

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110620\
 Data File : BG047245.D
 Acq On : 6 Nov 2020 15:49
 Operator : CG/JU
 Sample : L4534-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0BP1

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:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.394	7	9	12	rVB	1873	1898	0.17%	0.025%
2	3.470	20	22	25	rBV2	1378	1374	0.12%	0.018%
3	3.647	48	52	53	rBV3	3529	3557	0.32%	0.046%
4	3.670	53	56	62	rVB2	14888	19121	1.72%	0.249%
5	3.852	86	87	90	rBV	1351	1437	0.13%	0.019%
6	3.888	92	93	99	rVB	1347	1665	0.15%	0.022%
7	4.064	120	123	125	rVB	1923	1426	0.13%	0.019%
8	4.152	133	138	143	rBV2	5300	7310	0.66%	0.095%
9	4.199	143	146	149	rBV	1497	1601	0.14%	0.021%
10	4.352	168	172	174	rVB	1076	1439	0.13%	0.019%
11	4.604	212	215	219	rBV3	1480	2414	0.22%	0.031%
12	4.710	229	233	237	rVB3	1801	2343	0.21%	0.031%
13	4.781	241	245	247	rBV2	1426	1801	0.16%	0.023%
14	4.810	249	250	254	rVB	1846	1387	0.12%	0.018%
15	4.910	265	267	268	rBV	1639	1440	0.13%	0.019%
16	5.016	280	285	287	rVB3	2755	3677	0.33%	0.048%
17	5.133	302	305	307	rBV	1270	1493	0.13%	0.019%
18	5.445	355	358	359	rVB	1470	1437	0.13%	0.019%
19	5.727	404	406	408	rVB2	1785	1423	0.13%	0.019%
20	5.962	444	446	450	rVB2	1290	1684	0.15%	0.022%
21	6.091	463	468	471	rVB2	1344	2175	0.20%	0.028%
22	6.320	505	507	510	rVB	1829	1692	0.15%	0.022%
23	6.485	533	535	539	rBV	1452	1710	0.15%	0.022%
24	6.708	570	573	575	rVB	1111	1361	0.12%	0.018%
25	6.996	619	622	624	rBV	1126	1607	0.14%	0.021%
26	7.389	685	689	692	rBV3	1961	3273	0.29%	0.043%
27	7.683	736	739	741	rVB	1222	1378	0.12%	0.018%
28	8.000	786	793	800	rBV	251808	429014	38.59%	5.589%
29	8.065	803	804	809	rVB2	1267	1440	0.13%	0.019%
30	8.136	814	816	821	rVB3	3320	3338	0.30%	0.043%
31	8.294	840	843	845	rBV	1393	1406	0.13%	0.018%
32	8.541	882	885	887	rBV	1641	1482	0.13%	0.019%
33	10.380	1195	1198	1200	rVB	1812	1502	0.14%	0.020%
34	10.815	1263	1272	1280	rBV	308449	560729	50.43%	7.305%

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35	10.874	1280	1282	1285	rVB2	2233	2390	0.21%	0.031%
36	11.226	1338	1342	1343	rBV	1811	1502	0.14%	0.020%
37	11.514	1387	1391	1393	rBV2	1709	1471	0.13%	0.019%
38	12.483	1552	1556	1560	rVB	1266	2332	0.21%	0.030%
39	12.701	1589	1593	1595	rBV2	1414	1634	0.15%	0.021%
40	12.730	1595	1598	1601	rVB	1239	1615	0.15%	0.021%
41	13.970	1806	1809	1812	rBV3	1157	1883	0.17%	0.025%
42	14.640	1916	1923	1938	rBV	511725	767922	69.07%	10.004%
43	15.033	1988	1990	1992	rBV	1148	1398	0.13%	0.018%
44	15.433	2054	2058	2059	rBV2	1015	1536	0.14%	0.020%
45	15.562	2077	2080	2082	rVB2	1872	1602	0.14%	0.021%
46	15.592	2082	2085	2089	rVB2	1216	1998	0.18%	0.026%
47	15.691	2098	2102	2105	rVB2	2199	2117	0.19%	0.028%
48	15.727	2105	2108	2111	rBV	1179	1591	0.14%	0.021%
49	15.885	2133	2135	2137	rVB2	1539	1496	0.13%	0.019%
50	16.144	2174	2179	2182	rBV	1514	2887	0.26%	0.038%
51	16.250	2194	2197	2199	rBV	1176	1384	0.12%	0.018%
52	16.344	2211	2213	2215	rBV2	2143	1699	0.15%	0.022%
53	16.373	2215	2218	2223	rVB3	1825	2488	0.22%	0.032%
54	17.072	2332	2337	2339	rBV	1928	2419	0.22%	0.032%
55	17.178	2353	2355	2357	rVB2	2268	1461	0.13%	0.019%
56	17.384	2382	2390	2397	rBV	589693	926010	83.29%	12.063%
57	17.483	2406	2407	2414	rVB2	4939	5383	0.48%	0.070%
58	17.648	2433	2435	2438	rVB2	1942	2070	0.19%	0.027%
59	17.865	2469	2472	2473	rBV2	1439	1429	0.13%	0.019%
60	17.977	2488	2491	2492	rBV2	5938	5750	0.52%	0.075%
61	18.053	2500	2504	2507	rBV2	1850	2576	0.23%	0.034%
62	18.100	2508	2512	2516	rVV3	3198	3606	0.32%	0.047%
63	18.271	2538	2541	2545	rBV6	4188	6599	0.59%	0.086%
64	18.377	2557	2559	2562	rBV	3555	4535	0.41%	0.059%
65	18.453	2564	2572	2577	rVV2	95164	134429	12.09%	1.751%
66	18.512	2578	2582	2586	rVV4	26208	35465	3.19%	0.462%
67	18.559	2586	2590	2593	rVB4	7640	9393	0.84%	0.122%
68	18.676	2609	2610	2612	rBV	2515	1454	0.13%	0.019%
69	18.735	2616	2620	2624	rBV2	38826	48541	4.37%	0.632%
70	18.829	2632	2636	2639	rBV3	3507	4772	0.43%	0.062%
71	18.870	2639	2643	2646	rBV3	4792	6173	0.56%	0.080%

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72	18.911	2646	2650	2653	rVB4	14175	19354	1.74%	0.252%
73	19.011	2665	2667	2668	rBV	2730	1664	0.15%	0.022%
74	19.058	2673	2675	2677	rBV2	3275	2436	0.22%	0.032%
75	19.211	2699	2701	2706	rBV3	15783	20609	1.85%	0.268%
76	19.270	2706	2711	2712	rVV2	47008	63404	5.70%	0.826%
77	19.293	2712	2715	2722	rVB3	135400	191177	17.19%	2.490%
78	19.375	2726	2729	2735	rVB6	20354	26260	2.36%	0.342%
79	19.434	2735	2739	2744	rBV	36299	52143	4.69%	0.679%
80	19.499	2746	2750	2756	rVV8	28926	51335	4.62%	0.669%
81	19.575	2758	2763	2769	rVB	220133	287425	25.85%	3.744%
82	19.640	2769	2774	2778	rBV2	81088	103691	9.33%	1.351%
83	19.763	2793	2795	2798	rBV4	13001	15296	1.38%	0.199%
84	19.798	2798	2801	2803	rVV	9725	11913	1.07%	0.155%
85	19.834	2803	2807	2813	rVV2	58538	71928	6.47%	0.937%
86	19.910	2815	2820	2825	rVB3	101912	138466	12.45%	1.804%
87	19.963	2825	2829	2832	rBV3	36392	50685	4.56%	0.660%
88	20.016	2832	2838	2847	rVB	214443	295156	26.55%	3.845%
89	20.139	2856	2859	2860	rBV2	19916	17602	1.58%	0.229%
90	20.280	2878	2883	2887	rBV4	88296	122871	11.05%	1.601%
91	20.327	2887	2891	2897	rVB	183596	226401	20.36%	2.949%
92	20.556	2925	2930	2934	rBV	106656	134382	12.09%	1.751%
93	20.609	2935	2939	2941	rVV3	50216	69892	6.29%	0.910%
94	20.633	2941	2943	2949	rVV	76624	89457	8.05%	1.165%
95	20.703	2953	2955	2961	rVB5	24099	31329	2.82%	0.408%
96	20.850	2975	2980	2981	rBV3	55855	72831	6.55%	0.949%
97	20.874	2981	2984	2992	rVB3	132430	186528	16.78%	2.430%
98	21.203	3037	3040	3045	rVB6	41608	52085	4.68%	0.679%
99	21.643	3109	3115	3129	rVB2	654483	1111845	100.00%	14.484%
100	24.840	3650	3659	3673	rVB2	387557	1083117	97.42%	14.110%

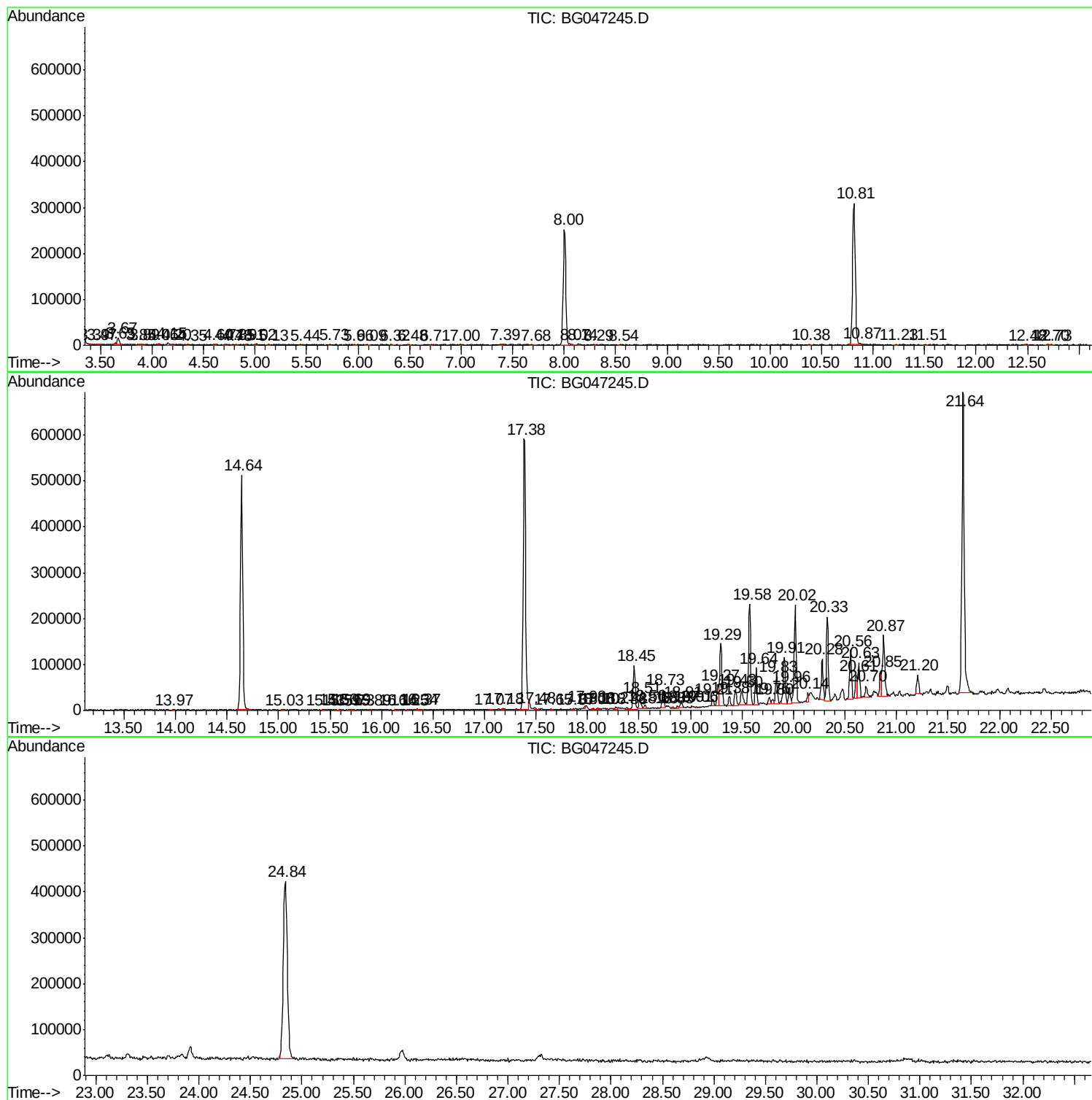
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

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



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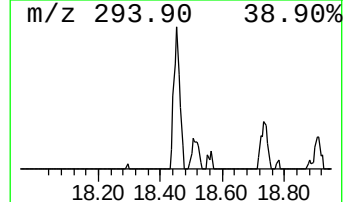
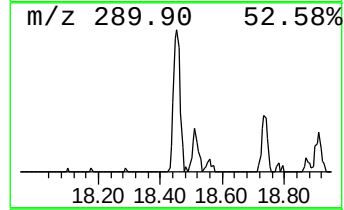
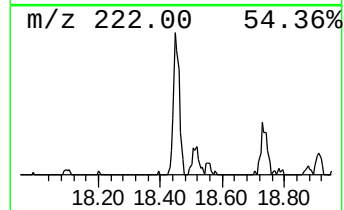
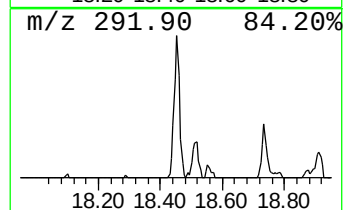
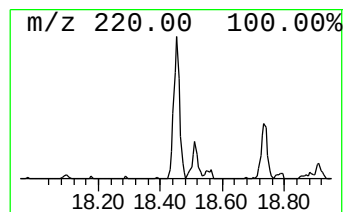
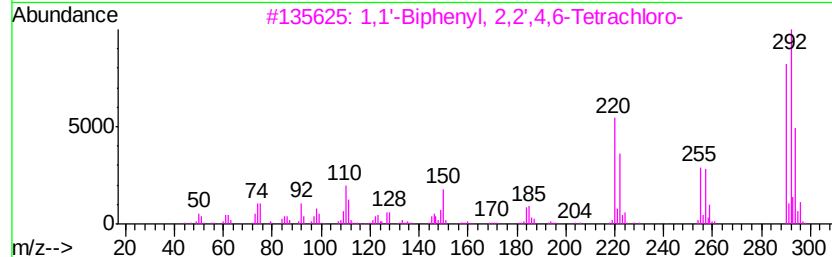
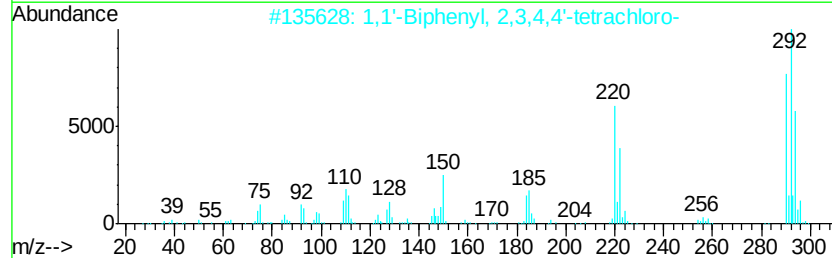
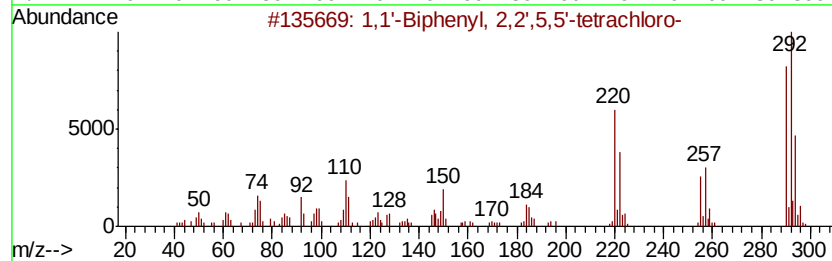
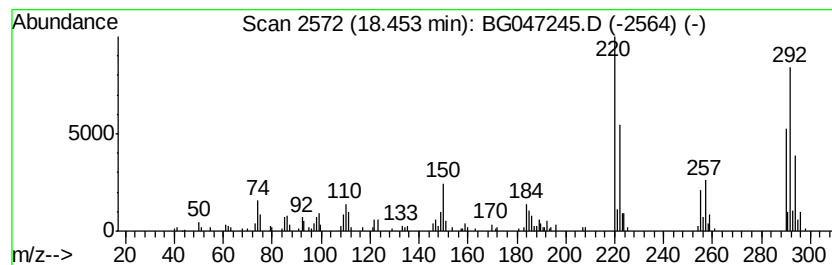
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.45	2.90 ng/ul	134429	Phenanthrene-d10	17.38	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	99
2	1,1'-Biphenyl, 2,3,4,4'-tetrachl...	290	C12H6Cl4	033025-41-1	99
3	1,1'-Biphenyl, 2,2',4,6-Tetrachl...	290	C12H6Cl4	062796-65-0	99
4	1,1'-Biphenyl, 3,3',5,5'-tetrach...	290	C12H6Cl4	033284-52-5	99
5	1,1'-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	99



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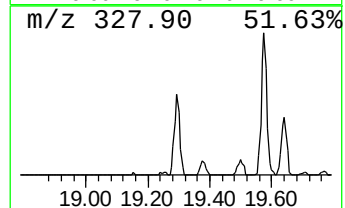
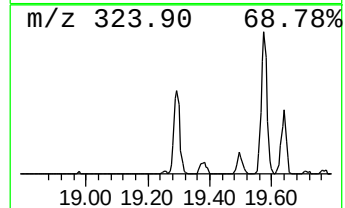
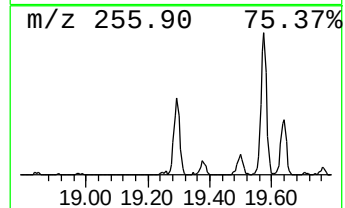
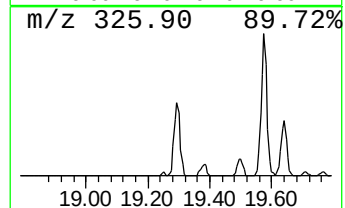
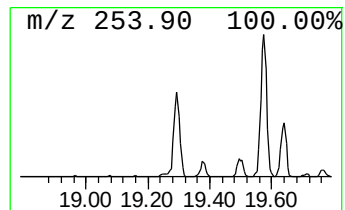
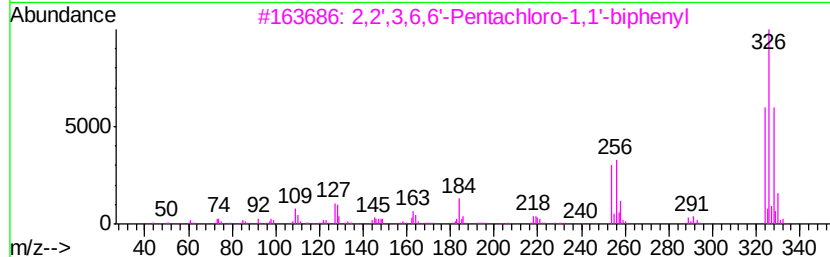
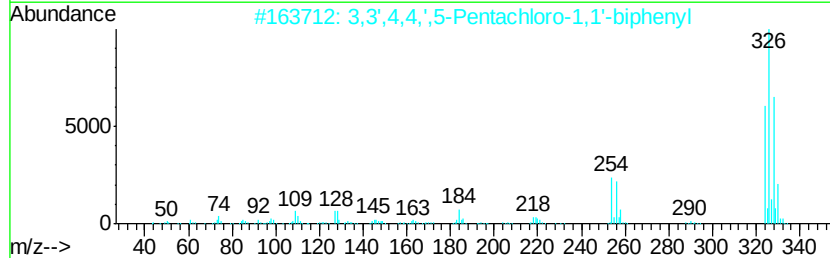
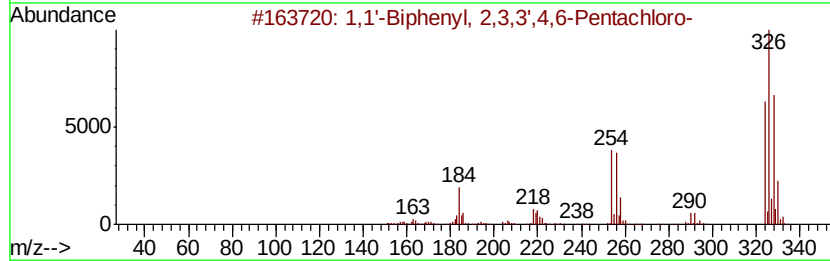
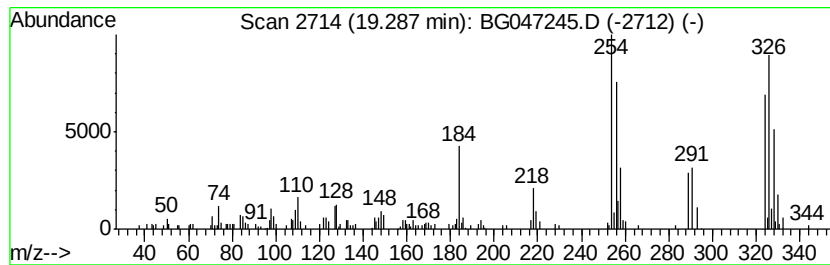
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
19.29	4.13 ng/ul	191177	Phenanthrene-d10	17.38		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,3',4,6-Pentac...	324	C12H5Cl5	074472-35-8	96	
2	3,3',4,4',5-Pentachloro-1,1'-bi...	324	C12H5Cl5	057465-28-8	95	
3	2,2',3,6,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-54-9	95	
4	1,1'-Biphenyl, 2,2',4,5',6-Penta...	324	C12H5Cl5	060145-21-3	95	
5	2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	070362-41-3	95	



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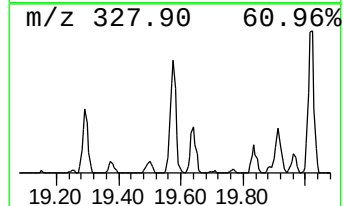
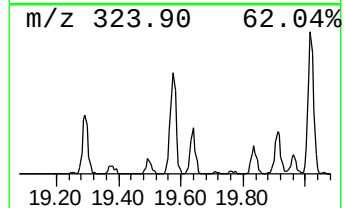
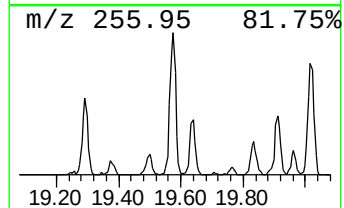
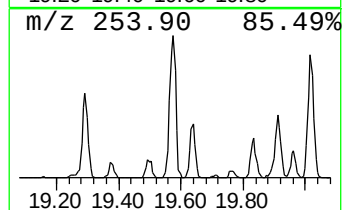
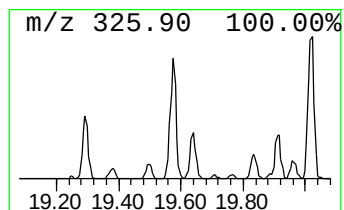
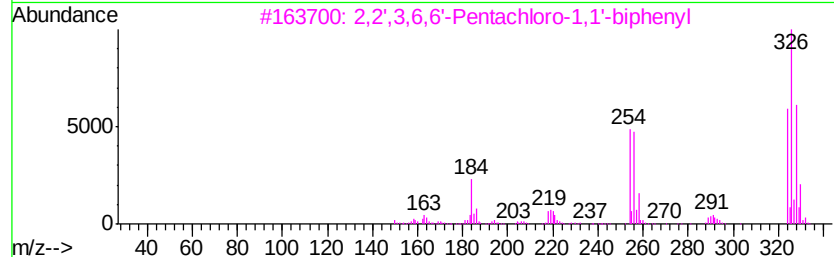
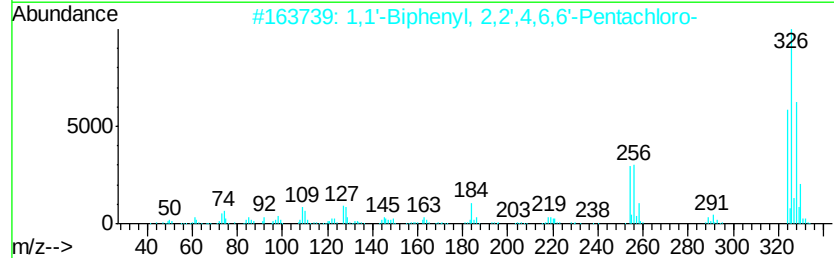
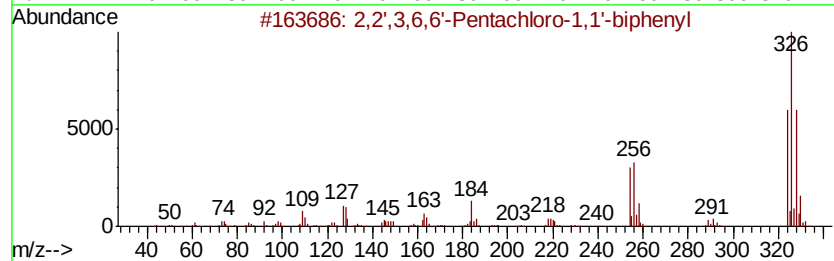
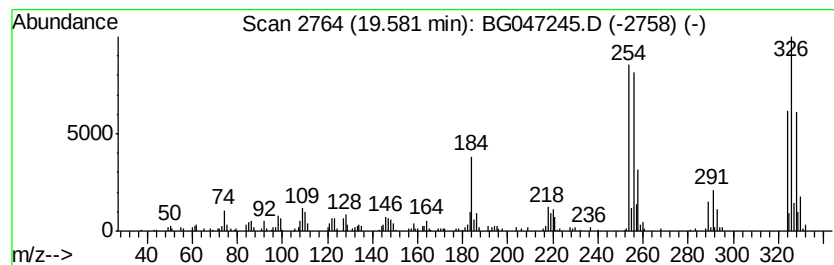
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
19.58	5.17 ng/ul	287425	Chrysene-d12	21.65		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,6,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-54-9	98	
2	1,1'-Biphenyl, 2,2',4,6,6'-Penta...	324	C12H5Cl5	056558-16-8	98	
3	2,2',3,6,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-54-9	96	
4	1,1'-Biphenyl, 2,2',4,5,5'-penta...	324	C12H5Cl5	037680-73-2	95	
5	1,1'-Biphenyl, 2,3,4,4',5-Pentac...	324	C12H5Cl5	074472-37-0	95	



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110620\
Data File : BG047245.D
Acq On : 6 Nov 2020 15:49
Operator : CG/JU
Sample : L4534-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

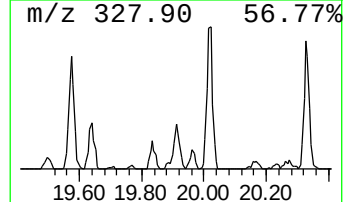
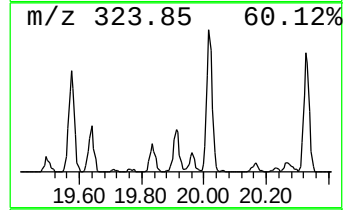
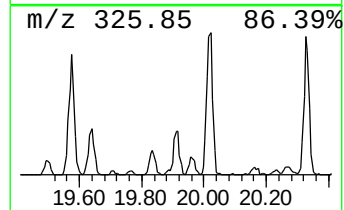
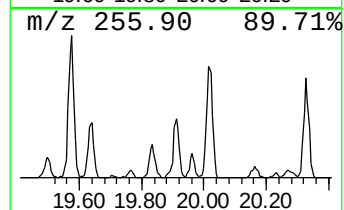
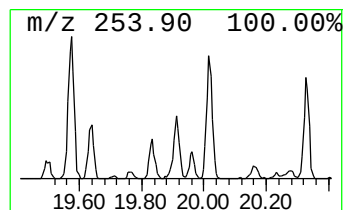
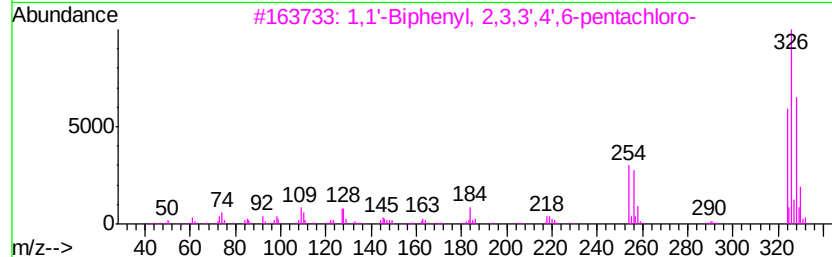
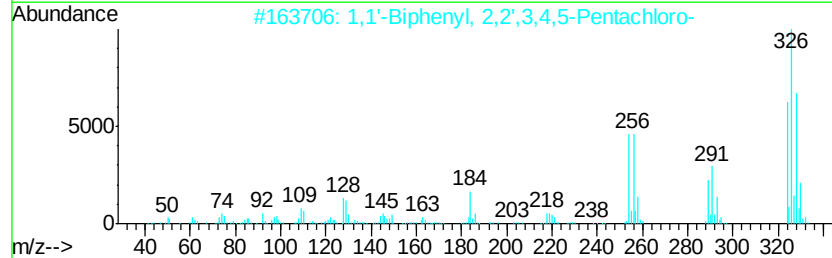
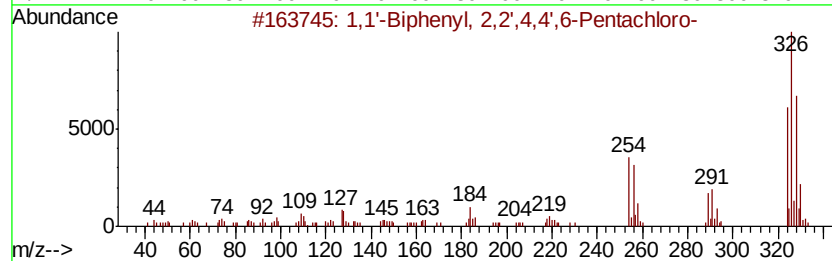
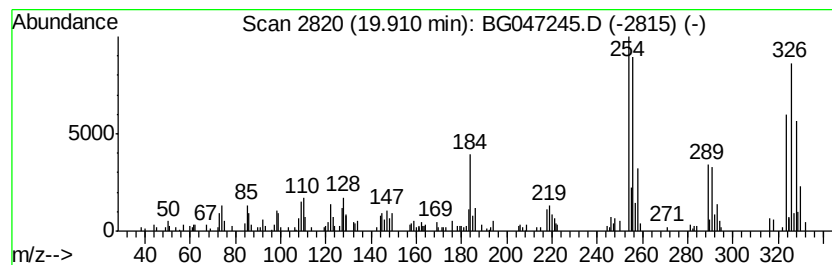
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 1,1'-Biphenyl, 2,2',4,4',6-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
19.91	2.49 ng/ul	138466	Chrysene-d12	21.65		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,4',6-Penta...	324	C12H5Cl5	039485-83-1	99	
2	1,1'-Biphenyl, 2,2',3,4,5-Pentac...	324	C12H5Cl5	055312-69-1	98	
3	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	98	
4	1,1'-Biphenyl, 2,2',4,6,6'-Penta...	324	C12H5Cl5	056558-16-8	98	
5	2,2',3,4,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-57-2	98	



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110620\
Data File : BG047245.D
Acq On : 6 Nov 2020 15:49
Operator : CG/JU
Sample : L4534-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

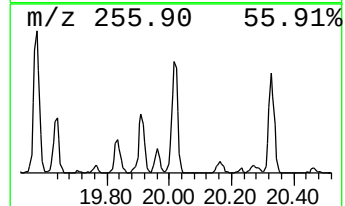
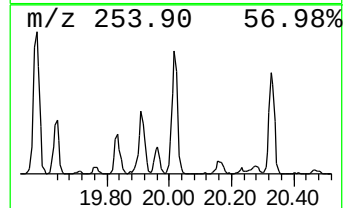
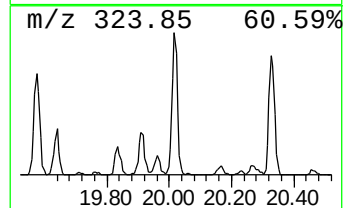
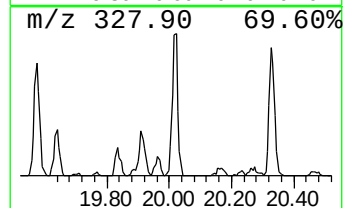
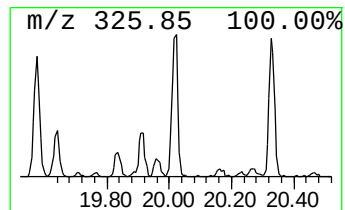
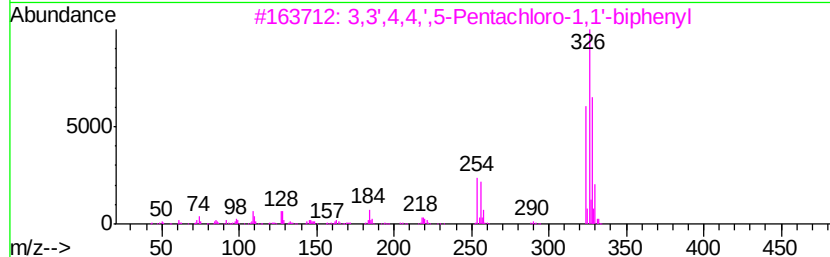
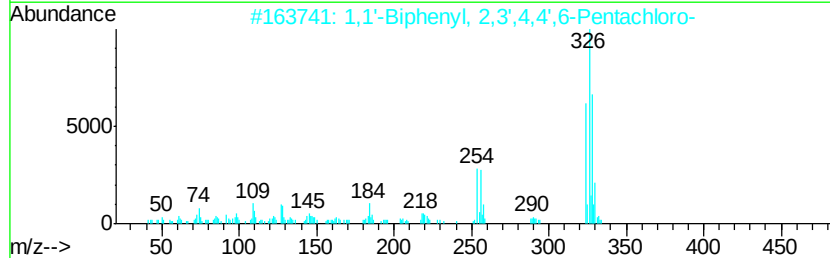
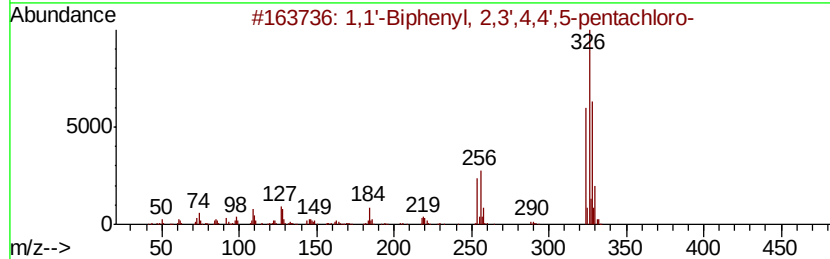
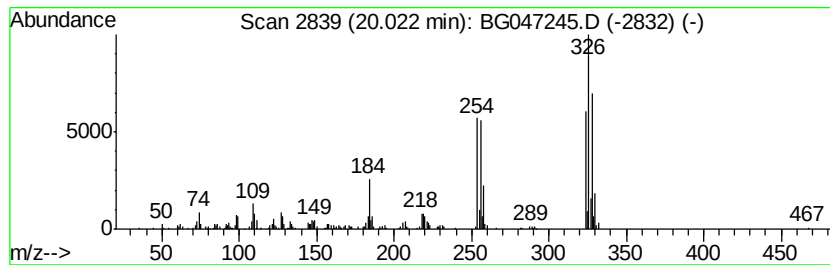
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.02	5.31 ng/ul	295156	Chrysene-d12		21.65	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'	-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	99
2	1,1'	-Biphenyl, 2,3',4,4',6-Penta...	324	C12H5Cl5	056558-17-9	97
3	3,3',4,4',	5-Pentachloro-1,1'-bi...	324	C12H5Cl5	057465-28-8	97
4	3,3',4,5,5'	-Pentachloro-1,1'-bip...	324	C12H5Cl5	039635-33-1	97
5	1,1'-Biphenyl,	2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	97



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110620\
Data File : BG047245.D
Acq On : 6 Nov 2020 15:49
Operator : CG/JU
Sample : L4534-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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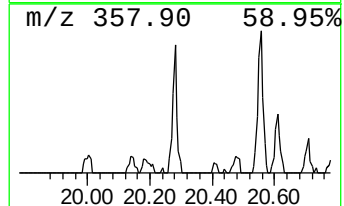
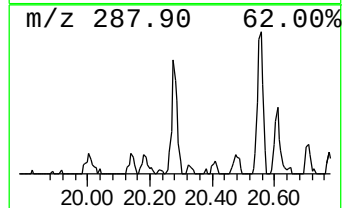
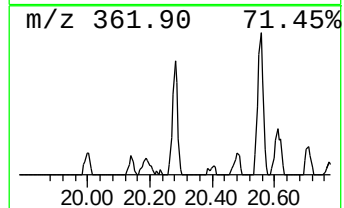
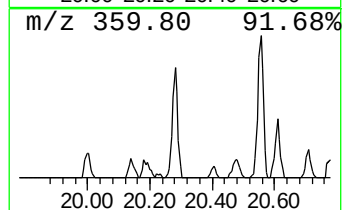
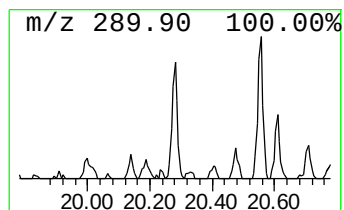
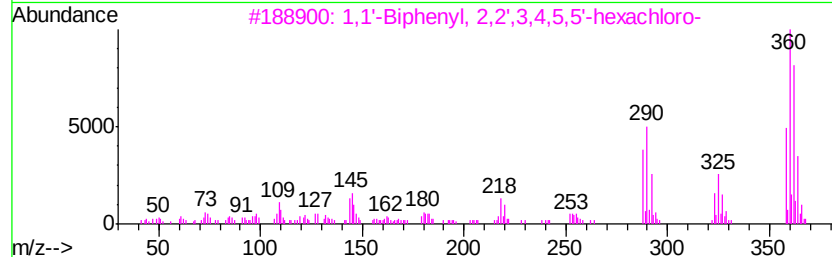
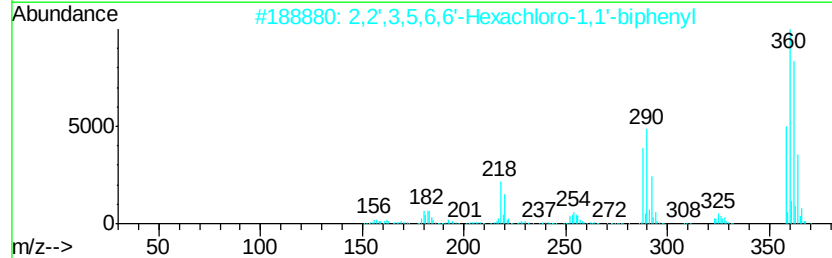
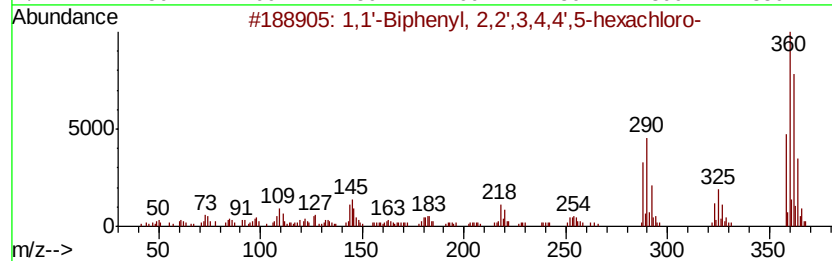
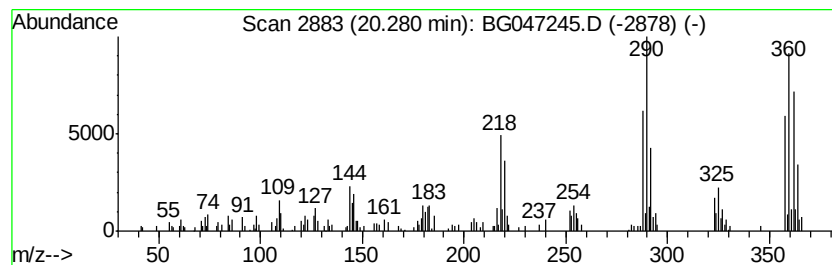
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1,1'-Biphenyl, 2,2',3,4,4',... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.28	2.21 ng/ul	122871	Chrysene-d12	21.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',3,4,4',5-hex...	358	C12H4Cl6	035694-06-5	99
2		2,2',3,5,6,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-09-2	99
3		1,1'-Biphenyl, 2,2',3,4,5,5'-hex...	358	C12H4Cl6	052712-04-6	99
4		2,2',3,4,6,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-40-5	98
5		2,2',3,4',6,6'-Hexachloro-1,1'-b...	358	C12H4Cl6	068194-08-1	98



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110620\
Data File : BG047245.D
Acq On : 6 Nov 2020 15:49
Operator : CG/JU
Sample : L4534-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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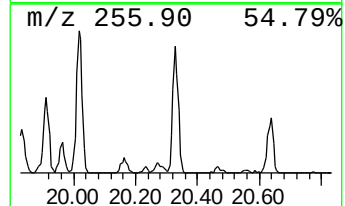
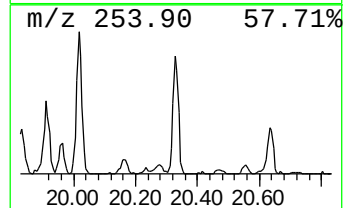
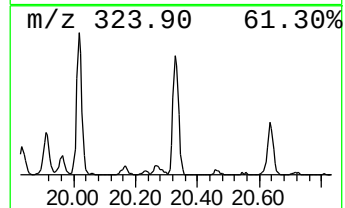
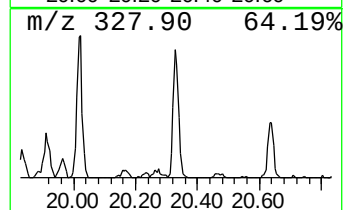
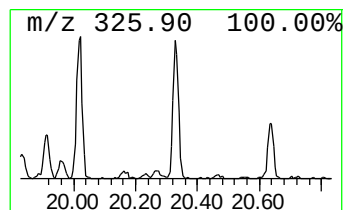
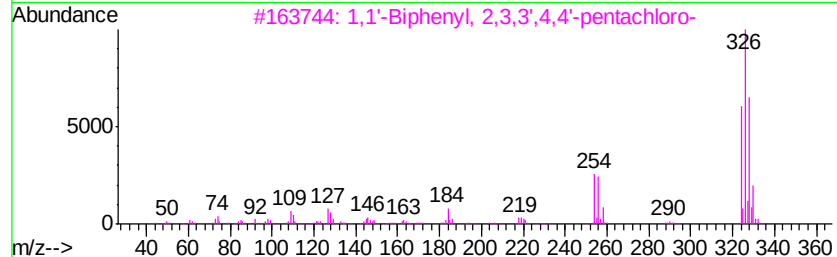
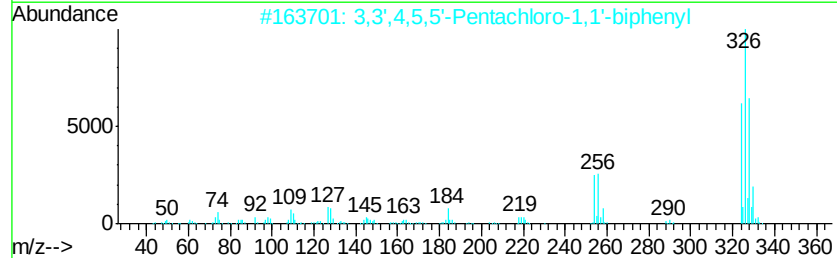
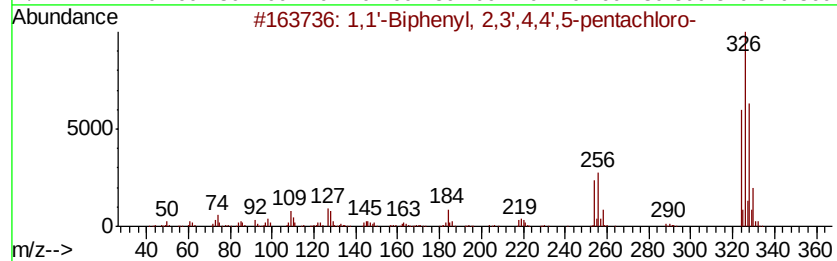
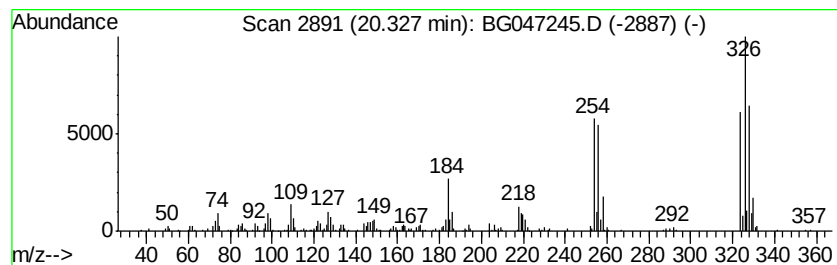
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.33	4.07 ng/ul	226401	Chrysene-d12	21.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	97
2		3,3',4,5,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	039635-33-1	97
3		1,1'-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	96
4		2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	070362-41-3	96
5		2,2',3,4',6-Pentachloro-1,1'-bip...	324	C12H5Cl5	068194-05-8	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110620\
Data File : BG047245.D
Acq On : 6 Nov 2020 15:49
Operator : CG/JU
Sample : L4534-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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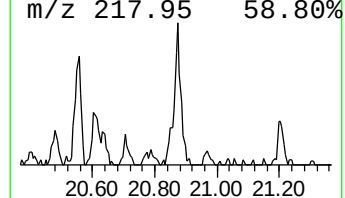
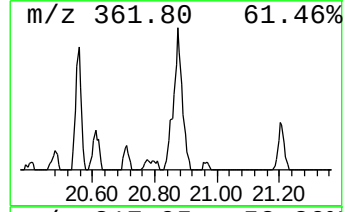
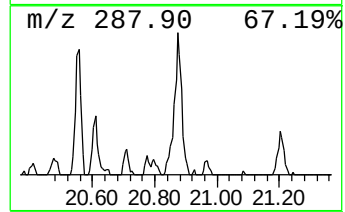
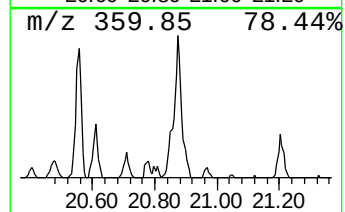
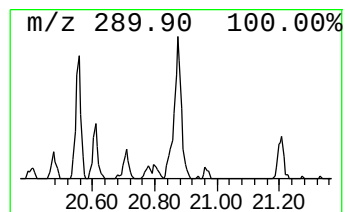
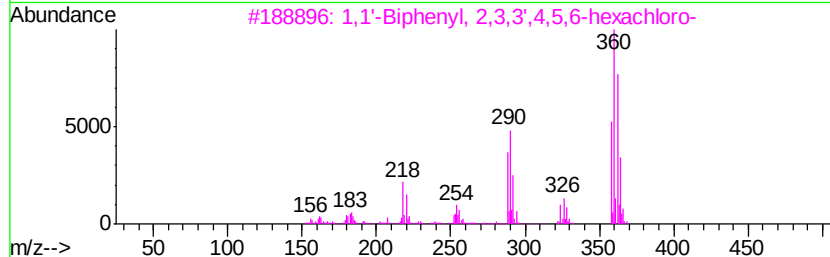
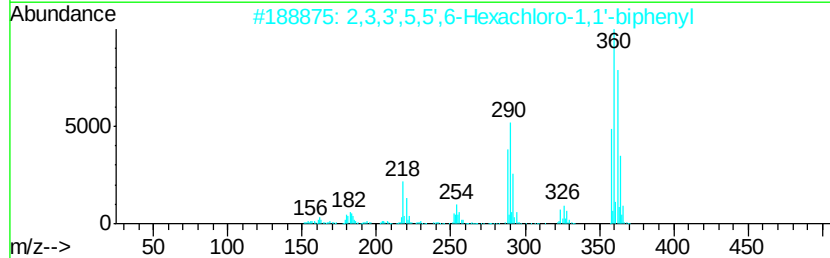
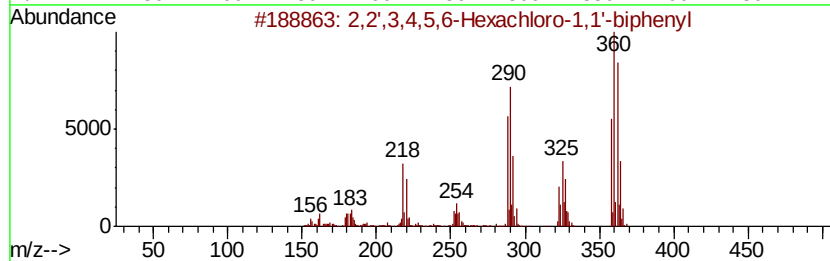
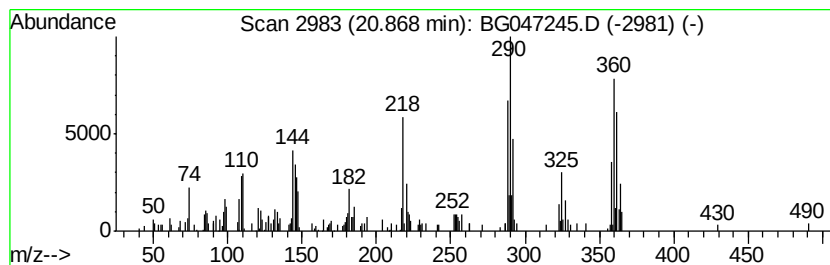
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 2,2',3,4,5,6-Hexachloro-1,1... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.87	3.36 ng/ul	186528	Chrysene-d12	21.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2',3,4,5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-61-4	96
2		2,3,3',5,5',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-46-1	96
3		1,1'-Biphenyl, 2,3,3',4,5,6-hexa...	358	C12H4Cl6	041411-62-5	96
4		1,1'-Biphenyl, 2,3,3',4,5,6-hexa...	358	C12H4Cl6	041411-62-5	96
5		1,1'-Biphenyl, 2,2',3,4',5',6-he...	358	C12H4Cl6	038380-04-0	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG110620\
Data File : BG047245.D
Acq On : 6 Nov 2020 15:49
Operator : CG/JU
Sample : L4534-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.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
1,1'-Biphenyl, 2,...	18.45	2.9	ng/ul	134429	4	17.38	926010	20.0
1,1'-Biphenyl, 2,...	19.29	4.1	ng/ul	191177	4	17.38	926010	20.0
2,2',3,6,6'-Penta...	19.58	5.2	ng/ul	287425	5	21.65	1111850	20.0
1,1'-Biphenyl, 2,...	19.91	2.5	ng/ul	138466	5	21.65	1111850	20.0
1,1'-Biphenyl, 2,...	20.02	5.3	ng/ul	295156	5	21.65	1111850	20.0
1,1'-Biphenyl, 2,...	20.28	2.2	ng/ul	122871	5	21.65	1111850	20.0
3,3',4,5,5'-Penta...	20.33	4.1	ng/ul	226401	5	21.65	1111850	20.0
2,3,4,4',5,6-Hexa...	20.56	2.4	ng/ul	134382	5	21.65	1111850	20.0
2,2',3,4,5,6-Hexa...	20.87	3.4	ng/ul	186528	5	21.65	1111850	20.0