

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071023\
 Data File : BG058149.D
 Acq On : 10 Jul 2023 19:52
 Operator : CG/JU
 Sample : AR1660-50PPM
 Misc : GCMS CONFIRMATION
 ALS Vial : 14 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-BG070123.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG058149.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.238	21	25	28	rBV	13722	14568	1.52%	0.221%
2	3.267	28	30	32	rVV2	4397	4418	0.46%	0.067%
3	3.291	32	34	36	rVV	7381	6876	0.72%	0.104%
4	3.326	36	40	46	rVB	147801	160878	16.77%	2.439%
5	3.632	91	92	95	rBV2	1253	1101	0.11%	0.017%
6	3.667	95	98	102	rVB4	5047	6699	0.70%	0.102%
7	3.749	110	112	115	rVB3	1126	1109	0.12%	0.017%
8	3.961	146	148	151	rBV	1174	1224	0.13%	0.019%
9	4.149	175	180	184	rBV2	3144	4165	0.43%	0.063%
10	4.278	195	202	206	rBV2	5128	7141	0.74%	0.108%
11	4.331	206	211	220	rVB	7140	10623	1.11%	0.161%
12	4.436	224	229	232	rBV	10896	16517	1.72%	0.250%
13	4.472	232	235	240	rVV2	9803	15112	1.58%	0.229%
14	4.548	244	248	251	rVB	810	1271	0.13%	0.019%
15	5.465	402	404	409	rBV	695	1064	0.11%	0.016%
16	7.968	823	830	838	rBV	165304	281519	29.35%	4.267%
17	10.776	1300	1308	1317	rBV	232502	398534	41.55%	6.041%
18	14.601	1952	1959	1968	rBV	364545	575956	60.04%	8.731%
19	14.719	1974	1979	1983	rVB	4168	4961	0.52%	0.075%
20	15.529	2114	2117	2121	rBV2	985	1150	0.12%	0.017%
21	15.847	2166	2171	2178	rVB2	34343	46101	4.81%	0.699%
22	16.287	2242	2246	2252	rVB2	7099	10021	1.04%	0.152%
23	16.452	2269	2274	2281	rBV2	15122	20601	2.15%	0.312%
24	16.569	2289	2294	2304	rBV	80155	112770	11.76%	1.709%
25	16.869	2340	2345	2349	rBV3	8096	10863	1.13%	0.165%
26	17.122	2386	2388	2392	rVB2	763	963	0.10%	0.015%
27	17.216	2399	2404	2407	rBV	104089	145661	15.18%	2.208%
28	17.251	2407	2410	2416	rVB	40612	49804	5.19%	0.755%
29	17.345	2416	2426	2437	rBV	486346	783118	81.64%	11.871%
30	17.433	2439	2441	2449	rVV3	2040	3645	0.38%	0.055%
31	17.509	2449	2454	2463	rVB2	59823	89928	9.37%	1.363%
32	17.703	2485	2487	2492	rBV3	1335	1148	0.12%	0.017%
33	17.791	2497	2502	2504	rVV2	13907	18828	1.96%	0.285%
34	17.815	2504	2506	2511	rVB2	6905	8859	0.92%	0.134%
35	17.932	2518	2526	2537	rBV2	119781	251162	26.18%	3.807%
36	18.068	2540	2549	2554	rVV2	69399	108011	11.26%	1.637%

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37	18.138	2556	2561	2563	rVV4	2612	3436	0.36%	0.052%
38	18.185	2564	2569	2575	rVV2	32523	46579	4.86%	0.706%
39	18.244	2575	2579	2585	rVB2	12214	14792	1.54%	0.224%
40	18.350	2591	2597	2601	rBV4	4144	5054	0.53%	0.077%
41	18.402	2601	2606	2612	rVV	50161	70650	7.36%	1.071%
42	18.467	2612	2617	2621	rVV	34738	46654	4.86%	0.707%
43	18.508	2621	2624	2629	rVB2	28427	37015	3.86%	0.561%
44	18.690	2648	2655	2659	rBV	46197	63523	6.62%	0.963%
45	18.737	2659	2663	2668	rVB2	15087	20973	2.19%	0.318%
46	18.790	2668	2672	2675	rBV	9204	12413	1.29%	0.188%
47	18.826	2675	2678	2681	rVV2	10770	13034	1.36%	0.198%
48	18.861	2681	2684	2689	rVB	25846	35717	3.72%	0.541%
49	18.925	2692	2695	2697	rBV2	1693	1900	0.20%	0.029%
50	18.967	2697	2702	2705	rBV3	7521	11284	1.18%	0.171%
51	19.031	2712	2713	2714	rBV	2143	1011	0.11%	0.015%
52	19.090	2717	2723	2724	rVV5	3732	5591	0.58%	0.085%
53	19.119	2727	2728	2730	rVV2	3526	1987	0.21%	0.030%
54	19.149	2730	2733	2734	rVV3	2669	2797	0.29%	0.042%
55	19.172	2734	2737	2741	rVV3	4793	7251	0.76%	0.110%
56	19.219	2741	2745	2746	rVV3	9497	10287	1.07%	0.156%
57	19.249	2746	2750	2760	rVB3	31185	49688	5.18%	0.753%
58	19.395	2774	2775	2777	rBV2	2190	1419	0.15%	0.022%
59	19.454	2782	2785	2788	rVV4	2966	3396	0.35%	0.051%
60	19.531	2794	2798	2802	rBV2	34082	41522	4.33%	0.629%
61	19.583	2805	2807	2811	rBV5	1565	2899	0.30%	0.044%
62	19.630	2814	2815	2819	rVB4	3456	3276	0.34%	0.050%
63	19.666	2819	2821	2823	rBV3	2551	1342	0.14%	0.020%
64	19.689	2823	2825	2827	rVB3	1938	1359	0.14%	0.021%
65	19.719	2827	2830	2831	rBV2	1828	1615	0.17%	0.024%
66	19.736	2831	2833	2835	rBV2	1910	1293	0.13%	0.020%
67	19.865	2851	2855	2857	rBV4	5545	7295	0.76%	0.111%
68	19.965	2866	2872	2877	rBV5	16172	36527	3.81%	0.554%
69	20.030	2882	2883	2885	rBV2	2697	1728	0.18%	0.026%
70	20.065	2888	2889	2890	rBV	2534	1669	0.17%	0.025%
71	20.095	2890	2894	2898	rVV2	35819	40400	4.21%	0.612%
72	20.136	2898	2901	2908	rVB7	12711	20207	2.11%	0.306%
73	20.236	2913	2918	2922	rVV	86576	109298	11.39%	1.657%
74	20.283	2923	2926	2931	rVB5	7406	9555	1.00%	0.145%
75	20.359	2937	2939	2940	rBV2	3396	2210	0.23%	0.034%
76	20.430	2946	2951	2956	rBV8	12859	18416	1.92%	0.279%

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77	20.506	2960	2964	2969	rVV	99760	121899	12.71%	1.848%
78	20.565	2970	2974	2980	rVB6	24368	34750	3.62%	0.527%
79	20.670	2987	2992	2997	rBV7	36790	62664	6.53%	0.950%
80	20.759	3004	3007	3011	rVB6	7107	10362	1.08%	0.157%
81	20.829	3011	3019	3025	rBV4	53749	115376	12.03%	1.749%
82	20.882	3025	3028	3032	rVB5	9164	12347	1.29%	0.187%
83	20.982	3041	3045	3051	rVV2	64186	79403	8.28%	1.204%
84	21.046	3053	3056	3062	rVB8	27552	35961	3.75%	0.545%
85	21.281	3090	3096	3100	rBV4	46265	68754	7.17%	1.042%
86	21.346	3104	3107	3112	rVB6	26542	39250	4.09%	0.595%
87	21.411	3115	3118	3122	rVB6	8530	12082	1.26%	0.183%
88	21.440	3122	3123	3126	rBV3	4998	4784	0.50%	0.073%
89	21.528	3136	3138	3142	rBV5	7447	7484	0.78%	0.113%
90	21.599	3144	3150	3163	rBV2	489726	959269	100.00%	14.541%
91	22.028	3219	3223	3232	rVB7	36606	68519	7.14%	1.039%
92	22.110	3234	3237	3244	rVB8	21411	33251	3.47%	0.504%
93	22.204	3249	3253	3257	rVB7	20563	33833	3.53%	0.513%
94	22.991	3385	3387	3394	rVB8	16572	27408	2.86%	0.415%
95	24.795	3684	3694	3713	rVB	310397	909653	94.83%	13.789%
96	24.918	3713	3715	3718	rVB4	5567	3924	0.41%	0.059%
97	27.210	4101	4105	4106	rBV3	6676	7231	0.75%	0.110%
98	31.922	4906	4907	4909	rVB2	4622	2078	0.22%	0.031%
99	32.415	4989	4991	4994	rVB4	4475	3178	0.33%	0.048%

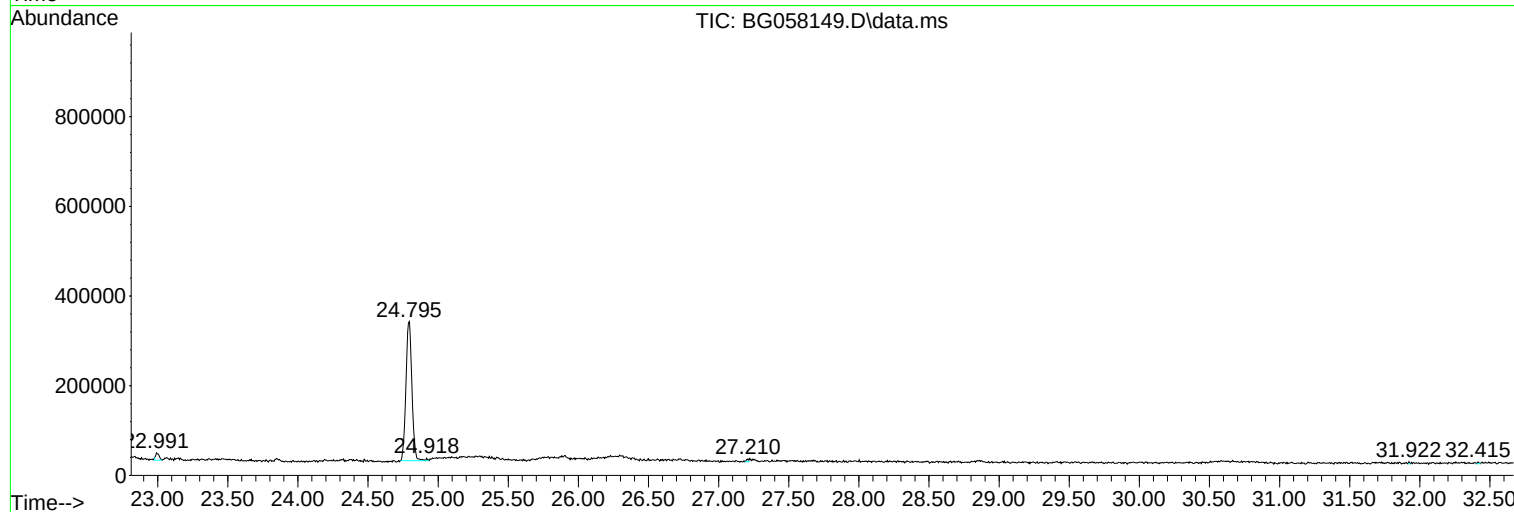
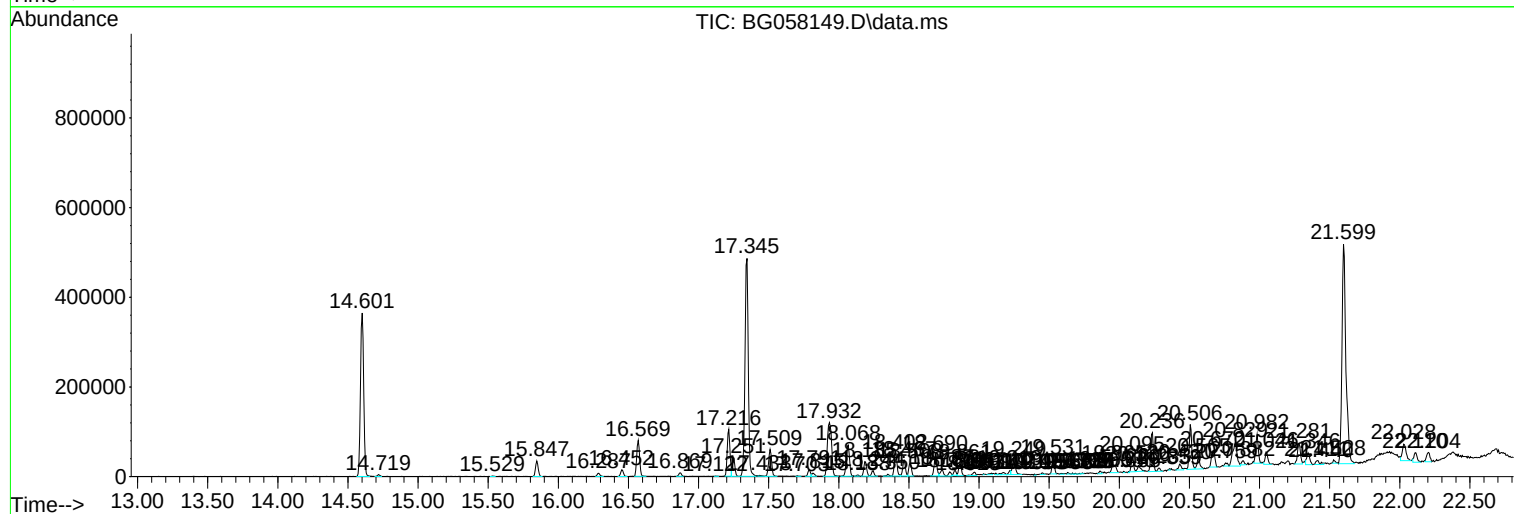
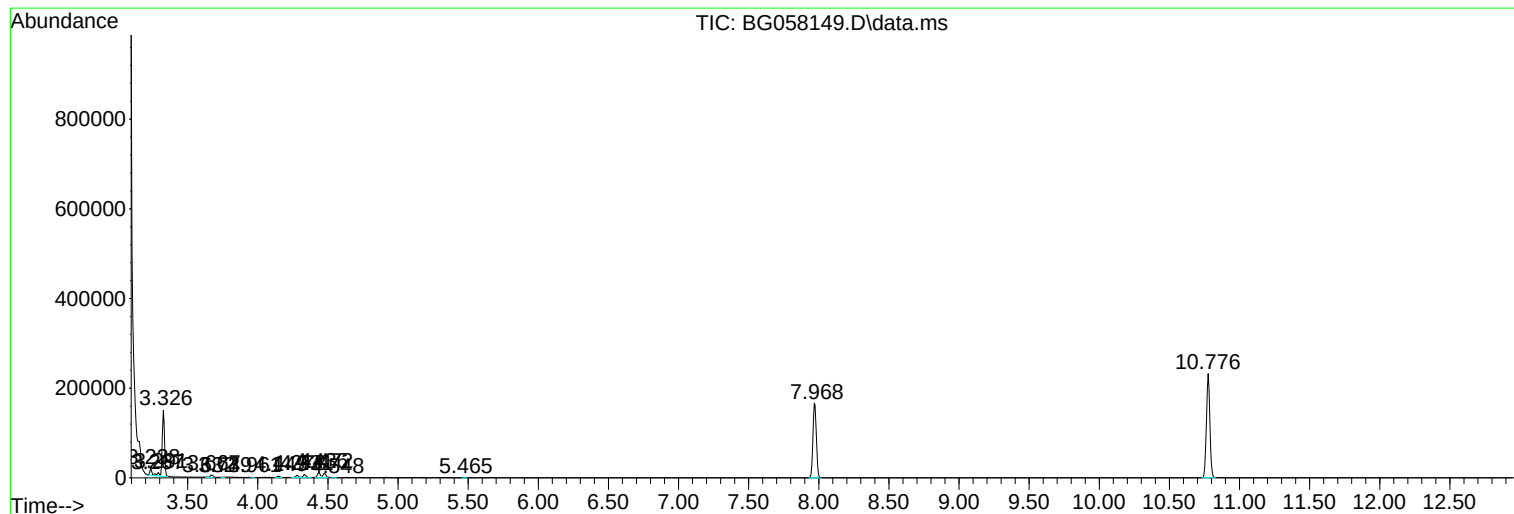
Sum of corrected areas: 6596891

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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



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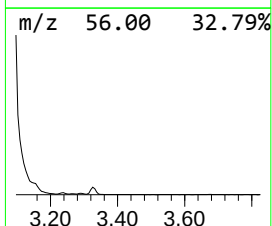
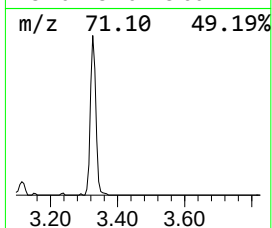
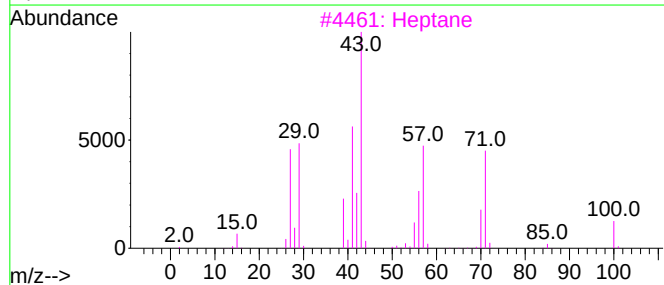
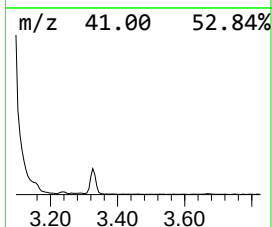
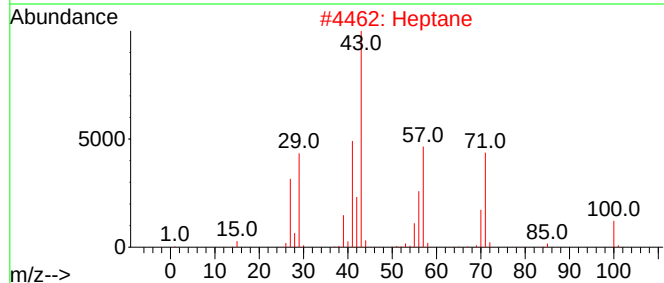
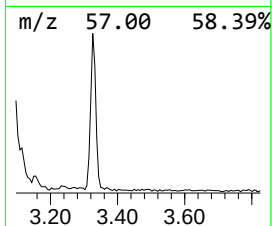
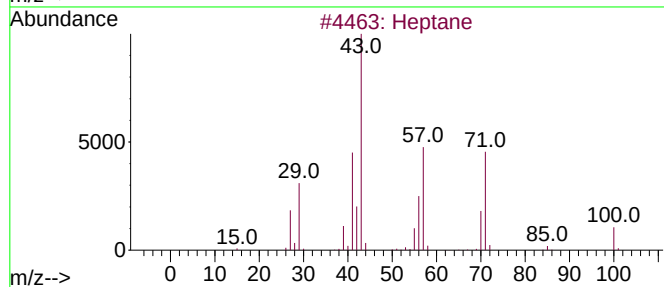
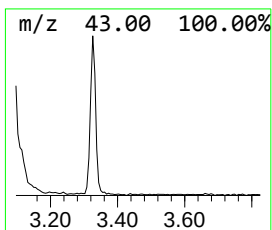
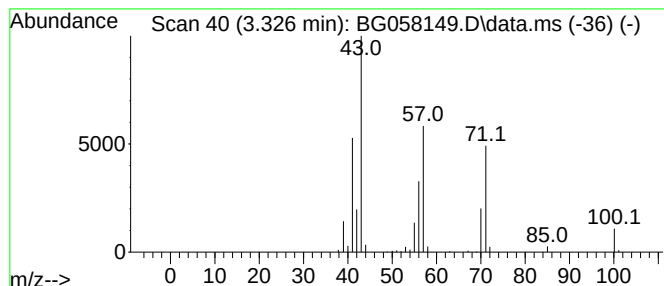
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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.
3.326	11.43 ng/ul	160878	1,4-Dichlorobenzene-d4	7.974

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	90
4			Heptane	100	C7H16	000142-82-5	80
5			Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	64



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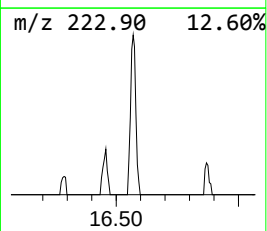
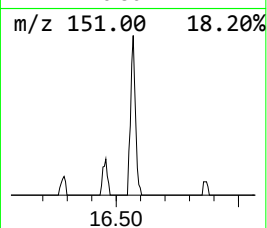
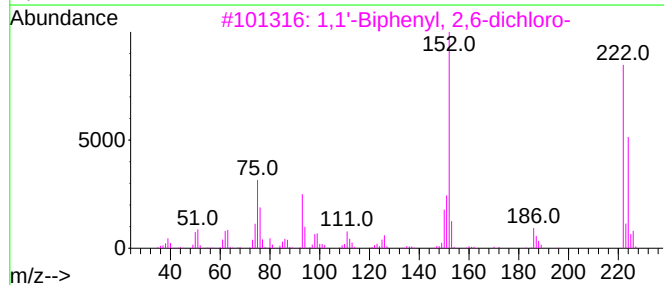
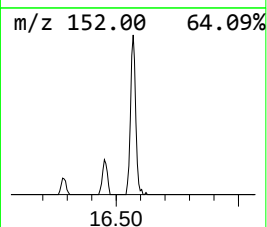
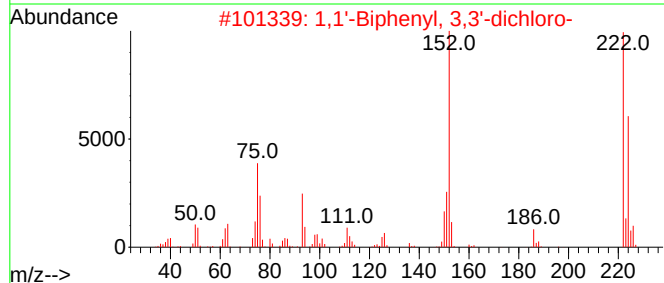
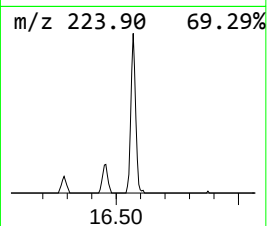
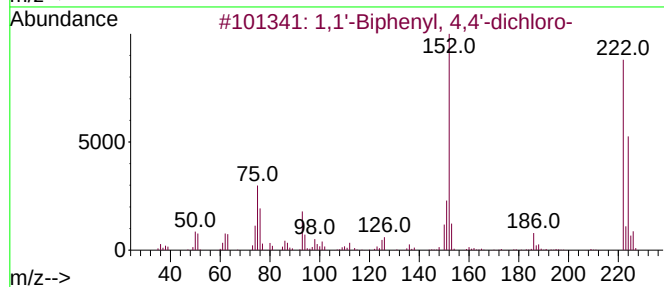
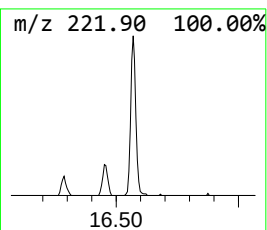
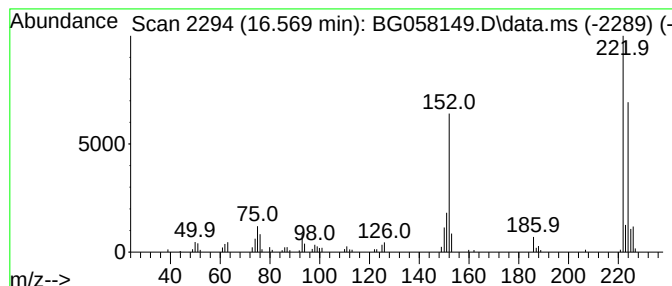
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 1,1'-Biphenyl, 4,4'-dichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.569	2.88 ng/ul	112770	Phenanthrene-d10	17.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 4,4'-dichloro-	222	C12H8Cl2	002050-68-2	98		
2	1,1'-Biphenyl, 3,3'-dichloro-	222	C12H8Cl2	002050-67-1	97		
3	1,1'-Biphenyl, 2,6-dichloro-	222	C12H8Cl2	033146-45-1	97		
4	1,1'-Biphenyl, 4,4'-dichloro-	222	C12H8Cl2	002050-68-2	96		
5	1,1'-Biphenyl, 2,4'-dichloro-	222	C12H8Cl2	034883-43-7	96		



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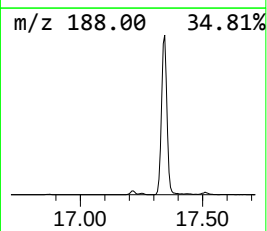
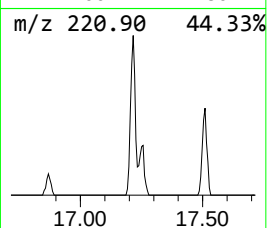
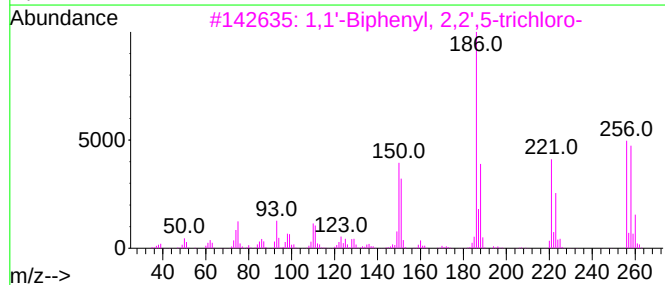
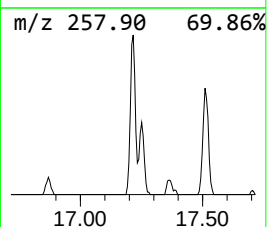
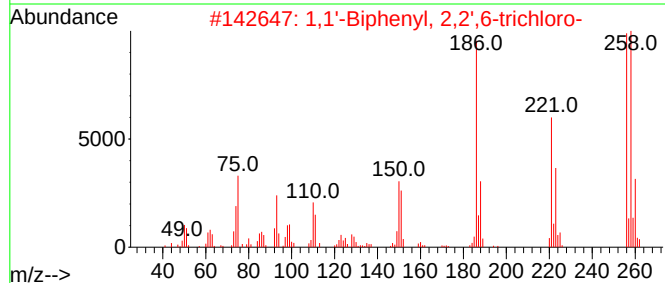
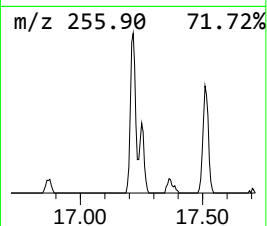
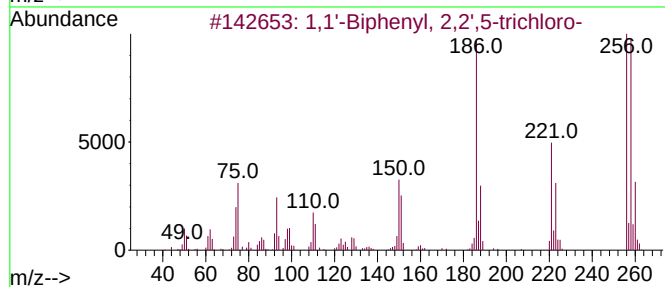
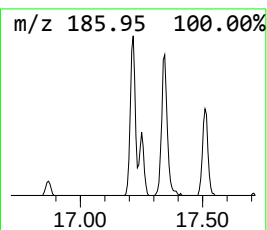
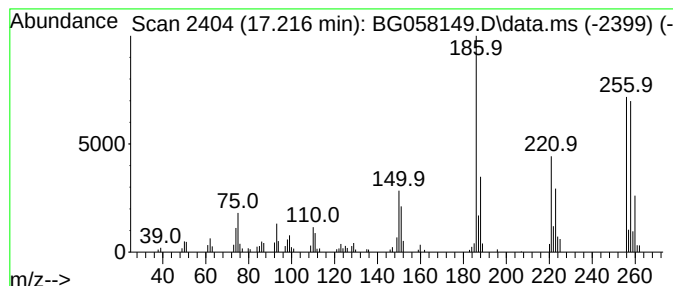
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.216	3.72 ng/ul	145661	Phenanthrene-d10	17.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99		
2	1,1'-Biphenyl, 2,2',6-trichloro-	256	C12H7Cl3	038444-73-4	99		
3	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99		
4	2,2',3-Trichloro-1,1'-biphenyl	256	C12H7Cl3	038444-78-9	99		
5	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071023\
Data File : BG058149.D
Acq On : 10 Jul 2023 19:52
Operator : CG/JU
Sample : AR1660-50PPM
Misc : GCMS CONFIRMATION
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AR1660-50PPM

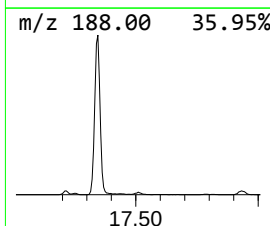
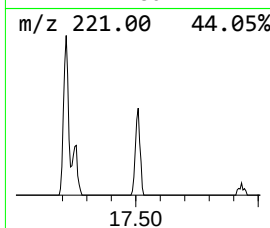
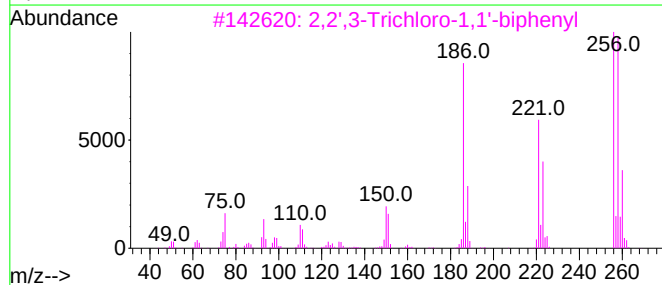
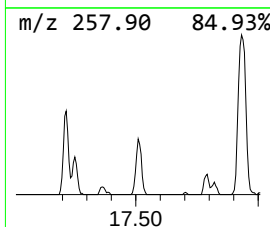
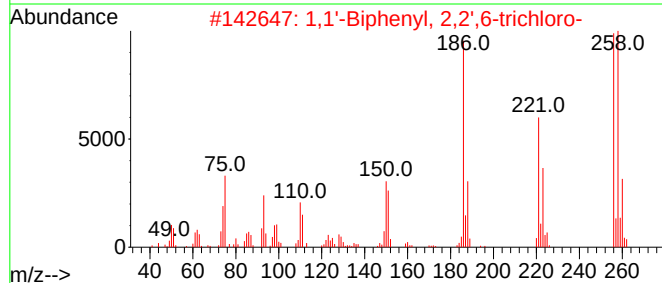
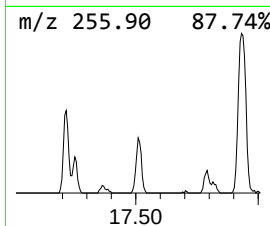
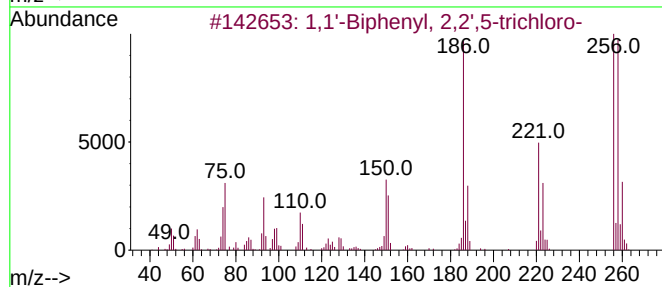
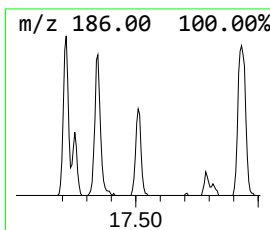
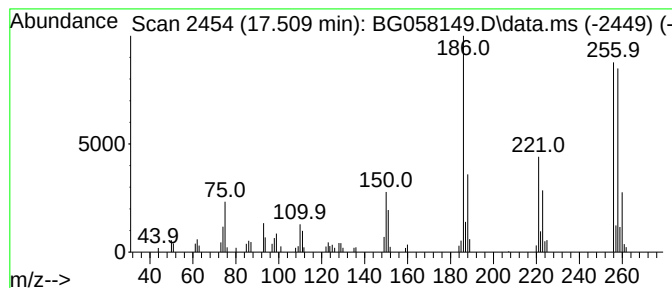
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.509	2.30 ng/ul	89928	Phenanthrene-d10	17.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99		
2	1,1'-Biphenyl, 2,2',6-trichloro-	256	C12H7Cl3	038444-73-4	99		
3	2,2',3-Trichloro-1,1'-biphenyl	256	C12H7Cl3	038444-78-9	99		
4	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99		
5	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	98		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071023\
Data File : BG058149.D
Acq On : 10 Jul 2023 19:52
Operator : CG/JU
Sample : AR1660-50PPM
Misc : GCMS CONFIRMATION
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AR1660-50PPM

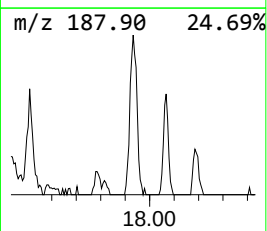
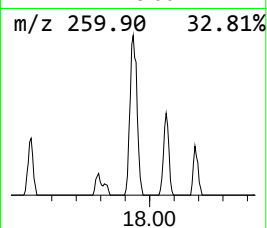
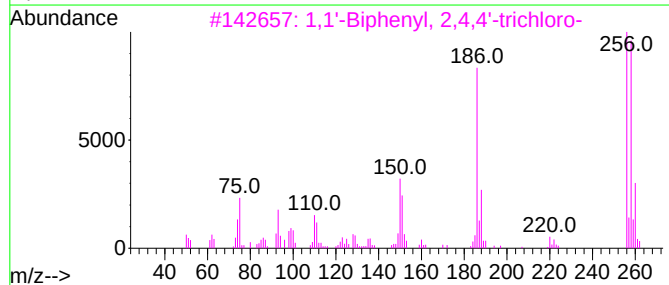
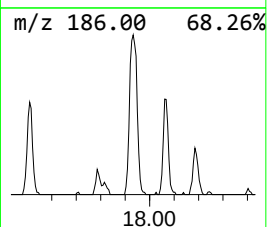
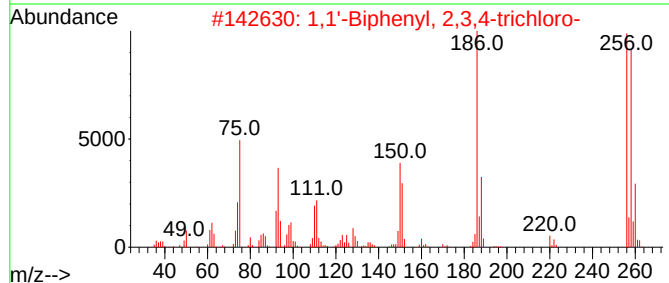
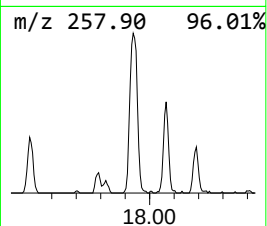
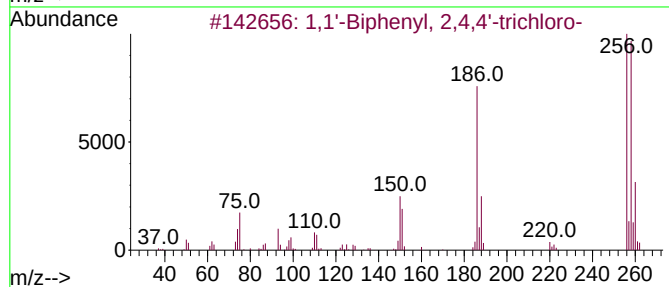
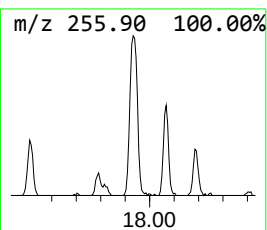
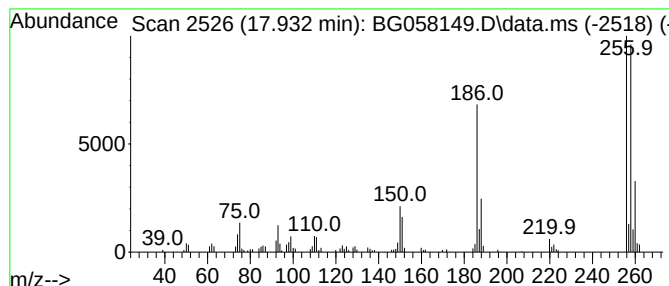
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.932	6.41 ng/ul	251162	Phenanthrene-d10	17.345

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	055702-46-0	99		
3	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99		
4	1,1'-Biphenyl, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	99		
5	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071023\
Data File : BG058149.D
Acq On : 10 Jul 2023 19:52
Operator : CG/JU
Sample : AR1660-50PPM
Misc : GCMS CONFIRMATION
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AR1660-50PPM

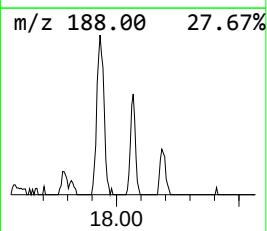
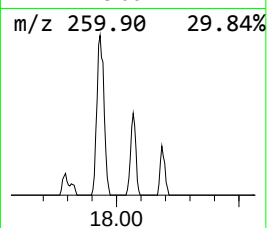
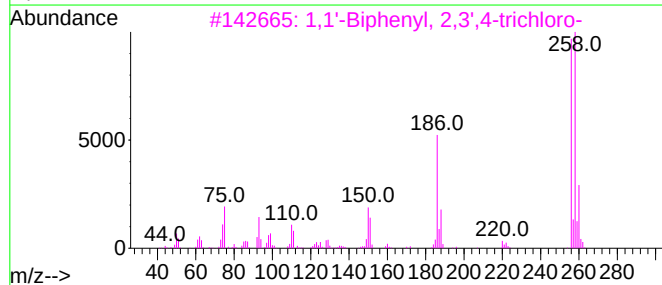
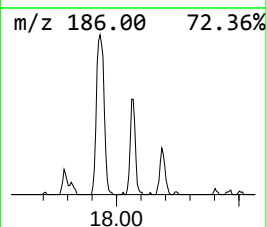
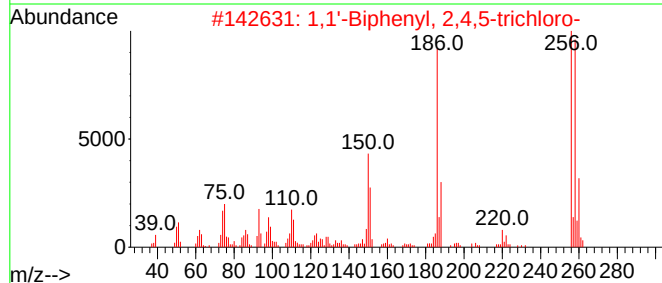
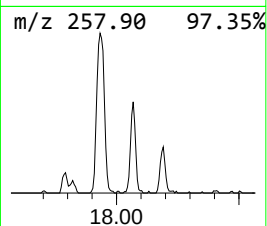
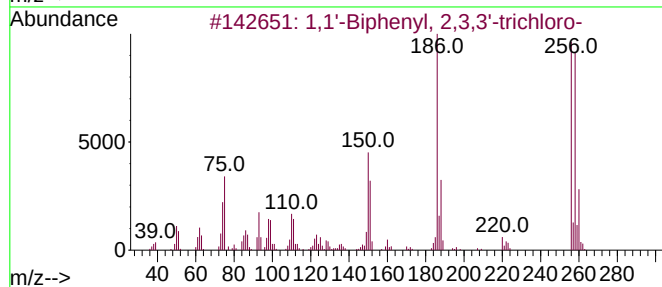
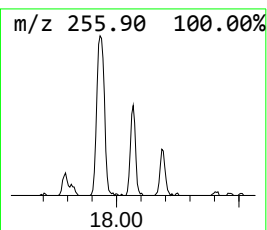
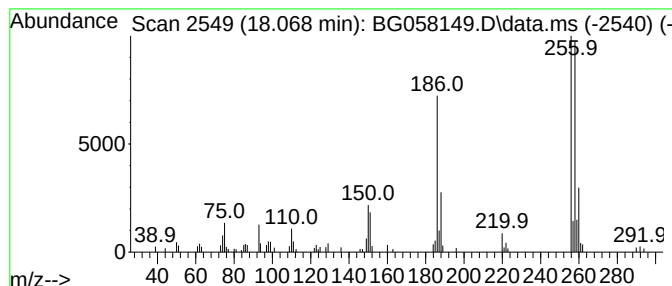
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.068	2.76 ng/ul	108011	Phenanthrene-d10	17.345

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',4-trichloro-	256	C12H7Cl3	055712-37-3	99	
4	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99	
5	1,1'-Biphenyl, 2,3,4'-Trichloro-	256	C12H7Cl3	038444-85-8	99	



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071023\
Data File : BG058149.D
Acq On : 10 Jul 2023 19:52
Operator : CG/JU
Sample : AR1660-50PPM
Misc : GCMS CONFIRMATION
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AR1660-50PPM

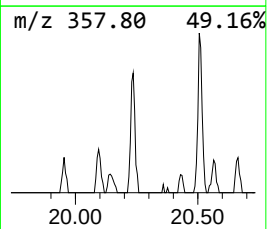
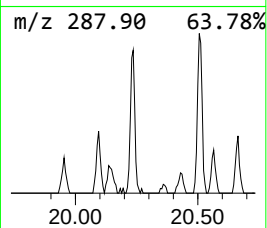
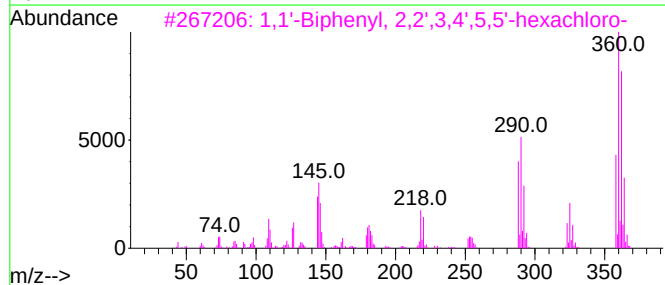
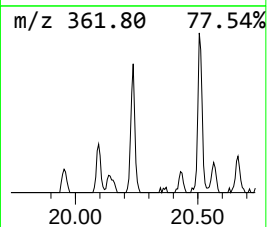
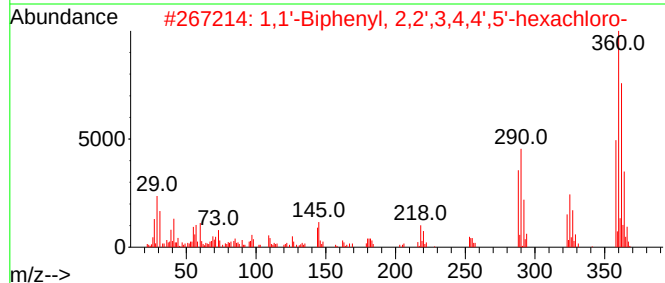
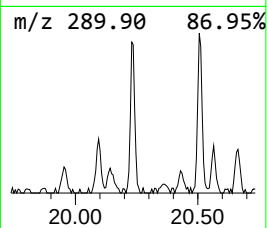
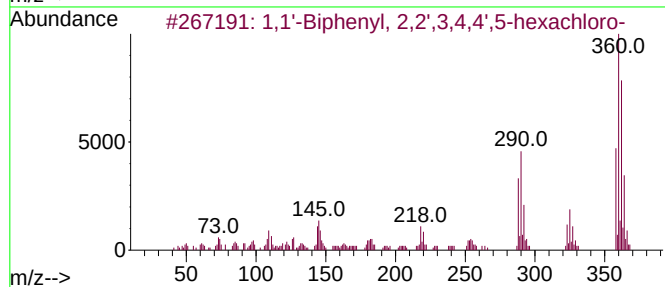
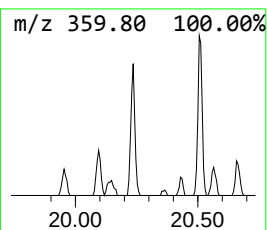
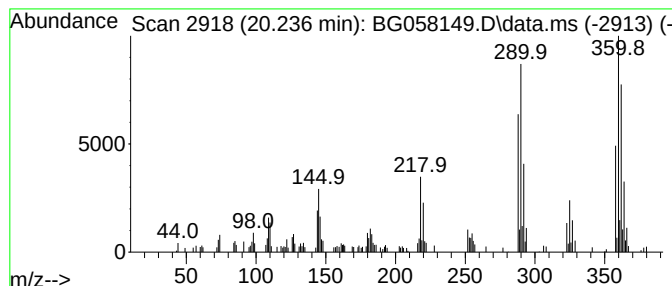
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.236	2.28 ng/ul	109298	Chrysene-d12	21.599

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,4,4',5-hex...	358	C12H4Cl6	035694-06-5	99		
2	1,1'-Biphenyl, 2,2',3,4,4',5'-he...	358	C12H4Cl6	035065-28-2	99		
3	1,1'-Biphenyl, 2,2',3,4',5,5'-he...	358	C12H4Cl6	051908-16-8	99		
4	1,1'-Biphenyl, 2,2',3,4,5,5'-hex...	358	C12H4Cl6	052712-04-6	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\BG071023\
Data File : BG058149.D
Acq On : 10 Jul 2023 19:52
Operator : CG/JU
Sample : AR1660-50PPM
Misc : GCMS CONFIRMATION
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AR1660-50PPM

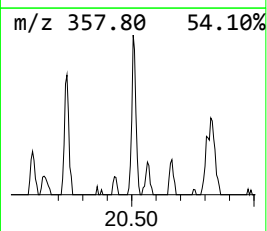
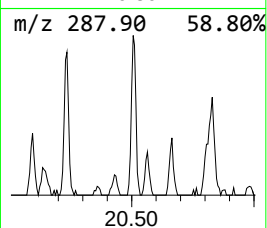
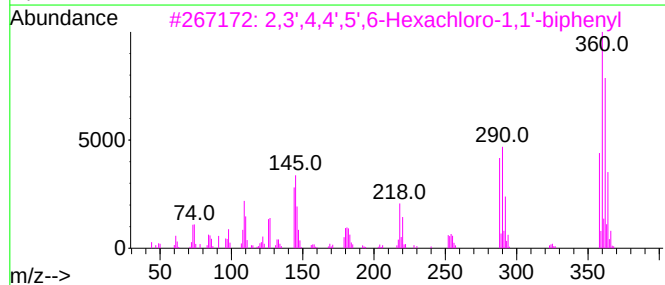
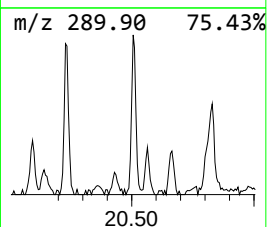
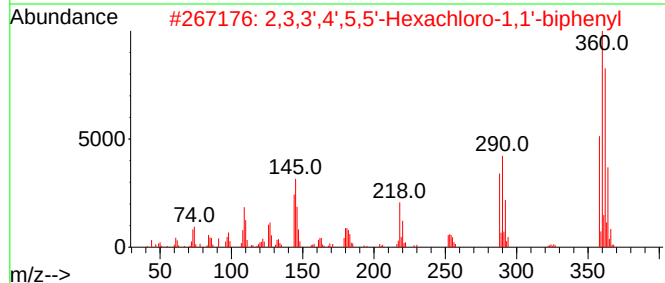
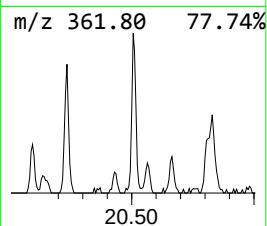
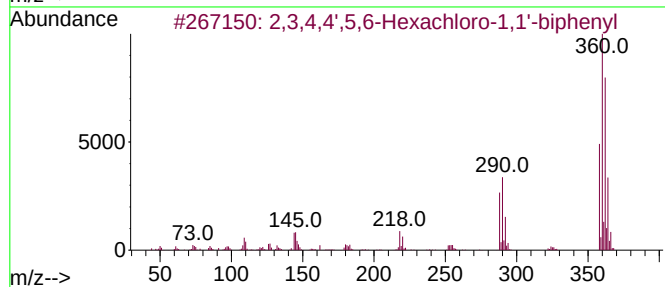
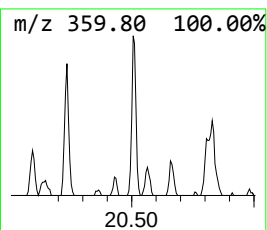
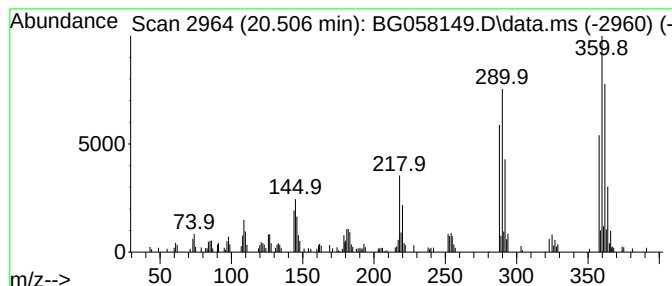
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L

TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.506	2.54 ng/ul	121899	Chrysene-d12	21.599	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,4,4',5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-63-6	99
2	2,3,3',4',5,5'-Hexachloro-1,1'-b...	358	C12H4Cl6	039635-34-2	99
3	2,3',4,4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-65-5	99
4	2,3,3',4,5',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-43-8	99
5	1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071023\
Data File : BG058149.D
Acq On : 10 Jul 2023 19:52
Operator : CG/JU
Sample : AR1660-50PPM
Misc : GCMS CONFIRMATION
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AR1660-50PPM

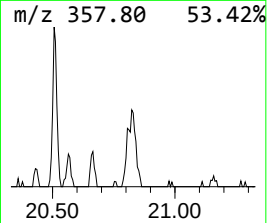
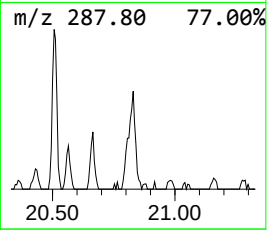
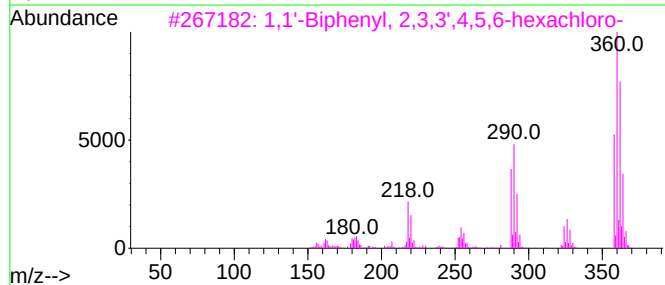
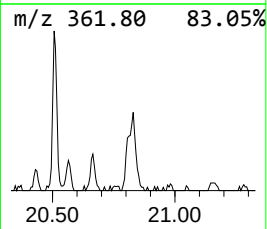
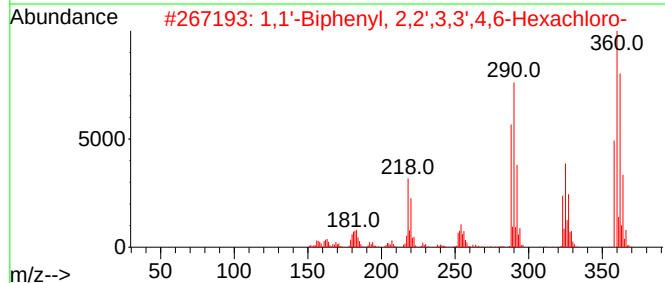
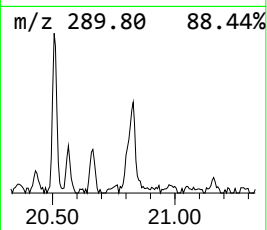
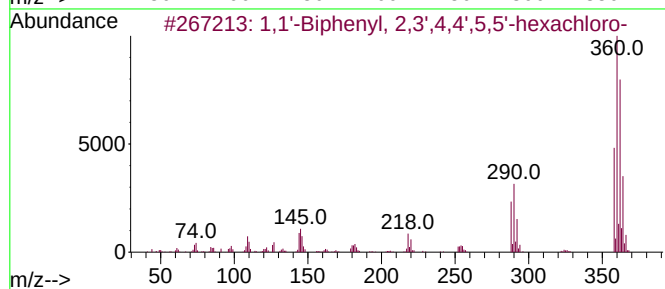
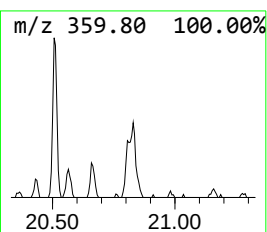
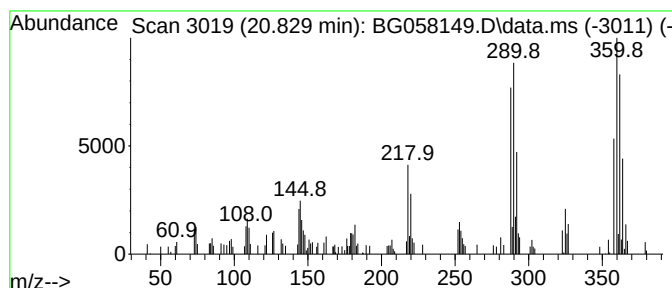
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L

TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.829	2.41 ng/ul	115376	Chrysene-d12	21.599	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3',4,4',5,5'-he...	358	C12H4Cl6	052663-72-6	96
2	1,1'-Biphenyl, 2,2',3,3',4,6-Hex...	358	C12H4Cl6	061798-70-7	95
3	1,1'-Biphenyl, 2,3,3',4,5,6-hexa...	358	C12H4Cl6	041411-62-5	95
4	1,1'-Biphenyl, 2,2',3,3',4,5-hex...	358	C12H4Cl6	055215-18-4	95
5	1,1'-Biphenyl, 2,2',3,3',4,4'-he...	358	C12H4Cl6	038380-07-3	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071023\
Data File : BG058149.D
Acq On : 10 Jul 2023 19:52
Operator : CG/JU
Sample : AR1660-50PPM
Misc : GCMS CONFIRMATION
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AR1660-50PPM

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.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: S...	3.326	11.4	ng/ul	160878	1	7.974	281519	20.0
1,1'-Biphenyl, ...	16.569	2.9	ng/ul	112770	4	17.345	783118	20.0
1,1'-Biphenyl, ...	17.216	3.7	ng/ul	145661	4	17.345	783118	20.0
1,1'-Biphenyl, ...	17.509	2.3	ng/ul	89928	4	17.345	783118	20.0
1,1'-Biphenyl, ...	17.932	6.4	ng/ul	251162	4	17.345	783118	20.0
1,1'-Biphenyl, ...	18.068	2.8	ng/ul	108011	4	17.345	783118	20.0
1,1'-Biphenyl, ...	20.236	2.3	ng/ul	109298	5	21.599	959269	20.0
2,3,4,4',5,6-He...	20.506	2.5	ng/ul	121899	5	21.599	959269	20.0
1,1'-Biphenyl, ...	20.829	2.4	ng/ul	115376	5	21.599	959269	20.0