

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
 Data File : BG047160.D
 Acq On : 30 Oct 2020 18:55
 Operator : CG/JU
 Sample : L4505-13
 Misc : GCMS CONFIRMATION
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0BH2

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	3.390	5	9	17	rVB	489957	531813	48.99%	6.907%
2	3.484	24	25	28	rVB	1656	1332	0.12%	0.017%
3	3.537	28	34	35	rBV2	1155	1602	0.15%	0.021%
4	3.578	38	41	44	rBV3	1696	2540	0.23%	0.033%
5	3.631	47	50	51	rVB	1775	1648	0.15%	0.021%
6	3.643	51	52	55	rBV	1688	1818	0.17%	0.024%
7	3.708	59	63	64	rBV2	1942	2077	0.19%	0.027%
8	3.737	64	68	74	rVB5	13993	16742	1.54%	0.217%
9	3.849	85	87	89	rBV	1562	1435	0.13%	0.019%
10	4.031	116	118	122	rBV2	1205	1661	0.15%	0.022%
11	4.219	144	150	154	rVB2	3580	4688	0.43%	0.061%
12	4.366	172	175	179	rBV2	1172	1845	0.17%	0.024%
13	4.665	222	226	228	rBV2	1510	1894	0.17%	0.025%
14	5.065	288	294	295	rBV2	1172	2225	0.20%	0.029%
15	5.200	315	317	320	rBV	1352	1398	0.13%	0.018%
16	5.441	356	358	361	rBV	1702	1523	0.14%	0.020%
17	6.263	491	498	499	rBV	1709	2131	0.20%	0.028%
18	7.444	697	699	702	rBV4	1786	1540	0.14%	0.020%
19	7.480	704	705	708	rVB3	1821	1340	0.12%	0.017%
20	7.527	708	713	715	rBV3	2741	3949	0.36%	0.051%
21	7.609	725	727	731	rVB2	2345	3090	0.28%	0.040%
22	7.709	740	744	749	rBV2	1194	1904	0.18%	0.025%
23	8.055	795	803	813	rBV	197773	338016	31.13%	4.390%
24	8.173	822	823	824	rBV	2379	1579	0.15%	0.021%
25	8.191	824	826	829	rVB2	2575	3216	0.30%	0.042%
26	8.802	927	930	934	rVB	1130	1377	0.13%	0.018%
27	10.006	1131	1135	1137	rBV	1013	1388	0.13%	0.018%
28	10.864	1273	1281	1292	rBV	249604	446112	41.09%	5.794%
29	11.493	1386	1388	1391	rVB2	1187	1319	0.12%	0.017%
30	12.421	1545	1546	1550	rBV2	1495	1548	0.14%	0.020%
31	13.167	1670	1673	1676	rVB2	1071	1486	0.14%	0.019%
32	13.772	1771	1776	1777	rBV2	1111	1649	0.15%	0.021%
33	14.336	1870	1872	1874	rBV	1784	1748	0.16%	0.023%
34	14.454	1888	1892	1895	rVB2	1092	1534	0.14%	0.020%

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

35	14.683	1925	1931	1941	rBV	445022	693159	63.85%	9.003%
36	14.795	1947	1950	1954	rVB2	1553	2088	0.19%	0.027%
37	14.830	1954	1956	1959	rBV	1319	1501	0.14%	0.019%
38	15.012	1985	1987	1990	rVB	1539	1391	0.13%	0.018%
39	15.053	1990	1994	1996	rBV	1794	2057	0.19%	0.027%
40	16.152	2178	2181	2183	rBV	1310	1606	0.15%	0.021%
41	16.240	2194	2196	2199	rVB	1489	1349	0.12%	0.018%
42	16.387	2218	2221	2223	rBV2	2479	2731	0.25%	0.035%
43	16.422	2225	2227	2231	rVB3	1221	1416	0.13%	0.018%
44	16.493	2236	2239	2241	rVV2	1515	1887	0.17%	0.025%
45	16.528	2241	2245	2248	rVB3	1903	1988	0.18%	0.026%
46	16.698	2270	2274	2277	rVB2	1179	1774	0.16%	0.023%
47	16.898	2304	2308	2311	rBV3	1716	3037	0.28%	0.039%
48	17.186	2353	2357	2358	rBV2	1922	1845	0.17%	0.024%
49	17.374	2384	2389	2390	rBV	1476	1721	0.16%	0.022%
50	17.427	2392	2398	2409	rVV	567294	881370	81.18%	11.447%
51	17.591	2425	2426	2431	rVB4	1681	1967	0.18%	0.026%
52	17.785	2458	2459	2463	rBV2	1922	2396	0.22%	0.031%
53	17.926	2478	2483	2486	rBV3	2527	3471	0.32%	0.045%
54	18.008	2495	2497	2498	rBV	2921	1881	0.17%	0.024%
55	18.032	2498	2501	2507	rVB3	8200	11319	1.04%	0.147%
56	18.091	2507	2511	2513	rBV3	2302	3680	0.34%	0.048%
57	18.273	2540	2542	2545	rBV3	2642	3489	0.32%	0.045%
58	18.373	2557	2559	2562	rVB3	1992	1936	0.18%	0.025%
59	18.414	2565	2566	2569	rBV2	1679	1636	0.15%	0.021%
60	18.490	2575	2579	2586	rVB	51389	72003	6.63%	0.935%
61	18.555	2586	2590	2594	rVB3	12604	17967	1.65%	0.233%
62	18.602	2594	2598	2602	rVB4	7724	11619	1.07%	0.151%
63	18.702	2613	2615	2618	rVB3	2086	1843	0.17%	0.024%
64	18.725	2618	2619	2622	rBV3	1936	1844	0.17%	0.024%
65	18.778	2623	2628	2632	rBV3	26886	33793	3.11%	0.439%
66	18.913	2648	2651	2654	rBV4	3411	3742	0.34%	0.049%
67	18.949	2654	2657	2663	rVB4	12007	15882	1.46%	0.206%
68	19.019	2667	2669	2670	rBV	2284	2296	0.21%	0.030%
69	19.095	2680	2682	2684	rBV3	1823	1914	0.18%	0.025%
70	19.254	2706	2709	2712	rBV	12957	17257	1.59%	0.224%
71	19.313	2714	2719	2720	rVV2	41284	61574	5.67%	0.800%

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72	19.336	2720	2723	2730	rVB3	88565	126848	11.68%	1.647%
73	19.419	2734	2737	2742	rVB4	13052	15996	1.47%	0.208%
74	19.536	2753	2757	2762	rBV6	21476	34825	3.21%	0.452%
75	19.612	2765	2770	2776	rVB2	165923	207751	19.14%	2.698%
76	19.677	2776	2781	2785	rBV2	59671	74760	6.89%	0.971%
77	19.724	2785	2789	2793	rBV	102461	111948	10.31%	1.454%
78	19.801	2799	2802	2806	rBV4	21483	25241	2.32%	0.328%
79	19.842	2806	2809	2811	rVV3	15843	16639	1.53%	0.216%
80	19.877	2811	2815	2821	rVB3	39506	51504	4.74%	0.669%
81	19.953	2824	2828	2832	rVV4	70305	91921	8.47%	1.194%
82	20.000	2833	2836	2840	rVV4	27530	30790	2.84%	0.400%
83	20.059	2840	2846	2852	rVB	157097	209639	19.31%	2.723%
84	20.182	2863	2867	2868	rBV3	11339	13195	1.22%	0.171%
85	20.318	2885	2890	2895	rBV4	73259	116993	10.78%	1.519%
86	20.370	2895	2899	2907	rVB	134294	176329	16.24%	2.290%
87	20.594	2933	2937	2942	rBV2	80551	100443	9.25%	1.305%
88	20.652	2943	2947	2948	rVV2	37940	47107	4.34%	0.612%
89	20.676	2948	2951	2956	rVV3	61581	81384	7.50%	1.057%
90	20.758	2960	2965	2970	rVV2	94193	121816	11.22%	1.582%
91	20.829	2973	2977	2981	rVV	48584	66829	6.16%	0.868%
92	20.917	2985	2992	2999	rVB3	96229	171185	15.77%	2.223%
93	21.252	3045	3049	3054	rVB7	34180	47800	4.40%	0.621%
94	21.334	3060	3063	3067	rVB3	31032	39732	3.66%	0.516%
95	21.698	3118	3125	3135	rBV	613105	1047726	96.51%	13.608%
96	21.880	3152	3156	3164	rVB2	46356	68141	6.28%	0.885%
97	22.491	3256	3260	3267	rVB2	41669	68374	6.30%	0.888%
98	23.185	3375	3378	3386	rVB2	45284	76502	7.05%	0.994%
99	23.996	3512	3516	3527	rVB4	53702	106796	9.84%	1.387%
100	24.930	3666	3675	3689	rVB3	369714	1085660	100.00%	14.100%

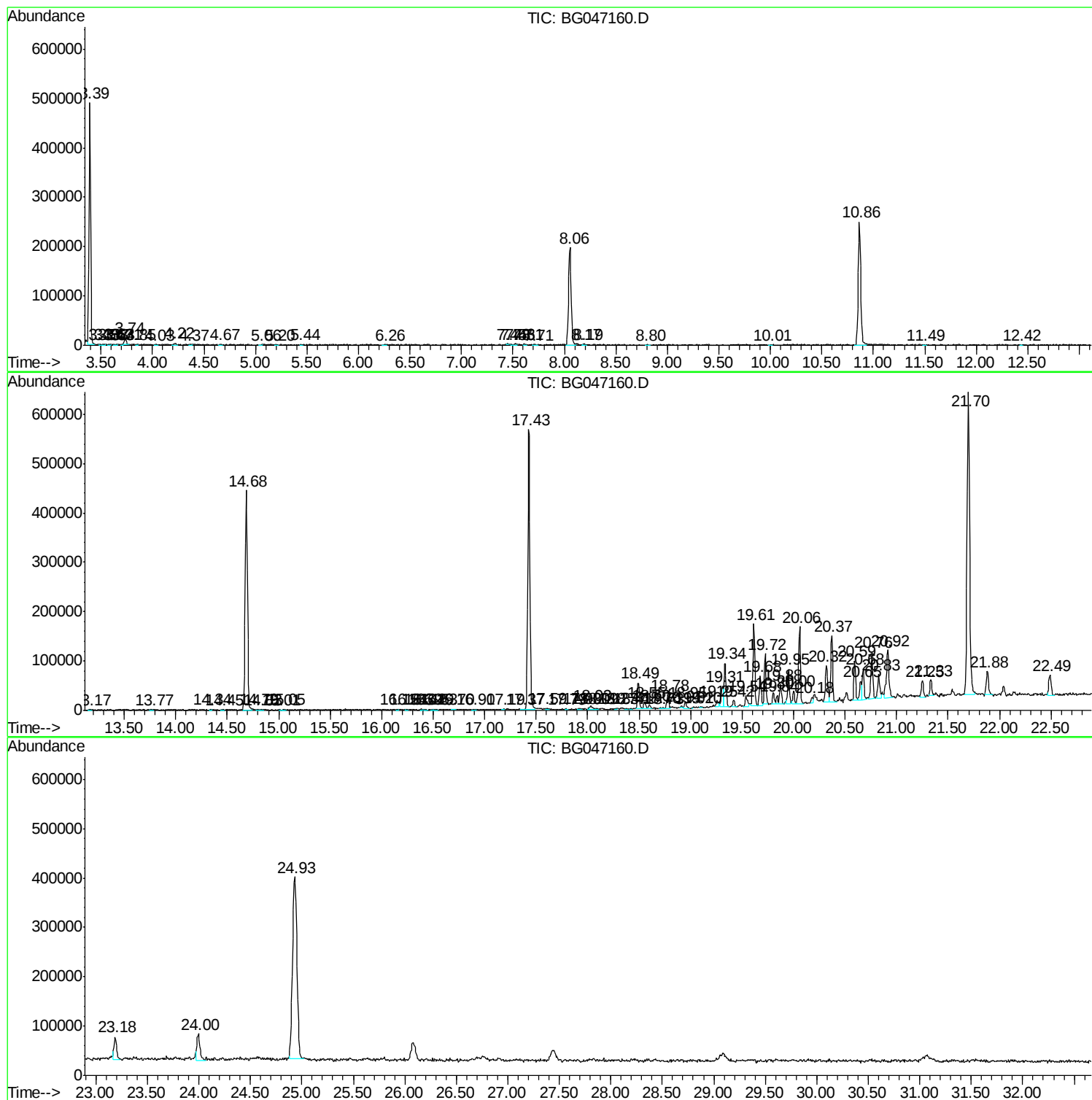
Sum of corrected areas: 7699470

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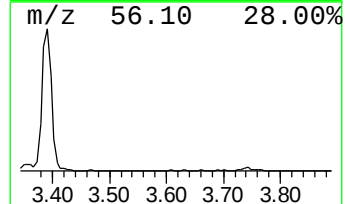
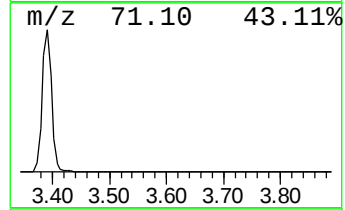
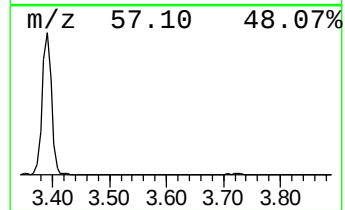
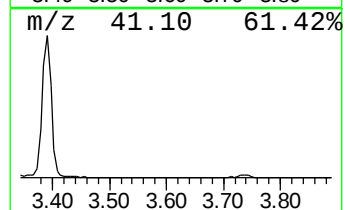
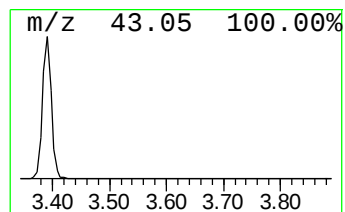
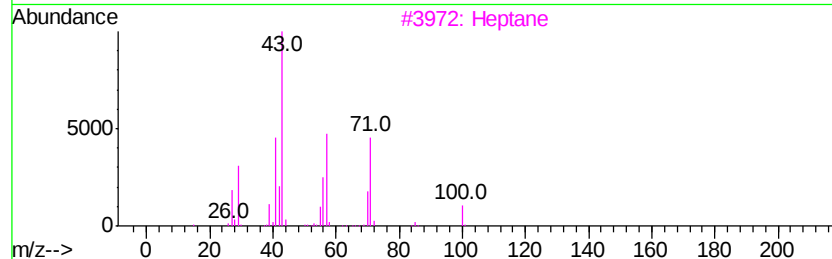
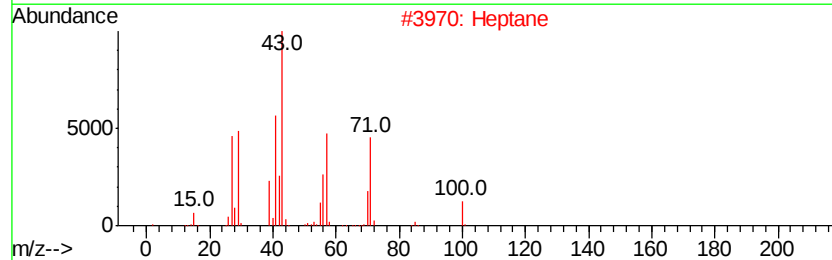
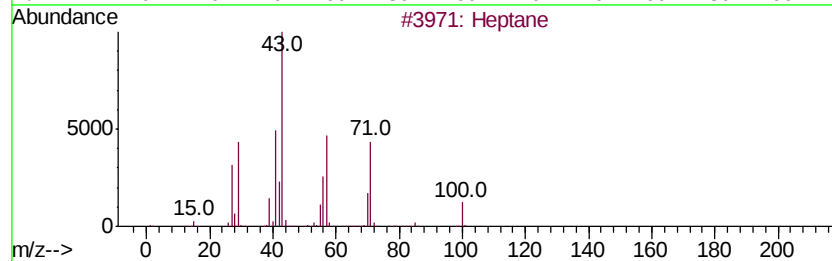
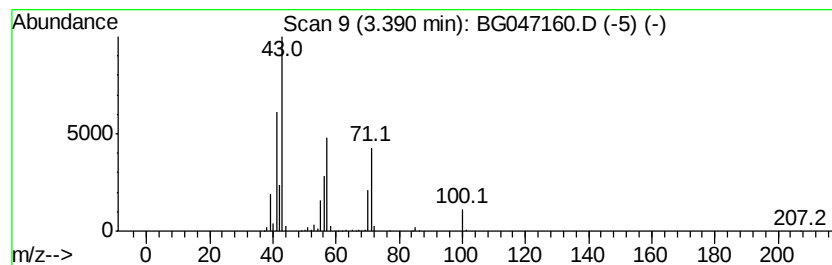
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

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.
3.39	31.47 ng/ul	531813	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane	100	C7H16	000142-82-5	91
2		Heptane	100	C7H16	000142-82-5	91
3		Heptane	100	C7H16	000142-82-5	91
4		Heptane	100	C7H16	000142-82-5	80
5		Hydroxylamine, O-(2-methylpropyl)-	89	C4H11NO	005618-62-2	42



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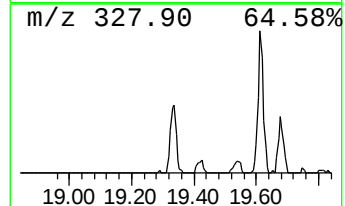
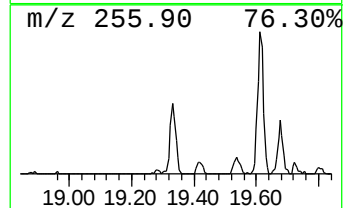
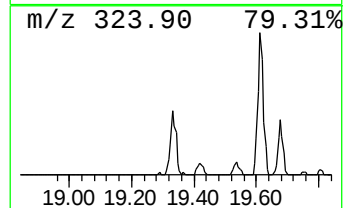
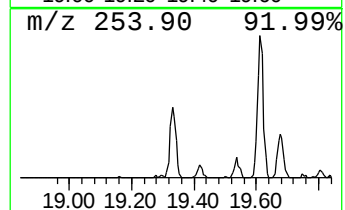
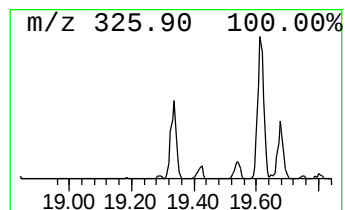
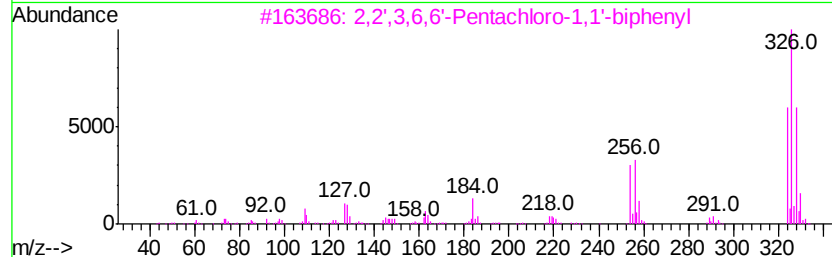
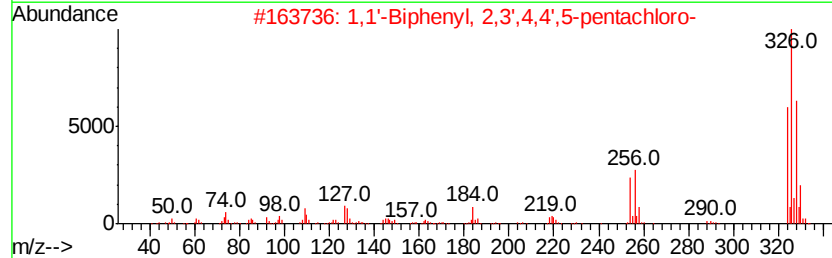
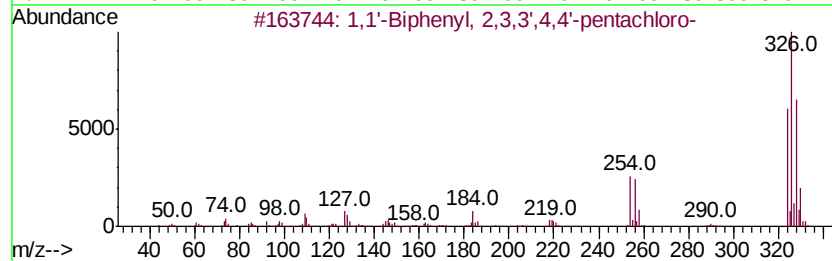
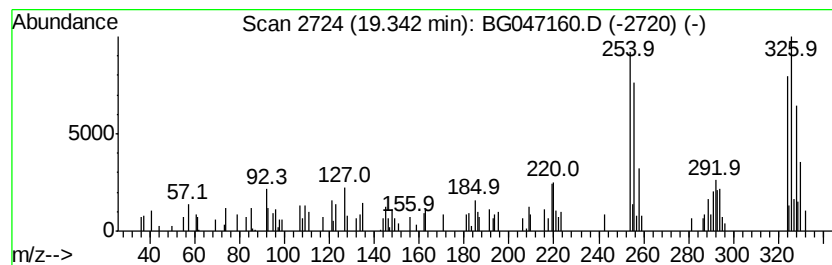
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
19.34	2.88 ng/ul	126848	Phenanthrene-d10	17.43		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'	-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	93
2	1,1'	-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	93
3	2,2',3,6,6'	-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-54-9	91
4	1,1'-Biphenyl, 2,2',3,3',4-penta...	324	C12H5Cl5	052663-62-4	91	
5	2,2',3,4',6-Pentachloro-1,1'-bip...	324	C12H5Cl5	068194-05-8	91	



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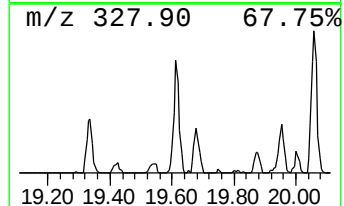
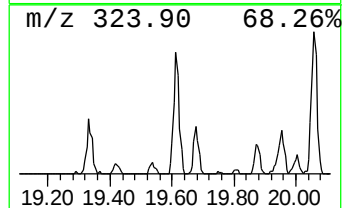
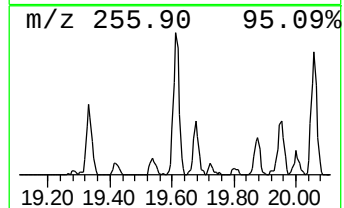
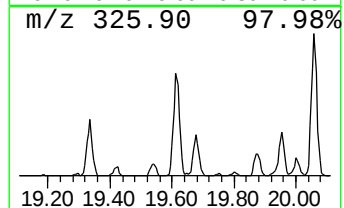
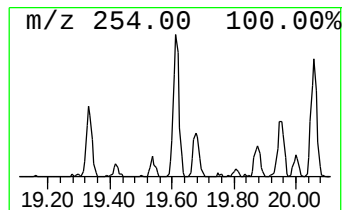
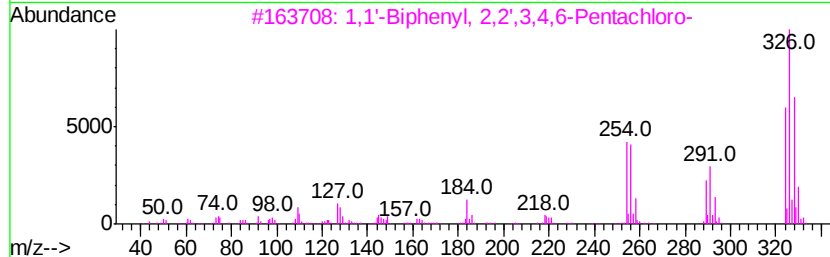
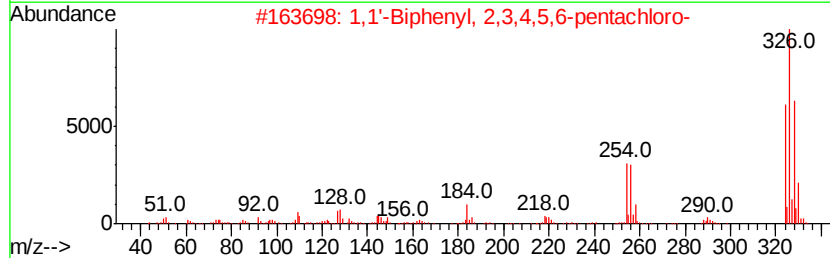
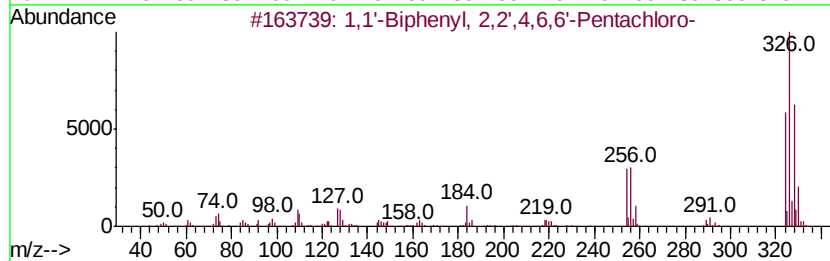
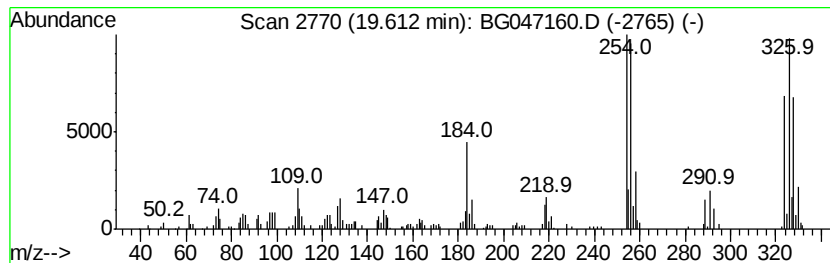
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.61	3.97 ng/ul	207751	Chrysene-d12	21.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',4,6,6'-Penta...	324	C12H5Cl5	056558-16-8	95
2		1,1'-Biphenyl, 2,3,4,5,6-pentach...	324	C12H5Cl5	018259-05-7	95
3		1,1'-Biphenyl, 2,2',3,4,6-Pentac...	324	C12H5Cl5	055215-17-3	95
4		2,2',3,6,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-54-9	95
5		1,1'-Biphenyl, 2,3,4',5,6-Pentac...	324	C12H5Cl5	068194-11-6	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
Data File : BG047160.D
Acq On : 30 Oct 2020 18:55
Operator : CG/JU
Sample : L4505-13
Misc : GCMS CONFIRMATION
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0BH2

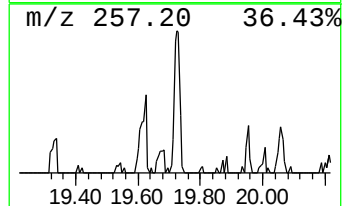
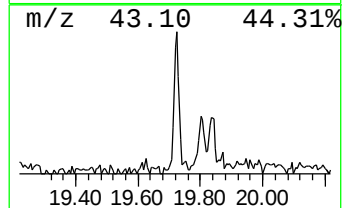
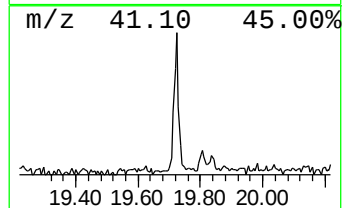
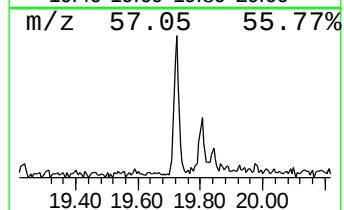
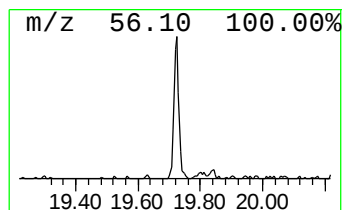
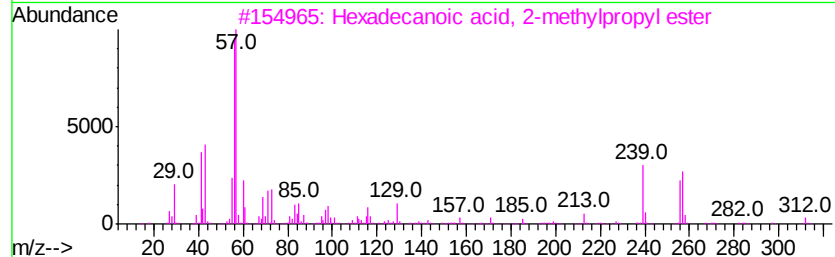
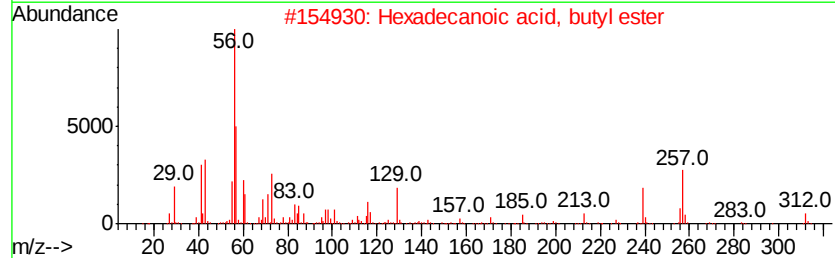
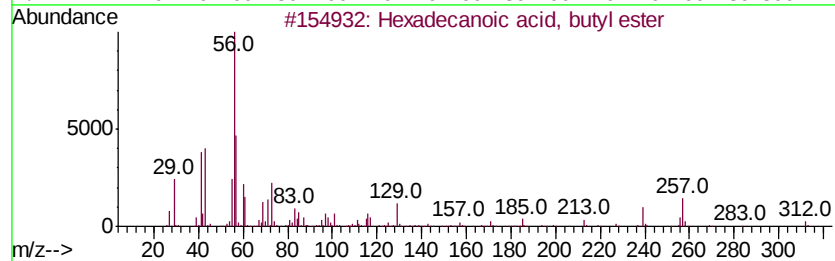
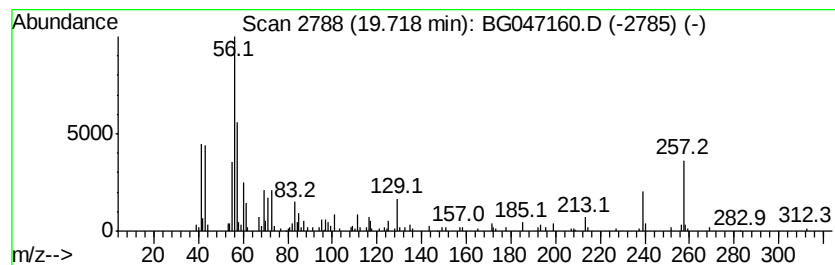
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Hexadecanoic acid, butyl ester Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.72	2.14 ng/ul	111948	Chrysene-d12	21.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	93
2		Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	91
3		Hexadecanoic acid, 2-methylpropyl...	312	C20H40O2	000110-34-9	87
4		Hexadecanoic acid, 1,1-dimethyl...	312	C20H40O2	031158-91-5	58
5		Cyclohexanol, 4-amino-, trans-	115	C6H13NO	027489-62-9	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
Data File : BG047160.D
Acq On : 30 Oct 2020 18:55
Operator : CG/JU
Sample : L4505-13
Misc : GCMS CONFIRMATION
ALS Vial : 50 Sample Multiplier: 1

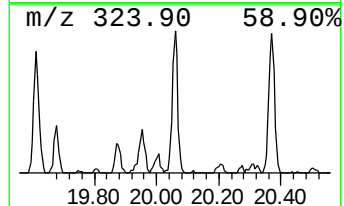
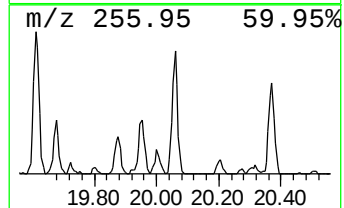
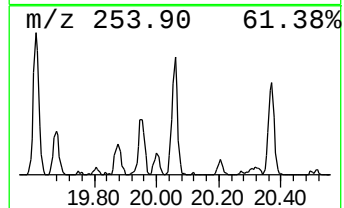
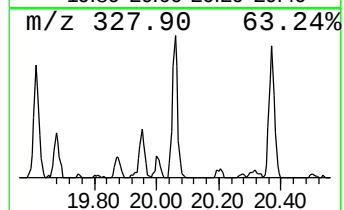
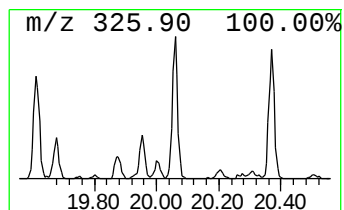
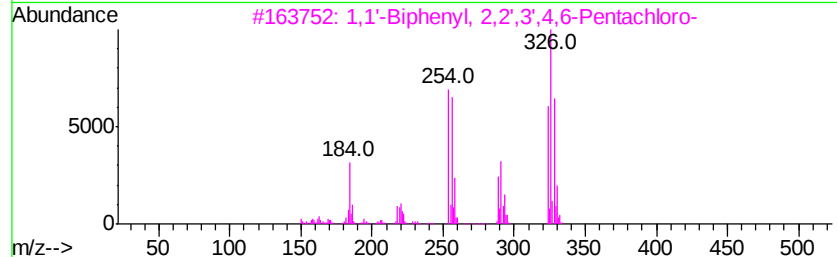
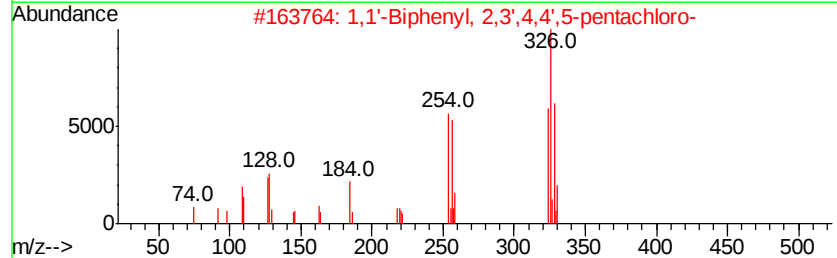
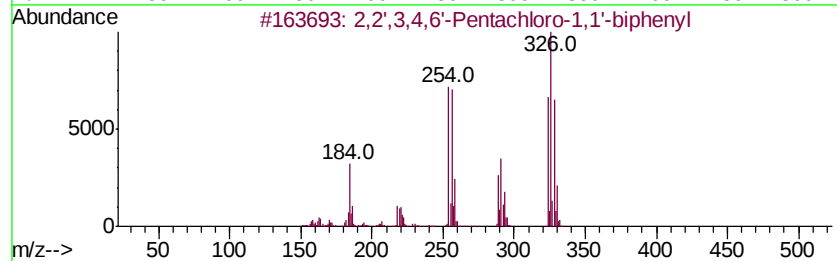
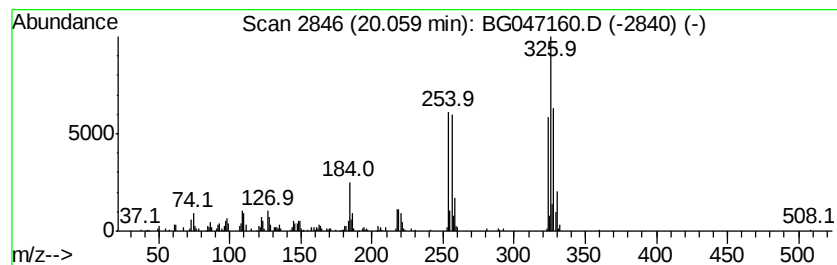
Instrument :
BNA_G
ClientSampleId :
C0BH2

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.06	4.00 ng/ul	209639	Chrysene-d12	21.70
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2,2',3,4,6'-Pentachloro-1,1'-bip...	324 C12H5Cl5	073575-57-2	99
2	1,1'-Biphenyl, 2,3',4,4',5-penta...	324 C12H5Cl5	031508-00-6	96
3	1,1'-Biphenyl, 2,2',3',4,6-Penta...	324 C12H5Cl5	060233-25-2	96
4	2,2',3,4,4'-Pentachloro-1,1'-bip...	324 C12H5Cl5	065510-45-4	96
5	2,3,3',4',5'- Pentachloro-1,1'-b...	324 C12H5Cl5	076842-07-4	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
Data File : BG047160.D
Acq On : 30 Oct 2020 18:55
Operator : CG/JU
Sample : L4505-13
Misc : GCMS CONFIRMATION
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0BH2

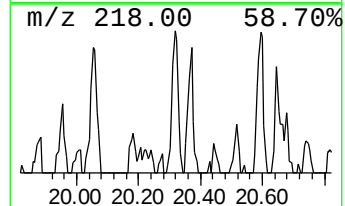
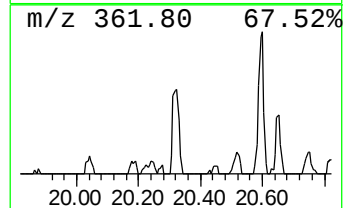
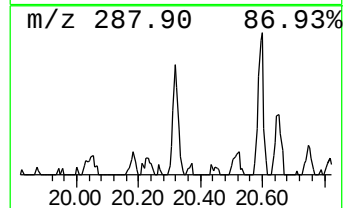
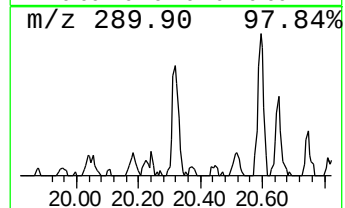
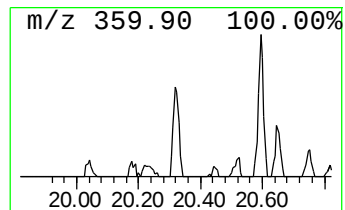
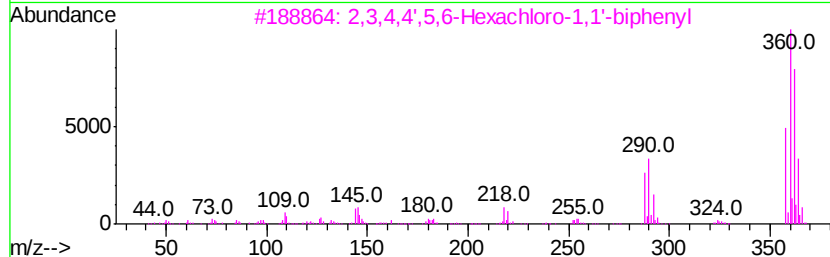
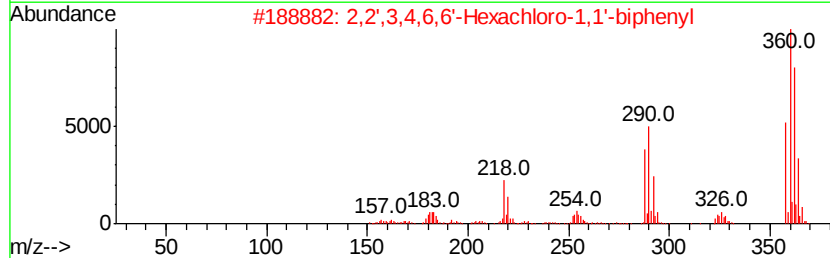
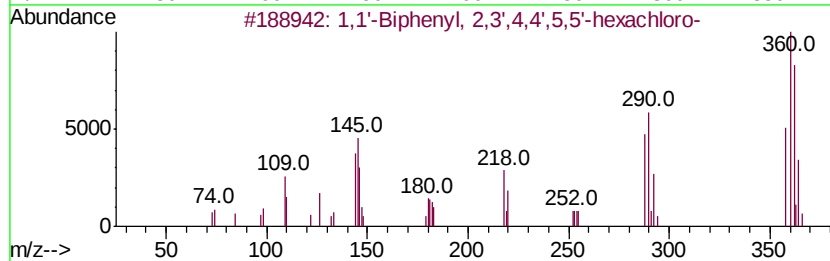
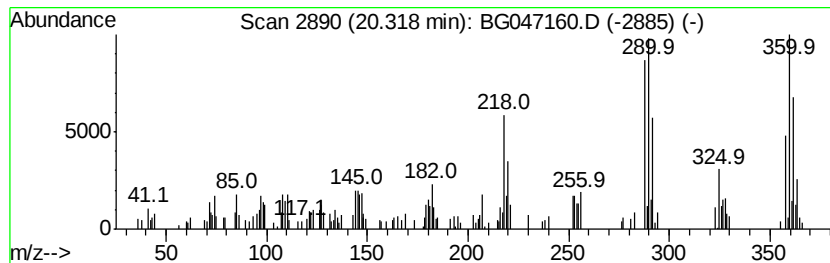
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.32	2.23 ng/ul	116993	Chrysene-d12	21.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3',4,4',5,5'-he...	358	C12H4Cl6	052663-72-6	96
2		2,2',3,4,6,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-40-5	95
3		2,3,4,4',5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-63-6	95
4		1,1'-Biphenyl, 2,2',4,4',5,5'-he...	358	C12H4Cl6	035065-27-1	95
5		1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
Data File : BG047160.D
Acq On : 30 Oct 2020 18:55
Operator : CG/JU
Sample : L4505-13
Misc : GCMS CONFIRMATION
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0BH2

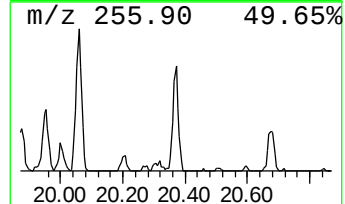
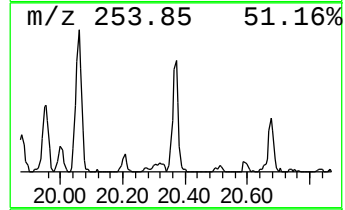
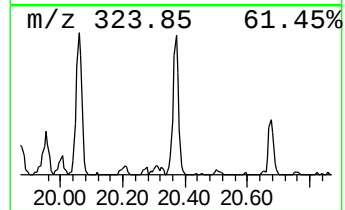
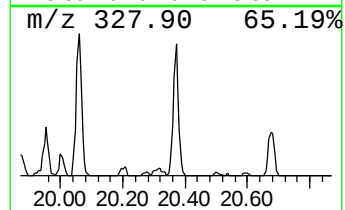
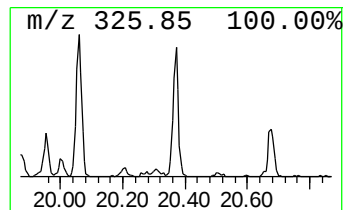
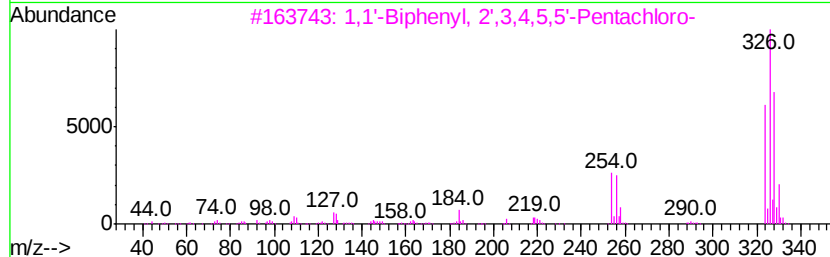
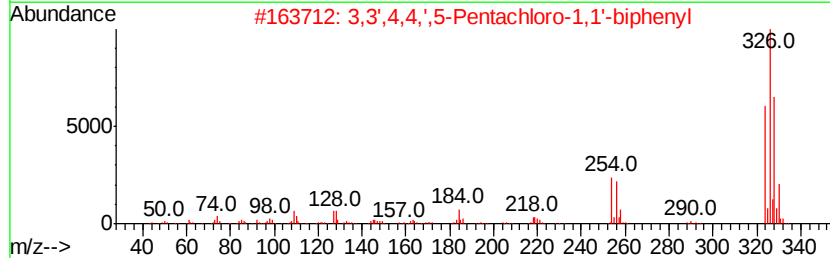
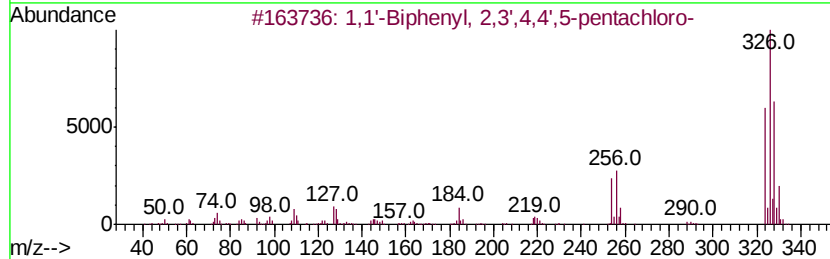
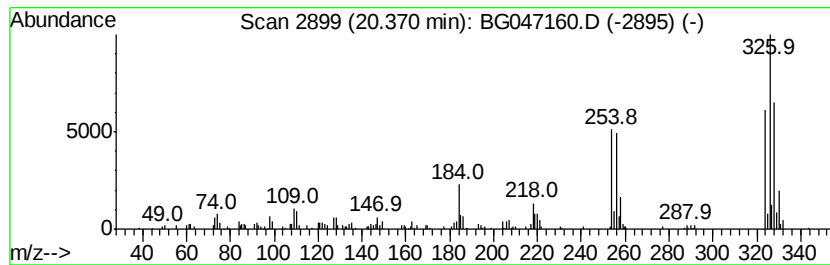
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.37	3.37 ng/ul	176329	Chrysene-d12	21.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	99
2		3,3',4,4',5-Pentachloro-1,1'-bi...	324	C12H5Cl5	057465-28-8	97
3		1,1'-Biphenyl, 2',3,4,5,5'-Penta...	324	C12H5Cl5	070424-70-3	97
4		1,1'-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	96
5		2,3,3',5,6-Pentachloro-1,1'-biph...	324	C12H5Cl5	074472-36-9	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
Data File : BG047160.D
Acq On : 30 Oct 2020 18:55
Operator : CG/JU
Sample : L4505-13
Misc : GCMS CONFIRMATION
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0BH2

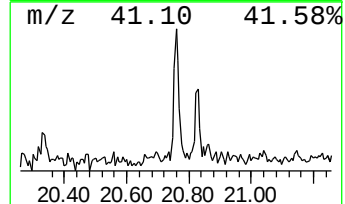
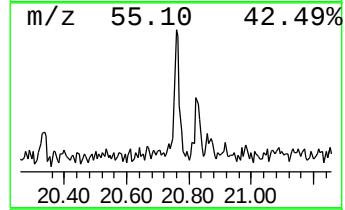
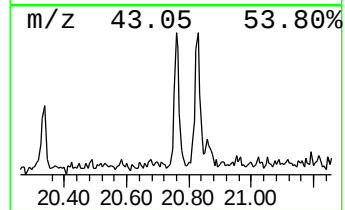
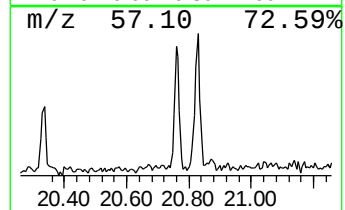
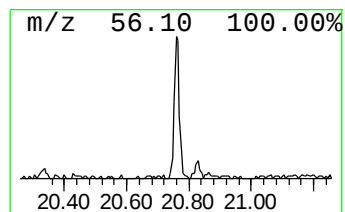
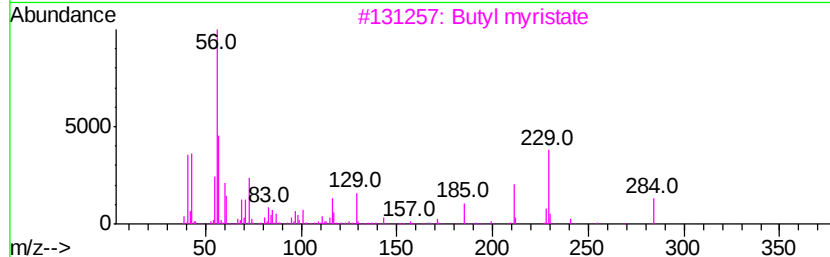
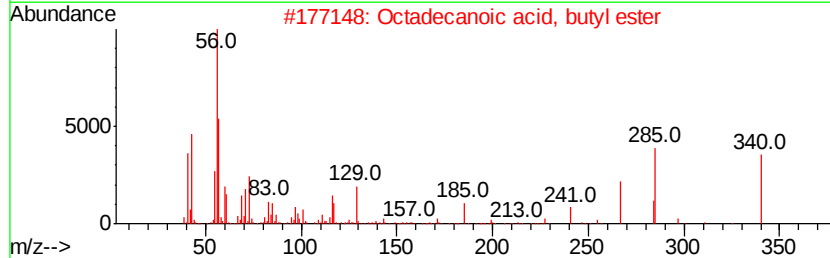
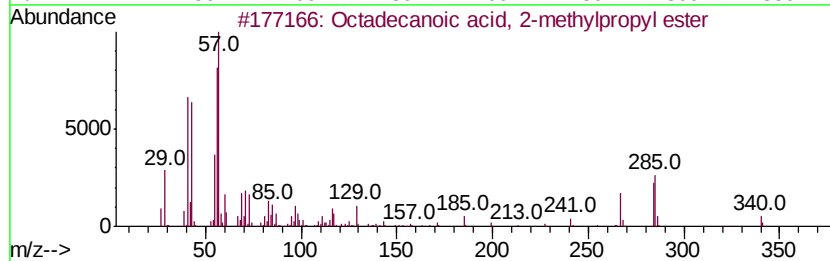
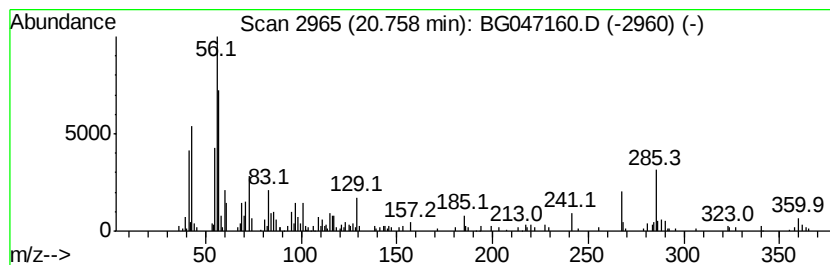
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Octadecanoic acid, 2-methyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.76	2.33 ng/ul	121816	Chrysene-d12	21.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid, 2-methylpropyl...	340	C22H44O2	000646-13-9	72
2		Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	59
3		Butyl myristate	284	C18H36O2	000110-36-1	52
4		Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	47
5		1,3-Hexanediol, 2-ethyl-	146	C8H18O2	000094-96-2	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102920\
Data File : BG047160.D
Acq On : 30 Oct 2020 18:55
Operator : CG/JU
Sample : L4505-13
Misc : GCMS CONFIRMATION
ALS Vial : 50 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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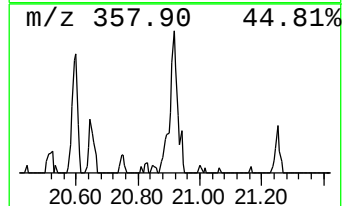
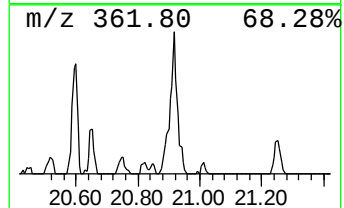
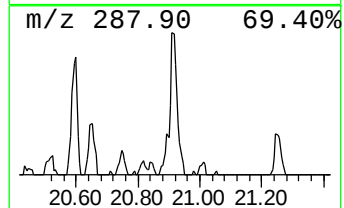
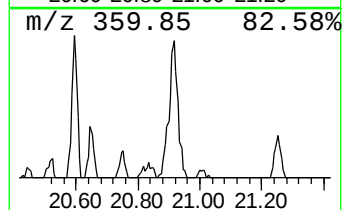
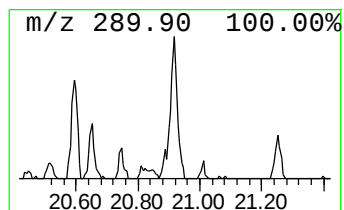
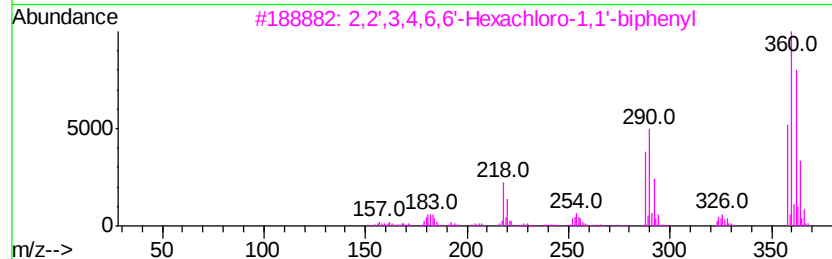
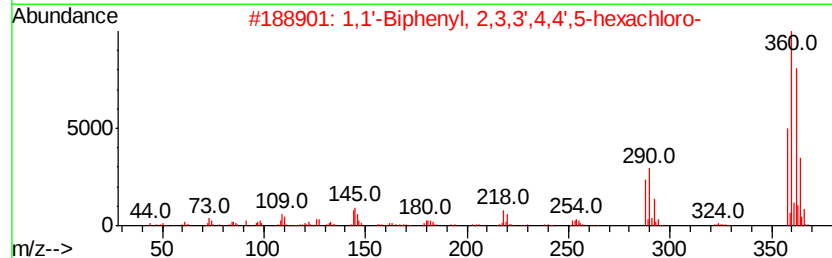
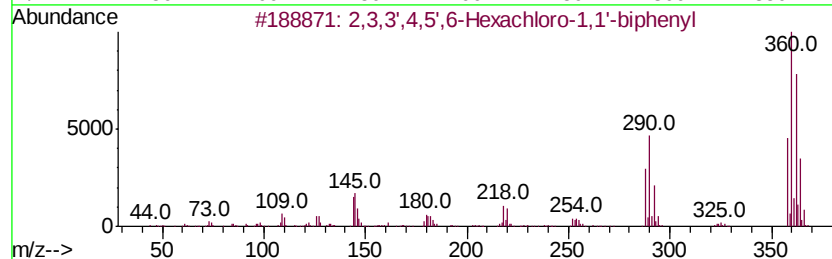
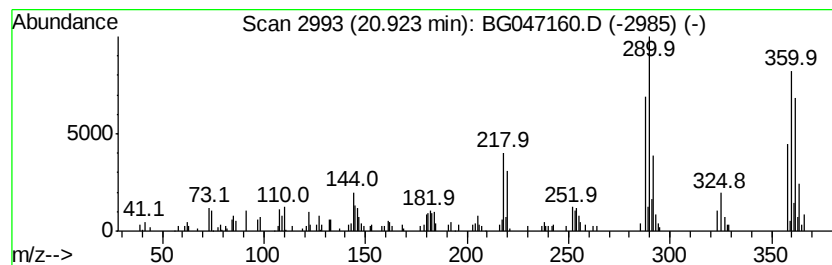
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.92	3.27 ng/ul	171185	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	98
2		1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	97
3		2,2',3,4,6,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-40-5	96
4		2,3,3',4',5,6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-44-9	95
5		1,1'-Biphenyl, 2,2',3,4,5,6'-Hex...	358	C12H4Cl6	068194-15-0	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG102920\
Data File : BG047160.D
Acq On : 30 Oct 2020 18:55
Operator : CG/JU
Sample : L4505-13
Misc : GCMS CONFIRMATION
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0BH2

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...	3.39	31.5	ng/ul	531813	1	8.06	338016	20.0
1,1'-Biphenyl, 2,...	19.34	2.9	ng/ul	126848	4	17.43	881370	20.0
1,1'-Biphenyl, 2,...	19.61	4.0	ng/ul	207751	5	21.70	1047730	20.0
Hexadecanoic acid...	19.72	2.1	ng/ul	111948	5	21.70	1047730	20.0
2,2',3,4,6'-Penta...	20.06	4.0	ng/ul	209639	5	21.70	1047730	20.0
1,1'-Biphenyl, 2,...	20.32	2.2	ng/ul	116993	5	21.70	1047730	20.0
1,1'-Biphenyl, 2,...	20.37	3.4	ng/ul	176329	5	21.70	1047730	20.0
Octadecanoic acid...	20.76	2.3	ng/ul	121816	5	21.70	1047730	20.0
2,3,3',4,5',6-Hex...	20.92	3.3	ng/ul	171185	5	21.70	1047730	20.0