

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110920\
 Data File : BG047261.D
 Acq On : 9 Nov 2020 18:47
 Operator : CG/JU
 Sample : L4534-05
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0BP2

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	3.452	16	19	21	rBV	1304	1585	0.15%	0.022%
2	3.552	35	36	41	rBV	1113	1704	0.16%	0.024%
3	3.640	45	51	52	rVV	6092	5576	0.53%	0.078%
4	3.669	52	56	60	rVB2	19719	24123	2.27%	0.338%
5	4.010	111	114	115	rVB2	2313	1759	0.17%	0.025%
6	4.063	115	123	129	rBV2	6702	11785	1.11%	0.165%
7	4.157	135	139	142	rBV3	8317	10551	0.99%	0.148%
8	4.380	174	177	180	rBV	1319	1638	0.15%	0.023%
9	4.592	210	213	215	rVB	1460	1724	0.16%	0.024%
10	4.786	243	246	248	rBV	1584	1572	0.15%	0.022%
11	4.815	248	251	255	rBV3	1656	2326	0.22%	0.033%
12	5.021	282	286	289	rVV3	1595	2288	0.22%	0.032%
13	5.297	330	333	335	rBV	1128	1370	0.13%	0.019%
14	6.231	487	492	494	rBV2	1164	1635	0.15%	0.023%
15	6.448	526	529	532	rBV2	1125	1805	0.17%	0.025%
16	6.495	535	537	542	rVB	1833	2094	0.20%	0.029%
17	6.548	545	546	551	rVB2	1436	1662	0.16%	0.023%
18	6.707	569	573	574	rBV2	1110	1434	0.14%	0.020%
19	6.760	579	582	585	rBV2	1050	1607	0.15%	0.023%
20	6.971	616	618	622	rVB	1394	1517	0.14%	0.021%
21	7.559	715	718	721	rBV2	1483	1462	0.14%	0.020%
22	8.005	787	794	801	rBV	227400	397219	37.40%	5.570%
23	8.052	801	802	807	rVB2	3131	2627	0.25%	0.037%
24	10.403	1200	1202	1208	rBV	1001	1675	0.16%	0.023%
25	10.502	1217	1219	1221	rVB	2576	2100	0.20%	0.029%
26	10.814	1264	1272	1285	rBV	268764	494782	46.59%	6.938%
27	11.437	1374	1378	1383	rBV2	1161	2379	0.22%	0.033%
28	11.507	1386	1390	1392	rBV3	1379	1713	0.16%	0.024%
29	12.812	1609	1612	1616	rVB	926	1410	0.13%	0.020%
30	13.822	1782	1784	1788	rBV2	1205	1382	0.13%	0.019%
31	13.910	1796	1799	1801	rBV2	1120	1352	0.13%	0.019%
32	14.022	1816	1818	1821	rBV2	1282	1702	0.16%	0.024%
33	14.181	1842	1845	1847	rVB2	1587	1670	0.16%	0.023%
34	14.227	1850	1853	1856	rBV	1017	1849	0.17%	0.026%

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35	14.557	1908	1909	1913	rBV2	1440	1417	0.13%	0.020%
36	14.639	1917	1923	1932	rBV	424463	668043	62.90%	9.367%
37	14.768	1942	1945	1949	rVB3	1078	1537	0.14%	0.022%
38	14.850	1958	1959	1964	rVB2	1356	1357	0.13%	0.019%
39	15.479	2065	2066	2070	rBV2	2048	1809	0.17%	0.025%
40	15.532	2073	2075	2078	rBV2	1631	2080	0.20%	0.029%
41	15.649	2093	2095	2098	rBV2	1254	1552	0.15%	0.022%
42	15.685	2098	2101	2103	rBV	2484	2948	0.28%	0.041%
43	15.978	2149	2151	2154	rBV	1117	1624	0.15%	0.023%
44	16.096	2168	2171	2173	rVB	2033	1629	0.15%	0.023%
45	16.343	2208	2213	2215	rBV2	2273	2491	0.23%	0.035%
46	16.425	2224	2227	2230	rVB2	1952	1909	0.18%	0.027%
47	16.965	2318	2319	2323	rVB	2721	2270	0.21%	0.032%
48	17.124	2344	2346	2351	rBV2	1834	3319	0.31%	0.047%
49	17.265	2367	2370	2371	rVB	2196	1767	0.17%	0.025%
50	17.383	2384	2390	2395	rBV	511250	793218	74.69%	11.122%
51	17.424	2395	2397	2405	rVV	52137	78720	7.41%	1.104%
52	17.488	2406	2408	2409	rVB2	3306	1891	0.18%	0.027%
53	17.559	2417	2420	2425	rVB5	6007	7046	0.66%	0.099%
54	17.829	2462	2466	2468	rBV2	2138	2820	0.27%	0.040%
55	17.888	2474	2476	2477	rBV	2548	2522	0.24%	0.035%
56	17.988	2486	2493	2498	rBV2	27822	53661	5.05%	0.752%
57	18.099	2509	2512	2513	rBV2	2237	2106	0.20%	0.030%
58	18.170	2520	2524	2525	rBV2	1786	2134	0.20%	0.030%
59	18.205	2528	2530	2531	rVB2	2058	1364	0.13%	0.019%
60	18.229	2531	2534	2537	rBV4	9180	14099	1.33%	0.198%
61	18.311	2546	2548	2553	rVB4	8706	9521	0.90%	0.133%
62	18.452	2565	2572	2577	rBV2	87410	123993	11.68%	1.739%
63	18.511	2578	2582	2586	rVV3	29388	42214	3.97%	0.592%
64	18.552	2586	2589	2593	rVV4	13389	15998	1.51%	0.224%
65	18.663	2606	2608	2610	rBV2	1549	1521	0.14%	0.021%
66	18.734	2615	2620	2623	rBV	43612	54745	5.15%	0.768%
67	18.840	2634	2638	2640	rBV3	5124	7022	0.66%	0.098%
68	18.869	2640	2643	2646	rVV4	11263	13893	1.31%	0.195%
69	18.910	2646	2650	2654	rVB2	24921	32279	3.04%	0.453%
70	18.969	2658	2660	2663	rBV3	2963	3236	0.30%	0.045%
71	19.010	2663	2667	2669	rVB4	6894	7357	0.69%	0.103%

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72	19.134	2686	2688	2689	rBV2	4207	2684	0.25%	0.038%
73	19.169	2691	2694	2696	rVV3	6316	9481	0.89%	0.133%
74	19.216	2698	2702	2706	rVV4	25939	37947	3.57%	0.532%
75	19.263	2706	2710	2712	rVV2	54293	76521	7.21%	1.073%
76	19.292	2712	2715	2724	rVV3	111413	173184	16.31%	2.428%
77	19.374	2726	2729	2733	rVV5	20922	27266	2.57%	0.382%
78	19.433	2733	2739	2745	rVV	96996	151896	14.30%	2.130%
79	19.504	2746	2751	2755	rVV8	27872	59210	5.58%	0.830%
80	19.574	2758	2763	2768	rVV	164673	218429	20.57%	3.063%
81	19.633	2769	2773	2779	rVB	61911	79234	7.46%	1.111%
82	19.680	2779	2781	2783	rBV3	4674	3024	0.28%	0.042%
83	19.762	2792	2795	2798	rBV3	17729	25509	2.40%	0.358%
84	19.797	2798	2801	2804	rVV	32122	40690	3.83%	0.571%
85	19.833	2804	2807	2811	rVB2	43165	51908	4.89%	0.728%
86	19.909	2816	2820	2825	rVB3	76023	98693	9.29%	1.384%
87	19.962	2825	2829	2832	rBV6	28598	38112	3.59%	0.534%
88	20.015	2833	2838	2843	rVB2	164181	218991	20.62%	3.071%
89	20.279	2877	2883	2887	rBV4	80017	116241	10.95%	1.630%
90	20.326	2887	2891	2896	rVB	143141	171409	16.14%	2.403%
91	20.549	2926	2929	2934	rVB2	86453	109301	10.29%	1.533%
92	20.608	2935	2939	2941	rVV5	41451	58691	5.53%	0.823%
93	20.632	2941	2943	2947	rVB	59780	69254	6.52%	0.971%
94	20.708	2953	2956	2960	rVB5	20347	24988	2.35%	0.350%
95	20.773	2965	2967	2970	rBV4	12174	16773	1.58%	0.235%
96	20.873	2975	2984	2993	rVV3	107220	225360	21.22%	3.160%
97	20.967	2996	3000	3003	rBV6	13272	18651	1.76%	0.262%
98	21.202	3037	3040	3049	rVB10	30061	41848	3.94%	0.587%
99	21.642	3109	3115	3129	rVB	595079	1062017	100.00%	14.891%
100	24.833	3650	3658	3671	rVB2	337384	944466	88.93%	13.243%

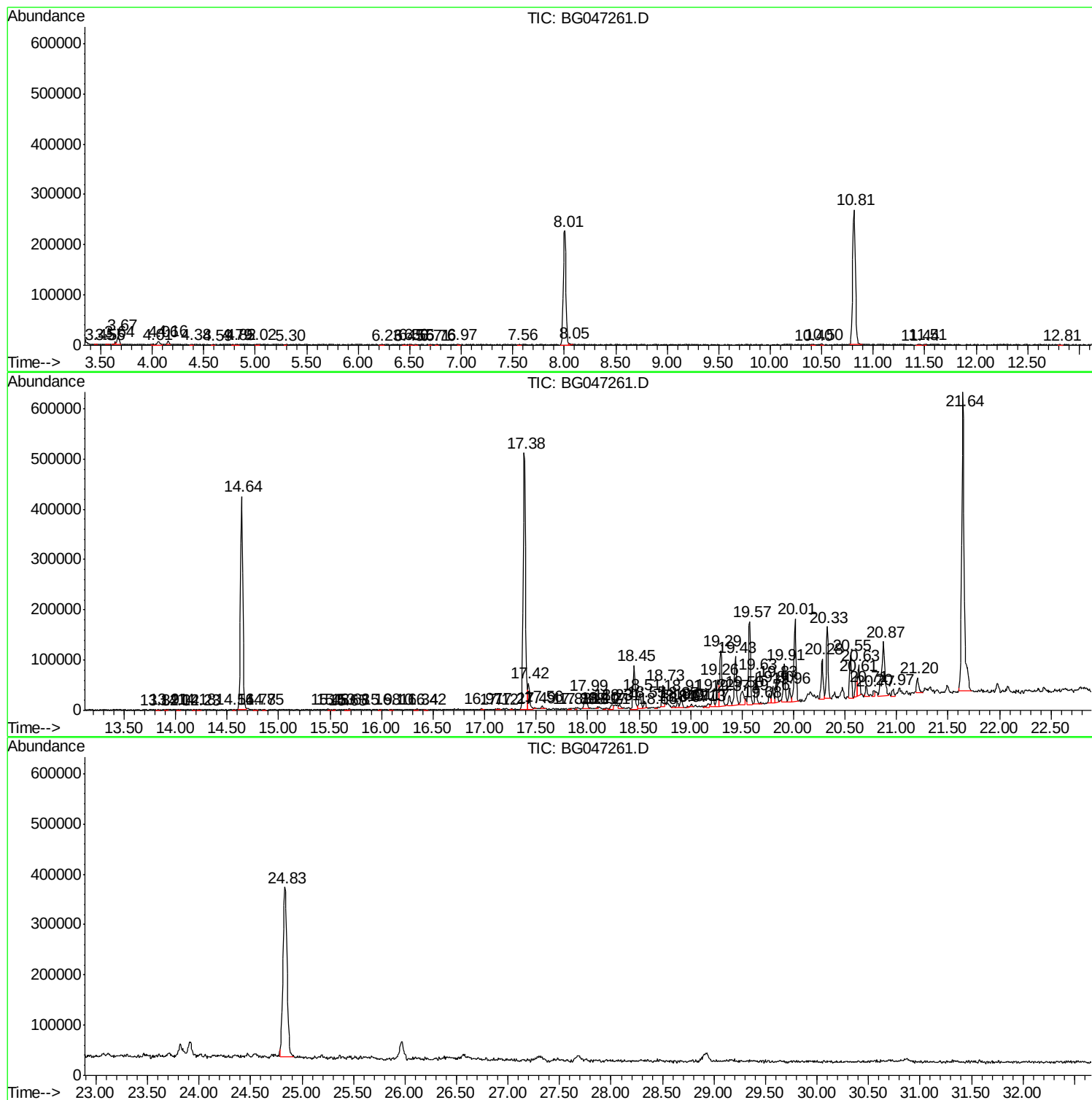
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

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



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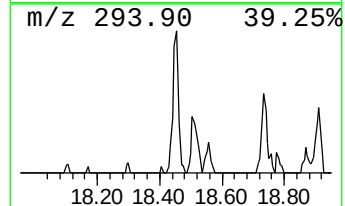
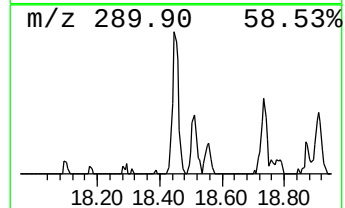
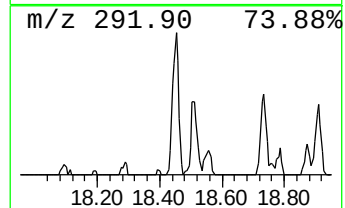
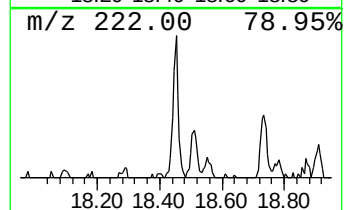
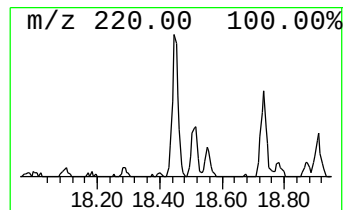
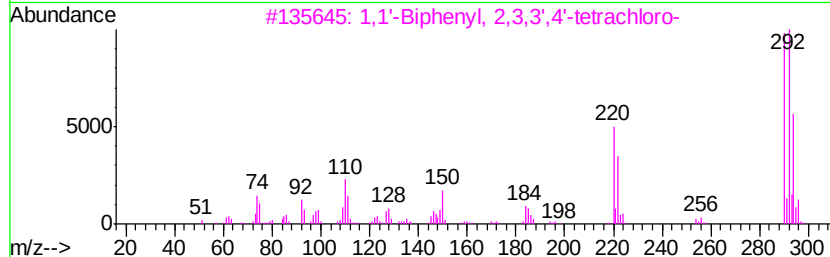
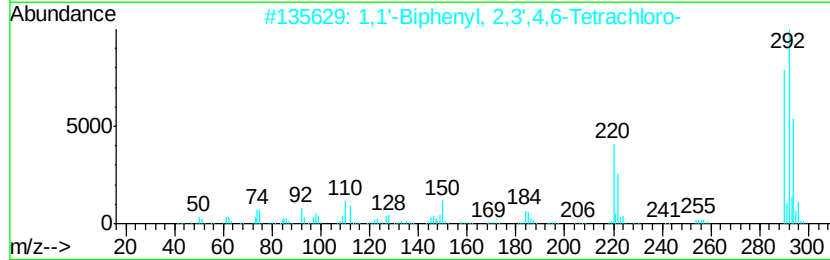
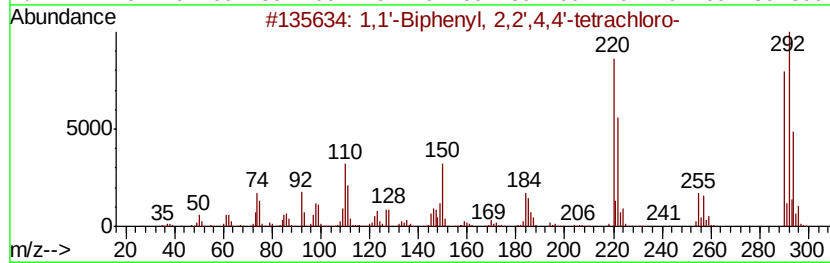
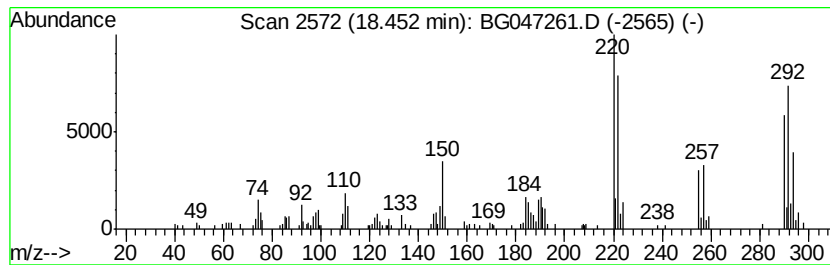
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.45	3.13 ng/ul	123993	Phenanthrene-d10	17.38	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
2	1,1'-Biphenyl, 2,3',4,6-Tetrachl...	290	C12H6Cl4	060233-24-1	99
3	1,1'-Biphenyl, 2,3,3',4'-tetrach...	290	C12H6Cl4	041464-43-1	99
4	1,1'-Biphenyl, 2,3,3',5'-tetrach...	290	C12H6Cl4	041464-49-7	99
5	2,3,3',6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074472-33-6	99



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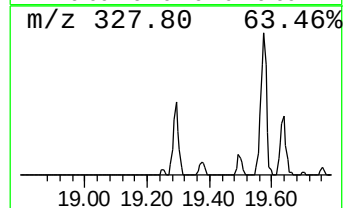
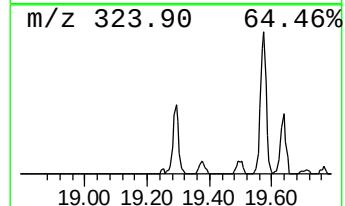
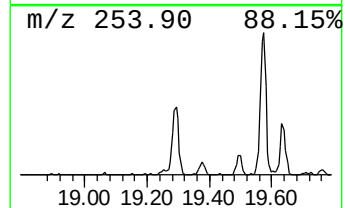
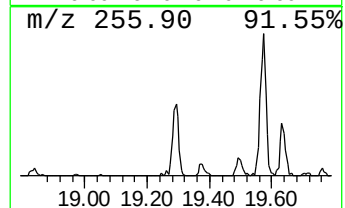
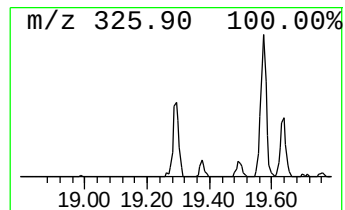
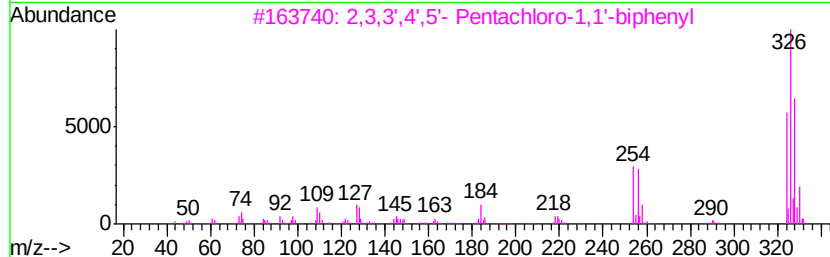
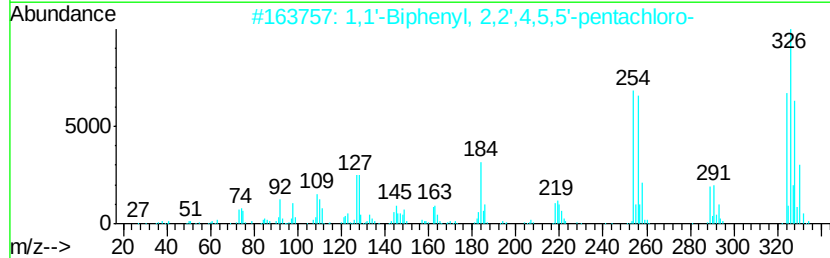
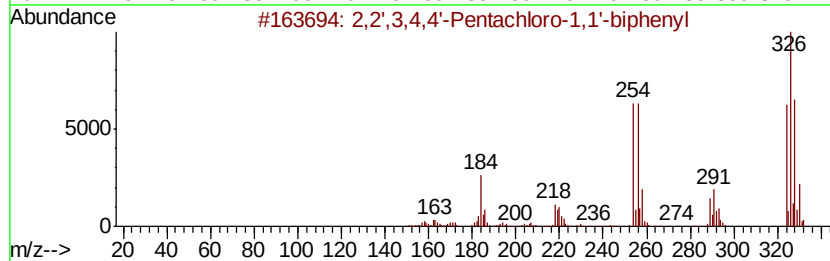
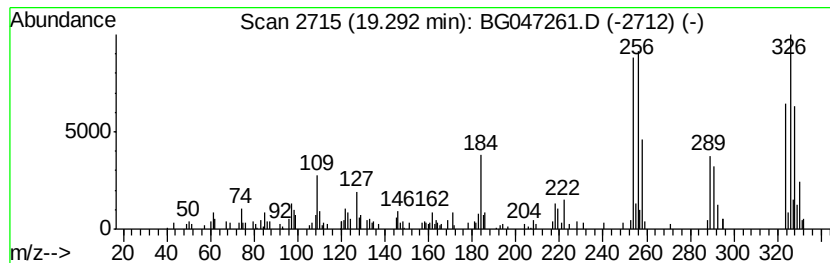
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.29	4.37 ng/ul	173184	Phenanthrene-d10	17.38	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,4,4'-Pentachloro-1,1'-bip...	324	C12H5Cl5	065510-45-4	99
2	1,1'-Biphenyl, 2,2',4,5,5'-penta...	324	C12H5Cl5	037680-73-2	98
3	2,3,3',4',5'- Pentachloro-1,1'-b...	324	C12H5Cl5	076842-07-4	98
4	2,2',3,4,4'-Pentachloro-1,1'-bip...	324	C12H5Cl5	065510-45-4	98
5	1,1'-Biphenyl, 2,2',3,4,5-Pentac...	324	C12H5Cl5	055312-69-1	98



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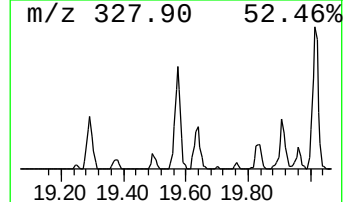
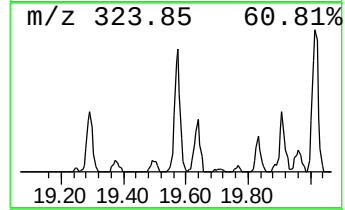
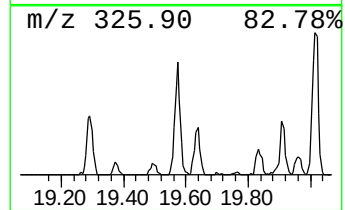
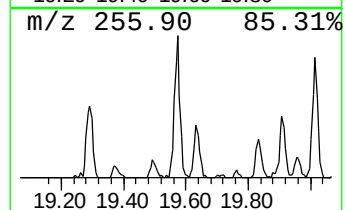
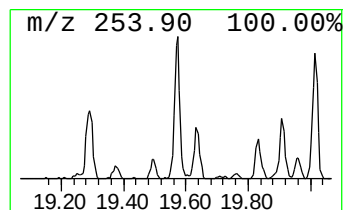
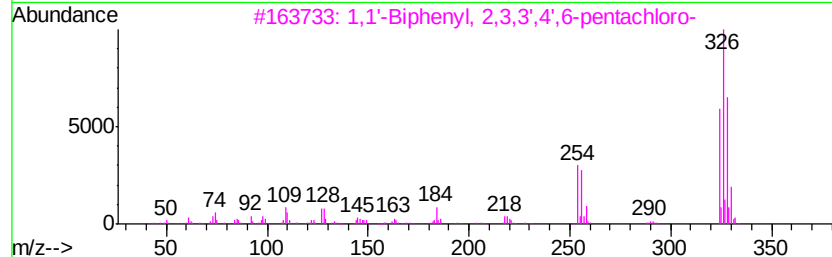
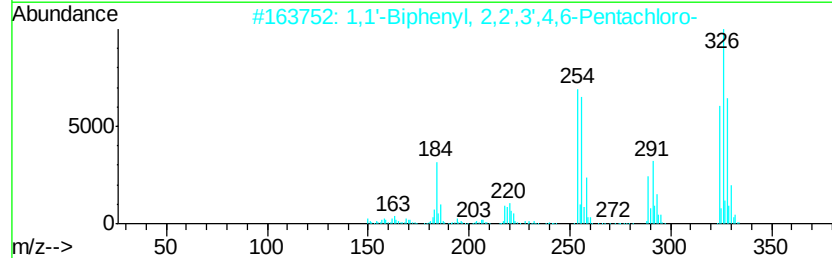
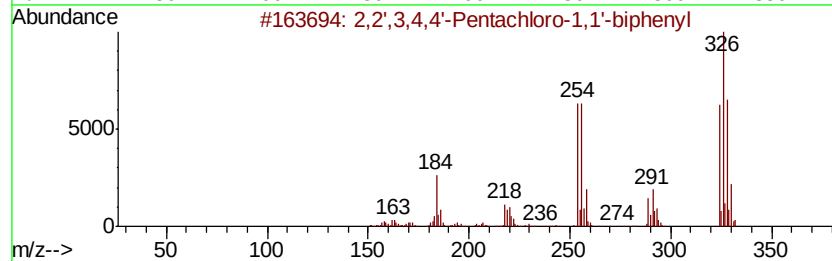
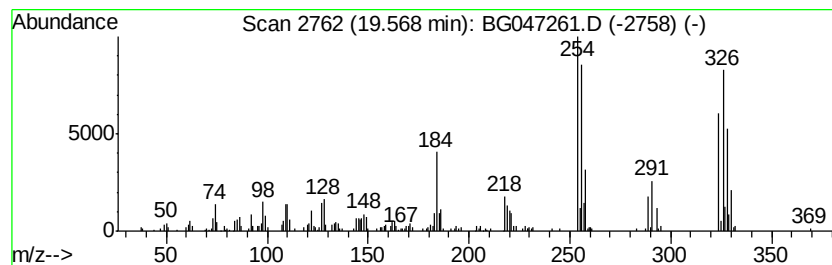
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
19.57	4.11 ng/ul	218429	Chrysene-d12	21.64		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,4,4'-Pentachloro-1,1'-bip...	324	C12H5Cl5	065510-45-4	95	
2	1,1'-Biphenyl, 2,2',3',4,6-Penta...	324	C12H5Cl5	060233-25-2	95	
3	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	95	
4	1,1'-Biphenyl, 2',3,4,5,5'-Penta...	324	C12H5Cl5	070424-70-3	95	
5	2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	070362-41-3	94	



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110920\
Data File : BG047261.D
Acq On : 9 Nov 2020 18:47
Operator : CG/JU
Sample : L4534-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

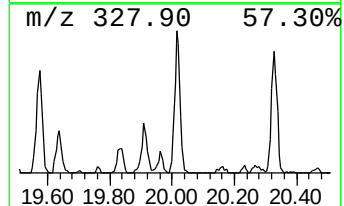
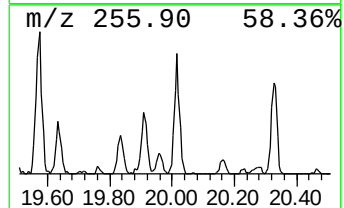
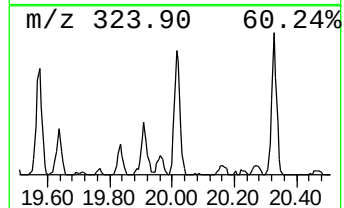
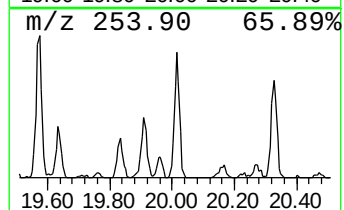
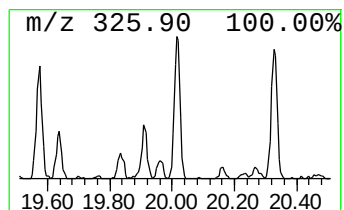
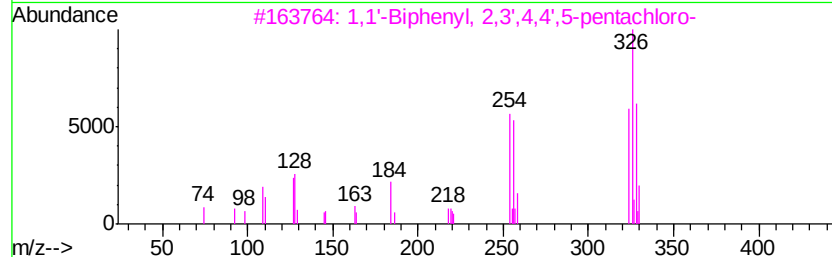
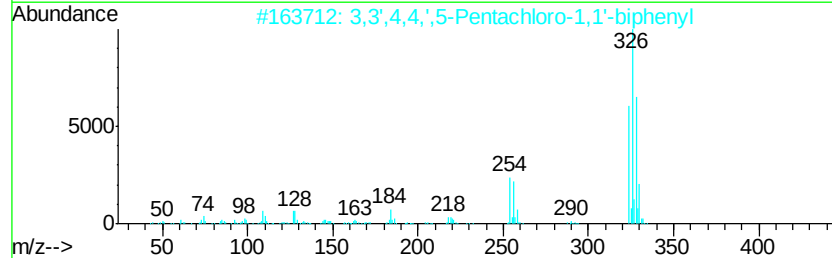
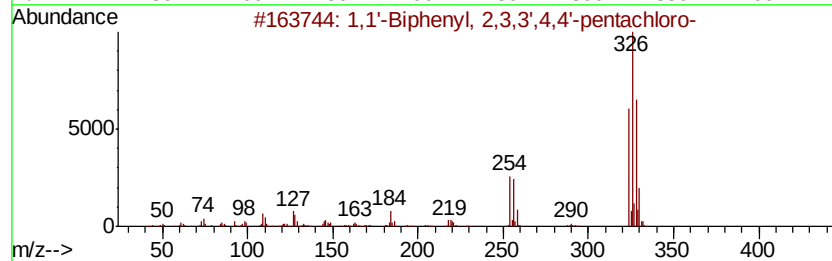
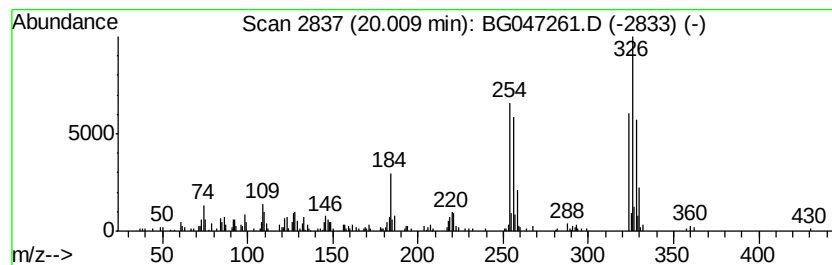
Instrument :
BNA_G
ClientSampleId :
C0BP2

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, 2,3,3',4,4'-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.01	4.12 ng/ul	218991	Chrysene-d12	21.64	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	97
2	3,3',4,4',5-Pentachloro-1,1'-bi...	324	C12H5Cl5	057465-28-8	96
3	1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	96
4	1,1'-Biphenyl, 2,3,4,5,6-pentach...	324	C12H5Cl5	018259-05-7	96
5	2,3,3',5,6-Pentachloro-1,1'-biph...	324	C12H5Cl5	074472-36-9	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110920\
Data File : BG047261.D
Acq On : 9 Nov 2020 18:47
Operator : CG/JU
Sample : L4534-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0BP2

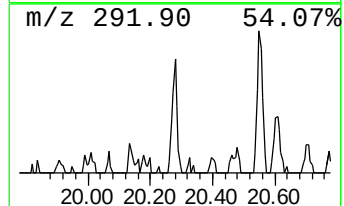
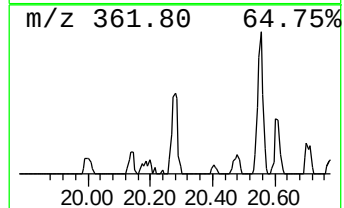
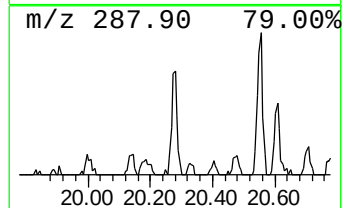
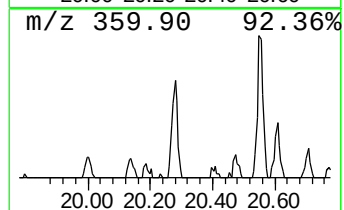
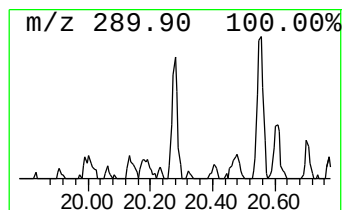
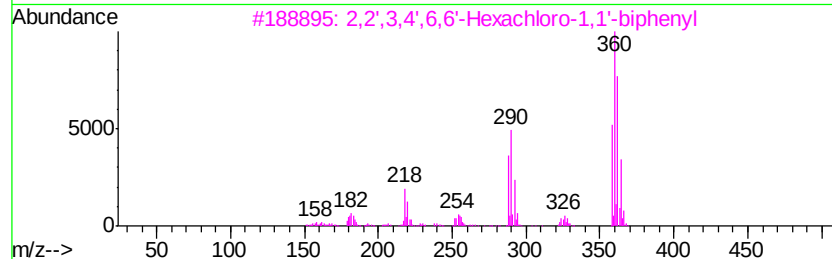
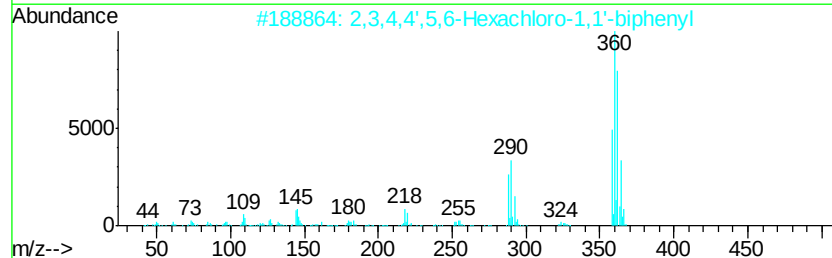
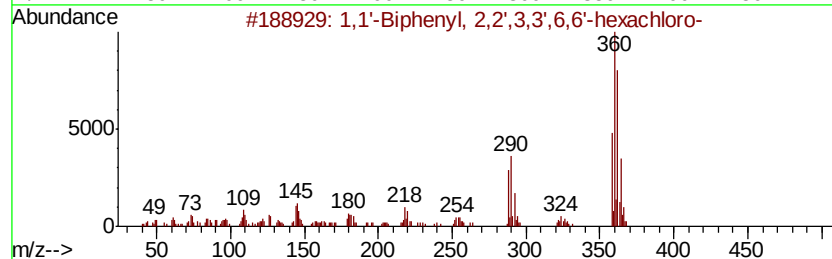
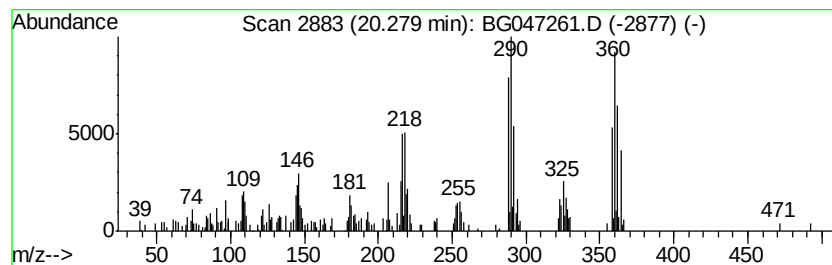
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,1'-Biphenyl, 2,2',3,3',6,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.28	2.19 ng/ul	116241	Chrysene-d12	21.64

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		2,3,4,4',5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-63-6	99
3		2,2',3,4',6,6'-Hexachloro-1,1'-b...	358	C12H4Cl6	068194-08-1	99
4		1,1'-Biphenyl, 2,2',4,4',5,5'-he...	358	C12H4Cl6	035065-27-1	99
5		1,1'-Biphenyl, 2,2',3,4,4',6-Hex...	358	C12H4Cl6	056030-56-9	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110920\
Data File : BG047261.D
Acq On : 9 Nov 2020 18:47
Operator : CG/JU
Sample : L4534-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

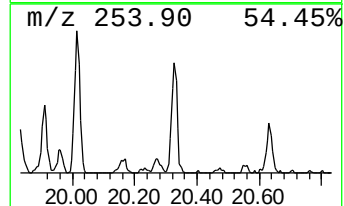
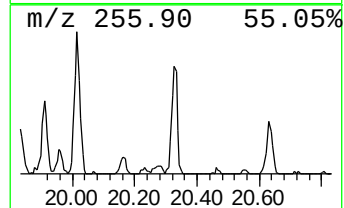
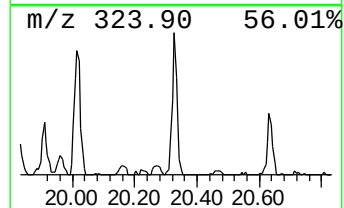
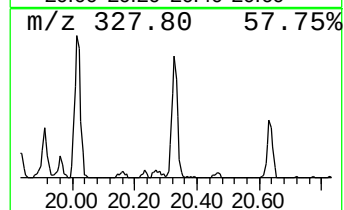
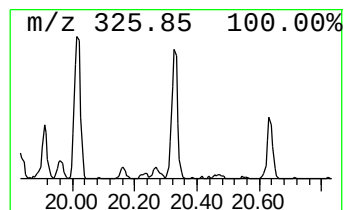
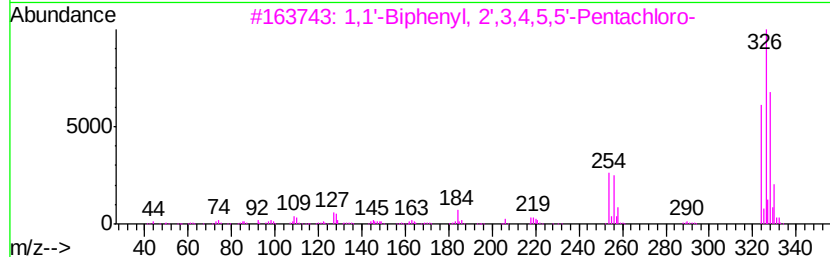
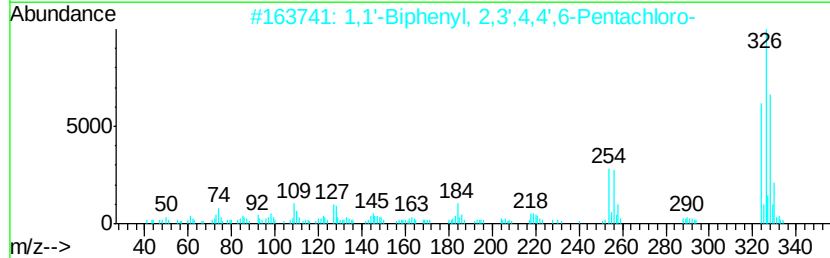
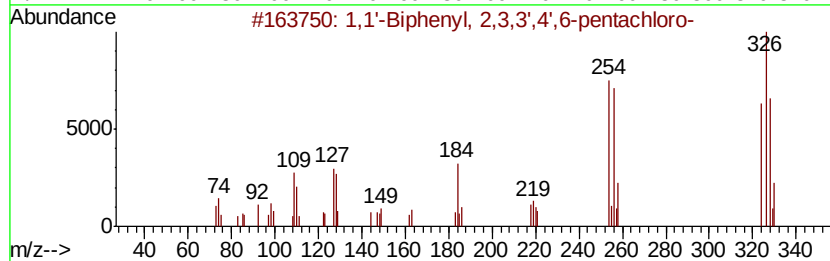
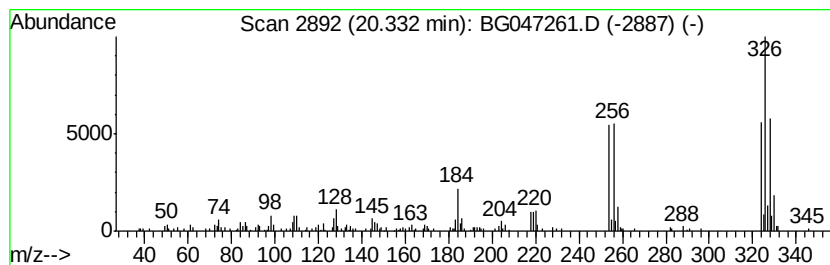
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.33	3.23 ng/ul	171409	Chrysene-d12	21.64	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99
2	1,1'-Biphenyl, 2,3',4,4',6-Penta...	324	C12H5Cl5	056558-17-9	98
3	1,1'-Biphenyl, 2',3,4,5,5'-Penta...	324	C12H5Cl5	070424-70-3	97
4	2,3,3',4,5-Pentachloro-1,1'-biph...	324	C12H5Cl5	070424-69-0	97
5	1,1'-Biphenyl, 2,2',4,4',5-penta...	324	C12H5Cl5	038380-01-7	97



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110920\
Data File : BG047261.D
Acq On : 9 Nov 2020 18:47
Operator : CG/JU
Sample : L4534-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

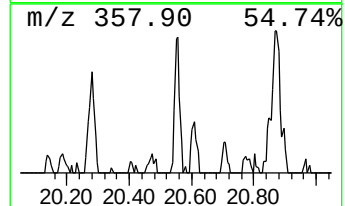
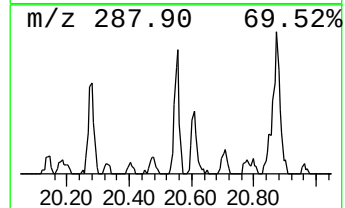
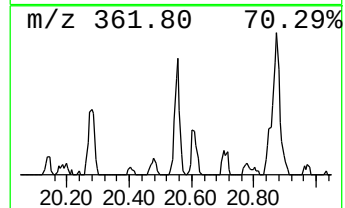
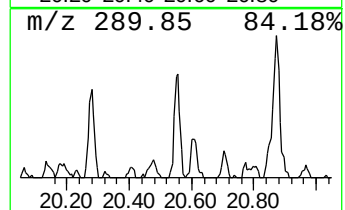
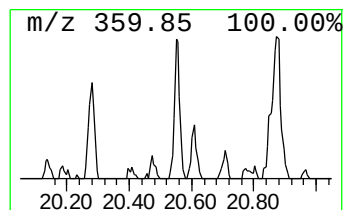
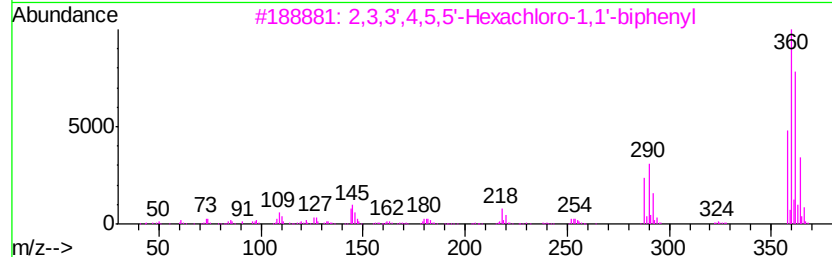
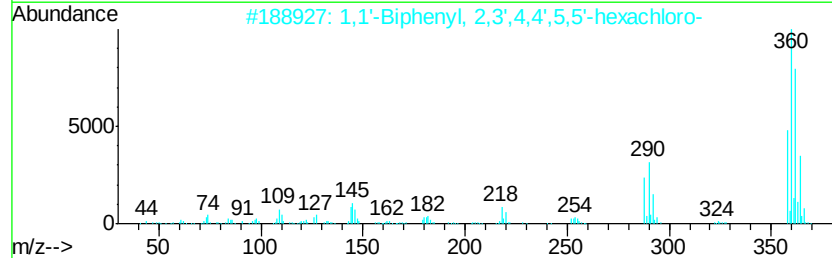
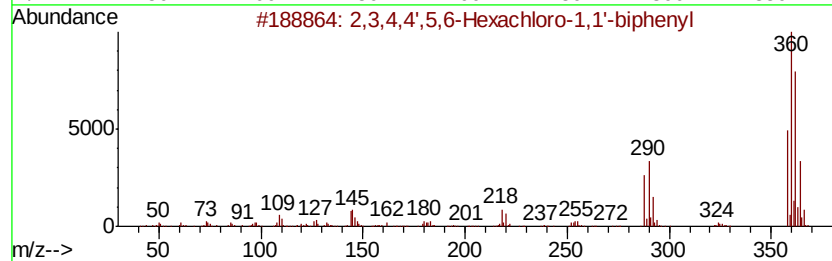
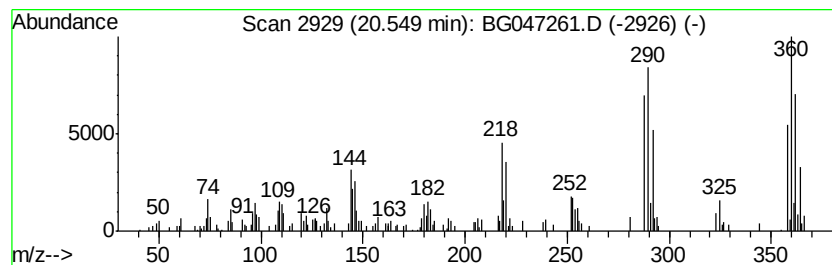
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 2,3,4,4',5,6-Hexachloro-1,1... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.55	2.06 ng/ul	109301	Chrysene-d12	21.64	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,4,4',5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-63-6	99
2	1,1'-Biphenyl, 2,3',4,4',5,5'-he...	358	C12H4Cl6	052663-72-6	98
3	2,3,3',4,5,5'-Hexachloro-1,1'-bi...	358	C12H4Cl6	039635-35-3	98
4	1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	97
5	2,3,3',4,4',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-42-7	97



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110920\
Data File : BG047261.D
Acq On : 9 Nov 2020 18:47
Operator : CG/JU
Sample : L4534-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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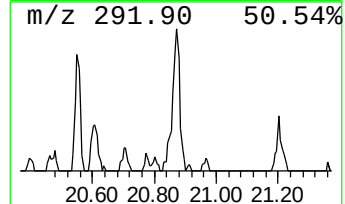
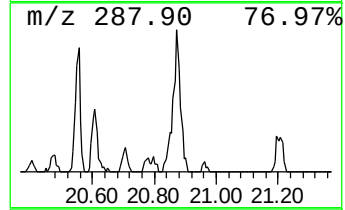
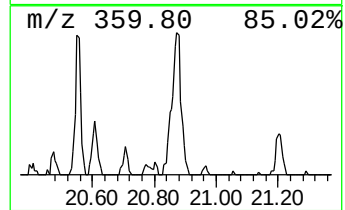
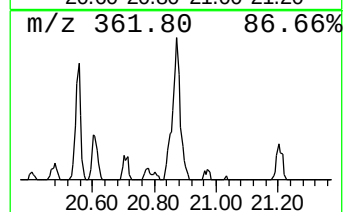
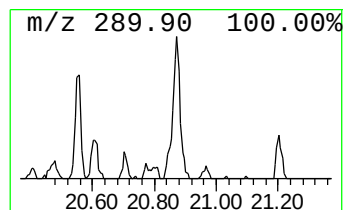
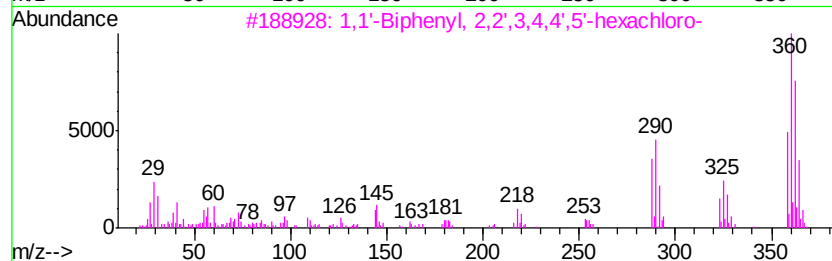
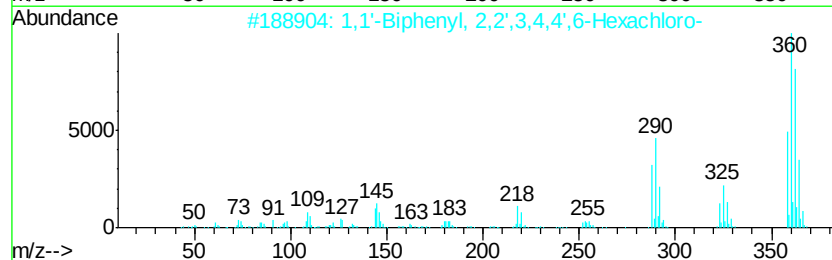
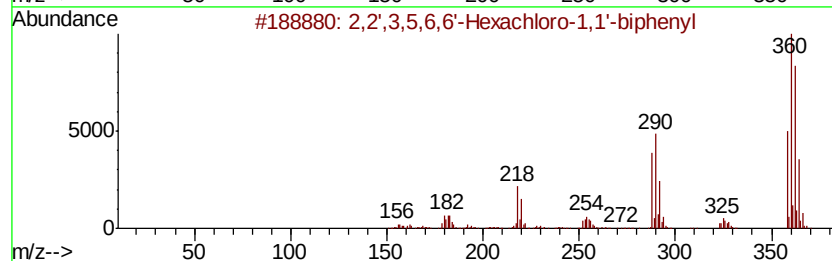
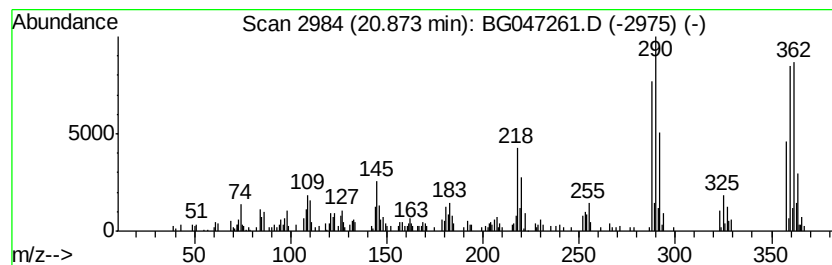
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 2,2',3,5,6,6'-Hexachloro-1,... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.87	4.24 ng/ul	225360	Chrysene-d12	21.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2',3,5,6,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-09-2	98
2		1,1'-Biphenyl, 2,2',3,4,4',6-Hex...	358	C12H4Cl6	056030-56-9	97
3		1,1'-Biphenyl, 2,2',3,4,4',5'-he...	358	C12H4Cl6	035065-28-2	97
4		1,1'-Biphenyl, 2,2',3,4,5,6'-Hex...	358	C12H4Cl6	068194-15-0	96
5		1,1'-Biphenyl, 2,2',3,3',4,4'-he...	358	C12H4Cl6	038380-07-3	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG110920\
Data File : BG047261.D
Acq On : 9 Nov 2020 18:47
Operator : CG/JU
Sample : L4534-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0BP2

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	3.1	ng/ul	123993	4	17.38	793218	20.0
2,2',3,4,4'-Penta...	19.29	4.4	ng/ul	173184	4	17.38	793218	20.0
1,1'-Biphenyl, 2,...	19.57	4.1	ng/ul	218429	5	21.64	1062020	20.0
1,1'-Biphenyl, 2,...	20.01	4.1	ng/ul	218991	5	21.64	1062020	20.0
1,1'-Biphenyl, 2,...	20.28	2.2	ng/ul	116241	5	21.64	1062020	20.0
1,1'-Biphenyl, 2,...	20.33	3.2	ng/ul	171409	5	21.64	1062020	20.0
2,3,4,4',5,6-Hexa...	20.55	2.1	ng/ul	109301	5	21.64	1062020	20.0
2,2',3,5,6,6'-Hex...	20.87	4.2	ng/ul	225360	5	21.64	1062020	20.0