

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
 Data File : BG062982.D
 Acq On : 20 Sep 2024 17:25
 Operator : JU/RC
 Sample : AR1660-50PPM
 Misc : GCMS CONFIRMATION
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 AR1660-50PPM

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-BG091724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG062982.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.163	9	13	15	rBV	58209	60518	6.00%	0.956%
2	3.252	25	28	34	rVB	238887	257606	25.55%	4.069%
3	3.457	61	63	66	rVB	4619	4585	0.45%	0.072%
4	3.498	66	70	73	rBV	3366	5198	0.52%	0.082%
5	3.586	83	85	89	rVB2	3982	3996	0.40%	0.063%
6	4.350	213	215	219	rVB2	4239	3748	0.37%	0.059%
7	4.867	302	303	305	rBV2	2439	2378	0.24%	0.038%
8	6.013	497	498	501	rBV2	3482	2606	0.26%	0.041%
9	6.066	504	507	513	rVB	3053	6238	0.62%	0.099%
10	6.219	531	533	535	rBV	2403	2489	0.25%	0.039%
11	6.947	654	657	661	rVB	2199	3468	0.34%	0.055%
12	7.852	804	811	817	rBV	183372	298358	29.59%	4.712%
13	8.252	874	879	880	rBV	2117	2739	0.27%	0.043%
14	8.868	982	984	989	rBV	2195	3696	0.37%	0.058%
15	9.527	1093	1096	1097	rBV	2515	2796	0.28%	0.044%
16	10.431	1248	1250	1253	rBV	1836	2543	0.25%	0.040%
17	10.643	1279	1286	1293	rBV	230056	407376	40.40%	6.434%
18	11.048	1351	1355	1356	rBV	2957	2729	0.27%	0.043%
19	11.160	1372	1374	1378	rBV	1481	2438	0.24%	0.039%
20	11.671	1457	1461	1464	rBV	1669	2440	0.24%	0.039%
21	13.199	1718	1721	1724	rBV	1561	2407	0.24%	0.038%
22	14.485	1934	1940	1947	rBV	358263	563331	55.87%	8.897%
23	14.615	1960	1962	1967	rVB2	5589	6280	0.62%	0.099%
24	14.750	1982	1985	1988	rVB2	2640	2471	0.25%	0.039%
25	15.314	2075	2081	2084	rBV4	7877	13098	1.30%	0.207%
26	15.696	2143	2146	2148	rVB4	2410	2786	0.28%	0.044%
27	15.743	2148	2154	2160	rBV3	30308	48594	4.82%	0.767%
28	16.183	2225	2229	2233	rBV2	10927	13318	1.32%	0.210%
29	16.348	2254	2257	2264	rVB2	14554	22558	2.24%	0.356%
30	16.465	2272	2277	2286	rVB3	71094	100295	9.95%	1.584%
31	16.601	2299	2300	2306	rVB3	3099	4989	0.49%	0.079%
32	16.665	2309	2311	2315	rVV	2651	2781	0.28%	0.044%
33	16.771	2325	2329	2332	rVB3	9180	12747	1.26%	0.201%
34	17.006	2367	2369	2371	rBV	3783	4017	0.40%	0.063%
35	17.118	2382	2388	2391	rBV	94260	133761	13.27%	2.113%
36	17.147	2391	2393	2398	rVB3	34535	45517	4.51%	0.719%

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Integration Parameters: LSCINT.P

Integrator: RTE

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Filtering: 5

Sampling : 1

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
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37	17.235	2402	2408	2417	rVB	517352	783597	77.72%	12.376%
38	17.411	2433	2438	2443	rBV3	51139	79660	7.90%	1.258%
39	17.535	2456	2459	2460	rBV	2718	2892	0.29%	0.046%
40	17.693	2482	2486	2489	rBV3	13494	19334	1.92%	0.305%
41	17.834	2503	2510	2517	rBV2	109751	228678	22.68%	3.612%
42	17.964	2525	2532	2540	rBV2	65638	103551	10.27%	1.635%
43	18.028	2540	2543	2547	rVB3	2584	3591	0.36%	0.057%
44	18.087	2549	2553	2559	rVB3	30516	43828	4.35%	0.692%
45	18.140	2559	2562	2566	rBV4	12759	14994	1.49%	0.237%
46	18.252	2577	2581	2585	rBV3	5643	8079	0.80%	0.128%
47	18.310	2585	2591	2597	rBV4	44104	68368	6.78%	1.080%
48	18.369	2597	2601	2605	rVV4	31886	46432	4.61%	0.733%
49	18.410	2605	2608	2613	rVB4	28037	37438	3.71%	0.591%
50	18.487	2620	2621	2624	rVB	4080	3221	0.32%	0.051%
51	18.592	2634	2639	2643	rBV3	41274	61741	6.12%	0.975%
52	18.698	2654	2657	2660	rBV4	5889	8114	0.80%	0.128%
53	18.733	2660	2663	2666	rVV3	11414	15063	1.49%	0.238%
54	18.775	2666	2670	2675	rVB4	27085	35829	3.55%	0.566%
55	18.874	2684	2687	2689	rBV3	6282	5167	0.51%	0.082%
56	19.156	2727	2735	2740	rBV7	31353	50366	5.00%	0.795%
57	19.444	2780	2784	2790	rVB4	32209	42150	4.18%	0.666%
58	19.879	2853	2858	2859	rBV4	13894	21469	2.13%	0.339%
59	19.944	2868	2869	2872	rBV2	4682	4127	0.41%	0.065%
60	20.008	2877	2880	2885	rVB4	30835	36415	3.61%	0.575%
61	20.149	2900	2904	2909	rBV3	76204	98353	9.75%	1.553%
62	20.426	2947	2951	2957	rVB2	92196	110480	10.96%	1.745%
63	20.484	2957	2961	2965	rBV5	22534	30309	3.01%	0.479%
64	20.584	2975	2978	2985	rVB7	35387	60709	6.02%	0.959%
65	20.743	2999	3005	3011	rVB6	51425	93590	9.28%	1.478%
66	20.896	3027	3031	3037	rVB3	60122	76497	7.59%	1.208%
67	21.183	3078	3080	3085	rVB4	43784	53240	5.28%	0.841%
68	21.489	3126	3132	3136	rBV	606315	953976	94.62%	15.067%
69	21.912	3200	3204	3211	rVB8	31281	52953	5.25%	0.836%
70	24.532	3640	3650	3662	rVB2	359495	1008234	100.00%	15.924%
71	25.960	3891	3893	3897	rBV4	8323	12128	1.20%	0.192%

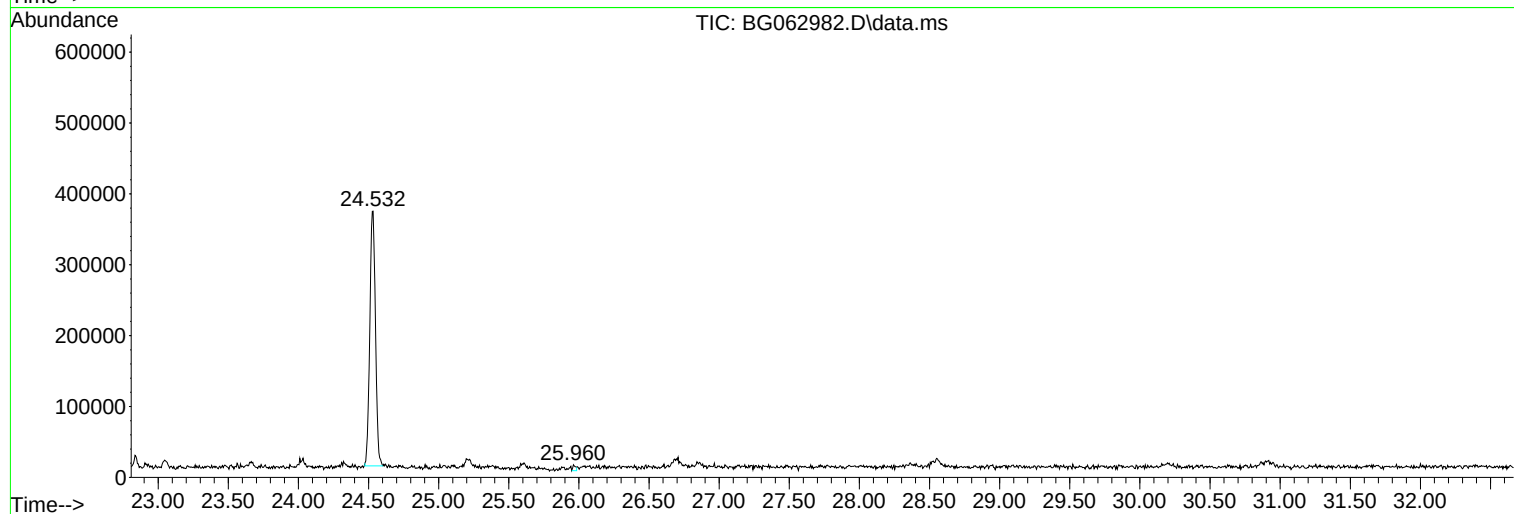
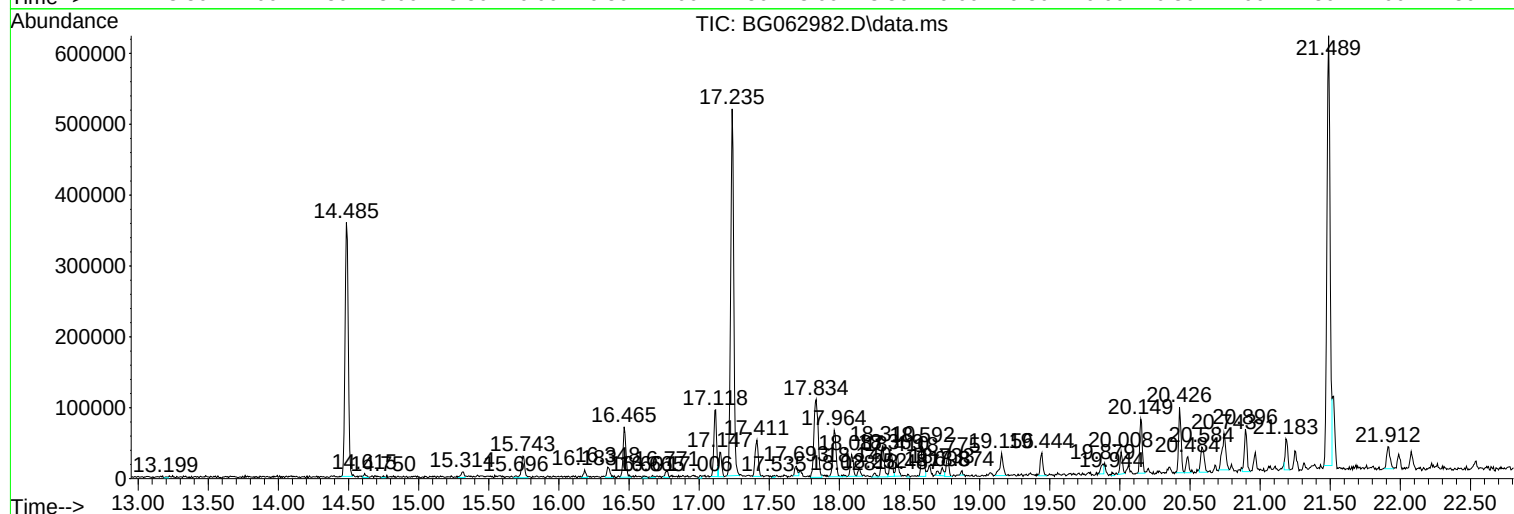
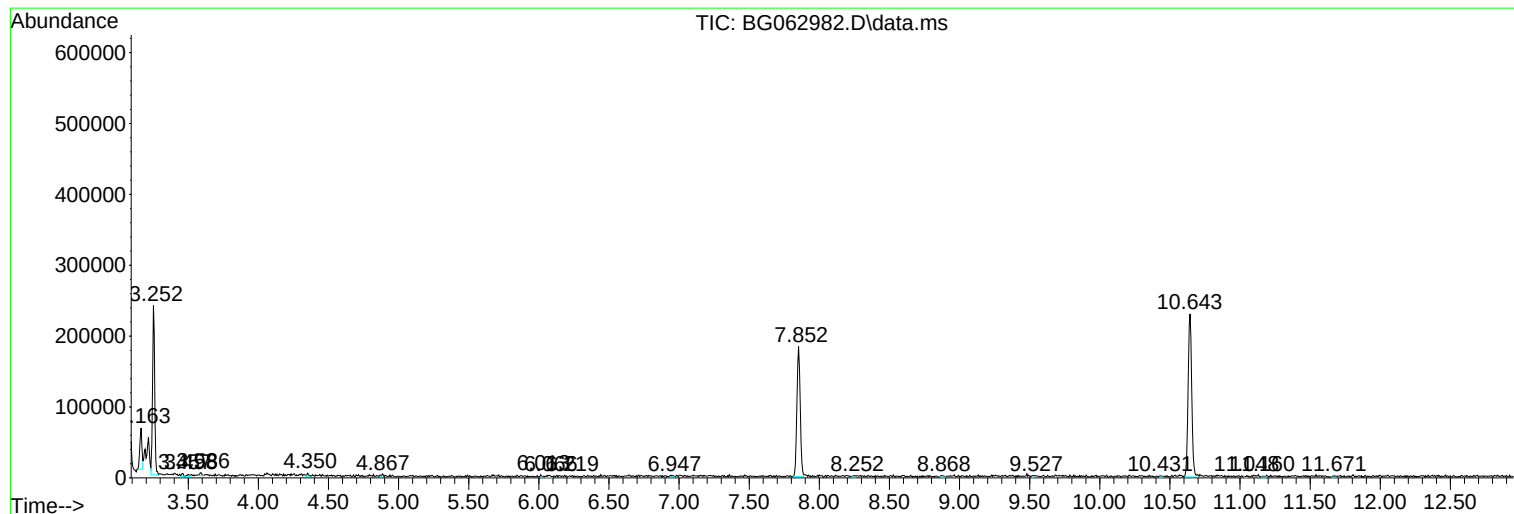
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Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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



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Operator : JU/RC
Sample : AR1660-50PPM
Misc : GCMS CONFIRMATION
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AR1660-50PPM

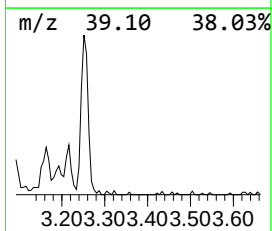
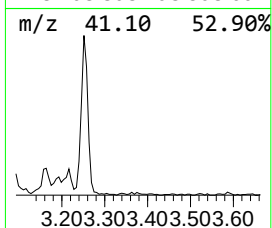
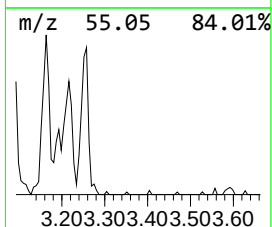
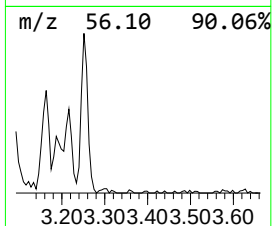
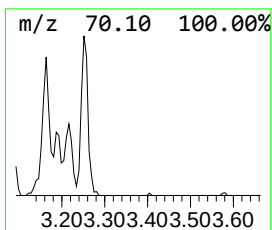
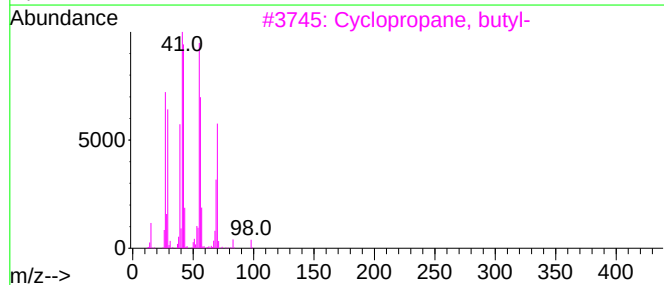
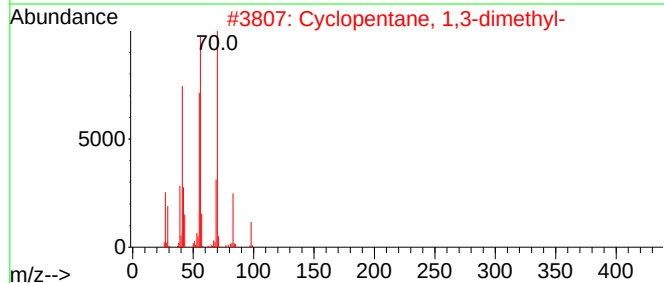
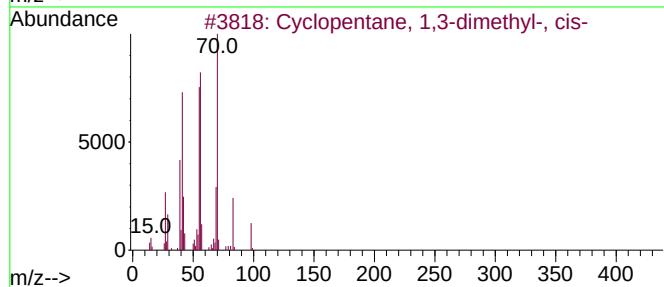
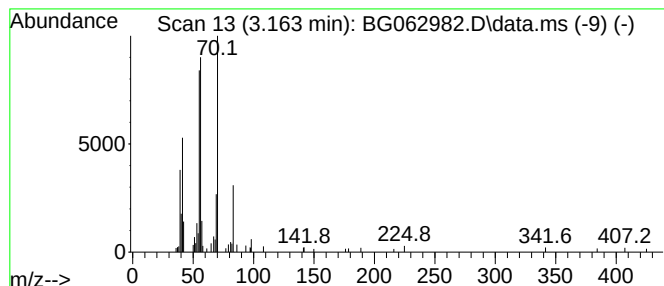
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic3.163 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.163	4.06 ng/ul	60518	1,4-Dichlorobenzene-d4	7.852

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	86
2		Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	80
3		Cyclopropane, butyl-	98	C7H14	000930-57-4	72
4		Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	72
5		Cyclopentane, 1,2,3-trimethyl-	112	C8H16	002815-57-8	72



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Misc : GCMS CONFIRMATION
ALS Vial : 46 Sample Multiplier: 1

Instrument :
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AR1660-50PPM

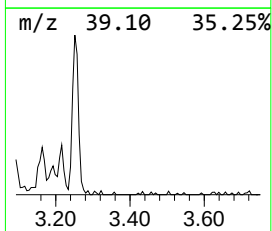
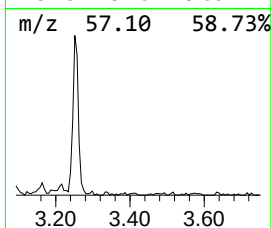
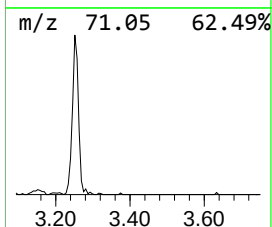
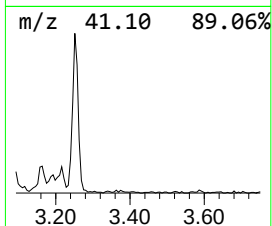
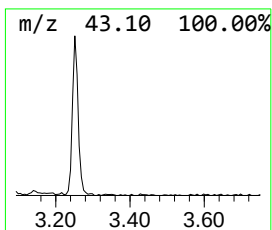
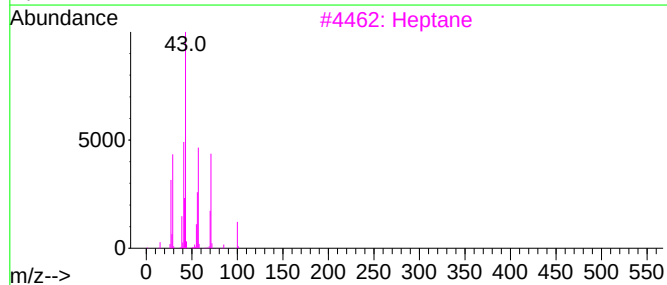
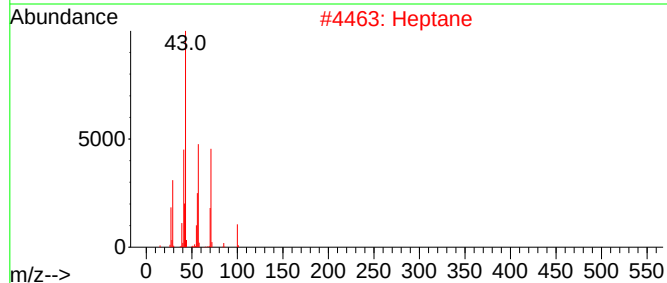
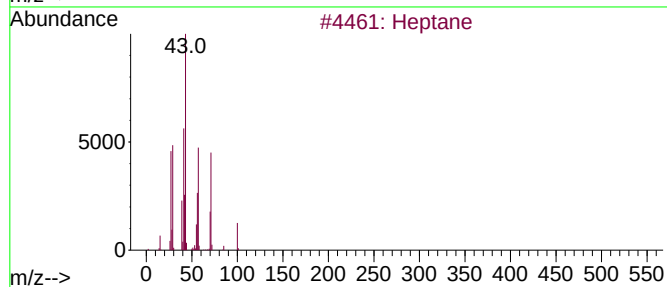
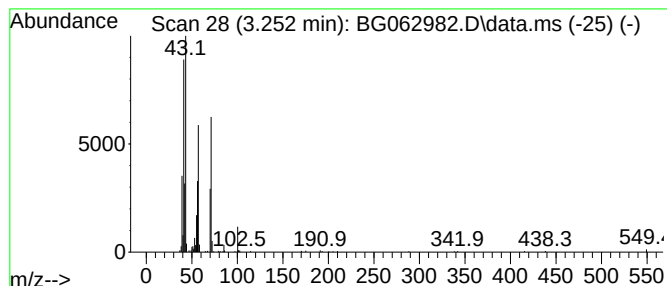
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.252	17.27 ng/ul	257606	1,4-Dichlorobenzene-d4	7.852

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane	100	C7H16	000142-82-5	80
2			Heptane	100	C7H16	000142-82-5	80
3			Heptane	100	C7H16	000142-82-5	80
4			Heptane	100	C7H16	000142-82-5	72
5			Isobutyl pentyl carbonate	188	C10H20O3	178442-32-5	50



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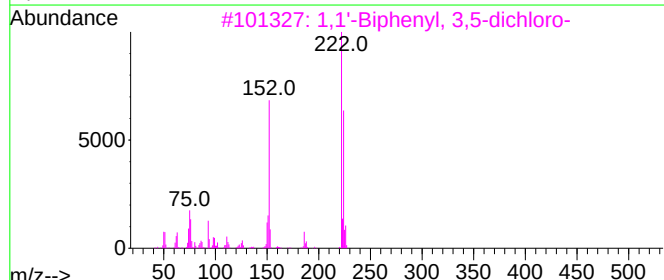
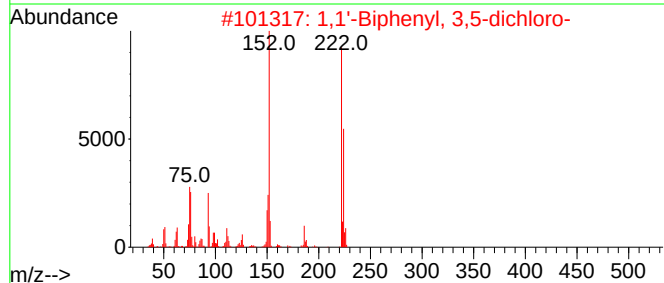
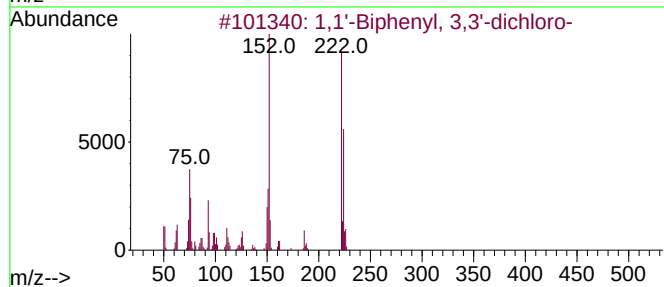
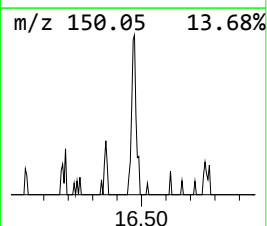
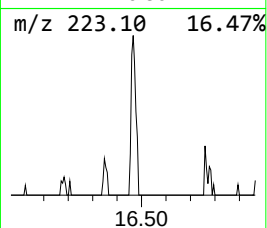
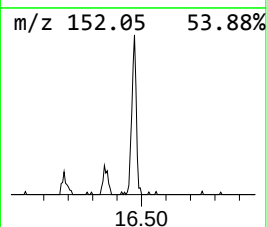
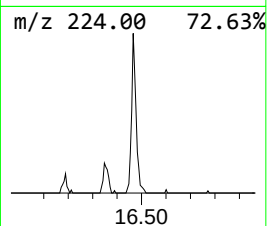
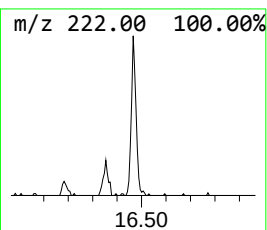
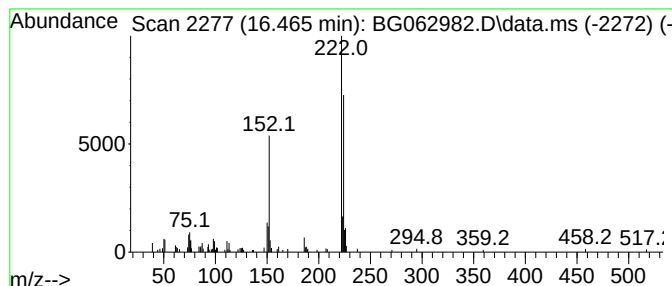
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1,1'-Biphenyl, 3,3'-dichloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.465	2.56 ng/ul	100295	Phenanthrene-d10	17.235

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 3,3'-dichloro-	222	C12H8Cl2	002050-67-1	97		
2	1,1'-Biphenyl, 3,5-dichloro-	222	C12H8Cl2	034883-41-5	97		
3	1,1'-Biphenyl, 3,5-dichloro-	222	C12H8Cl2	034883-41-5	97		
4	1,1'-Biphenyl, 4,4'-dichloro-	222	C12H8Cl2	002050-68-2	96		
5	1,1'-Biphenyl 2,3'-dichloro-	222	C12H8Cl2	025569-80-6	96		



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ClientSampleId :
AR1660-50PPM

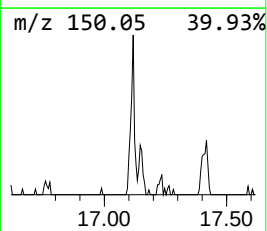
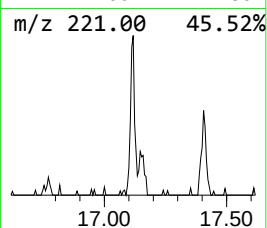
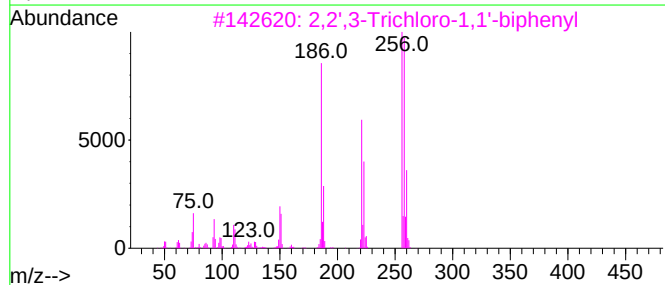
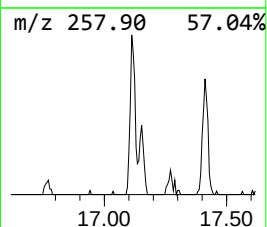
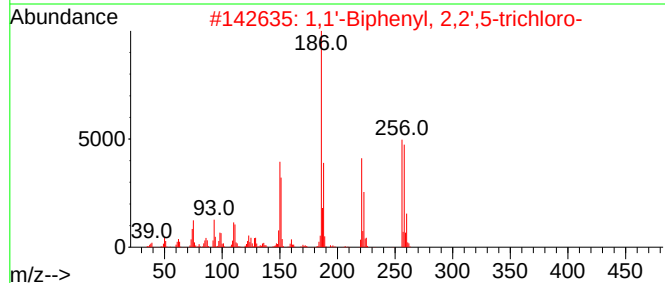
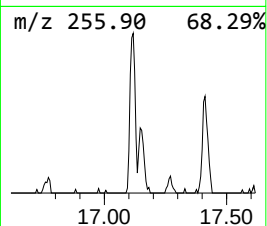
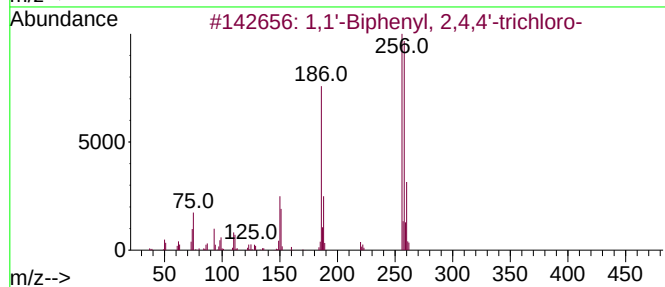
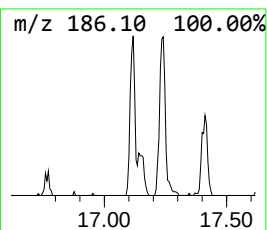
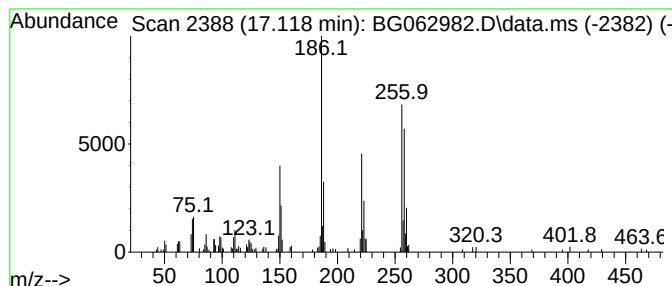
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.118	3.41 ng/ul	133761	Phenanthrene-d10	17.235

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	98		
2	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	97		
3	2,2',3-Trichloro-1,1'-biphenyl	256	C12H7Cl3	038444-78-9	97		
4	2,4',6-Trichloro-1,1'-biphenyl	256	C12H7Cl3	038444-77-8	97		
5	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	95		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062982.D
Acq On : 20 Sep 2024 17:25
Operator : JU/RC
Sample : AR1660-50PPM
Misc : GCMS CONFIRMATION
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AR1660-50PPM

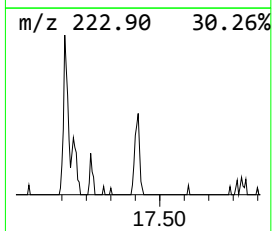
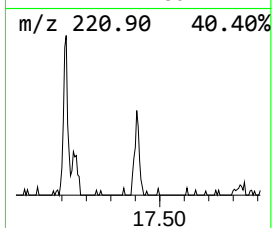
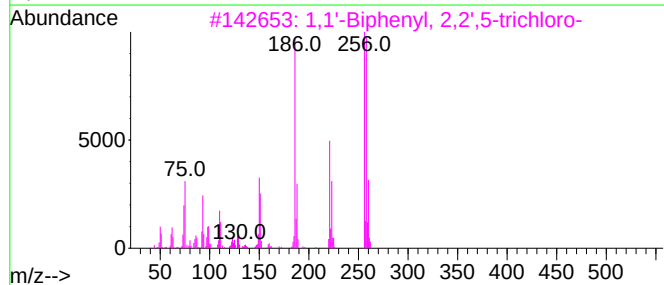
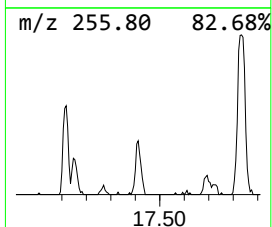
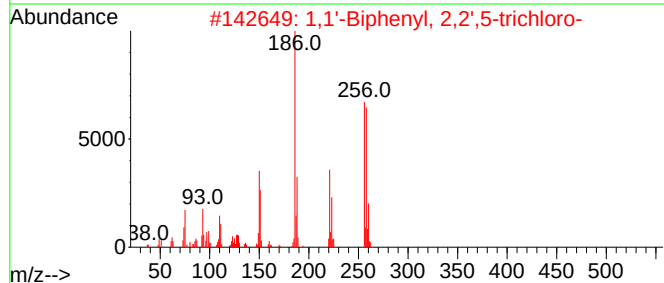
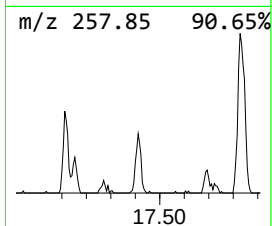
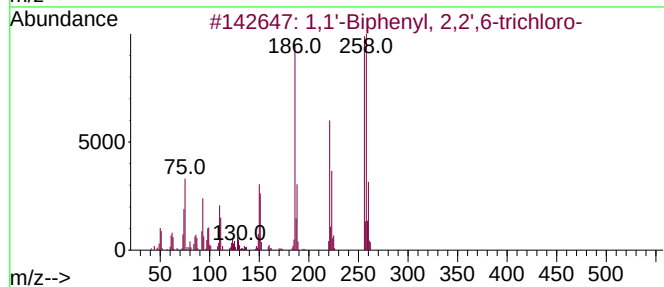
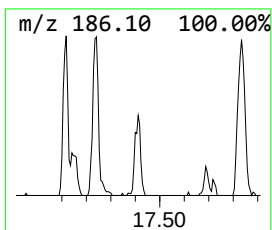
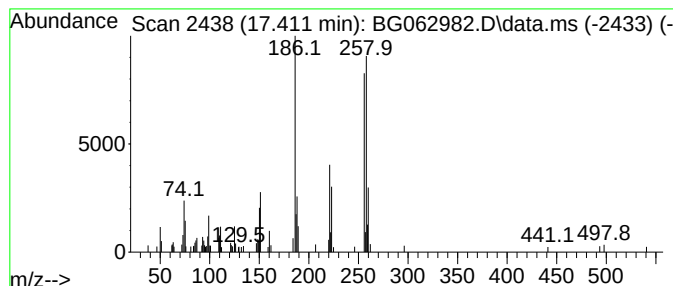
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.411	2.03 ng/ul	79660	Phenanthrene-d10	17.235

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',6-trichloro-	256	C12H7Cl3	038444-73-4	94		
2	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	94		
3	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	94		
4	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	93		
5	2,2',3-Trichloro-1,1'-biphenyl	256	C12H7Cl3	038444-78-9	93		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062982.D
Acq On : 20 Sep 2024 17:25
Operator : JU/RC
Sample : AR1660-50PPM
Misc : GCMS CONFIRMATION
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AR1660-50PPM

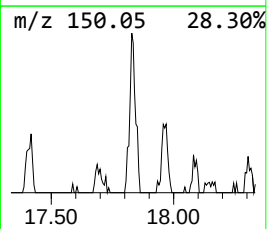
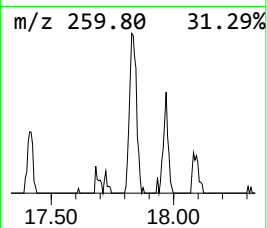
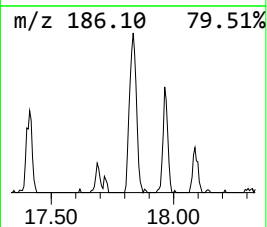
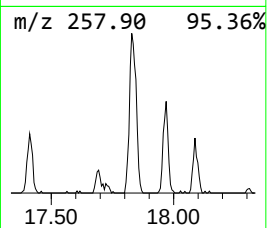
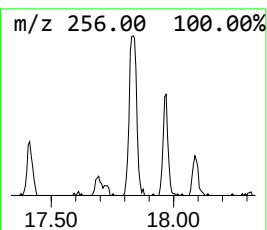
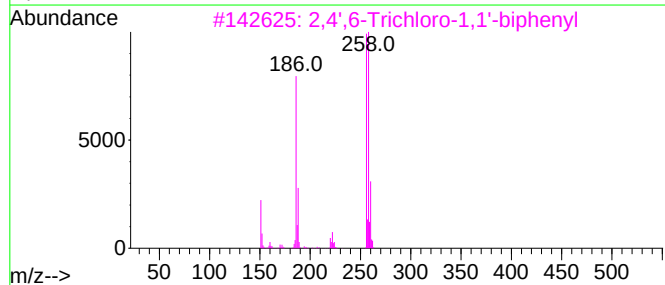
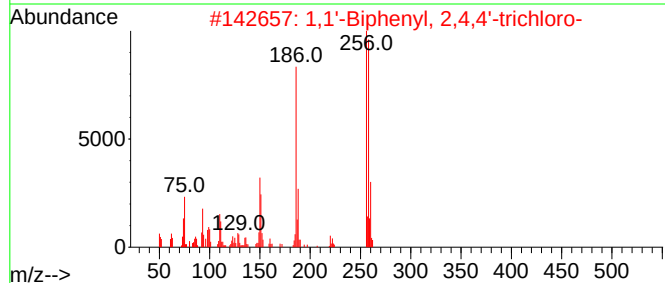
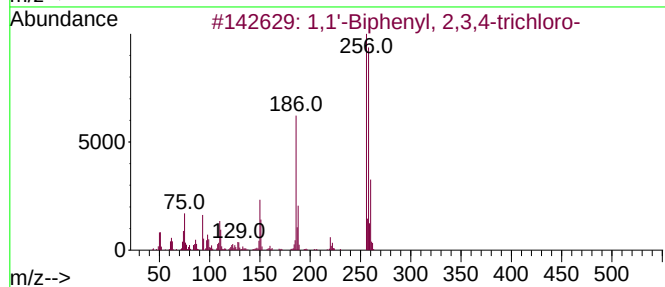
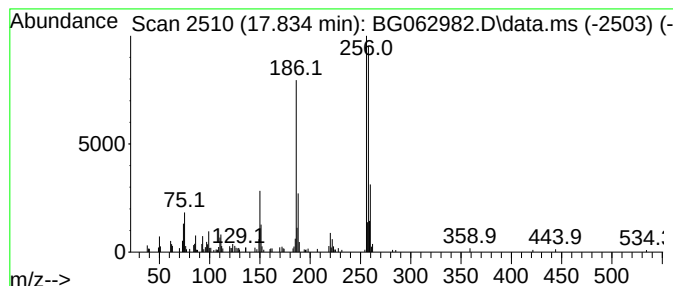
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.834	5.84 ng/ul	228678	Phenanthrene-d10	17.235

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,4-trichloro-	256	C12H7Cl3	055702-46-0	96		
2	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	96		
3	2,4',6-Trichloro-1,1'-biphenyl	256	C12H7Cl3	038444-77-8	95		
4	1,1'-Biphenyl, 2,3,6-trichloro-	256	C12H7Cl3	055702-45-9	95		
5	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	95		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062982.D
Acq On : 20 Sep 2024 17:25
Operator : JU/RC
Sample : AR1660-50PPM
Misc : GCMS CONFIRMATION
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AR1660-50PPM

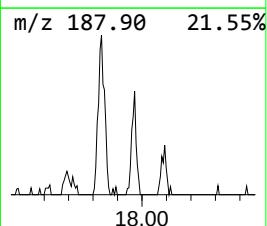
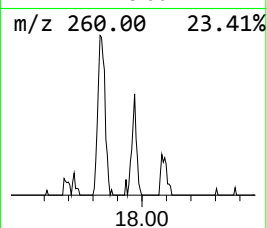
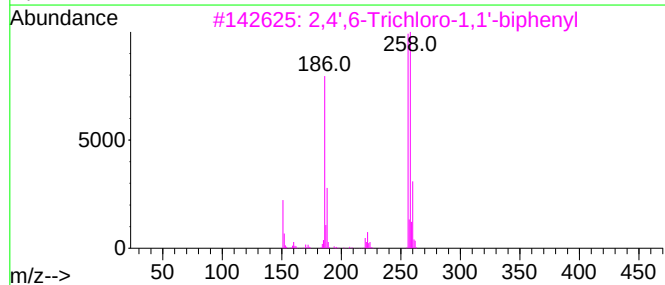
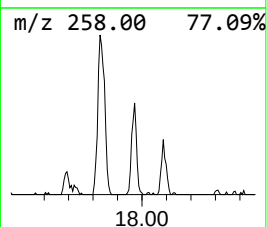
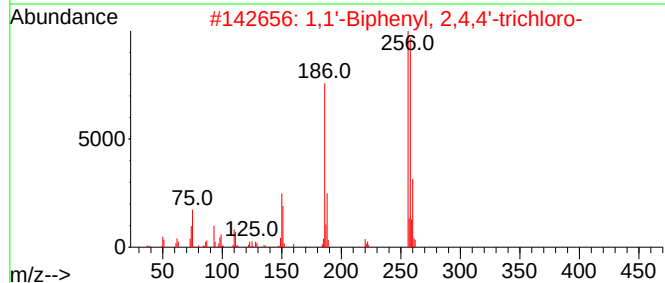
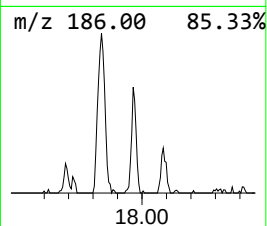
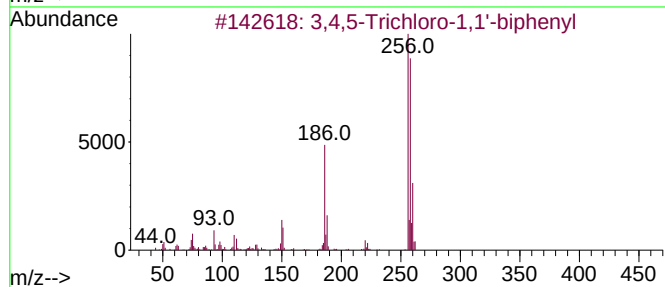
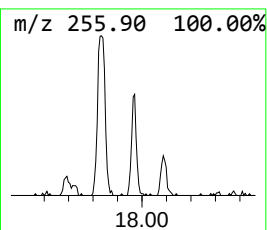
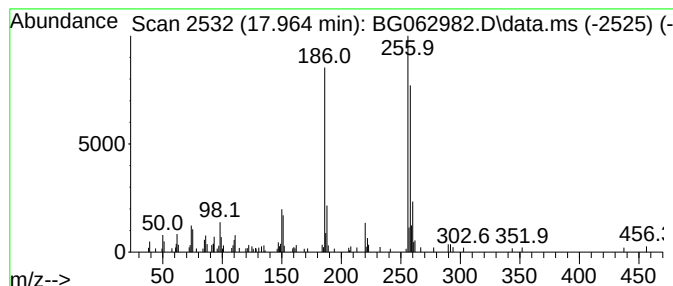
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 3,4,5-Trichloro-1,1'-biphenyl Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.964	2.64 ng/ul	103551	Phenanthrene-d10	17.235

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3,4,5-Trichloro-1,1'-biphenyl	256	C12H7Cl3	053555-66-1	97
2		1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	96
3		2,4',6-Trichloro-1,1'-biphenyl	256	C12H7Cl3	038444-77-8	95
4		1,1'-Biphenyl, 2,4,6-trichloro-	256	C12H7Cl3	035693-92-6	95
5		1,1'-Biphenyl, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062982.D
Acq On : 20 Sep 2024 17:25
Operator : JU/RC
Sample : AR1660-50PPM
Misc : GCMS CONFIRMATION
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AR1660-50PPM

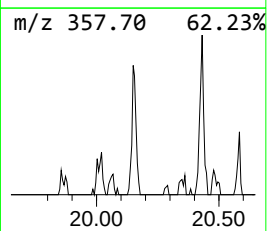
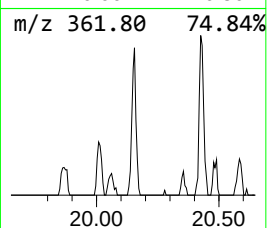
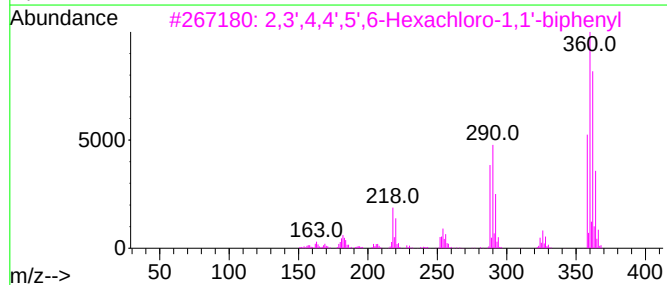
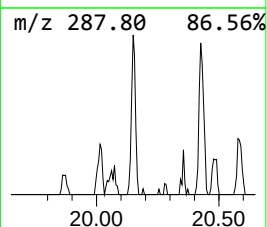
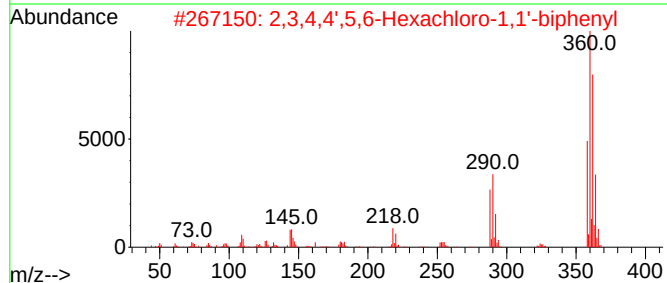
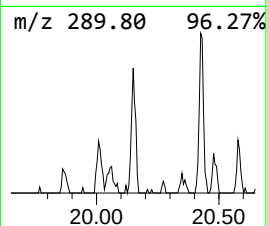
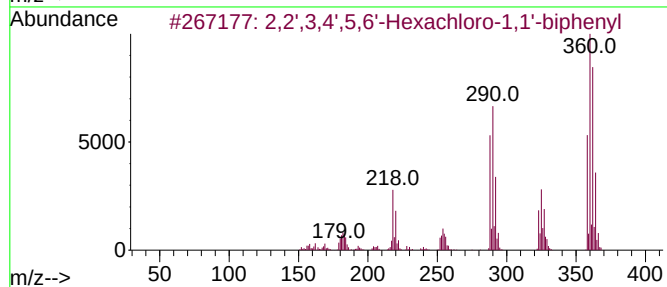
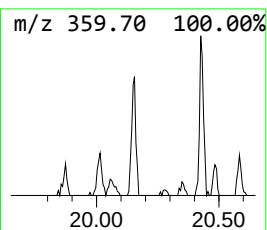
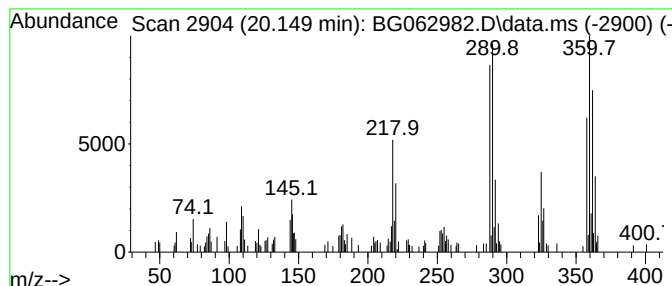
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.149	2.06 ng/ul	98353	Chrysene-d12	21.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,4',5,6'-Hexachloro-1,1'-b...	358	C12H4Cl6	074472-41-6	98		
2	2,3,4,4',5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-63-6	98		
3	2,3',4,4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-65-5	98		
4	2,3,3',4,4',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-42-7	97		
5	2,2',3,5,6,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-09-2	96		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062982.D
Acq On : 20 Sep 2024 17:25
Operator : JU/RC
Sample : AR1660-50PPM
Misc : GCMS CONFIRMATION
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AR1660-50PPM

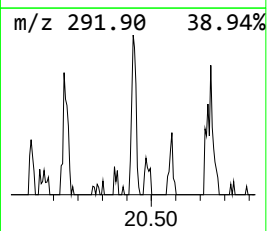
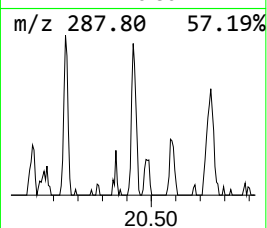
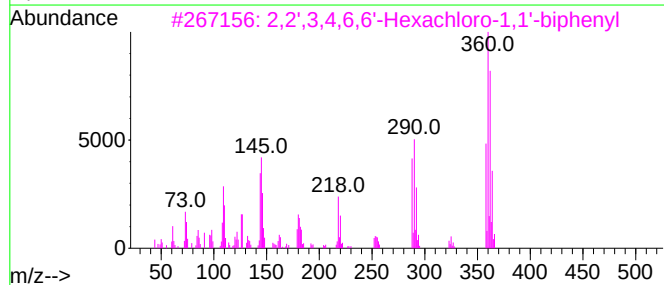
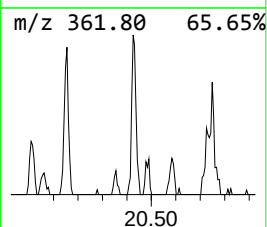
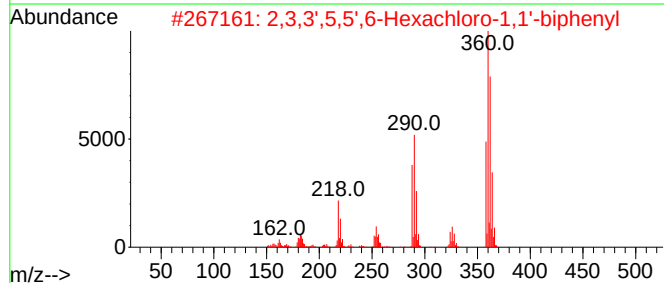
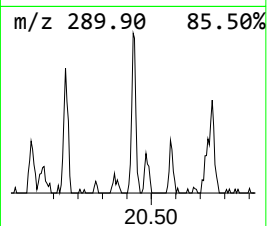
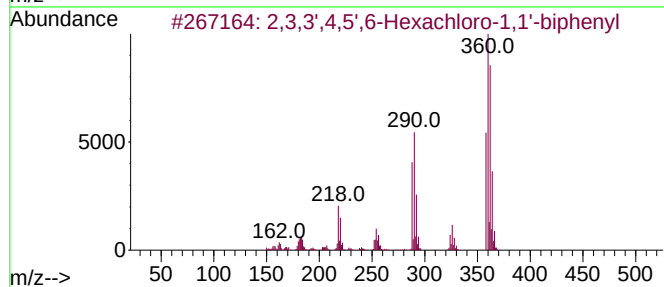
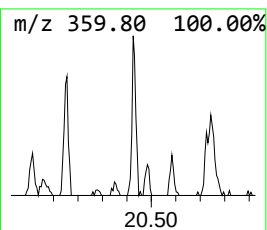
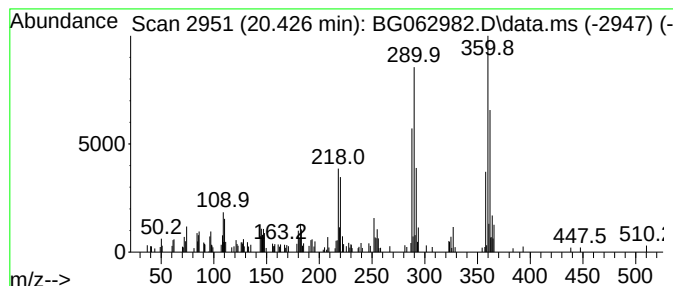
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 2,3,3',4,5',6-Hexachloro-1,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.426	2.32 ng/ul	110480	Chrysene-d12	21.489

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	96		
2	2,3,3',5,5',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-46-1	95		
3	2,2',3,4,6,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-40-5	94		
4	1,1'-Biphenyl, 2,2',4,4',6,6'-he...	358	C12H4Cl6	033979-03-2	93		
5	1,1'-Biphenyl, 2,2',3,4,5,5'-hex...	358	C12H4Cl6	052712-04-6	91		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062982.D
Acq On : 20 Sep 2024 17:25
Operator : JU/RC
Sample : AR1660-50PPM
Misc : GCMS CONFIRMATION
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AR1660-50PPM

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.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
(DEL) Alkane: C...	3.163	4.1	ng/ul	60518	1	7.852	298358	20.0
(DEL) Alkane: S...	3.252	17.3	ng/ul	257606	1	7.852	298358	20.0
1,1'-Biphenyl, ...	16.465	2.6	ng/ul	100295	4	17.235	783597	20.0
1,1'-Biphenyl, ...	17.118	3.4	ng/ul	133761	4	17.235	783597	20.0
1,1'-Biphenyl, ...	17.411	2.0	ng/ul	79660	4	17.235	783597	20.0
1,1'-Biphenyl, ...	17.834	5.8	ng/ul	228678	4	17.235	783597	20.0
3,4,5-Trichloro...	17.964	2.6	ng/ul	103551	4	17.235	783597	20.0
2,2',3,4',5,6'-...	20.149	2.1	ng/ul	98353	5	21.489	953976	20.0
2,3,3',4,5',6-H...	20.426	2.3	ng/ul	110480	5	21.489	953976	20.0