

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042523\
 Data File : BM039663.D
 Acq On : 26 Apr 2023 05:18
 Operator : CG/JU
 Sample : 02419-27
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW924

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_M\Methods\SFAM-EPA-BM041823.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM039663.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.084	6	8	11	rVB	75981	69231	3.14%	0.327%
2	3.125	11	15	26	rVB	1432310	1462224	66.35%	6.914%
3	3.437	64	68	70	rBV	12909	17503	0.79%	0.083%
4	3.466	70	73	78	rVB	56367	64026	2.91%	0.303%
5	3.796	127	129	133	rVB4	4484	5607	0.25%	0.027%
6	3.848	134	138	146	rVV2	71325	105574	4.79%	0.499%
7	3.954	151	156	161	rBV2	69416	96116	4.36%	0.454%
8	4.396	227	231	233	rVB2	5321	8421	0.38%	0.040%
9	4.495	243	248	252	rBV8	5256	9593	0.44%	0.045%
10	4.601	264	266	272	rVB6	4273	6299	0.29%	0.030%
11	4.707	279	284	288	rVB7	5468	8368	0.38%	0.040%
12	4.801	295	300	306	rBV3	9459	15516	0.70%	0.073%
13	4.966	326	328	333	rVB6	3669	5474	0.25%	0.026%
14	5.295	380	384	390	rVV	16247	26759	1.21%	0.127%
15	5.566	429	430	437	rVB5	3729	6067	0.28%	0.029%
16	6.042	506	511	514	rBV6	3571	6003	0.27%	0.028%
17	6.295	547	554	557	rBV6	4022	8358	0.38%	0.040%
18	6.331	557	560	564	rBV5	3138	4808	0.22%	0.023%
19	6.531	590	594	601	rVB7	4434	8100	0.37%	0.038%
20	6.742	626	630	635	rBV8	3302	4728	0.21%	0.022%
21	6.919	656	660	664	rBV6	5634	9584	0.43%	0.045%
22	7.284	717	722	725	rVV7	3528	5376	0.24%	0.025%
23	7.413	740	744	748	rBV6	4127	7245	0.33%	0.034%
24	7.807	799	811	828	rBV	836223	1618153	73.43%	7.652%
25	8.342	900	902	909	rVB8	4702	7030	0.32%	0.033%
26	9.054	1017	1023	1024	rBV4	4257	7266	0.33%	0.034%
27	9.978	1178	1180	1185	rBV4	2758	4889	0.22%	0.023%
28	10.283	1226	1232	1237	rVB7	3393	6482	0.29%	0.031%
29	10.383	1243	1249	1253	rVB7	3794	8098	0.37%	0.038%
30	10.554	1271	1278	1280	rBV8	3625	7777	0.35%	0.037%
31	10.613	1280	1288	1310	rBV	1106075	2162051	98.11%	10.223%
32	11.083	1361	1368	1372	rBV9	5123	12389	0.56%	0.059%
33	11.272	1394	1400	1401	rBV5	2877	4042	0.18%	0.019%
34	11.430	1424	1427	1431	rBV6	2298	4493	0.20%	0.021%
35	11.624	1456	1460	1471	rBV6	7791	15093	0.68%	0.071%
36	12.036	1527	1530	1531	rBV3	4297	5114	0.23%	0.024%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
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37	12.330	1572	1580	1581	rBV7	3729	7314	0.33%	0.035%
38	12.519	1606	1612	1618	rBV9	5115	9516	0.43%	0.045%
39	12.807	1656	1661	1664	rVB7	2359	4480	0.20%	0.021%
40	12.848	1664	1668	1670	rBV4	3514	4332	0.20%	0.020%

41	12.983	1687	1691	1696	rVV3	10108	14914	0.68%	0.071%
42	13.283	1734	1742	1747	rVV5	13015	25130	1.14%	0.119%
43	13.377	1755	1758	1764	rBV8	3861	6100	0.28%	0.029%
44	13.518	1778	1782	1783	rVV3	3013	4232	0.19%	0.020%
45	13.554	1783	1788	1796	rVB2	46542	73100	3.32%	0.346%

46	13.789	1824	1828	1834	rVB7	6703	10856	0.49%	0.051%
47	13.860	1834	1840	1847	rBV7	5394	14670	0.67%	0.069%
48	13.936	1847	1853	1857	rVV5	13034	20725	0.94%	0.098%
49	14.065	1870	1875	1879	rVV6	5882	8857	0.40%	0.042%
50	14.195	1893	1897	1898	rBV4	5070	6163	0.28%	0.029%

51	14.283	1906	1912	1919	rVB6	7359	13787	0.63%	0.065%
52	14.360	1919	1925	1931	rBV2	15043	28516	1.29%	0.135%
53	14.465	1936	1943	1959	rBV	1358228	2203702	100.00%	10.420%
54	14.871	2008	2012	2014	rBV5	3438	6012	0.27%	0.028%
55	14.971	2027	2029	2037	rVB8	5348	10318	0.47%	0.049%

56	15.095	2045	2050	2055	rBV8	5196	10079	0.46%	0.048%
57	15.165	2057	2062	2068	rBV9	6192	12186	0.55%	0.058%
58	15.242	2072	2075	2078	rBV4	7559	10123	0.46%	0.048%
59	15.307	2082	2086	2091	rBV3	14700	21102	0.96%	0.100%
60	15.507	2118	2120	2121	rBV2	4795	4017	0.18%	0.019%

61	15.530	2121	2124	2127	rVV5	4180	7187	0.33%	0.034%
62	15.718	2150	2156	2162	rVV4	33893	62757	2.85%	0.297%
63	15.942	2189	2194	2198	rBV2	41094	64566	2.93%	0.305%
64	16.189	2229	2236	2243	rBV2	47643	110689	5.02%	0.523%
65	16.448	2276	2280	2287	rBV	34471	60606	2.75%	0.287%

66	16.742	2326	2330	2336	rBV4	29975	49801	2.26%	0.235%
67	16.965	2365	2368	2372	rBV3	21647	25514	1.16%	0.121%
68	17.018	2372	2377	2385	rBV3	32141	61354	2.78%	0.290%
69	17.095	2385	2390	2393	rBV	93272	130759	5.93%	0.618%
70	17.130	2393	2396	2402	rVB2	49495	72767	3.30%	0.344%

71	17.218	2405	2411	2416	rBV	1377188	2182484	99.04%	10.320%
72	17.395	2436	2441	2451	rVB2	140955	217437	9.87%	1.028%
73	17.671	2484	2488	2491	rBV	42343	59071	2.68%	0.279%
74	17.706	2491	2494	2498	rVB3	41948	51580	2.34%	0.244%
75	17.818	2506	2513	2528	rBV	496879	1028347	46.66%	4.863%

76	17.948	2528	2535	2542	rBV3	90210	194469	8.82%	0.920%
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 Title : SVOA CALIBRATION

77	18.071	2551	2556	2560	rBV	128286	201097	9.13%	0.951%
78	18.118	2562	2564	2568	rVB2	46931	58133	2.64%	0.275%
79	18.289	2588	2593	2599	rBV2	341577	476873	21.64%	2.255%
80	18.348	2599	2603	2607	rVV2	198687	266750	12.10%	1.261%
81	18.395	2607	2611	2618	rVB2	109382	161825	7.34%	0.765%
82	18.571	2637	2641	2647	rBV	236636	380667	17.27%	1.800%
83	18.618	2647	2649	2654	rVV	79865	112650	5.11%	0.533%
84	18.683	2657	2660	2663	rVV	87651	131699	5.98%	0.623%
85	18.712	2663	2665	2668	rVV	82817	104632	4.75%	0.495%
86	18.753	2668	2672	2680	rVB	156743	219531	9.96%	1.038%
87	18.853	2686	2689	2695	rVB3	48053	65278	2.96%	0.309%
88	19.065	2721	2725	2729	rVB2	89748	114724	5.21%	0.542%
89	19.112	2729	2733	2735	rBV	235296	296503	13.45%	1.402%
90	19.283	2758	2762	2768	rBV	371167	518331	23.52%	2.451%
91	19.365	2773	2776	2782	rVV	92701	190250	8.63%	0.900%
92	19.418	2782	2785	2793	rVV	231885	326865	14.83%	1.546%
93	19.483	2793	2796	2802	rVB3	78181	100339	4.55%	0.474%
94	19.618	2816	2819	2821	rBV2	65436	83010	3.77%	0.393%
95	19.647	2821	2824	2828	rVV	134000	178858	8.12%	0.846%
96	19.759	2841	2843	2848	rVB	74864	82051	3.72%	0.388%
97	19.871	2858	2862	2867	rVB	197448	244710	11.10%	1.157%
98	20.183	2912	2915	2923	rVB	138110	168460	7.64%	0.797%
99	21.412	3119	3124	3129	rBV	1287319	1952576	88.60%	9.233%
100	23.771	3520	3525	3539	rVB2	941007	1978614	89.79%	9.356%

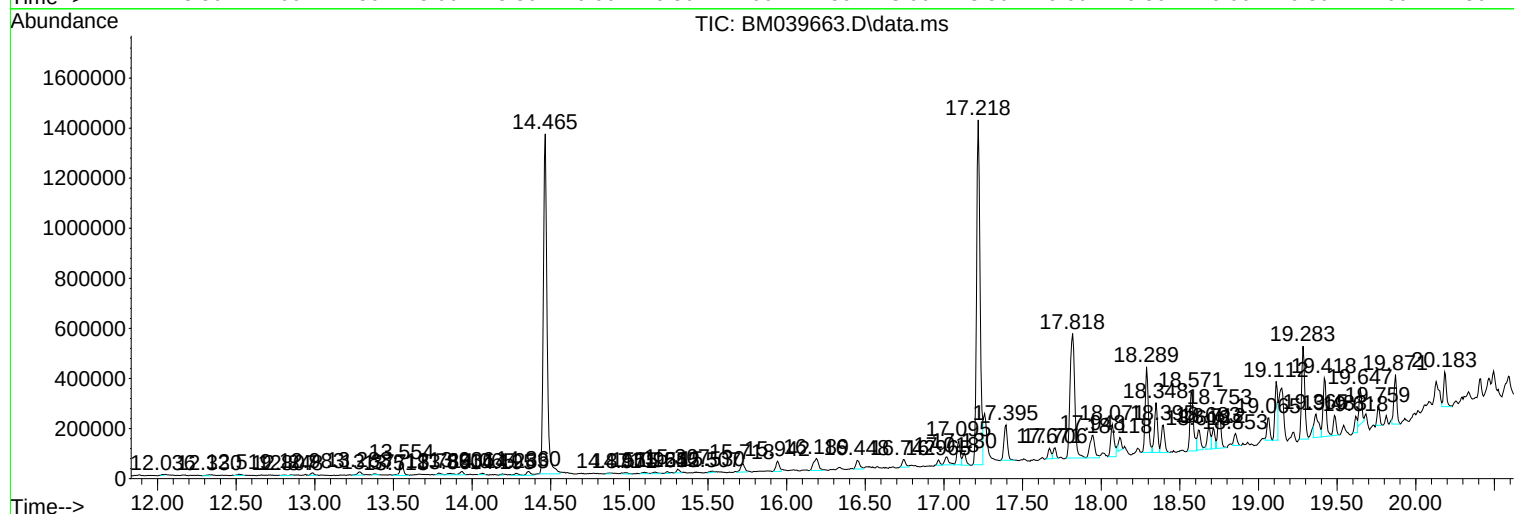
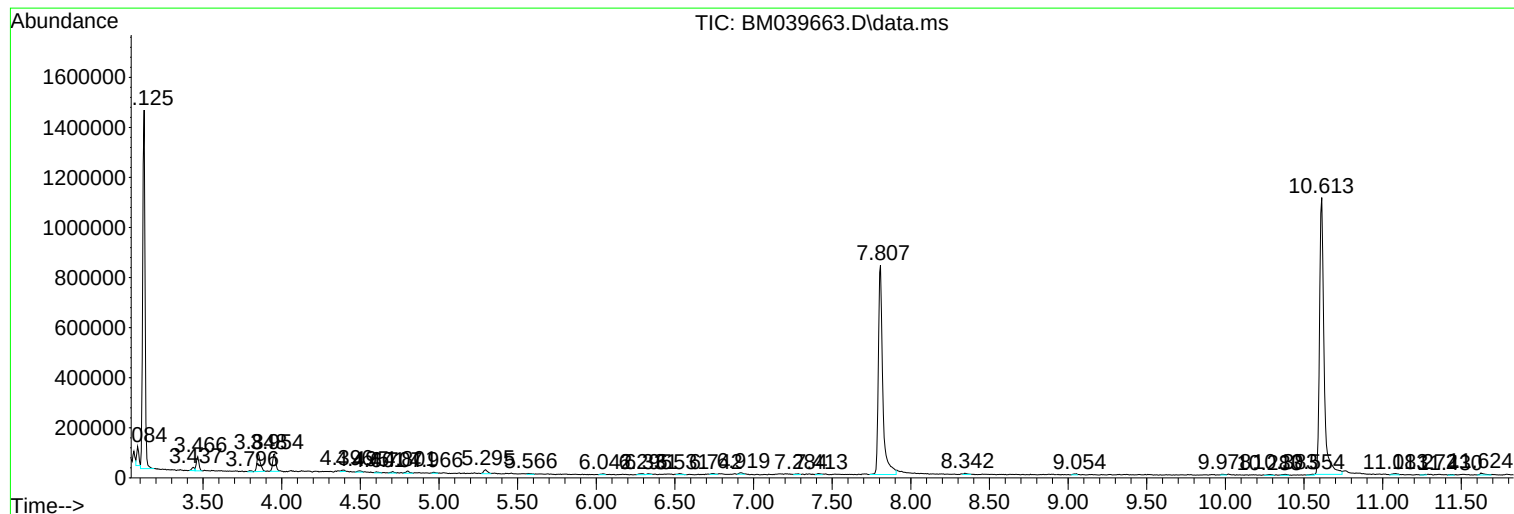
Sum of corrected areas: 21147902

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.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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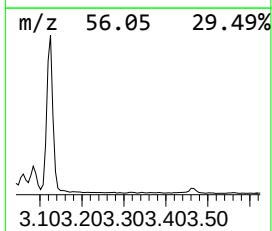
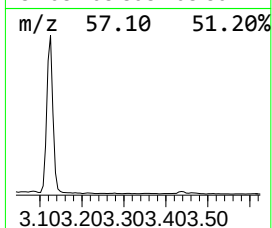
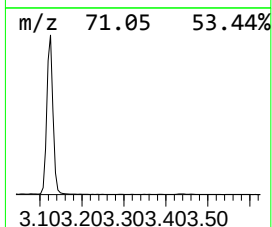
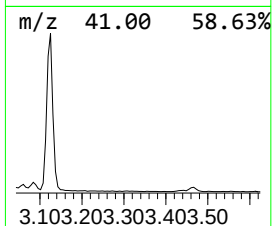
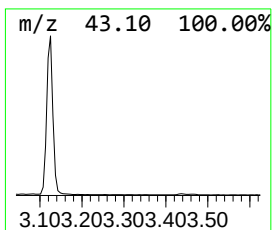
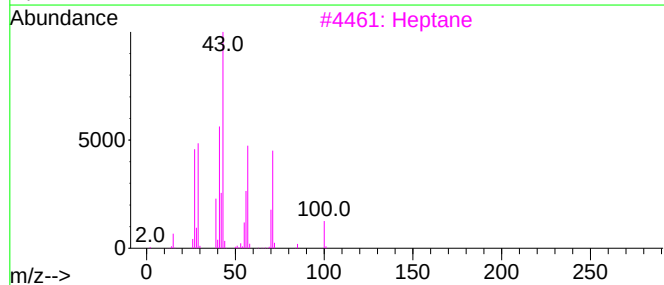
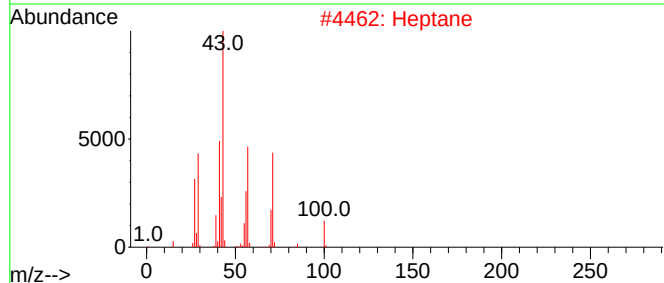
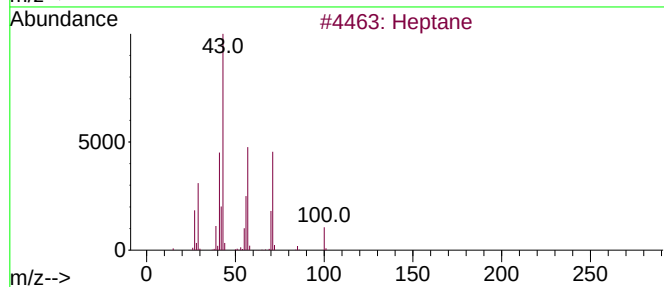
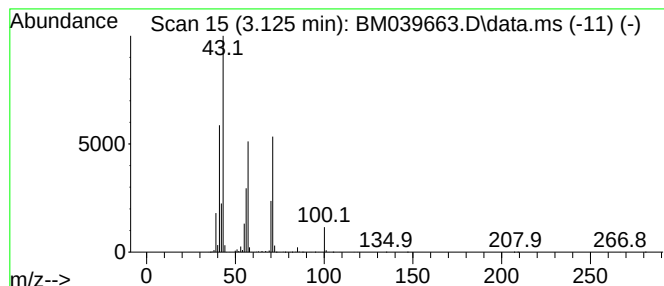
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.125	18.07 ng/ul	1462220	1,4-Dichlorobenzene-d4	7.807

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane			100	C7H16	000142-82-5	95
2	Heptane			100	C7H16	000142-82-5	94
3	Heptane			100	C7H16	000142-82-5	87
4	Heptane			100	C7H16	000142-82-5	83
5	Oxalic acid, isobutyl pentyl ester			216	C11H20O4	1000309-37-0	64



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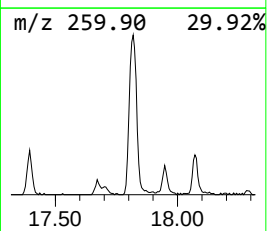
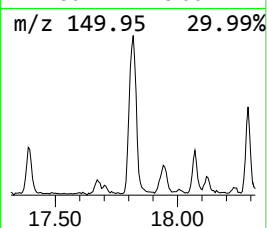
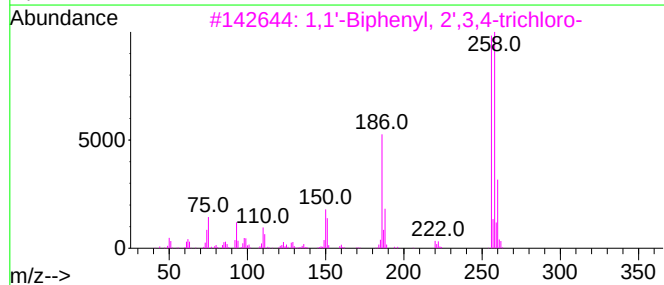
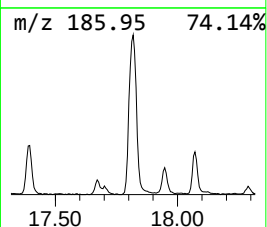
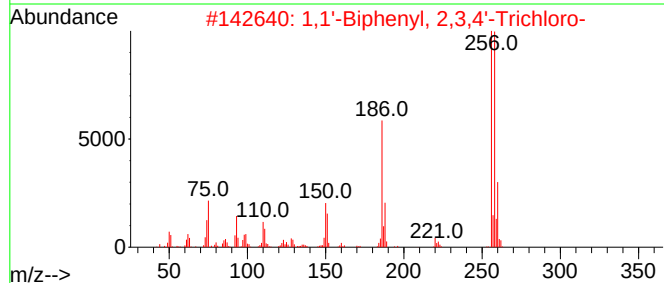
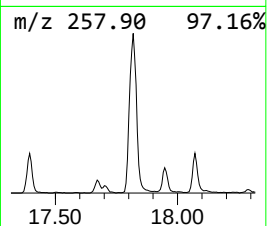
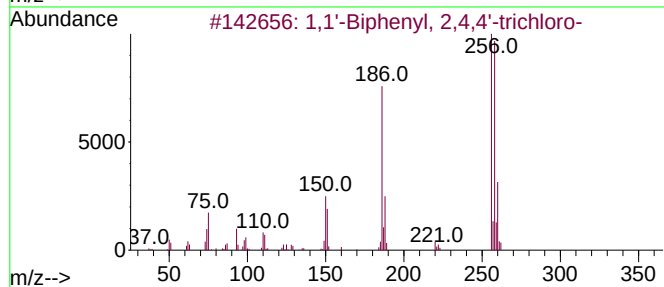
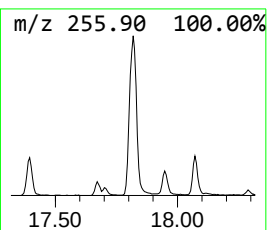
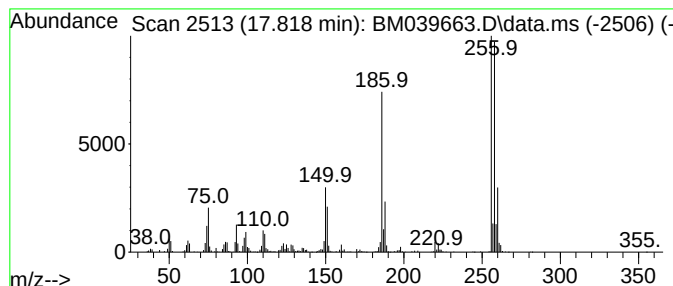
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.818	9.42 ng/ul	1028350	Phenanthrene-d10	17.218

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, 2,3,4'-Trichloro-	256	C12H7Cl3	038444-85-8	99		
3	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99		
4	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99		
5	1,1'-Biphenyl, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	99		



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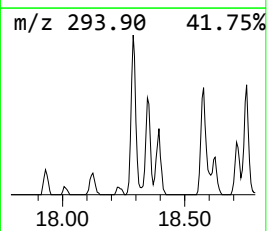
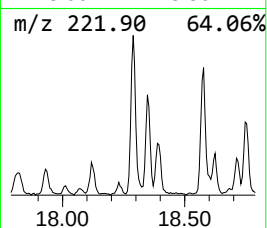
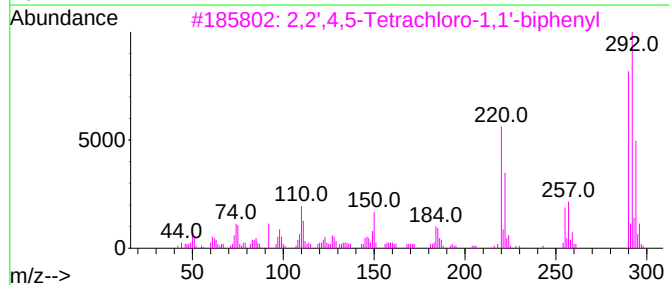
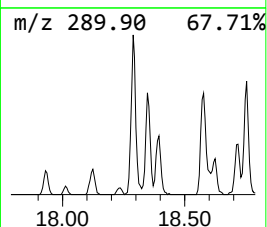
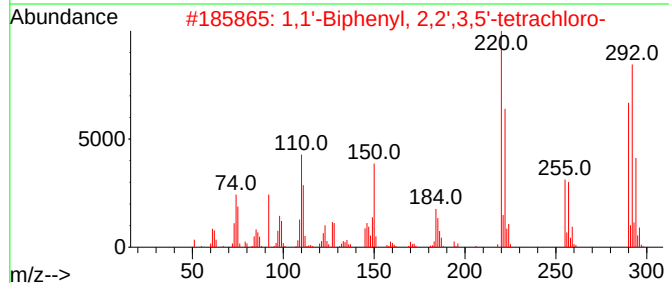
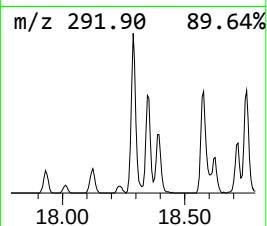
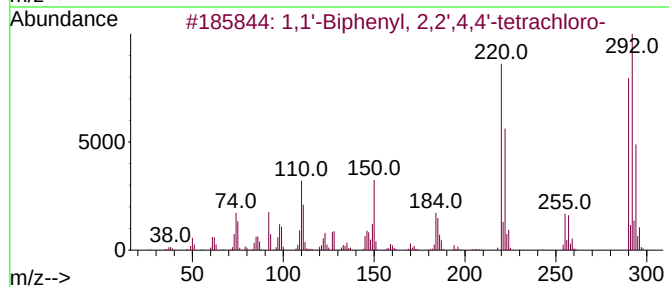
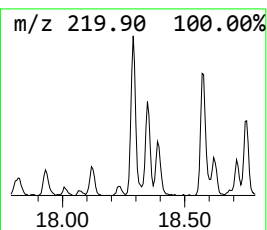
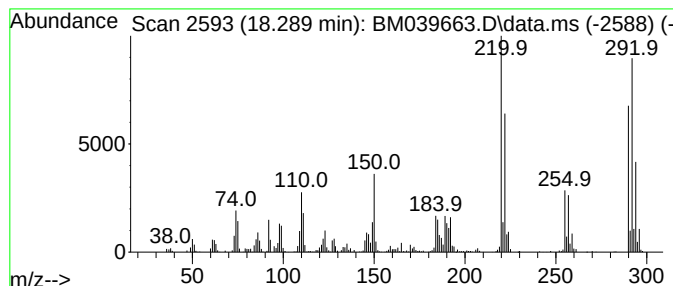
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.289	4.37 ng/ul	476873	Phenanthrene-d10	17.218

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',3,5'-tetrach...	290	C12H6Cl4	041464-39-5	99		
3	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99		
4	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99		
5	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042523\
Data File : BM039663.D
Acq On : 26 Apr 2023 05:18
Operator : CG/JU
Sample : 02419-27
Misc :
ALS Vial : 32 Sample Multiplier: 1

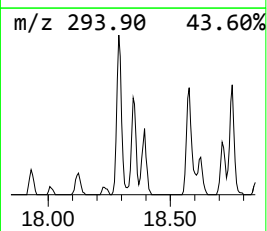
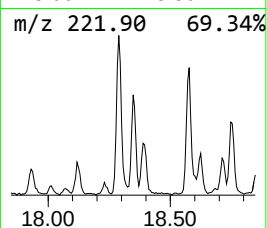
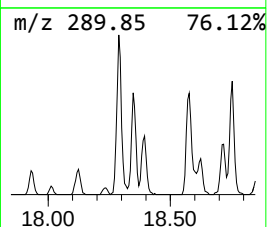
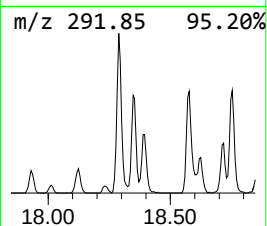
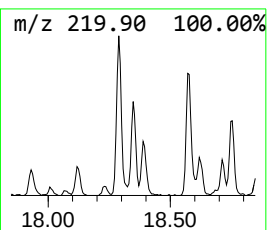
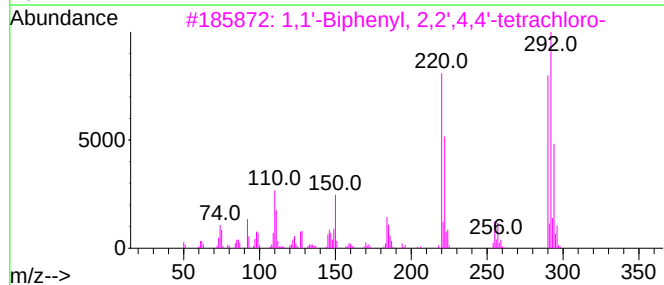
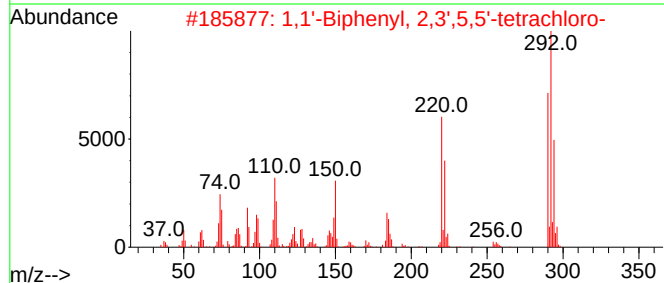
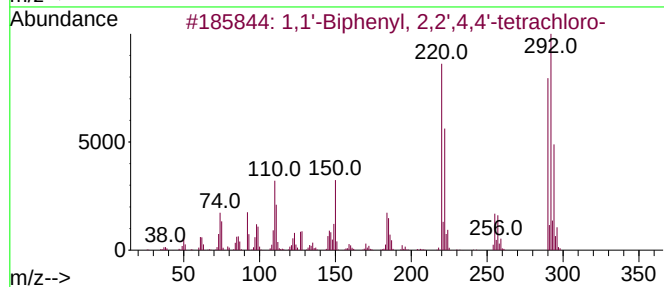
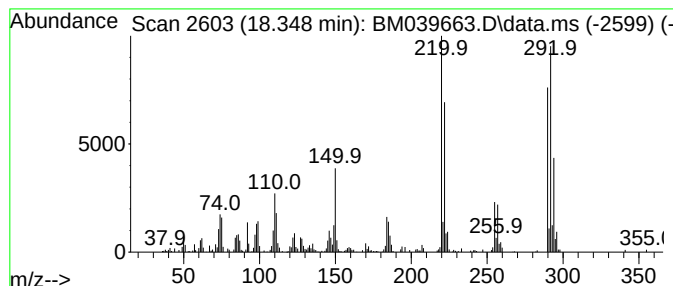
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.348	2.44 ng/ul	266750	Phenanthrene-d10	17.218	
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',5,5'-tetrach...	290	C12H6Cl4	041464-42-0	99
3	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
4	1,1'-Biphenyl, 2,3,3',4'-tetrach...	290	C12H6Cl4	041464-43-1	99
5	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042523\
Data File : BM039663.D
Acq On : 26 Apr 2023 05:18
Operator : CG/JU
Sample : 02419-27
Misc :
ALS Vial : 32 Sample Multiplier: 1

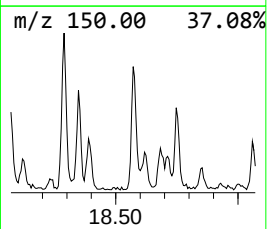
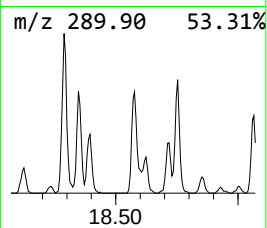
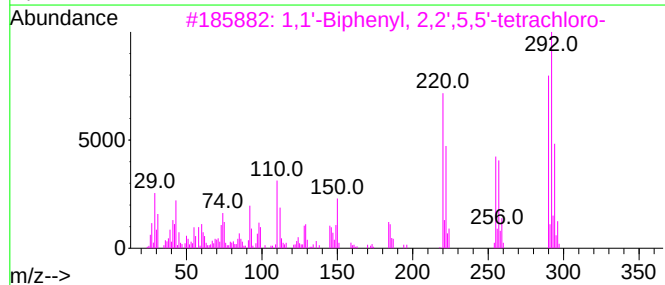
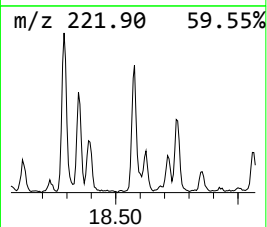
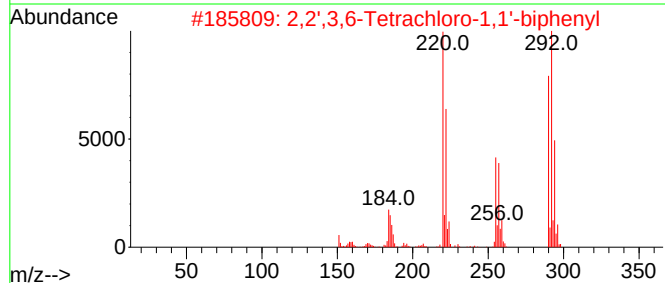
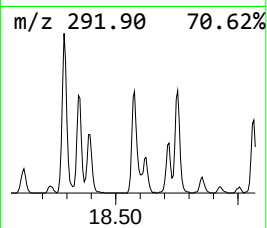
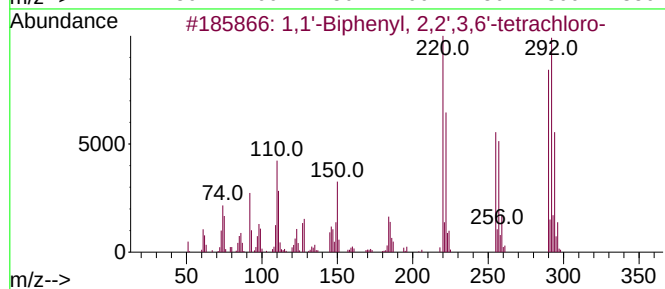
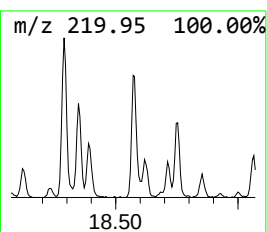
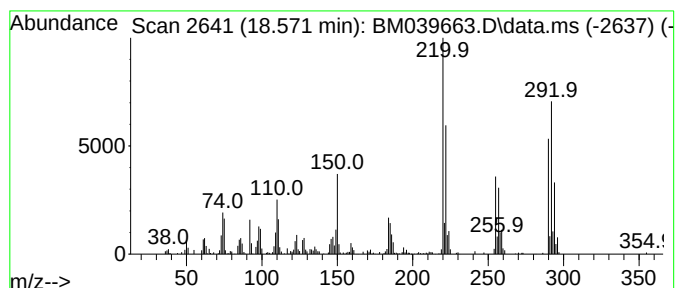
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.571	3.49 ng/ul	380667	Phenanthrene-d10	17.218	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,6'-tetrach...	290	C12H6Cl4	041464-47-5	99
2	2,2',3,6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-45-7	99
3	1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	99
4	2,2',3,6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-45-7	99
5	1,1'-Biphenyl, 2,2',4,6-Tetrachl...	290	C12H6Cl4	062796-65-0	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042523\
Data File : BM039663.D
Acq On : 26 Apr 2023 05:18
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Sample : 02419-27
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BNA_M
ClientSampleId :
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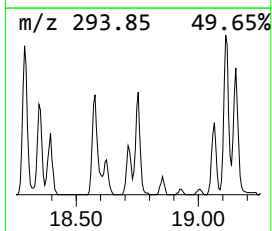
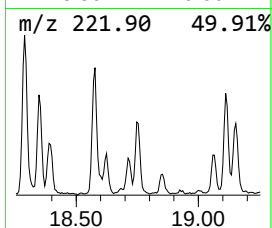
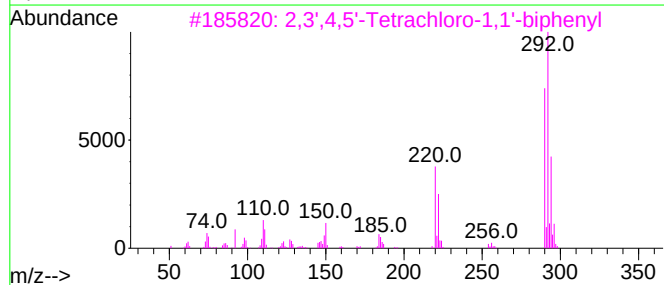
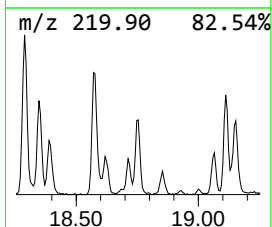
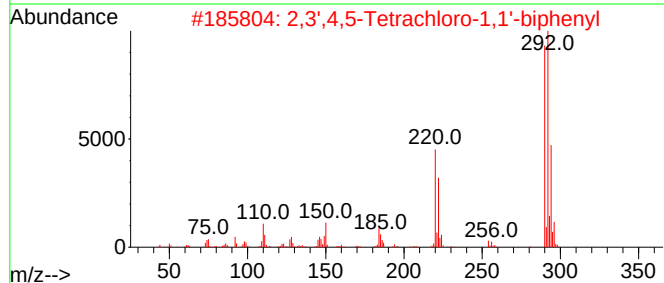
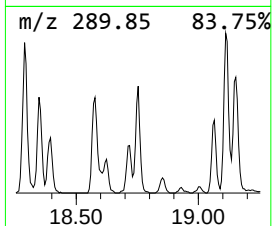
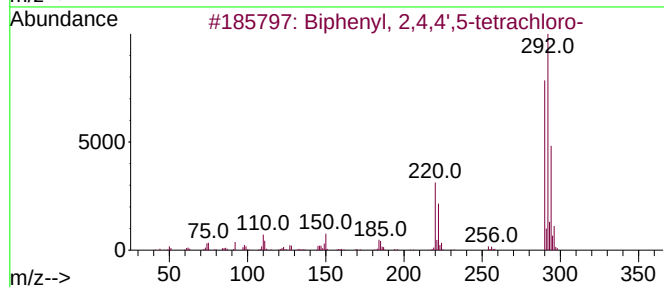
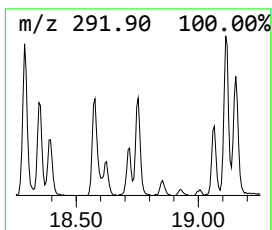
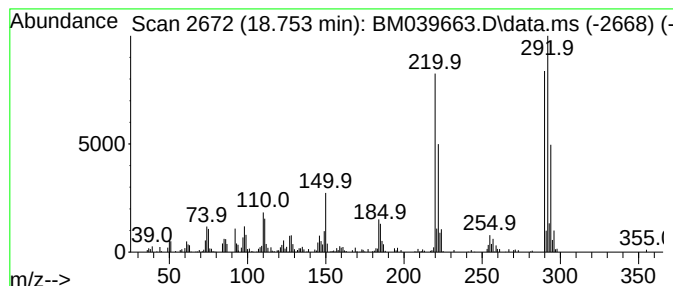
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.753	2.01 ng/ul	219531	Phenanthrene-d10	17.218

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	99
2			2,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-53-8	99
3			2,3',4,5'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-52-7	99
4			1,1'-Biphenyl, 2,3',4,6-Tetrachl...	290	C12H6Cl4	060233-24-1	99
5			3,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-49-1	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042523\
Data File : BM039663.D
Acq On : 26 Apr 2023 05:18
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ClientSampleId :
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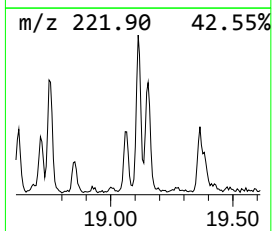
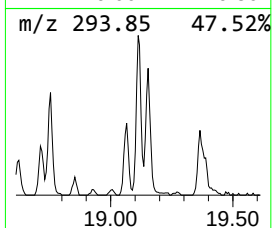
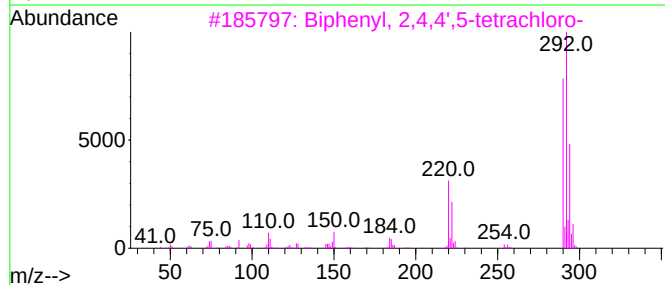
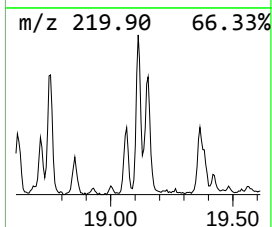
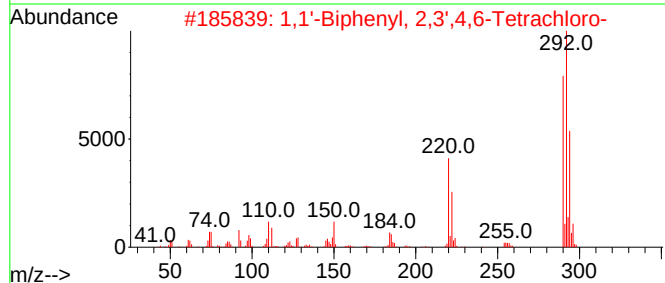
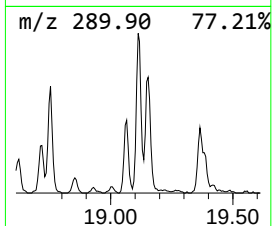
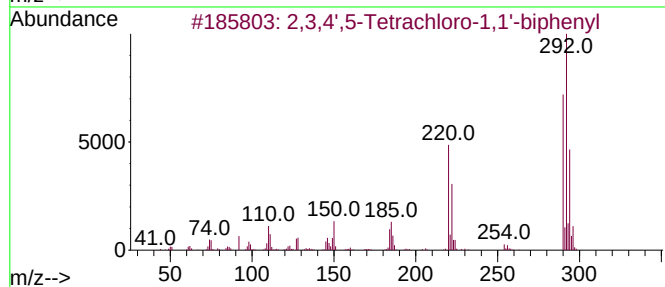
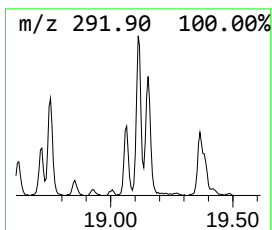
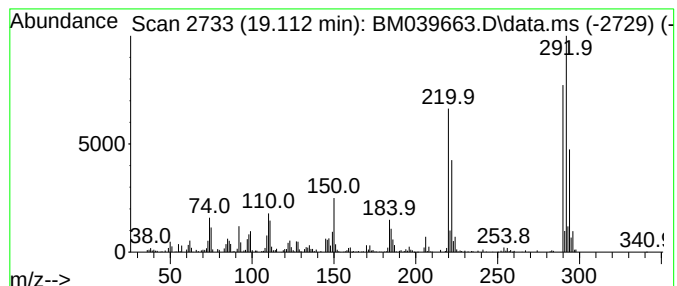
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 2,3,4',5-Tetrachloro-1,1'-b... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.112	2.72 ng/ul	296503	Phenanthrene-d10	17.218

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	1,1'-Biphenyl, 2,3',4,6-Tetrachl...	290	C12H6Cl4	060233-24-1	99		
3	Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	99		
4	1,1'-Biphenyl, 3,3',4,5'-tetrach...	290	C12H6Cl4	041464-48-6	99		
5	1,1'-Biphenyl, 2,3,3',5'-tetrach...	290	C12H6Cl4	041464-49-7	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042523\
Data File : BM039663.D
Acq On : 26 Apr 2023 05:18
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Sample : 02419-27
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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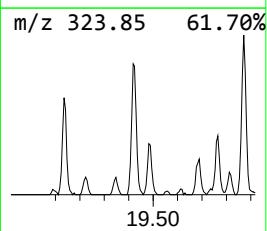
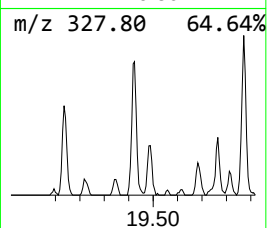
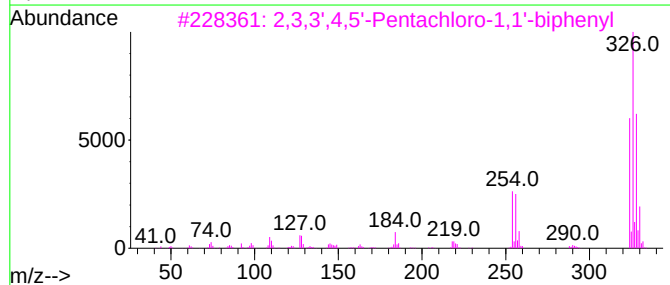
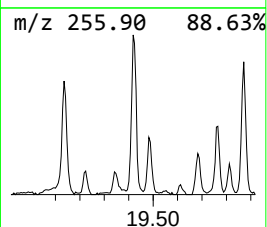
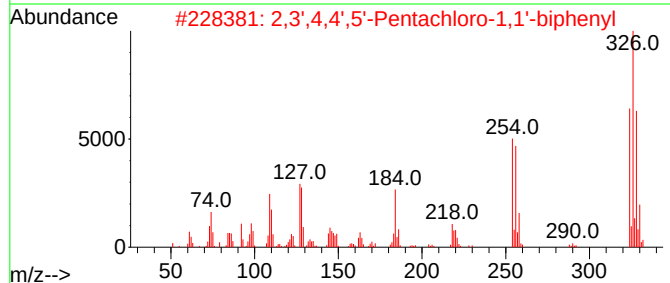
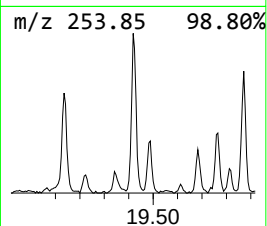
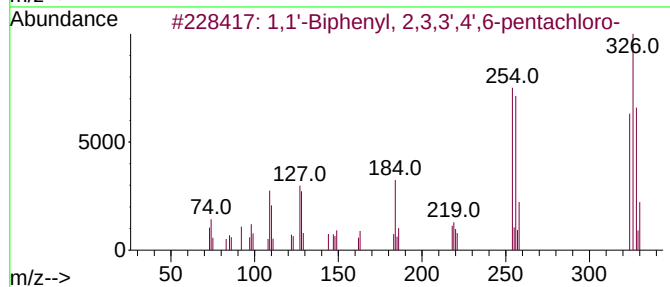
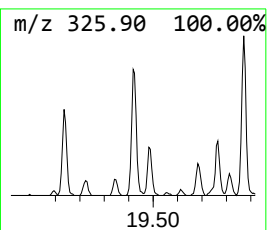
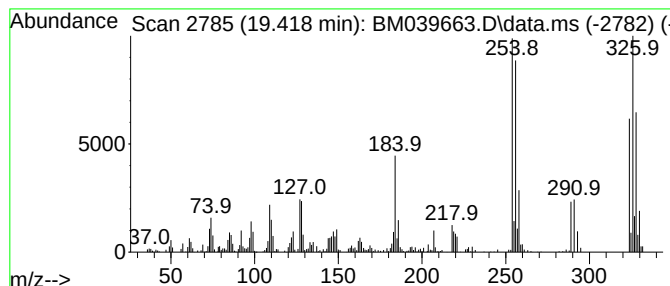
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.418	3.35 ng/ul	326865	Chrysene-d12	21.412

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	2,3,3',4,4',5'-Pentachloro-1,1'-bi...	324	C12H5Cl5	065510-44-3	99		
3	2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	070362-41-3	99		
4	1,1'-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	99		
5	1,1'-Biphenyl, 2',3,4,5,5'-Penta...	324	C12H5Cl5	070424-70-3	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042523\
Data File : BM039663.D
Acq On : 26 Apr 2023 05:18
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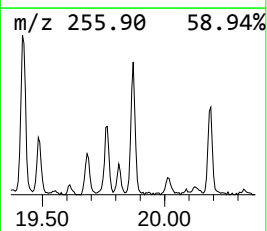
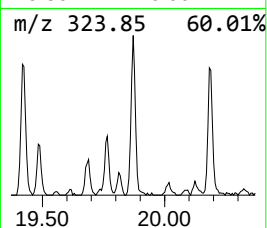
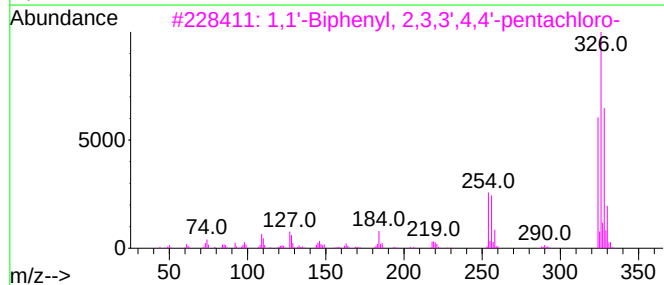
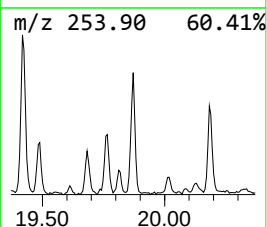
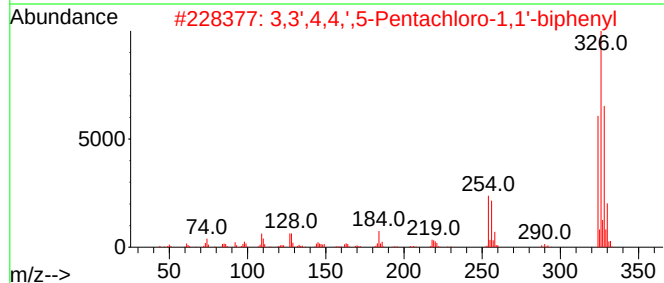
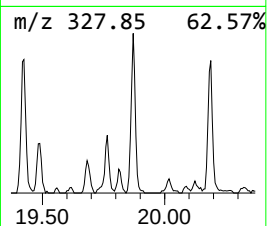
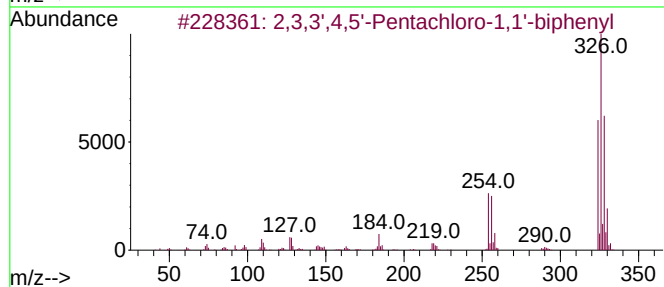
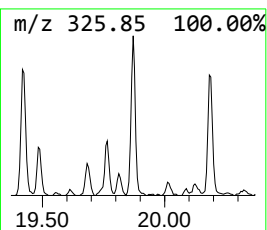
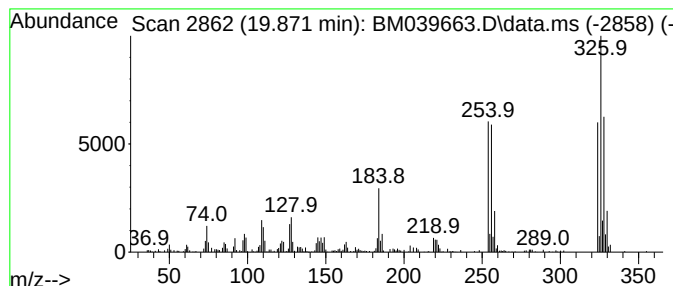
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 9 2,3,3',4,5'-Pentachloro-1,1... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.871	2.51 ng/ul	244710	Chrysene-d12	21.412	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	070362-41-3	99
2	3,3',4,4',5-Pentachloro-1,1'-bi...	324	C12H5Cl5	057465-28-8	99
3	1,1'-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	99
4	1,1'-Biphenyl, 2,2',3,3',6-penta...	324	C12H5Cl5	052663-60-2	99
5	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	98



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042523\
Data File : BM039663.D
Acq On : 26 Apr 2023 05:18
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Sample : 02419-27
Misc :
ALS Vial : 32 Sample Multiplier: 1

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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
(DEL) Alkane: S...	3.125	18.1	ng/ul	1462220	1	7.807	1618150	20.0
1,1'-Biphenyl, ...	17.818	9.4	ng/ul	1028350	4	17.218	2182480	20.0
1,1'-Biphenyl, ...	18.289	4.4	ng/ul	476873	4	17.218	2182480	20.0
1,1'-Biphenyl, ...	18.348	2.4	ng/ul	266750	4	17.218	2182480	20.0
1,1'-Biphenyl, ...	18.571	3.5	ng/ul	380667	4	17.218	2182480	20.0
Biphenyl, 2,4,4...	18.753	2.0	ng/ul	219531	4	17.218	2182480	20.0
2,3,4',5-Tetrac...	19.112	2.7	ng/ul	296503	4	17.218	2182480	20.0
1,1'-Biphenyl, ...	19.418	3.4	ng/ul	326865	5	21.412	1952580	20.0
2,3,3',4,5'-Pen...	19.871	2.5	ng/ul	244710	5	21.412	1952580	20.0