

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050423\
 Data File : BG057299.D
 Acq On : 4 May 2023 11:55
 Operator : CG/JU
 Sample : 02419-36DL 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW931DL

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG057299.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.177	277	279	284	rBV3	674	776	0.23%	0.021%
2	5.747	373	376	378	rBV	502	758	0.22%	0.020%
3	5.806	384	386	389	rBV3	435	536	0.16%	0.014%
4	6.088	430	434	436	rVB2	416	499	0.15%	0.013%
5	6.293	467	469	472	rBV	679	583	0.17%	0.016%
6	6.845	561	563	565	rBV	376	479	0.14%	0.013%
7	6.928	575	577	582	rVB2	442	727	0.21%	0.020%
8	7.444	661	665	666	rBV	570	701	0.21%	0.019%
9	7.521	673	678	686	rBV4	3596	7034	2.08%	0.189%
10	7.674	701	704	710	rBV2	3255	5097	1.51%	0.137%
11	7.832	727	731	734	rVB	390	744	0.22%	0.020%
12	7.903	738	743	747	rBV4	3641	5465	1.61%	0.147%
13	8.373	818	823	830	rVB	43553	72716	21.47%	1.951%
14	9.078	940	943	948	rBV2	2880	4495	1.33%	0.121%
15	9.207	959	965	972	rBV2	3319	6583	1.94%	0.177%
16	9.553	1018	1024	1030	rBV2	3974	6825	2.02%	0.183%
17	10.282	1144	1148	1153	rBV2	3076	4617	1.36%	0.124%
18	10.828	1238	1241	1251	rVB2	3686	7158	2.11%	0.192%
19	11.216	1300	1307	1314	rVB	69061	119672	35.34%	3.210%
20	11.345	1323	1329	1332	rBV2	1232	2065	0.61%	0.055%
21	12.308	1489	1493	1496	rBV	315	620	0.18%	0.017%
22	12.837	1580	1583	1586	rBV	477	628	0.19%	0.017%
23	13.431	1678	1684	1685	rBV	305	687	0.20%	0.018%
24	13.777	1741	1743	1747	rVB	868	716	0.21%	0.019%
25	13.848	1750	1755	1760	rVB2	2499	3592	1.06%	0.096%
26	14.370	1840	1844	1850	rBV	14961	20660	6.10%	0.554%
27	14.423	1850	1853	1855	rVB	632	564	0.17%	0.015%
28	14.688	1893	1898	1904	rBV	14722	22452	6.63%	0.602%
29	14.834	1919	1923	1926	rBV3	1243	1773	0.52%	0.048%
30	14.993	1941	1950	1956	rBV	139512	218718	64.58%	5.867%
31	15.052	1958	1960	1962	rBV2	745	606	0.18%	0.016%
32	15.193	1980	1984	1996	rVB4	2226	4121	1.22%	0.111%
33	15.369	2011	2014	2016	rBV	717	732	0.22%	0.020%
34	15.393	2016	2018	2019	rVB	1014	571	0.17%	0.015%
35	15.457	2024	2029	2031	rVB	601	745	0.22%	0.020%
36	15.581	2048	2050	2053	rBV	482	522	0.15%	0.014%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
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37	15.628	2056	2058	2061	rBV	411	470	0.14%	0.013%
38	15.704	2068	2071	2073	rBV	607	603	0.18%	0.016%
39	15.739	2073	2077	2079	rVB	913	1334	0.39%	0.036%
40	15.774	2079	2083	2088	rBV5	2808	4424	1.31%	0.119%

41	15.839	2090	2094	2095	rVV2	971	929	0.27%	0.025%
42	15.880	2099	2101	2105	rVB	423	488	0.14%	0.013%
43	15.974	2111	2117	2125	rBV	23288	34974	10.33%	0.938%
44	16.027	2125	2126	2129	rBV2	890	911	0.27%	0.024%
45	16.068	2129	2133	2135	rVB3	668	732	0.22%	0.020%

46	16.097	2135	2138	2142	rBV2	2605	3289	0.97%	0.088%
47	16.180	2144	2152	2157	rBV5	2009	4490	1.33%	0.120%
48	16.221	2157	2159	2162	rBV2	3051	3458	1.02%	0.093%
49	16.280	2167	2169	2170	rBV	674	623	0.18%	0.017%
50	16.327	2172	2177	2183	rBV2	8828	12132	3.58%	0.325%

51	16.368	2183	2184	2185	rVB	1533	540	0.16%	0.014%
52	16.403	2185	2190	2191	rBV2	1244	1339	0.40%	0.036%
53	16.421	2191	2193	2195	rVB	1038	694	0.20%	0.019%
54	16.462	2199	2200	2201	rBV	1913	1050	0.31%	0.028%
55	16.509	2207	2208	2210	rBV2	801	641	0.19%	0.017%

56	16.532	2210	2212	2213	rVB2	1206	705	0.21%	0.019%
57	16.579	2218	2220	2223	rVV3	1974	2260	0.67%	0.061%
58	16.661	2227	2234	2240	rVV5	5616	13510	3.99%	0.362%
59	16.703	2240	2241	2243	rVV2	1474	905	0.27%	0.024%
60	16.738	2246	2247	2248	rBV	1875	681	0.20%	0.018%

61	16.773	2248	2253	2255	rBV6	1659	2444	0.72%	0.066%
62	16.826	2260	2262	2265	rBV3	1348	1251	0.37%	0.034%
63	16.885	2270	2272	2273	rBV	1044	611	0.18%	0.016%
64	16.943	2278	2282	2288	rVB	7207	10604	3.13%	0.284%
65	17.043	2297	2299	2300	rVB2	1678	840	0.25%	0.023%

66	17.079	2303	2305	2306	rBV2	1193	790	0.23%	0.021%
67	17.478	2371	2373	2376	rVB3	5920	5992	1.77%	0.161%
68	17.590	2388	2392	2395	rBV2	21710	29562	8.73%	0.793%
69	17.625	2396	2398	2403	rVB3	10971	12773	3.77%	0.343%
70	17.731	2406	2416	2421	rBV	204973	338666	100.00%	9.085%

71	17.772	2421	2423	2428	rVV	16847	19054	5.63%	0.511%
72	17.830	2428	2433	2437	rVV	24734	36961	10.91%	0.991%
73	17.889	2438	2443	2450	rVB3	35413	52392	15.47%	1.405%
74	18.165	2485	2490	2493	rBV5	13215	19101	5.64%	0.512%
75	18.312	2508	2515	2525	rVB2	120711	264423	78.08%	7.093%

76	18.435	2531	2536	2541	rVB4	22379	43247	12.77%	1.160%
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 Title : SVOA CALIBRATION

77	18.565	2553	2558	2562	rBV	36774	57392	16.95%	1.540%
78	18.647	2569	2572	2575	rVB	15300	17144	5.06%	0.460%
79	18.770	2588	2593	2598	rBV	113504	157318	46.45%	4.220%
80	18.829	2600	2603	2607	rVV2	53003	76310	22.53%	2.047%
81	18.876	2607	2611	2617	rVB4	30998	45381	13.40%	1.217%
82	19.052	2636	2641	2645	rBV	77803	110657	32.67%	2.968%
83	19.152	2656	2658	2662	rBV	24296	26655	7.87%	0.715%
84	19.229	2667	2671	2677	rVB	50218	63633	18.79%	1.707%
85	19.528	2719	2722	2726	rBV	32794	43469	12.84%	1.166%
86	19.581	2726	2731	2733	rVV	87475	137328	40.55%	3.684%
87	19.604	2733	2735	2742	rVB3	106633	165896	48.99%	4.450%
88	19.763	2758	2762	2766	rVB	42295	54686	16.15%	1.467%
89	19.881	2778	2782	2789	rVB	135701	180991	53.44%	4.855%
90	19.939	2789	2792	2797	rBV4	43295	62917	18.58%	1.688%
91	20.215	2836	2839	2844	rVB2	56880	70315	20.76%	1.886%
92	20.321	2854	2857	2863	rVB	108950	142077	41.95%	3.811%
93	20.633	2906	2910	2915	rVB	102401	130717	38.60%	3.506%
94	20.856	2945	2948	2954	rVB4	52320	69324	20.47%	1.860%
95	21.713	3090	3094	3098	rVB	102136	145085	42.84%	3.892%
96	21.866	3116	3120	3126	rVB	154270	218543	64.53%	5.862%
97	22.048	3146	3151	3156	rVB2	166410	296924	87.67%	7.965%

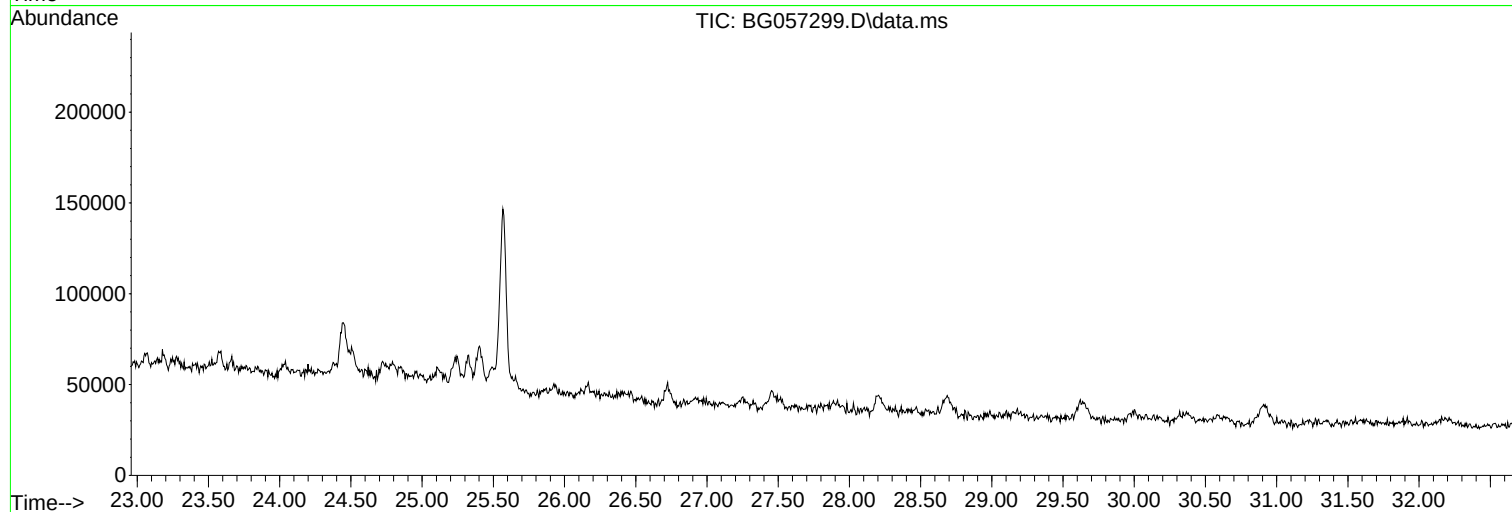
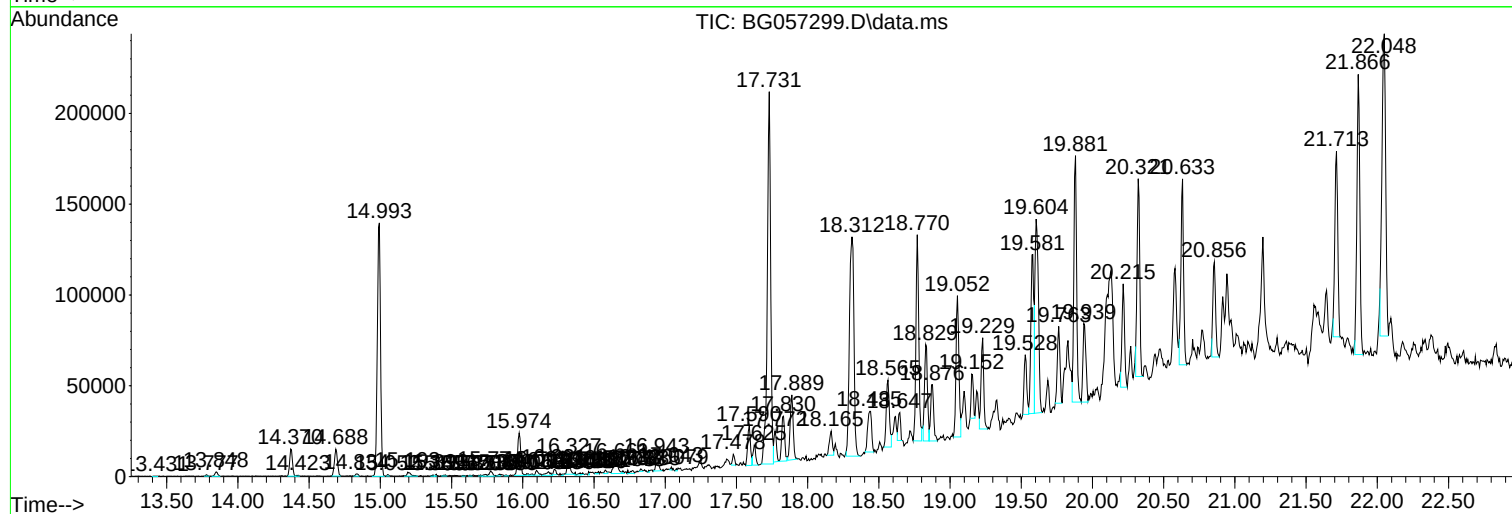
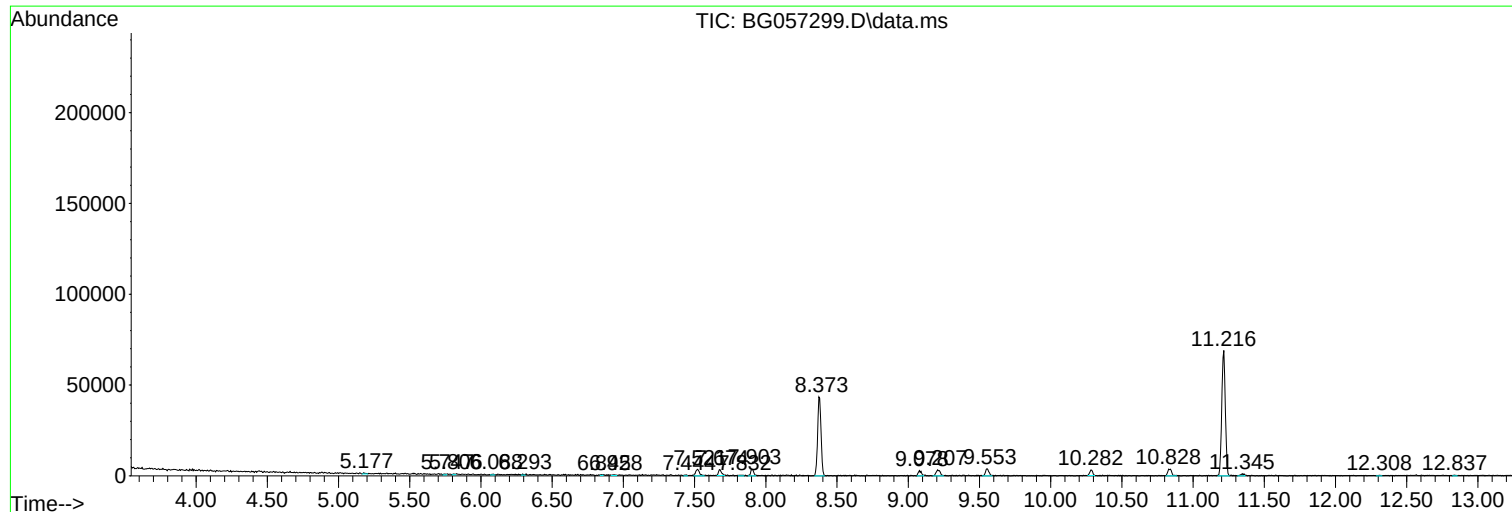
Sum of corrected areas: 3727892

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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Instrument :
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EW931DL

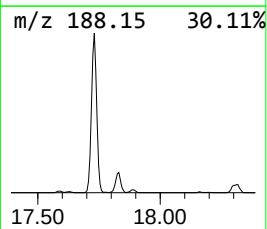
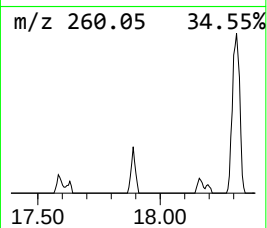
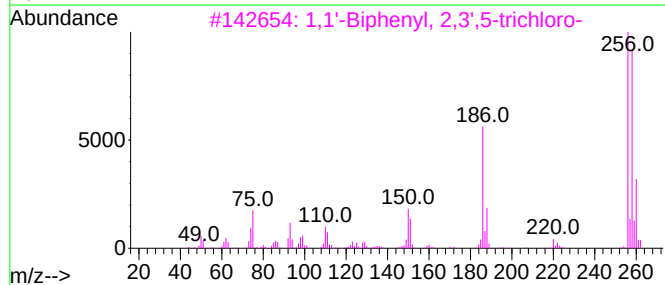
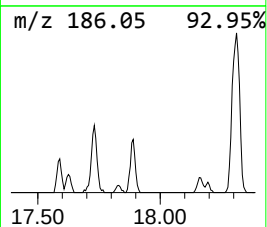
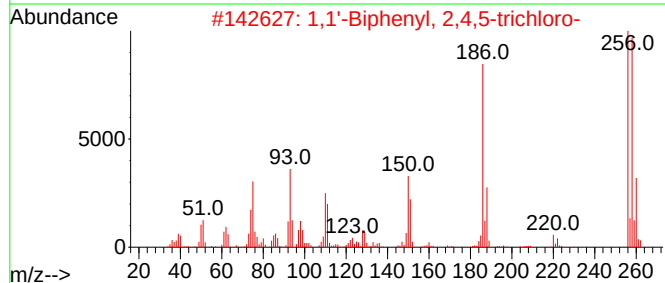
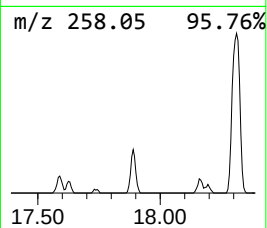
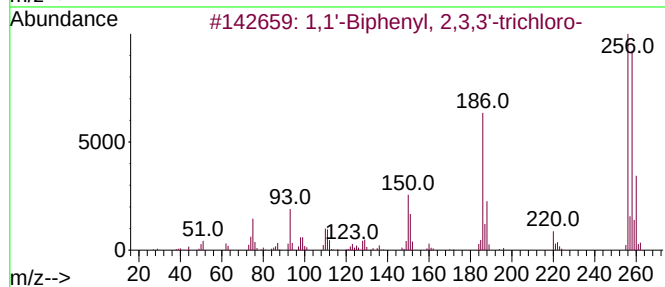
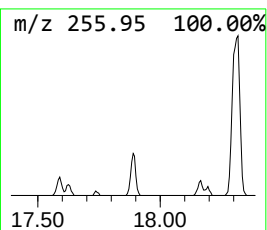
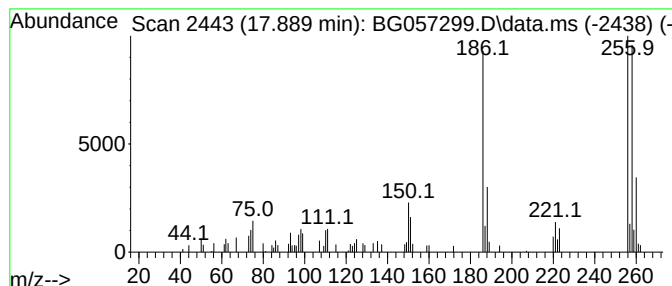
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.889	3.09 ng/ul	52392	Phenanthrene-d10	17.731

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99		
2	1,1'-Biphenyl, 2,4,5-trichloro-	256	C12H7Cl3	015862-07-4	99		
3	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	98		
4	1,1'-Biphenyl, 2',3,5-trichloro-	256	C12H7Cl3	037680-68-5	98		
5	1,1'-Biphenyl, 2,3,6-trichloro-	256	C12H7Cl3	055702-45-9	98		



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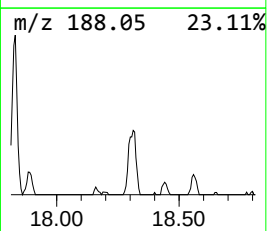
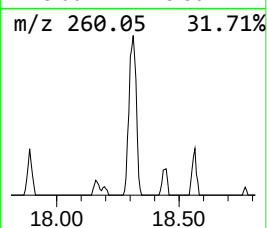
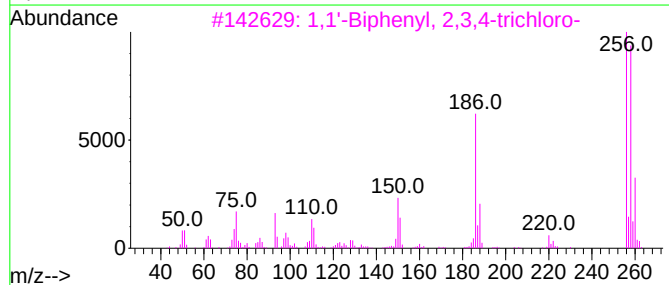
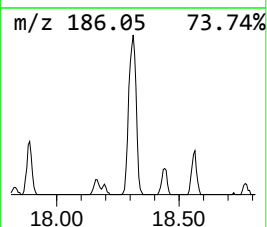
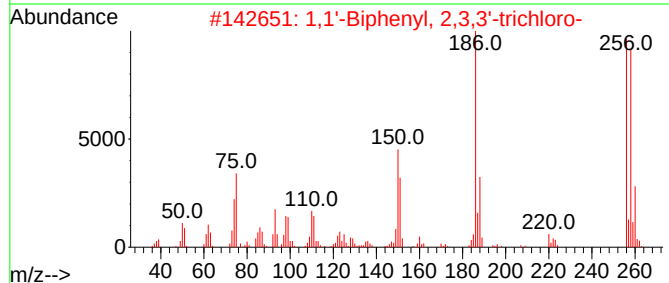
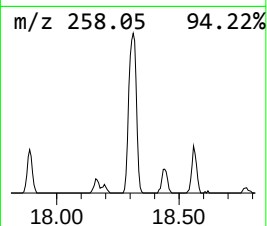
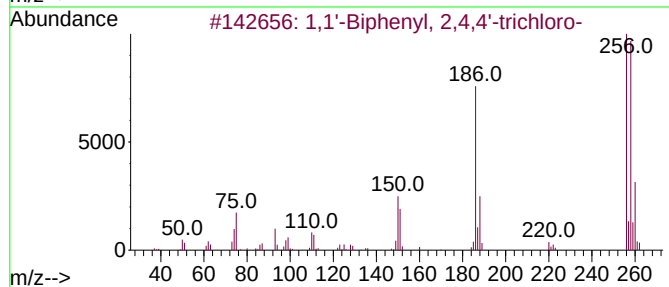
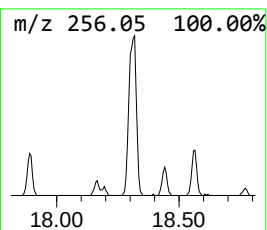
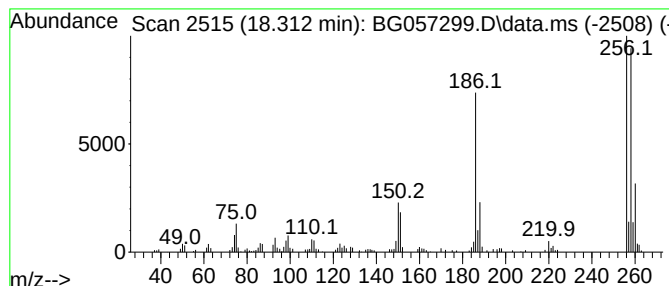
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.312	15.62 ng/ul	264423	Phenanthrene-d10	17.731

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99		
2	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99		
3	1,1'-Biphenyl, 2,3,4-trichloro-	256	C12H7Cl3	055702-46-0	98		
4	1,1'-Biphenyl, 2',3,5-trichloro-	256	C12H7Cl3	037680-68-5	98		
5	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	97		



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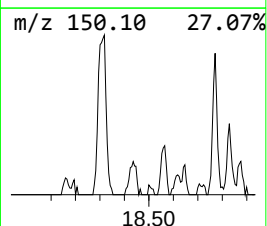
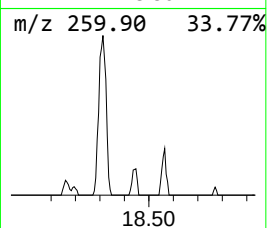
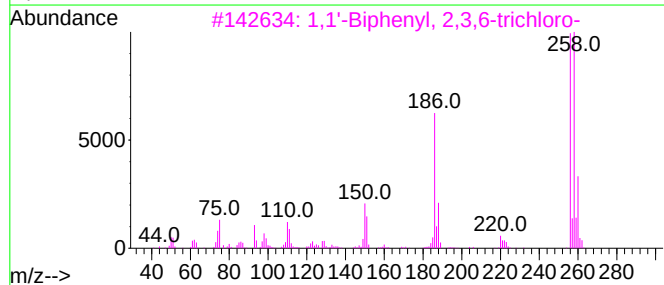
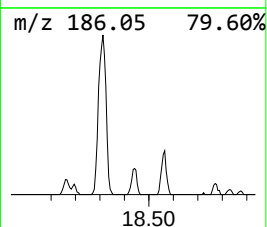
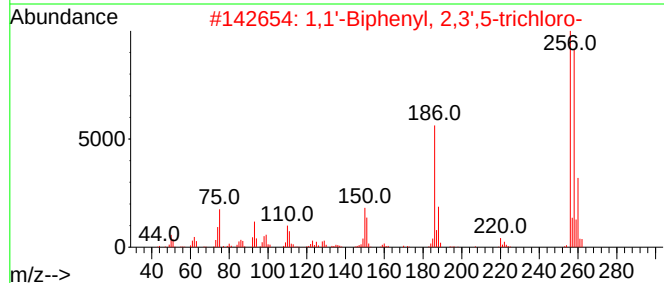
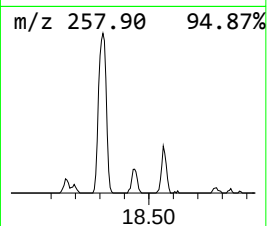
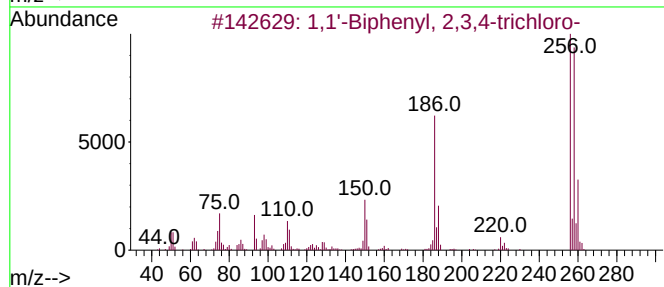
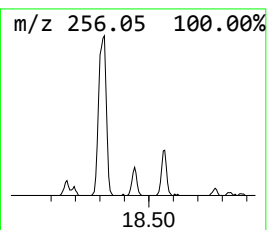
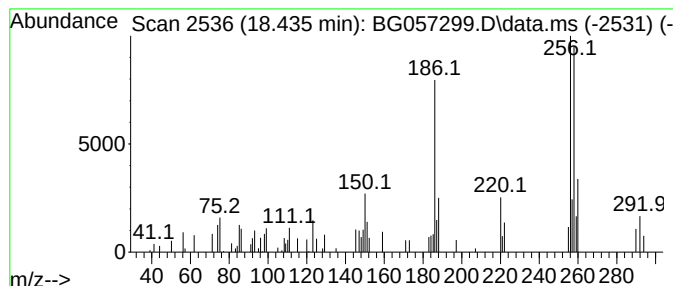
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.436	2.55 ng/ul	43247	Phenanthrene-d10	17.731

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,4-trichloro-	256	C12H7Cl3	055702-46-0	93		
2	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	93		
3	1,1'-Biphenyl, 2,3,6-trichloro-	256	C12H7Cl3	055702-45-9	93		
4	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	93		
5	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	93		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050423\
Data File : BG057299.D
Acq On : 4 May 2023 11:55
Operator : CG/JU
Sample : 02419-36DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW931DL

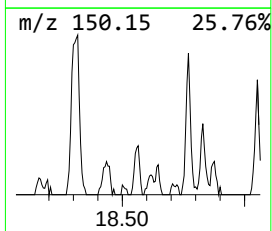
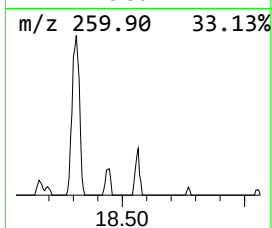
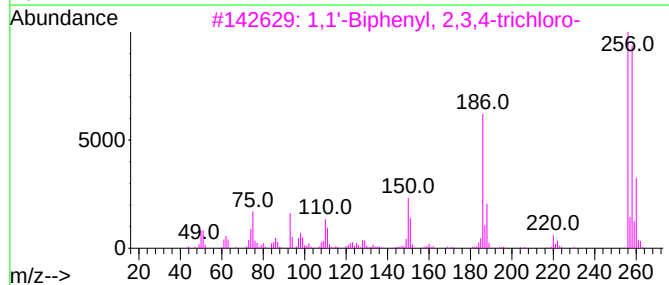
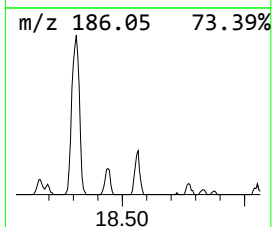
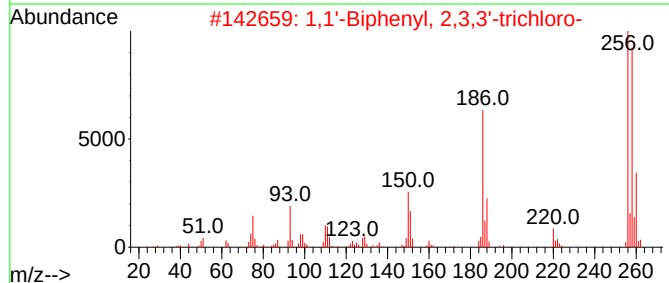
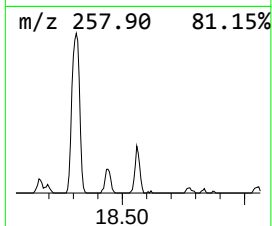
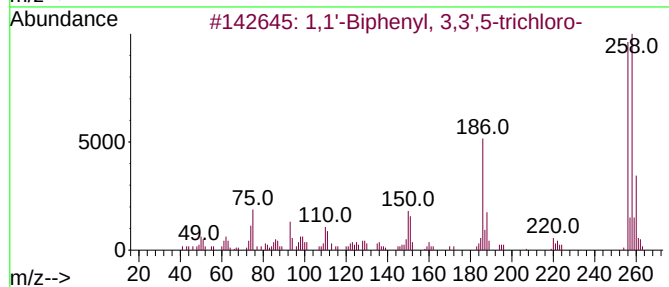
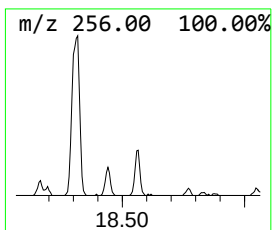
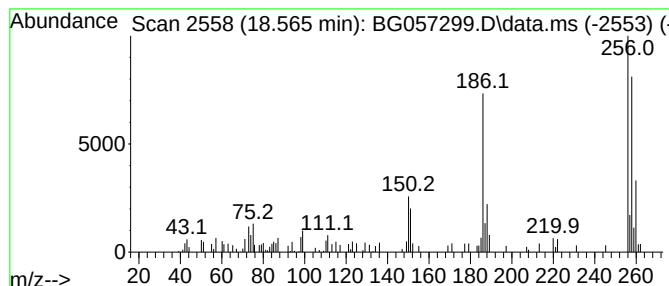
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1,1'-Biphenyl, 3,3',5-trich... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.565	3.39 ng/ul	57392	Phenanthrene-d10	17.731

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	97		
2	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	97		
3	1,1'-Biphenyl, 2,3,4-trichloro-	256	C12H7Cl3	055702-46-0	96		
4	3,4,5-Trichloro-1,1'-biphenyl	256	C12H7Cl3	053555-66-1	96		
5	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	96		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050423\
Data File : BG057299.D
Acq On : 4 May 2023 11:55
Operator : CG/JU
Sample : 02419-36DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

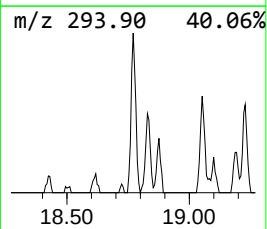
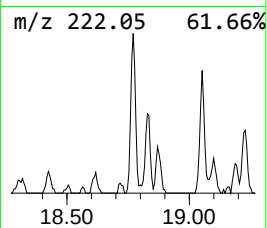
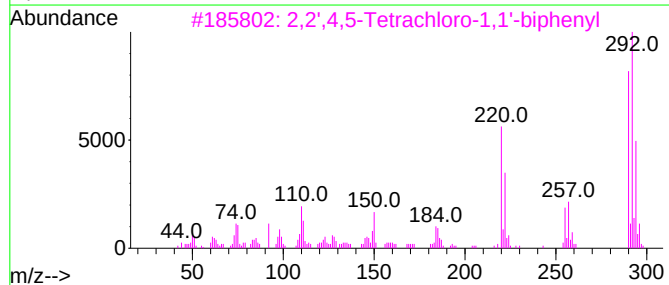
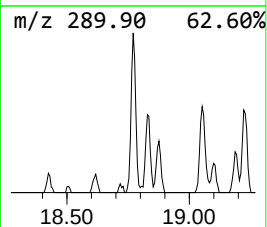
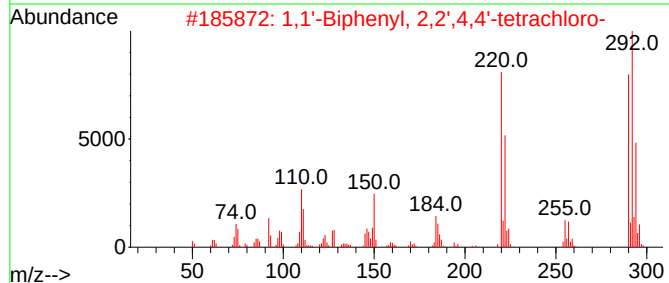
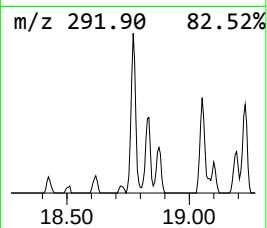
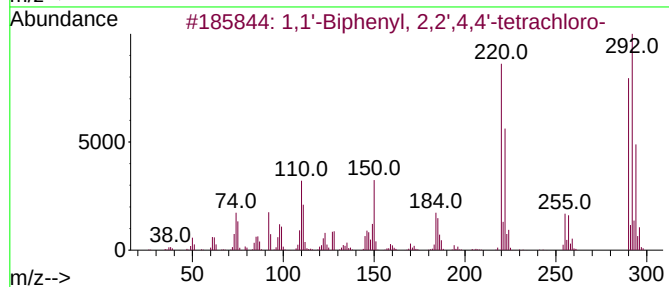
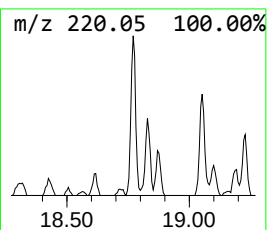
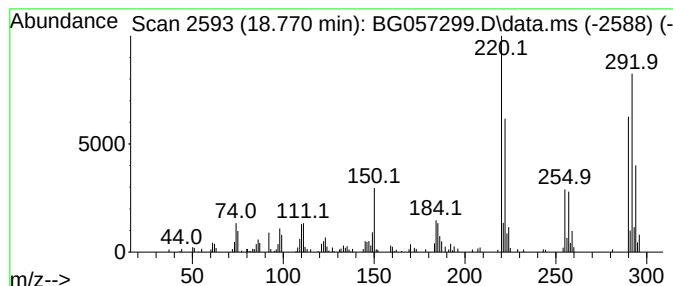
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.770	9.29 ng/ul	157318	Phenanthrene-d10	17.731	
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, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
3	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99
4	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
5	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050423\
Data File : BG057299.D
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Operator : CG/JU
Sample : 02419-36DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

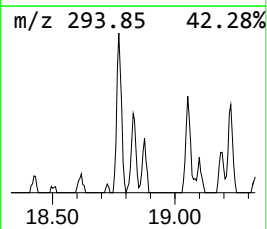
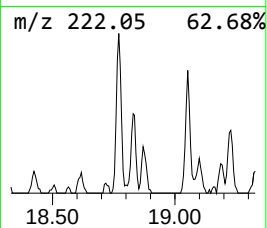
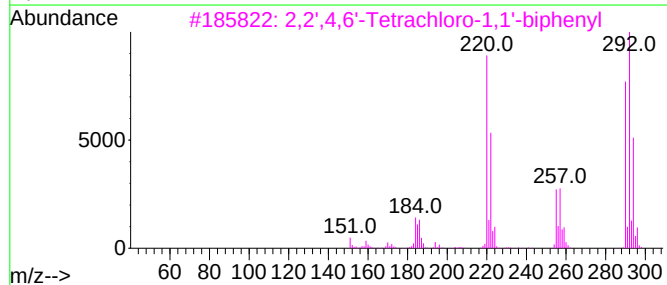
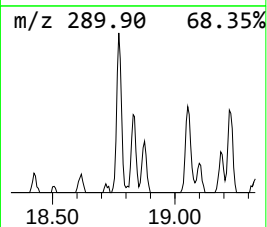
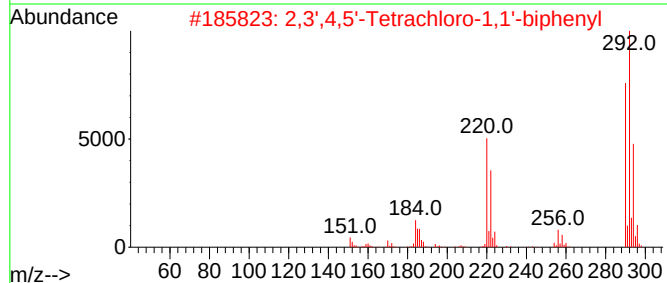
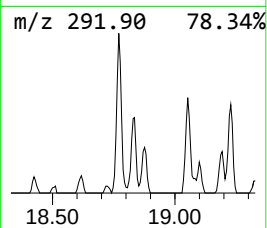
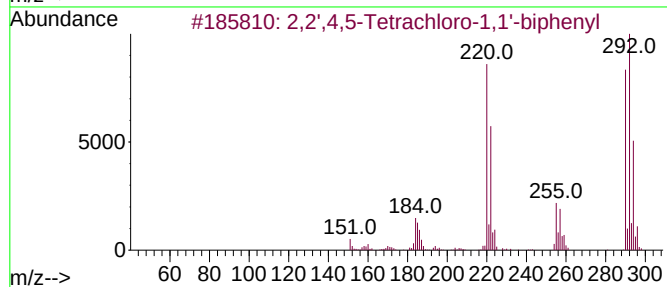
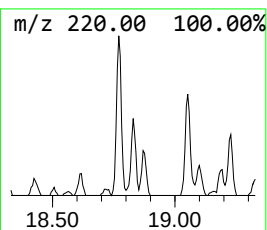
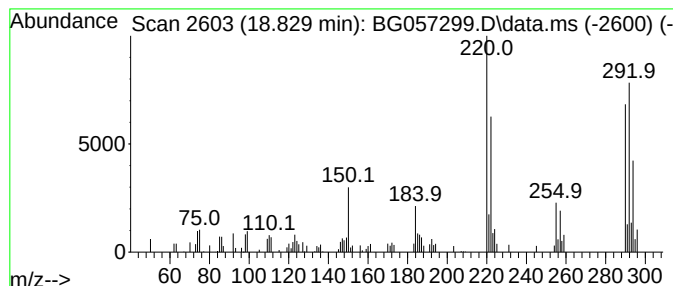
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 2,2',4,5-Tetrachloro-1,1'-b... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.829	4.51 ng/ul	76310	Phenanthrene-d10	17.731	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	98
2	2,3',4,5'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-52-7	97
3	2,2',4,6'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	068194-04-7	97
4	2,3,3',6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074472-33-6	97
5	2,3,4',5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074472-34-7	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050423\
Data File : BG057299.D
Acq On : 4 May 2023 11:55
Operator : CG/JU
Sample : 02419-36DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

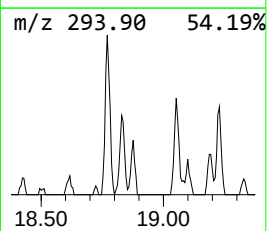
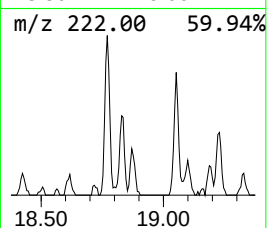
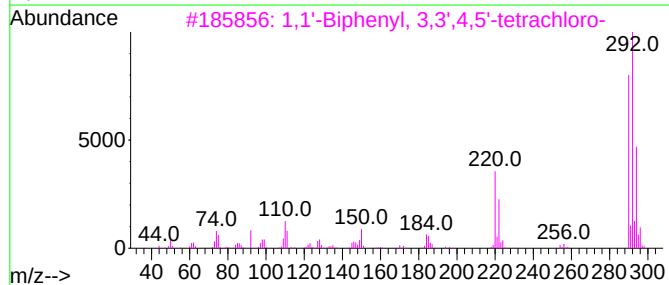
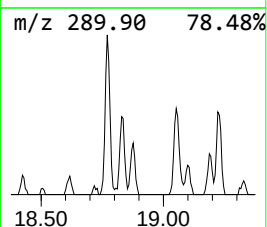
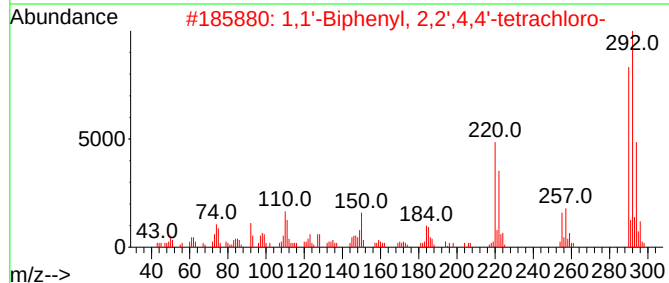
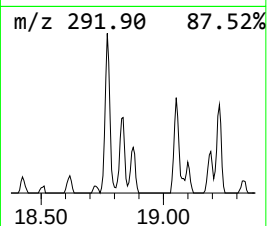
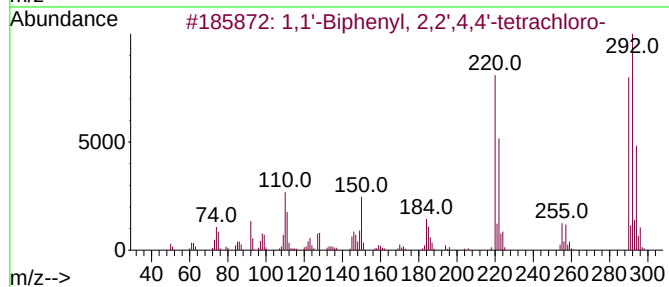
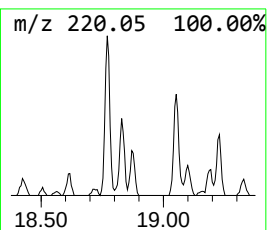
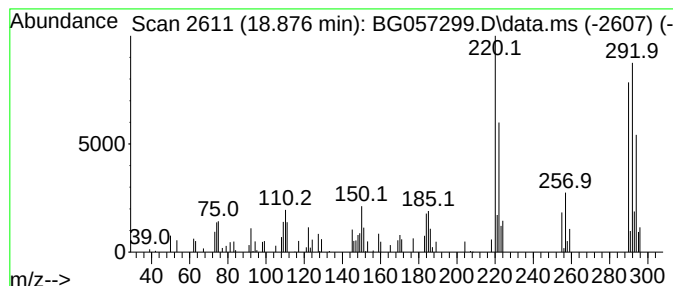
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 1,1'-Biphenyl, 3,3',4,5'-te... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.876	2.68 ng/ul	45381	Phenanthrene-d10	17.731		
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,4'-tetrachloro-	290	C12H6Cl4	002437-79-8	99	
2	1,1'-Biphenyl, 2,2',4,4'-tetrachloro-	290	C12H6Cl4	002437-79-8	99	
3	1,1'-Biphenyl, 3,3',4,5'-tetrachloro-	290	C12H6Cl4	041464-48-6	99	
4	2,3',4',5'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-48-0	99	
5	3,4,4',5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-50-4	99	



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050423\
Data File : BG057299.D
Acq On : 4 May 2023 11:55
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Sample : 02419-36DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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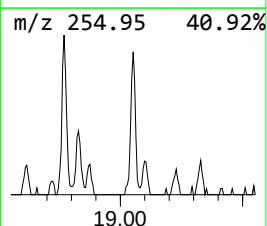
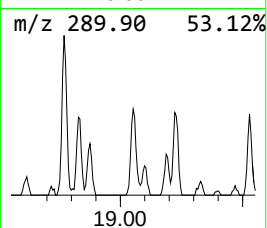
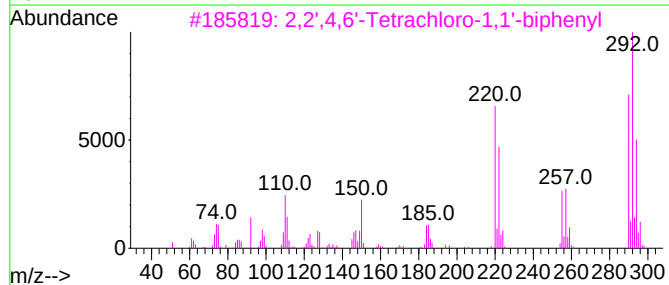
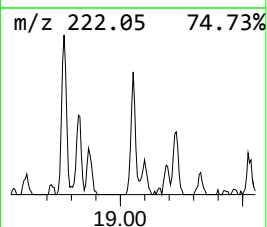
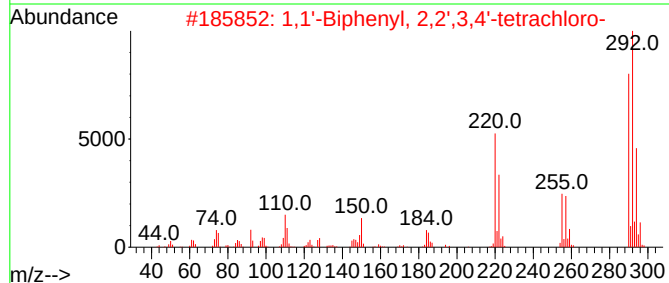
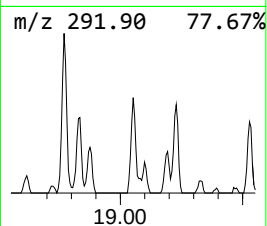
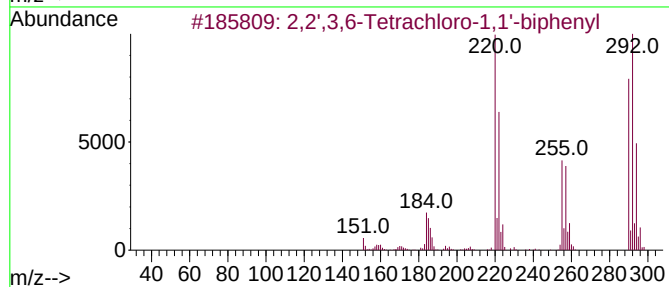
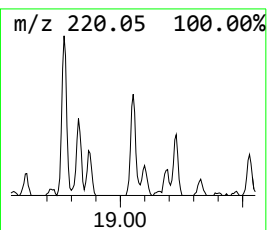
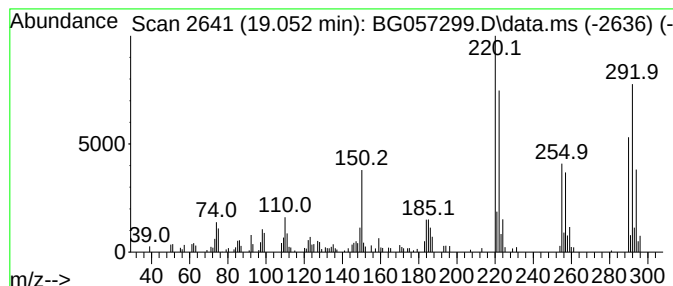
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 2,2',3,6-Tetrachloro-1,1'-b... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.052	6.53 ng/ul	110657	Phenanthrene-d10	17.731

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-45-7	99		
2	1,1'-Biphenyl, 2,2',3,4'-tetrach...	290	C12H6Cl4	036559-22-5	99		
3	2,2',4,6'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	068194-04-7	99		
4	1,1'-Biphenyl, 2,2',3,4-tetrachl...	290	C12H6Cl4	052663-59-9	99		
5	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	98		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050423\
Data File : BG057299.D
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Operator : CG/JU
Sample : 02419-36DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

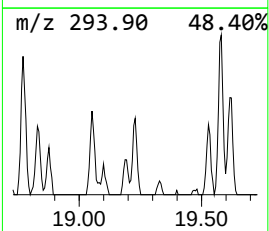
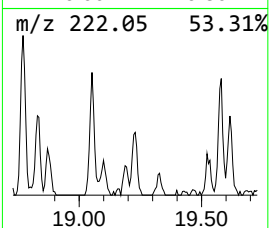
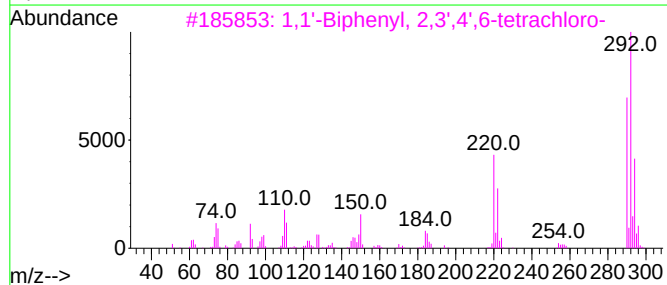
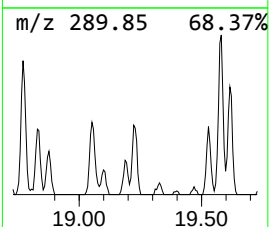
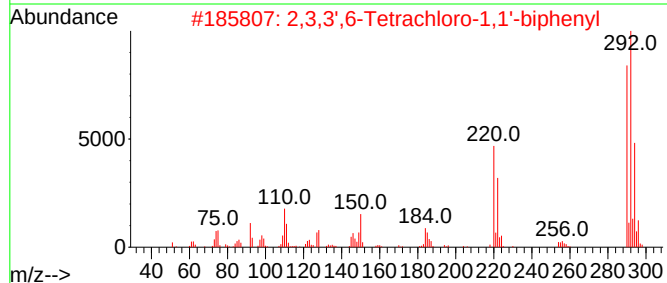
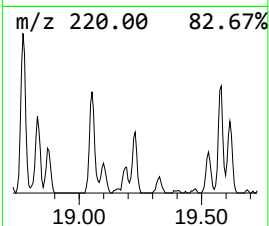
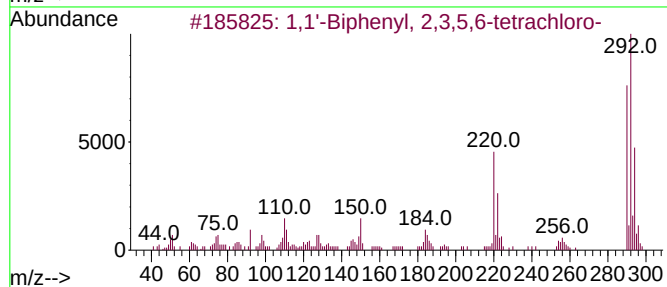
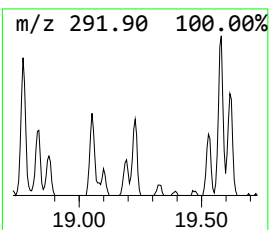
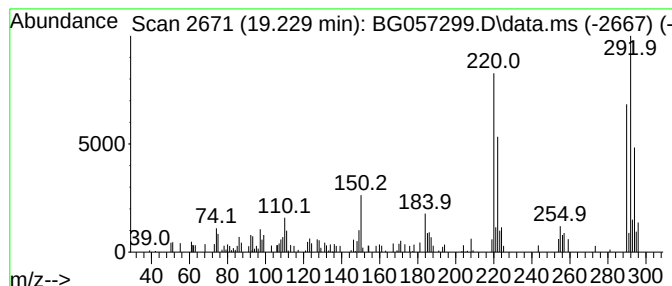
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 1,1'-Biphenyl, 2,3,5,6-tetr... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.229	3.76 ng/ul	63633	Phenanthrene-d10	17.731	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,5,6-tetrachloro-	290	C12H6Cl4	033284-54-7	99
2	2,3,3',6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074472-33-6	99
3	1,1'-Biphenyl, 2,3',4',6-tetrachl...	290	C12H6Cl4	041464-46-4	99
4	1,1'-Biphenyl, 2,3,4',6-tetrachl...	290	C12H6Cl4	052663-58-8	99
5	1,1'-Biphenyl, 2,3,4,4'-tetrachl...	290	C12H6Cl4	033025-41-1	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050423\
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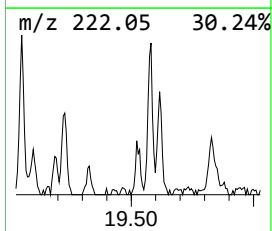
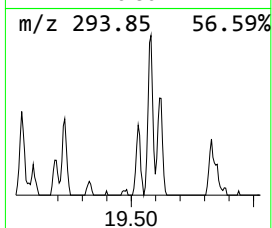
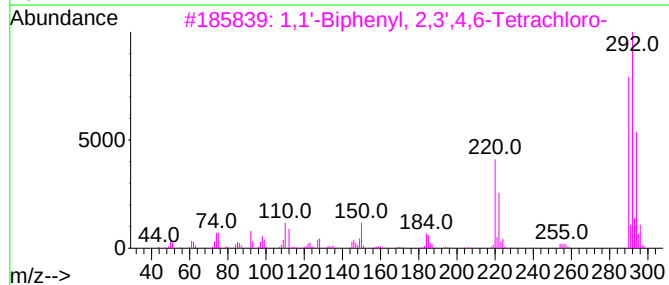
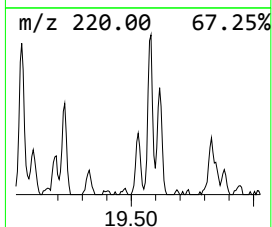
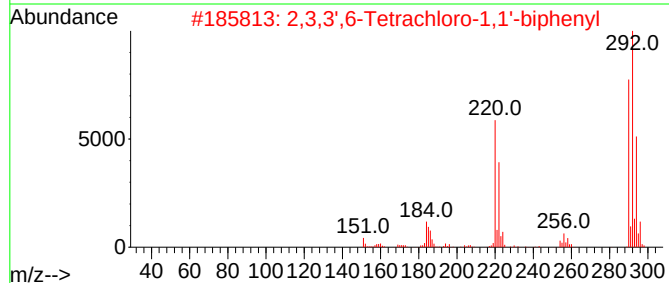
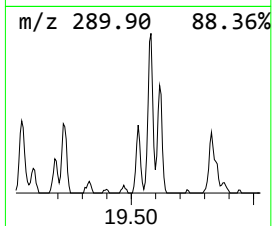
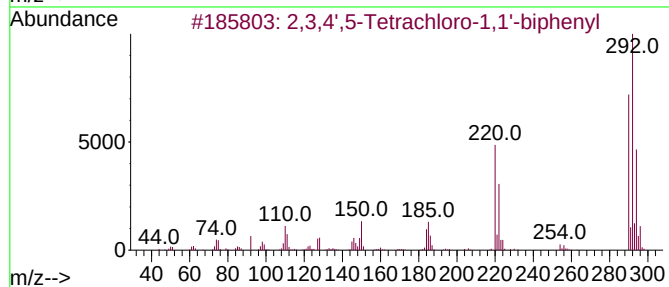
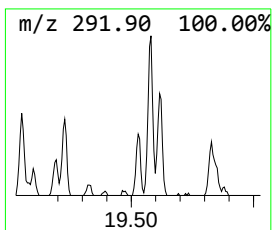
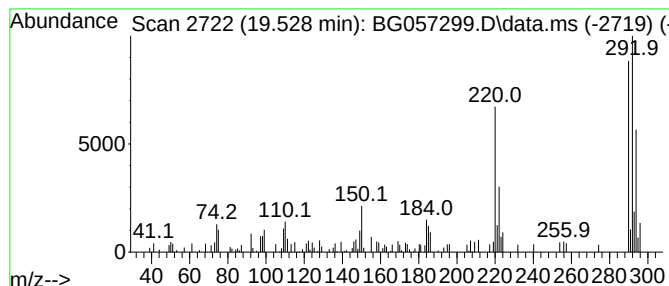
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 2,3,4',5-Tetrachloro-1,1'-b... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.528	2.57 ng/ul	43469	Phenanthrene-d10	17.731	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,4',5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074472-34-7	99
2	2,3,3',6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074472-33-6	99
3	1,1'-Biphenyl, 2,3',4,6-Tetrachl...	290	C12H6Cl4	060233-24-1	98
4	2,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-53-8	98
5	2,3',5',6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074338-23-1	98



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050423\
Data File : BG057299.D
Acq On : 4 May 2023 11:55
Operator : CG/JU
Sample : 02419-36DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

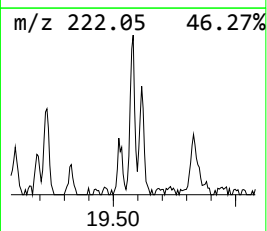
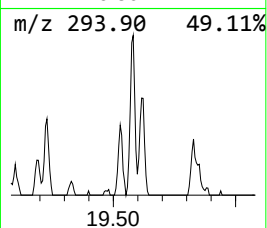
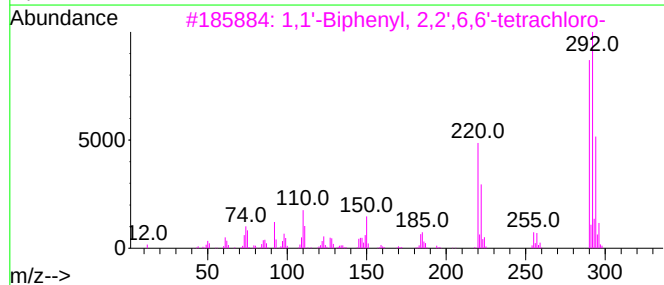
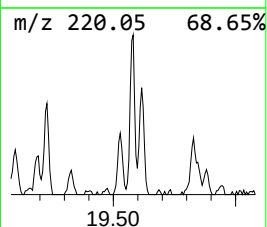
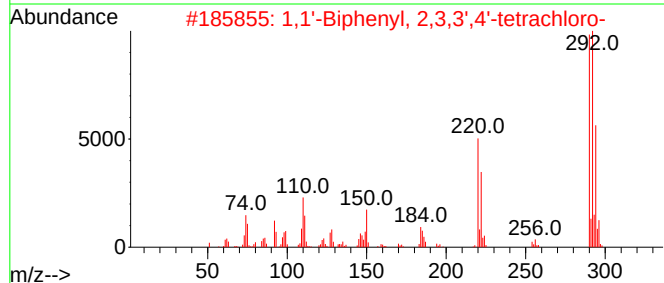
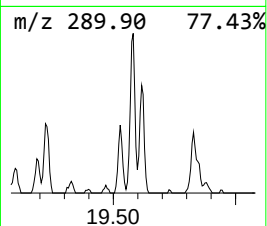
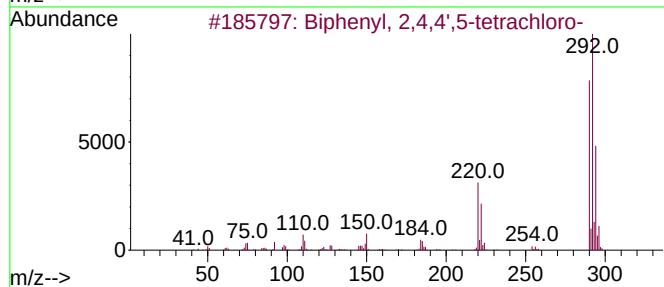
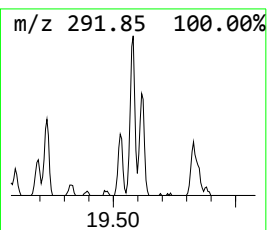
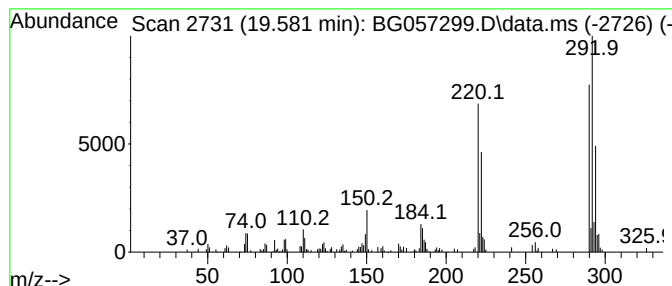
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Biphenyl, 2,4,4',5-tetrachl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.581	8.11 ng/ul	137328	Phenanthrene-d10	17.731	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	99
2	1,1'-Biphenyl, 2,3,3',4'-tetrach...	290	C12H6Cl4	041464-43-1	99
3	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99
4	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99
5	3,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-49-1	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050423\
Data File : BG057299.D
Acq On : 4 May 2023 11:55
Operator : CG/JU
Sample : 02419-36DL 5X
Misc :
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Instrument :
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ClientSampleId :
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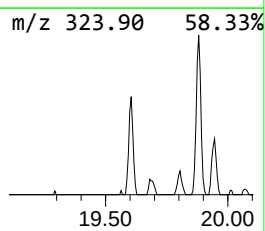
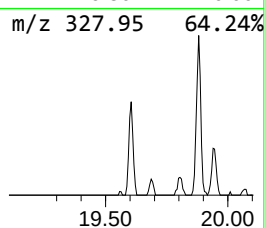
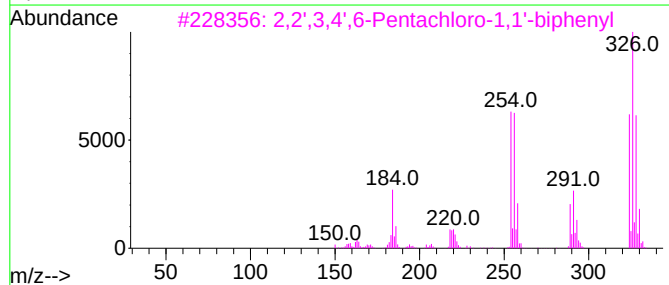
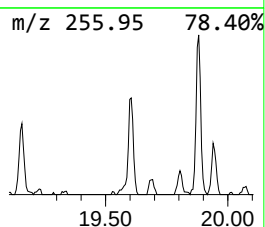
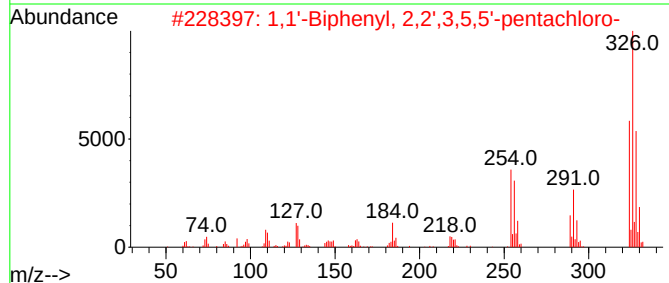
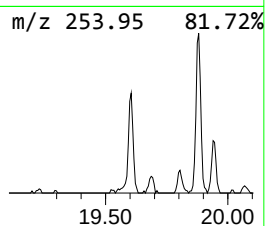
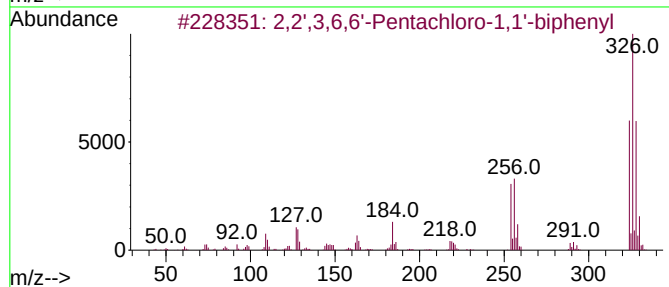
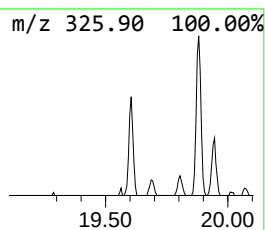
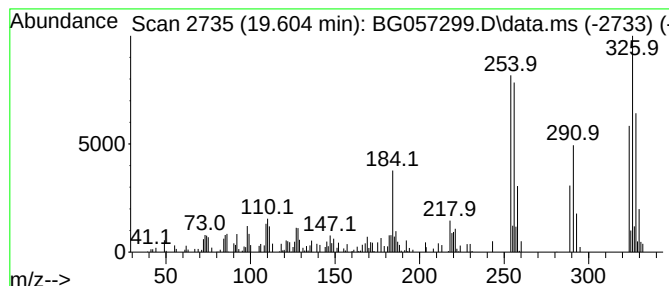
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 2,2',3,6,6'-Pentachloro-1,1... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.605	9.80 ng/ul	165896	Phenanthrene-d10	17.731

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',3,5,5'-penta...	324	C12H5Cl5	052663-61-3	98		
3	2,2',3,4',6-Pentachloro-1,1'-bip...	324	C12H5Cl5	068194-05-8	97		
4	1,1'-Biphenyl, 2,2',4,5,5'-penta...	324	C12H5Cl5	037680-73-2	97		
5	2,2',3,4,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-57-2	97		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050423\
Data File : BG057299.D
Acq On : 4 May 2023 11:55
Operator : CG/JU
Sample : 02419-36DL 5X
Misc :
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Instrument :
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ClientSampleId :
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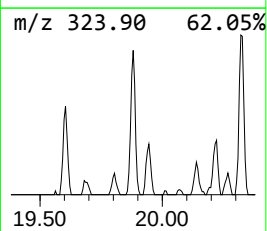
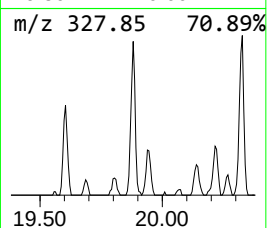
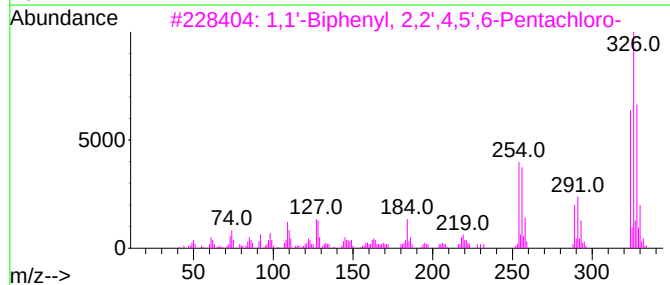
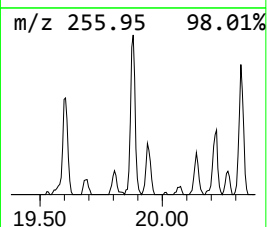
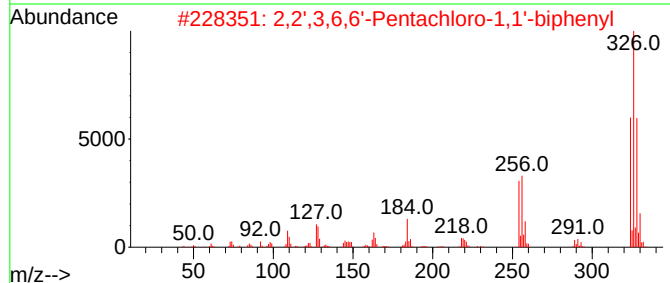
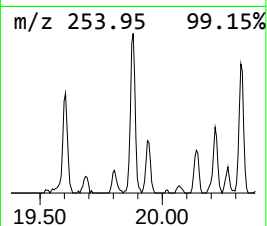
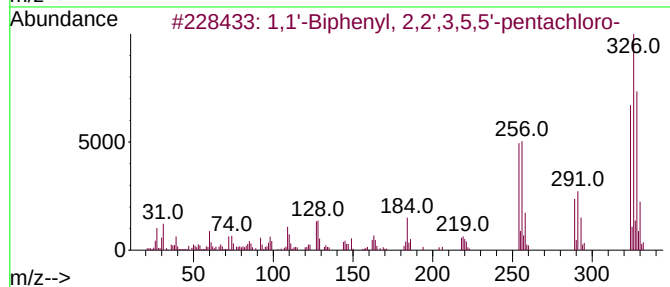
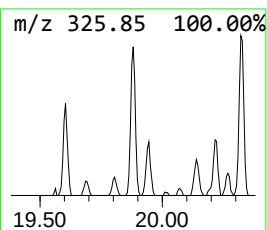
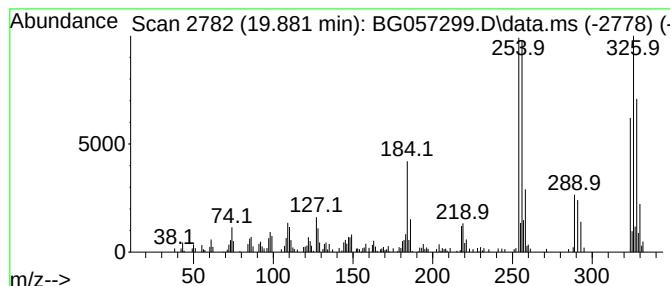
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 1,1'-Biphenyl, 2,2',3,5,5'-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.881	10.69 ng/ul	180991	Phenanthrene-d10	17.731

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,5,5'-penta...	324	C12H5Cl5	052663-61-3	99		
2	2,2',3,6,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-54-9	97		
3	1,1'-Biphenyl, 2,2',4,5',6-Penta...	324	C12H5Cl5	060145-21-3	97		
4	1,1'-Biphenyl, 2,2',3',4,5-Penta...	324	C12H5Cl5	041464-51-1	96		
5	1,1'-Biphenyl, 2,2',3,4,6-Pentac...	324	C12H5Cl5	055215-17-3	96		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050423\
Data File : BG057299.D
Acq On : 4 May 2023 11:55
Operator : CG/JU
Sample : 02419-36DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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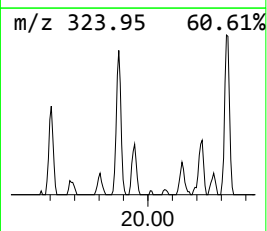
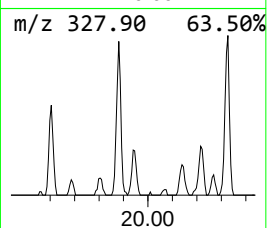
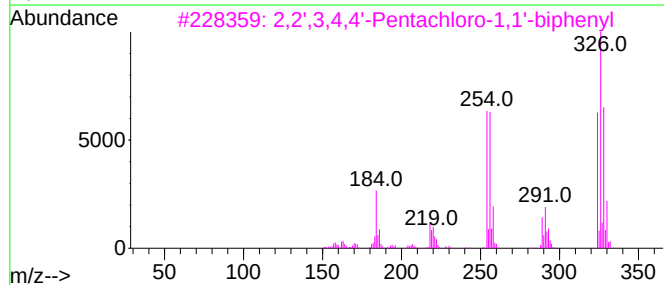
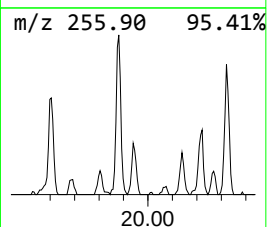
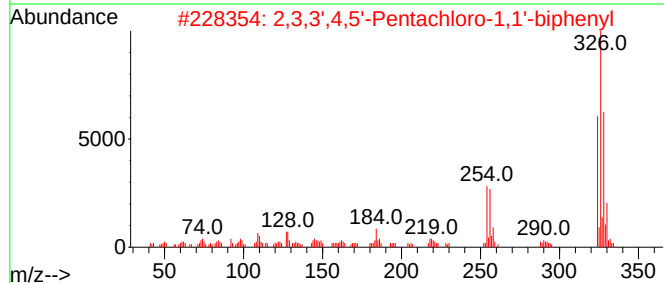
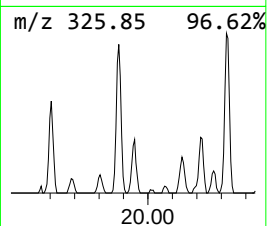
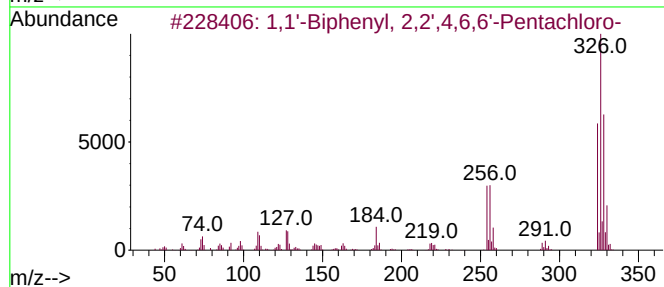
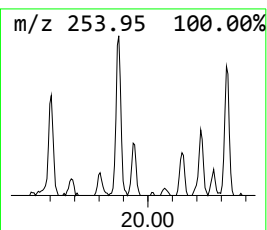
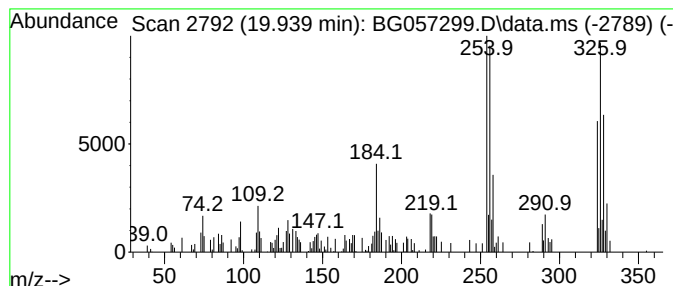
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 1,1'-Biphenyl, 2,2',4,6,6'-... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.939	4.24 ng/ul	62917	Chrysene-d12	22.048

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,6,6'-Penta...	324	C12H5Cl5	056558-16-8	99		
2	2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	070362-41-3	98		
3	2,2',3,4,4'-Pentachloro-1,1'-bip...	324	C12H5Cl5	065510-45-4	98		
4	2,2',3,6,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-54-9	97		
5	2,3,3',5,6-Pentachloro-1,1'-biph...	324	C12H5Cl5	074472-36-9	97		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050423\
Data File : BG057299.D
Acq On : 4 May 2023 11:55
Operator : CG/JU
Sample : 02419-36DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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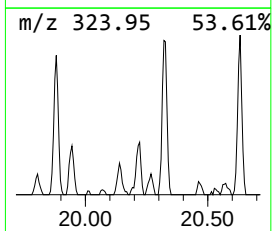
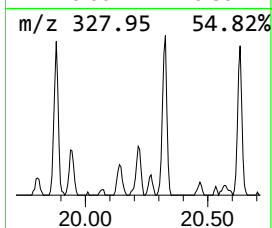
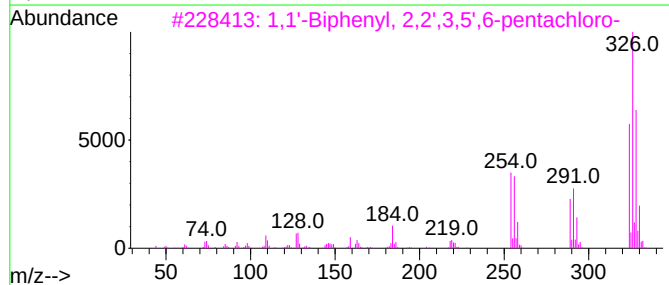
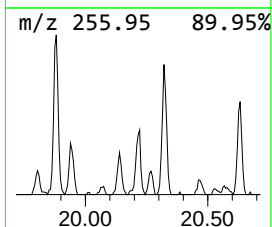
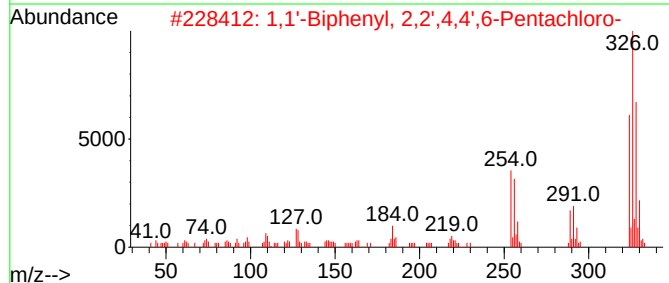
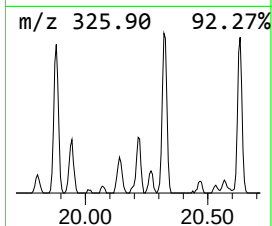
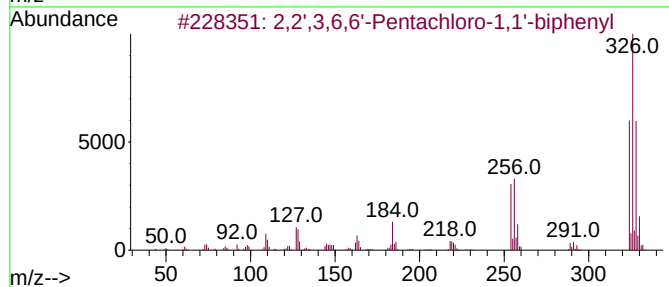
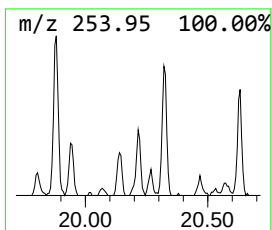
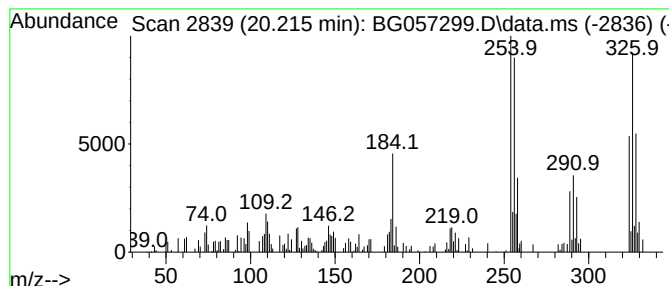
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 1,1'-Biphenyl, 2,2',4,4',6-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.215	4.74 ng/ul	70315	Chrysene-d12	22.048

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,6,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-54-9	99		
2	1,1'-Biphenyl, 2,2',4,4',6-Penta...	324	C12H5Cl5	039485-83-1	99		
3	1,1'-Biphenyl, 2,2',3,5',6-penta...	324	C12H5Cl5	038379-99-6	99		
4	2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	070362-41-3	97		
5	1,1'-Biphenyl, 2,2',4,5',6-Penta...	324	C12H5Cl5	060145-21-3	96		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050423\
Data File : BG057299.D
Acq On : 4 May 2023 11:55
Operator : CG/JU
Sample : 02419-36DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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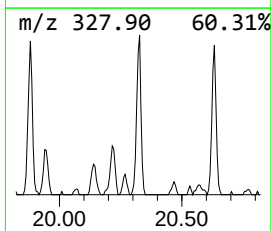
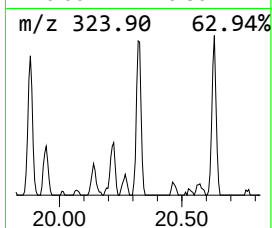
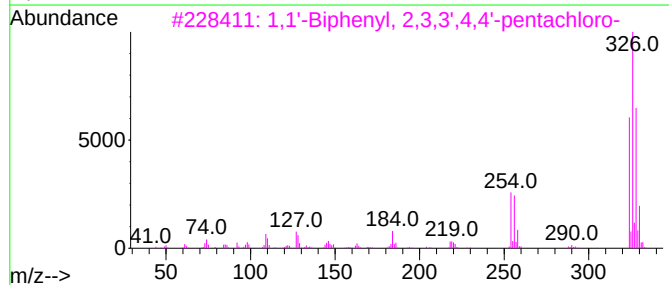
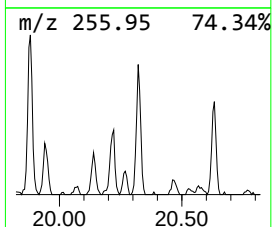
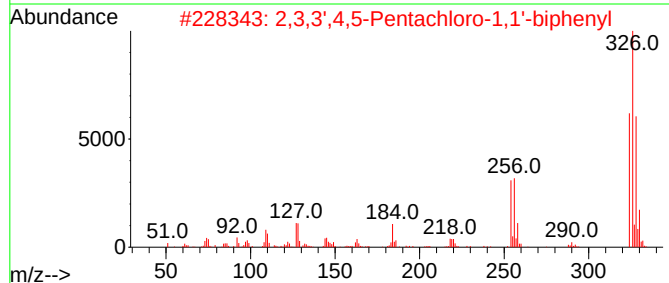
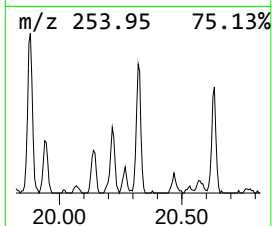
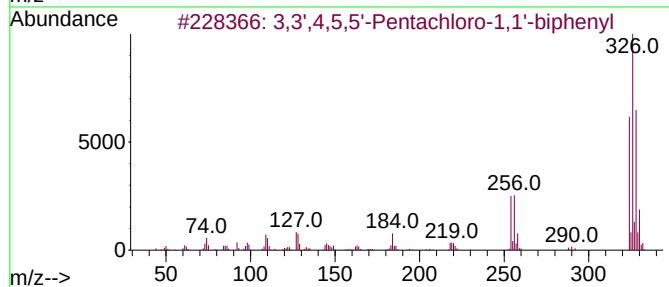
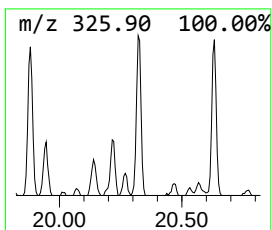
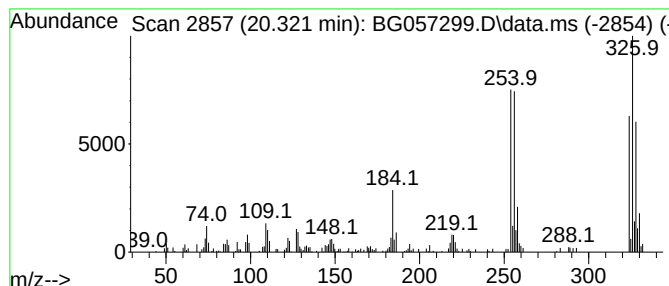
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 3,3',4,5,5'-Pentachloro-1,1... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.321	9.57 ng/ul	142077	Chrysene-d12	22.048

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,3',4,5,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	039635-33-1	99		
2	2,3,3',4,5-Pentachloro-1,1'-biph...	324	C12H5Cl5	070424-69-0	99		
3	1,1'-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	98		
4	2,3,3',4,5-Pentachloro-1,1'-biph...	324	C12H5Cl5	070424-69-0	98		
5	1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	97		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050423\
Data File : BG057299.D
Acq On : 4 May 2023 11:55
Operator : CG/JU
Sample : 02419-36DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW931DL

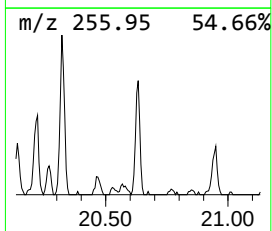
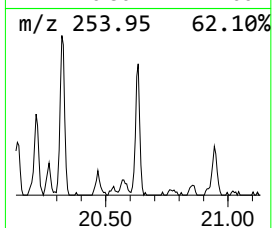
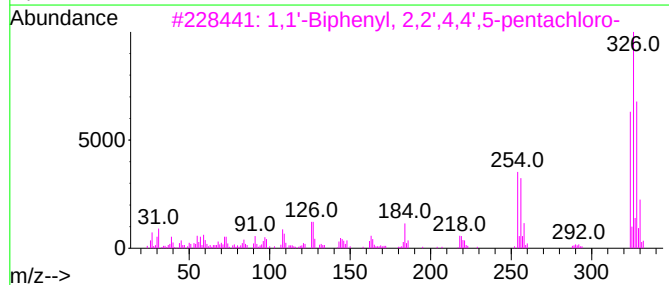
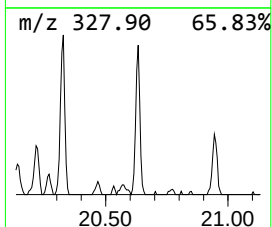
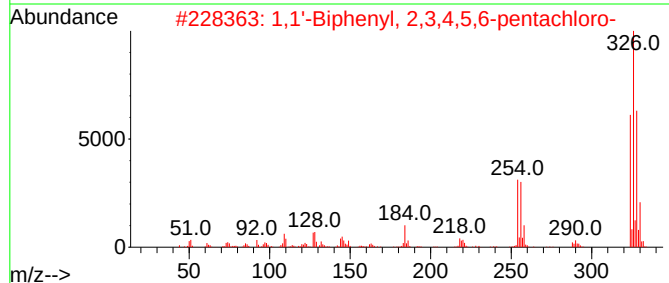
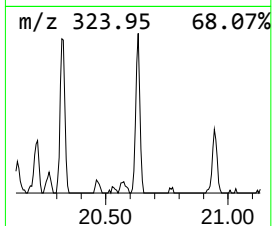
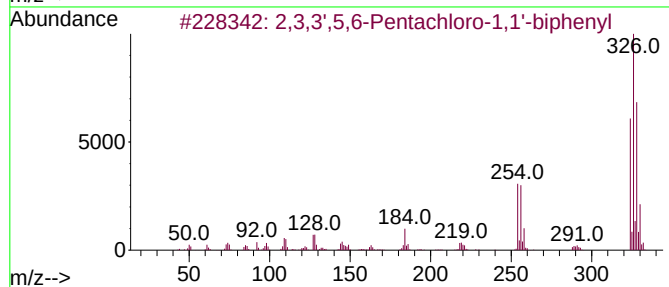
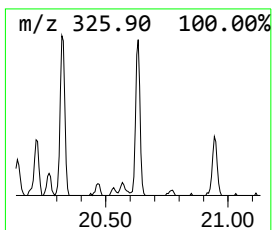
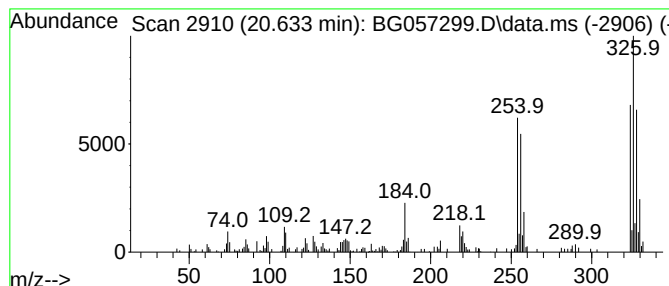
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 2,3,3',5,6-Pentachloro-1,1'... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.633	8.80 ng/ul	130717	Chrysene-d12	22.048

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,3',5,6-Pentachloro-1,1'-biph...	324	C12H5Cl5	074472-36-9	99		
2	1,1'-Biphenyl, 2,3,4,5,6-pentach...	324	C12H5Cl5	018259-05-7	99		
3	1,1'-Biphenyl, 2,2',4,4',5-penta...	324	C12H5Cl5	038380-01-7	99		
4	2,3,3',4',5'- Pentachloro-1,1'-b...	324	C12H5Cl5	076842-07-4	98		
5	1,1'-Biphenyl, 2,2',4,6,6'-Penta...	324	C12H5Cl5	056558-16-8	98		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050423\
Data File : BG057299.D
Acq On : 4 May 2023 11:55
Operator : CG/JU
Sample : 02419-36DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW931DL

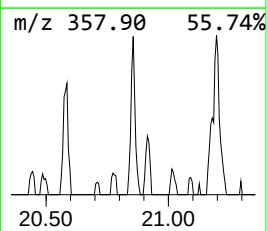
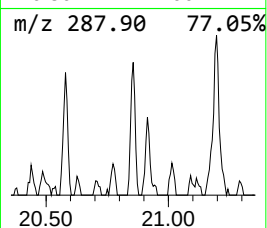
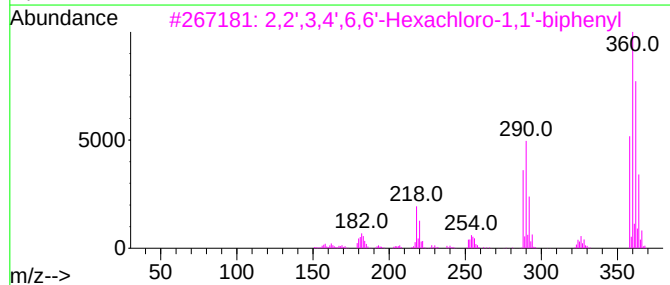
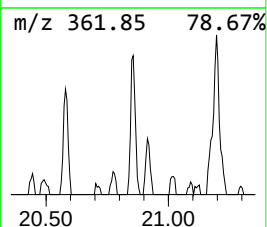
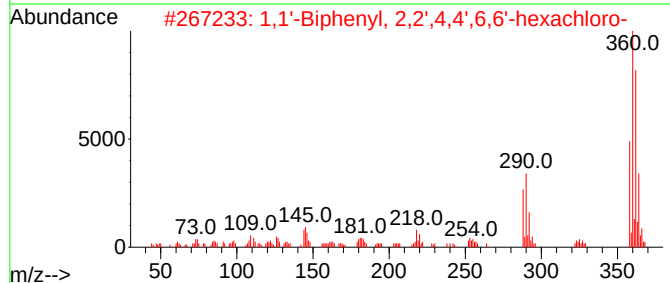
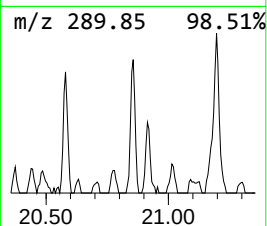
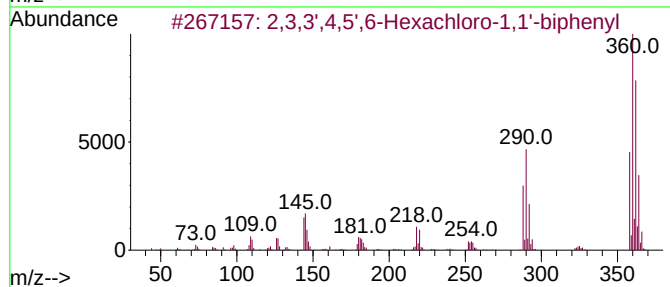
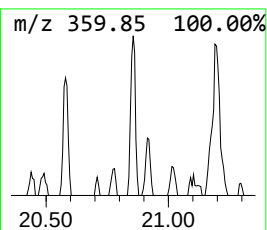
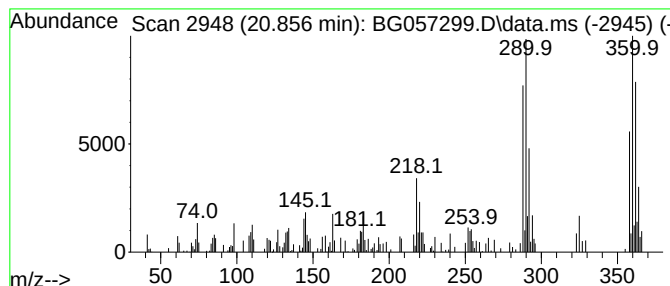
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 2,3,3',4,5',6-Hexachloro-1,... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.856	4.67 ng/ul	69324	Chrysene-d12	22.048

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,3',4,5',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-43-8	99		
2	1,1'-Biphenyl, 2,2',4,4',6,6'-he...	358	C12H4Cl6	033979-03-2	99		
3	2,2',3,4',6,6'-Hexachloro-1,1'-b...	358	C12H4Cl6	068194-08-1	99		
4	1,1'-Biphenyl, 2,2',3,3',6,6'-he...	358	C12H4Cl6	038411-22-2	99		
5	2,3,4,4',5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-63-6	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050423\
Data File : BG057299.D
Acq On : 4 May 2023 11:55
Operator : CG/JU
Sample : 02419-36DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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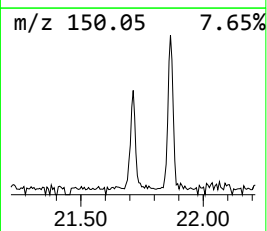
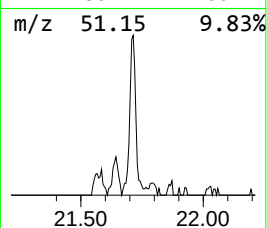
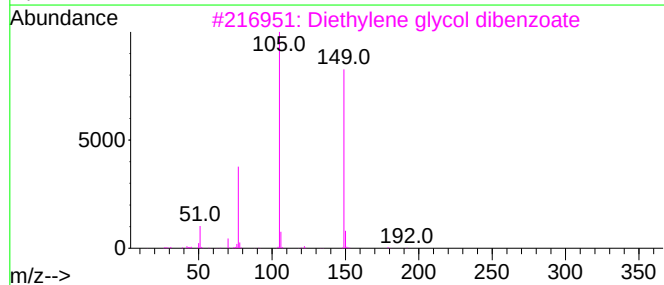
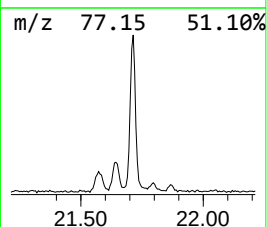
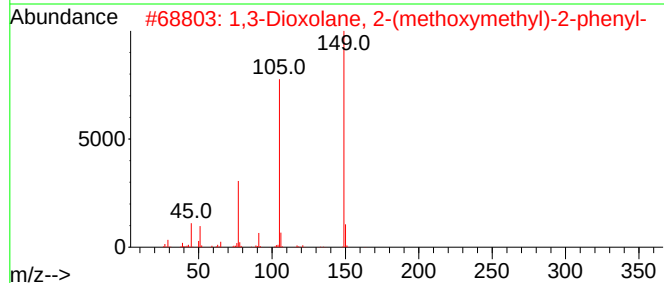
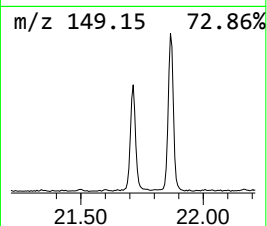
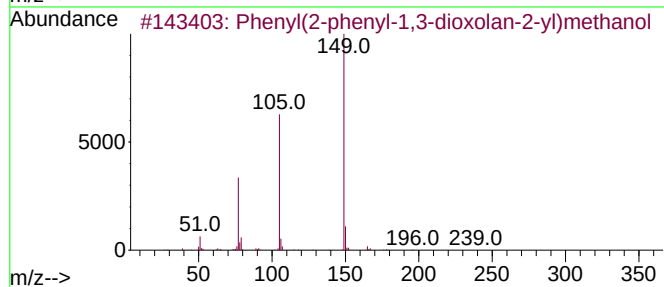
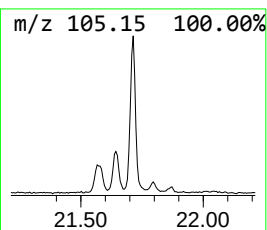
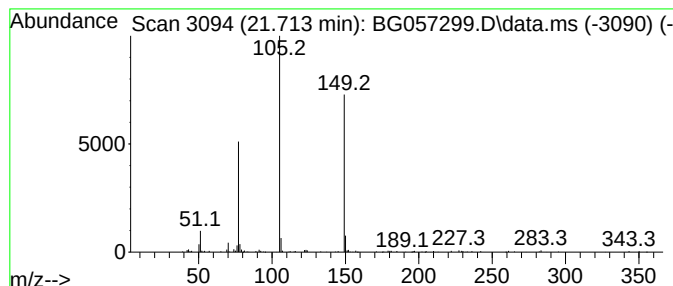
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 Phenyl(2-phenyl-1,3-dioxola... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.713	9.77 ng/ul	145085	Chrysene-d12	22.048

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenyl(2-phenyl-1,3-dioxolan-2-y...	256	C16H16O3	005694-69-9	72
2			1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	72
3			Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	59
4			Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	53
5			Benzoic acid, 3,5-difluophenyl e...	234	C13H8F2O2	1000357-79-5	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050423\
Data File : BG057299.D
Acq On : 4 May 2023 11:55
Operator : CG/JU
Sample : 02419-36DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW931DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1,1'-Biphenyl, ...	17.889	3.1	ng/ul	52392	4	17.731	338666	20.0
1,1'-Biphenyl, ...	18.312	15.6	ng/ul	264423	4	17.731	338666	20.0
1,1'-Biphenyl, ...	18.436	2.5	ng/ul	43247	4	17.731	338666	20.0
1,1'-Biphenyl, ...	18.565	3.4	ng/ul	57392	4	17.731	338666	20.0
1,1'-Biphenyl, ...	18.770	9.3	ng/ul	157318	4	17.731	338666	20.0
2,2',4,5-Tetrac...	18.829	4.5	ng/ul	76310	4	17.731	338666	20.0
1,1'-Biphenyl, ...	18.876	2.7	ng/ul	45381	4	17.731	338666	20.0
2,2',3,6-Tetrac...	19.052	6.5	ng/ul	110657	4	17.731	338666	20.0
1,1'-Biphenyl, ...	19.229	3.8	ng/ul	63633	4	17.731	338666	20.0
2,3,4',5-Tetrac...	19.528	2.6	ng/ul	43469	4	17.731	338666	20.0
Biphenyl, 2,4,4...	19.581	8.1	ng/ul	137328	4	17.731	338666	20.0
2,2',3,6,6'-Pen...	19.605	9.8	ng/ul	165896	4	17.731	338666	20.0
1,1'-Biphenyl, ...	19.881	10.7	ng/ul	180991	4	17.731	338666	20.0
1,1'-Biphenyl, ...	19.939	4.2	ng/ul	62917	5	22.048	296924	20.0
1,1'-Biphenyl, ...	20.215	4.7	ng/ul	70315	5	22.048	296924	20.0
3,3',4,5,5'-Pen...	20.321	9.6	ng/ul	142077	5	22.048	296924	20.0
2,3,3',5,6-Pent...	20.633	8.8	ng/ul	130717	5	22.048	296924	20.0
2,3,3',4,5',6-H...	20.856	4.7	ng/ul	69324	5	22.048	296924	20.0
Phenyl(2-phenyl...	21.713	9.8	ng/ul	145085	5	22.048	296924	20.0