

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110620\
 Data File : BG047247.D
 Acq On : 6 Nov 2020 17:10
 Operator : CG/JU
 Sample : L4534-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0BP3

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	3.508	26	29	34	rBV	1793	2156	0.20%	0.029%
2	3.596	43	44	47	rVB2	1754	1393	0.13%	0.019%
3	3.637	47	51	53	rBV3	4692	6009	0.54%	0.081%
4	3.666	53	56	64	rVB3	17082	22424	2.03%	0.302%
5	3.948	102	104	107	rVB2	1618	1386	0.13%	0.019%
6	4.013	111	115	117	rBV	1480	1604	0.15%	0.022%
7	4.054	119	122	130	rVV4	8677	14946	1.35%	0.201%
8	4.154	133	139	145	rBV3	9819	14854	1.35%	0.200%
9	4.601	209	215	221	rBV	3108	5222	0.47%	0.070%
10	4.712	230	234	235	rBV	1376	1846	0.17%	0.025%
11	5.006	281	284	288	rBV	1662	2283	0.21%	0.031%
12	5.570	376	380	382	rBV2	1150	1526	0.14%	0.021%
13	5.629	386	390	392	rVV2	1676	1539	0.14%	0.021%
14	6.187	482	485	488	rVB	1070	1421	0.13%	0.019%
15	6.322	504	508	511	rVB	1212	1822	0.17%	0.025%
16	6.493	530	537	540	rVB2	2086	3554	0.32%	0.048%
17	6.951	611	615	617	rBV2	1265	1568	0.14%	0.021%
18	7.391	688	690	693	rBV2	1826	1586	0.14%	0.021%
19	7.568	717	720	722	rVB	1815	1655	0.15%	0.022%
20	7.644	729	733	735	rBV2	2036	2314	0.21%	0.031%
21	7.873	767	772	773	rBV	1333	1673	0.15%	0.023%
22	8.003	787	794	807	rBV	256237	423599	38.40%	5.702%
23	8.132	814	816	821	rBV3	1999	2252	0.20%	0.030%
24	8.379	855	858	860	rBV3	2622	3149	0.29%	0.042%
25	8.843	933	937	939	rBV2	1204	1622	0.15%	0.022%
26	8.872	939	942	947	rVB2	1010	1679	0.15%	0.023%
27	9.248	1002	1006	1009	rVB2	1792	2428	0.22%	0.033%
28	10.212	1166	1170	1172	rBV	1247	1626	0.15%	0.022%
29	10.500	1217	1219	1222	rBV	1936	1680	0.15%	0.023%
30	10.817	1266	1273	1284	rBV	297600	546002	49.50%	7.349%
31	10.887	1284	1285	1289	rVB2	2499	1815	0.16%	0.024%
32	13.044	1650	1652	1655	rBV	1206	1389	0.13%	0.019%
33	13.760	1772	1774	1776	rVB2	2493	1571	0.14%	0.021%
34	13.849	1786	1789	1793	rBV2	1032	1434	0.13%	0.019%

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 Title : SVOA CALIBRATION

35	14.089	1826	1830	1831	rVB	1263	1471	0.13%	0.020%
36	14.107	1831	1833	1836	rBV	1748	2210	0.20%	0.030%
37	14.642	1917	1924	1932	rBV	481016	747639	67.78%	10.063%
38	14.936	1971	1974	1976	rBV2	2207	2161	0.20%	0.029%
39	15.200	2017	2019	2022	rBV2	1263	1454	0.13%	0.020%
40	15.247	2025	2027	2032	rVV	1775	2681	0.24%	0.036%
41	15.306	2035	2037	2039	rVB2	1675	1496	0.14%	0.020%
42	15.782	2114	2118	2121	rBV2	1485	1897	0.17%	0.026%
43	16.023	2156	2159	2162	rBV3	1440	1993	0.18%	0.027%
44	16.346	2211	2214	2216	rBV2	2330	2050	0.19%	0.028%
45	16.446	2230	2231	2234	rBV2	1833	1570	0.14%	0.021%
46	16.904	2306	2309	2312	rBV	1067	1428	0.13%	0.019%
47	16.992	2322	2324	2327	rBV	1293	1473	0.13%	0.020%
48	17.139	2346	2349	2352	rBV	1807	2234	0.20%	0.030%
49	17.180	2352	2356	2359	rBV2	2007	3139	0.28%	0.042%
50	17.262	2365	2370	2372	rVB2	2398	3162	0.29%	0.043%
51	17.280	2372	2373	2376	rBV	1377	1479	0.13%	0.020%
52	17.386	2384	2391	2404	rBV	579267	927703	84.10%	12.487%
53	17.550	2418	2419	2424	rBV3	1912	2909	0.26%	0.039%
54	17.638	2432	2434	2437	rVB2	1980	1388	0.13%	0.019%
55	17.668	2437	2439	2441	rVV3	2134	1566	0.14%	0.021%
56	17.709	2444	2446	2449	rVB3	1894	1542	0.14%	0.021%
57	17.856	2469	2471	2475	rBV3	2142	3929	0.36%	0.053%
58	17.985	2487	2493	2500	rBV4	13057	24940	2.26%	0.336%
59	18.032	2500	2501	2503	rBV2	2315	1695	0.15%	0.023%
60	18.155	2518	2522	2524	rBV2	2940	3553	0.32%	0.048%
61	18.238	2533	2536	2540	rBV4	4382	7321	0.66%	0.099%
62	18.290	2543	2545	2547	rVV3	4363	3837	0.35%	0.052%
63	18.314	2547	2549	2553	rVV3	3189	4608	0.42%	0.062%
64	18.349	2553	2555	2559	rVB4	3657	4266	0.39%	0.057%
65	18.390	2559	2562	2563	rBV	2767	2672	0.24%	0.036%
66	18.449	2568	2572	2579	rBV	81534	111192	10.08%	1.497%
67	18.508	2579	2582	2586	rVB4	21986	30365	2.75%	0.409%
68	18.555	2586	2590	2594	rBV6	6777	9275	0.84%	0.125%
69	18.737	2616	2621	2625	rBV2	38753	53099	4.81%	0.715%
70	18.778	2625	2628	2633	rVB6	4653	7235	0.66%	0.097%
71	18.902	2647	2649	2656	rVB5	14644	21662	1.96%	0.292%

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72	18.972	2659	2661	2663	rBV3	2981	2508	0.23%	0.034%
73	19.048	2673	2674	2676	rBV2	2442	1979	0.18%	0.027%
74	19.131	2687	2688	2689	rBV	3595	1965	0.18%	0.026%
75	19.213	2699	2702	2705	rBV4	16662	20320	1.84%	0.274%
76	19.266	2707	2711	2712	rVV	50097	50690	4.60%	0.682%
77	19.295	2712	2716	2723	rVB2	126161	194019	17.59%	2.611%
78	19.348	2723	2725	2727	rBV3	5152	6601	0.60%	0.089%
79	19.377	2727	2730	2734	rVB4	18739	21068	1.91%	0.284%
80	19.436	2736	2740	2746	rVV2	27205	38687	3.51%	0.521%
81	19.495	2746	2750	2756	rVV7	26844	48954	4.44%	0.659%
82	19.571	2758	2763	2769	rVV	186877	255924	23.20%	3.445%
83	19.636	2770	2774	2778	rVB2	72408	86973	7.88%	1.171%
84	19.759	2793	2795	2799	rBV4	12730	18111	1.64%	0.244%
85	19.836	2804	2808	2813	rVB3	52953	72468	6.57%	0.975%
86	19.912	2816	2821	2825	rVV2	89576	118239	10.72%	1.591%
87	19.959	2825	2829	2832	rVV3	31532	39226	3.56%	0.528%
88	20.018	2834	2839	2845	rVB	189725	254837	23.10%	3.430%
89	20.141	2857	2860	2861	rBV2	14922	15436	1.40%	0.208%
90	20.276	2879	2883	2887	rBV3	82007	116392	10.55%	1.567%
91	20.329	2888	2892	2900	rVB2	167184	205788	18.66%	2.770%
92	20.553	2926	2930	2935	rVB	100656	125286	11.36%	1.686%
93	20.611	2936	2940	2942	rVV4	47185	70171	6.36%	0.944%
94	20.635	2942	2944	2949	rVV	73252	82985	7.52%	1.117%
95	20.711	2953	2957	2962	rVB7	28397	43624	3.95%	0.587%
96	20.852	2976	2981	2982	rBV3	50183	68033	6.17%	0.916%
97	20.876	2982	2985	2992	rVB2	119478	171686	15.56%	2.311%
98	21.205	3037	3041	3045	rBV4	35606	46425	4.21%	0.625%
99	21.645	3110	3116	3130	rVB	660700	1103074	100.00%	14.847%
100	24.842	3650	3660	3673	rVB2	362367	1060729	96.16%	14.277%

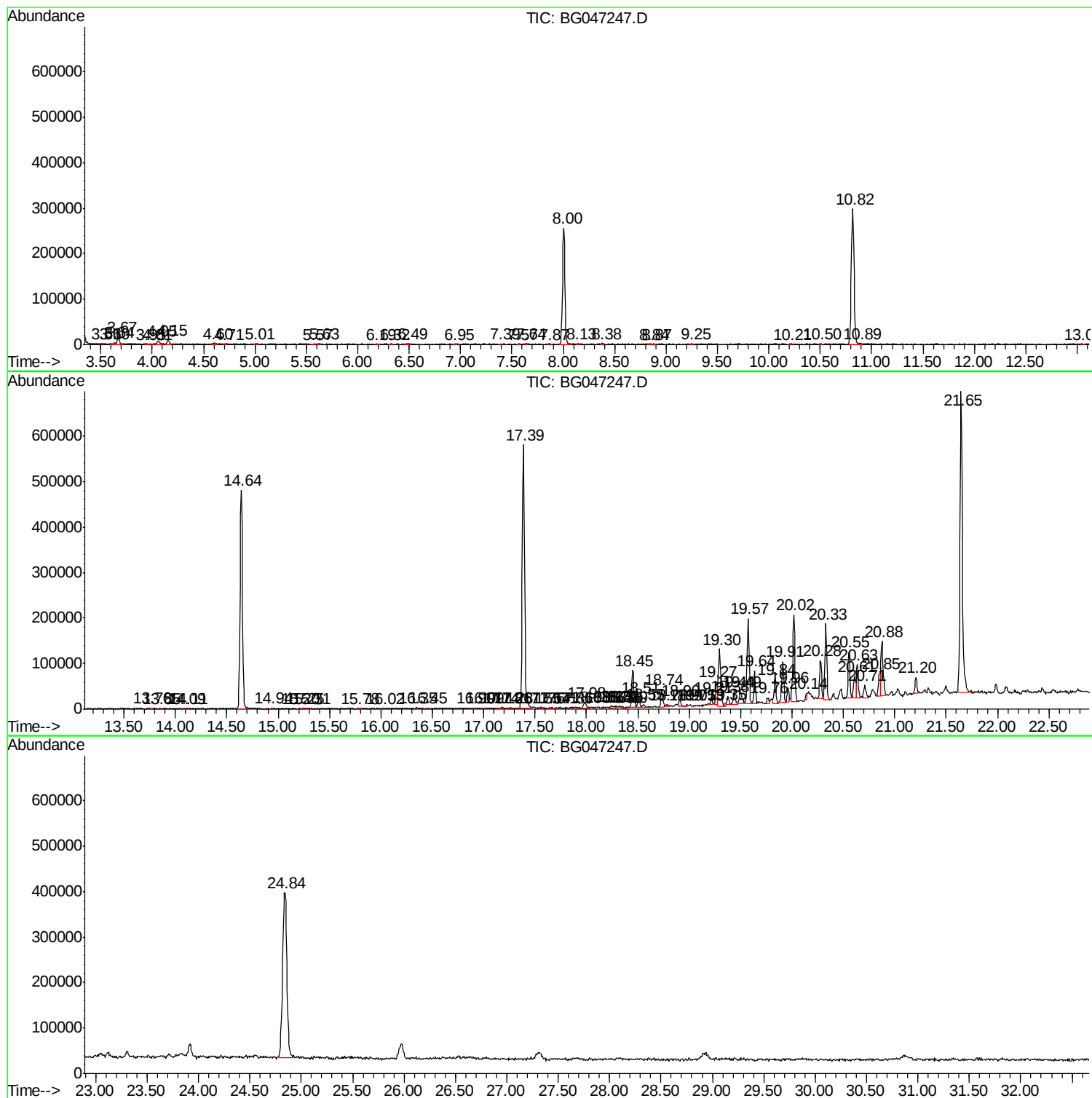
Sum of corrected areas: 7429519

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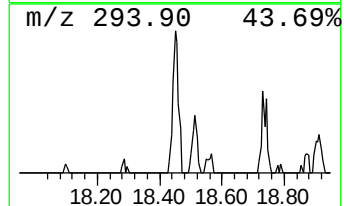
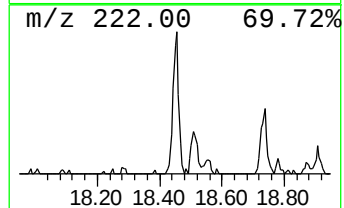
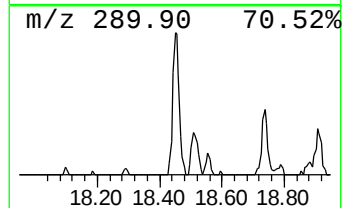
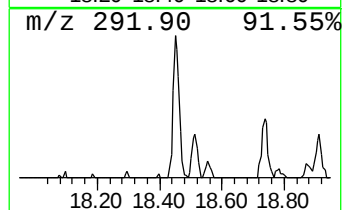
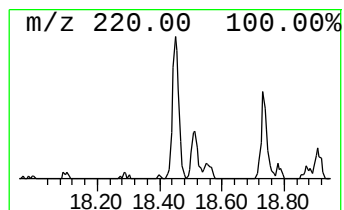
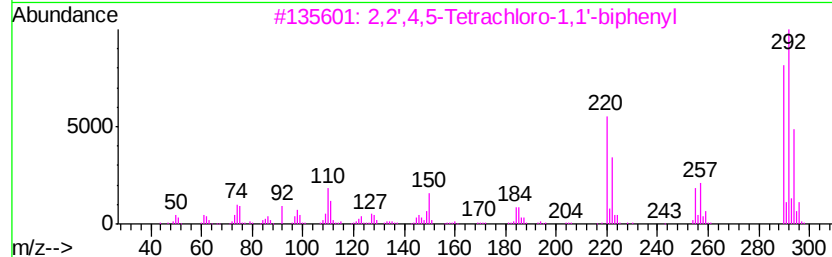
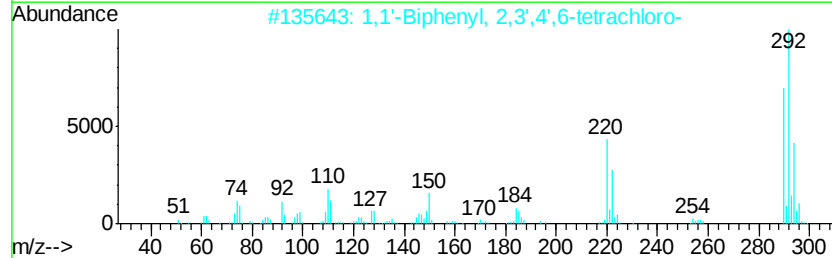
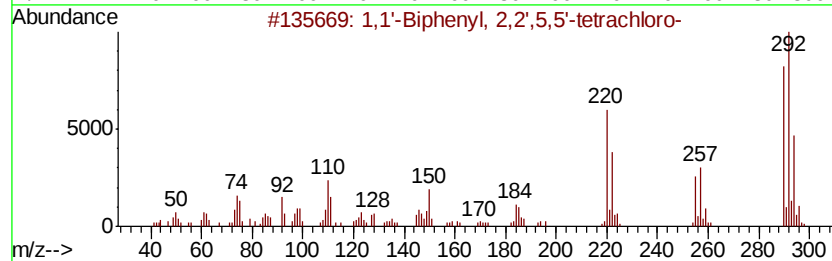
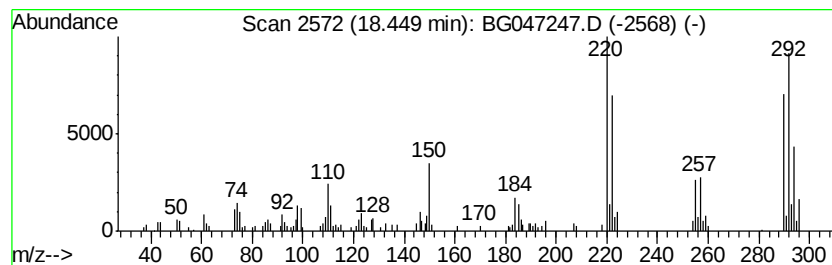
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.45	2.40 ng/ul	111192	Phenanthrene-d10	17.39	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5,5'-tetrachloro-1,1'-biphenyl	290	C12H6Cl4	035693-99-3	99
2	1,1'-Biphenyl, 2,3',4',6-tetrachloro-1,1'-biphenyl	290	C12H6Cl4	041464-46-4	99
3	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99
4	1,1'-Biphenyl, 2,3,3',5'-tetrachloro-1,1'-biphenyl	290	C12H6Cl4	041464-49-7	99
5	2,3',5',6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074338-23-1	99



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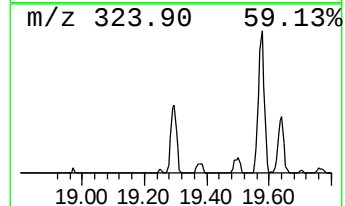
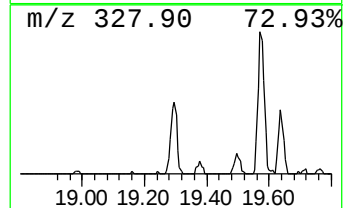
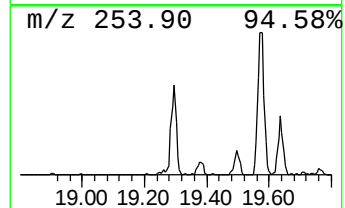
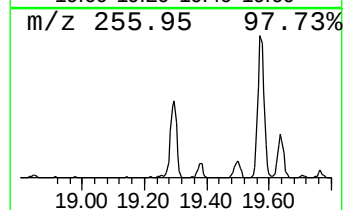
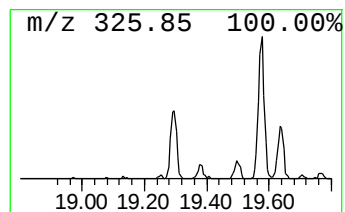
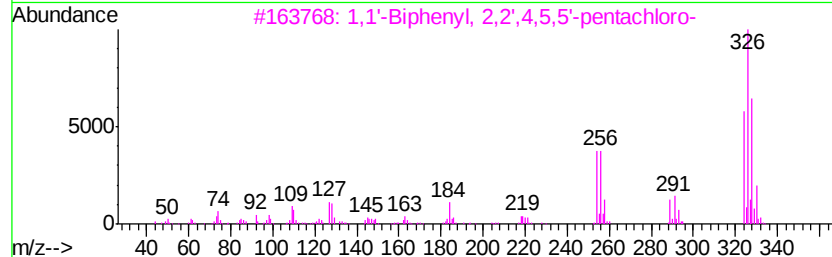
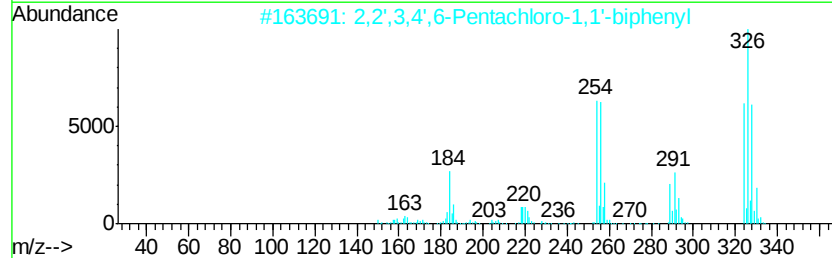
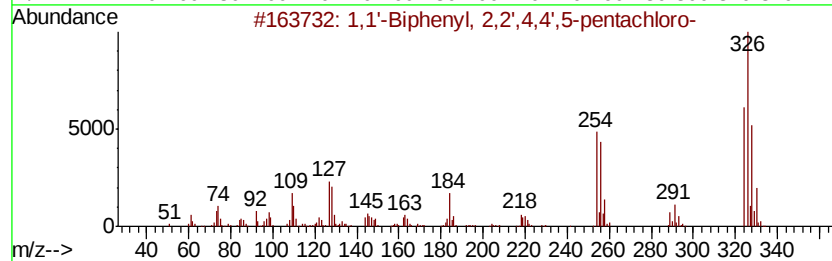
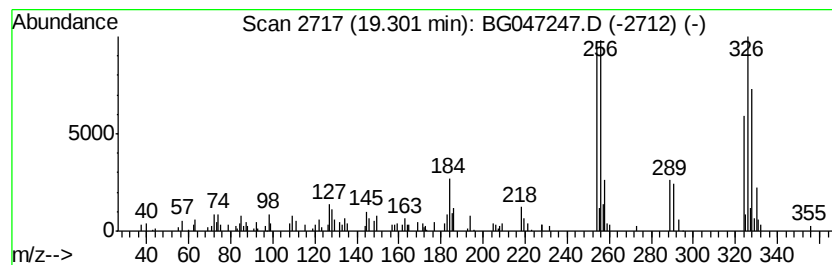
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
19.30	4.18 ng/ul	194019	Phenanthrene-d10	17.39		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,4',5-penta...	324	C12H5Cl5	038380-01-7	98	
2	2,2',3,4',6-Pentachloro-1,1'-bip...	324	C12H5Cl5	068194-05-8	98	
3	1,1'-Biphenyl, 2,2',4,5,5'-penta...	324	C12H5Cl5	037680-73-2	97	
4	1,1'-Biphenyl, 2,2',3',4,6-Penta...	324	C12H5Cl5	060233-25-2	97	
5	2,2',3,4,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-57-2	97	



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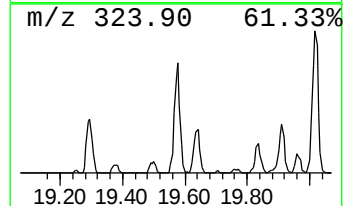
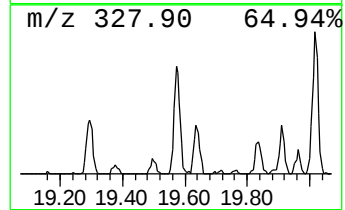
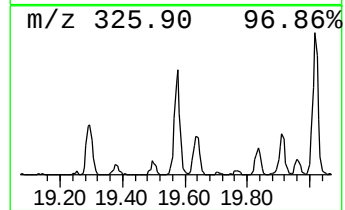
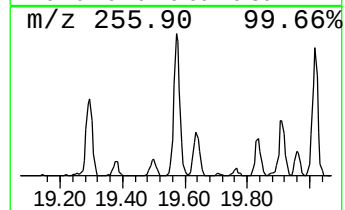
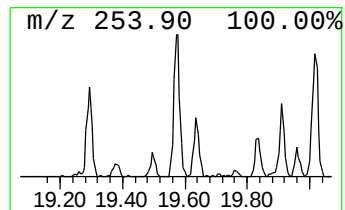
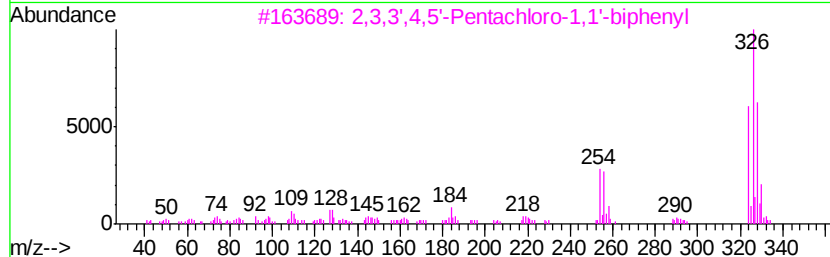
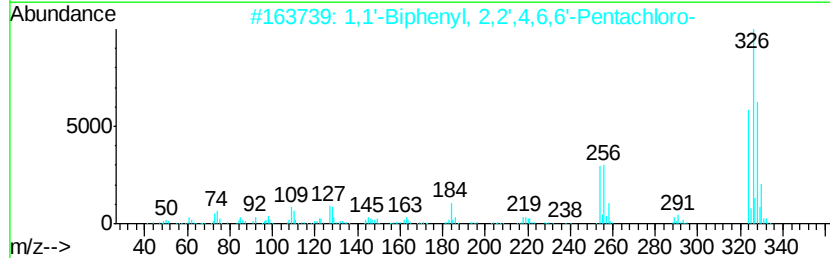
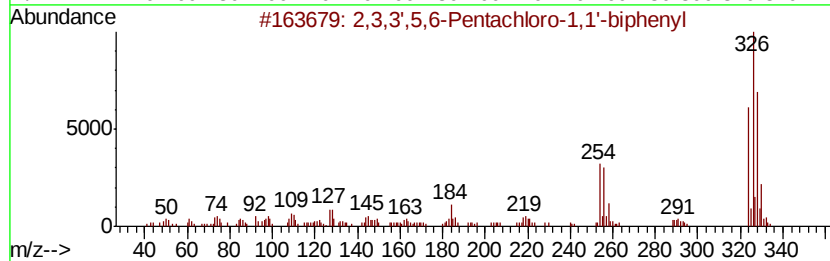
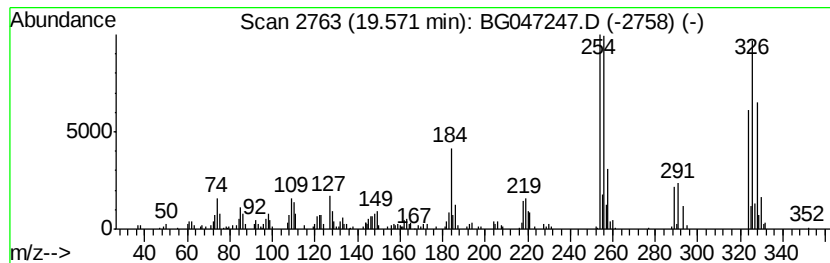
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.57	4.64 ng/ul	255924	Chrysene-d12	21.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,3,3',5,6-Pentachloro-1,1'-biph...	324	C12H5Cl5	074472-36-9	99
2		1,1'-Biphenyl, 2,2',4,6,6'-Penta...	324	C12H5Cl5	056558-16-8	98
3		2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	070362-41-3	96
4		2,2',3,4',5-Pentachloro-1,1'-bip...	324	C12H5Cl5	068194-07-0	96
5		1,1'-Biphenyl, 2,2',4,5,5'-penta...	324	C12H5Cl5	037680-73-2	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110620\
Data File : BG047247.D
Acq On : 6 Nov 2020 17:10
Operator : CG/JU
Sample : L4534-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0BP3

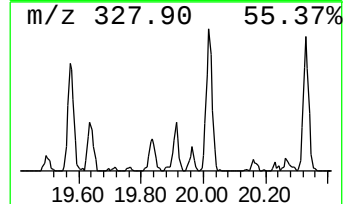
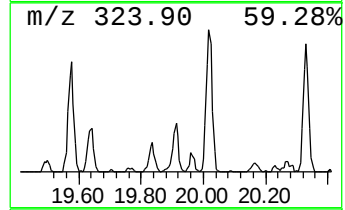
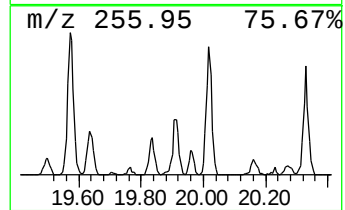
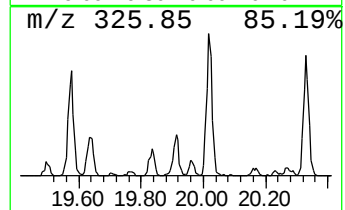
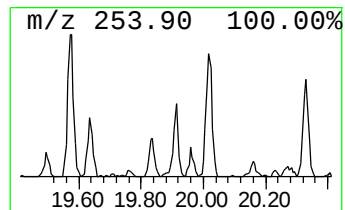
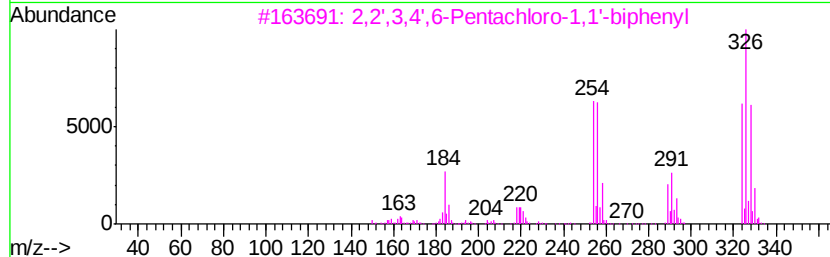
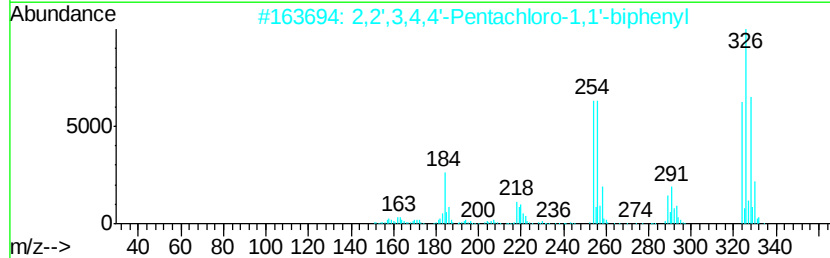
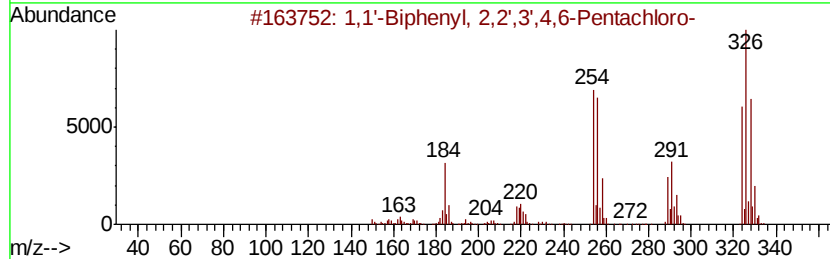
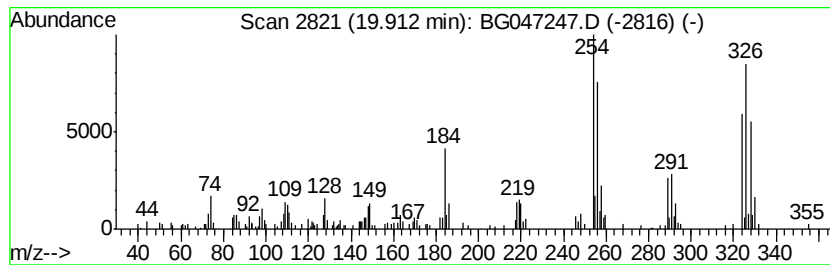
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1,1'-Biphenyl, 2,2',3',4,6-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.91	2.14 ng/ul	118239	Chrysene-d12	21.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1'-Biphenyl, 2,2',3',4,6-Penta...	324	C12H5Cl5	060233-25-2	99
2			2,2',3,4,4'-Pentachloro-1,1'-bip...	324	C12H5Cl5	065510-45-4	99
3			2,2',3,4',6-Pentachloro-1,1'-bip...	324	C12H5Cl5	068194-05-8	99
4			1,1'-Biphenyl, 2',3,4,5,5'-Penta...	324	C12H5Cl5	070424-70-3	99
5			2,2',3,6,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-54-9	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110620\
Data File : BG047247.D
Acq On : 6 Nov 2020 17:10
Operator : CG/JU
Sample : L4534-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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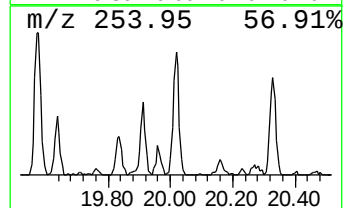
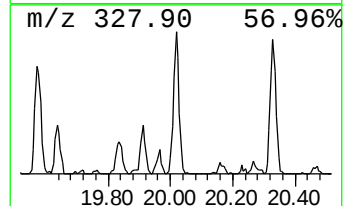
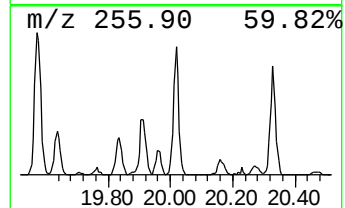
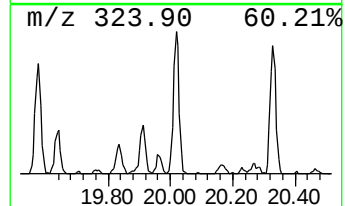
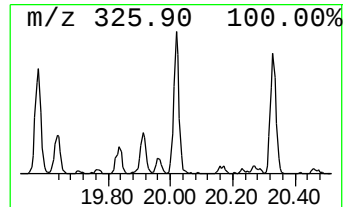
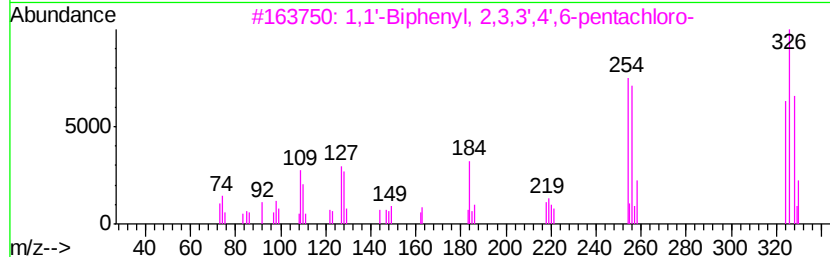
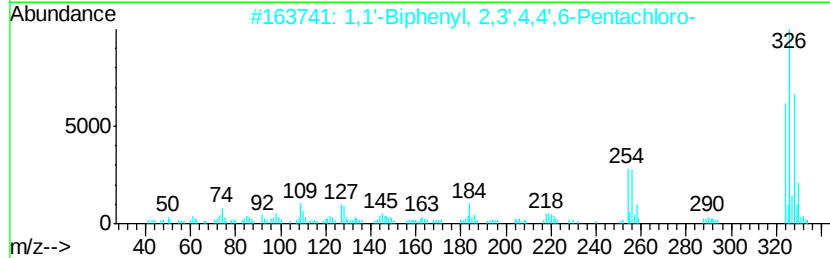
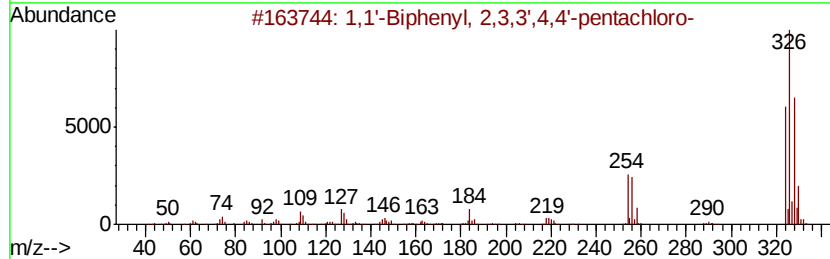
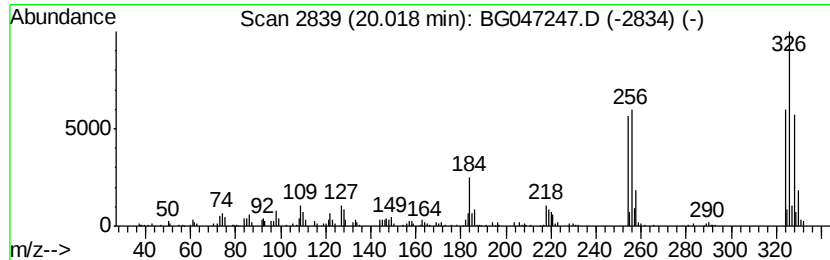
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.02	4.62 ng/ul	254837	Chrysene-d12	21.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'	-Biphenyl,	2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	99
2	1,1'	-Biphenyl,	2,3',4,4',6-Penta...	324	C12H5Cl5	056558-17-9	99
3	1,1'	-Biphenyl,	2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99
4	2,3,3'	,4,5-Pentachloro-1,1'-biph...		324	C12H5Cl5	070424-69-0	99
5	2,3,3'	,4',5'- Pentachloro-1,1'-b...		324	C12H5Cl5	076842-07-4	97



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110620\
Data File : BG047247.D
Acq On : 6 Nov 2020 17:10
Operator : CG/JU
Sample : L4534-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

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ClientSampleId :
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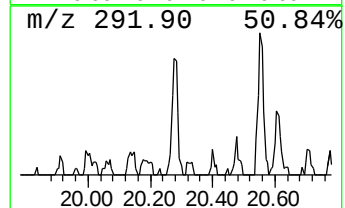
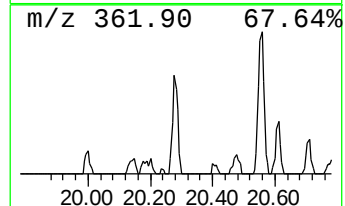
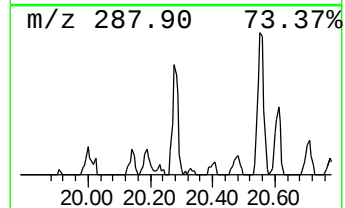
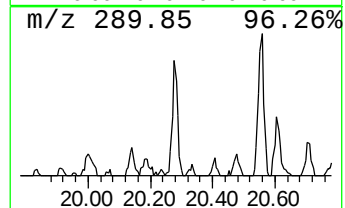
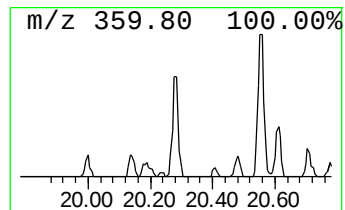
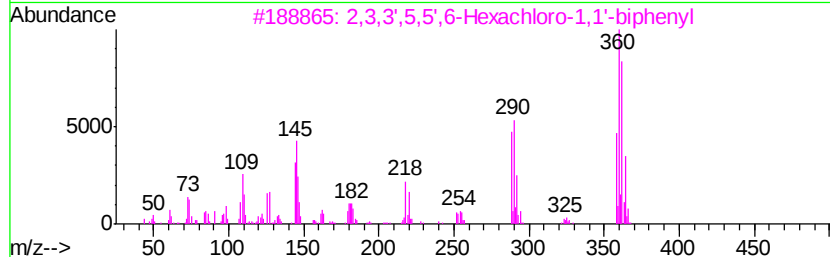
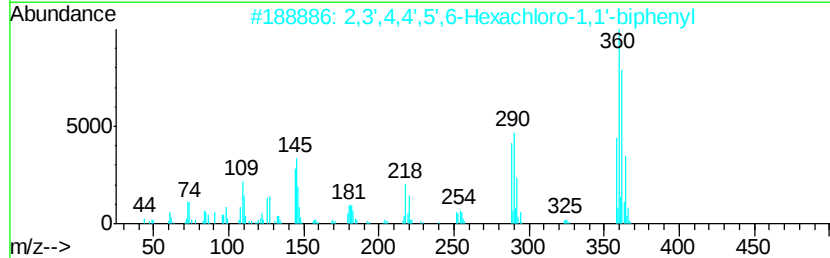
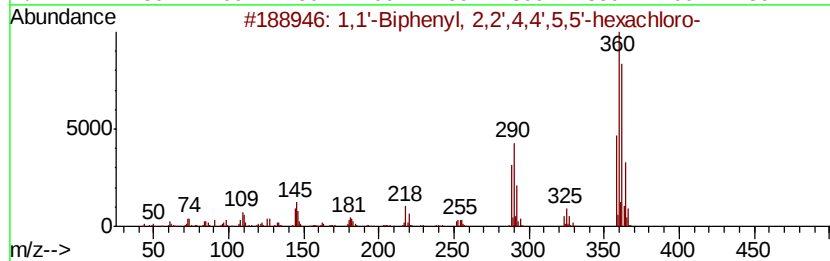
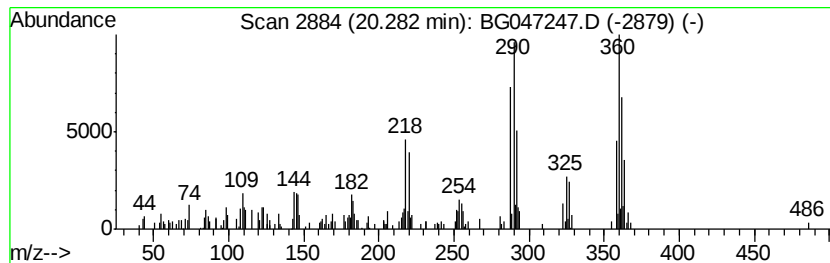
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.28	2.11 ng/ul	116392	Chrysene-d12	21.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',4,4',5,5'-he...	358	C12H4Cl6	035065-27-1	99
2		2,3',4,4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-65-5	99
3		2,3,3',5,5',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-46-1	99
4		2,3,3',4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	074472-45-0	99
5		1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110620\
Data File : BG047247.D
Acq On : 6 Nov 2020 17:10
Operator : CG/JU
Sample : L4534-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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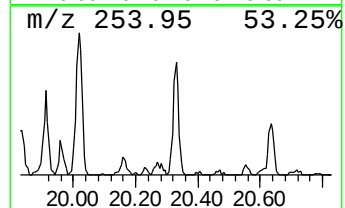
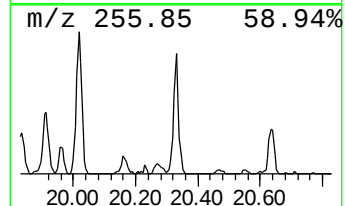
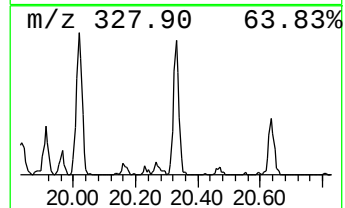
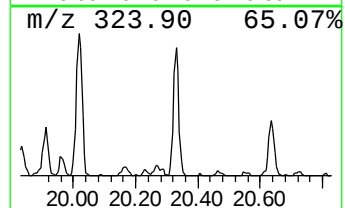
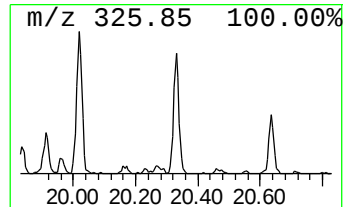
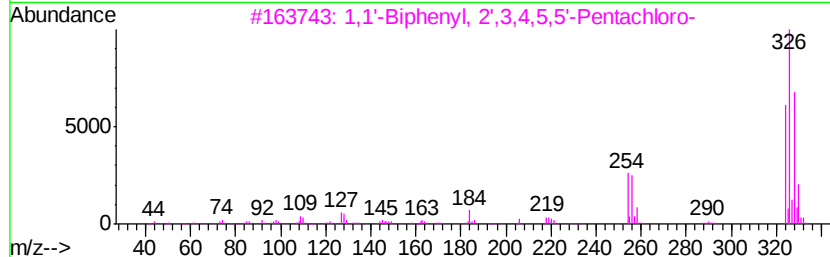
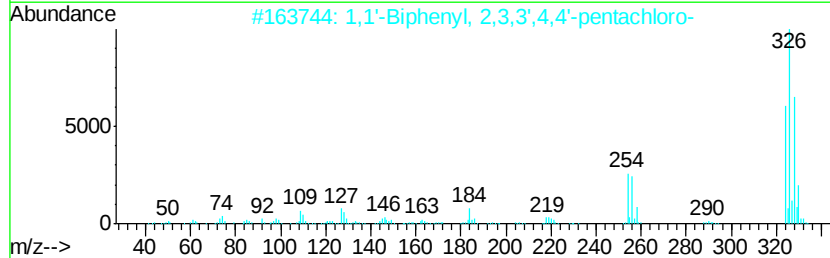
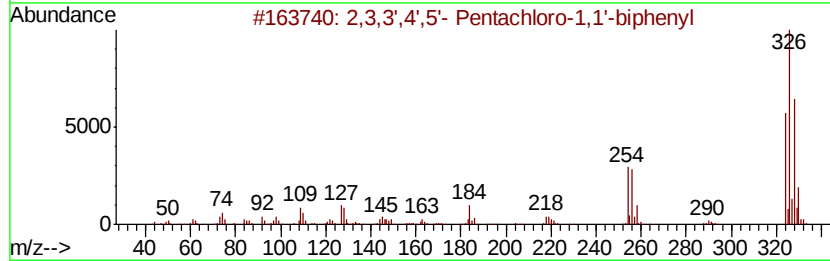
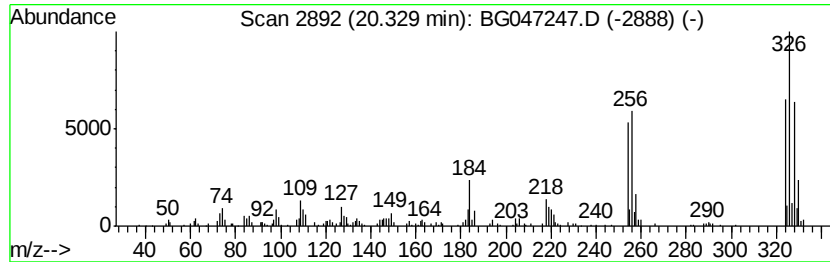
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.33	3.73 ng/ul	205788	Chrysene-d12	21.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,3,3',4',5'- Pentachloro-1,1'-b...	324	C12H5Cl5	076842-07-4	98
2		1,1'-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	97
3		1,1'-Biphenyl, 2',3,4,5,5'-Penta...	324	C12H5Cl5	070424-70-3	96
4		2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	070362-41-3	96
5		2,2',3,4,4'-Pentachloro-1,1'-bip...	324	C12H5Cl5	065510-45-4	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110620\
Data File : BG047247.D
Acq On : 6 Nov 2020 17:10
Operator : CG/JU
Sample : L4534-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

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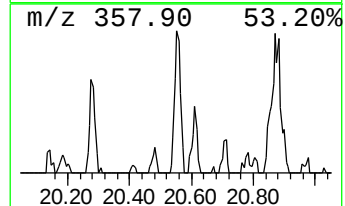
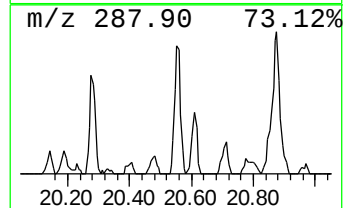
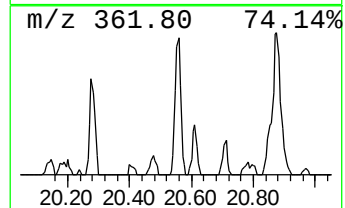
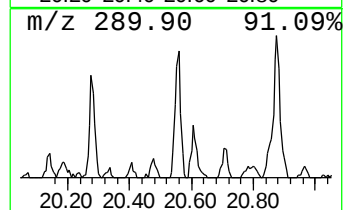
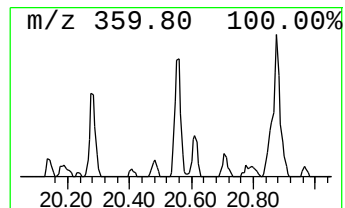
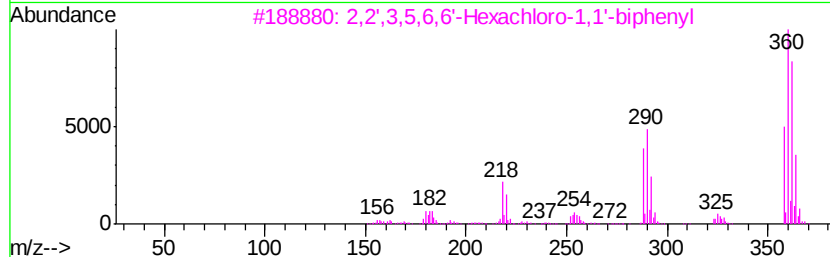
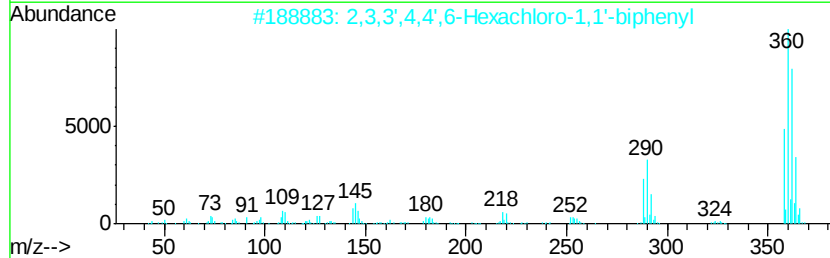
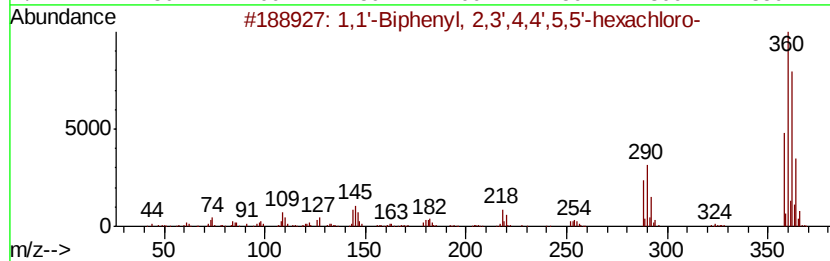
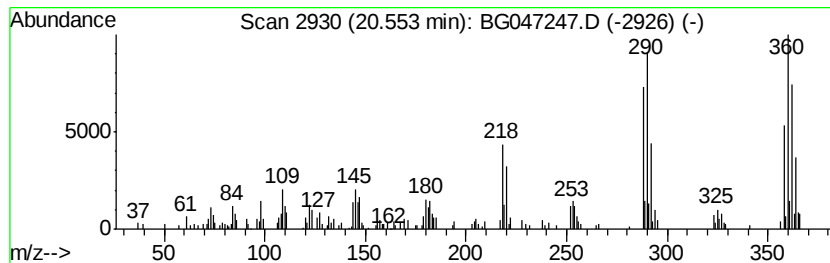
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.55	2.27 ng/ul	125286	Chrysene-d12	21.65

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		2,3,3',4,4',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-42-7	99
3		2,2',3,5,6,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-09-2	99
4		2,3,4,4',5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-63-6	99
5		1,1'-Biphenyl, 2,2',4,4',6,6'-he...	358	C12H4Cl6	033979-03-2	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110620\
Data File : BG047247.D
Acq On : 6 Nov 2020 17:10
Operator : CG/JU
Sample : L4534-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

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ClientSampleId :
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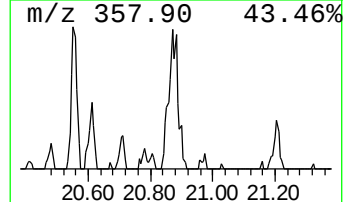
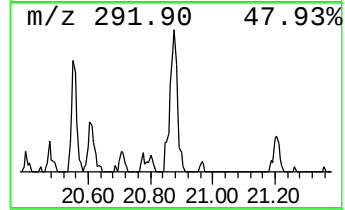
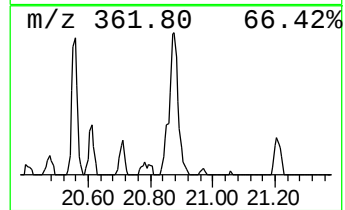
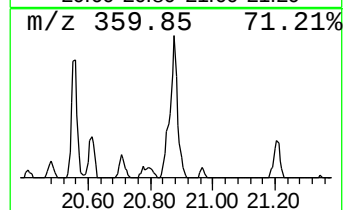
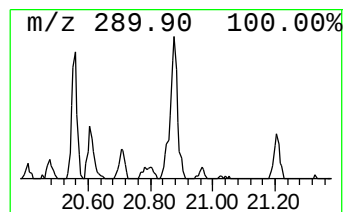
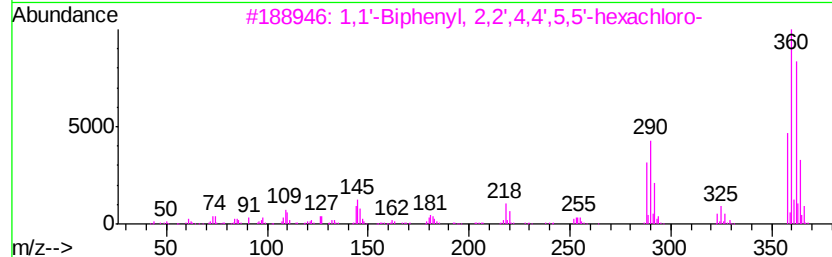
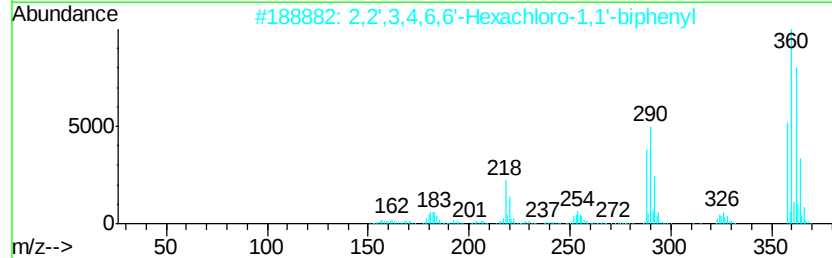
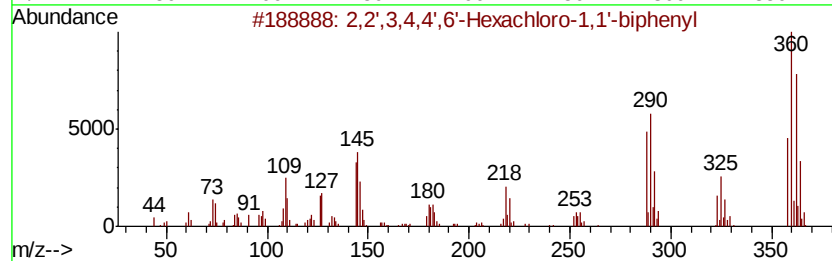
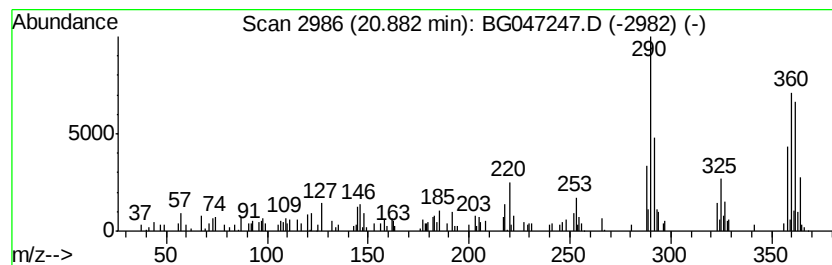
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.88	3.11 ng/ul	171686	Chrysene-d12	21.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2',3,4,4',6'-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-64-4	93
2		2,2',3,4,6,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-40-5	93
3		1,1'-Biphenyl, 2,2',4,4',5,5'-he...	358	C12H4Cl6	035065-27-1	93
4		1,1'-Biphenyl, 2,2',3,3',4,5-hex...	358	C12H4Cl6	055215-18-4	92
5		2,3,3',5,5',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-46-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG110620\
 Data File : BG047247.D
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 Operator : CG/JU
 Sample : L4534-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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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	2.4	ng/ul	111192	4	17.39	927703	20.0
1,1'-Biphenyl, 2,...	19.30	4.2	ng/ul	194019	4	17.39	927703	20.0
2,3,3',5,6-Pentac...	19.57	4.6	ng/ul	255924	5	21.65	1103070	20.0
1,1'-Biphenyl, 2,...	19.91	2.1	ng/ul	118239	5	21.65	1103070	20.0
1,1'-Biphenyl, 2,...	20.02	4.6	ng/ul	254837	5	21.65	1103070	20.0
1,1'-Biphenyl, 2,...	20.28	2.1	ng/ul	116392	5	21.65	1103070	20.0
2,3,3',4',5'- Pen...	20.33	3.7	ng/ul	205788	5	21.65	1103070	20.0
1,1'-Biphenyl, 2,...	20.55	2.3	ng/ul	125286	5	21.65	1103070	20.0
2,2',3,4,4',6'-He...	20.88	3.1	ng/ul	171686	5	21.65	1103070	20.0