

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111920\
 Data File : BG047364.D
 Acq On : 19 Nov 2020 22:45
 Operator : CG/JU
 Sample : L4710-20
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AY3

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-BG111920MA.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	3.615	43	47	52	rBV6	6363	11172	0.50%	0.061%
2	4.162	136	140	144	rVB6	4408	7434	0.33%	0.041%
3	4.279	158	160	161	rBV2	3658	3364	0.15%	0.018%
4	4.303	161	164	170	rVB4	6762	10100	0.45%	0.055%
5	4.391	177	179	183	rVB3	3220	3756	0.17%	0.021%
6	5.037	286	289	290	rVB3	3622	3290	0.15%	0.018%
7	5.331	333	339	345	rBV3	7463	13003	0.58%	0.071%
8	5.654	391	394	396	rBV	2757	3247	0.14%	0.018%
9	5.707	401	403	408	rVB2	4103	5455	0.24%	0.030%
10	6.330	507	509	512	rBV3	2547	3249	0.14%	0.018%
11	6.776	583	585	588	rBV3	2632	3137	0.14%	0.017%
12	6.835	591	595	599	rVB	2209	3573	0.16%	0.020%
13	6.923	608	610	613	rVB2	2957	3264	0.15%	0.018%
14	7.399	684	691	701	rVB	147464	265615	11.82%	1.455%
15	7.587	716	723	731	rVB	88017	155741	6.93%	0.853%
16	7.798	753	759	765	rVB	141986	238840	10.63%	1.308%
17	8.274	834	840	848	rBV	196788	339705	15.11%	1.860%
18	8.962	951	957	967	rVB	163813	289021	12.86%	1.583%
19	9.097	977	980	982	rBV2	4921	5210	0.23%	0.029%
20	9.173	989	993	996	rBV3	3514	4299	0.19%	0.024%
21	9.438	1031	1038	1046	rVB2	131340	242539	10.79%	1.328%
22	9.602	1061	1066	1068	rBV2	2723	4250	0.19%	0.023%
23	9.679	1076	1079	1082	rVV2	2626	3872	0.17%	0.021%
24	9.714	1082	1085	1088	rVB2	3705	4742	0.21%	0.026%
25	9.861	1105	1110	1113	rBV4	6125	7621	0.34%	0.042%
26	10.113	1147	1153	1156	rBV	2723	5368	0.24%	0.029%
27	10.166	1156	1162	1169	rVB2	123961	211548	9.41%	1.159%
28	10.307	1183	1186	1189	rBV2	2580	3733	0.17%	0.020%
29	10.707	1246	1254	1263	rBV	193457	350148	15.58%	1.918%
30	10.865	1275	1281	1282	rBV	2383	3397	0.15%	0.019%
31	11.100	1313	1321	1329	rBV	252064	450463	20.04%	2.467%
32	11.224	1335	1342	1350	rVB2	92444	174689	7.77%	0.957%
33	11.594	1400	1405	1407	rBV	2070	3088	0.14%	0.017%
34	12.663	1585	1587	1590	rVB3	3677	3835	0.17%	0.021%

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35	12.728	1595	1598	1602	rBV2	2668	4319	0.19%	0.024%
36	12.939	1633	1634	1640	rVB3	3780	4899	0.22%	0.027%
37	13.492	1725	1728	1733	rBV2	4669	5731	0.25%	0.031%
38	13.556	1736	1739	1742	rBV	2499	3707	0.16%	0.020%
39	13.715	1762	1766	1771	rVB4	5480	9165	0.41%	0.050%
40	13.762	1771	1774	1782	rBV4	7629	14062	0.63%	0.077%
41	14.279	1855	1862	1866	rBV	327821	492805	21.93%	2.699%
42	14.320	1866	1869	1879	rVB	178445	261337	11.63%	1.431%
43	14.514	1900	1902	1906	rBV2	4518	4512	0.20%	0.025%
44	14.579	1906	1913	1921	rVB	334113	546226	24.30%	2.991%
45	14.884	1957	1965	1972	rBV	388242	636241	28.31%	3.484%
46	15.031	1984	1990	1996	rBV2	150514	244805	10.89%	1.341%
47	15.202	2017	2019	2022	rVB3	4719	4515	0.20%	0.025%
48	15.272	2026	2031	2036	rBV5	6143	12195	0.54%	0.067%
49	15.489	2063	2068	2070	rBV	3035	5003	0.22%	0.027%
50	15.672	2095	2099	2103	rBV5	4968	10601	0.47%	0.058%
51	15.713	2104	2106	2111	rVB3	4778	5685	0.25%	0.031%
52	15.865	2126	2132	2144	rBV	439520	675484	30.05%	3.699%
53	15.965	2144	2149	2160	rVB2	109232	173092	7.70%	0.948%
54	16.106	2166	2173	2177	rBV7	8218	17454	0.78%	0.096%
55	16.529	2244	2245	2249	rVB4	5402	5056	0.22%	0.028%
56	16.565	2249	2251	2255	rBV4	3764	4330	0.19%	0.024%
57	16.647	2260	2265	2267	rBV4	4001	5391	0.24%	0.030%
58	16.853	2294	2300	2305	rVB4	12240	20586	0.92%	0.113%
59	16.894	2305	2307	2309	rBV3	4115	4032	0.18%	0.022%
60	17.076	2335	2338	2341	rBV5	5983	6849	0.30%	0.038%
61	17.140	2345	2349	2356	rBV7	6226	9223	0.41%	0.051%
62	17.264	2366	2370	2372	rBV4	3385	4443	0.20%	0.024%
63	17.328	2378	2381	2384	rBV3	3526	4768	0.21%	0.026%
64	17.411	2390	2395	2398	rBV5	6862	12395	0.55%	0.068%
65	17.516	2411	2413	2416	rVB4	4802	3777	0.17%	0.021%
66	17.616	2424	2430	2436	rBV	519595	800363	35.61%	4.383%
67	17.657	2436	2437	2441	rVV2	37862	34898	1.55%	0.191%
68	17.716	2441	2447	2460	rVB	475755	738326	32.85%	4.043%
69	17.875	2472	2474	2479	rVB5	4394	8169	0.36%	0.045%
70	17.910	2479	2480	2482	rBV2	5386	3811	0.17%	0.021%
71	17.939	2484	2485	2487	rBV2	6015	4367	0.19%	0.024%

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72	18.186	2523	2527	2532	rBV3	52644	75449	3.36%	0.413%
73	18.257	2535	2539	2543	rVB6	12636	19602	0.87%	0.107%
74	18.363	2553	2557	2558	rBV4	5625	7148	0.32%	0.039%
75	18.439	2568	2570	2573	rBV3	4970	5757	0.26%	0.032%
76	18.486	2573	2578	2584	rBV3	196213	313861	13.96%	1.719%
77	18.880	2639	2645	2654	rBV3	1168177	2247610	100.00%	12.309%
78	18.950	2654	2657	2661	rVV6	49581	73406	3.27%	0.402%
79	19.009	2662	2667	2675	rVB4	98770	166524	7.41%	0.912%
80	19.162	2690	2693	2697	rVB5	26961	32302	1.44%	0.177%
81	19.261	2705	2710	2716	rBV2	87123	155399	6.91%	0.851%
82	19.355	2721	2726	2730	rBV3	106461	150251	6.68%	0.823%
83	19.408	2732	2735	2741	rVB6	44446	62801	2.79%	0.344%
84	19.497	2745	2750	2755	rBV3	146170	214950	9.56%	1.177%
85	19.585	2761	2765	2772	rBV6	145093	288606	12.84%	1.581%
86	19.649	2772	2776	2779	rVV	91764	125242	5.57%	0.686%
87	19.685	2779	2782	2787	rVB5	96490	129102	5.74%	0.707%
88	19.761	2788	2795	2797	rBV5	151075	315012	14.02%	1.725%
89	19.796	2797	2801	2807	rVB	811469	1150286	51.18%	6.299%
90	19.855	2807	2811	2813	rBV5	19035	29946	1.33%	0.164%
91	19.955	2822	2828	2831	rBV2	1090541	1726839	76.83%	9.457%
92	20.090	2847	2851	2858	rVB5	137208	200980	8.94%	1.101%
93	20.548	2925	2929	2935	rBV9	28320	50142	2.23%	0.275%
94	20.707	2953	2956	2961	rBV7	51794	70585	3.14%	0.387%
95	20.924	2988	2993	3000	rVB2	330513	458342	20.39%	2.510%
96	21.518	3089	3094	3103	rBV3	159758	250203	11.13%	1.370%
97	21.776	3136	3138	3142	rVB	47193	52571	2.34%	0.288%
98	21.900	3152	3159	3165	rBV2	453646	807366	35.92%	4.422%
99	25.061	3690	3697	3707	rVB	238408	684563	30.46%	3.749%
100	25.301	3729	3738	3747	rBV	234790	750753	33.40%	4.111%

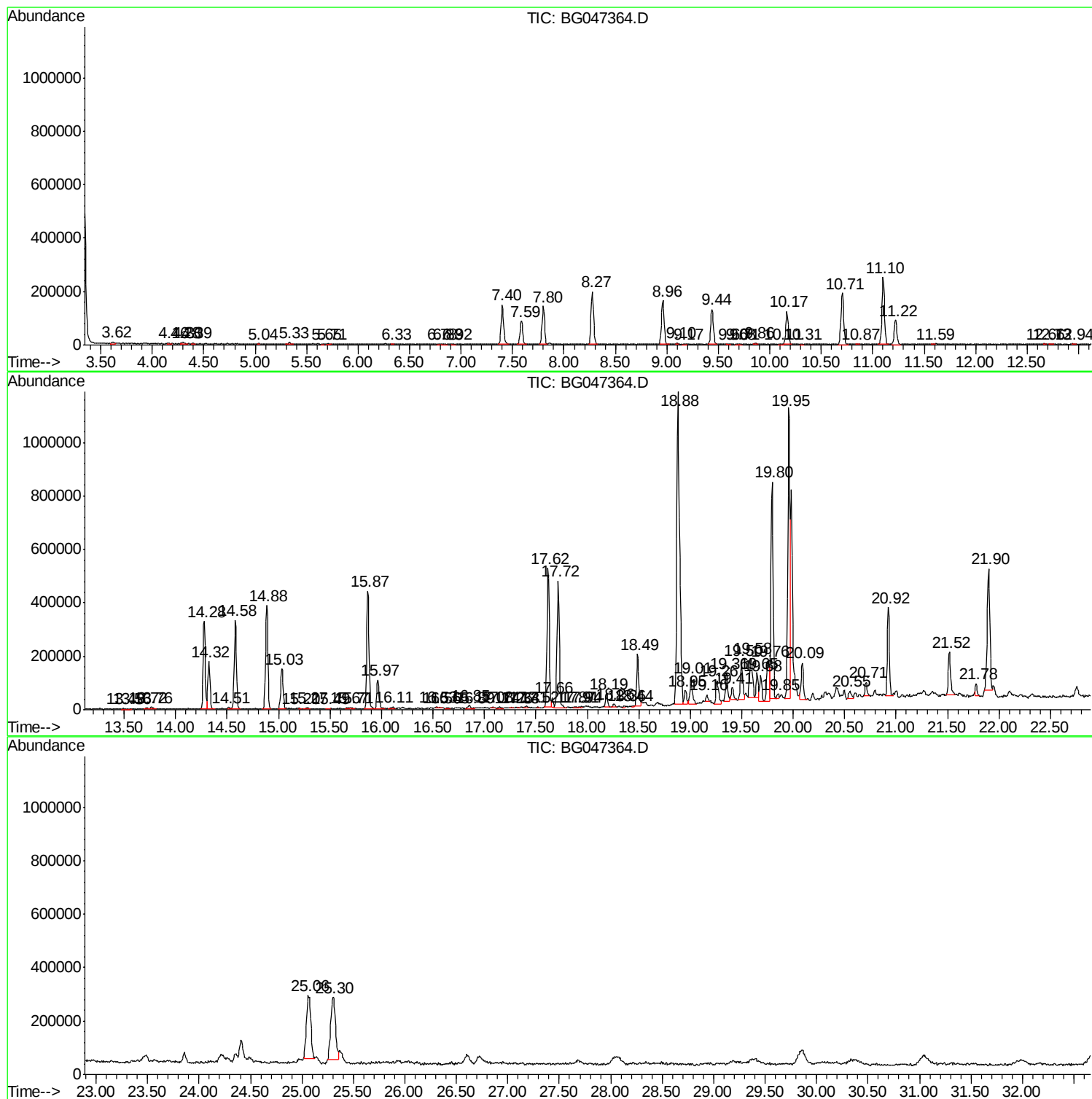
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

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



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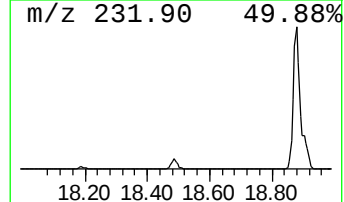
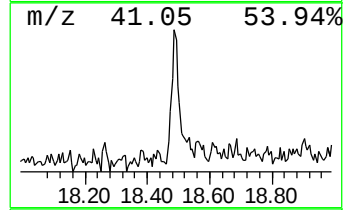
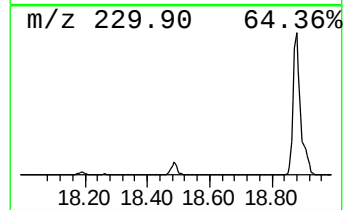
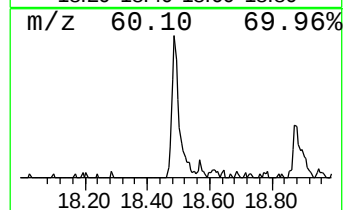
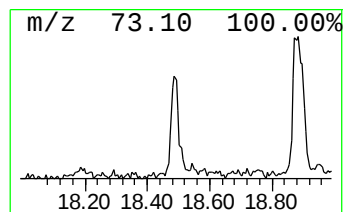
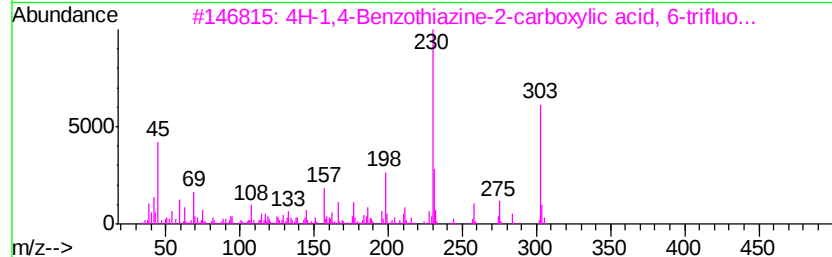
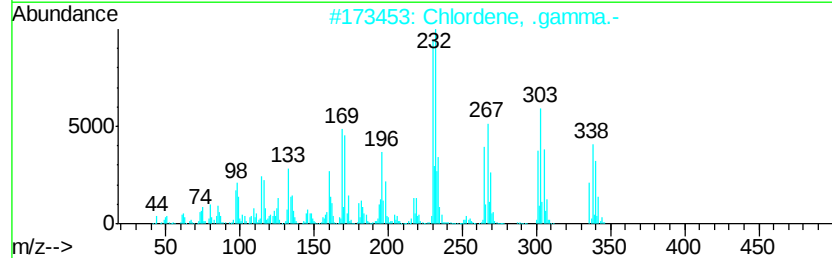
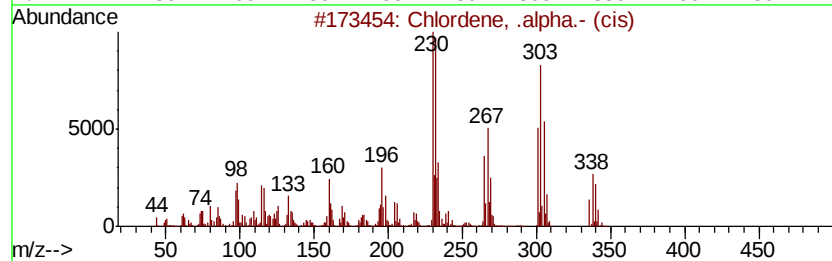
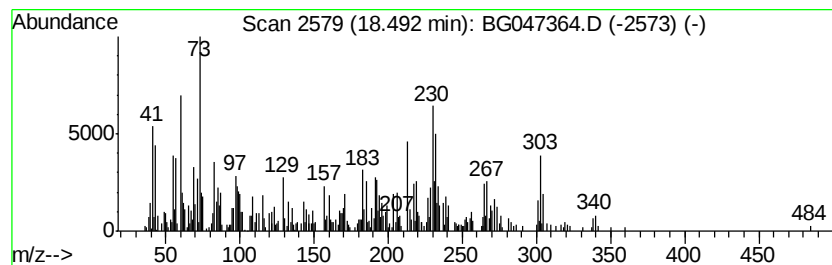
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Chlordene, .alpha.- (cis) Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.49	7.84 ng/ul	313861	Phenanthrene-d10	17.62		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Chlordene, .alpha.- (cis)	336	C10H6Cl6	1000378-50-5	86
2		Chlordene, .gamma.-	336	C10H6Cl6	1000378-50-6	38
3		4H-1,4-Benzothiazine-2-carboxyli...	303	C13H12F3N02S	101767-70-8	38
4		Benzaldehyde, 3-bromo-4-hydroxyv...	230	C8H7BrO3	002973-76-4	38
5		Phenol, 2-(3,5-dichlorophenylimi...	265	C13H9Cl2NO	303769-52-0	35



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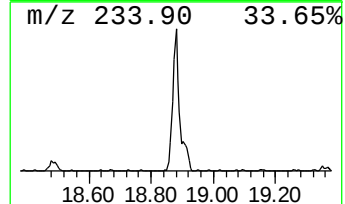
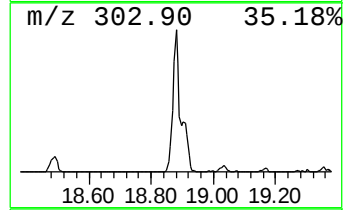
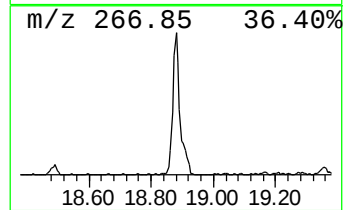
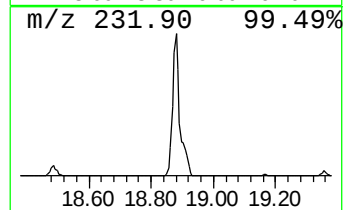
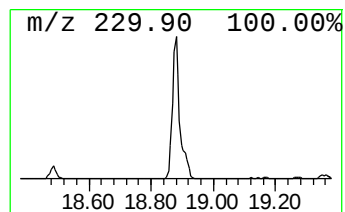
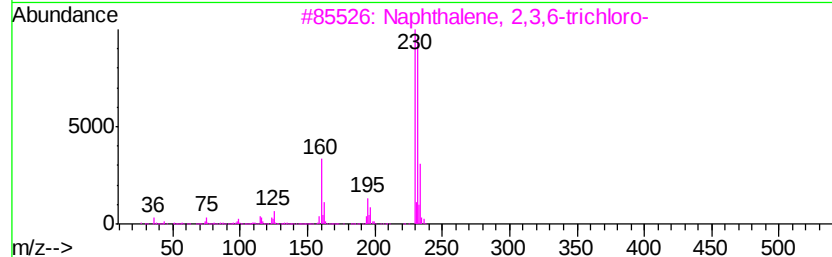
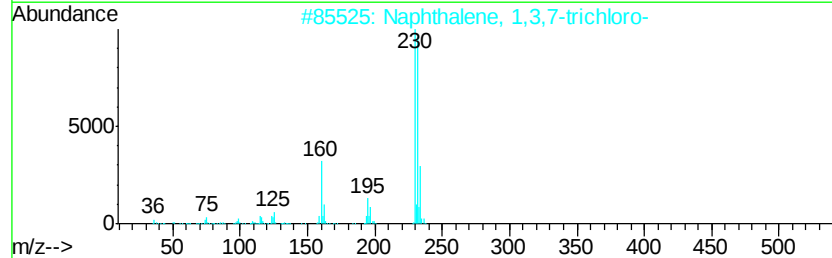
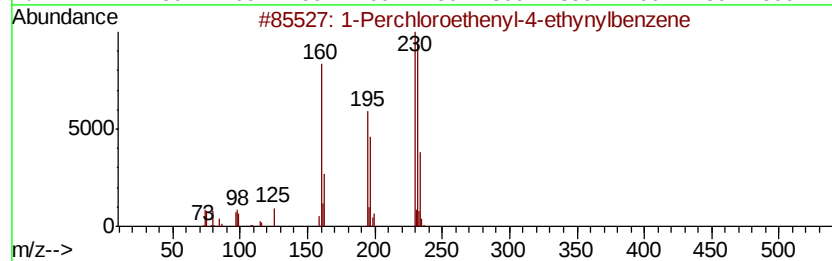
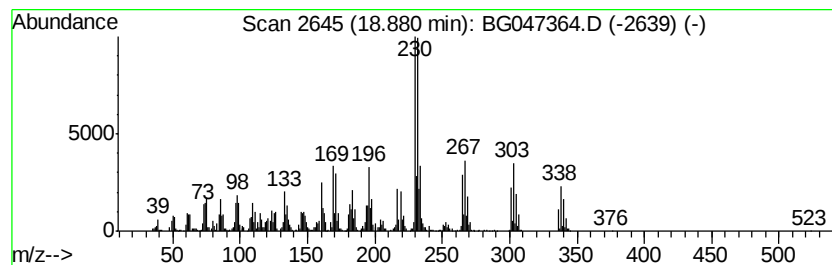
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 1-Perchloroethenyl-4-ethyny... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.88	56.16 ng/ul	2247610	Phenanthrene-d10	17.62

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Perchloroethenyl-4-ethynylbenzene	230	C10H5Cl3	1000283-76-9	91
2		Naphthalene, 1,3,7-trichloro-	230	C10H5Cl3	055720-37-1	91
3		Naphthalene, 2,3,6-trichloro-	230	C10H5Cl3	055720-40-6	90
4		Chlordene, .alpha.- (cis)	336	C10H6Cl6	1000378-50-5	68
5		Chlordene, .gamma.-	336	C10H6Cl6	1000378-50-6	58



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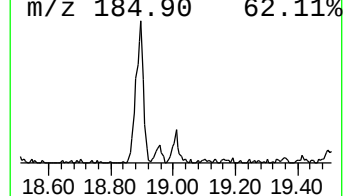
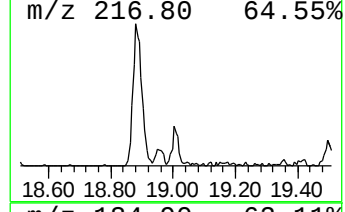
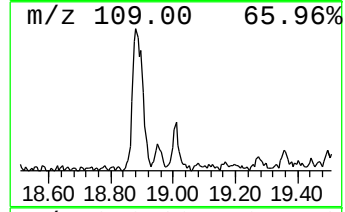
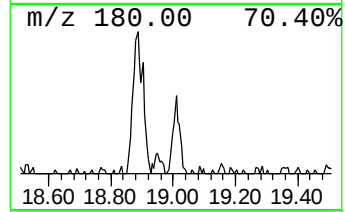
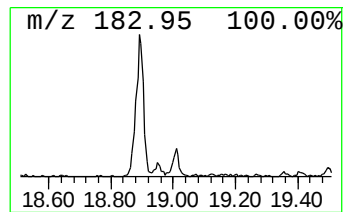
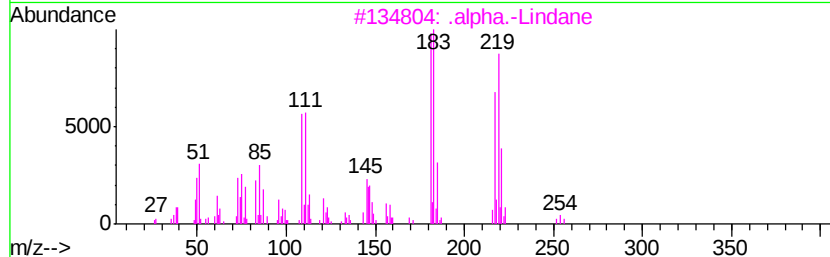
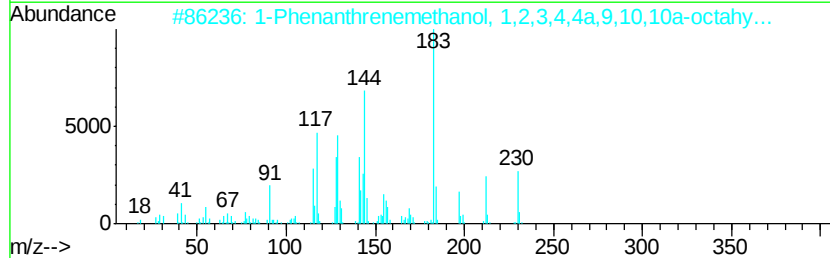
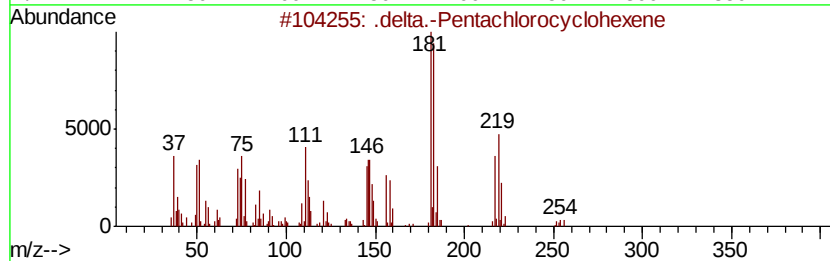
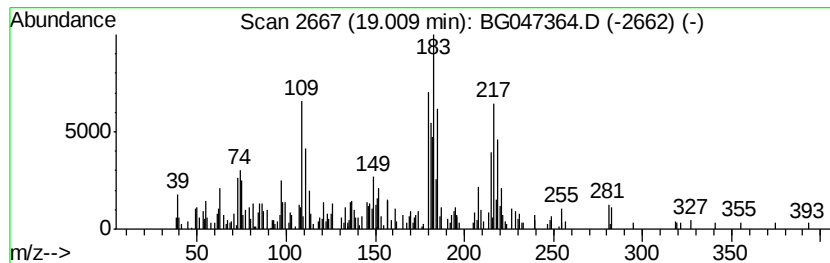
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TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.01	4.16 ng/ul	166524	Phenanthrene-d10	17.62

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	delta.-Pentachlorocyclohexene	252	C6H5Cl5	000643-15-2	30
2	1	Phenanthrenemethanol, 1,2,3,4,...	230	C16H22O	057378-57-1	25
3	.	alpha.-Lindane	288	C6H6Cl6	000319-84-6	25
4	.	alpha.-Lindane	288	C6H6Cl6	000319-84-6	22
5	2	Methyl-5-p-tolylpyridine	183	C13H13N	030456-56-5	22



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Data File : BG047364.D
Acq On : 19 Nov 2020 22:45
Operator : CG/JU
Sample : L4710-20
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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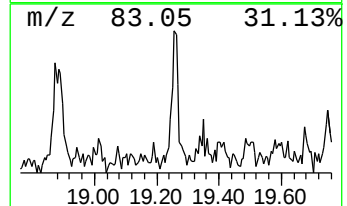
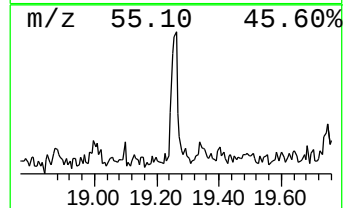
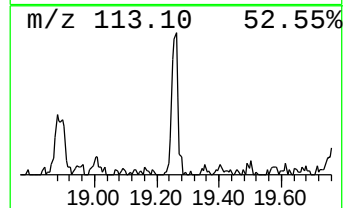
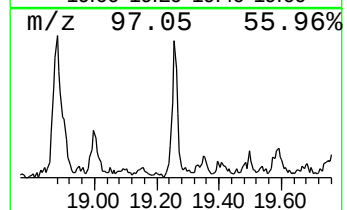
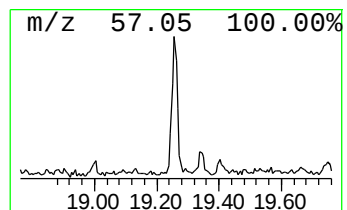
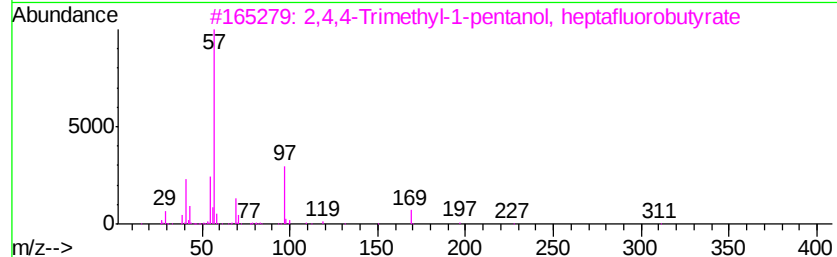
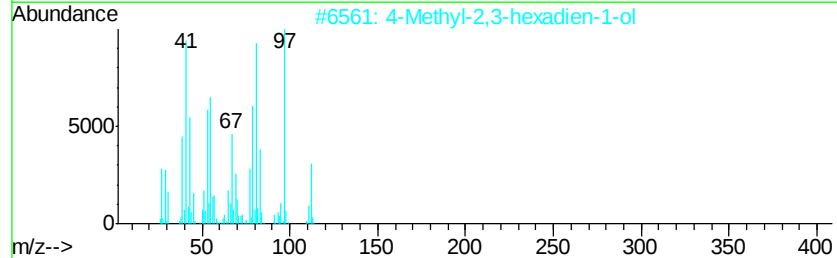
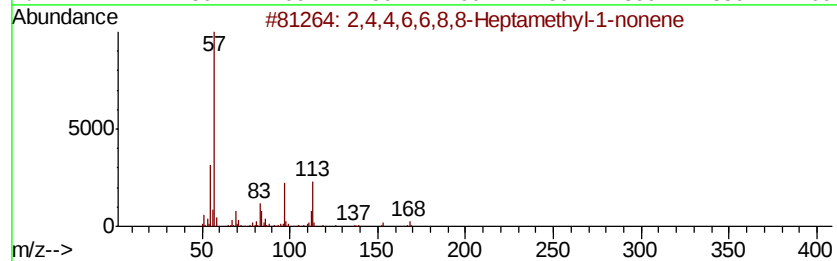
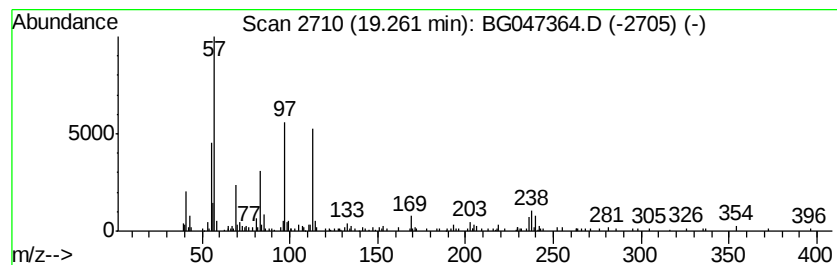
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.26	3.88 ng/ul	155399	Phenanthrene-d10	17.62

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	64		
2	4-Methyl-2,3-hexadien-1-ol	112	C7H12O	1000196-09-6	27		
3	2,4,4-Trimethyl-1-pentanol, hept...	326	C12H17F7O2	1000365-01-4	27		
4	Methyl-3,4-O-furvlidene.beta.l-a...	242	C11H14O6	053547-58-3	22		
5	2,5-Pyrrolidinedione, 1-methyl-	113	C5H7N02	001121-07-9	22		



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111920\
Data File : BG047364.D
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ClientSampleId :
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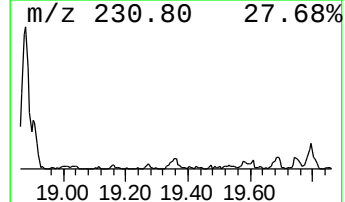
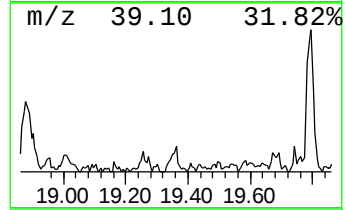
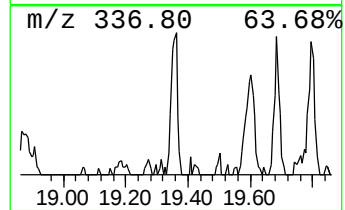
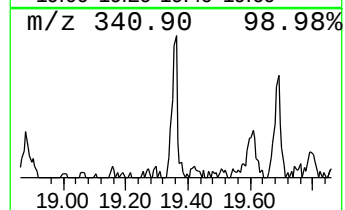
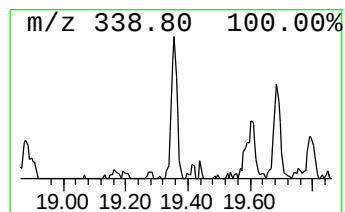
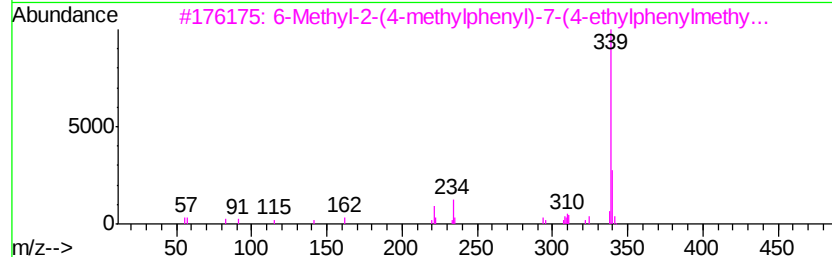
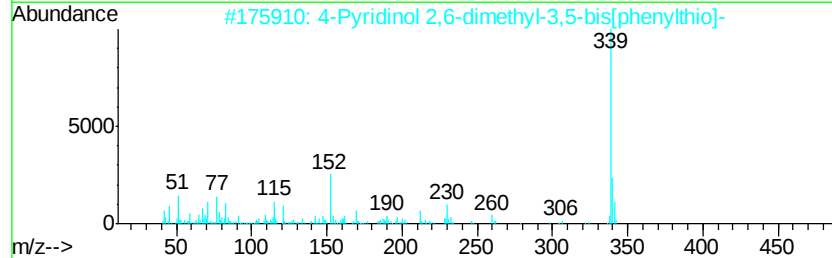
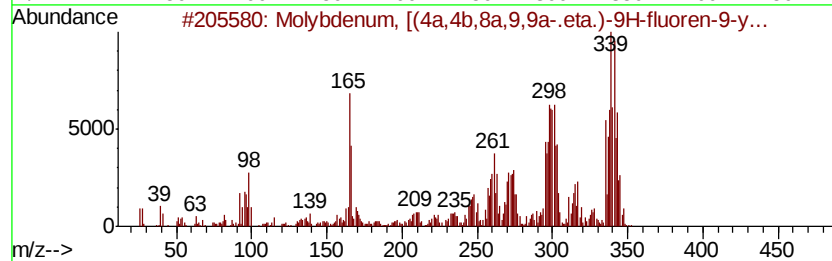
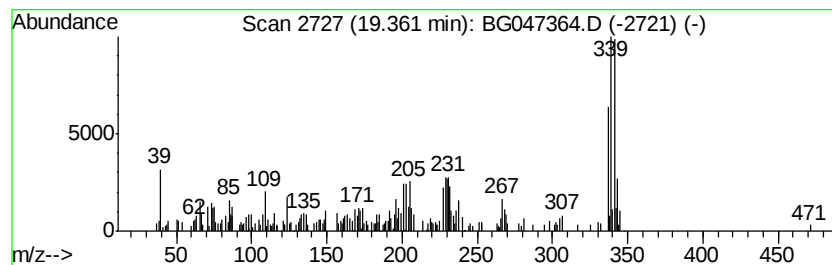
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-02 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.36	3.75 ng/ul	150251	Phenanthrene-d10	17.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Molybdenum, [(4a,4b,8a,9,9a-eta...	386	C22H24Mo	125312-19-8	30
2		4-Pyridinol 2,6-dimethyl-3,5-bis...	339	C19H17NOS2	1000252-21-8	25
3		6-Methyl-2-(4-methylphenyl)-7-(4...	339	C25H25N	080193-45-9	14
4		1-Ethyl-1-hexadecyloxv-1-silacyc...	368	C23H48OSi	1000279-38-6	14
5		Butanamide, N-(1-naphthyl)-2,2,3...	339	C14H8F7NO	1000307-28-3	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111920\
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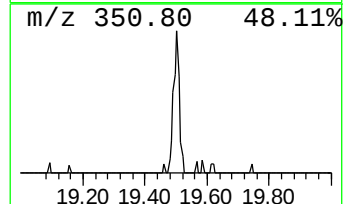
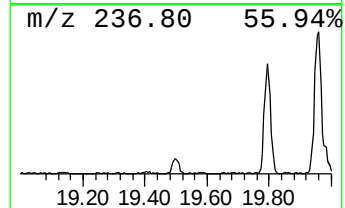
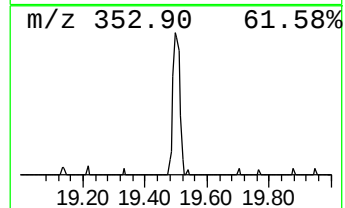
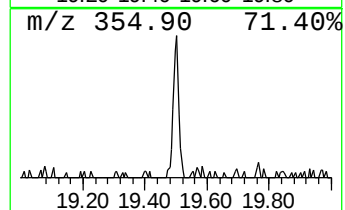
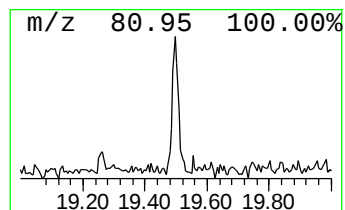
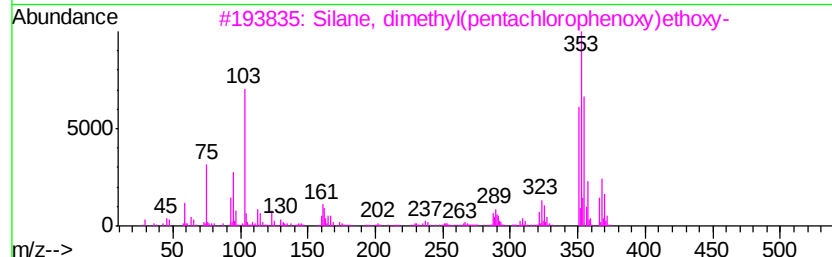
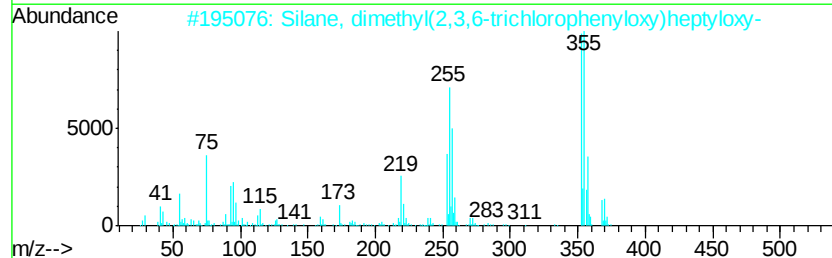
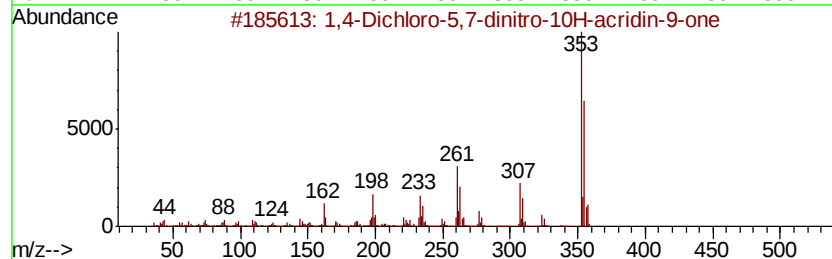
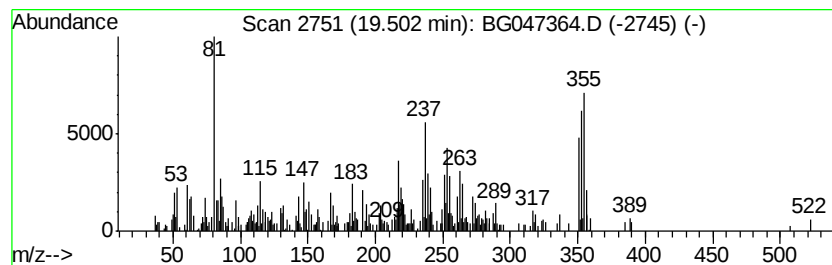
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-03 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.50	5.37 ng/ul	214950	Phenanthrene-d10	17.62

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Dichloro-5,7-dinitro-10H-acr...	353	C13H5Cl2N3O5	148877-13-8	18
2			Silane, dimethyl(2,3,6-trichloro...	368	C15H23Cl3O2Si	1000346-82-6	14
3			Silane, dimethyl(pentachlorophen...	366	C10H11Cl5O2Si	1000347-47-6	14
4			Nickel(II), [O-methyl-N(1),N(4)-...	353	C16H13N3NiO3	1000164-81-5	11
5			Silane, diethylisohexyloxy(2,4,5...	382	C16H25Cl3O2Si	1000363-35-3	10



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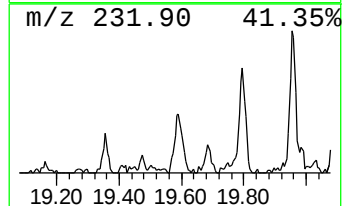
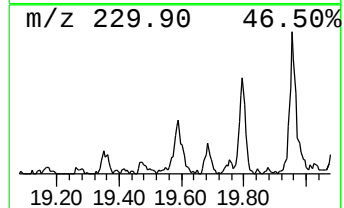
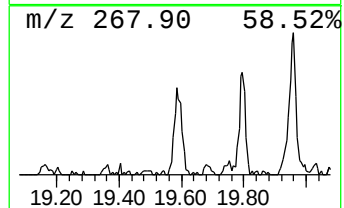
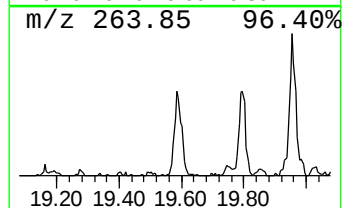
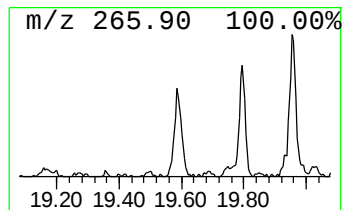
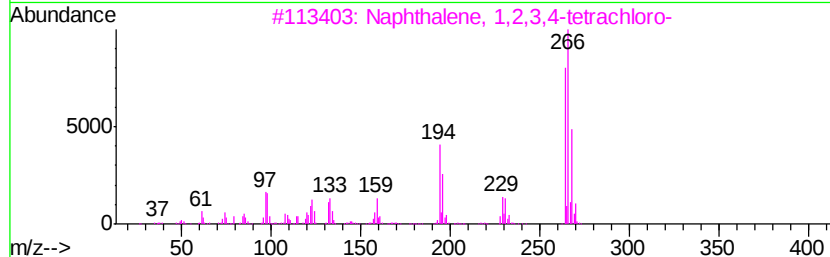
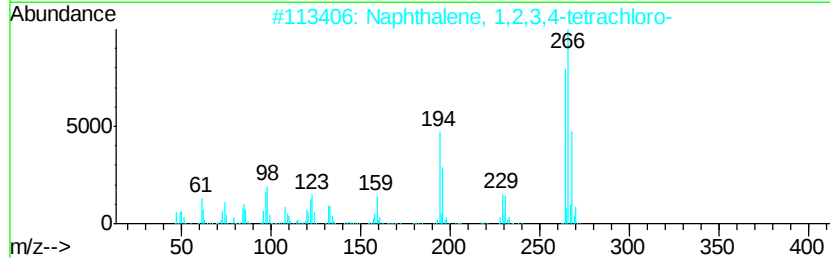
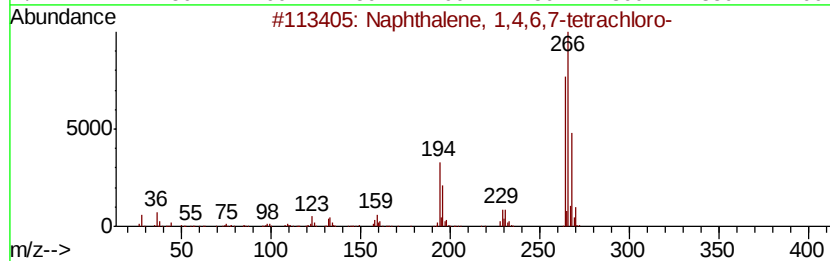
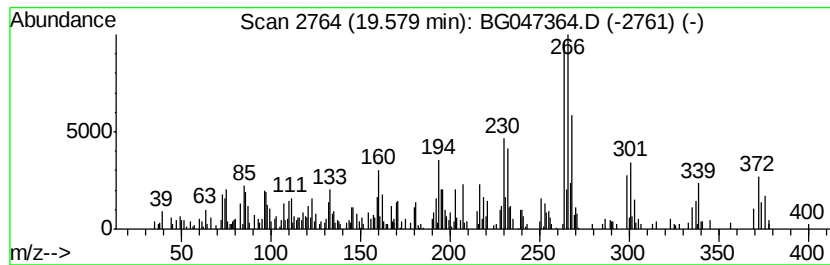
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Naphthalene, 1,4,6,7-tetrac... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.58	7.21 ng/ul	288606	Phenanthrene-d10	17.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,4,6,7-tetrachloro-	264	C10H4Cl4	055720-43-9	90
2		Naphthalene, 1,2,3,4-tetrachloro-	264	C10H4Cl4	020020-02-4	78
3		Naphthalene, 1,2,3,4-tetrachloro-	264	C10H4Cl4	020020-02-4	78
4		Naphthalene, 1,2,3,4-tetrachloro-	264	C10H4Cl4	020020-02-4	78
5		Tetrachloroisophthalonitrile	264	C8Cl4N2	001897-45-6	60



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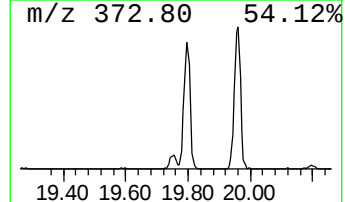
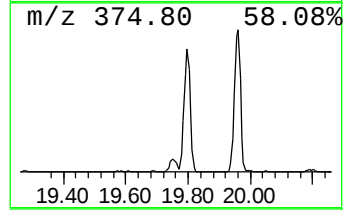
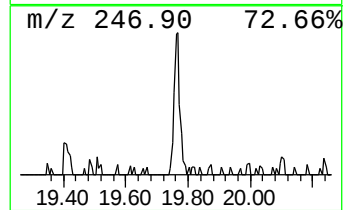
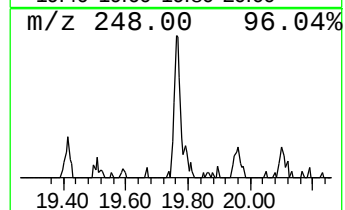
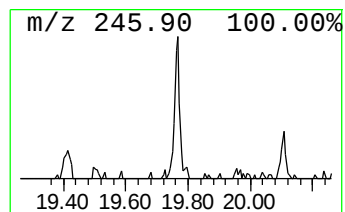
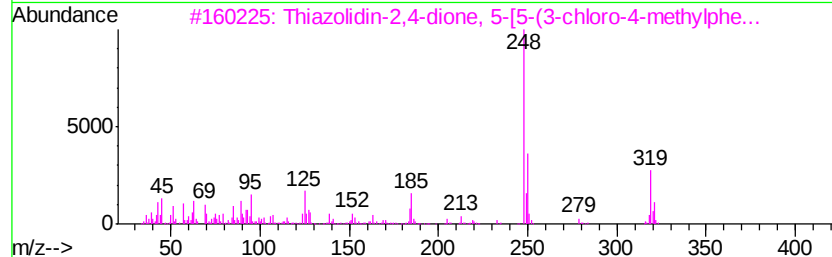
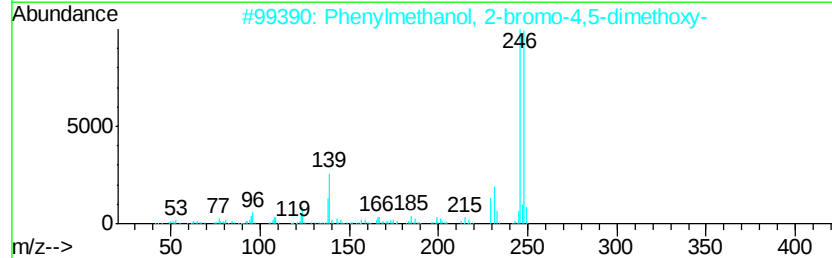
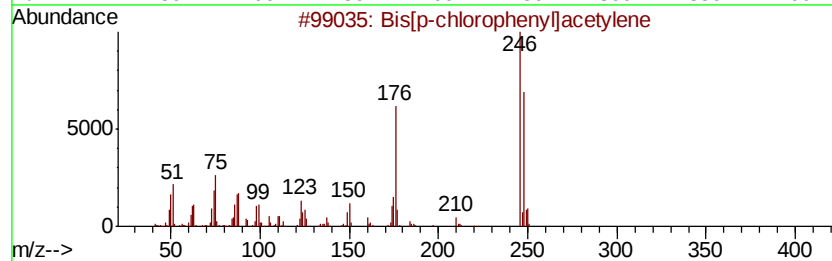
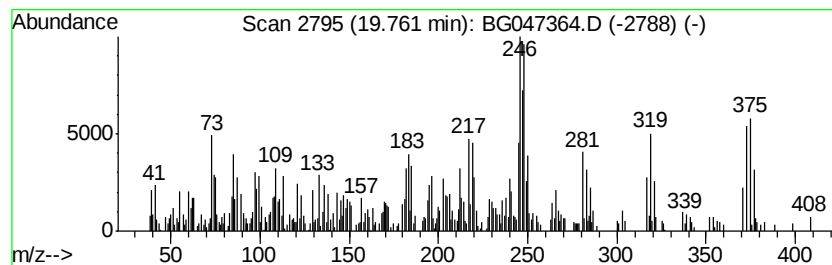
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TIC Integration Parameters: LSCINT.P

Peak Number 8 unknown-04 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.76	7.80 ng/ul	315012	Chrysene-d12	21.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bis[p-chlorophenyl]acetylene	246	C14H8Cl2	001820-42-4	30
2		Phenylmethanol, 2-bromo-4,5-dime...	246	C9H11BrO3	054370-00-2	30
3		Thiazolidin-2,4-dione, 5-[5-(3-c...	319	C15H10ClN03S	1000297-44-2	30
4		Pyrimethamine	248	C12H13ClN4	000058-14-0	25
5		Lumisantonin	246	C15H18O3	000467-41-4	25



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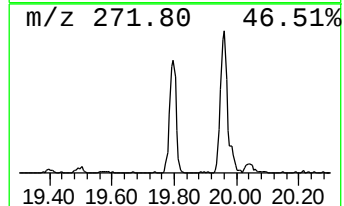
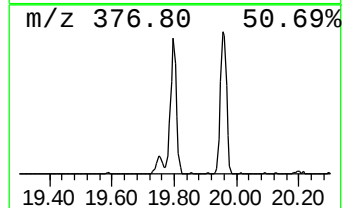
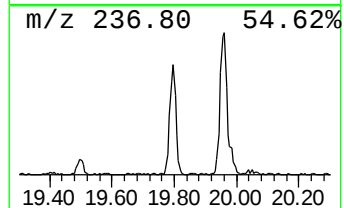
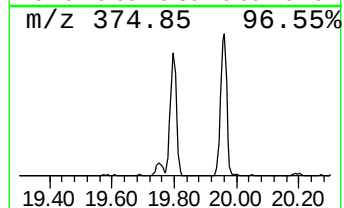
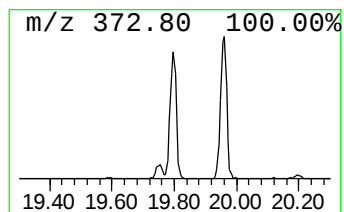
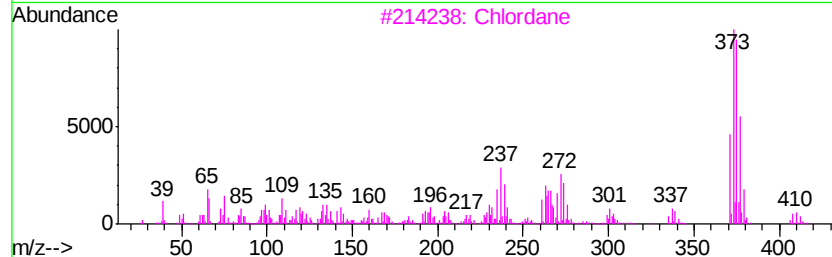
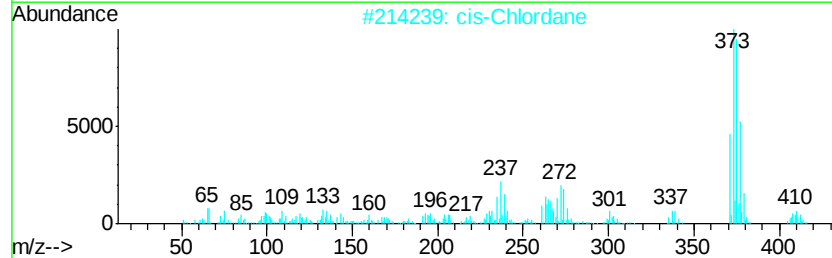
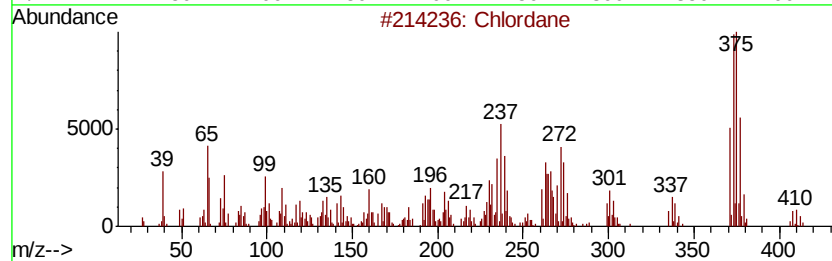
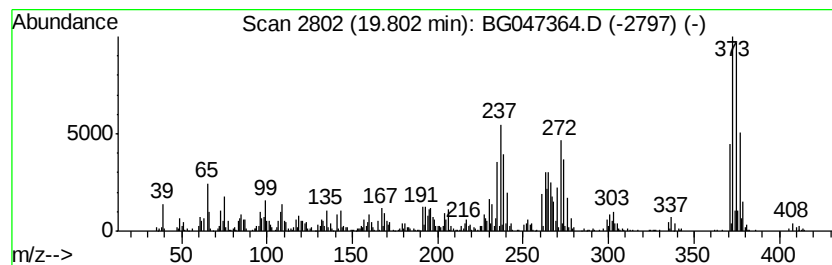
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Chlordane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.80	28.49 ng/ul	1150290	Chrysene-d12	21.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Chlordane	406	C10H6Cl8	000057-74-9	94
2		cis-Chlordane	406	C10H6Cl8	005103-71-9	91
3		Chlordane	406	C10H6Cl8	000057-74-9	91
4		Chlordane	406	C10H6Cl8	000057-74-9	91
5		trans-Chlordane	406	C10H6Cl8	005103-74-2	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111920\
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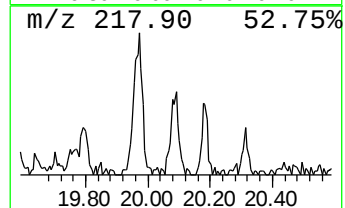
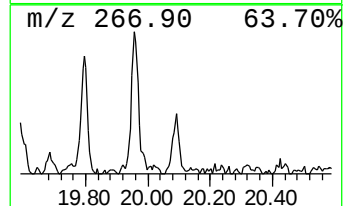
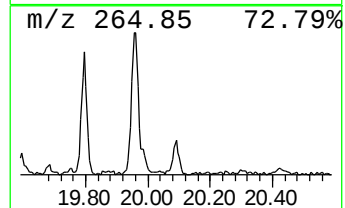
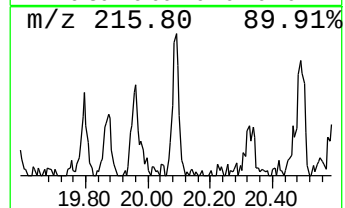
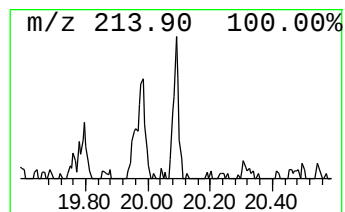
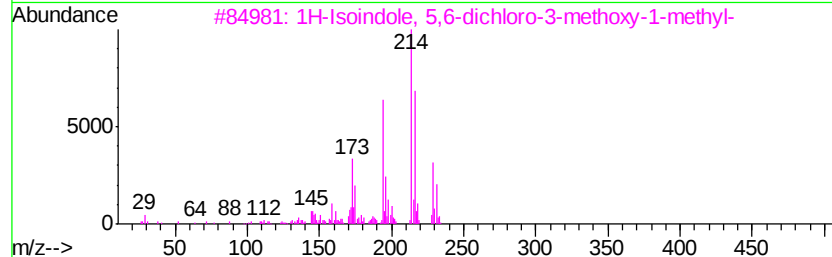
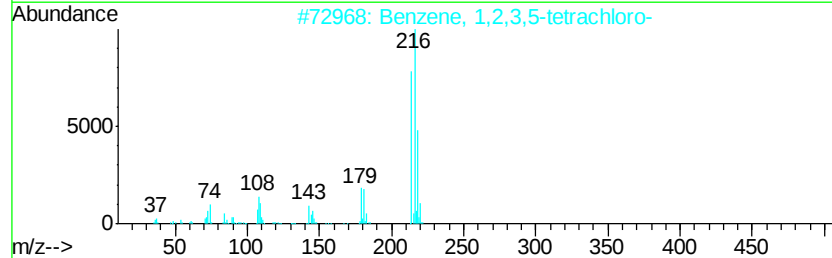
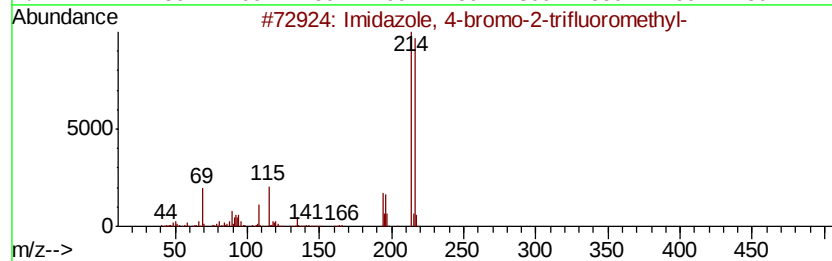
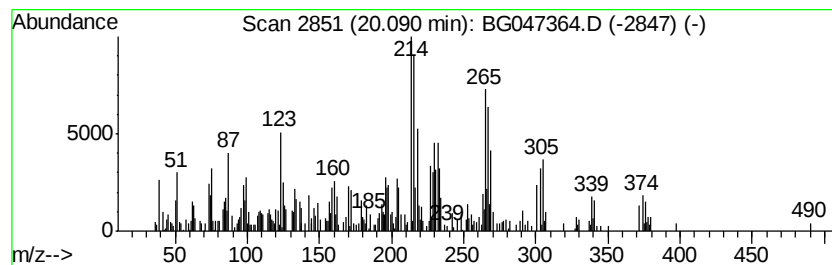
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-05 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.09	4.98 ng/ul	200980	Chrysene-d12	21.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Imidazole, 4-bromo-2-trifluorome...	214	C4H2BrF3N2	219534-98-2	42
2		Benzene, 1,2,3,5-tetrachloro-	214	C6H2Cl4	000634-90-2	35
3		1H-Isoindole, 5,6-dichloro-3-met...	229	C10H9Cl2N0	100813-57-8	30
4		Quinoxaline-2-carbonitrile, 3-am...	216	C10H8N4O2	094728-06-0	25
5		Benzene, 1,2,3,4-tetrachloro-	214	C6H2Cl4	000634-66-2	20



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111920\
Data File : BG047364.D
Acq On : 19 Nov 2020 22:45
Operator : CG/JU
Sample : L4710-20
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AY3

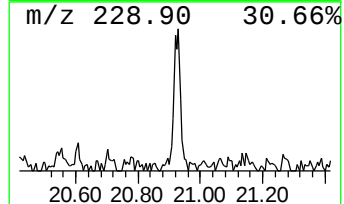
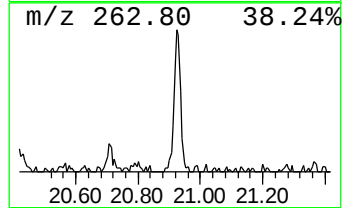
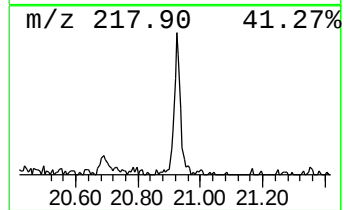
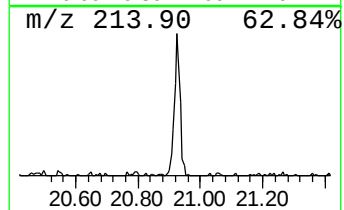
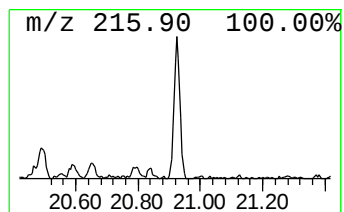
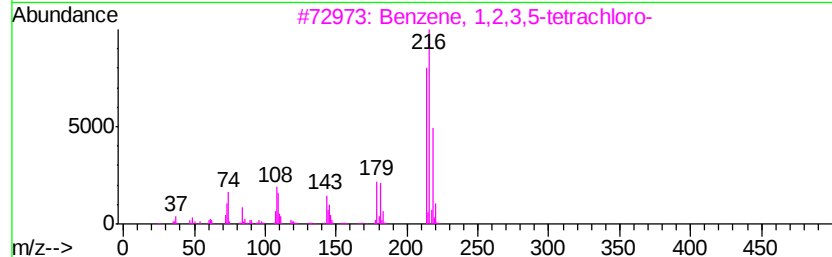
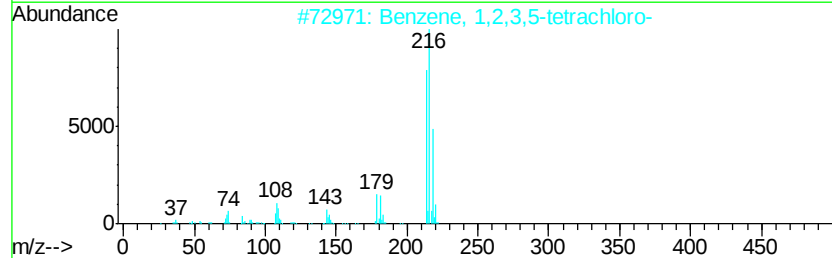
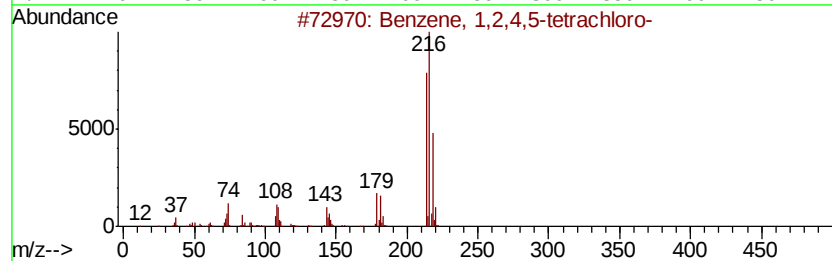
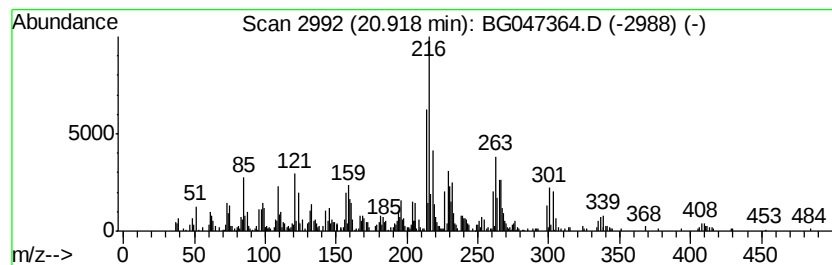
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-06 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.92	11.35 ng/ul	458342	Chrysene-d12	21.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetrachloro-	214	C6H2Cl4	000095-94-3	46
2		Benzene, 1,2,3,5-tetrachloro-	214	C6H2Cl4	000634-90-2	46
3		Benzene, 1,2,3,5-tetrachloro-	214	C6H2Cl4	000634-90-2	41
4		Benzene, 1,2,3,4-tetrachloro-	214	C6H2Cl4	000634-66-2	41
5		Benzene, 1,2,3,4-tetrachloro-	214	C6H2Cl4	000634-66-2	41



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111920\
Data File : BG047364.D
Acq On : 19 Nov 2020 22:45
Operator : CG/JU
Sample : L4710-20
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AY3

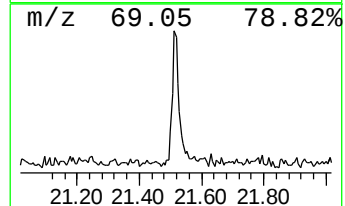
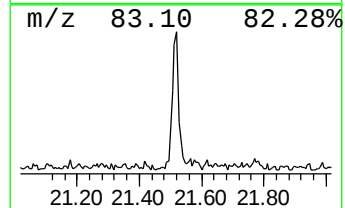
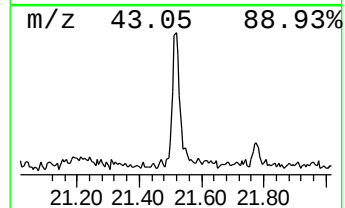
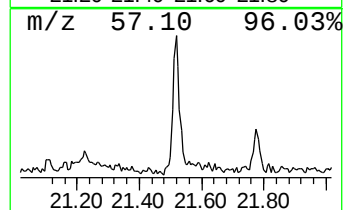
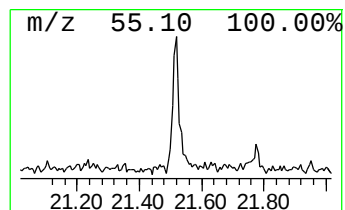
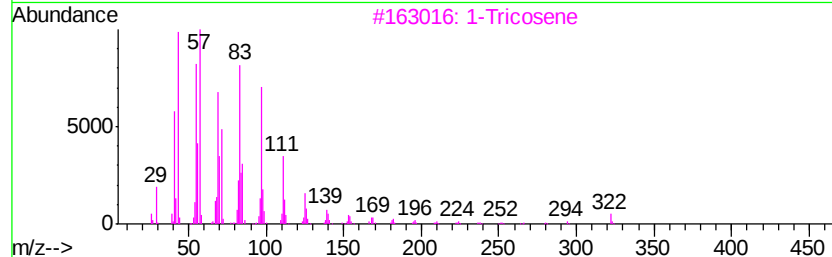
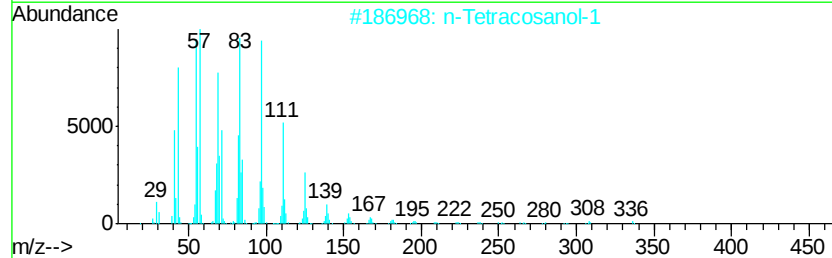
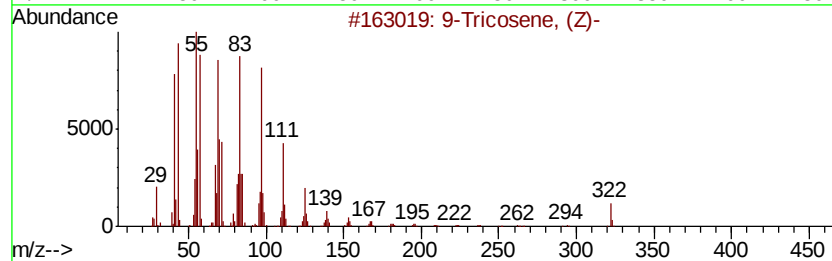
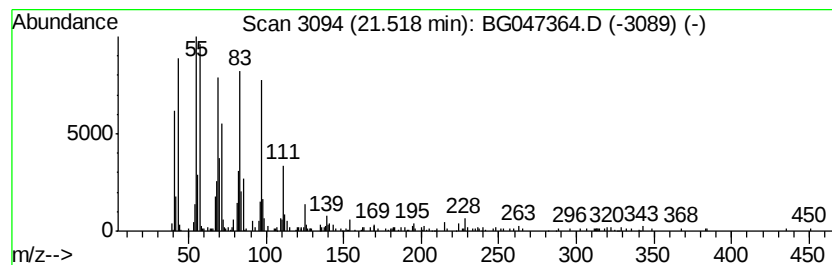
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.52	6.20 ng/ul	250203	Chrysene-d12	21.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Tricosene, (Z)-	322	C23H46	027519-02-4	93
2		n-Tetracosanol-1	354	C24H50O	000506-51-4	90
3		1-Tricosene	322	C23H46	018835-32-0	90
4		1-Heneicosanol	312	C21H44O	015594-90-8	87
5		Dichloroacetic acid, tridecyl ester	310	C15H28Cl2O2	1000280-48-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG111920\
 Data File : BG047364.D
 Acq On : 19 Nov 2020 22:45
 Operator : CG/JU
 Sample : L4710-20
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AY3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.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
Chlordene, .alpha...	18.49	7.8	ng/ul	313861	4	17.62	800363	20.0
1-Perchloroetheny...	18.88	56.2	ng/ul	2247610	4	17.62	800363	20.0
unknown-01	19.01	4.2	ng/ul	166524	4	17.62	800363	20.0
(DEL) Alkane: Str...	19.26	3.9	ng/ul	155399	4	17.62	800363	20.0
unknown-02	19.36	3.8	ng/ul	150251	4	17.62	800363	20.0
unknown-03	19.50	5.4	ng/ul	214950	4	17.62	800363	20.0
Naphthalene, 1,4,...	19.58	7.2	ng/ul	288606	4	17.62	800363	20.0
unknown-04	19.76	7.8	ng/ul	315012	5	21.90	807366	20.0
Chlordane	19.80	28.5	ng/ul	1150290	5	21.90	807366	20.0
unknown-05	20.09	5.0	ng/ul	200980	5	21.90	807366	20.0
unknown-06	20.92	11.4	ng/ul	458342	5	21.90	807366	20.0
(DEL) Alkane: Str...	21.52	6.2	ng/ul	250203	5	21.90	807366	20.0