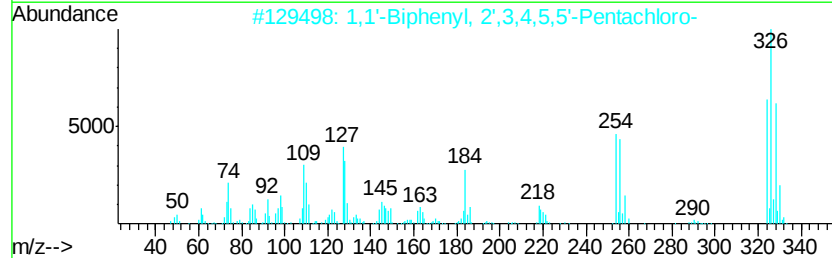
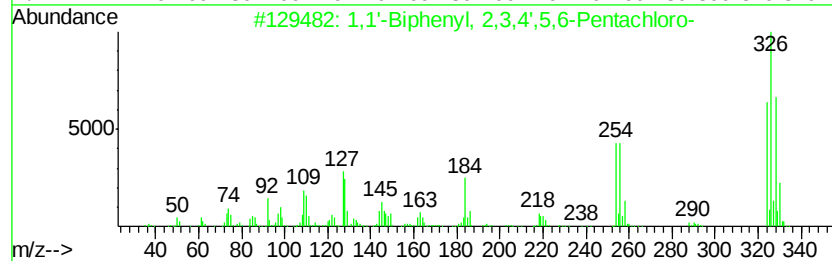
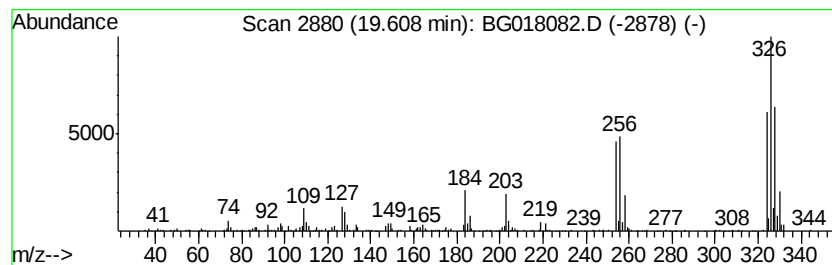
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 1,1'-Biphenyl, 2,3,4',5,6-P... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.61	6.20 ng/ul	455228	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3,4',5,6-Pentac...	324	C12H5Cl5	068194-11-6	98
2		1,1'-Biphenyl, 2',3,4,5,5'-Penta...	324	C12H5Cl5	070424-70-3	97
3		1,1'-Biphenyl, 2,3,4,4',6-Pentac...	324	C12H5Cl5	074472-38-1	96
4		1,1'-Biphenyl, 2,3,3',4,6-Pentac...	324	C12H5Cl5	074472-35-8	96
5		1,1'-Biphenyl, 2,2',4,5,5'-penta...	324	C12H5Cl5	037680-73-2	96



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
Data File : BG018082.D
Acq On : 24 Jul 2015 1:33
Operator : TP/UM
Sample : G3035-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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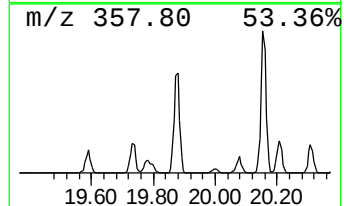
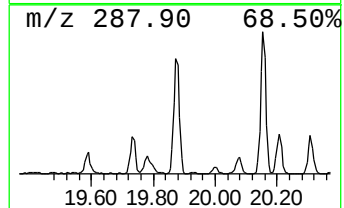
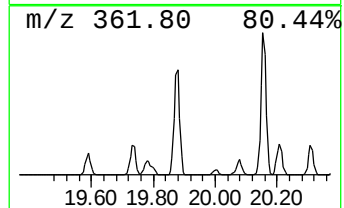
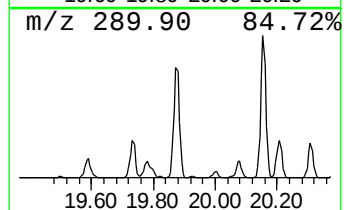
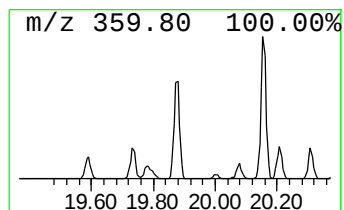
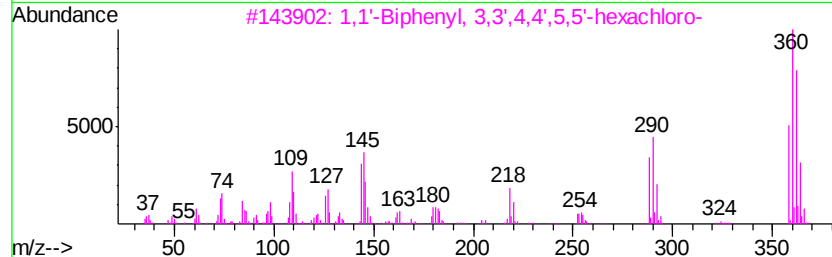
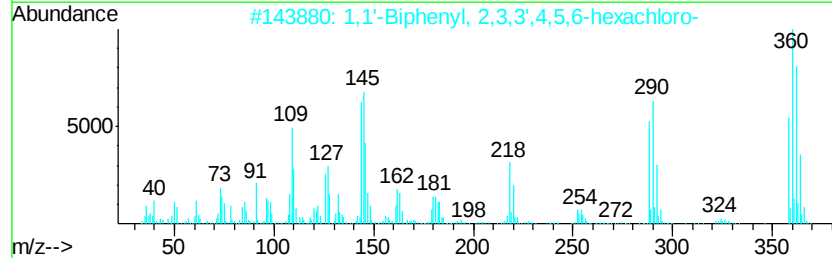
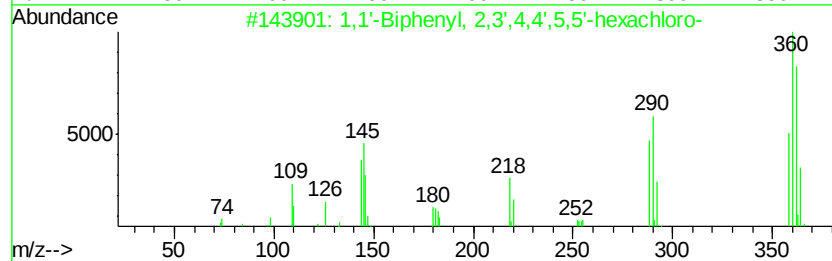
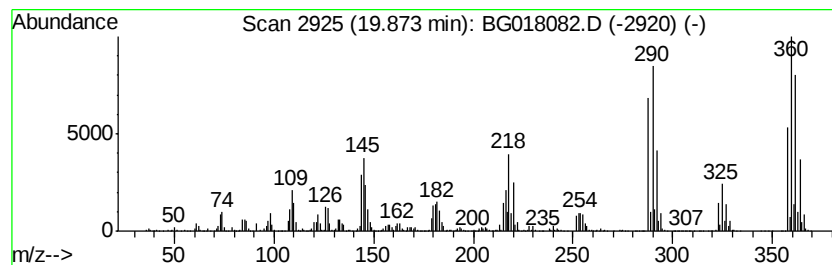
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 1,1'-Biphenyl, 3,3',4,4',5,... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.87	24.94 ng/ul	1832120	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3',4,4',5,5'-he...	358	C12H4Cl6	052663-72-6	99
2		1,1'-Biphenyl, 2,3,3',4,5,6-hexa...	358	C12H4Cl6	041411-62-5	99
3		1,1'-Biphenyl, 3,3',4,4',5,5'-he...	358	C12H4Cl6	032774-16-6	99
4		1,1'-Biphenyl, 2,2',3,4,4',5'-he...	358	C12H4Cl6	035065-28-2	99
5		1,1'-Biphenyl, 2,2',4,4',6,6'-he...	358	C12H4Cl6	033979-03-2	99



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
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Acq On : 24 Jul 2015 1:33
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Sample : G3035-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C01K4

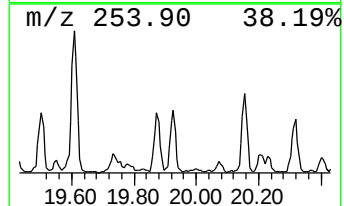
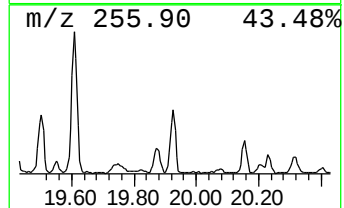
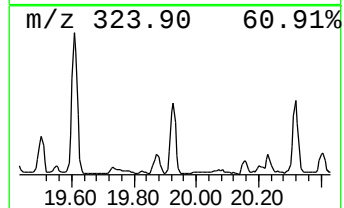
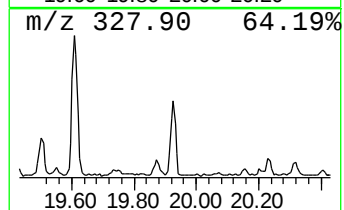
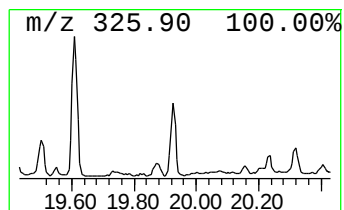
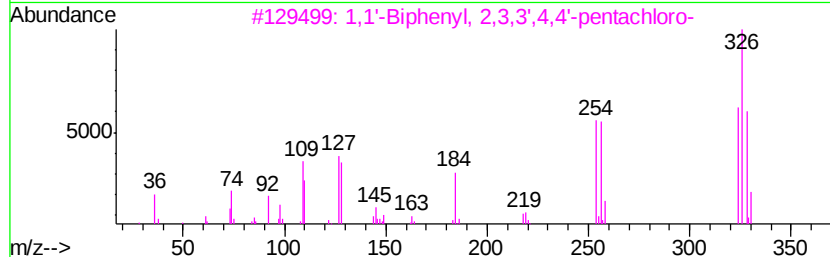
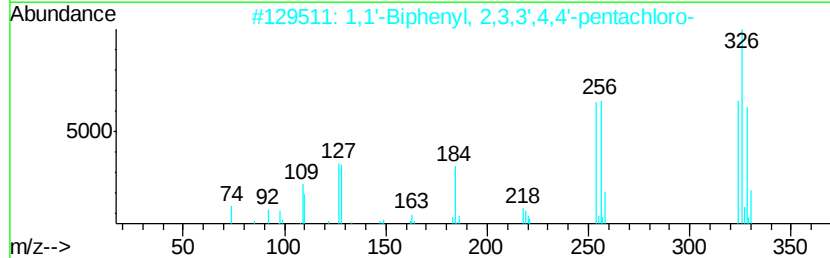
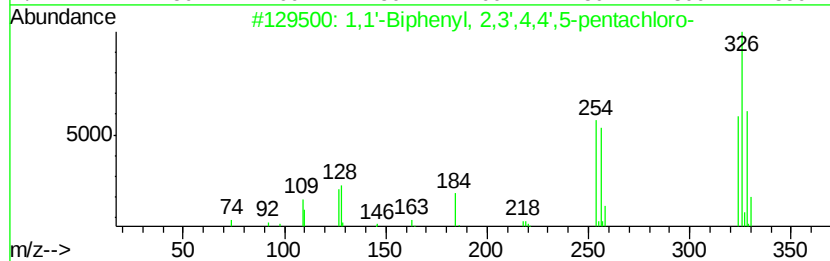
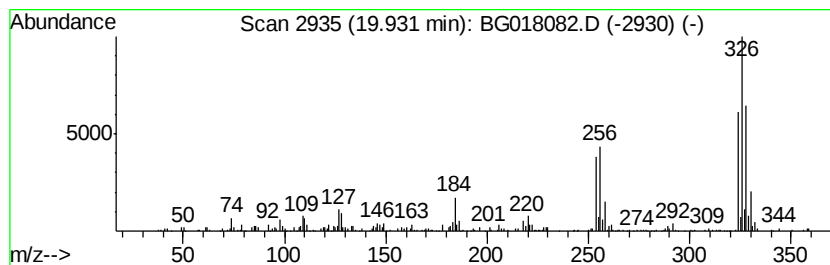
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 1,1'-Biphenyl, 2,3',4,4',5-... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.93	2.98 ng/ul	218673	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	99
2		1,1'-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	99
3		1,1'-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	98
4		1,1'-Biphenyl, 2,2',4,5,5'-penta...	324	C12H5Cl5	037680-73-2	97
5		1,1'-Biphenyl, 2,2',3,5,5'-penta...	324	C12H5Cl5	052663-61-3	96



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
Data File : BG018082.D
Acq On : 24 Jul 2015 1:33
Operator : TP/UM
Sample : G3035-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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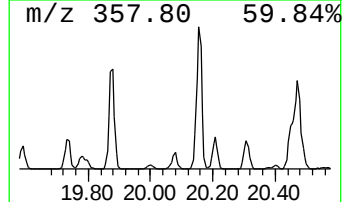
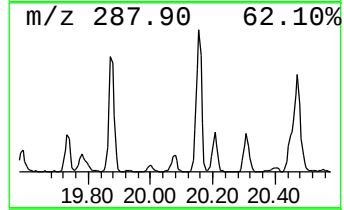
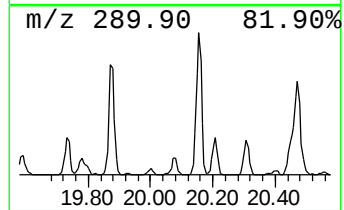
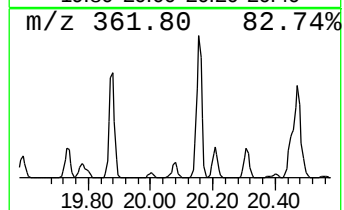
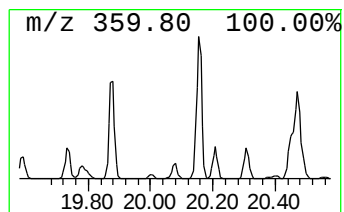
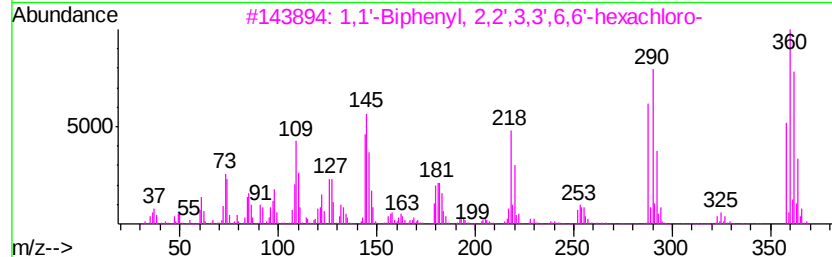
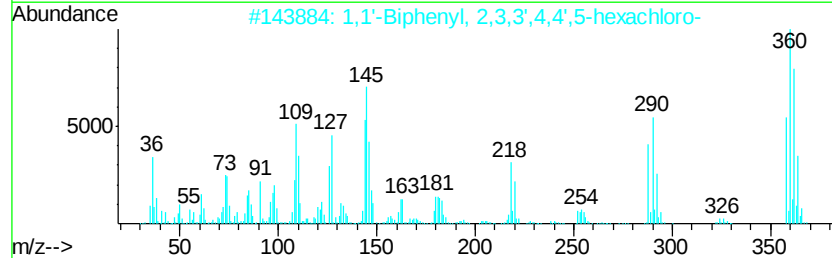
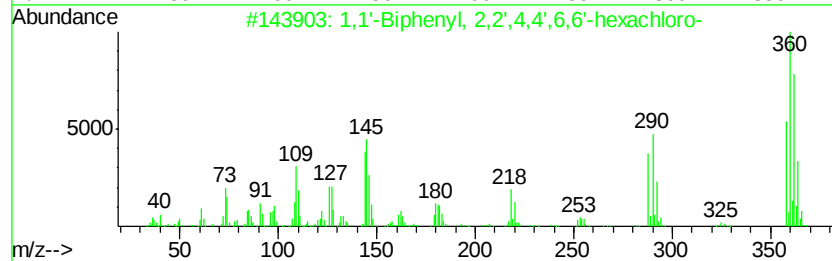
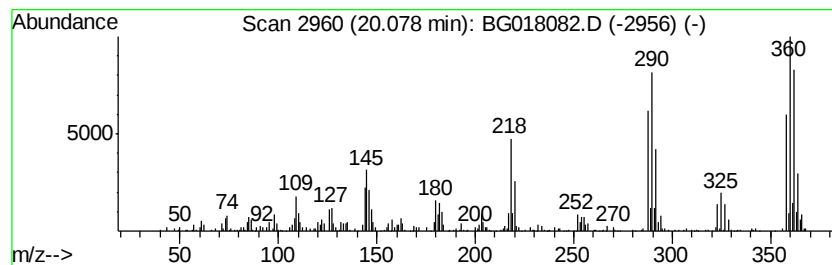
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 1,1'-Biphenyl, 2,2',4,4',6,... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.08	3.54 ng/ul	259755	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',4,4',6,6'-he...	358	C12H4Cl6	033979-03-2	99
2		1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	98
3		1,1'-Biphenyl, 2,2',3,3',6,6'-he...	358	C12H4Cl6	038411-22-2	96
4		1,1'-Biphenyl, 2,2',3,4,4',6-Hex...	358	C12H4Cl6	056030-56-9	95
5		1,1'-Biphenyl, 2,2',4,4',5,5'-he...	358	C12H4Cl6	035065-27-1	95



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
Data File : BG018082.D
Acq On : 24 Jul 2015 1:33
Operator : TP/UM
Sample : G3035-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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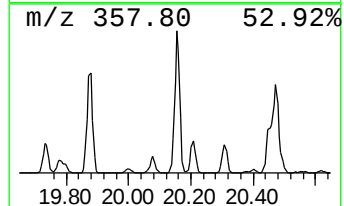
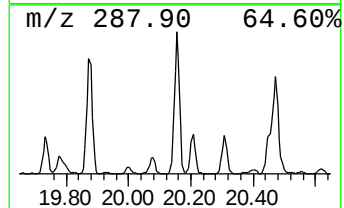
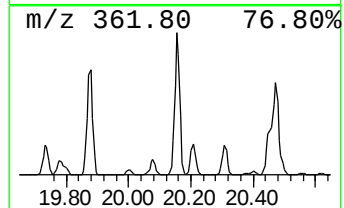
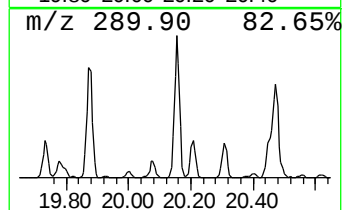
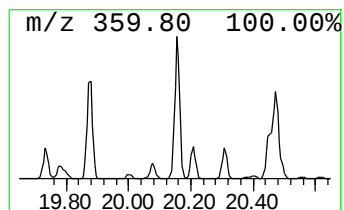
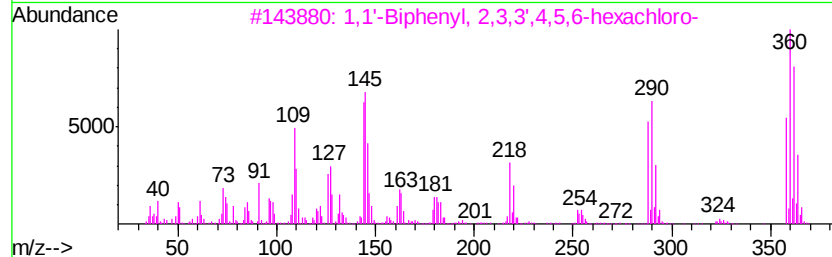
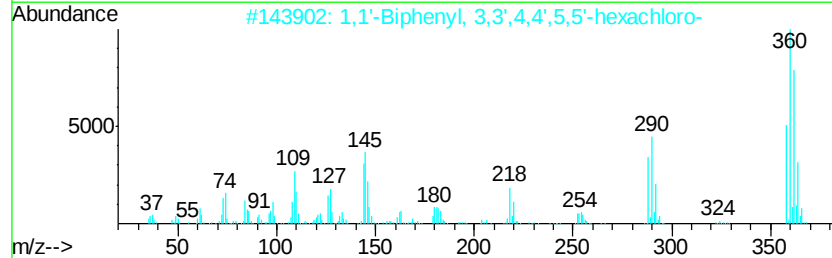
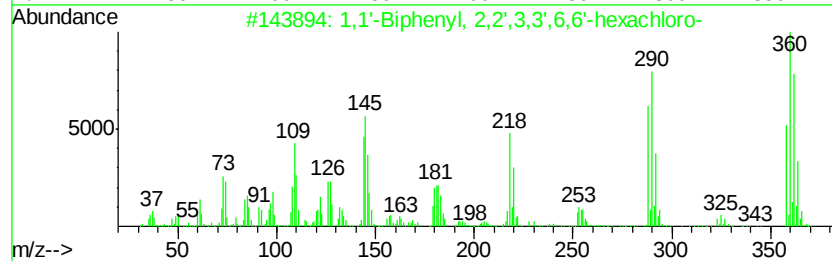
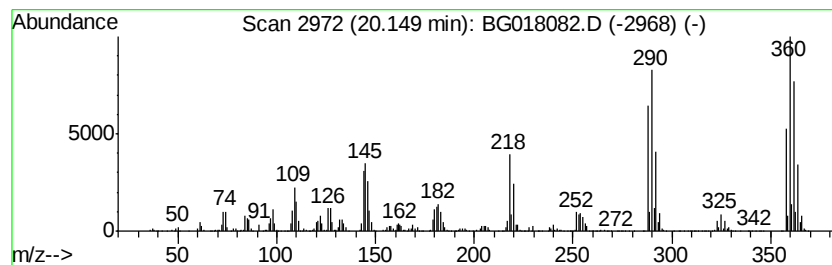
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 1,1'-Biphenyl, 2,2',3,3',6,... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.15	27.61 ng/ul	2027780	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',3,3',6,6'-he...	358	C12H4Cl6	038411-22-2	99
2		1,1'-Biphenyl, 3,3',4,4',5,5'-he...	358	C12H4Cl6	032774-16-6	99
3		1,1'-Biphenyl, 2,3,3',4,5,6-hexa...	358	C12H4Cl6	041411-62-5	99
4		1,1'-Biphenyl, 2,2',4,4',6,6'-he...	358	C12H4Cl6	033979-03-2	99
5		1,1'-Biphenyl, 2,3',4,4',5,5'-he...	358	C12H4Cl6	052663-72-6	99



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
Data File : BG018082.D
Acq On : 24 Jul 2015 1:33
Operator : TP/UM
Sample : G3035-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C01K4

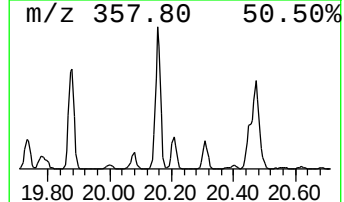
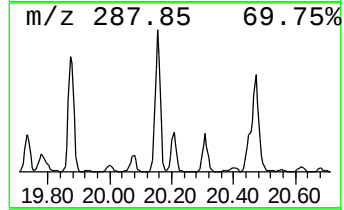
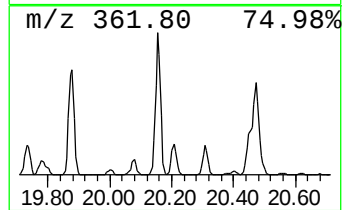
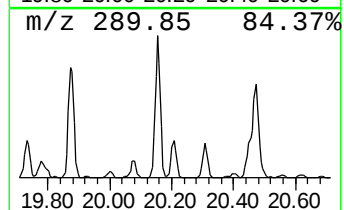
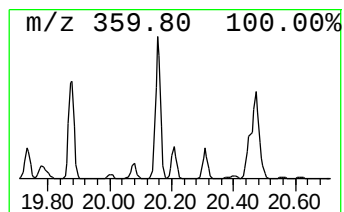
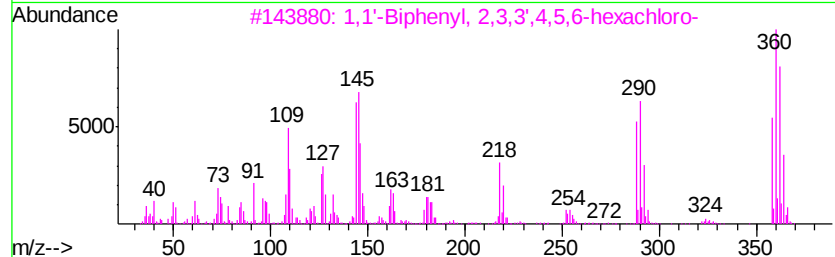
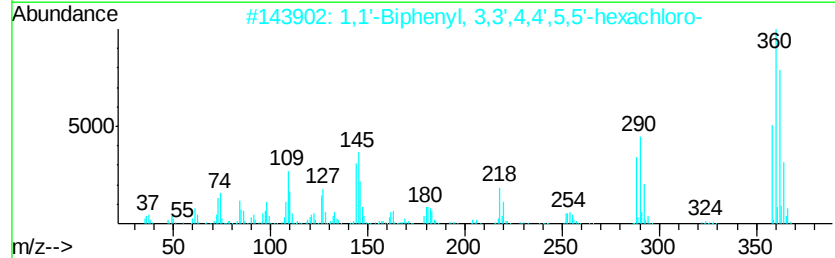
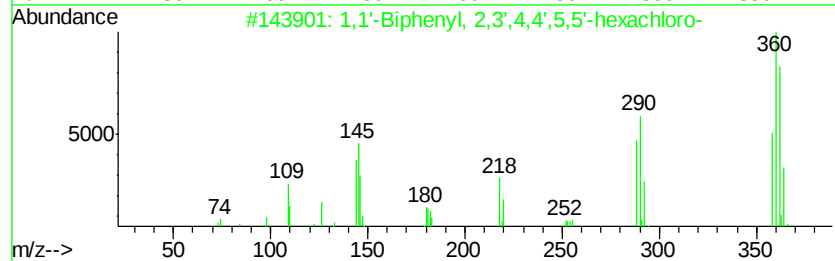
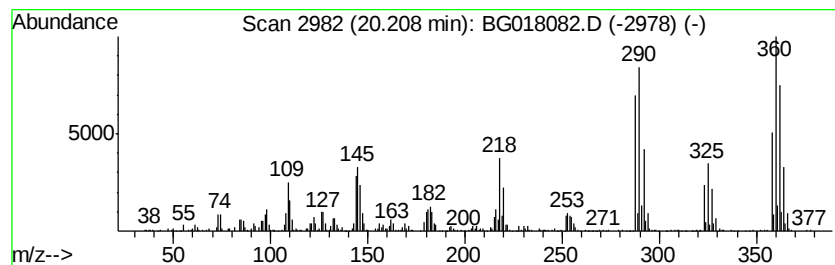
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 1,1'-Biphenyl, 2,3',4,4',5,... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.21	9.23 ng/ul	678224	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3',4,4',5,5'-he...	358	C12H4Cl6	052663-72-6	99
2		1,1'-Biphenyl, 3,3',4,4',5,5'-he...	358	C12H4Cl6	032774-16-6	99
3		1,1'-Biphenyl, 2,3,3',4,5,6-hexa...	358	C12H4Cl6	041411-62-5	99
4		1,1'-Biphenyl, 2,2',3,3',4,6'-he...	358	C12H4Cl6	038380-05-1	99
5		1,1'-Biphenyl, 2,2',3,4,4',6-Hex...	358	C12H4Cl6	056030-56-9	99



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
Data File : BG018082.D
Acq On : 24 Jul 2015 1:33
Operator : TP/UM
Sample : G3035-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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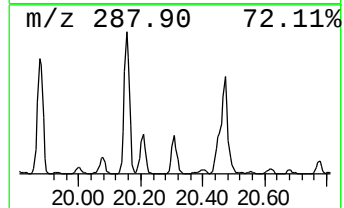
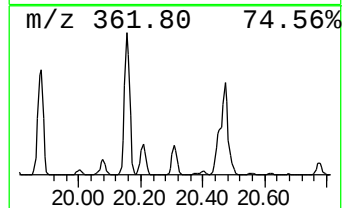
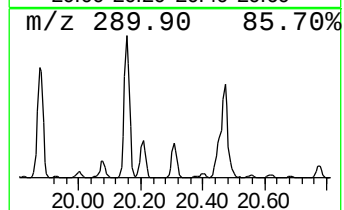
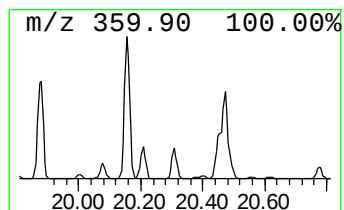
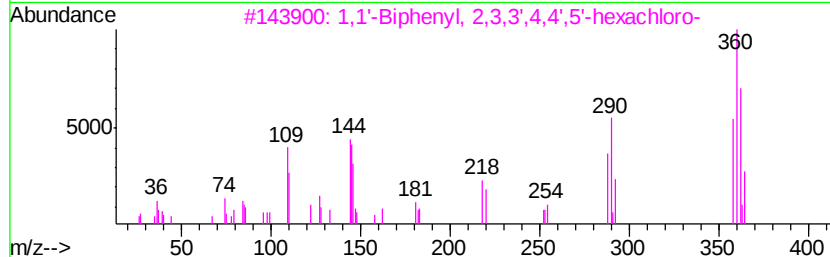
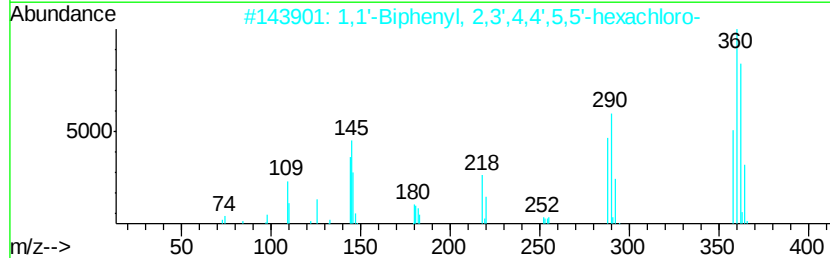
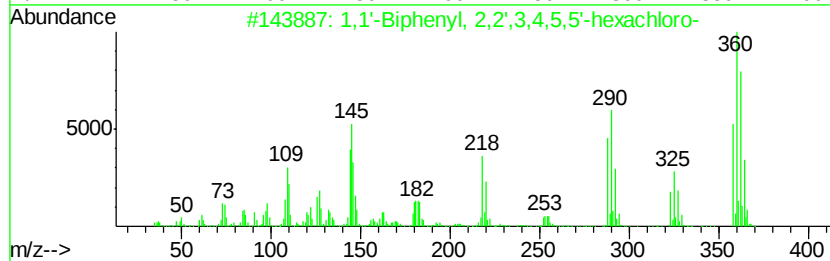
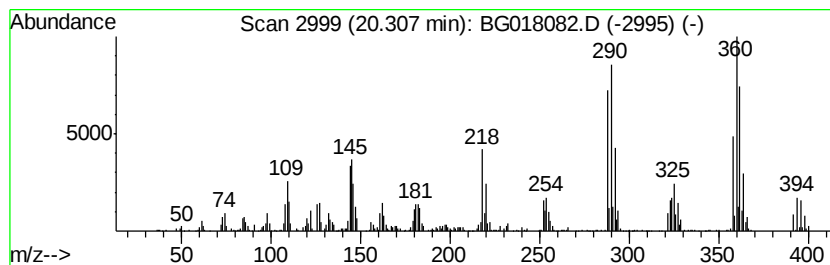
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 1,1'-Biphenyl, 2,2',3,4,5,5... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.31	10.65 ng/ul	782548	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',3,4,5,5'-hex...	358	C12H4Cl6	052712-04-6	99
2		1,1'-Biphenyl, 2,3',4,4',5,5'-he...	358	C12H4Cl6	052663-72-6	99
3		1,1'-Biphenyl, 2,3,3',4,4',5'-he...	358	C12H4Cl6	069782-90-7	99
4		1,1'-Biphenyl, 2,3,3',4,5,6-hexa...	358	C12H4Cl6	041411-62-5	99
5		1,1'-Biphenyl, 3,3',4,4',5,5'-he...	358	C12H4Cl6	032774-16-6	99



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
Data File : BG018082.D
Acq On : 24 Jul 2015 1:33
Operator : TP/UM
Sample : G3035-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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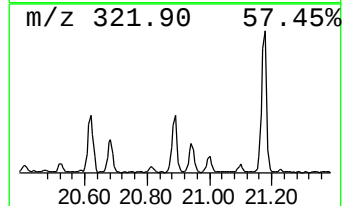
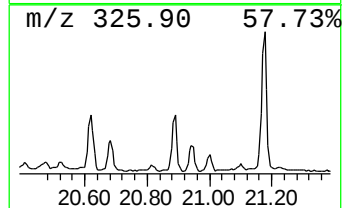
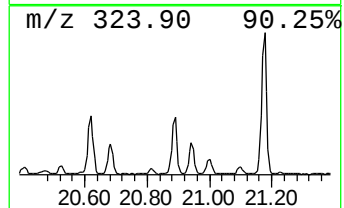
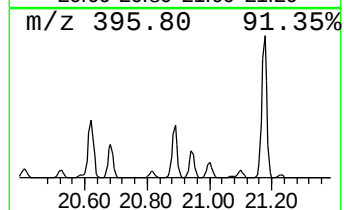
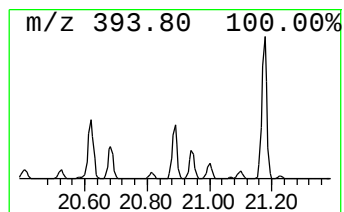
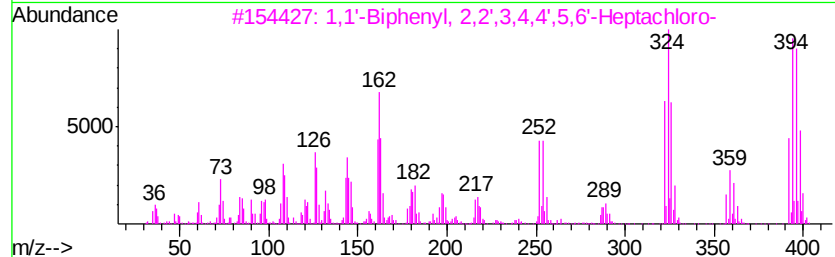
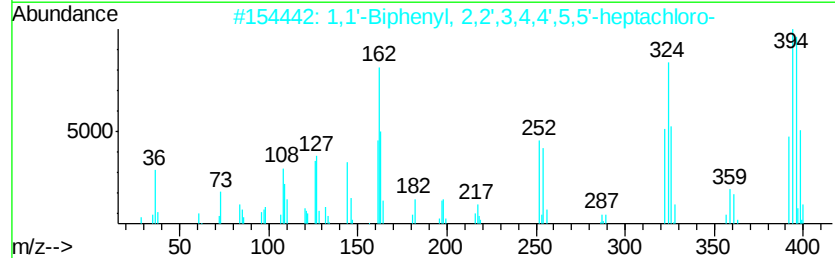
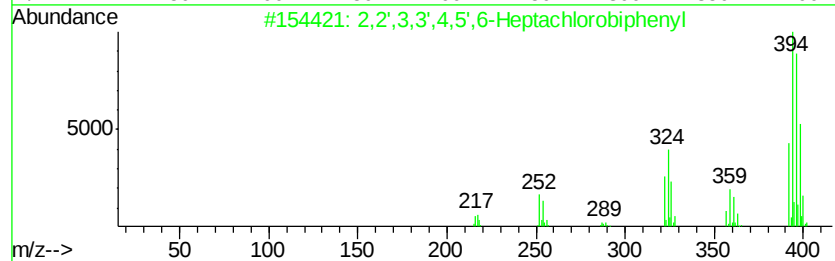
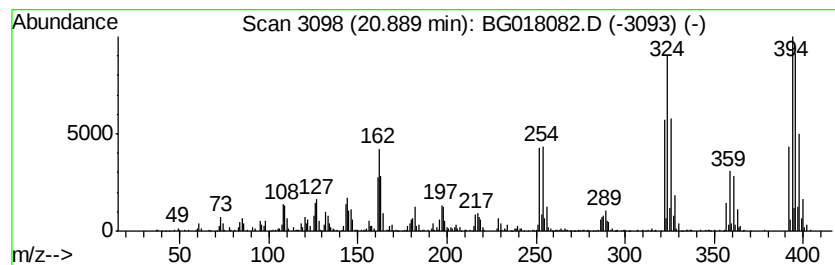
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 2,2',3,3',4,5',6-Heptachlor... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.89	9.44 ng/ul	693347	Chrysene-d12	21.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,3',4,5',6-Heptachlorobiph...	392	C12H3Cl7	040186-70-7	99		
2	1,1'-Biphenyl, 2,2',3,4,4',5,5'-...	392	C12H3Cl7	035065-29-3	99		
3	1,1'-Biphenyl, 2,2',3,4,4',5,6'-...	392	C12H3Cl7	060145-23-5	99		
4	1,1'-Biphenyl, 2,2',3,4,5,5',6-h...	392	C12H3Cl7	052712-05-7	99		
5	1,1'-Biphenyl, 2,3,3',4,5,5',6-h...	392	C12H3Cl7	074472-51-8	99		



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
Data File : BG018082.D
Acq On : 24 Jul 2015 1:33
Operator : TP/UM
Sample : G3035-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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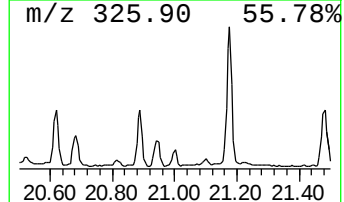
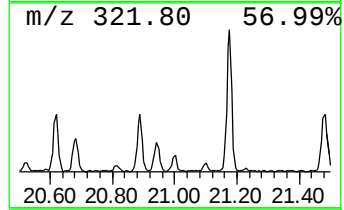
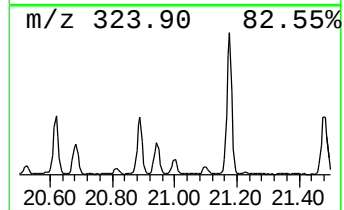
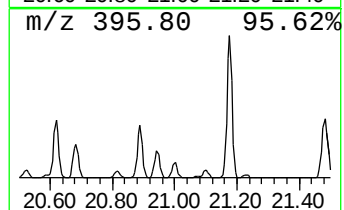
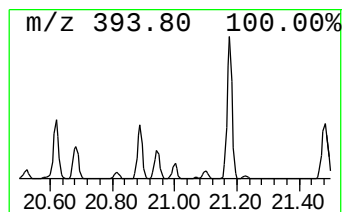
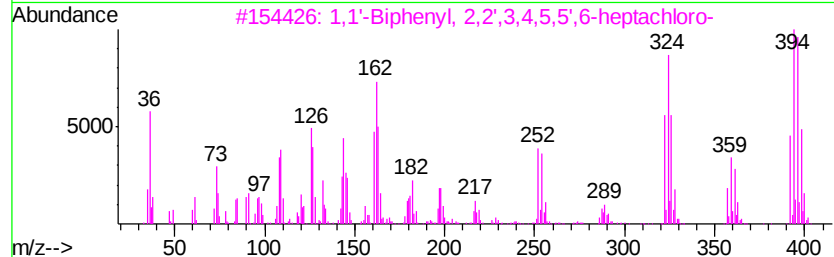
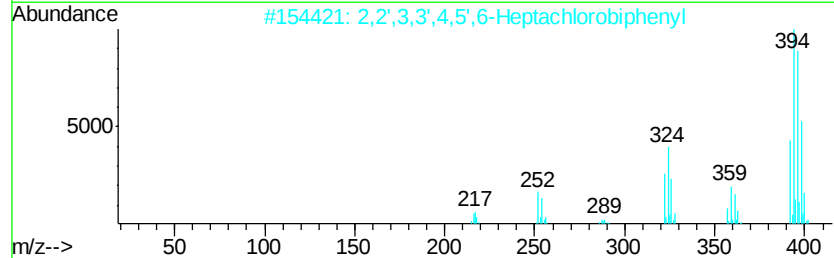
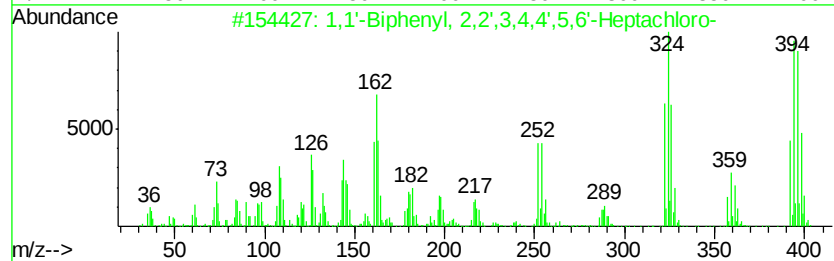
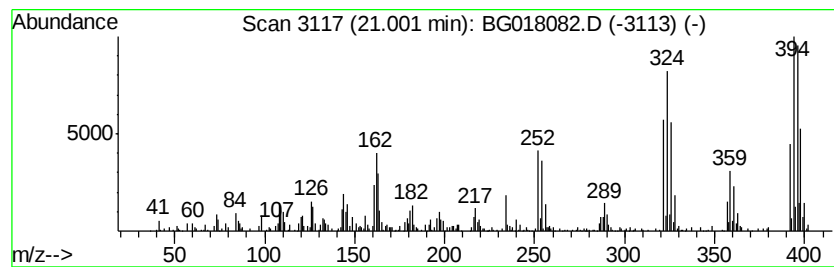
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 1,1'-Biphenyl, 2,2',3,4,4',... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.00	2.74 ng/ul	201101	Chrysene-d12	21.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,4,4',5,6'-...	392	C12H3Cl7	060145-23-5	99		
2	2,2',3,3',4,5',6-Heptachlorobiph...	392	C12H3Cl7	040186-70-7	99		
3	1,1'-Biphenyl, 2,2',3,4,5,5',6-h...	392	C12H3Cl7	052712-05-7	99		
4	1,1'-Biphenyl, 2,2',3,4,4',5',6-...	392	C12H3Cl7	052663-69-1	99		
5	1,1'-Biphenyl, 2,2',3,4,4',5,5',6-...	392	C12H3Cl7	052663-68-0	96		



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
Data File : BG018082.D
Acq On : 24 Jul 2015 1:33
Operator : TP/UM
Sample : G3035-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C01K4

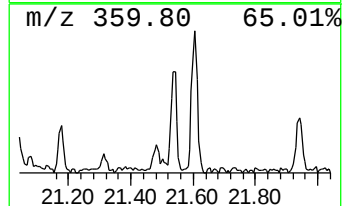
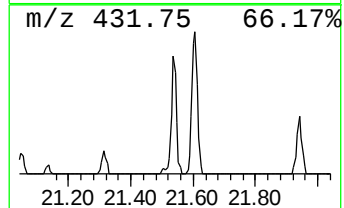
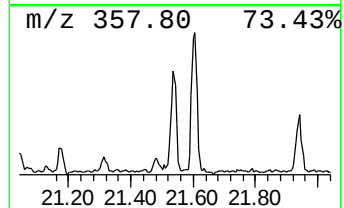
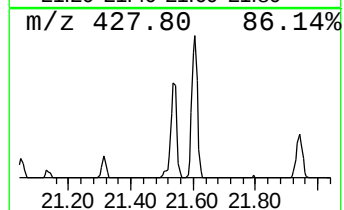
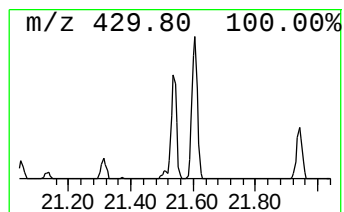
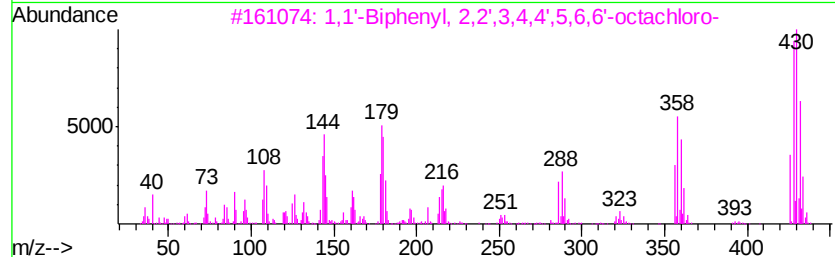
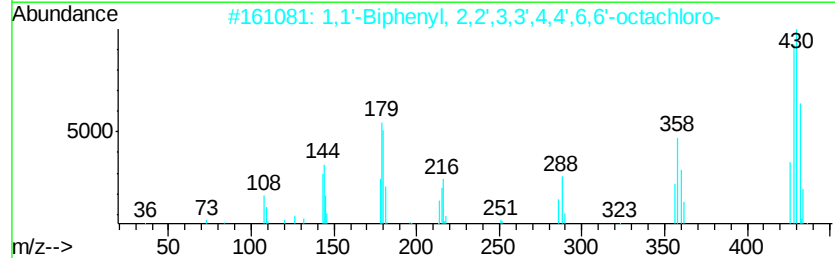
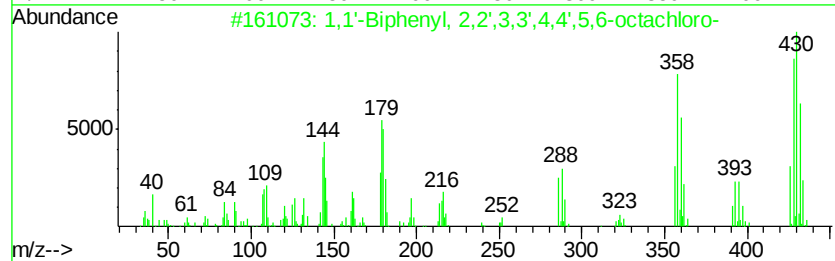
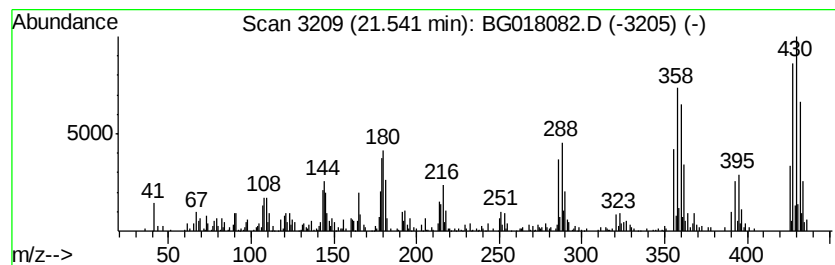
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 1,1'-Biphenyl, 2,2',3,3',4,... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.54	3.63 ng/ul	266670	Chrysene-d12	21.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,3',4,4',5,...	426	C12H2Cl8	052663-78-2	99		
2	1,1'-Biphenyl, 2,2',3,3',4,4',6,...	426	C12H2Cl8	033091-17-7	99		
3	1,1'-Biphenyl, 2,2',3,4,4',5,6,6...	426	C12H2Cl8	074472-52-9	99		
4	1,1'-Biphenyl, 2,2',3,3',5,5',6,...	426	C12H2Cl8	002136-99-4	99		
5	1,1'-Biphenyl, 2,2',3,3',4,4',5,...	426	C12H2Cl8	042740-50-1	99		



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
Data File : BG018082.D
Acq On : 24 Jul 2015 1:33
Operator : TP/UM
Sample : G3035-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C01K4

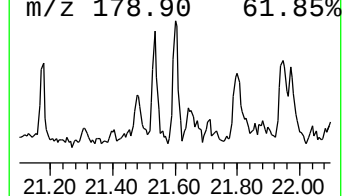
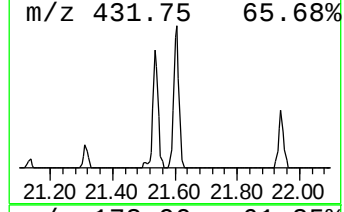
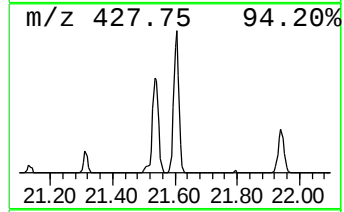
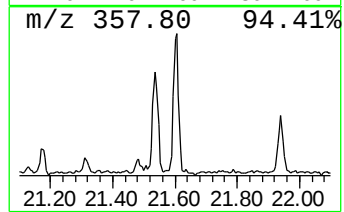
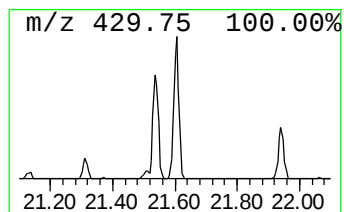
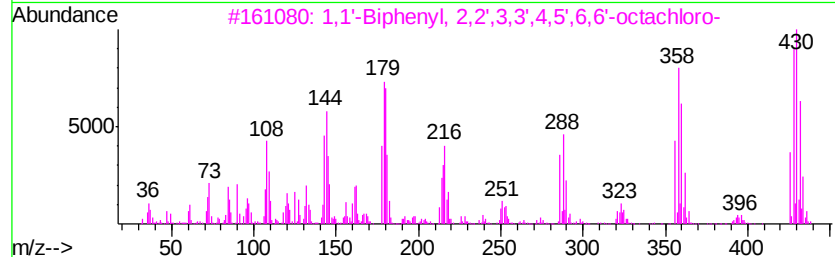
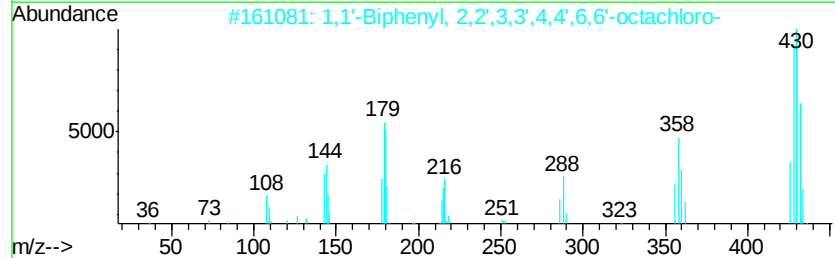
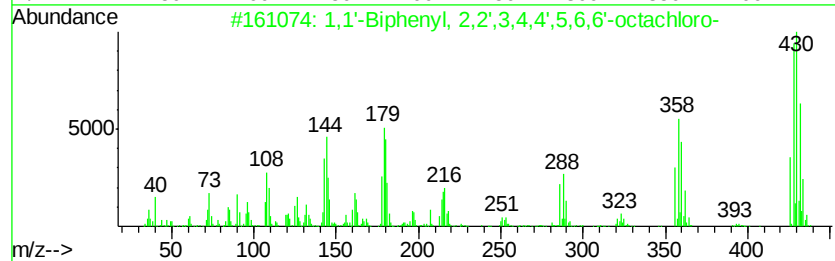
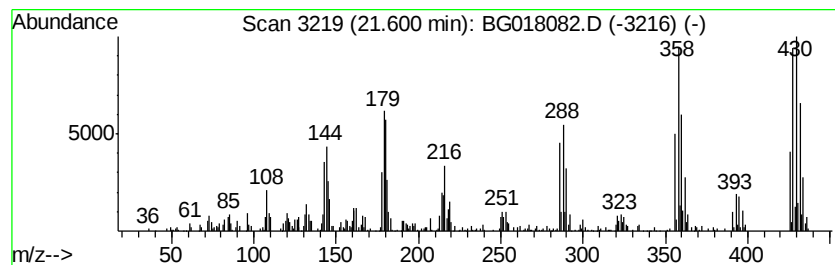
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 1,1'-Biphenyl, 2,2',3,4,4',... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.60	4.12 ng/ul	302474	Chrysene-d12	21.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,4,4',5,6,6...	426	C12H2Cl8	074472-52-9	99		
2	1,1'-Biphenyl, 2,2',3,3',4,4',6,...	426	C12H2Cl8	033091-17-7	99		
3	1,1'-Biphenyl, 2,2',3,3',4,5',6,...	426	C12H2Cl8	040186-71-8	99		
4	1,1'-Biphenyl, 2,2',3,3',5,5',6,...	426	C12H2Cl8	002136-99-4	99		
5	1,1'-Biphenyl, 2,2',3,3',4,4',5,...	426	C12H2Cl8	035694-08-7	99		



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
Data File : BG018082.D
Acq On : 24 Jul 2015 1:33
Operator : TP/UM
Sample : G3035-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C01K4

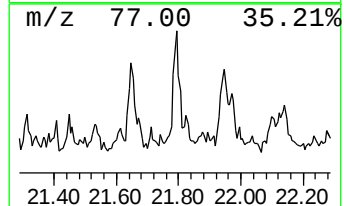
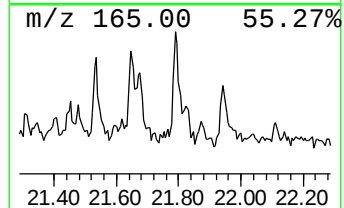
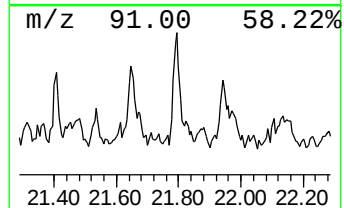
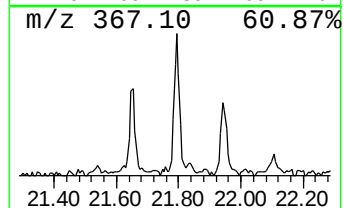
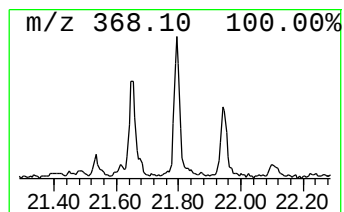
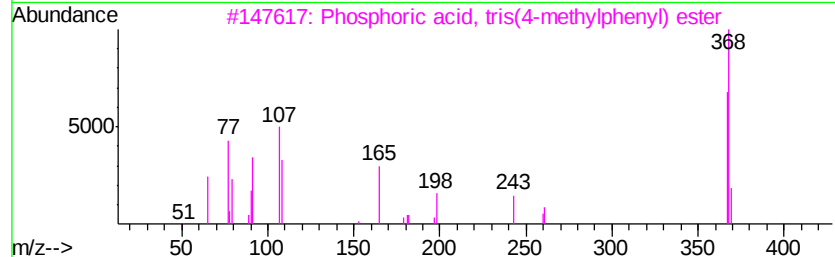
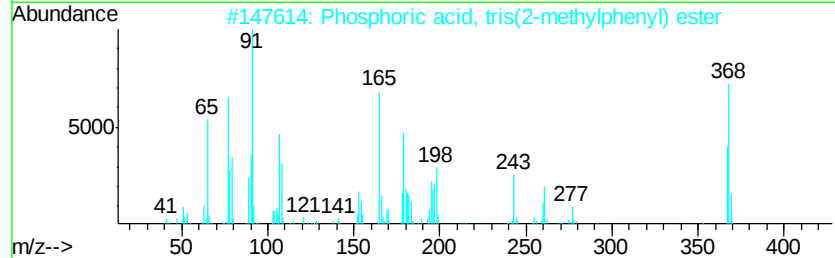
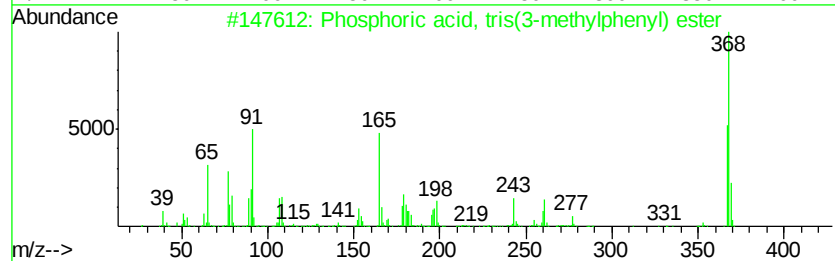
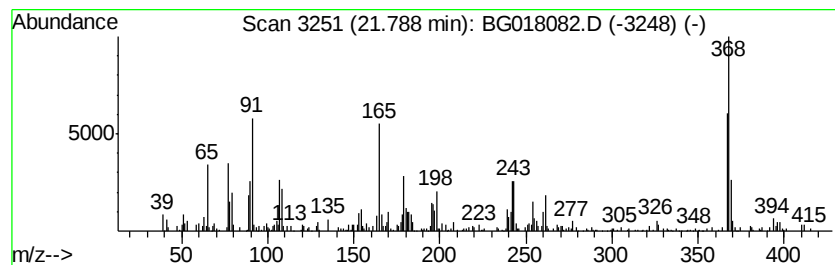
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Quant Title : SVOA CALIBRATION

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

Peak Number 29 Phosphoric acid, tris(3-met... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.79	2.20 ng/ul	161427	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phosphoric acid, tris(3-methylph...	368	C21H21O4P	000563-04-2	98
2		Phosphoric acid, tris(2-methylph...	368	C21H21O4P	000078-30-8	95
3		Phosphoric acid, tris(4-methylph...	368	C21H21O4P	000078-32-0	89
4		Phosphoric acid, tris(2-methylph...	368	C21H21O4P	000078-30-8	83
5		Phosphoric acid, tris(3-methylph...	368	C21H21O4P	000563-04-2	76



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
Data File : BG018082.D
Acq On : 24 Jul 2015 1:33
Operator : TP/UM
Sample : G3035-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C01K4

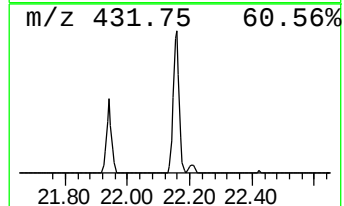
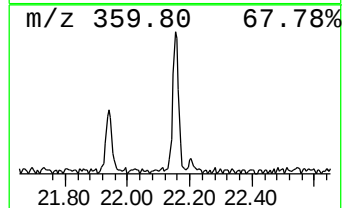
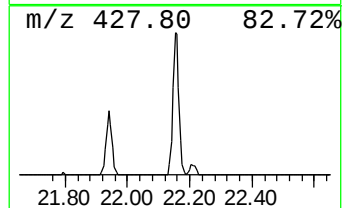
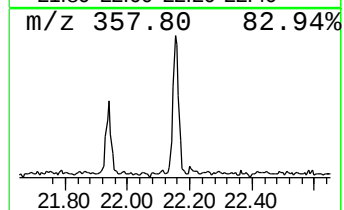
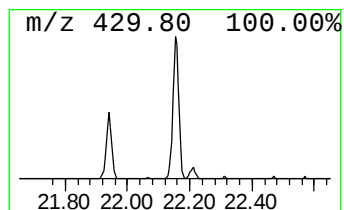
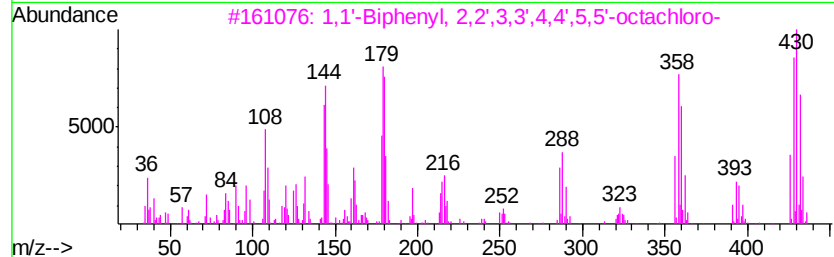
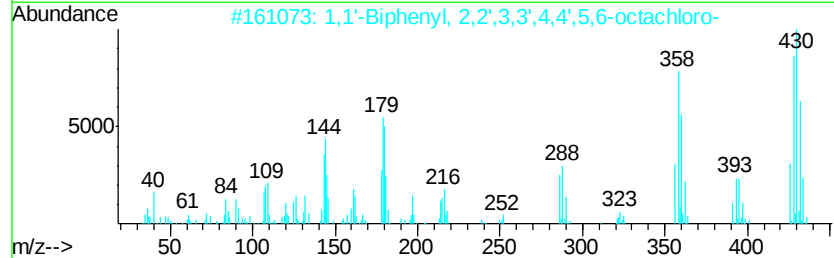
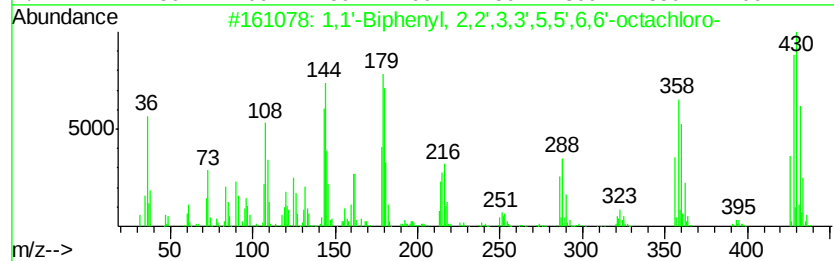
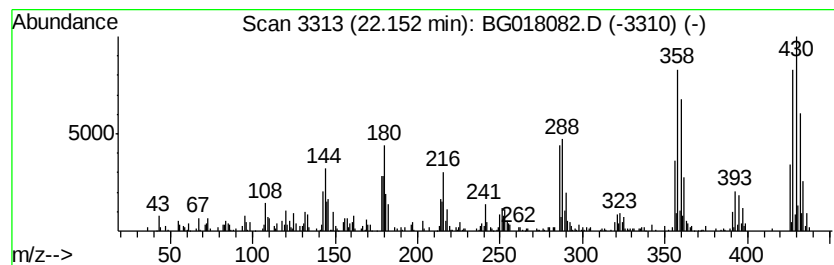
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Quant Title : SVOA CALIBRATION

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

Peak Number 31 1,1'-Biphenyl, 2,2',3,3',5,... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.15	2.33 ng/ul	170856	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',3,3',5,5',6,...	426	C12H2Cl8	002136-99-4	96
2		1,1'-Biphenyl, 2,2',3,3',4,4',5,...	426	C12H2Cl8	052663-78-2	93
3		1,1'-Biphenyl, 2,2',3,3',4,4',5,...	426	C12H2Cl8	035694-08-7	91
4		1,1'-Biphenyl, 2,2',3,3',4,5,5',...	426	C12H2Cl8	068194-17-2	91
5		1,1'-Biphenyl, 2,2',3,3',4,4',5,...	426	C12H2Cl8	042740-50-1	91



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
Data File : BG018082.D
Acq On : 24 Jul 2015 1:33
Operator : TP/UM
Sample : G3035-16
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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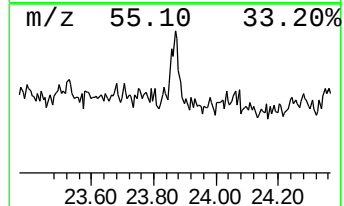
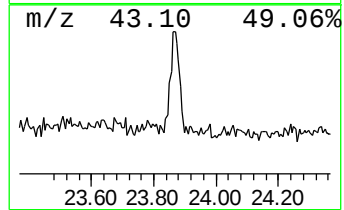
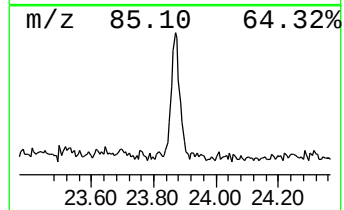
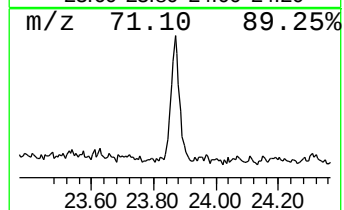
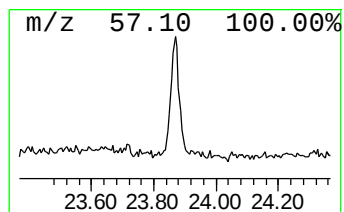
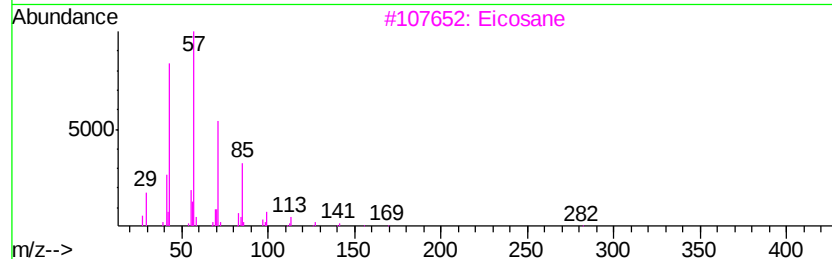
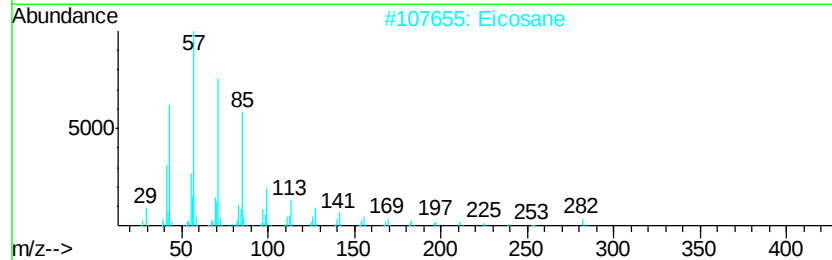
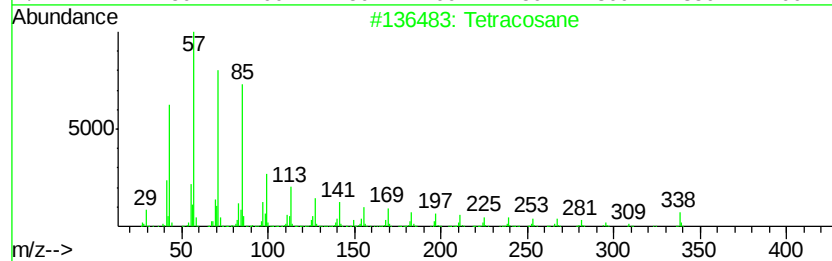
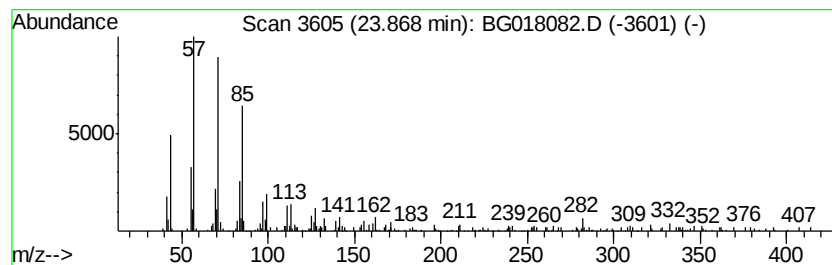
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG072315.M
Quant Title : SVOA CALIBRATION

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

Peak Number 33 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.87	3.32 ng/ul	165767	Perylene-d12	23.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	94
2			Eicosane	282	C20H42	000112-95-8	93
3			Eicosane	282	C20H42	000112-95-8	93
4			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
5			Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072315\
 Data File : BG018082.D
 Acq On : 24 Jul 2015 1:33
 Operator : TP/UM
 Sample : G3035-16
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C01K4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG072315.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
Phenanthrene, 1-m...	17.86	2.7	ng/ul	123115	4	16.93	902295	20.0
1,1'-Biphenyl, 2,...	18.02	5.3	ng/ul	240762	4	16.93	902295	20.0
1,1'-Biphenyl, 2,...	18.30	2.7	ng/ul	121249	4	16.93	902295	20.0
1,1'-Biphenyl, 2,...	18.87	17.2	ng/ul	777491	4	16.93	902295	20.0
1,1'-Biphenyl, 2,...	19.09	2.1	ng/ul	153745	5	21.14	1469040	20.0
1,1'-Biphenyl, 2,...	19.16	12.3	ng/ul	905034	5	21.14	1469040	20.0
1,1'-Biphenyl, 2,...	19.50	2.0	ng/ul	148992	5	21.14	1469040	20.0
1,1'-Biphenyl, 2,...	19.61	6.2	ng/ul	455228	5	21.14	1469040	20.0
1,1'-Biphenyl, 3,...	19.87	24.9	ng/ul	1832120	5	21.14	1469040	20.0
1,1'-Biphenyl, 2,...	19.93	3.0	ng/ul	218673	5	21.14	1469040	20.0
1,1'-Biphenyl, 2,...	20.08	3.5	ng/ul	259755	5	21.14	1469040	20.0
1,1'-Biphenyl, 2,...	20.15	27.6	ng/ul	2027780	5	21.14	1469040	20.0
1,1'-Biphenyl, 2,...	20.21	9.2	ng/ul	678224	5	21.14	1469040	20.0
1,1'-Biphenyl, 2,...	20.31	10.7	ng/ul	782548	5	21.14	1469040	20.0
2,2',3,3',4,5',6-...	20.89	9.4	ng/ul	693347	5	21.14	1469040	20.0
1,1'-Biphenyl, 2,...	21.00	2.7	ng/ul	201101	5	21.14	1469040	20.0
1,1'-Biphenyl, 2,...	21.54	3.6	ng/ul	266670	5	21.14	1469040	20.0
1,1'-Biphenyl, 2,...	21.60	4.1	ng/ul	302474	5	21.14	1469040	20.0
Phosphoric acid, ...	21.79	2.2	ng/ul	161427	5	21.14	1469040	20.0
1,1'-Biphenyl, 2,...	22.15	2.3	ng/ul	170856	5	21.14	1469040	20.0
(DEL) Alkane: Str...	23.87	3.3	ng/ul	165767	6	23.33	999356	20.0