

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102220\
 Data File : BG047054.D
 Acq On : 23 Oct 2020 6:30
 Operator : CG/JU
 Sample : L4451-03
 Misc : GCMS CONFIRMATION
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AE1

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

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

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.399	6	10	20	rVB	640216	689105	59.02%	8.381%
2	3.716	60	64	65	rBV2	3449	2895	0.25%	0.035%
3	3.746	65	69	74	rVB	23665	32339	2.77%	0.393%
4	3.834	79	84	87	rVB	1682	2618	0.22%	0.032%
5	3.904	95	96	102	rVB	1311	1939	0.17%	0.024%
6	4.128	131	134	138	rVB3	2523	3170	0.27%	0.039%
7	4.227	145	151	156	rBV5	5498	8005	0.69%	0.097%
8	4.345	170	171	174	rVB5	1937	1654	0.14%	0.020%
9	4.410	181	182	186	rVB	2087	2121	0.18%	0.026%
10	4.680	222	228	229	rBV	2052	1969	0.17%	0.024%
11	4.880	260	262	264	rBV2	2411	1818	0.16%	0.022%
12	5.085	292	297	299	rBV2	1915	2419	0.21%	0.029%
13	5.191	313	315	320	rBV2	1773	1936	0.17%	0.024%
14	5.520	369	371	376	rVB2	1915	2286	0.20%	0.028%
15	5.914	433	438	442	rBV2	1384	2258	0.19%	0.027%
16	6.454	528	530	536	rVB3	962	1697	0.15%	0.021%
17	6.771	581	584	587	rVB2	1650	1968	0.17%	0.024%
18	7.053	628	632	635	rBV3	2608	2709	0.23%	0.033%
19	7.682	734	739	745	rBV	7944	13119	1.12%	0.160%
20	8.058	794	803	812	rBV	249491	430515	36.88%	5.236%
21	8.352	851	853	856	rVB2	2142	1908	0.16%	0.023%
22	8.587	890	893	894	rBV	1778	1682	0.14%	0.020%
23	8.652	899	904	908	rVB4	2734	4118	0.35%	0.050%
24	8.716	912	915	919	rBV2	4006	5343	0.46%	0.065%
25	8.822	930	933	934	rBV2	2048	1643	0.14%	0.020%
26	9.286	1007	1012	1019	rBV3	13779	24434	2.09%	0.297%
27	9.439	1035	1038	1043	rVB	1283	1845	0.16%	0.022%
28	9.598	1060	1065	1068	rBV2	1014	1655	0.14%	0.020%
29	9.709	1082	1084	1087	rVB	1692	1877	0.16%	0.023%
30	10.291	1180	1183	1188	rVB2	1589	2328	0.20%	0.028%
31	10.532	1221	1224	1227	rVB2	1498	1687	0.14%	0.021%
32	10.649	1240	1244	1245	rBV2	1369	2108	0.18%	0.026%
33	10.737	1256	1259	1262	rBV2	1893	1905	0.16%	0.023%
34	10.873	1267	1282	1291	rBV	308183	568317	48.68%	6.912%

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35	11.002	1301	1304	1305	rBV2	1694	1947	0.17%	0.024%
36	11.014	1305	1306	1314	rVB2	3482	4624	0.40%	0.056%
37	11.131	1322	1326	1332	rVB4	6563	10122	0.87%	0.123%
38	11.225	1339	1342	1346	rVB2	3016	3119	0.27%	0.038%
39	11.284	1349	1352	1356	rVV3	3212	4396	0.38%	0.053%
40	11.343	1356	1362	1367	rVV6	6590	10920	0.94%	0.133%
41	11.566	1395	1400	1407	rBV3	3342	6553	0.56%	0.080%
42	11.642	1407	1413	1418	rVV5	3017	5733	0.49%	0.070%
43	11.683	1418	1420	1424	rVB	2532	1947	0.17%	0.024%
44	12.124	1492	1495	1498	rVB	1468	1605	0.14%	0.020%
45	12.764	1601	1604	1606	rBV2	1713	1920	0.16%	0.023%
46	13.099	1656	1661	1665	rBV2	1021	1767	0.15%	0.021%
47	13.569	1739	1741	1749	rVB3	1581	3311	0.28%	0.040%
48	13.634	1749	1752	1753	rBV	1329	1601	0.14%	0.019%
49	13.658	1753	1756	1758	rVB2	1469	1633	0.14%	0.020%
50	13.693	1758	1762	1765	rBV	1149	1547	0.13%	0.019%
51	14.692	1925	1932	1939	rBV	489445	759809	65.08%	9.241%
52	15.156	2006	2011	2014	rVB3	1558	2270	0.19%	0.028%
53	16.654	2263	2266	2270	rBV	1005	1660	0.14%	0.020%
54	16.724	2276	2278	2282	rBV2	1379	1968	0.17%	0.024%
55	17.136	2347	2348	2351	rBV2	1701	1964	0.17%	0.024%
56	17.218	2359	2362	2366	rBV3	3523	4179	0.36%	0.051%
57	17.435	2392	2399	2405	rBV	558576	854305	73.18%	10.391%
58	17.770	2454	2456	2458	rBV3	2070	2061	0.18%	0.025%
59	17.853	2469	2470	2473	rVB2	3043	1885	0.16%	0.023%
60	17.882	2473	2475	2479	rBV4	2011	3100	0.27%	0.038%
61	17.988	2490	2493	2495	rVB3	2475	1945	0.17%	0.024%
62	18.041	2495	2502	2506	rBV3	13997	28111	2.41%	0.342%
63	18.140	2517	2519	2521	rVB2	3167	2358	0.20%	0.029%
64	18.240	2534	2536	2537	rVB2	3405	1674	0.14%	0.020%
65	18.276	2540	2542	2545	rVV3	3372	3646	0.31%	0.044%
66	18.434	2567	2569	2572	rBV3	2006	1966	0.17%	0.024%
67	18.499	2575	2580	2584	rVV2	43873	55401	4.75%	0.674%
68	18.558	2587	2590	2594	rVV6	12836	17098	1.46%	0.208%
69	18.599	2595	2597	2602	rVB6	8354	7808	0.67%	0.095%
70	18.781	2625	2628	2633	rBV4	19947	24713	2.12%	0.301%
71	18.822	2633	2635	2639	rVV4	5821	6729	0.58%	0.082%

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Integration Parameters: LSCINT.P

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72	19.045	2671	2673	2676	rBV3	3656	3990	0.34%	0.049%
73	19.216	2700	2702	2706	rBV4	5102	6660	0.57%	0.081%
74	19.257	2707	2709	2714	rVV5	12274	16343	1.40%	0.199%
75	19.310	2714	2718	2719	rVV2	31020	36742	3.15%	0.447%
76	19.339	2719	2723	2728	rVB2	98262	144704	12.39%	1.760%
77	19.480	2743	2747	2753	rBV4	17364	27464	2.35%	0.334%
78	19.545	2754	2758	2764	rVV9	19568	43035	3.69%	0.523%
79	19.621	2766	2771	2776	rVV2	144464	203282	17.41%	2.472%
80	19.680	2777	2781	2786	rVB5	35291	42878	3.67%	0.522%
81	19.880	2811	2815	2820	rVB7	31005	39327	3.37%	0.478%
82	19.956	2824	2828	2832	rVB3	53355	68392	5.86%	0.832%
83	20.062	2840	2846	2852	rVB2	129206	188196	16.12%	2.289%
84	20.179	2863	2866	2870	rBV3	44578	55268	4.73%	0.672%
85	20.320	2886	2890	2895	rVB	146860	196924	16.87%	2.395%
86	20.373	2895	2899	2904	rVB2	94641	114984	9.85%	1.399%
87	20.520	2921	2924	2929	rVB6	33291	39380	3.37%	0.479%
88	20.596	2933	2937	2942	rBV	215829	253291	21.70%	3.081%
89	20.655	2943	2947	2949	rBV4	58049	75488	6.47%	0.918%
90	20.749	2959	2963	2970	rVB6	52403	85762	7.35%	1.043%
91	20.920	2984	2992	2999	rBV2	166967	313142	26.82%	3.809%
92	21.078	3015	3019	3025	rVB3	58930	77796	6.66%	0.946%
93	21.143	3028	3030	3038	rVB6	39112	54480	4.67%	0.663%
94	21.255	3046	3049	3053	rVB5	27405	34565	2.96%	0.420%
95	21.378	3066	3070	3076	rVB5	55172	81333	6.97%	0.989%
96	21.442	3079	3081	3089	rVB9	28683	40940	3.51%	0.498%
97	21.701	3118	3125	3137	rBV	577383	1167480	100.00%	14.200%
98	22.142	3196	3200	3207	rVB7	54033	95701	8.20%	1.164%
99	22.318	3226	3230	3236	rBV9	16901	35349	3.03%	0.430%
100	24.933	3666	3675	3688	rVB2	358782	1004146	86.01%	12.213%

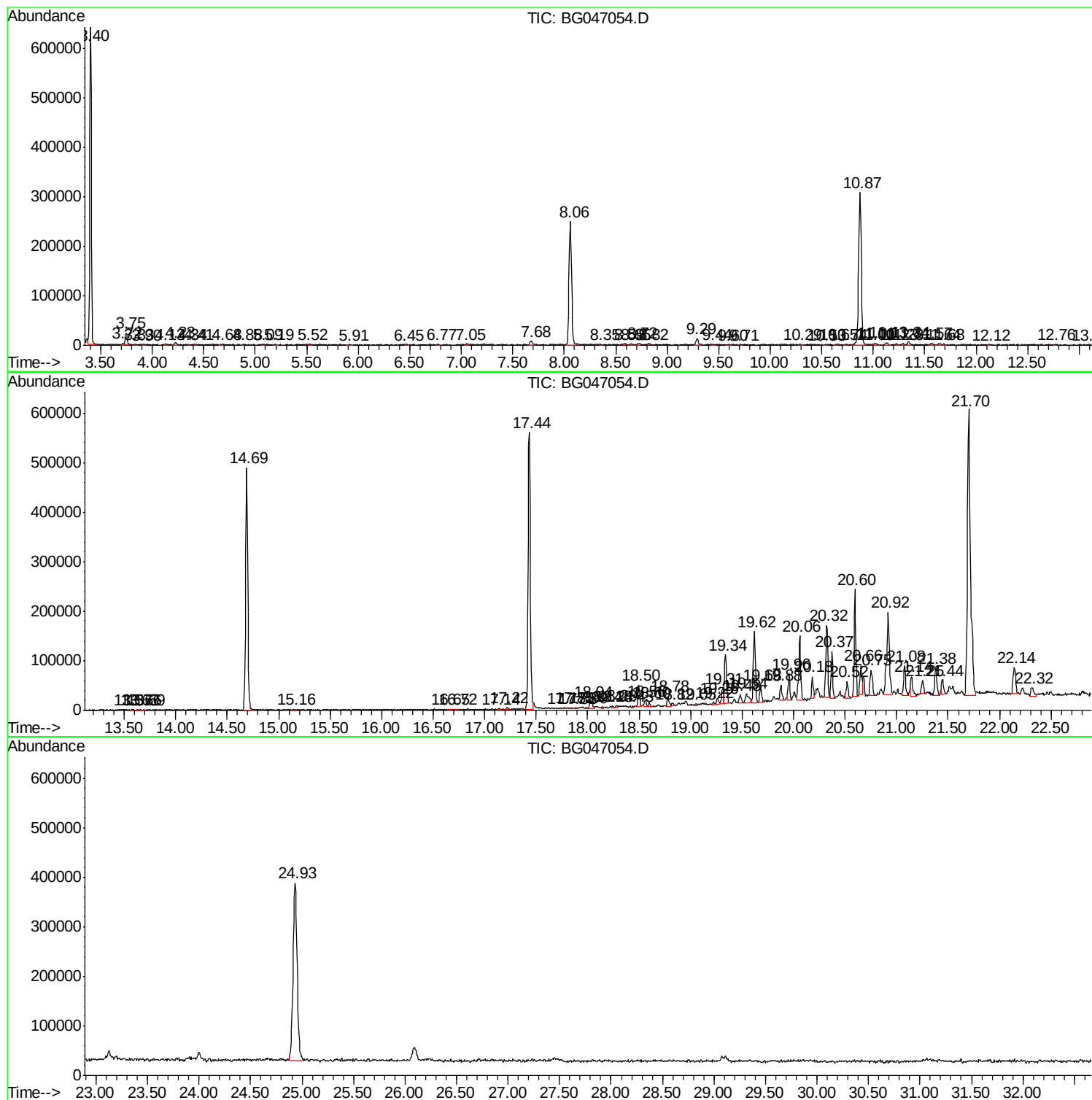
Sum of corrected areas: 8221827

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

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



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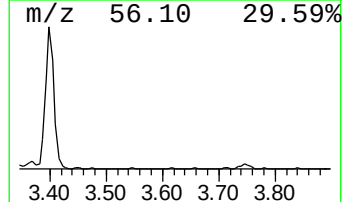
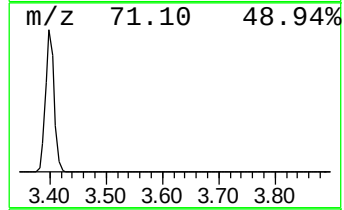
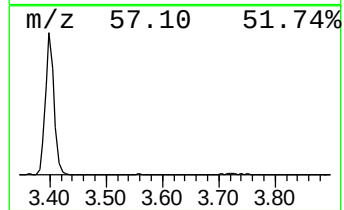
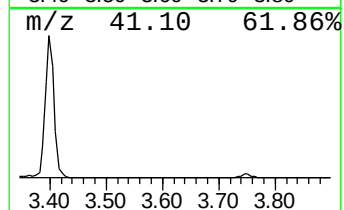
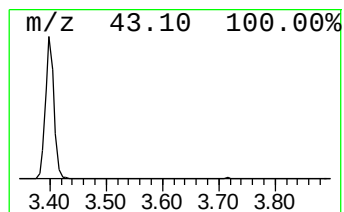
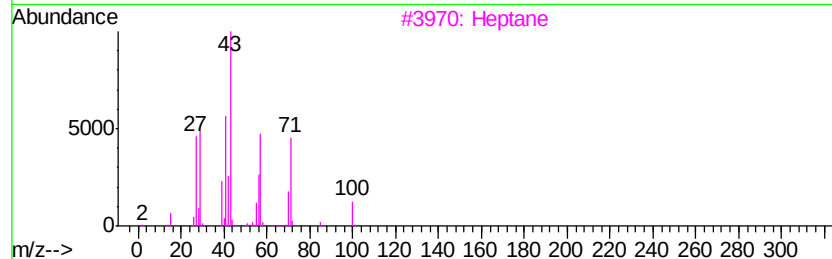
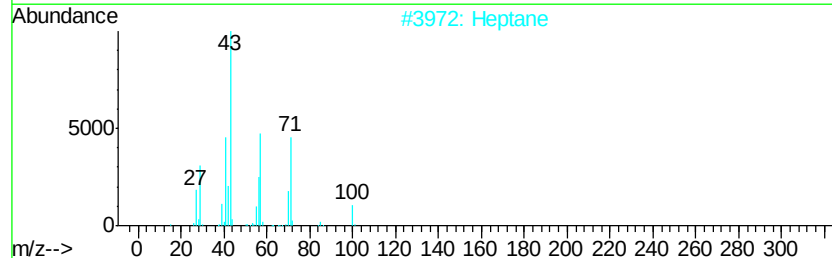
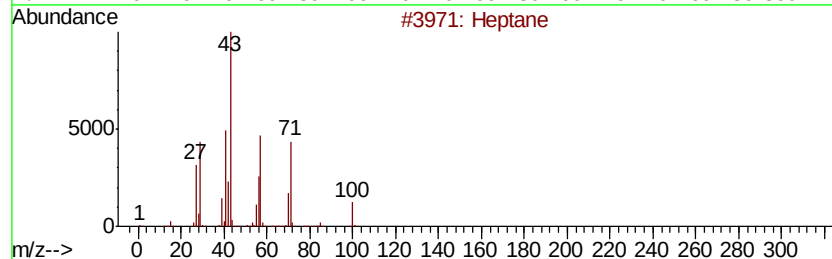
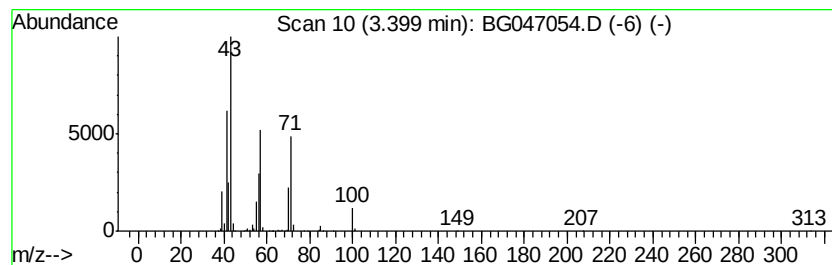
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.40	32.01 ng/ul	689105	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane	100	C7H16	000142-82-5	94
2		Heptane	100	C7H16	000142-82-5	91
3		Heptane	100	C7H16	000142-82-5	91
4		Heptane	100	C7H16	000142-82-5	87
5		Hexane, 3-methyl-	100	C7H16	000589-34-4	43



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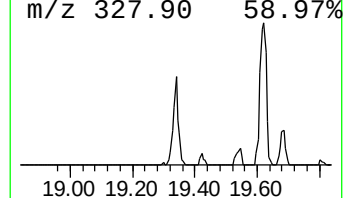
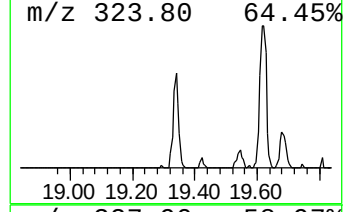
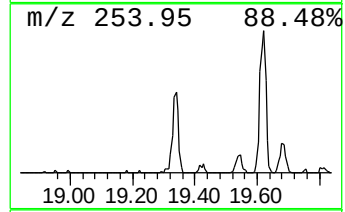
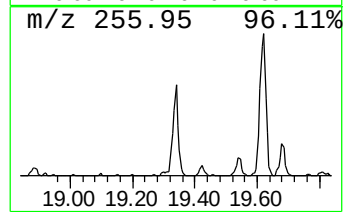
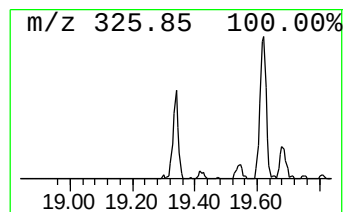
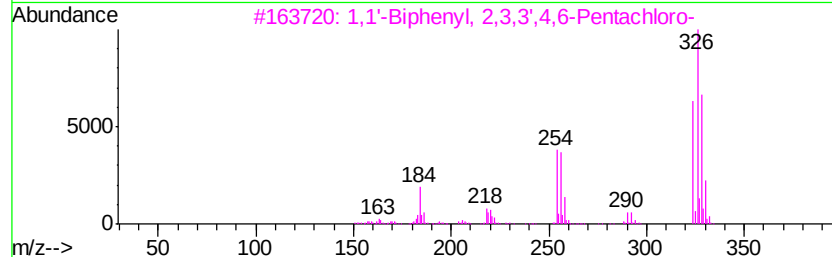
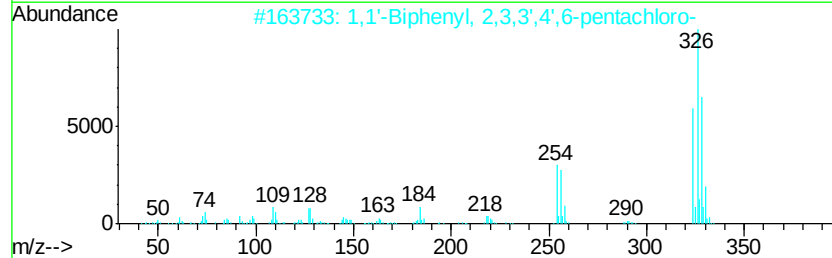
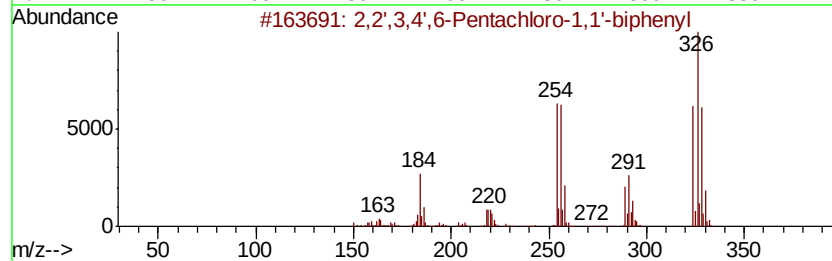
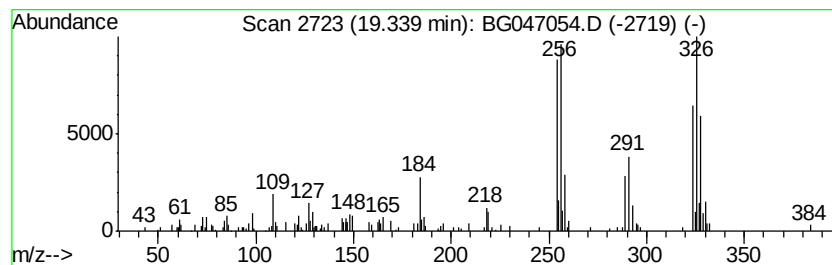
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.34	3.39 ng/ul	144704	Phenanthrene-d10	17.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2',3,4',6-Pentachloro-1,1'-bip...	324	C12H5Cl5	068194-05-8	99
2		1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99
3		1,1'-Biphenyl, 2,3,3',4,6-Pentac...	324	C12H5Cl5	074472-35-8	99
4		1,1'-Biphenyl, 2,2',3',4,6-Penta...	324	C12H5Cl5	060233-25-2	99
5		2,3,3',5,6-Pentachloro-1,1'-biph...	324	C12H5Cl5	074472-36-9	98



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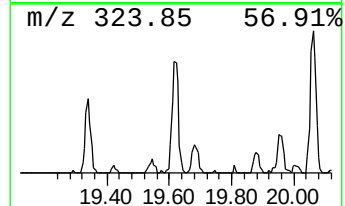
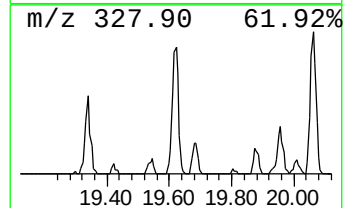
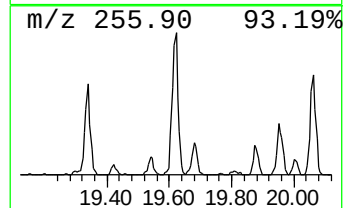
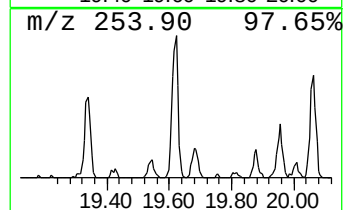
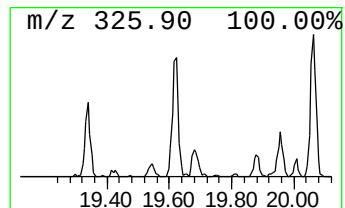
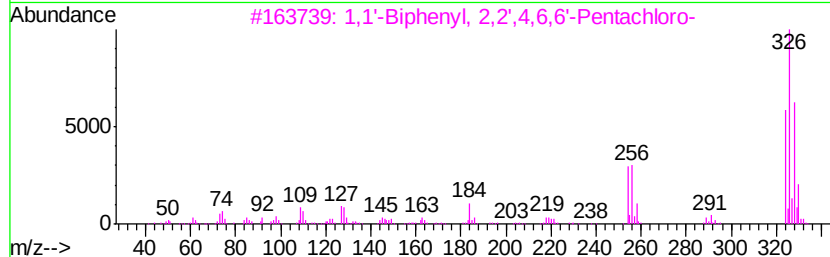
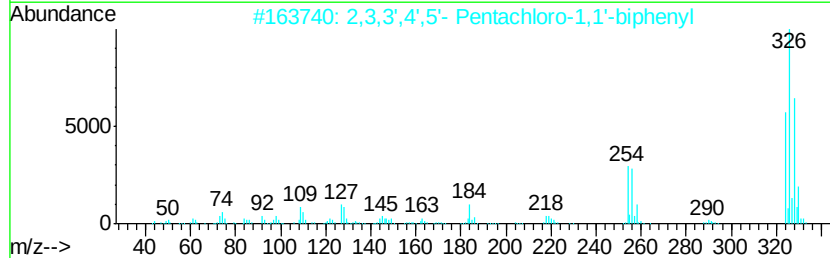
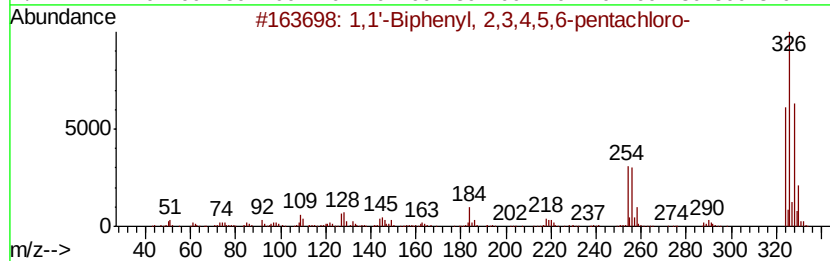
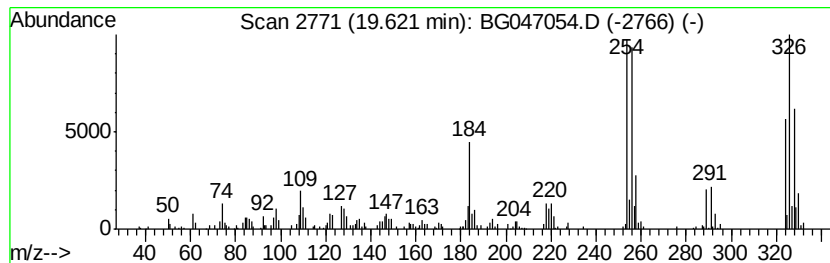
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Peak Number 3 1,1'-Biphenyl, 2,3,4,5,6-pe... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
19.62	3.48 ng/ul	203282	Chrysene-d12	21.70		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,4,5,6-pentachloro-	324	C12H5Cl5	018259-05-7	95	
2	2,3,3',4',5'-Pentachloro-1,1'-biphenyl	324	C12H5Cl5	076842-07-4	95	
3	1,1'-Biphenyl, 2,2',4,6,6'-Pentachloro-	324	C12H5Cl5	056558-16-8	95	
4	(2,3,4,5-Tetrachloro-2,4-cyclopentadienylidene)-	324	C12H5Cl5	029887-33-0	95	
5	2,2',3,4',5'-Pentachloro-1,1'-biphenyl	324	C12H5Cl5	068194-07-0	95	



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102220\
Data File : BG047054.D
Acq On : 23 Oct 2020 6:30
Operator : CG/JU
Sample : L4451-03
Misc : GCMS CONFIRMATION
ALS Vial : 28 Sample Multiplier: 1

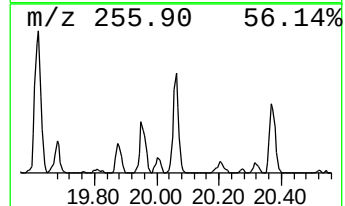
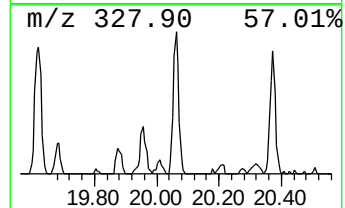
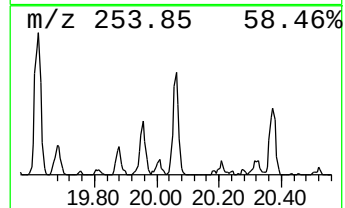
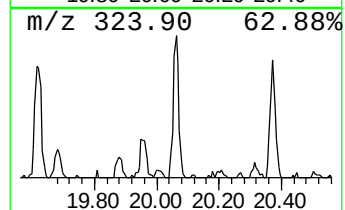
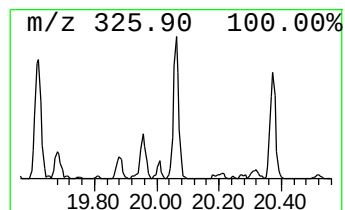
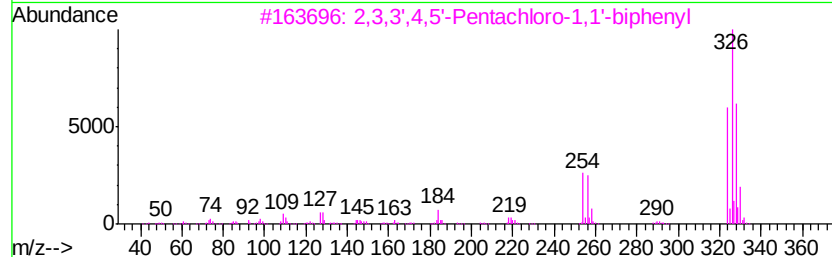
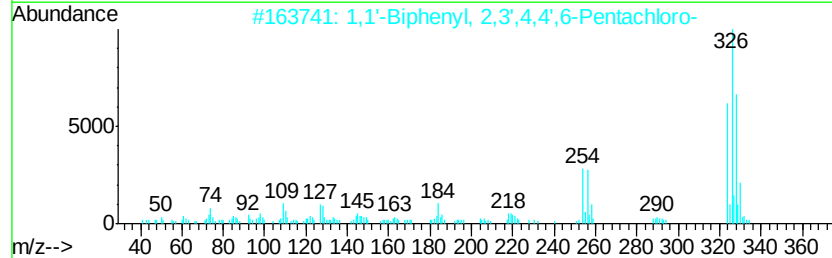
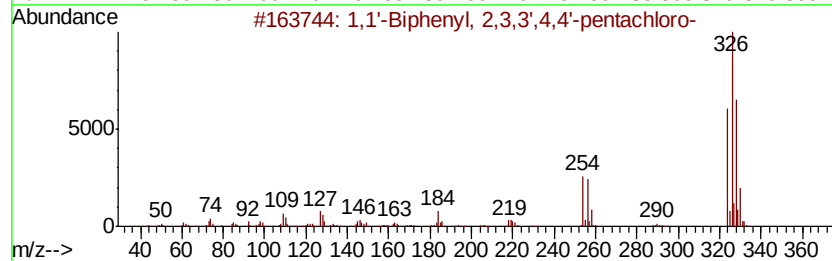
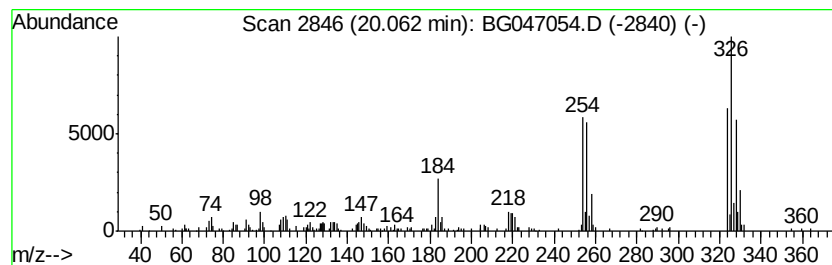
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BNA_G
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.06	3.22 ng/ul	188196	Chrysene-d12	21.70
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1,1'-Biphenyl, 2,3,3',4,4'-penta...	324 C12H5Cl5	032598-14-4	99
2	1,1'-Biphenyl, 2,3',4,4',6-Penta...	324 C12H5Cl5	056558-17-9	99
3	2,3,3',4,5'-Pentachloro-1,1'-bip...	324 C12H5Cl5	070362-41-3	99
4	3,3',4,4',5-Pentachloro-1,1'-bi...	324 C12H5Cl5	057465-28-8	99
5	1,1'-Biphenyl, 2,3,3',4',6-penta...	324 C12H5Cl5	038380-03-9	97



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102220\
Data File : BG047054.D
Acq On : 23 Oct 2020 6:30
Operator : CG/JU
Sample : L4451-03
Misc : GCMS CONFIRMATION
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AE1

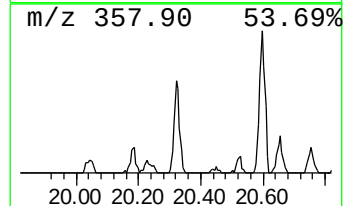
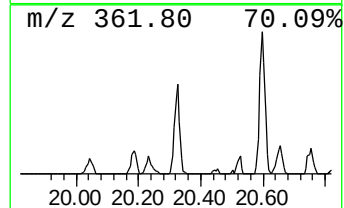
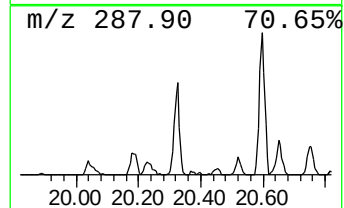
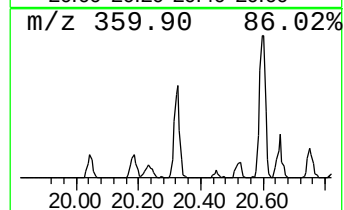
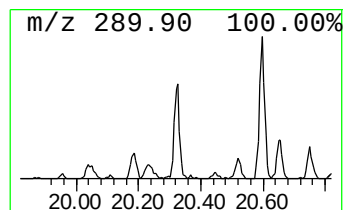
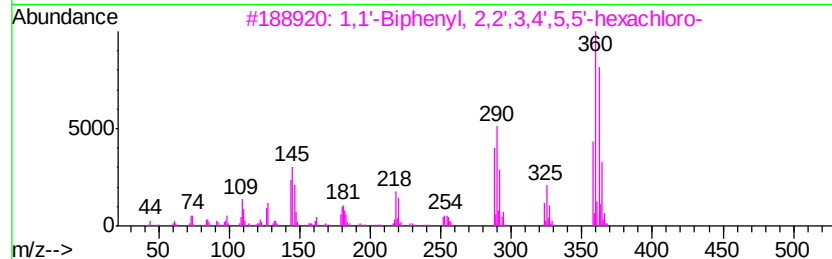
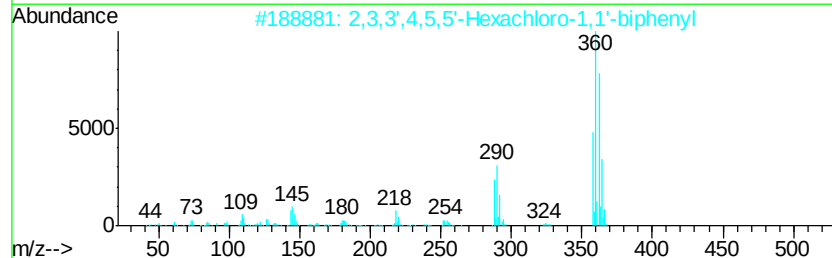
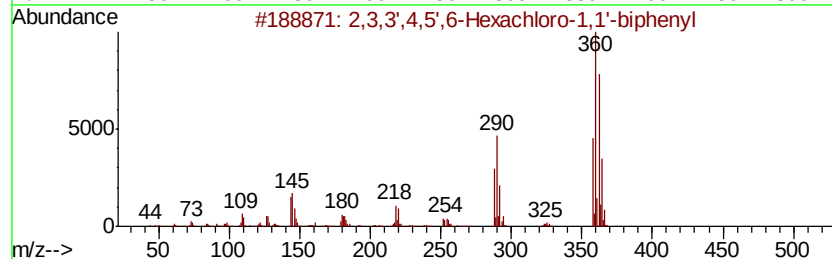
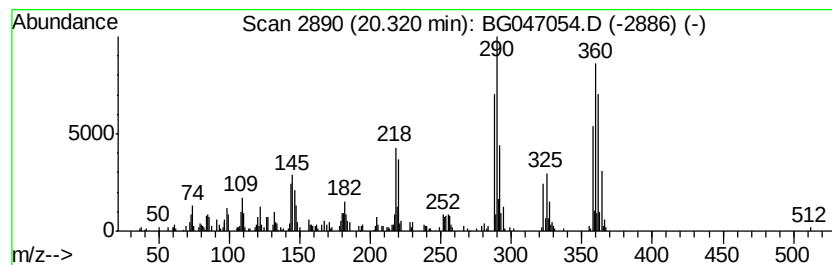
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.32	3.37 ng/ul	196924	Chrysene-d12	21.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,3,3',4,5',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-43-8	99
2		2,3,3',4,5,5'-Hexachloro-1,1'-bi...	358	C12H4Cl6	039635-35-3	99
3		1,1'-Biphenyl, 2,2',3,4',5,5'-he...	358	C12H4Cl6	051908-16-8	99
4		2,3,4,4',5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-63-6	99
5		1,1'-Biphenyl, 2,2',4,4',5',6-He...	358	C12H4Cl6	060145-22-4	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102220\
Data File : BG047054.D
Acq On : 23 Oct 2020 6:30
Operator : CG/JU
Sample : L4451-03
Misc : GCMS CONFIRMATION
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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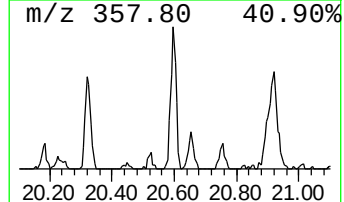
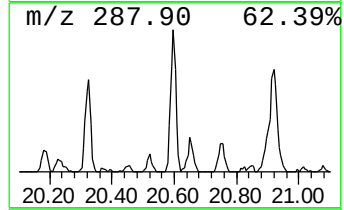
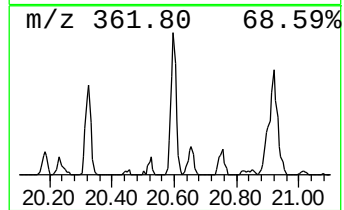
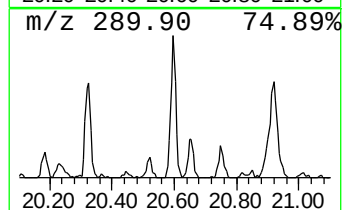
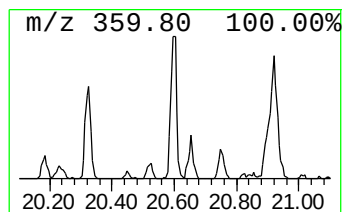
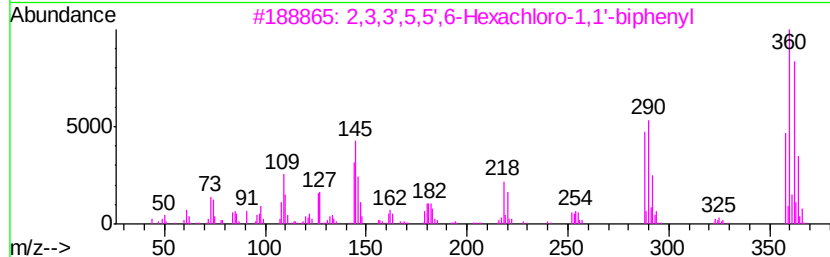
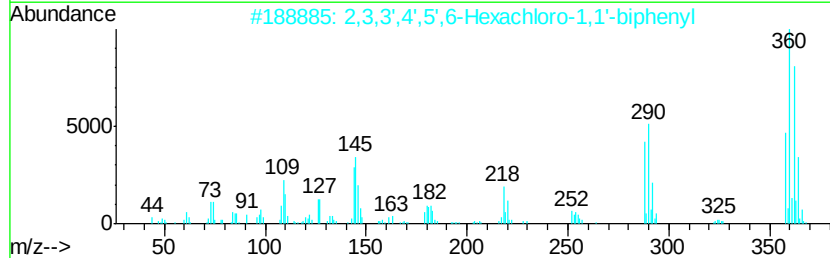
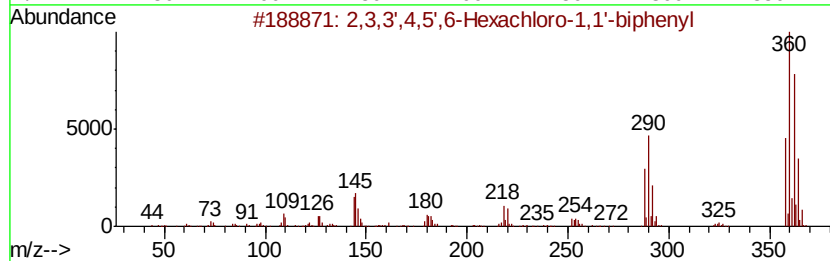
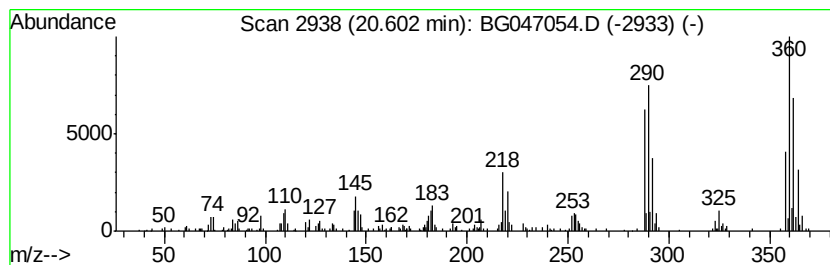
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.60	4.34 ng/ul	253291	Chrysene-d12	21.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,3,3',4,5',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-43-8	99
2		2,3,3',4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	074472-45-0	99
3		2,3,3',5,5',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-46-1	99
4		2,3',4,4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-65-5	99
5		2,2',3,5,6,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-09-2	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102220\
Data File : BG047054.D
Acq On : 23 Oct 2020 6:30
Operator : CG/JU
Sample : L4451-03
Misc : GCMS CONFIRMATION
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AE1

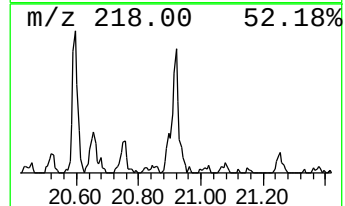
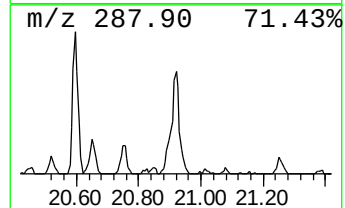
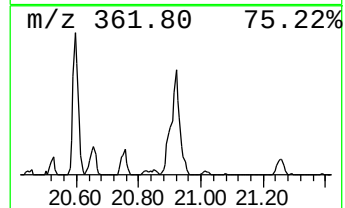
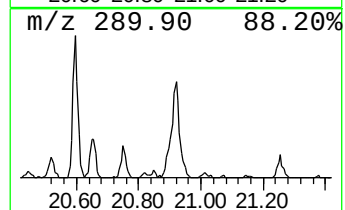
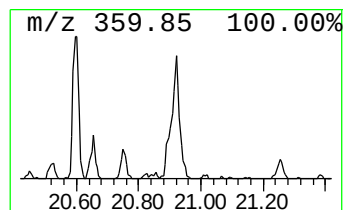
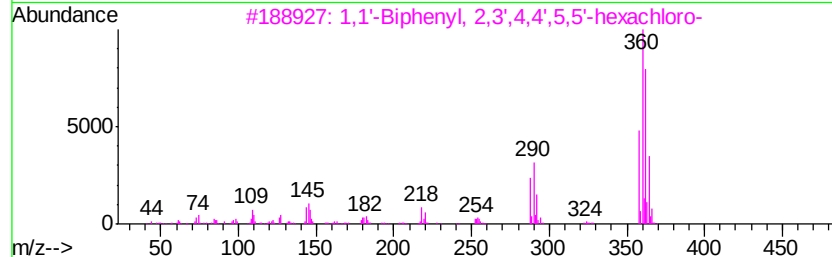
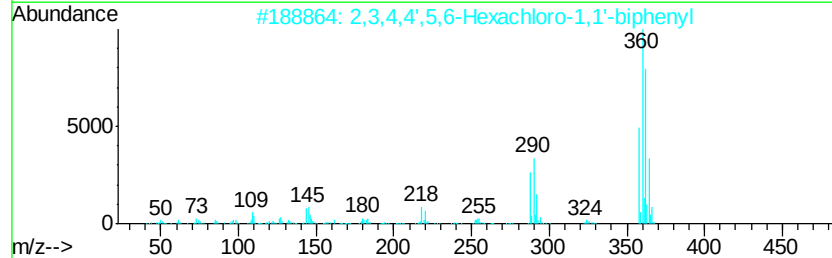
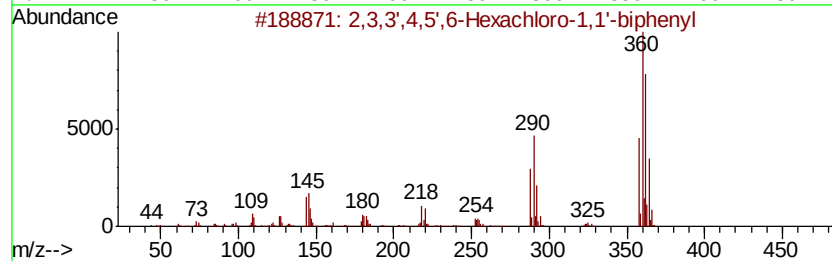
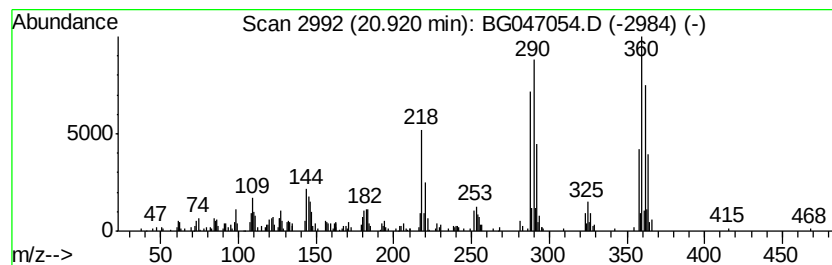
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.92	5.36 ng/ul	313142	Chrysene-d12	21.70

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	2,3,4,4',5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-63-6	99		
3	1,1'-Biphenyl, 2,3',4,4',5,5'-he...	358	C12H4Cl6	052663-72-6	99		
4	2,3,3',4,4',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-42-7	99		
5	2,3,3',4,5,5'-Hexachloro-1,1'-bi...	358	C12H4Cl6	039635-35-3	99		



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG102220\
Data File : BG047054.D
Acq On : 23 Oct 2020 6:30
Operator : CG/JU
Sample : L4451-03
Misc : GCMS CONFIRMATION
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AE1

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Str...	3.40	32.0	ng/ul	689105	1	8.06	430515	20.0
2,2',3,4',6-Penta...	19.34	3.4	ng/ul	144704	4	17.44	854305	20.0
1,1'-Biphenyl, 2,...	19.62	3.5	ng/ul	203282	5	21.70	1167480	20.0
1,1'-Biphenyl, 2,...	20.06	3.2	ng/ul	188196	5	21.70	1167480	20.0
2,3,3',4,5',6-Hex...	20.32	3.4	ng/ul	196924	5	21.70	1167480	20.0
2,3,3',4',5',6-He...	20.60	4.3	ng/ul	253291	5	21.70	1167480	20.0
1,1'-Biphenyl, 2,...	20.92	5.4	ng/ul	313142	5	21.70	1167480	20.0