

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
 Data File : BM029421.D
 Acq On : 08 Apr 2021 23:41
 Operator : CG/JU
 Sample : M1959-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B41

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_M\METHODS\SOM-EPA-BM040721MA.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	2.946	8	10	17	rVB	132321	145256	5.65%	0.697%
2	3.028	20	24	26	rVV	39904	41691	1.62%	0.200%
3	3.058	26	29	30	rVV	17215	16722	0.65%	0.080%
4	3.081	30	33	35	rVV	25856	25745	1.00%	0.124%
5	3.117	35	39	44	rVB	2645939	2570089	100.00%	12.337%
6	3.423	87	91	92	rBV2	6112	6943	0.27%	0.033%
7	3.446	92	95	101	rVB	76507	90856	3.54%	0.436%
8	3.540	108	111	113	rBV4	4539	4360	0.17%	0.021%
9	3.823	156	159	166	rVB5	4542	6314	0.25%	0.030%
10	3.917	170	175	181	rBV	22076	27837	1.08%	0.134%
11	4.093	201	205	208	rVB4	3797	4425	0.17%	0.021%
12	4.123	208	210	216	rBV5	2316	3661	0.14%	0.018%
13	4.458	263	267	270	rVB3	5281	6534	0.25%	0.031%
14	4.746	313	316	322	rVB6	2737	5408	0.21%	0.026%
15	4.822	324	329	330	rVV5	2430	3684	0.14%	0.018%
16	4.864	333	336	339	rVB4	3421	3949	0.15%	0.019%
17	5.717	476	481	485	rBV5	2944	5208	0.20%	0.025%
18	5.975	522	525	530	rVB4	2424	3631	0.14%	0.017%
19	6.193	557	562	568	rVB7	2861	4228	0.16%	0.020%
20	6.258	568	573	577	rBV6	2084	3465	0.13%	0.017%
21	6.916	682	685	692	rVB7	1767	2679	0.10%	0.013%
22	7.322	747	754	755	rBV4	3858	5528	0.22%	0.027%
23	7.687	809	816	823	rBV	553284	856836	33.34%	4.113%
24	8.322	921	924	929	rBV4	2027	3654	0.14%	0.018%
25	8.916	1021	1025	1029	rVB4	2416	3801	0.15%	0.018%
26	10.334	1261	1266	1271	rBV4	7724	13411	0.52%	0.064%
27	10.469	1280	1289	1296	rBV	678911	1152802	44.85%	5.534%
28	11.910	1528	1534	1538	rBV5	2270	3642	0.14%	0.017%
29	12.140	1566	1573	1576	rBV4	2301	4064	0.16%	0.020%
30	12.357	1606	1610	1615	rBV3	2890	5236	0.20%	0.025%
31	13.157	1743	1746	1752	rVB4	5015	7139	0.28%	0.034%
32	13.640	1824	1828	1829	rBV2	2786	2580	0.10%	0.012%
33	13.692	1834	1837	1841	rVB5	2525	3920	0.15%	0.019%
34	14.151	1913	1915	1921	rVB4	2498	3570	0.14%	0.017%

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 Title : SVOA CALIBRATION

35	14.240	1925	1930	1936	rBV6	4379	8267	0.32%	0.040%
36	14.322	1936	1944	1954	rBV	1034468	1487599	57.88%	7.141%
37	14.722	2007	2012	2016	rVB3	4720	6783	0.26%	0.033%
38	14.857	2031	2035	2040	rBV7	2325	4099	0.16%	0.020%
39	15.122	2077	2080	2082	rBV4	2502	3056	0.12%	0.015%
40	15.192	2088	2092	2096	rBV4	5238	8328	0.32%	0.040%
41	15.375	2119	2123	2127	rBV3	8833	10516	0.41%	0.050%
42	15.592	2154	2160	2163	rBV7	3498	7114	0.28%	0.034%
43	15.786	2191	2193	2195	rBV3	2440	2677	0.10%	0.013%
44	15.810	2195	2197	2200	rVV4	3878	4845	0.19%	0.023%
45	16.051	2235	2238	2240	rBV4	6583	8836	0.34%	0.042%
46	16.275	2274	2276	2278	rVB3	4202	3274	0.13%	0.016%
47	16.416	2298	2300	2305	rBV6	5050	7782	0.30%	0.037%
48	16.598	2327	2331	2336	rVB2	38632	52034	2.02%	0.250%
49	17.063	2404	2410	2414	rBV	1163366	1639024	63.77%	7.868%
50	17.104	2414	2417	2423	rVB	152618	210725	8.20%	1.012%
51	17.245	2436	2441	2448	rBV	105683	148858	5.79%	0.715%
52	17.669	2506	2513	2520	rBV	392757	544884	21.20%	2.616%
53	17.780	2528	2532	2540	rVB3	60015	89594	3.49%	0.430%
54	17.916	2550	2555	2560	rBV2	140432	218951	8.52%	1.051%
55	17.969	2561	2564	2574	rVB5	72428	129728	5.05%	0.623%
56	18.133	2587	2592	2596	rBV	257329	351526	13.68%	1.687%
57	18.192	2598	2602	2606	rVV	176097	242756	9.45%	1.165%
58	18.239	2606	2610	2615	rVB	104052	140398	5.46%	0.674%
59	18.422	2636	2641	2646	rBV2	226215	330694	12.87%	1.587%
60	18.469	2646	2649	2654	rVB	91284	116665	4.54%	0.560%
61	18.563	2661	2665	2667	rBV	67624	86214	3.35%	0.414%
62	18.598	2667	2671	2678	rVB	181722	231791	9.02%	1.113%
63	18.698	2685	2688	2691	rVB4	46582	57391	2.23%	0.275%
64	18.904	2719	2723	2728	rBV	151373	205984	8.01%	0.989%
65	18.957	2728	2732	2733	rVV3	69046	82521	3.21%	0.396%
66	18.992	2733	2738	2744	rVB3	372018	596159	23.20%	2.862%
67	19.122	2755	2760	2764	rVB	129491	182268	7.09%	0.875%
68	19.204	2769	2774	2781	rVV3	145922	334698	13.02%	1.607%
69	19.269	2781	2785	2792	rVV	278429	382413	14.88%	1.836%
70	19.333	2793	2796	2799	rVB4	44911	53058	2.06%	0.255%
71	19.527	2827	2829	2833	rVB3	42711	49912	1.94%	0.240%

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72	19.604	2839	2842	2847	rVB3	69113	89756	3.49%	0.431%
73	19.716	2854	2861	2865	rVB2	182883	310001	12.06%	1.488%
74	19.839	2878	2882	2886	rBV2	154656	200660	7.81%	0.963%
75	19.880	2887	2889	2896	rVB7	59992	93313	3.63%	0.448%
76	19.974	2901	2905	2910	rVV	457733	583507	22.70%	2.801%
77	20.027	2910	2914	2921	rVB	110868	136677	5.32%	0.656%
78	20.180	2937	2940	2944	rVB4	65643	74124	2.88%	0.356%
79	20.257	2948	2953	2957	rBV	583837	705837	27.46%	3.388%
80	20.310	2958	2962	2965	rBV2	153879	216737	8.43%	1.040%
81	20.410	2975	2979	2987	rVB3	170772	277818	10.81%	1.334%
82	20.569	2999	3006	3012	rBV	386644	673363	26.20%	3.232%
83	20.716	3027	3031	3036	rVB2	234222	273808	10.65%	1.314%
84	20.780	3038	3042	3047	rVB2	129209	149811	5.83%	0.719%
85	20.986	3073	3077	3081	rVB3	202185	258923	10.07%	1.243%
86	21.039	3083	3086	3092	rVB4	120479	148211	5.77%	0.711%
87	21.092	3093	3095	3098	rBV3	50867	62226	2.42%	0.299%
88	21.239	3115	3120	3123	rBV	1028743	1354115	52.69%	6.500%
89	21.268	3123	3125	3131	rVB2	568700	616732	24.00%	2.961%
90	21.574	3174	3177	3183	rBV2	208232	268406	10.44%	1.288%
91	23.445	3490	3495	3507	rVB2	651491	1247597	48.54%	5.989%

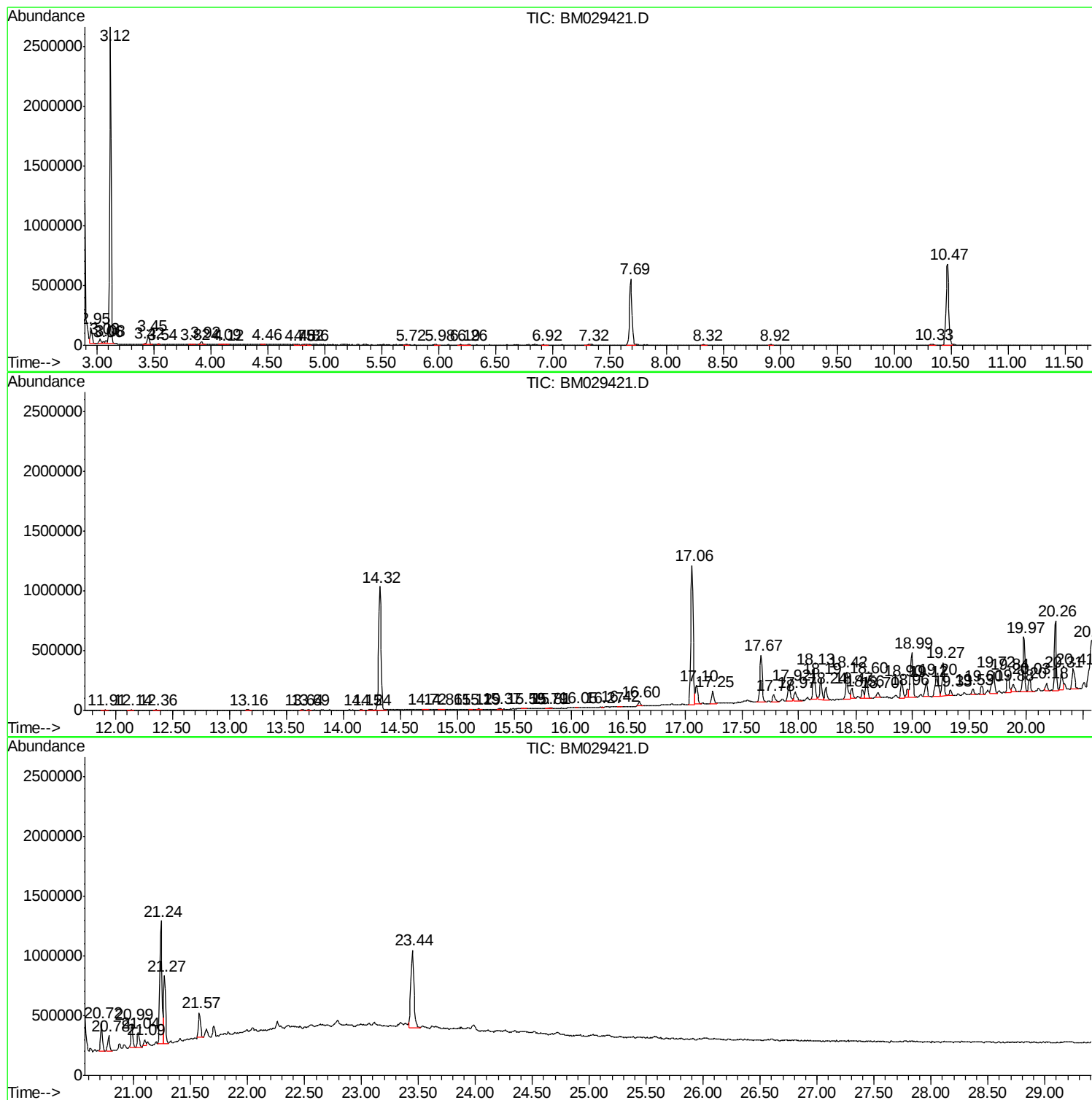
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM040721MA.M
Quant Title : SVOA CALIBRATION

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



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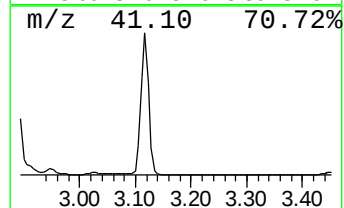
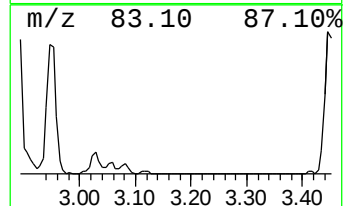
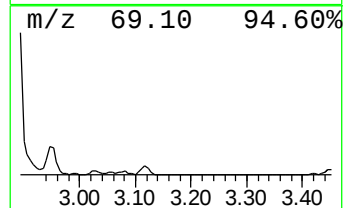
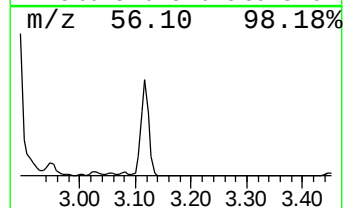
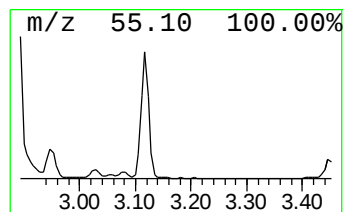
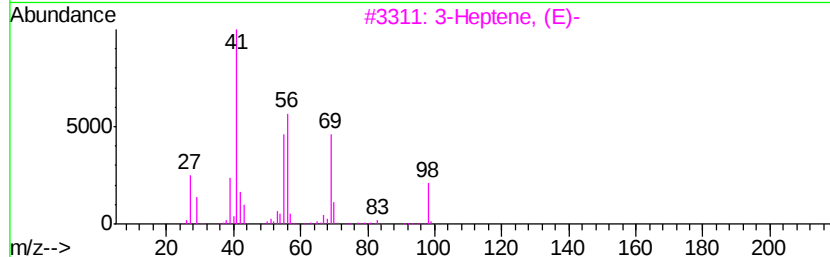
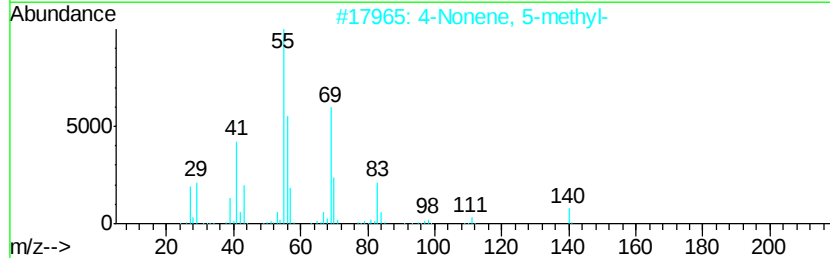
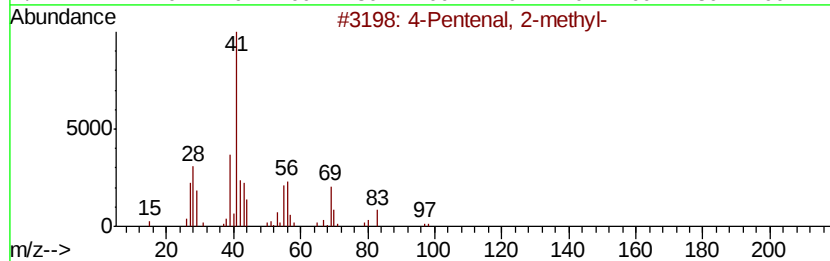
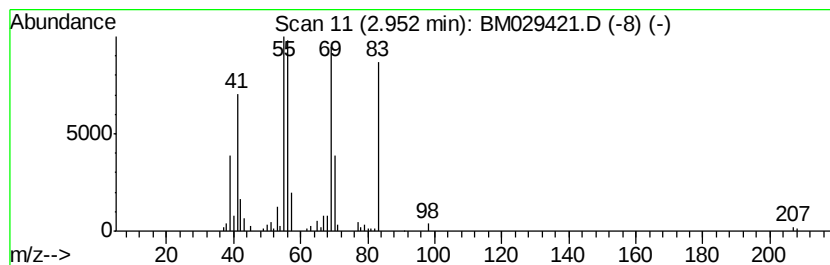
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM040721MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 4-Pentenal, 2-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.95	3.39 ng/ul	145256	1,4-Dichlorobenzene-d4	7.69

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Pentenal, 2-methyl-	98	C6H10O	005187-71-3	53
2			4-Nonene, 5-methyl-	140	C10H20	015918-07-7	50
3			3-Heptene, (E)-	98	C7H14	014686-14-7	50
4			Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	49
5			6-Dodecene, (E)-	168	C12H24	007206-17-9	47



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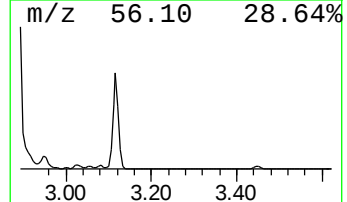
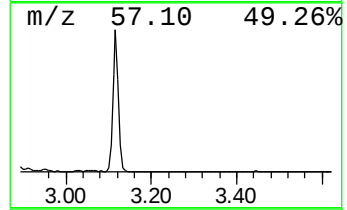
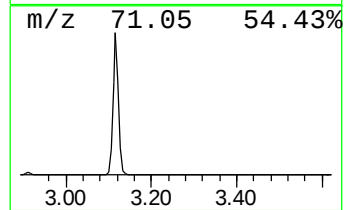
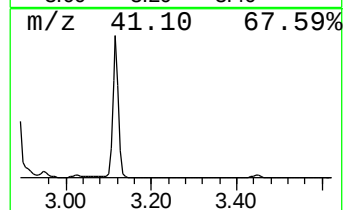
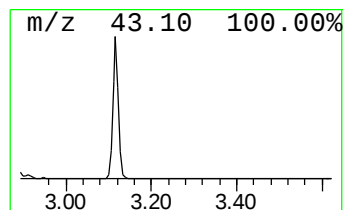
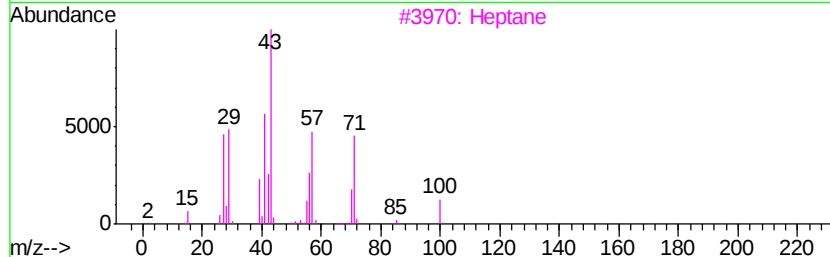
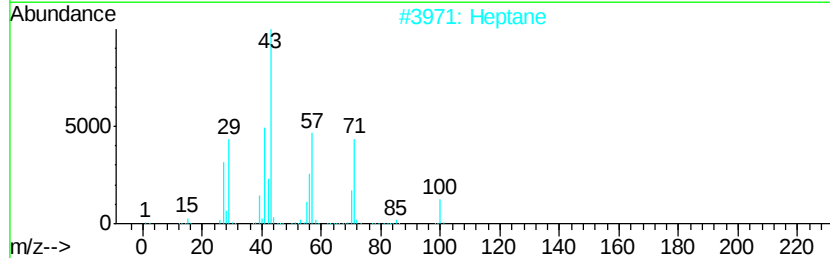
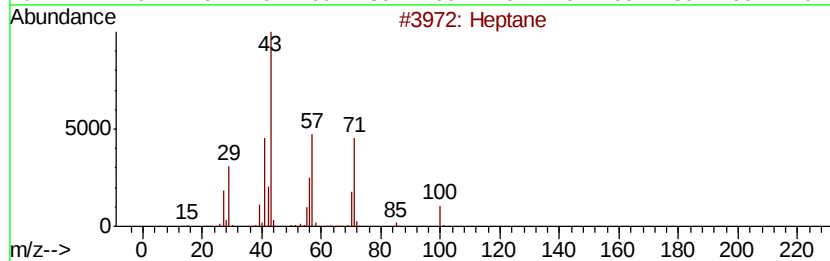
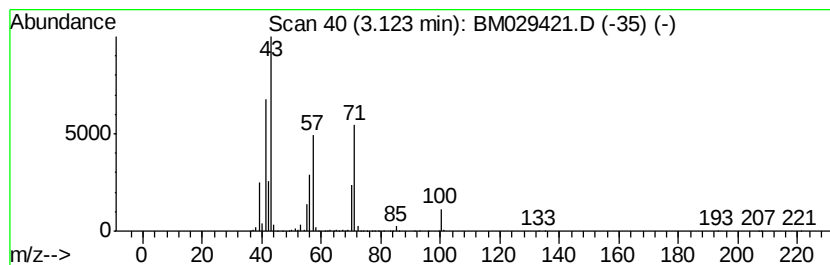
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.12	59.99 ng/ul	2570090	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane	100	C7H16	000142-82-5	94
2		Heptane	100	C7H16	000142-82-5	90
3		Heptane	100	C7H16	000142-82-5	87
4		Heptane	100	C7H16	000142-82-5	80
5		Hexane, 3-methyl-	100	C7H16	000589-34-4	45



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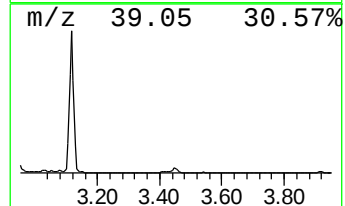
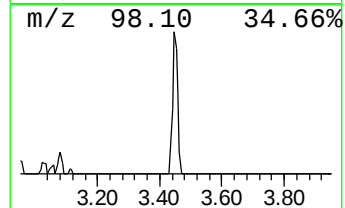
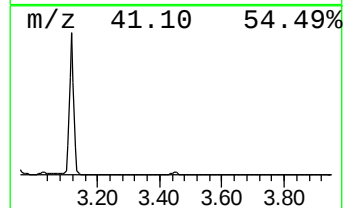
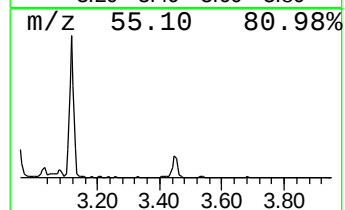
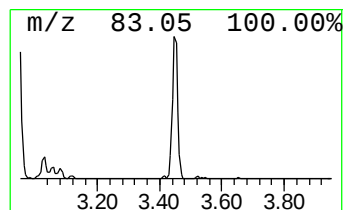
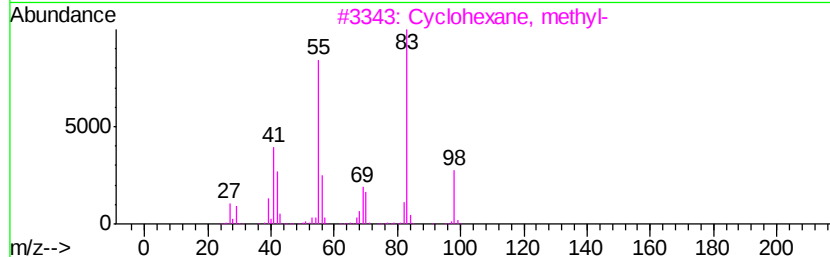
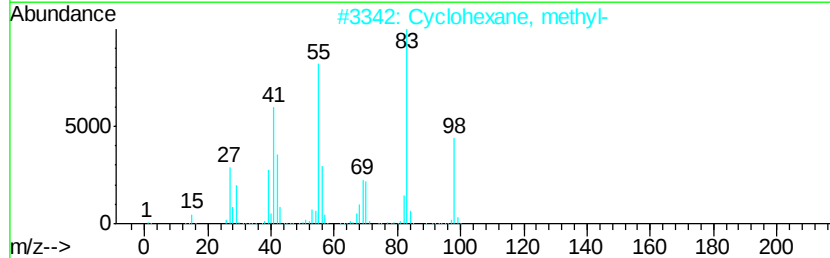
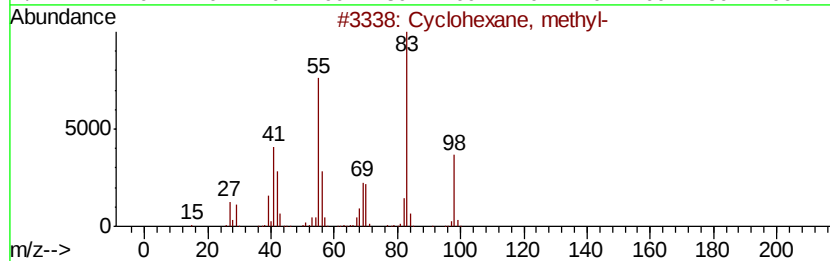
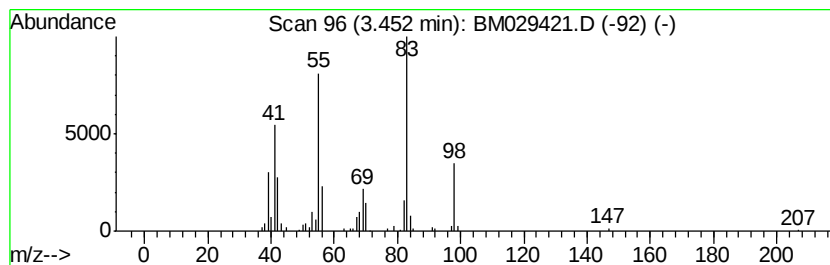
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Peak Number 3 (DEL) Alkane: Cyclic3.45 Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.45	2.12 ng/ul	90856	1,4-Dichlorobenzene-d4	7.69

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, methyl-	98	C7H14	000108-87-2	97
2			Cyclohexane, methyl-	98	C7H14	000108-87-2	97
3			Cyclohexane, methyl-	98	C7H14	000108-87-2	94
4			Cyclohexane, methyl-	98	C7H14	000108-87-2	90
5			2-Pentene, 2,3-dimethyl-	98	C7H14	010574-37-5	58



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C0B41

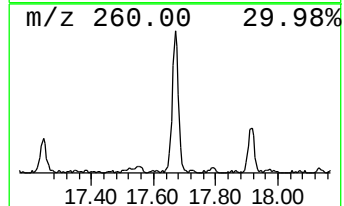
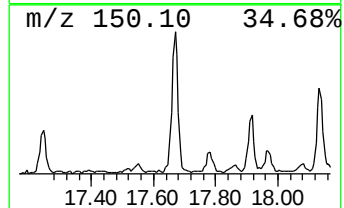
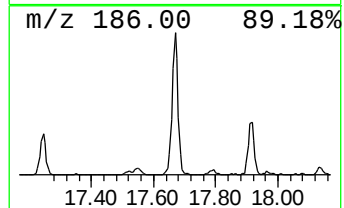
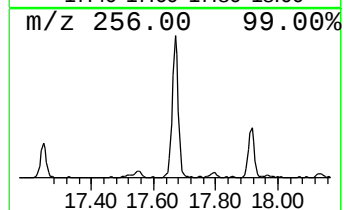
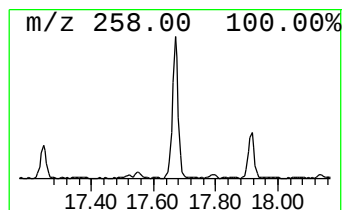
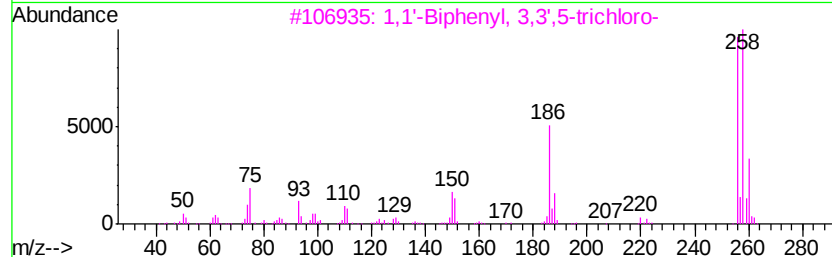
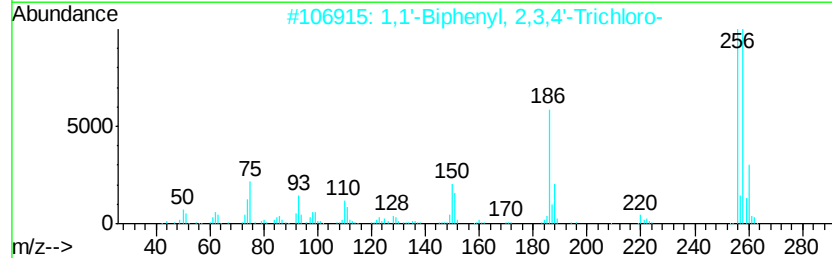
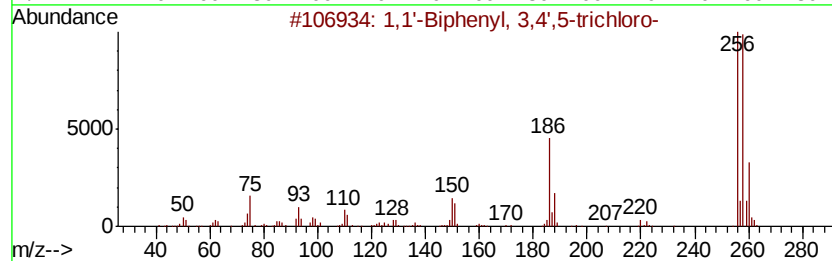
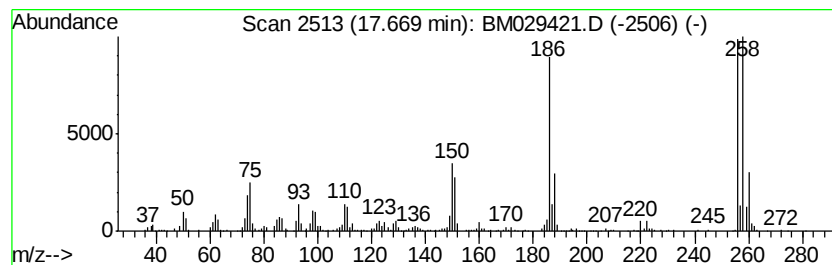
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.67	6.65 ng/ul	544884	Phenanthrene-d10	17.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1'-Biphenyl, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	99
2			1,1'-Biphenyl, 2,3,4'-Trichloro-	256	C12H7Cl3	038444-85-8	99
3			1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	99
4			1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99
5			1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
Data File : BM029421.D
Acq On : 08 Apr 2021 23:41
Operator : CG/JU
Sample : M1959-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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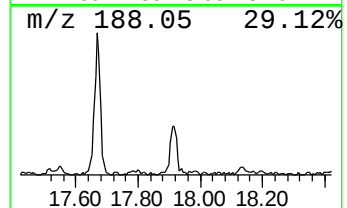
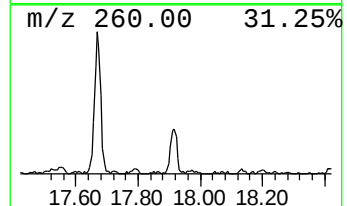
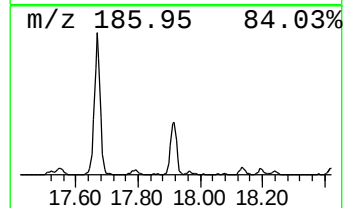
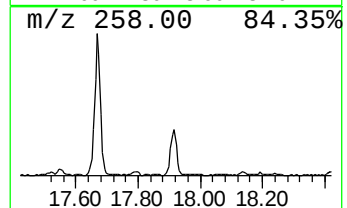
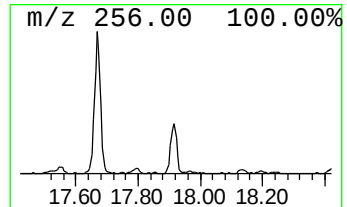
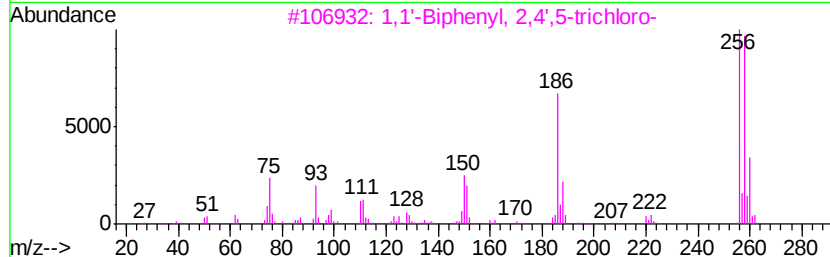
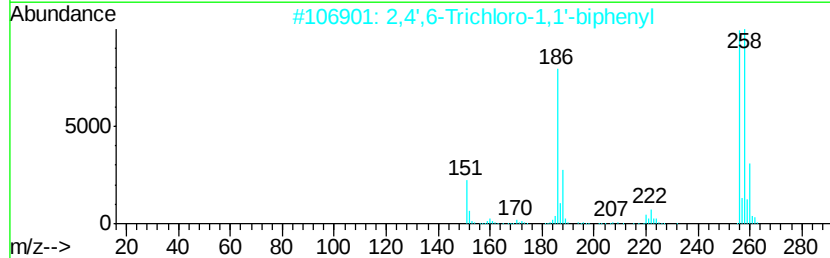
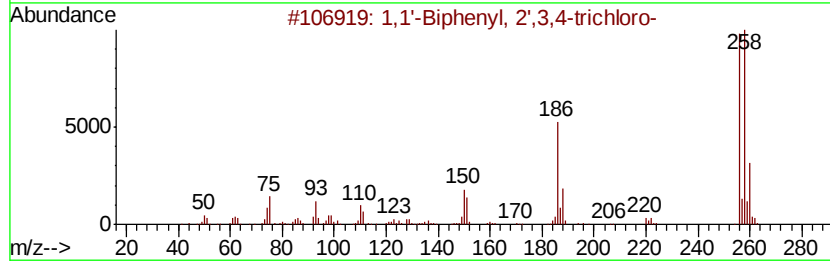
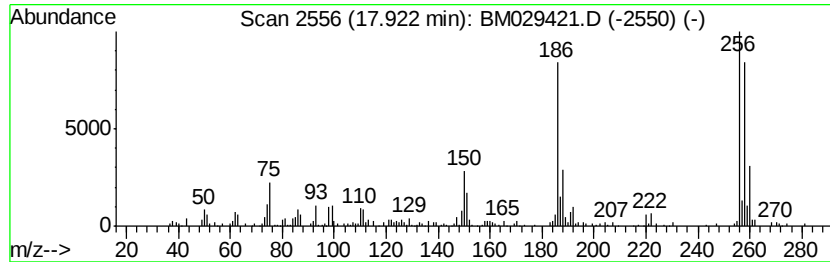
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,1'-Biphenyl, 2',3,4-trich... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.92	2.67 ng/ul	218951	Phenanthrene-d10	17.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99
2		2,4',6-Trichloro-1,1'-biphenyl	256	C12H7Cl3	038444-77-8	99
3		1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	99
4		1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	99
5		1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
Data File : BM029421.D
Acq On : 08 Apr 2021 23:41
Operator : CG/JU
Sample : M1959-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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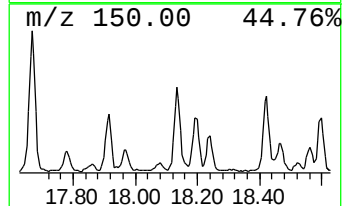
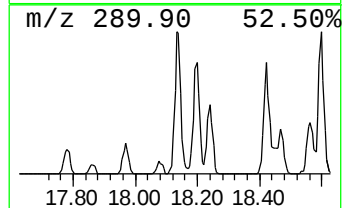
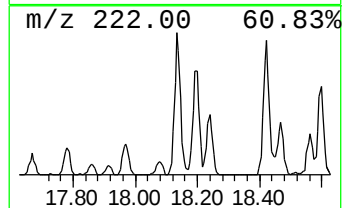
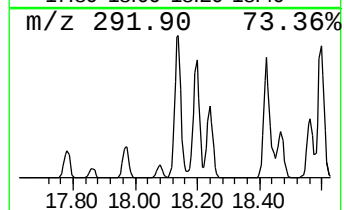
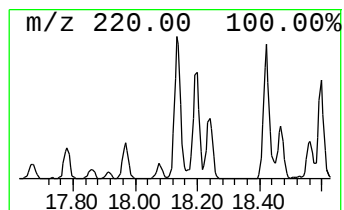
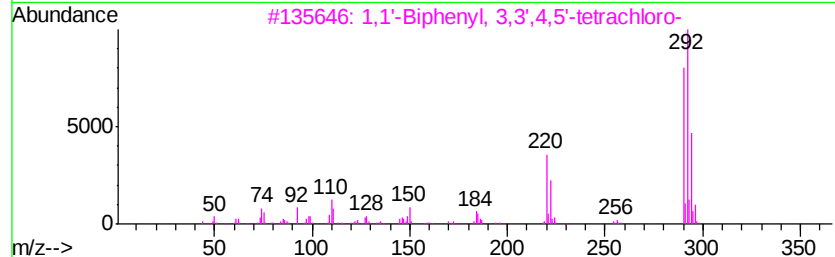
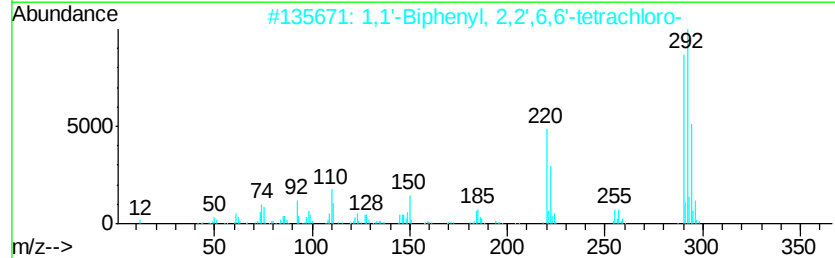
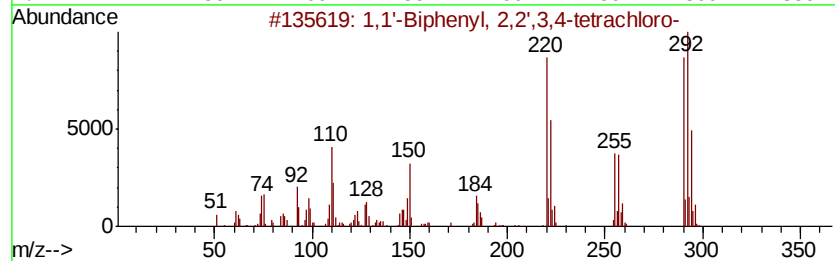
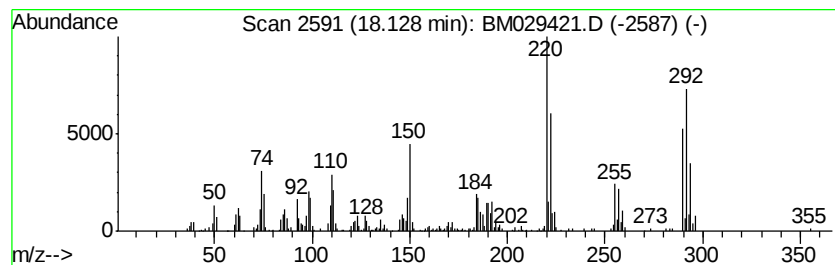
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1,1'-Biphenyl, 2,2',3,4-tet... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.13	4.29 ng/ul	351526	Phenanthrene-d10	17.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',3,4-tetrachl...	290	C12H6Cl4	052663-59-9	99
2		1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99
3		1,1'-Biphenyl, 3,3',4,5'-tetrach...	290	C12H6Cl4	041464-48-6	99
4		2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99
5		2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
Data File : BM029421.D
Acq On : 08 Apr 2021 23:41
Operator : CG/JU
Sample : M1959-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

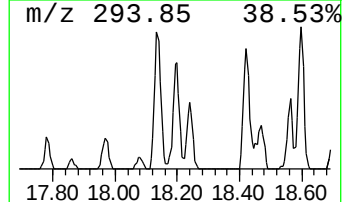
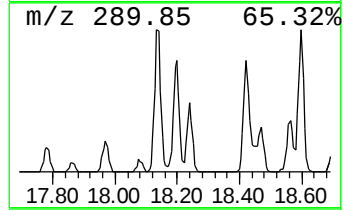
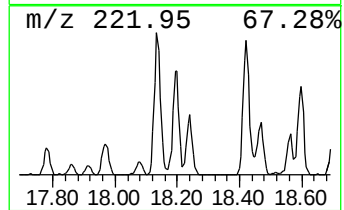
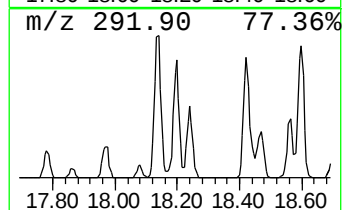
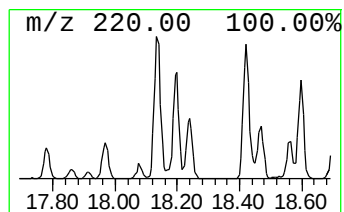
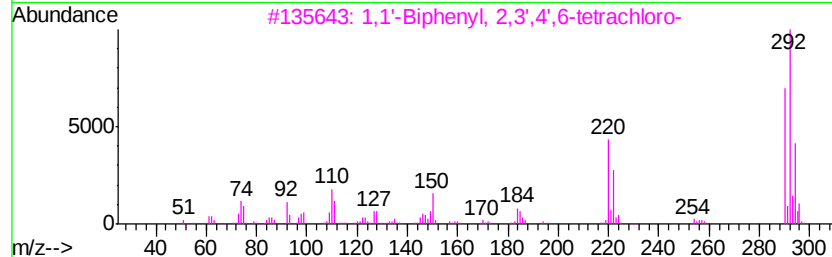
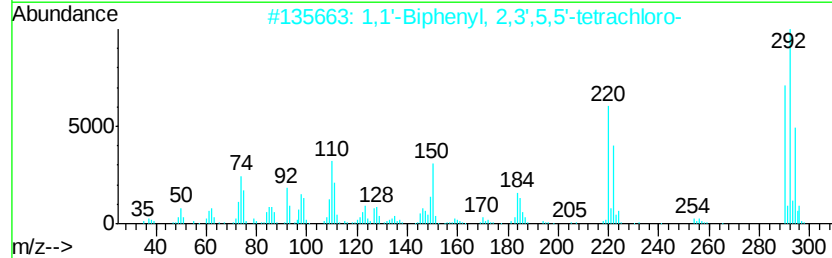
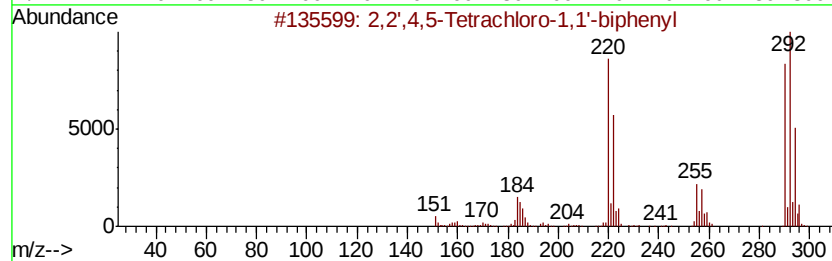
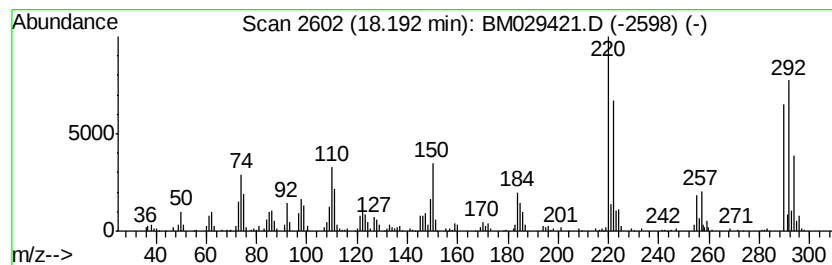
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 2,2',4,5-Tetrachloro-1,1'-b... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.19	2.96 ng/ul	242756	Phenanthrene-d10	17.06		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99	
2	1,1'-Biphenyl, 2,3',5,5'-tetrachloro-	290	C12H6Cl4	041464-42-0	99	
3	1,1'-Biphenyl, 2,3',4',6-tetrachloro-	290	C12H6Cl4	041464-46-4	99	
4	1,1'-Biphenyl, 2,2',4,4'-tetrachloro-	290	C12H6Cl4	002437-79-8	99	
5	1,1'-Biphenyl, 2,2',4,4'-tetrachloro-	290	C12H6Cl4	002437-79-8	99	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
Data File : BM029421.D
Acq On : 08 Apr 2021 23:41
Operator : CG/JU
Sample : M1959-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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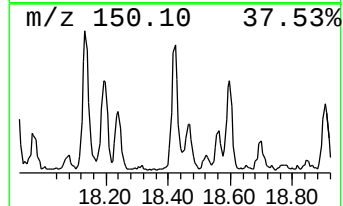
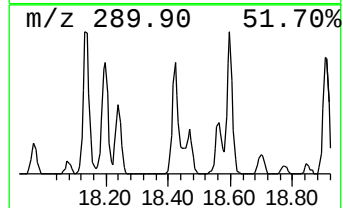
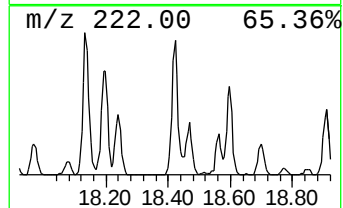
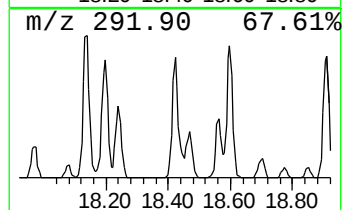
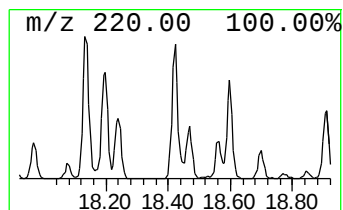
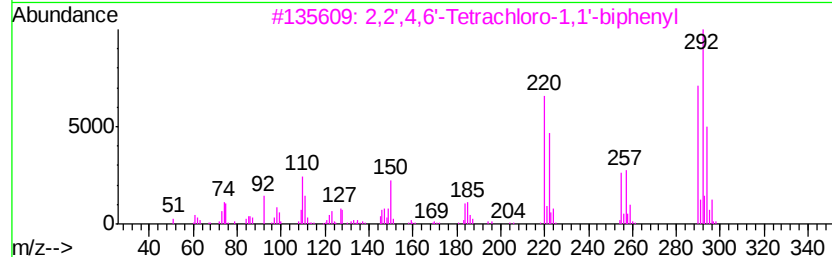
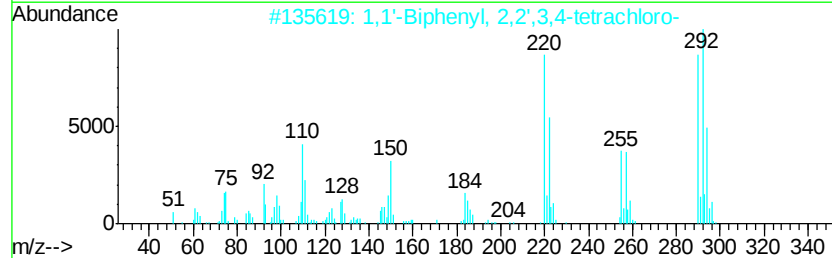
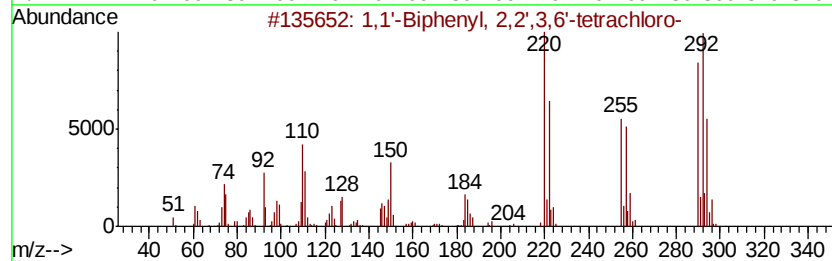
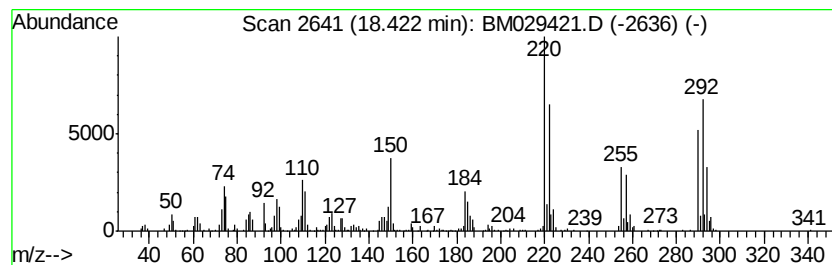
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1,1'-Biphenyl, 2,2',3,6'-te... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.42	4.04 ng/ul	330694	Phenanthrene-d10	17.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',3,6'-tetrach...	290	C12H6Cl4	041464-47-5	99
2		1,1'-Biphenyl, 2,2',3,4'-tetrachl...	290	C12H6Cl4	052663-59-9	99
3		2,2',4,6'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	068194-04-7	99
4		2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99
5		1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
Data File : BM029421.D
Acq On : 08 Apr 2021 23:41
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Sample : M1959-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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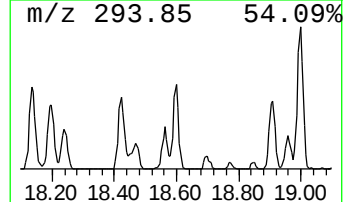
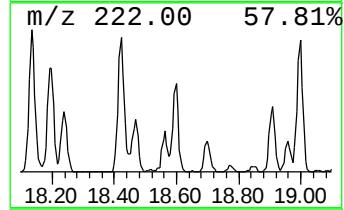
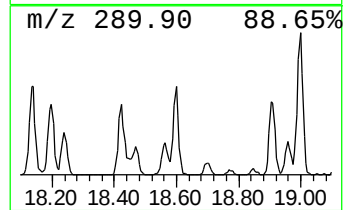
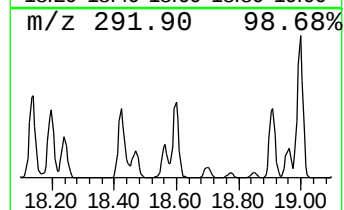
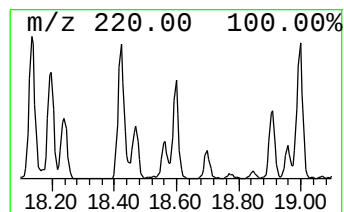
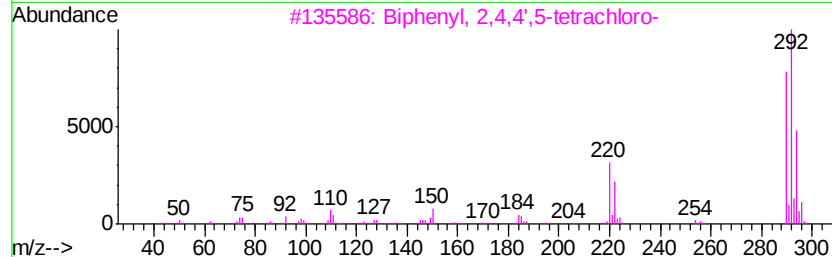
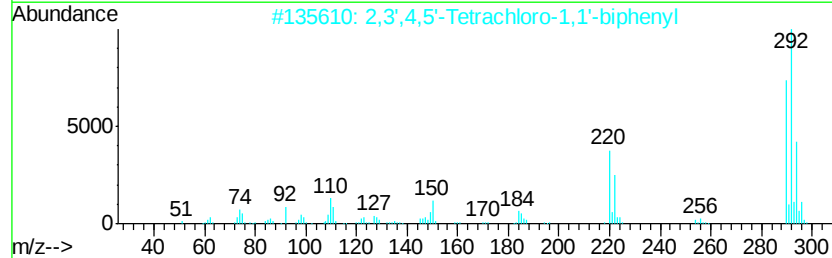
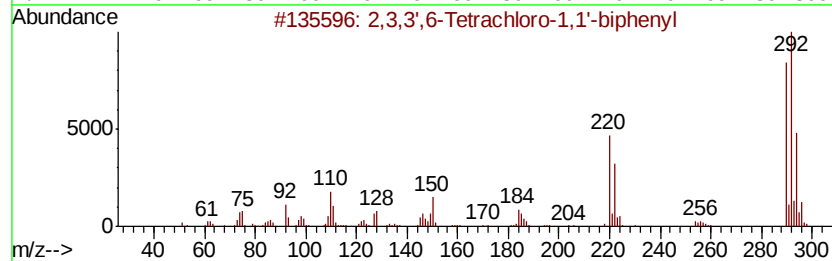
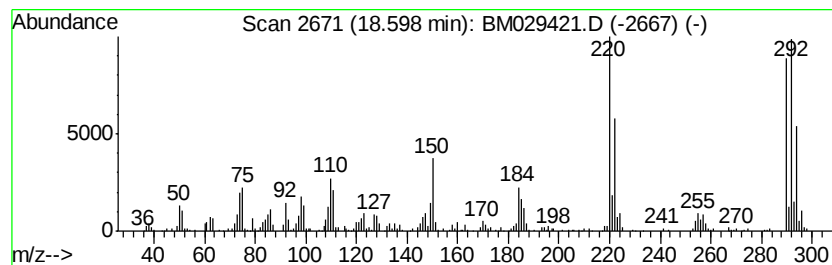
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 2,3,3',6-Tetrachloro-1,1'-b... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.60	2.83 ng/ul	231791	Phenanthrene-d10	17.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,3',6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074472-33-6	99		
2	2,3',4,5'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-52-7	99		
3	Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	99		
4	2,3,3',4-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074338-24-2	99		
5	1,1'-Biphenyl, 2,3,3',5'-tetrach...	290	C12H6Cl4	041464-49-7	99		



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
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Sample : M1959-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

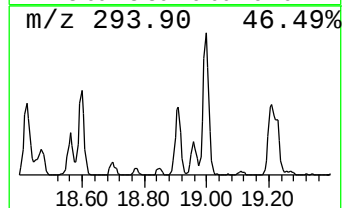
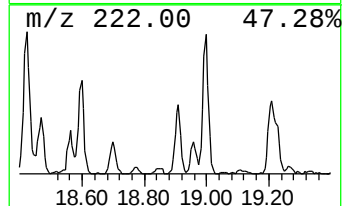
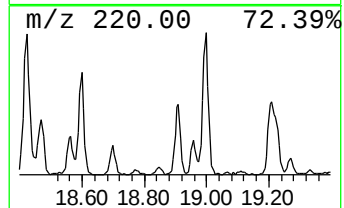
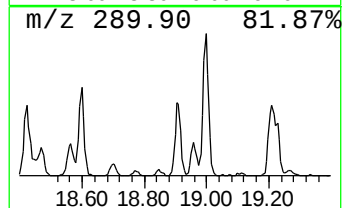
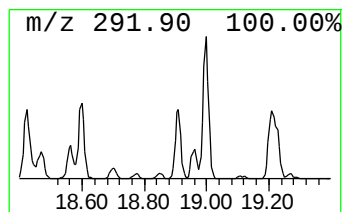
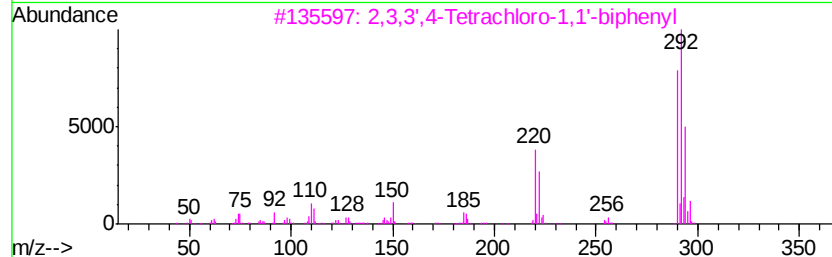
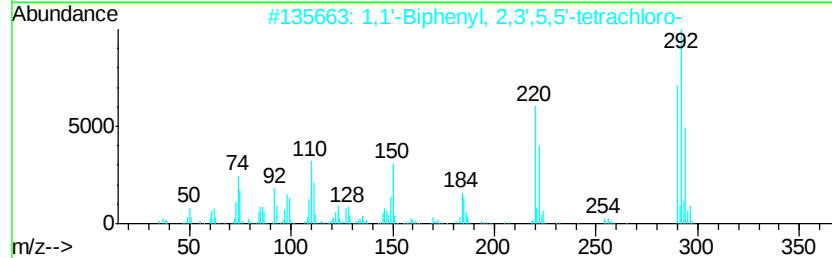
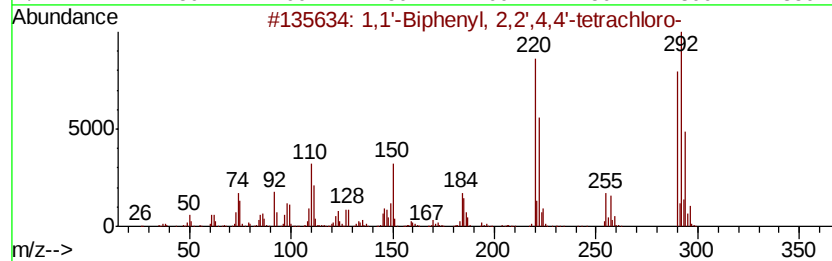
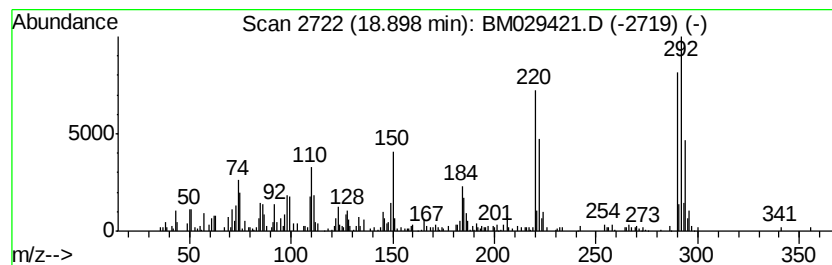
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ClientSampled :
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 1,1'-Biphenyl, 2,2',4,4'-te... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.90	2.51 ng/ul	205984	Phenanthrene-d10	17.06		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'	-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
2	1,1'	-Biphenyl, 2,3',5,5'-tetrach...	290	C12H6Cl4	041464-42-0	99
3	2,3,3',4-	Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074338-24-2	99
4	2,3',5',6-	Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074338-23-1	99
5	2,3,4',5-	Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074472-34-7	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
Data File : BM029421.D
Acq On : 08 Apr 2021 23:41
Operator : CG/JU
Sample : M1959-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B41

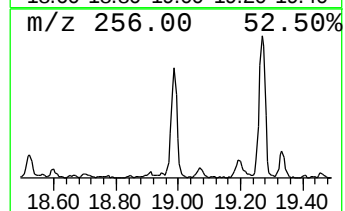
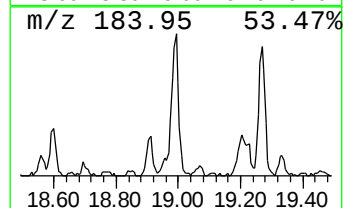
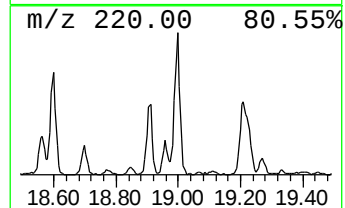
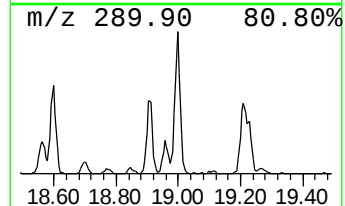
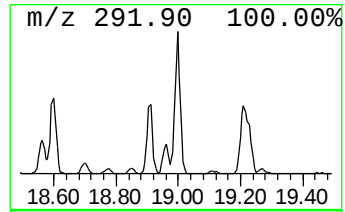
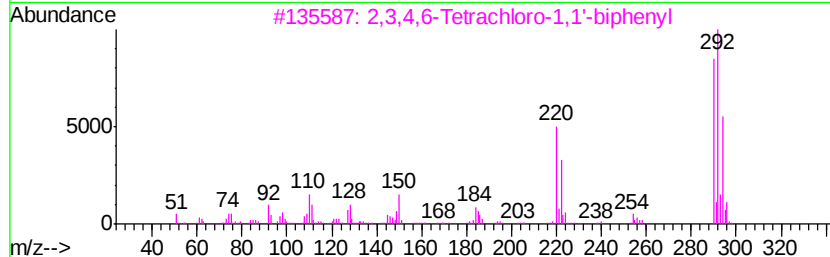
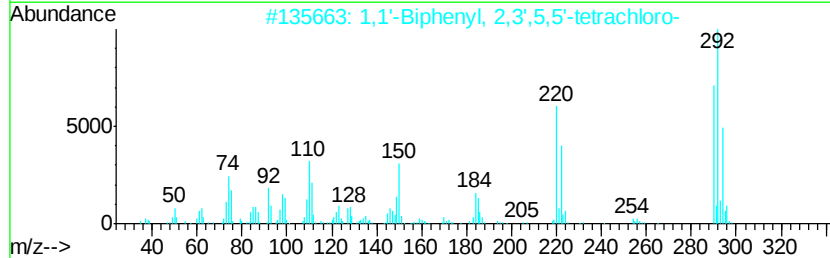
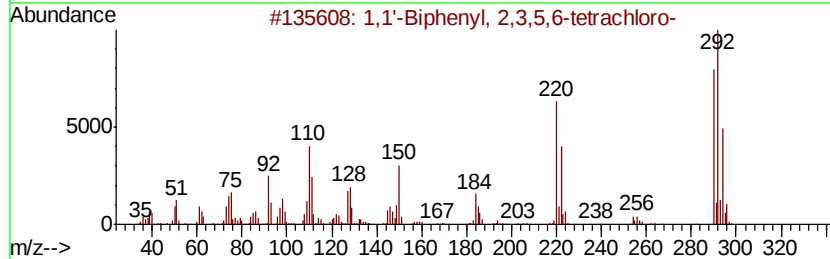
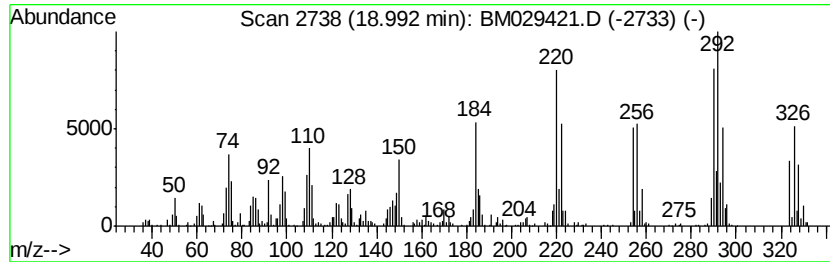
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 1,1'-Biphenyl, 2,3,5,6-tetr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.99	7.27 ng/ul	596159	Phenanthrene-d10	17.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3,5,6-tetrachloro-	290	C12H6Cl4	033284-54-7	99
2		1,1'-Biphenyl, 2,3',5,5'-tetrach...	290	C12H6Cl4	041464-42-0	99
3		2,3,4,6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	054230-22-7	99
4		1,1'-Biphenyl, 2,2',3,4-tetrachl...	290	C12H6Cl4	052663-59-9	99
5		1,1'-Biphenyl, 2,3',4,6-Tetrachl...	290	C12H6Cl4	060233-24-1	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
Data File : BM029421.D
Acq On : 08 Apr 2021 23:41
Operator : CG/JU
Sample : M1959-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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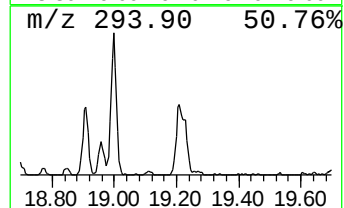
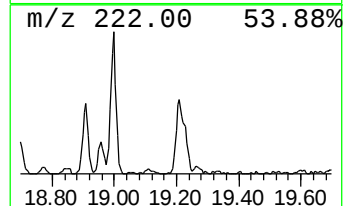
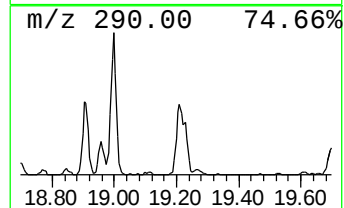
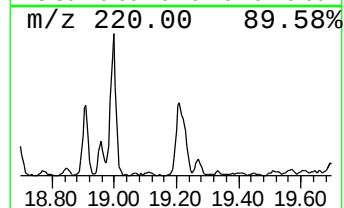
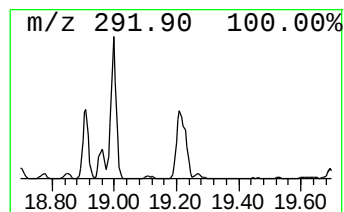
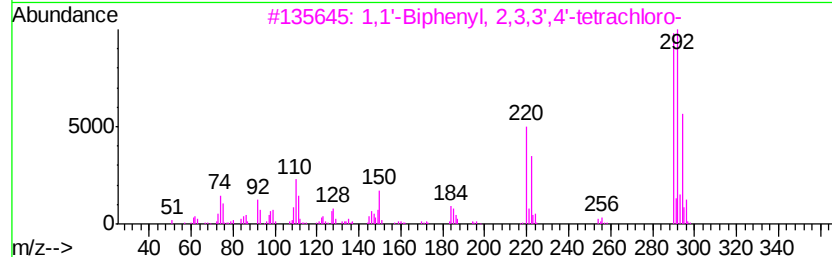
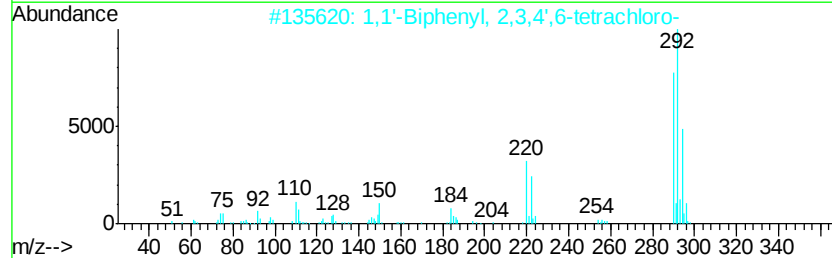
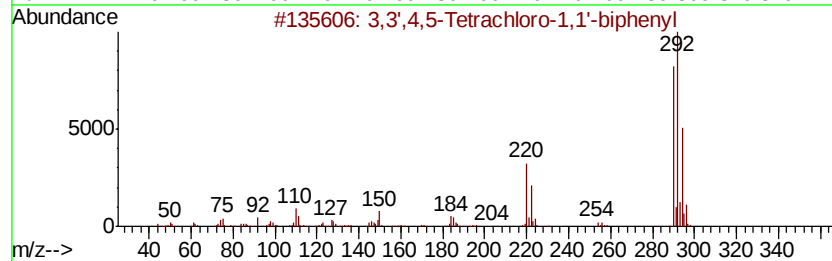
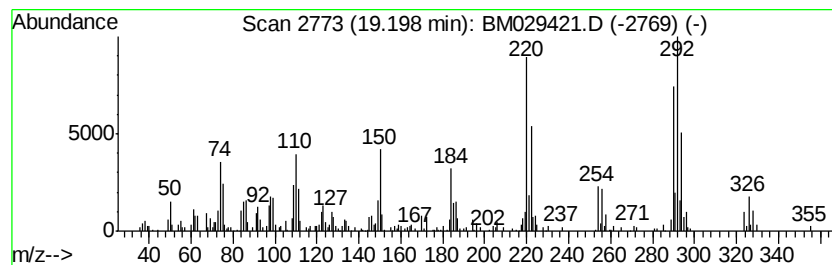
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 3,3',4,5-Tetrachloro-1,1'-b... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.20	4.94 ng/ul	334698	Chrysene-d12	21.24

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-49-1	99
2			1,1'-Biphenyl, 2,3,4',6-tetrachl...	290	C12H6Cl4	052663-58-8	99
3			1,1'-Biphenyl, 2,3,3',4'-tetrach...	290	C12H6Cl4	041464-43-1	99
4			2,3',4',5'-Tetrachloro-1,1'-biph...	290	C12H6Cl4	070362-48-0	99
5			1,1'-Biphenyl, 3,3',4,5'-tetrach...	290	C12H6Cl4	041464-48-6	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
Data File : BM029421.D
Acq On : 08 Apr 2021 23:41
Operator : CG/JU
Sample : M1959-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

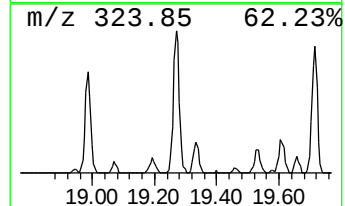
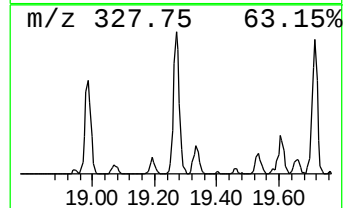
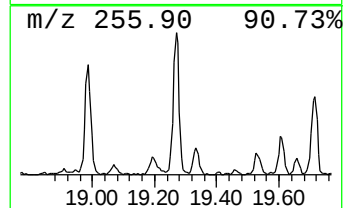
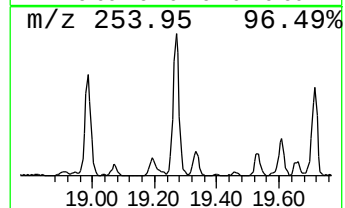
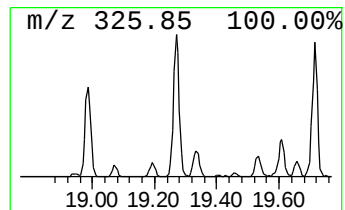
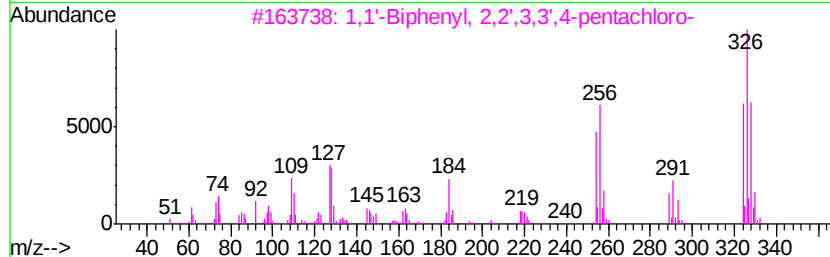
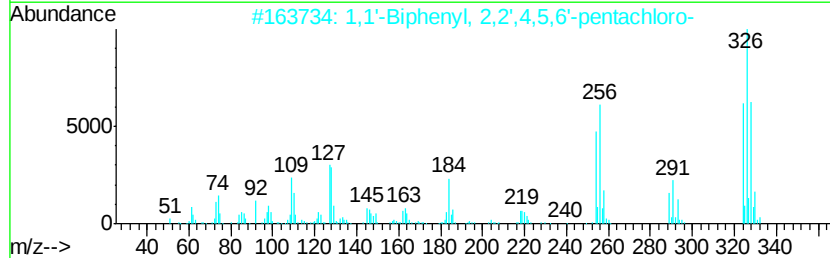
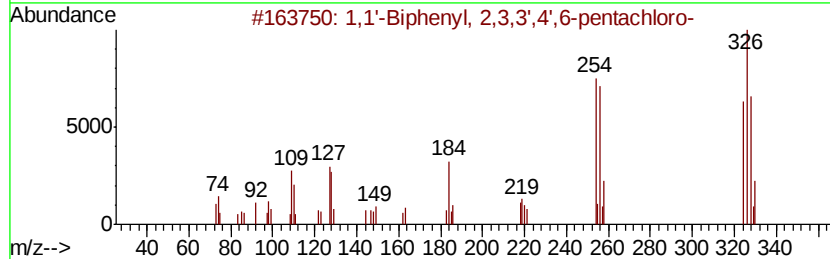
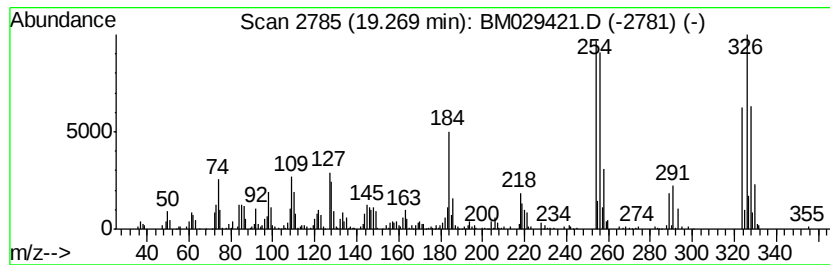
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM040721MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 13 1,1'-Biphenyl, 2,3,3',4',6-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.		
19.27	5.65 ng/ul	382413	Chrysene-d12		21.24		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'	-Biphenyl,	2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99
2	1,1'	-Biphenyl,	2,2',4,5,6'-penta...	324	C12H5Cl5	068194-06-9	99
3	1,1'	-Biphenyl,	2,2',3,3',4-penta...	324	C12H5Cl5	052663-62-4	99
4	1,1'	-Biphenyl,	2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	98
5	2,3,3',	4,5-Pentachloro-1,1'-biph...		324	C12H5Cl5	070424-69-0	97



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
Data File : BM029421.D
Acq On : 08 Apr 2021 23:41
Operator : CG/JU
Sample : M1959-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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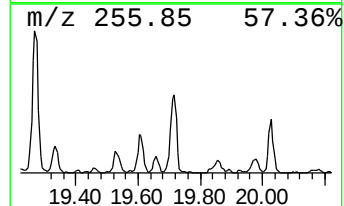
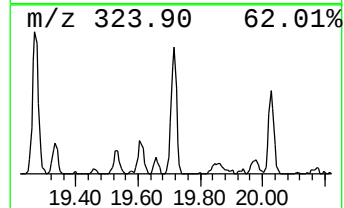
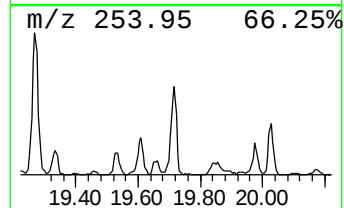
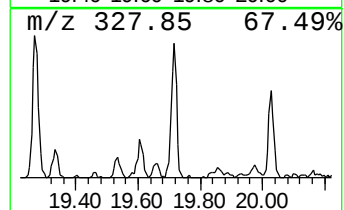
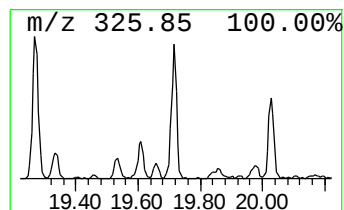
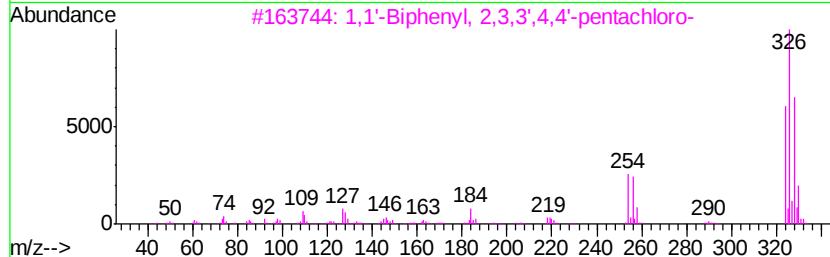
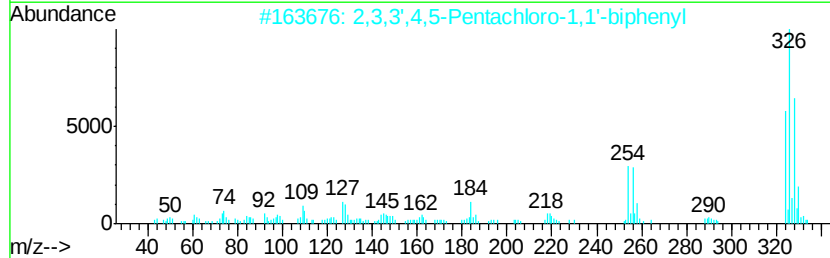
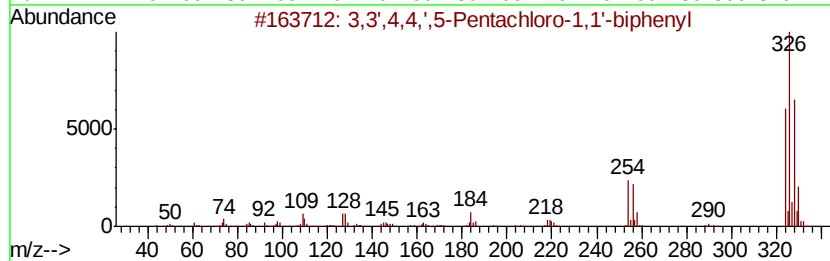
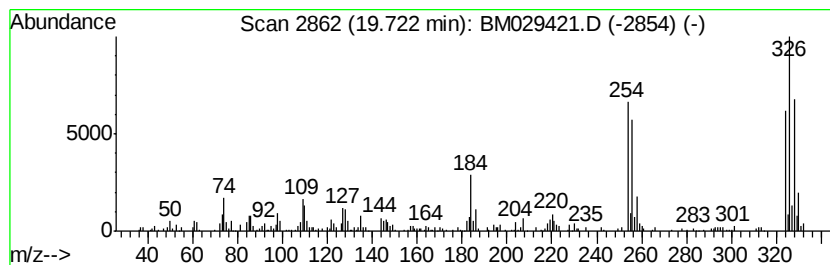
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 3,3',4,4',5-Pentachloro-1,... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.72	4.58 ng/ul	310001	Chrysene-d12	21.24

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,3',4,4',5-Pentachloro-1,1'-bi...	324	C12H5Cl5	057465-28-8	99		
2	2,3,3',4,5-Pentachloro-1,1'-biph...	324	C12H5Cl5	070424-69-0	98		
3	1,1'-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	98		
4	1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	97		
5	2,3,3',4,5-Pentachloro-1,1'-biph...	324	C12H5Cl5	070424-69-0	97		



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
Data File : BM029421.D
Acq On : 08 Apr 2021 23:41
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Sample : M1959-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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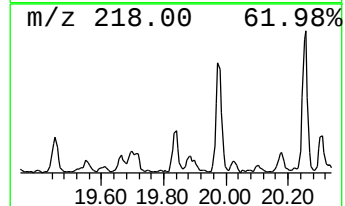
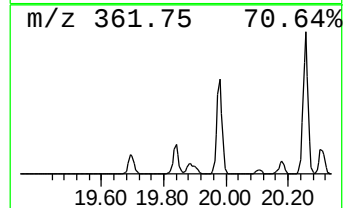
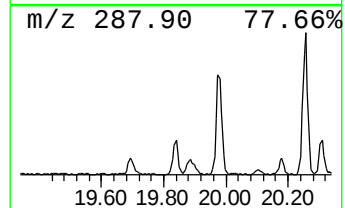
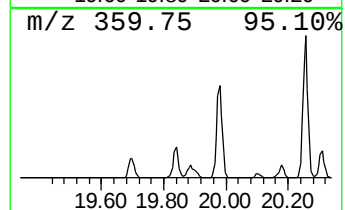
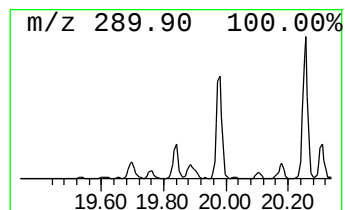
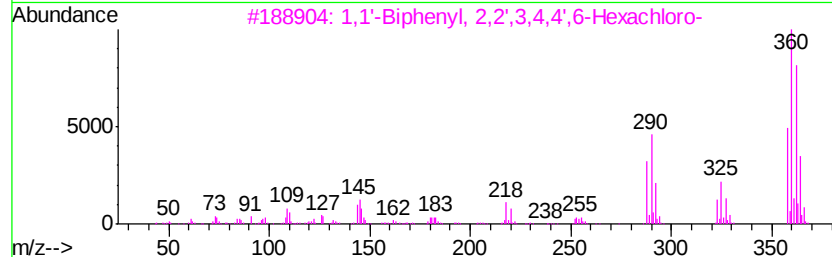
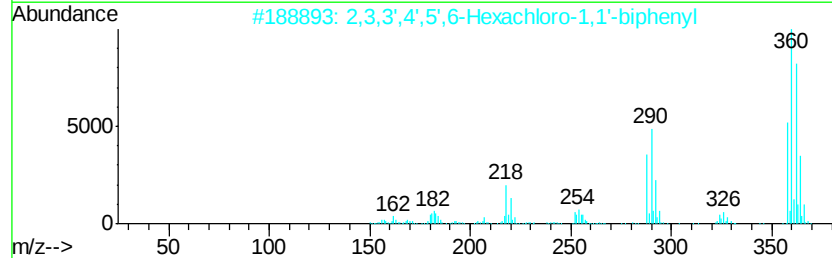
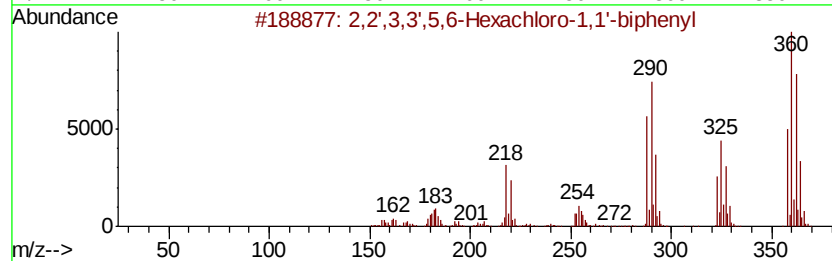
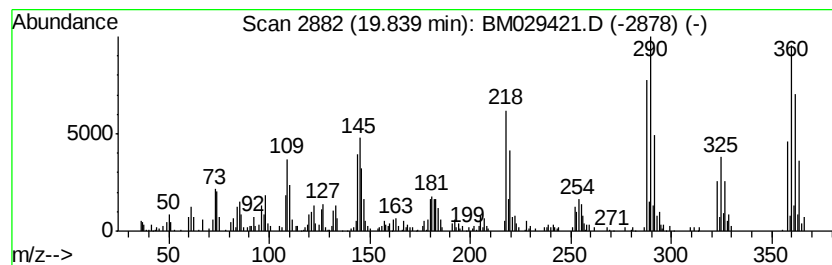
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 2,2',3,3',5,6-Hexachloro-1,... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.84	2.96 ng/ul	200660	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2',3,3',5,6-Hexachloro-1,1'-bi...	358	C12H4Cl6	052704-70-8	99
2		2,3,3',4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	074472-45-0	99
3		1,1'-Biphenyl, 2,2',3,4,4',6-Hex...	358	C12H4Cl6	056030-56-9	99
4		2,3',4,4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-65-5	99
5		2,2',3,3',5,6-Hexachloro-1,1'-bi...	358	C12H4Cl6	052704-70-8	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
Data File : BM029421.D
Acq On : 08 Apr 2021 23:41
Operator : CG/JU
Sample : M1959-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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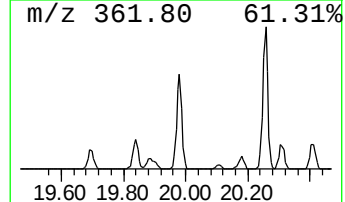
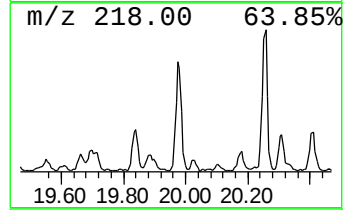
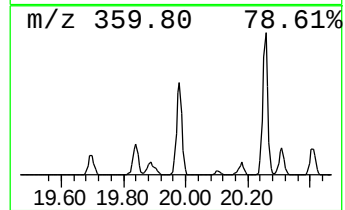
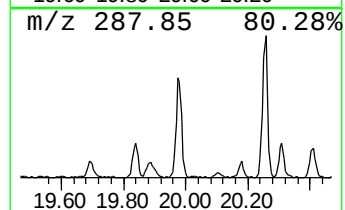
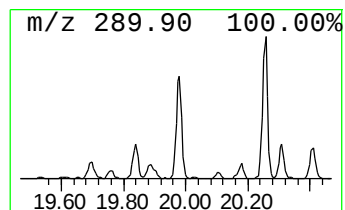
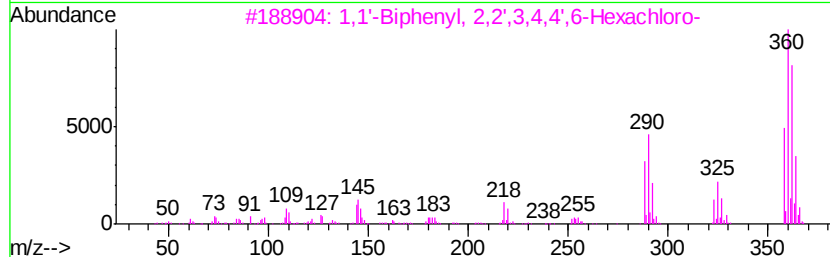
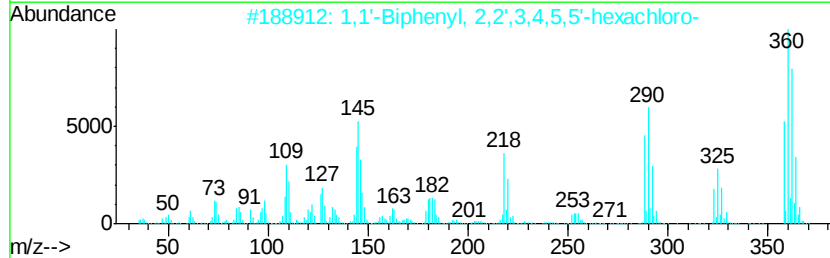
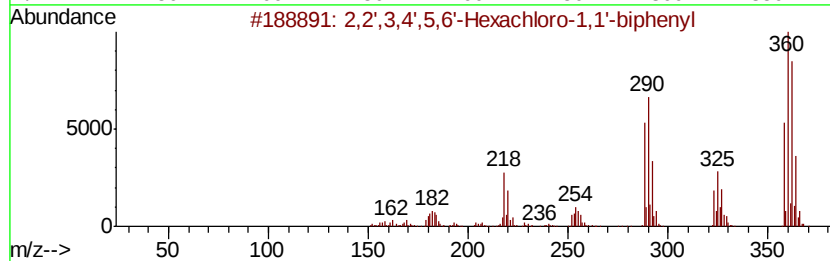
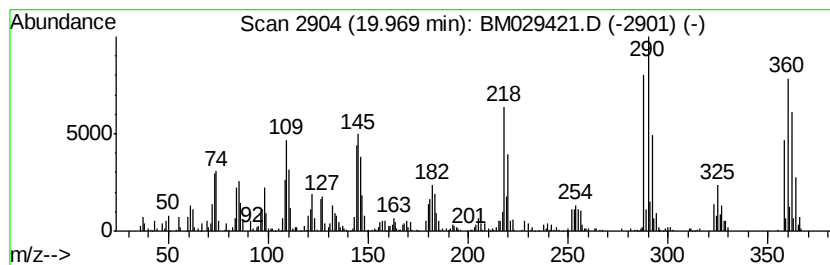
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 2,2',3,4',5,6'-Hexachloro-1... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.97	8.62 ng/ul	583507	Chrysene-d12	21.24

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,4',5,6'-Hexachloro-1,1'-b...	358	C12H4Cl6	074472-41-6	99		
2	1,1'-Biphenyl, 2,2',3,4,5,5'-hex...	358	C12H4Cl6	052712-04-6	99		
3	1,1'-Biphenyl, 2,2',3,4,4',6-Hex...	358	C12H4Cl6	056030-56-9	99		
4	2,2',3,4',6,6'-Hexachloro-1,1'-b...	358	C12H4Cl6	068194-08-1	99		
5	1,1'-Biphenyl, 2,2',4,4',5',6-He...	358	C12H4Cl6	060145-22-4	99		



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
Data File : BM029421.D
Acq On : 08 Apr 2021 23:41
Operator : CG/JU
Sample : M1959-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

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BNA_M
ClientSampleId :
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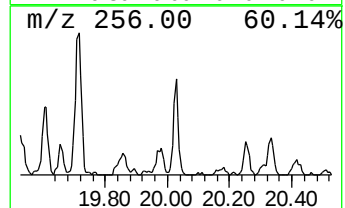
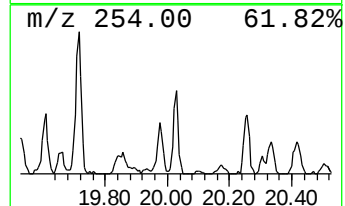
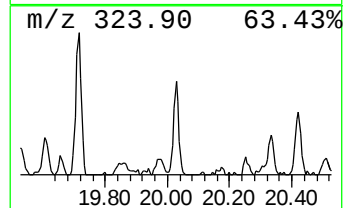
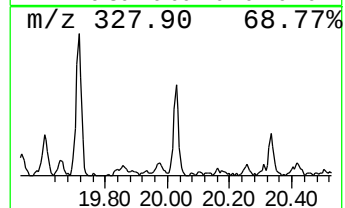
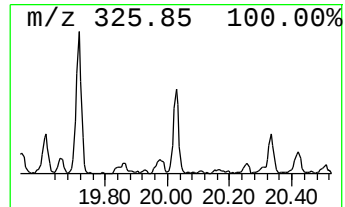
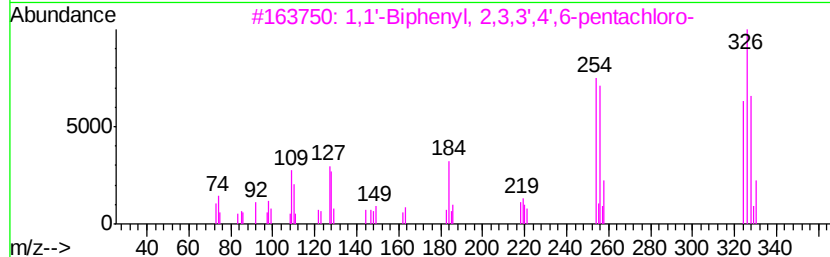
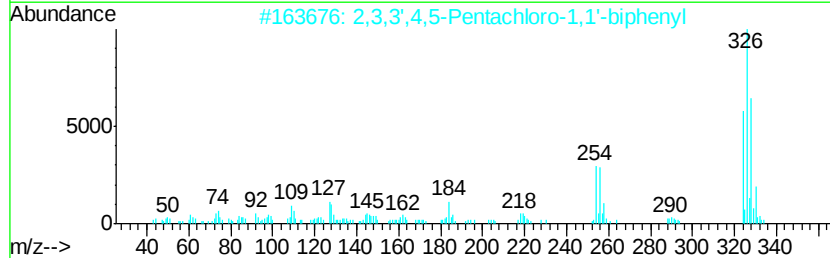
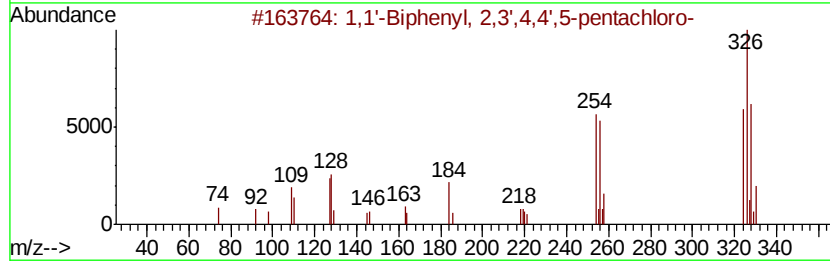
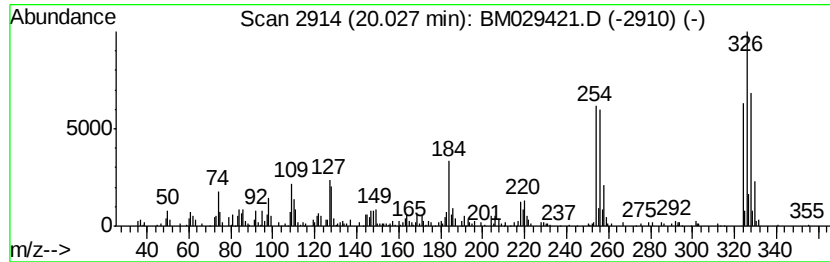
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 1,1'-Biphenyl, 2,3',4,4',5-... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.03	2.02 ng/ul	136677	Chrysene-d12	21.24

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'	-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	99	
2	2,3,3',4,5-Pentachloro-1,1'-biph...	324	C12H5Cl5	070424-69-0	99		
3	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99		
4	1,1'-Biphenyl, 2,3',4,4',6-Penta...	324	C12H5Cl5	056558-17-9	98		
5	2,3,3',5',6-Pentachloro-1,1'-bi...	324	C12H5Cl5	068194-10-5	98		



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
Data File : BM029421.D
Acq On : 08 Apr 2021 23:41
Operator : CG/JU
Sample : M1959-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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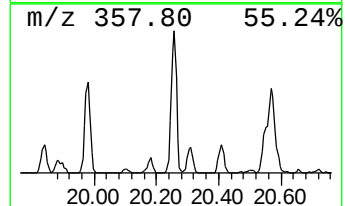
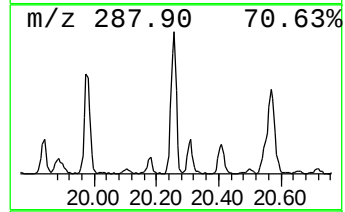
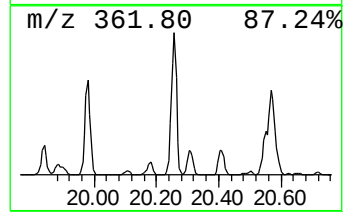
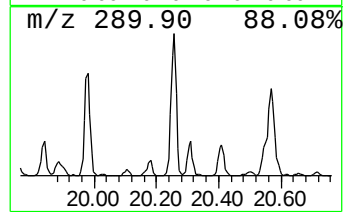
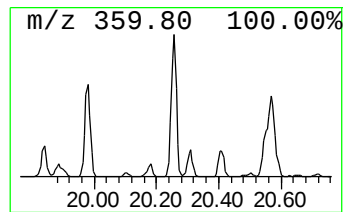
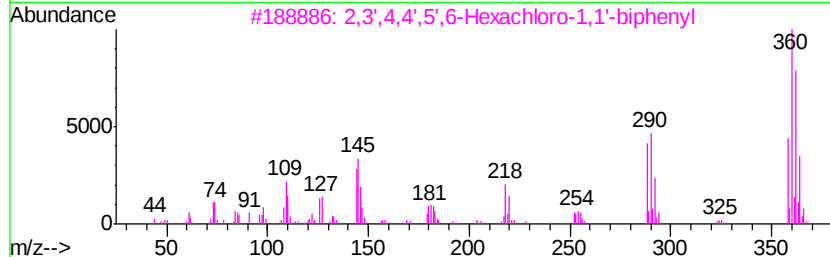
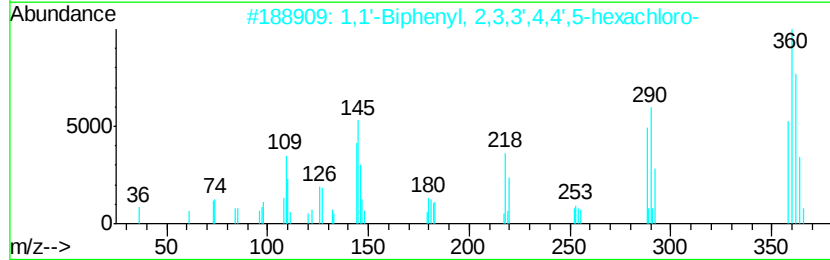
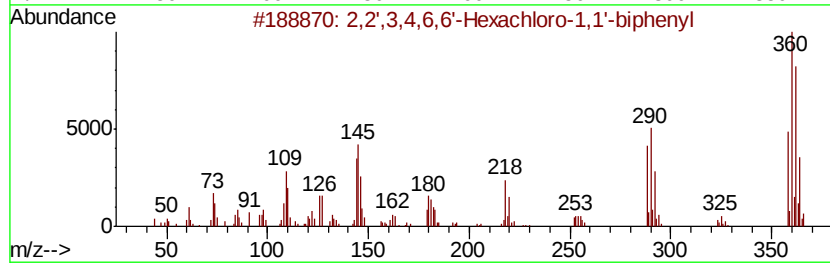
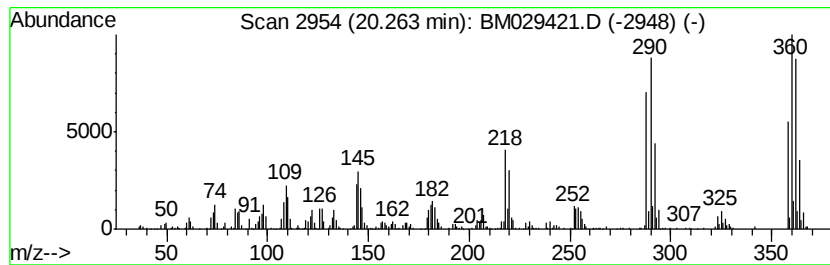
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 2,2',3,4,6,6'-Hexachloro-1,... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.26	10.43 ng/ul	705837	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2',3,4,6,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-40-5	99
2		1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	99
3		2,3',4,4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-65-5	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 M\DATA\BM040721\
Data File : BM029421.D
Acq On : 08 Apr 2021 23:41
Operator : CG/JU
Sample : M1959-01
Misc :
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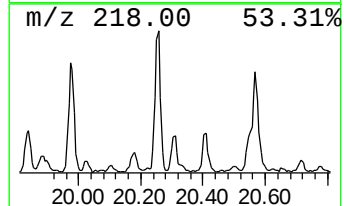
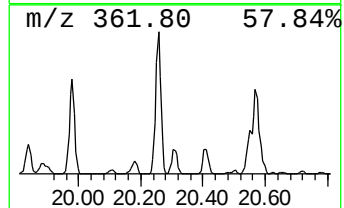
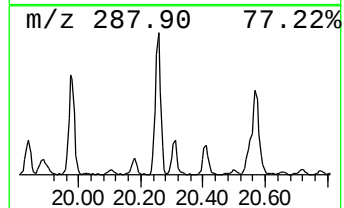
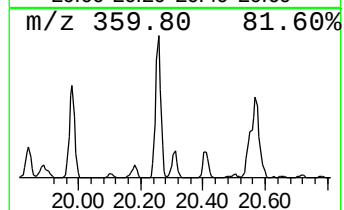
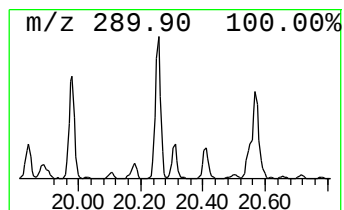
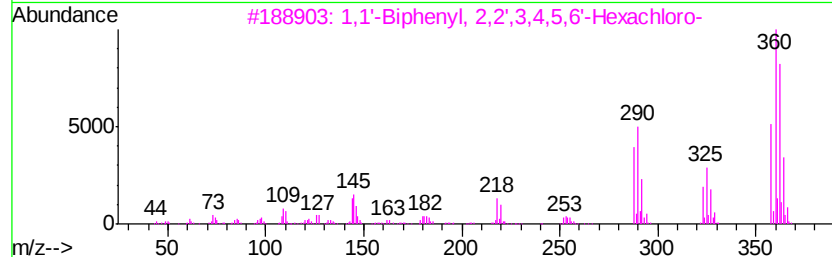
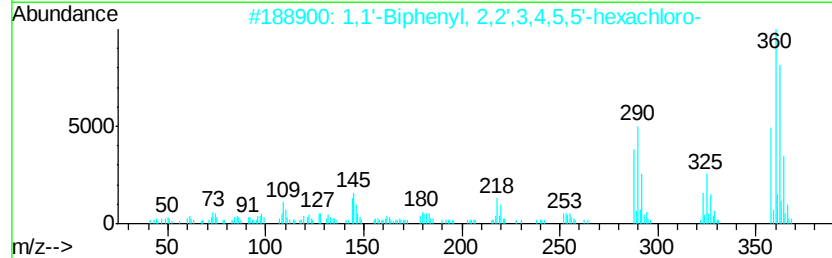
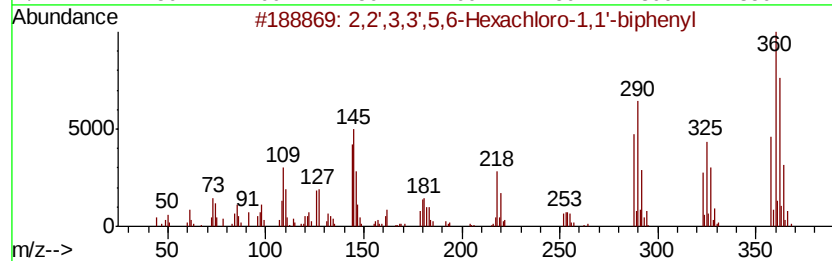
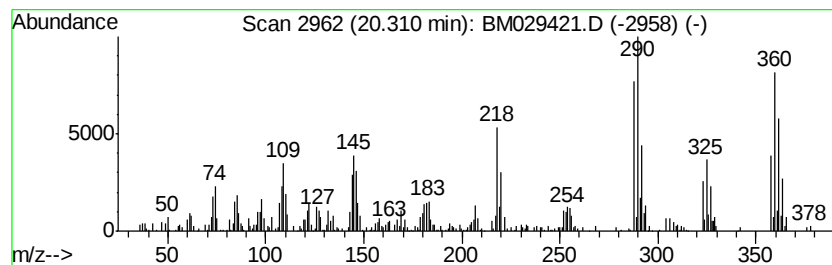
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 1,1'-Biphenyl, 2,2',3,4,5,5... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.31	3.20 ng/ul	216737	Chrysene-d12	21.24

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,3',5,6-Hexachloro-1,1'-bi...	358	C12H4Cl6	052704-70-8	99		
2	1,1'-Biphenyl, 2,2',3,4,5,5'-hex...	358	C12H4Cl6	052712-04-6	99		
3	1,1'-Biphenyl, 2,2',3,4,5,6'-Hex...	358	C12H4Cl6	068194-15-0	99		
4	2,2',3,4,5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-61-4	99		
5	1,1'-Biphenyl, 2,2',3,3',4,5-hex...	358	C12H4Cl6	055215-18-4	99		



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
Data File : BM029421.D
Acq On : 08 Apr 2021 23:41
Operator : CG/JU
Sample : M1959-01
Misc :
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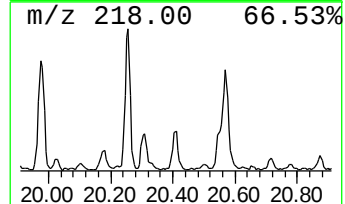
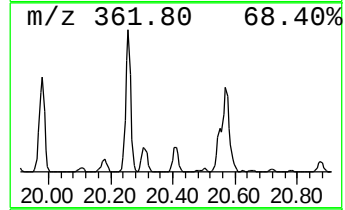
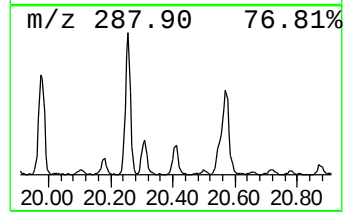
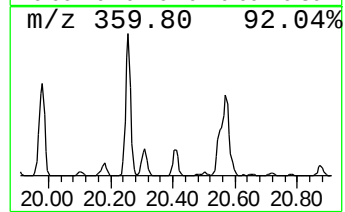
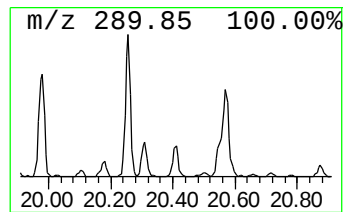
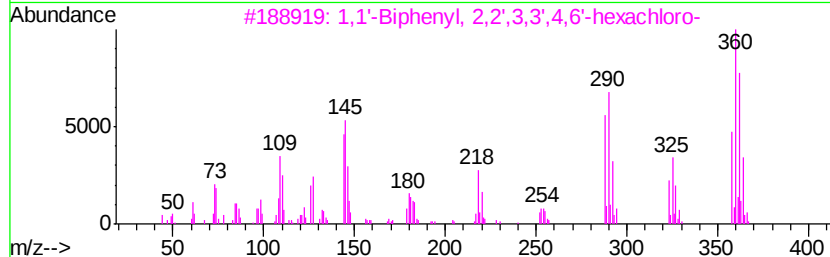
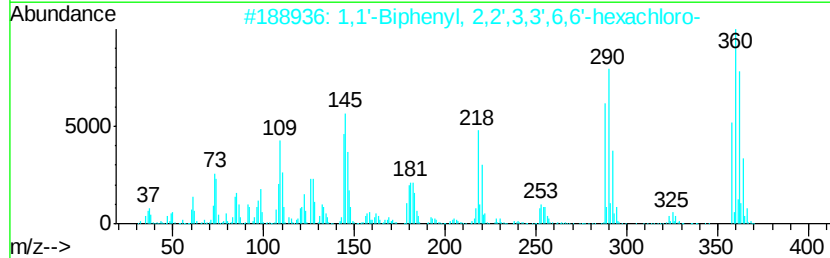
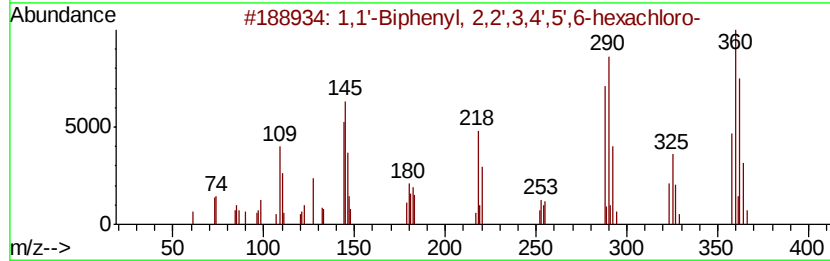
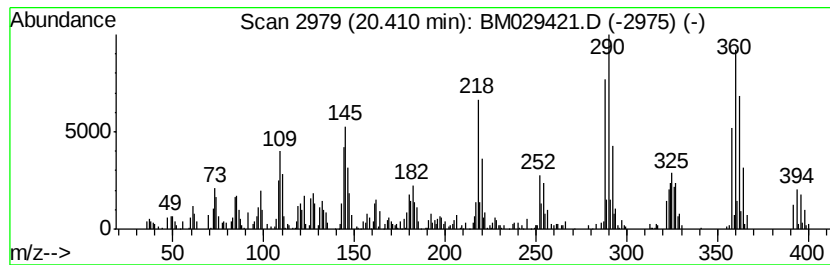
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 1,1'-Biphenyl, 2,2',3,4',5'... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.41	4.10 ng/ul	277818	Chrysene-d12	21.24

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'	-Biphenyl, 2,2',3,4',5',6-he...	358	C12H4Cl6	038380-04-0	99	
2	1,1'	-Biphenyl, 2,2',3,3',6,6'-he...	358	C12H4Cl6	038411-22-2	99	
3	1,1'	-Biphenyl, 2,2',3,3',4,6'-he...	358	C12H4Cl6	038380-05-1	99	
4	1,1'	-Biphenyl, 2,2',3,3',4,4'-he...	358	C12H4Cl6	038380-07-3	99	
5	1,1'	-Biphenyl, 2,2',3,3',4,6'-he...	358	C12H4Cl6	038380-05-1	99	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
Data File : BM029421.D
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Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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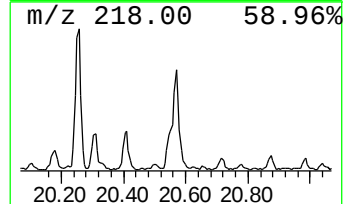
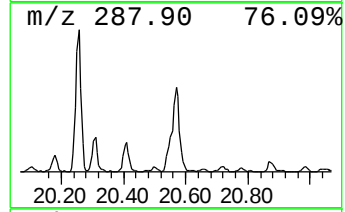
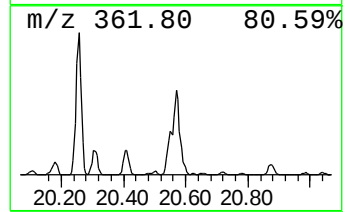
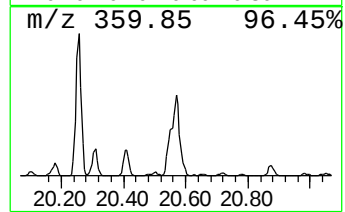
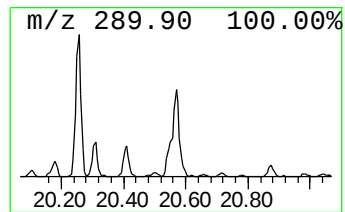
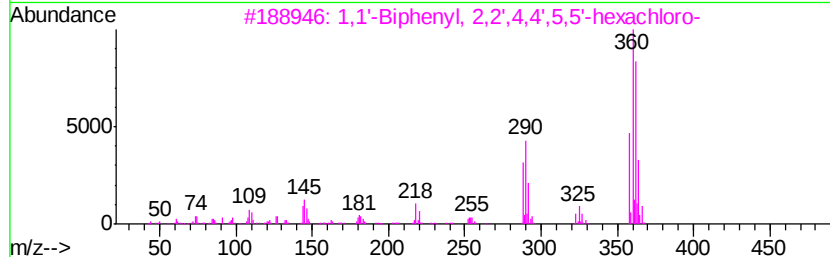
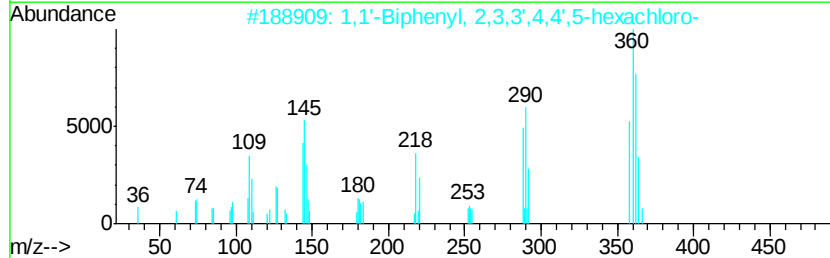
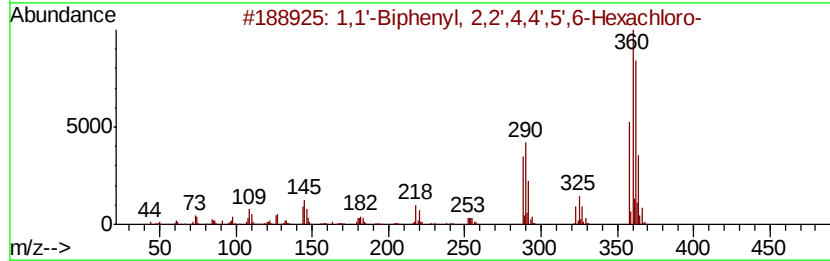
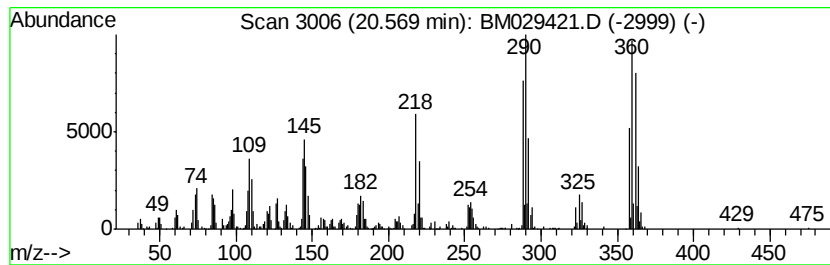
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 1,1'-Biphenyl, 2,2',4,4',5'... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.57	9.95 ng/ul	673363	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',4,4',5',6-He...	358	C12H4Cl6	060145-22-4	99
2		1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	99
3		1,1'-Biphenyl, 2,2',4,4',5,5'-he...	358	C12H4Cl6	035065-27-1	99
4		1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	99
5		2,2',3,5,6,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-09-2	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
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Misc :
ALS Vial : 19 Sample Multiplier: 1

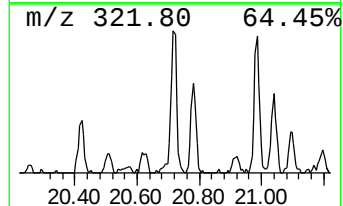
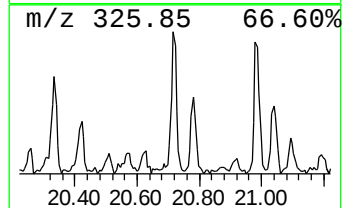
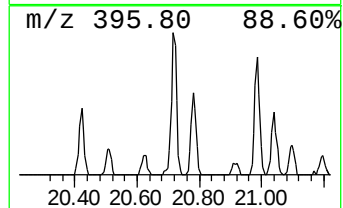
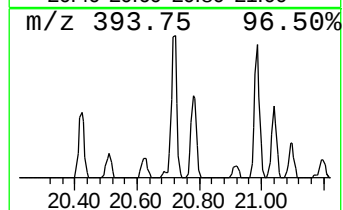
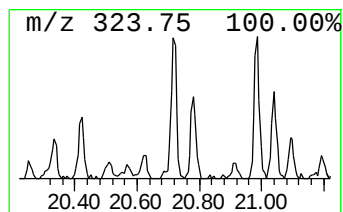
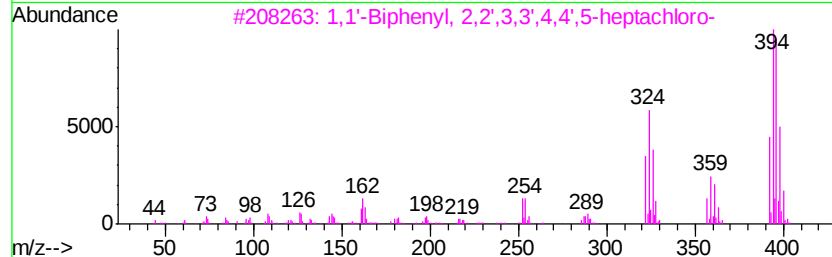
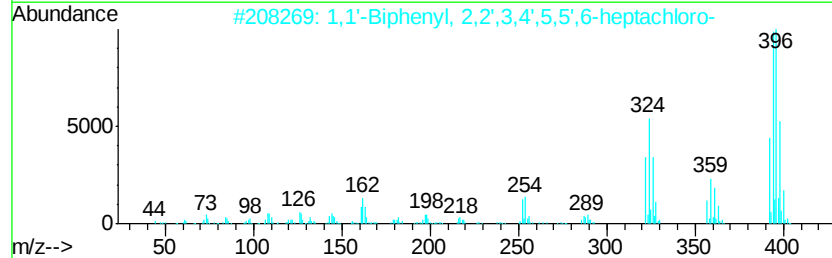
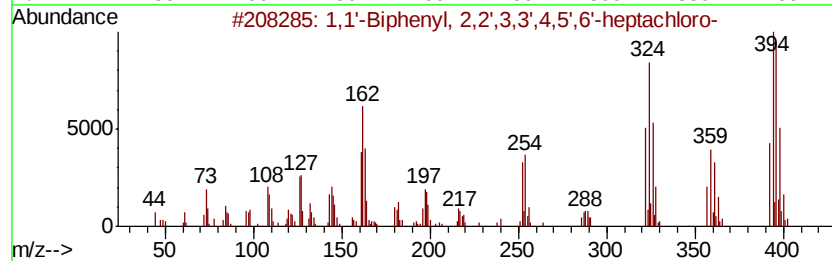
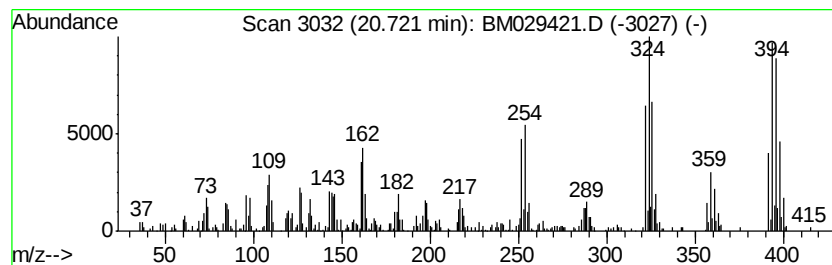
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 1,1'-Biphenyl, 2,2',3,3',4,... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.		
20.72	4.04 ng/ul	273808	Chrysene-d12	21.24		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl,	2,2',3,3',4,5',6'...	392	C12H3Cl7	052663-70-4	99
2	1,1'-Biphenyl,	2,2',3,4',5,5',6-...	392	C12H3Cl7	052663-68-0	99
3	1,1'-Biphenyl,	2,2',3,3',4,4',5-...	392	C12H3Cl7	035065-30-6	99
4	2,3,3',4,4',5',6-Heptachloro-1,1...		392	C12H3Cl7	074472-50-7	99
5	1,1'-Biphenyl,	2,2',3,3',4,6,6'-...	392	C12H3Cl7	052663-65-7	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
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ALS Vial : 19 Sample Multiplier: 1

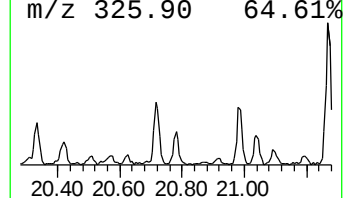
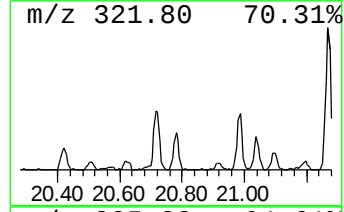
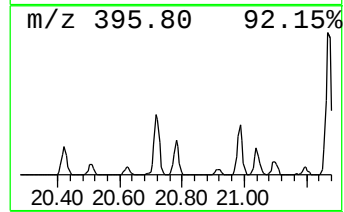
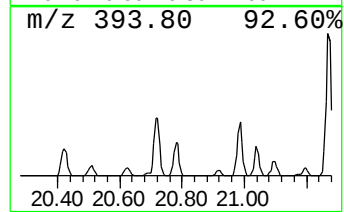
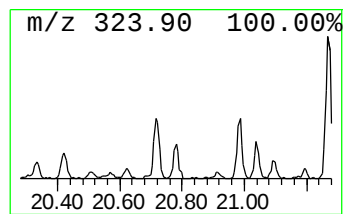
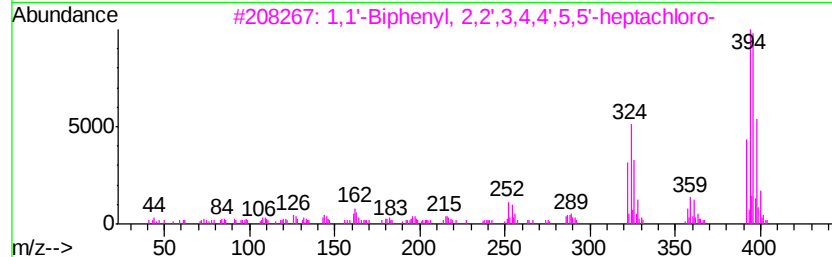
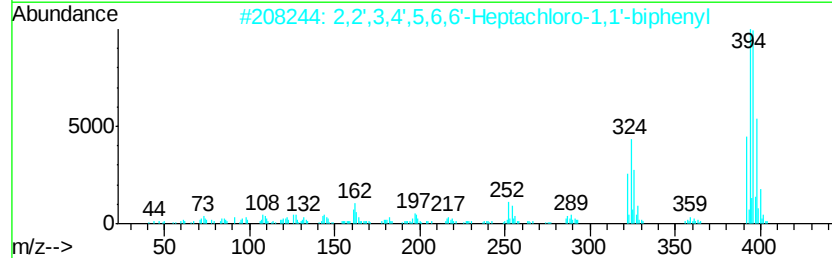
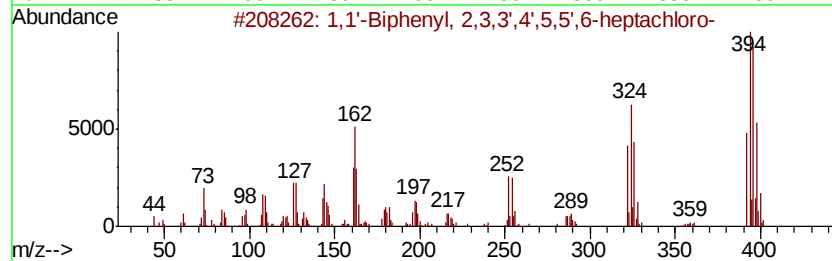
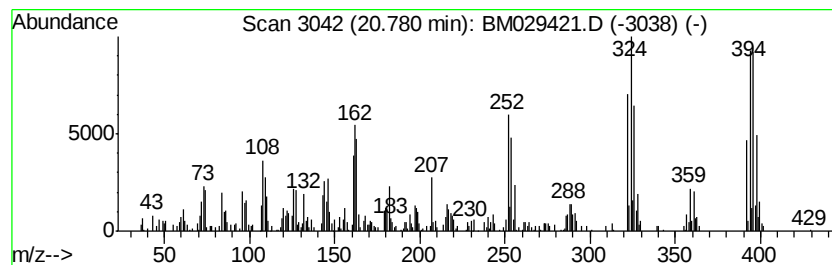
Instrument :
BNA_M
ClientSampleId :
C0B41

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM040721MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 23 1,1'-Biphenyl, 2,3,3',4',5,... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.		
20.78	2.21 ng/ul	149811	Chrysene-d12	21.24		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'	-Biphenyl, 2,3,3',4',5,5',6-...	392	C12H3Cl7	069782-91-8	99
2	2,2',3,4',5,6,6'	-Heptachloro-1,1...	392	C12H3Cl7	074487-85-7	99
3	1,1'-Biphenyl, 2,2',3,4,4',5,5'	-...	392	C12H3Cl7	035065-29-3	99
4	1,1'-Biphenyl, 2,2',3,4,5,6,6'	-H...	392	C12H3Cl7	074472-49-4	99
5	1,1'-Biphenyl, 2,2',3,4,4',5',6-...		392	C12H3Cl7	052663-69-1	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
Data File : BM029421.D
Acq On : 08 Apr 2021 23:41
Operator : CG/JU
Sample : M1959-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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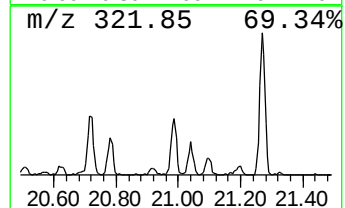
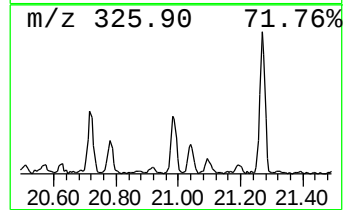
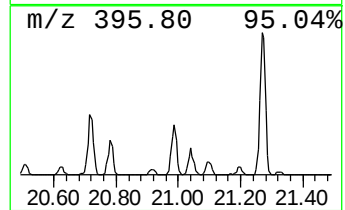
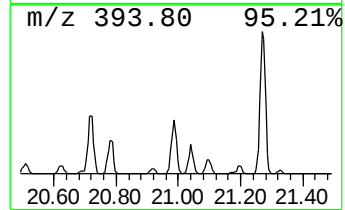
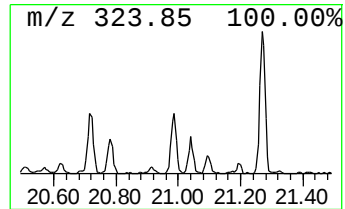
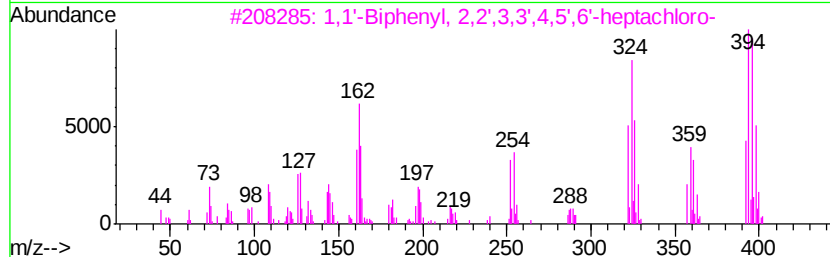
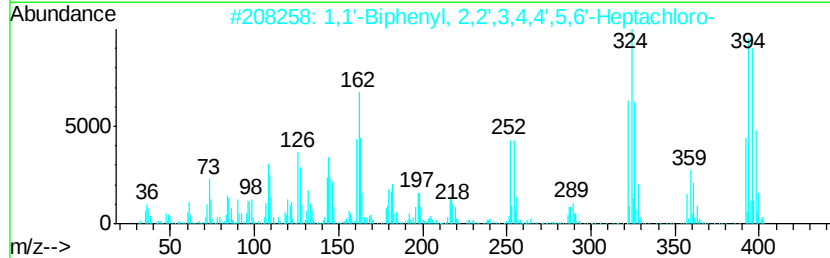
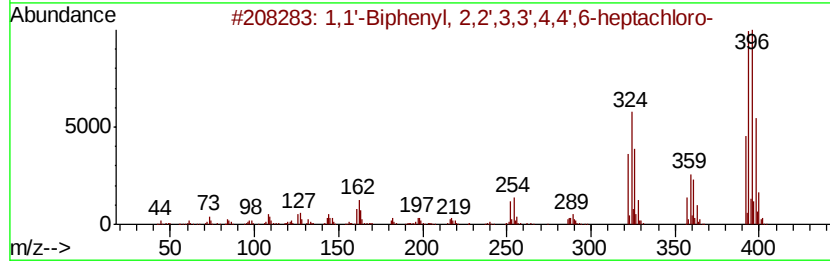
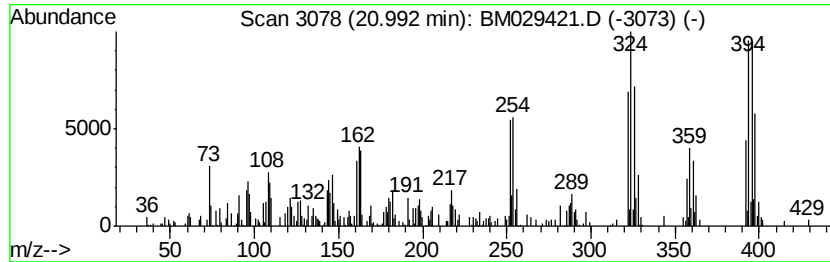
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 1,1'-Biphenyl, 2,2',3,3',4,... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.99	3.82 ng/ul	258923	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',3,3',4,4',6-...	392	C12H3Cl7	052663-71-5	99
2		1,1'-Biphenyl, 2,2',3,4,4',5,6'-...	392	C12H3Cl7	060145-23-5	99
3		1,1'-Biphenyl, 2,2',3,3',4,5',6'...	392	C12H3Cl7	052663-70-4	99
4		2,2',3,4',5,6,6'-Heptachloro-1,1...	392	C12H3Cl7	074487-85-7	99
5		1,1'-Biphenyl, 2,2',3,3',4,4',5-...	392	C12H3Cl7	035065-30-6	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
Data File : BM029421.D
Acq On : 08 Apr 2021 23:41
Operator : CG/JU
Sample : M1959-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

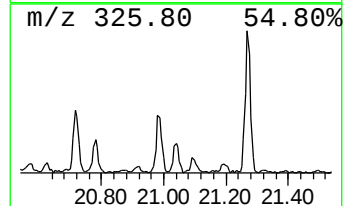
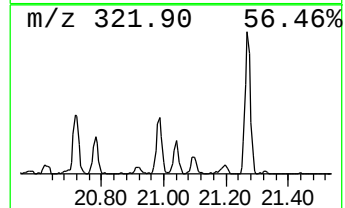
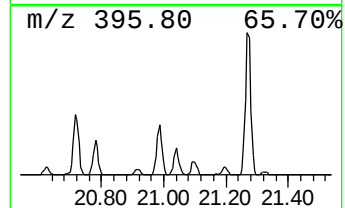
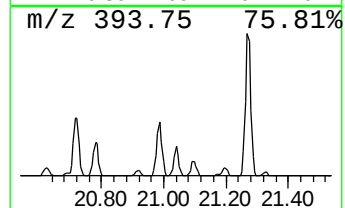
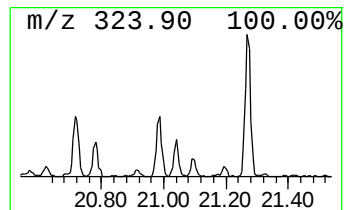
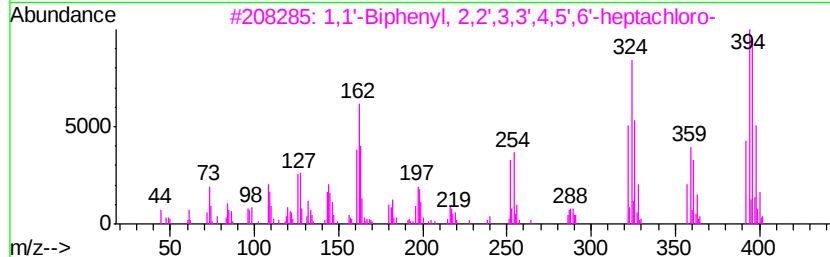
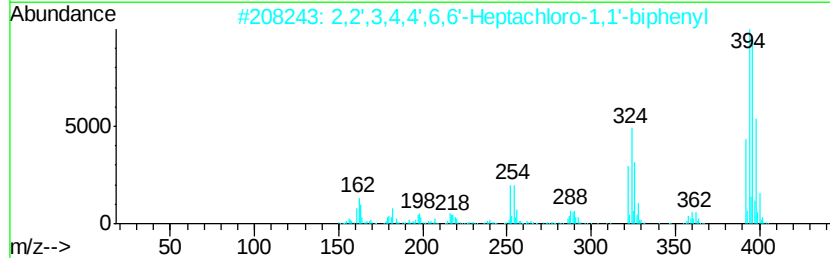
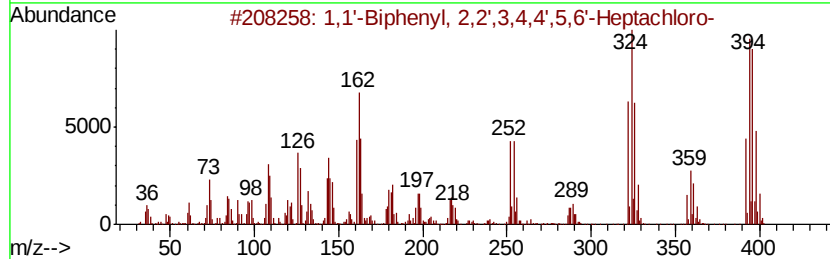
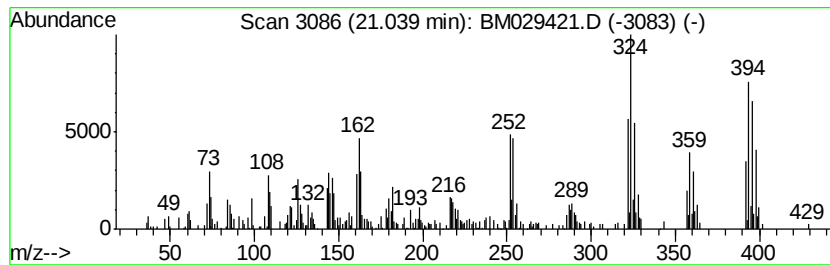
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM040721MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 25 1,1'-Biphenyl, 2,2',3,4,4',... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.		
21.04	2.19 ng/ul	148211	Chrysene-d12	21.24		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,4,4',5,6'-...	392	C12H3Cl7	060145-23-5	99	
2	2,2',3,4,4',6,6'-Heptachloro-1,1...	392	C12H3Cl7	074472-48-3	99	
3	1,1'-Biphenyl, 2,2',3,3',4,5',6'...	392	C12H3Cl7	052663-70-4	99	
4	1,1'-Biphenyl, 2,3,3',4,4',5,6-H...	392	C12H3Cl7	041411-64-7	99	
5	1,1'-Biphenyl, 2,2',3,4,5,6,6'-H...	392	C12H3Cl7	074472-49-4	99	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040721\
Data File : BM029421.D
Acq On : 08 Apr 2021 23:41
Operator : CG/JU
Sample : M1959-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

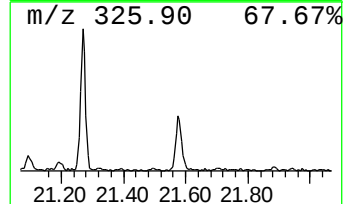
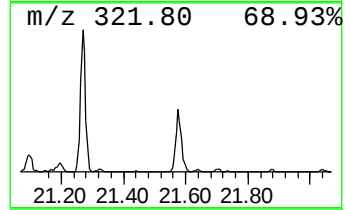
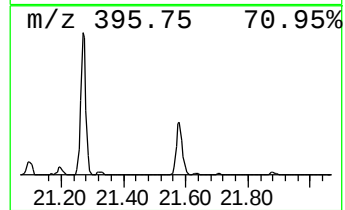
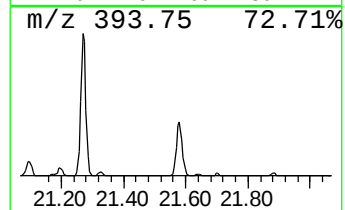
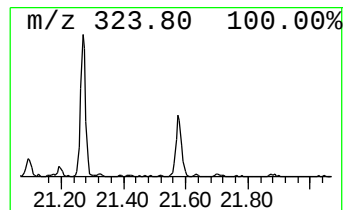
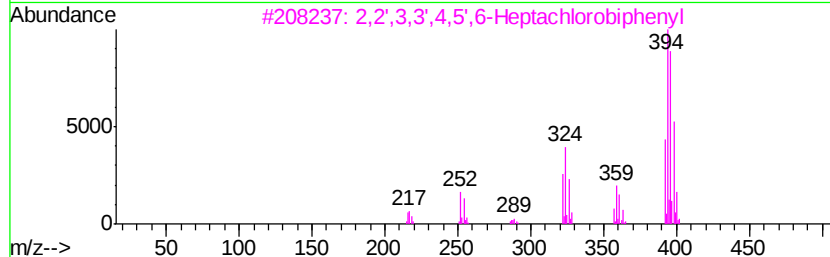
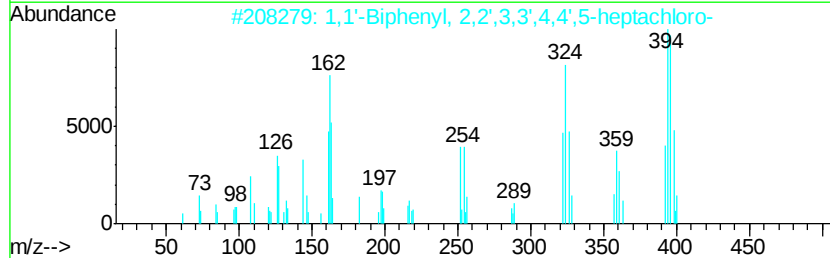
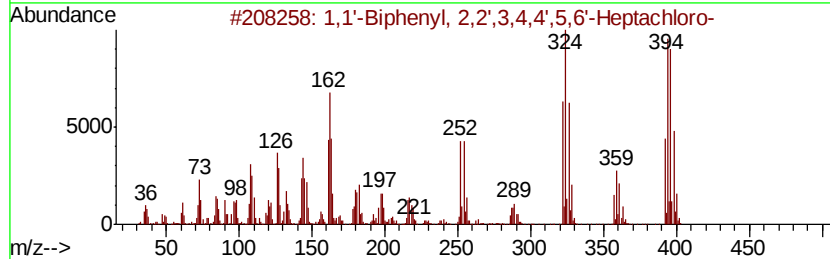
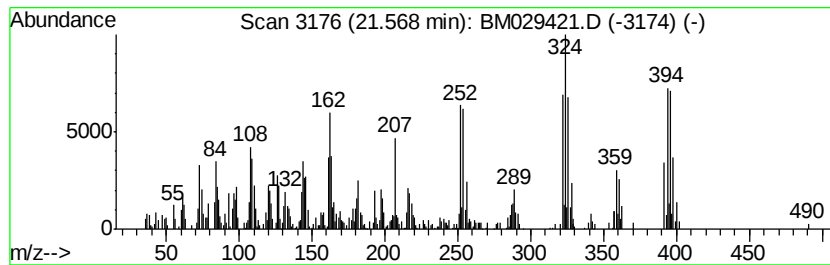
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM040721MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 26 1,1'-Biphenyl, 2,2',3,3',4,... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.		
21.57	3.96 ng/ul	268406	Chrysene-d12	21.24		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,4,4',5,6'-...	392	C12H3Cl7	060145-23-5	99	
2	1,1'-Biphenyl, 2,2',3,3',4,4',5-...	392	C12H3Cl7	035065-30-6	99	
3	2,2',3,3',4,5',6-Heptachlorobiph...	392	C12H3Cl7	040186-70-7	99	
4	1,1'-Biphenyl, 2,2',3,4,4',5',6-...	392	C12H3Cl7	052663-69-1	99	
5	1,1'-Biphenyl, 2,3,3',4,4',5,5'-...	392	C12H3Cl7	039635-31-9	99	



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM040721\
 Data File : BM029421.D
 Acq On : 08 Apr 2021 23:41
 Operator : CG/JU
 Sample : M1959-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B41

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM040721MA.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
4-Pentenal, 2-met...	2.95	3.4	ng/ul	145256	1	7.69	856836	20.0
(DEL) Alkane: Str...	3.12	60.0	ng/ul	2570090	1	7.69	856836	20.0
(DEL) Alkane: Cyc...	3.45	2.1	ng/ul	90856	1	7.69	856836	20.0
1,1'-Biphenyl, 3,...	17.67	6.7	ng/ul	544884	4	17.06	1639020	20.0
1,1'-Biphenyl, 2'...	17.92	2.7	ng/ul	218951	4	17.06	1639020	20.0
1,1'-Biphenyl, 2,...	18.13	4.3	ng/ul	351526	4	17.06	1639020	20.0
2,2',4,5-Tetrachl...	18.19	3.0	ng/ul	242756	4	17.06	1639020	20.0
1,1'-Biphenyl, 2,...	18.42	4.0	ng/ul	330694	4	17.06	1639020	20.0
2,3,3',6-Tetrachl...	18.60	2.8	ng/ul	231791	4	17.06	1639020	20.0
1,1'-Biphenyl, 2,...	18.90	2.5	ng/ul	205984	4	17.06	1639020	20.0
1,1'-Biphenyl, 2,...	18.99	7.3	ng/ul	596159	4	17.06	1639020	20.0
3,3',4,5-Tetrachl...	19.20	4.9	ng/ul	334698	5	21.24	1354120	20.0
1,1'-Biphenyl, 2,...	19.27	5.7	ng/ul	382413	5	21.24	1354120	20.0
3,3',4,4',5-Pent...	19.72	4.6	ng/ul	310001	5	21.24	1354120	20.0
2,2',3,3',5,6-Hex...	19.84	3.0	ng/ul	200660	5	21.24	1354120	20.0
2,2',3,4',5,6'-He...	19.97	8.6	ng/ul	583507	5	21.24	1354120	20.0
1,1'-Biphenyl, 2,...	20.03	2.0	ng/ul	136677	5	21.24	1354120	20.0
2,2',3,4,6,6'-Hex...	20.26	10.4	ng/ul	705837	5	21.24	1354120	20.0
1,1'-Biphenyl, 2,...	20.31	3.2	ng/ul	216737	5	21.24	1354120	20.0
1,1'-Biphenyl, 2,...	20.41	4.1	ng/ul	277818	5	21.24	1354120	20.0
1,1'-Biphenyl, 2,...	20.57	9.9	ng/ul	673363	5	21.24	1354120	20.0
1,1'-Biphenyl, 2,...	20.72	4.0	ng/ul	273808	5	21.24	1354120	20.0
1,1'-Biphenyl, 2,...	20.78	2.2	ng/ul	149811	5	21.24	1354120	20.0
1,1'-Biphenyl, 2,...	20.99	3.8	ng/ul	258923	5	21.24	1354120	20.0
1,1'-Biphenyl, 2,...	21.04	2.2	ng/ul	148211	5	21.24	1354120	20.0
1,1'-Biphenyl, 2,...	21.57	4.0	ng/ul	268406	5	21.24	1354120	20.0