

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052021\
 Data File : BM030151.D
 Acq On : 22 May 2021 15:39
 Operator : CG/JU
 Sample : M2455-01
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0BA1

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\SFAM-EPA-BM050721.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	2.940	5	9	23	rVB	3222589	3289719	100.00%	15.344%
2	3.263	61	64	70	rVB	86319	97958	2.98%	0.457%
3	3.740	141	145	150	rVB	13238	17715	0.54%	0.083%
4	4.169	214	218	223	rVB	6191	9925	0.30%	0.046%
5	4.263	229	234	240	rVB7	3826	7324	0.22%	0.034%
6	7.504	778	785	807	rBV	463768	901886	27.42%	4.207%
7	10.275	1249	1256	1278	rBV	611592	1240828	37.72%	5.788%
8	13.557	1810	1814	1822	rBV10	5750	13314	0.40%	0.062%
9	13.892	1866	1871	1872	rBV5	2910	3738	0.11%	0.017%
10	13.980	1882	1886	1891	rVB7	2469	4346	0.13%	0.020%
11	14.116	1905	1909	1910	rBV2	4367	4047	0.12%	0.019%
12	14.157	1910	1916	1926	rBV	1144005	1734906	52.74%	8.092%
13	14.804	2023	2026	2028	rBV4	4371	5990	0.18%	0.028%
14	14.863	2033	2036	2042	rVV7	3836	6978	0.21%	0.033%
15	14.939	2045	2049	2052	rBV6	6696	9447	0.29%	0.044%
16	15.039	2063	2066	2069	rBV4	2809	3827	0.12%	0.018%
17	15.157	2083	2086	2089	rBV5	2492	3467	0.11%	0.016%
18	15.639	2165	2168	2171	rBV5	4019	6318	0.19%	0.029%
19	15.680	2171	2175	2176	rVV4	3490	4127	0.13%	0.019%
20	16.398	2295	2297	2298	rBV2	4602	3496	0.11%	0.016%
21	16.910	2378	2384	2389	rBV	1297932	1971915	59.94%	9.197%
22	17.521	2482	2488	2493	rBV	78150	127924	3.89%	0.597%
23	17.763	2526	2529	2532	rBV3	19187	24622	0.75%	0.115%
24	17.980	2561	2566	2571	rBV	89193	135752	4.13%	0.633%
25	18.039	2572	2576	2580	rVB4	48388	65797	2.00%	0.307%
26	18.268	2611	2615	2618	rBV	55098	73109	2.22%	0.341%
27	18.445	2642	2645	2650	rVB	41381	53155	1.62%	0.248%
28	18.757	2695	2698	2703	rVB4	35313	53066	1.61%	0.248%
29	18.833	2704	2711	2719	rBV2	182732	346340	10.53%	1.615%
30	18.974	2731	2735	2743	rBV	133030	203853	6.20%	0.951%
31	19.051	2744	2748	2755	rVV7	37367	105855	3.22%	0.494%
32	19.121	2755	2760	2767	rVB	245246	334522	10.17%	1.560%
33	19.462	2815	2818	2822	rVB4	46732	57270	1.74%	0.267%
34	19.545	2829	2832	2833	rBV	81866	84969	2.58%	0.396%

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35	19.568	2833	2836	2842	rVB	145003	182867	5.56%	0.853%
36	19.692	2853	2857	2861	rBV	151039	193277	5.88%	0.901%
37	19.733	2861	2864	2871	rVB4	67402	112199	3.41%	0.523%
38	19.827	2876	2880	2885	rVV	495403	624362	18.98%	2.912%
39	19.886	2886	2890	2896	rVB3	77981	110583	3.36%	0.516%
40	20.033	2912	2915	2919	rVB	70567	84582	2.57%	0.395%
41	20.109	2924	2928	2933	rVB	728862	794304	24.15%	3.705%
42	20.162	2933	2937	2941	rBV	169568	216578	6.58%	1.010%
43	20.262	2950	2954	2960	rVB2	197164	274123	8.33%	1.279%
44	20.356	2968	2970	2974	rBV4	46168	56809	1.73%	0.265%
45	20.427	2974	2982	2988	rBV	449505	850315	25.85%	3.966%
46	20.574	3003	3007	3011	rVB	222107	273181	8.30%	1.274%
47	20.633	3014	3017	3021	rBV2	128475	146282	4.45%	0.682%
48	20.733	3031	3034	3036	rBV3	54825	56440	1.72%	0.263%
49	20.839	3048	3052	3058	rBV2	251917	312159	9.49%	1.456%
50	20.898	3058	3062	3067	rVB	137985	170190	5.17%	0.794%
51	20.951	3068	3071	3074	rBV4	71093	84984	2.58%	0.396%
52	21.098	3091	3096	3099	rBV	1781414	2334595	70.97%	10.889%
53	21.127	3099	3101	3109	rVB2	728580	835596	25.40%	3.897%
54	21.433	3149	3153	3157	rBV2	265175	348637	10.60%	1.626%
55	21.486	3160	3162	3168	rVB5	94262	115276	3.50%	0.538%
56	21.551	3171	3173	3177	rVB4	92815	101133	3.07%	0.472%
57	23.239	3455	3460	3471	rVB2	1185375	2153797	65.47%	10.046%

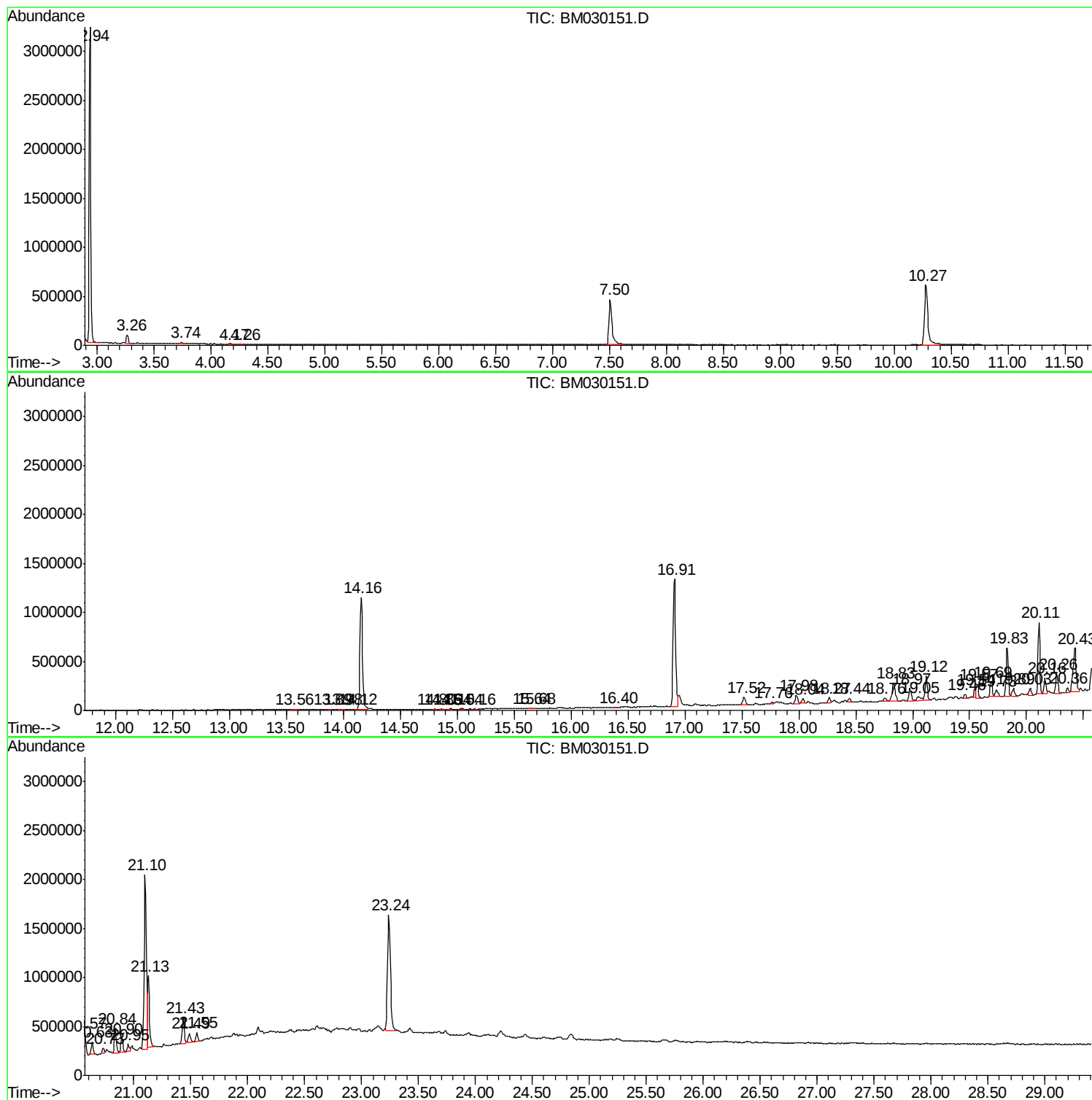
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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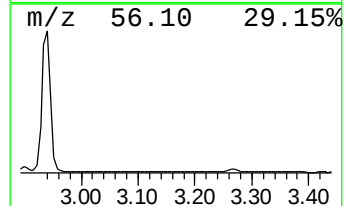
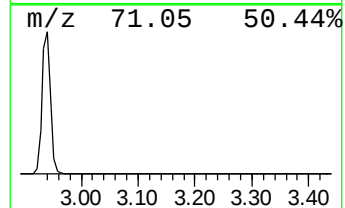
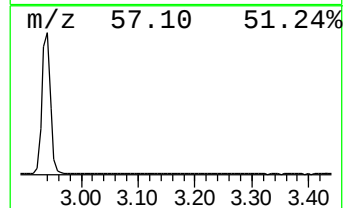
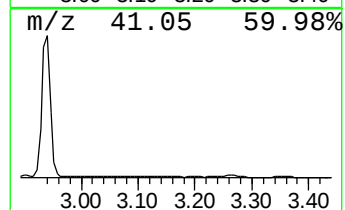
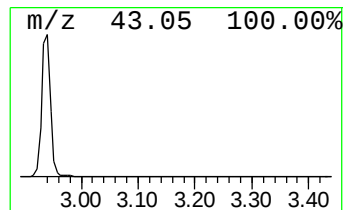
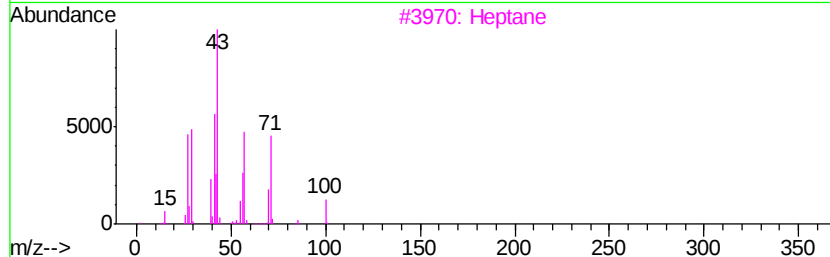
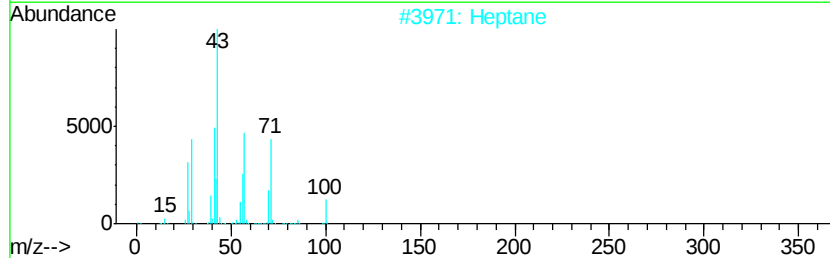
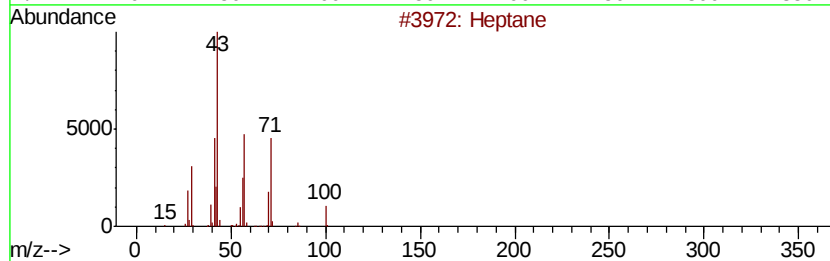
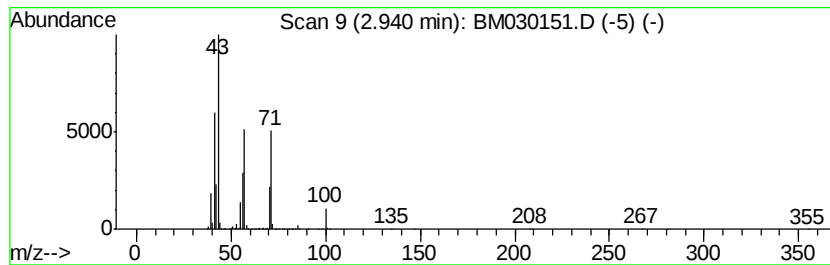
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.94	72.95 ng/ul	3289720	1,4-Dichlorobenzene-d4	7.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane	100	C7H16	000142-82-5	95
2		Heptane	100	C7H16	000142-82-5	94
3		Heptane	100	C7H16	000142-82-5	91
4		Heptane	100	C7H16	000142-82-5	83
5		Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	64



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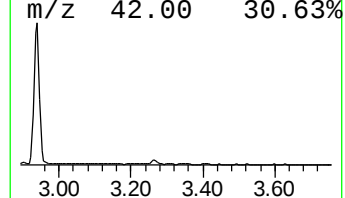
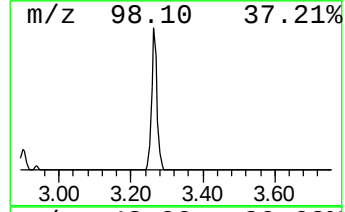
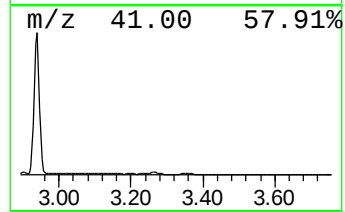
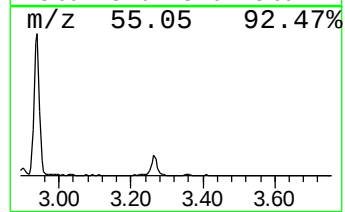
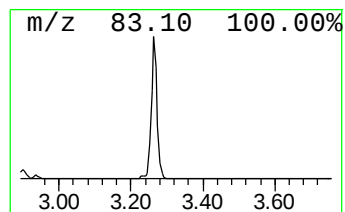
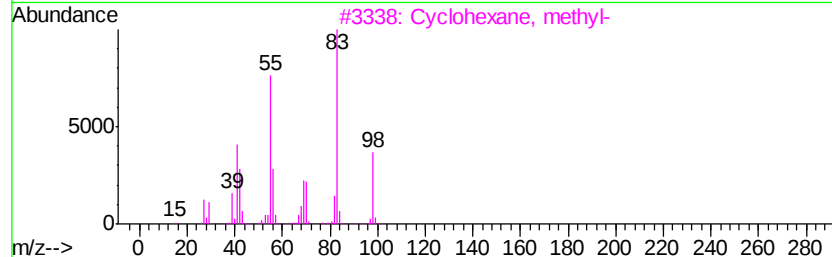
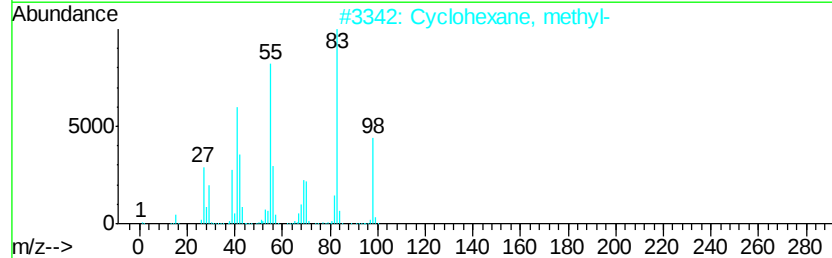
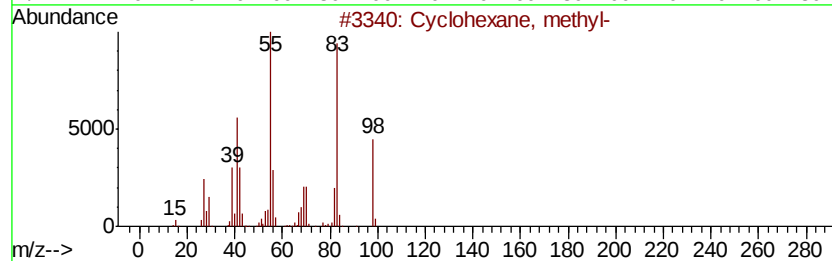
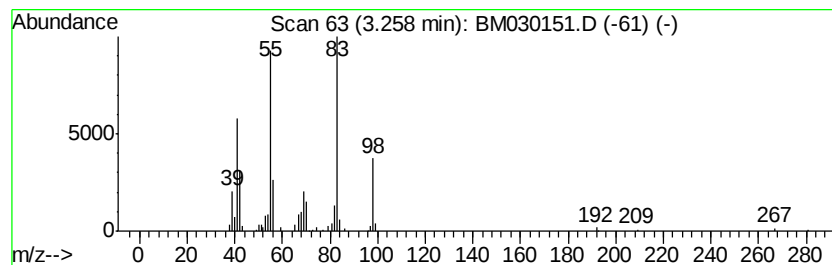
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Cyclic3.26 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.26	2.17 ng/ul	97958	1,4-Dichlorobenzene-d4	7.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, methyl-	98	C7H14	000108-87-2	94
2		Cyclohexane, methyl-	98	C7H14	000108-87-2	91
3		Cyclohexane, methyl-	98	C7H14	000108-87-2	90
4		2-Pentene, 4,4-dimethyl-, (Z)-	98	C7H14	000762-63-0	64
5		2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	64



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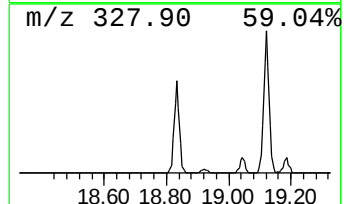
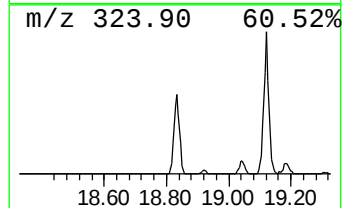
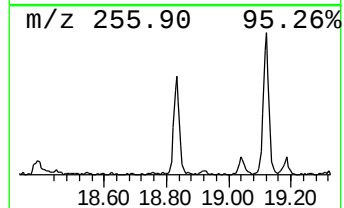
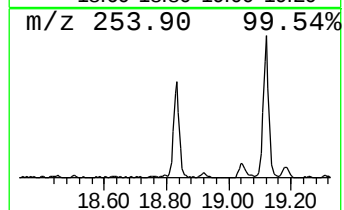
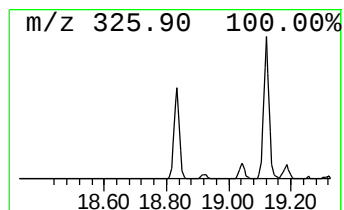
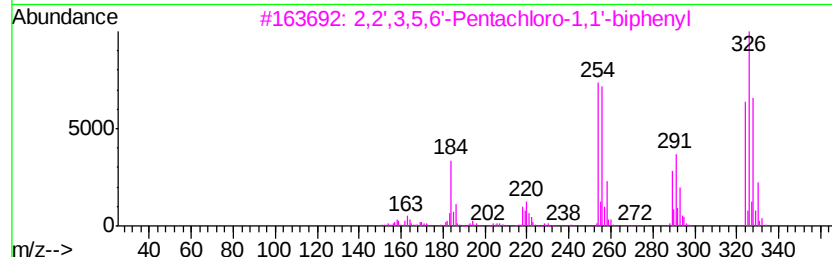
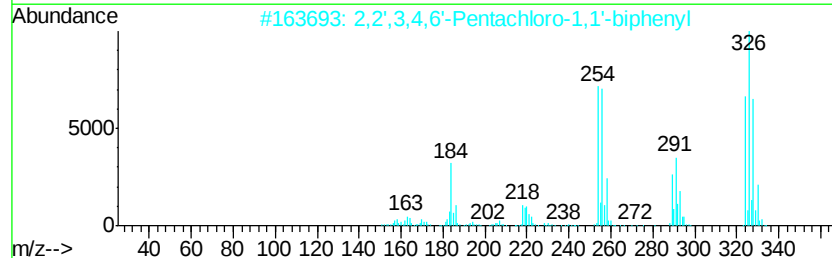
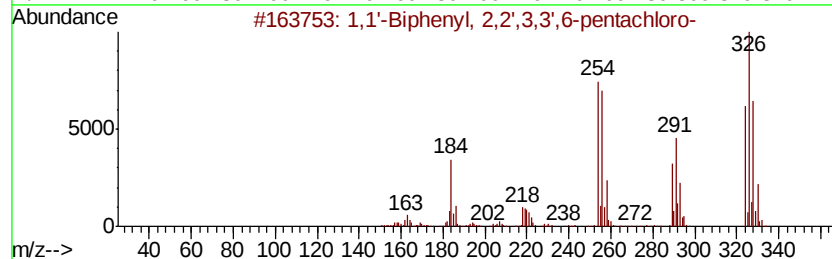
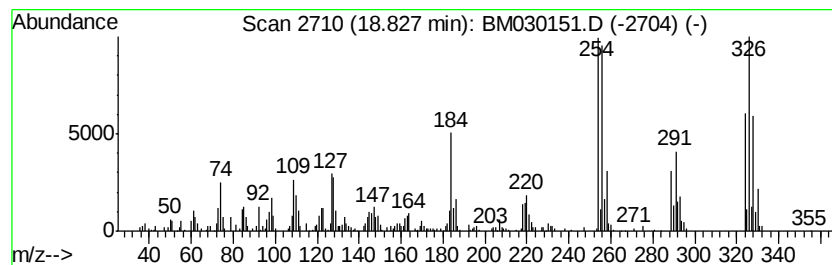
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.83	3.51 ng/ul	346340	Phenanthrene-d10	16.91	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,3',6-penta...	324	C12H5Cl5	052663-60-2	99
2	2,2',3,4,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-57-2	99
3	2,2',3,5,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-55-0	99
4	1,1'-Biphenyl, 2,2',3,3',5-penta...	324	C12H5Cl5	060145-20-2	99
5	1,1'-Biphenyl, 2,2',3,3',5-penta...	324	C12H5Cl5	060145-20-2	99



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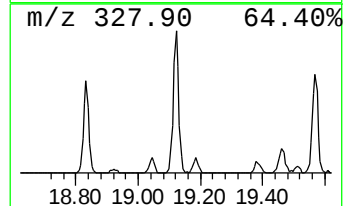
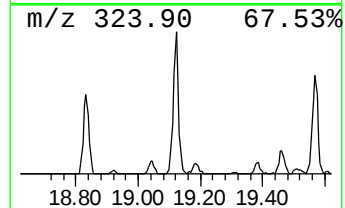
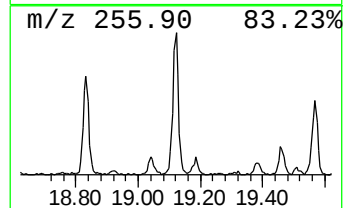
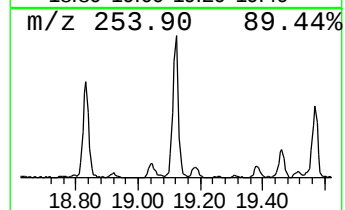
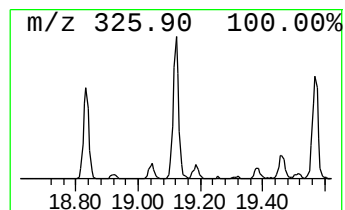
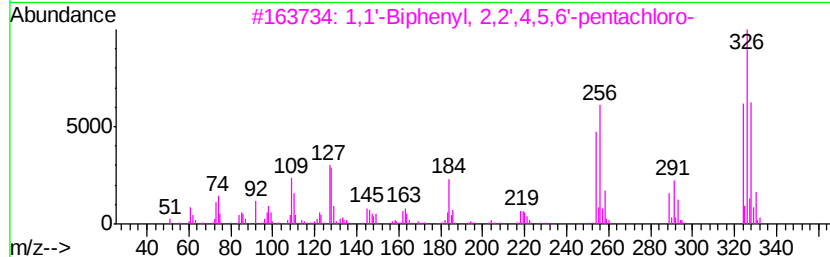
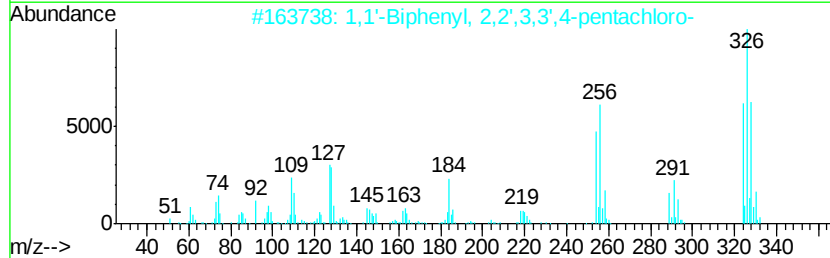
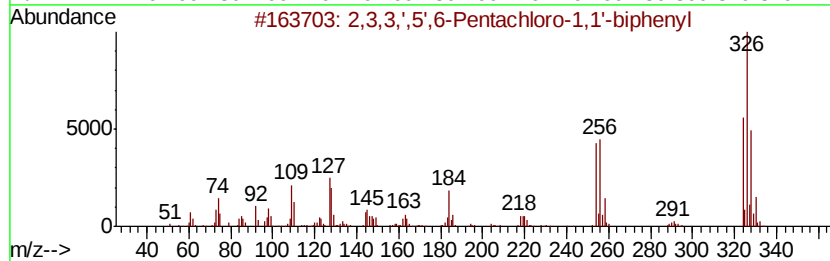
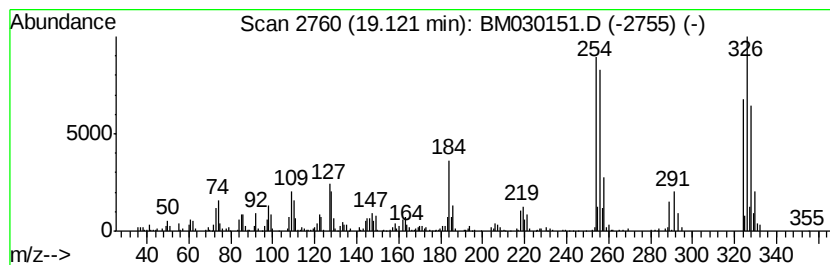
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 2,3,3',5',6-Pentachloro-1,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.12	2.87 ng/ul	334522	Chrysene-d12	21.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,3',5',6-Pentachloro-1,1'-bi...	324	C12H5Cl5	068194-10-5	99		
2	1,1'-Biphenyl, 2,2',3,3',4-penta...	324	C12H5Cl5	052663-62-4	98		
3	1,1'-Biphenyl, 2,2',4,5,6'-penta...	324	C12H5Cl5	068194-06-9	98		
4	2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	070362-41-3	98		
5	1,1'-Biphenyl, 2,2',4,4',5-penta...	324	C12H5Cl5	038380-01-7	98		



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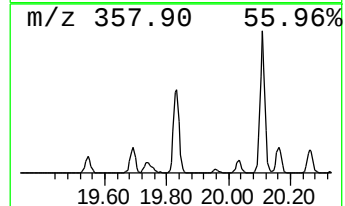
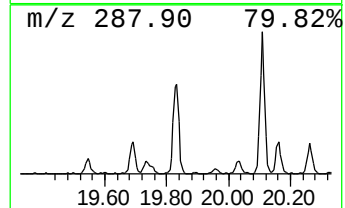
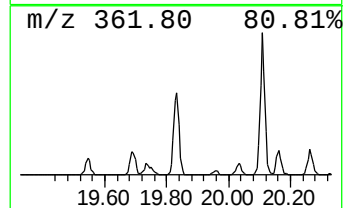
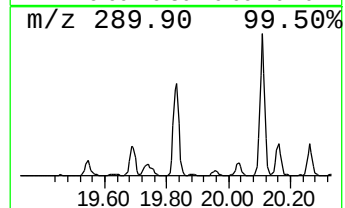
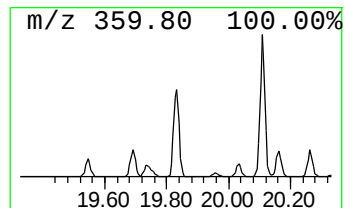
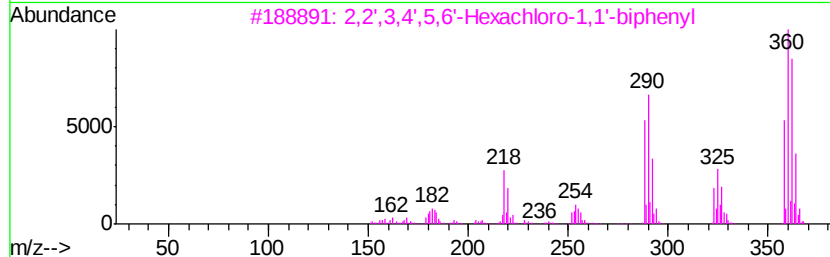
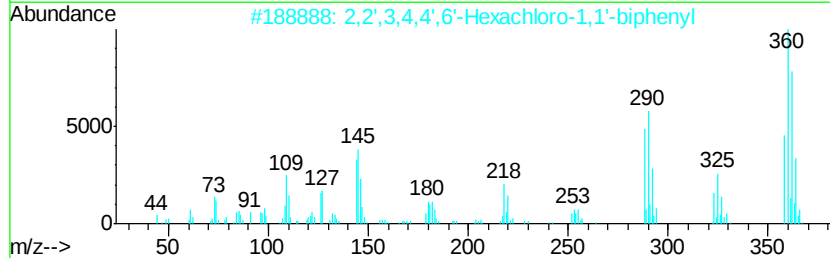
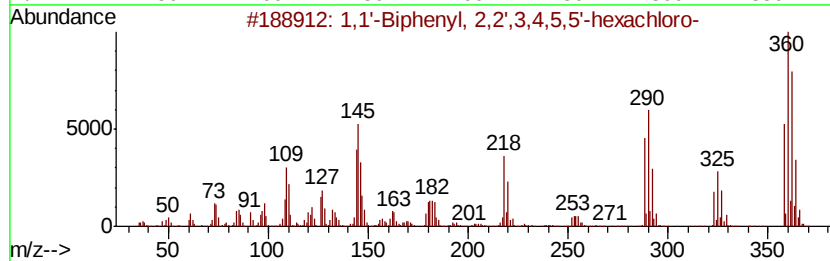
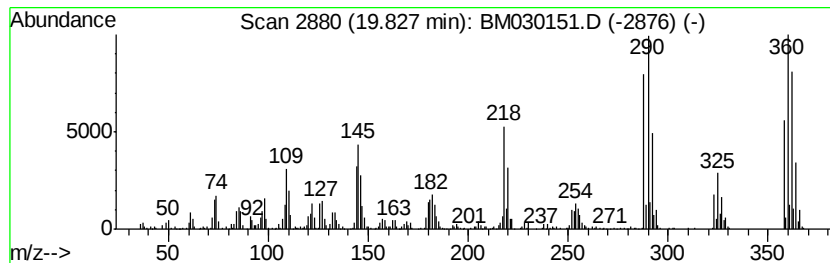
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.83	5.35 ng/ul	624362	Chrysene-d12	21.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',3,4,5,5'-hex...	358	C12H4Cl6	052712-04-6	99
2		2,2',3,4,4',6'-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-64-4	99
3		2,2',3,4',5,6'-Hexachloro-1,1'-b...	358	C12H4Cl6	074472-41-6	99
4		1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	99
5		1,1'-Biphenyl, 2,3',4,4',5,5'-he...	358	C12H4Cl6	052663-72-6	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052021\
Data File : BM030151.D
Acq On : 22 May 2021 15:39
Operator : CG/JU
Sample : M2455-01
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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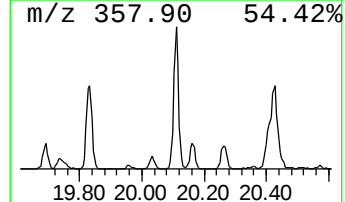
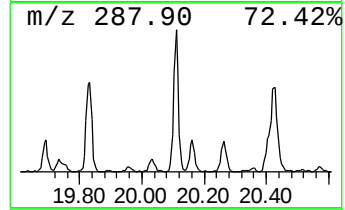
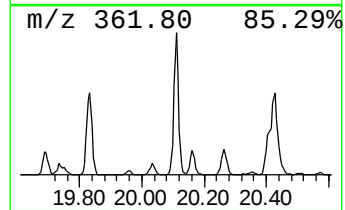
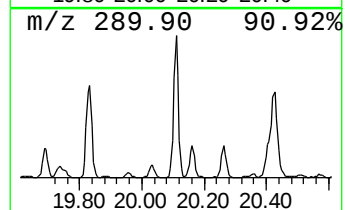
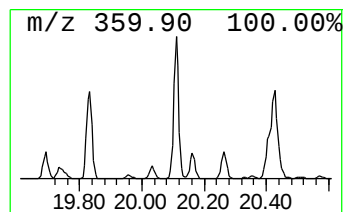
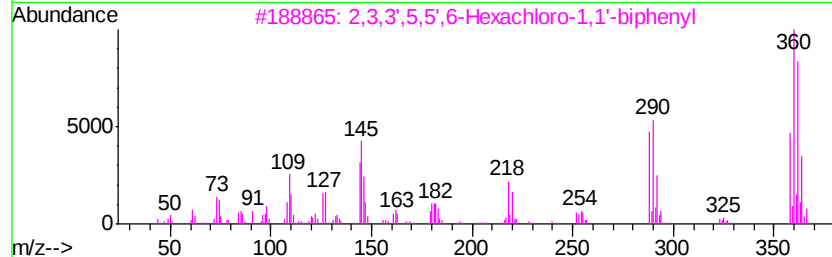
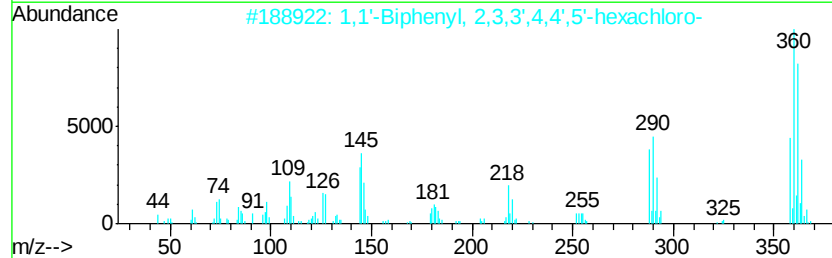
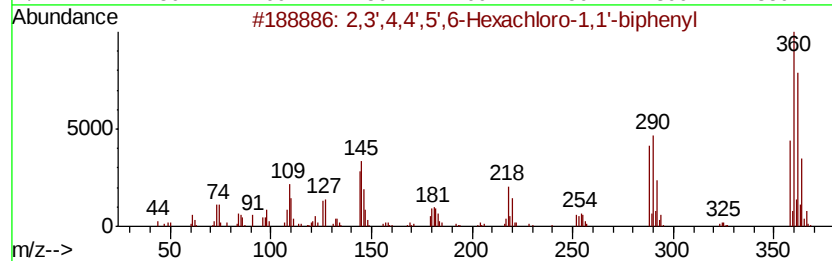
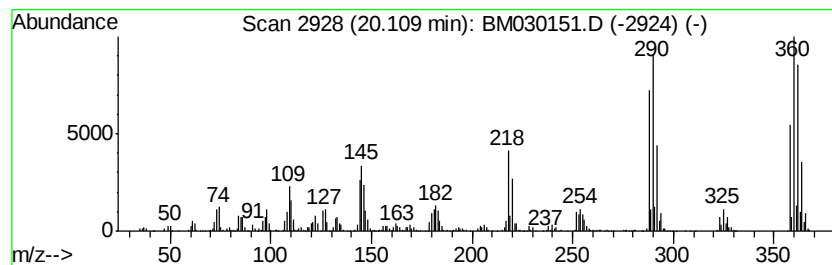
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.11	6.80 ng/ul	794304	Chrysene-d12	21.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,3',4,4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-65-5	99
2		1,1'-Biphenyl, 2,3,3',4,4',5'-he...	358	C12H4Cl6	069782-90-7	99
3		2,3,3',5,5',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-46-1	99
4		2,2',3,4,6,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-40-5	99
5		2,3,3',4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	074472-45-0	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052021\
Data File : BM030151.D
Acq On : 22 May 2021 15:39
Operator : CG/JU
Sample : M2455-01
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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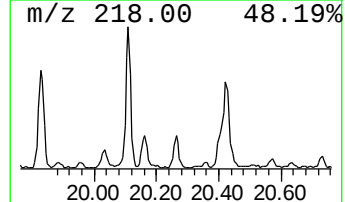
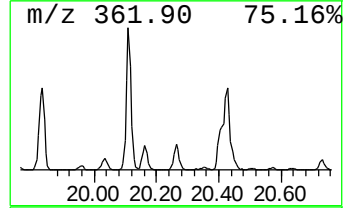
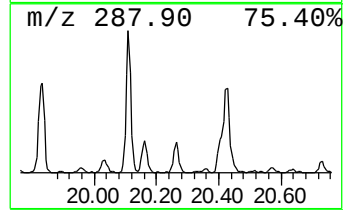
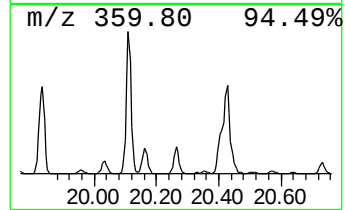
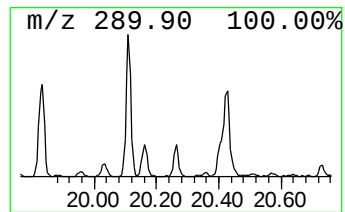
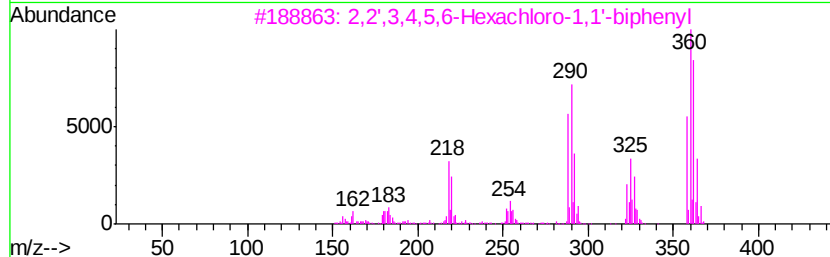
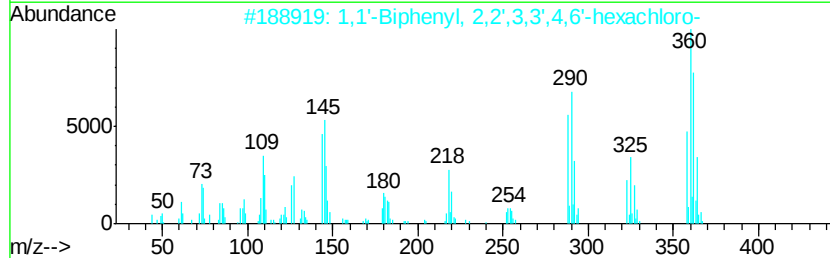
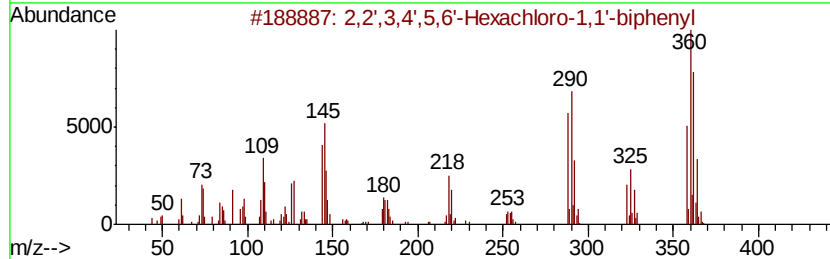
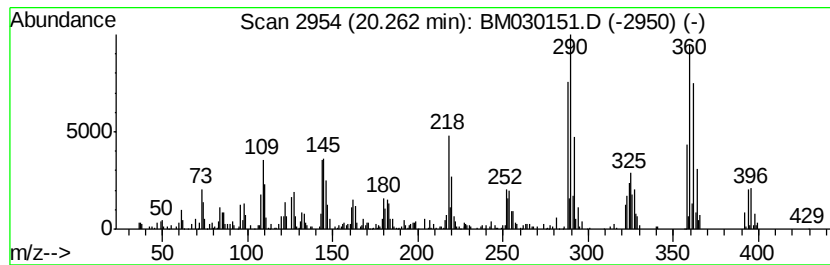
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 2,2',3,4',5,6'-Hexachloro-1... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.26	2.35 ng/ul	274123	Chrysene-d12	21.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2',3,4',5,6'-Hexachloro-1,1'-b...	358	C12H4Cl6	074472-41-6	99
2		1,1'-Biphenyl, 2,2',3,3',4,6'-he...	358	C12H4Cl6	038380-05-1	99
3		2,2',3,4,5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-61-4	99
4		1,1'-Biphenyl, 2,2',3,4,4',5'-he...	358	C12H4Cl6	035065-28-2	99
5		2,3,4,4',5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-63-6	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052021\
Data File : BM030151.D
Acq On : 22 May 2021 15:39
Operator : CG/JU
Sample : M2455-01
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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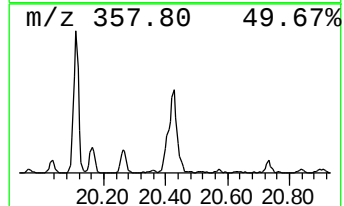
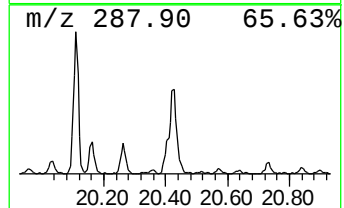
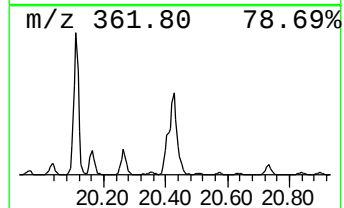
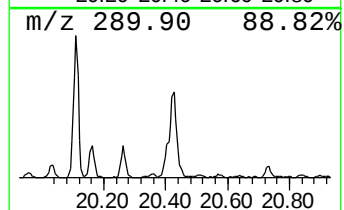
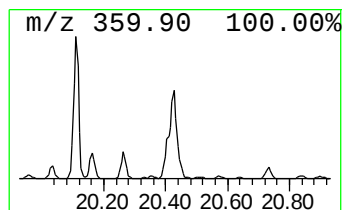
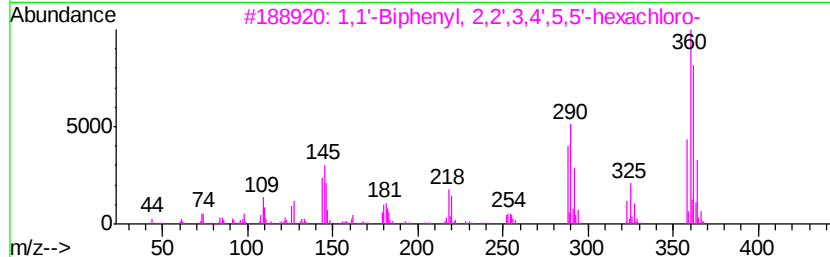
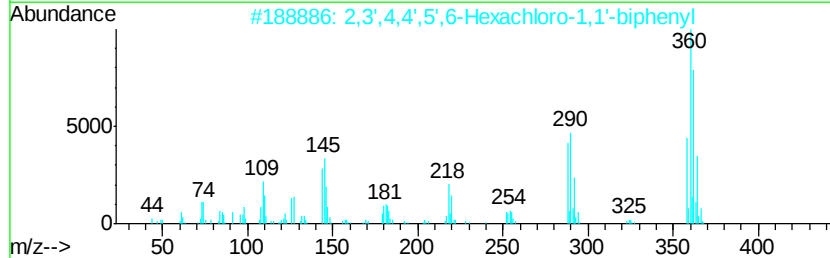
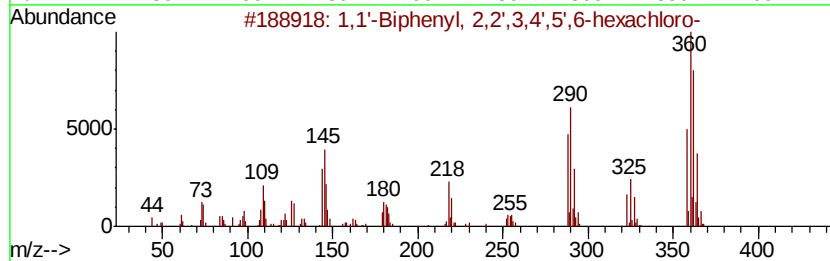
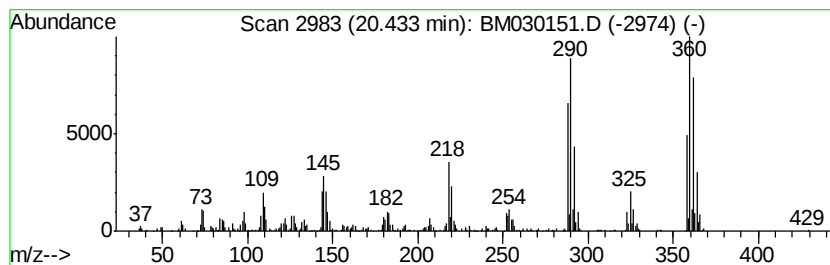
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.43	7.28 ng/ul	850315	Chrysene-d12	21.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',3,4',5',6-he...	358	C12H4Cl6	038380-04-0	99
2		2,3',4,4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-65-5	99
3		1,1'-Biphenyl, 2,2',3,4',5,5'-he...	358	C12H4Cl6	051908-16-8	99
4		1,1'-Biphenyl, 2,2',3,3',4,6-Hex...	358	C12H4Cl6	061798-70-7	99
5		2,2',3,4,4',6'-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-64-4	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052021\
Data File : BM030151.D
Acq On : 22 May 2021 15:39
Operator : CG/JU
Sample : M2455-01
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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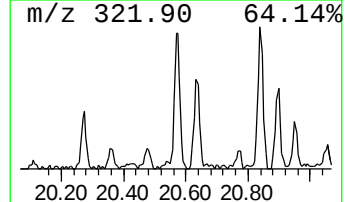
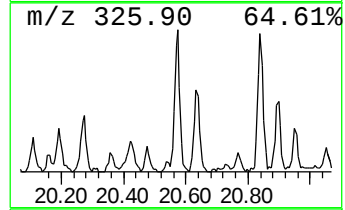
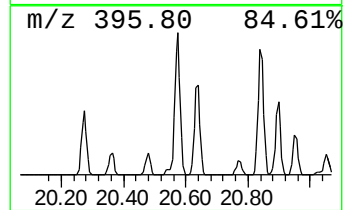
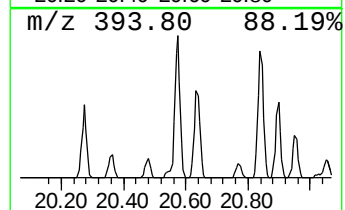
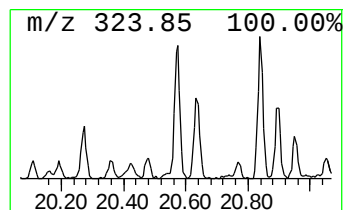
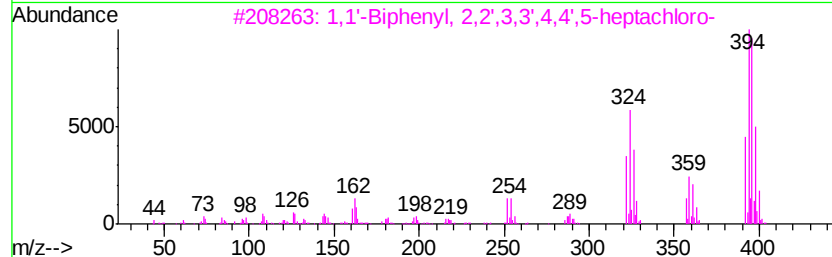
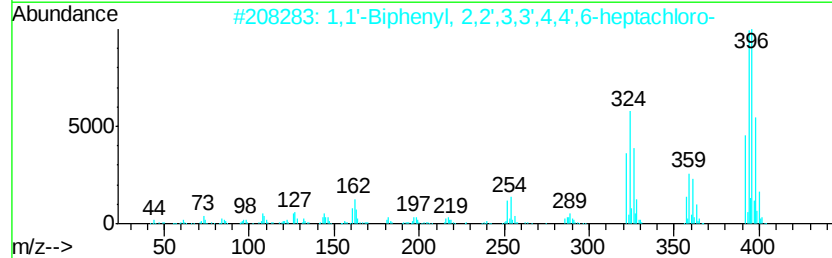
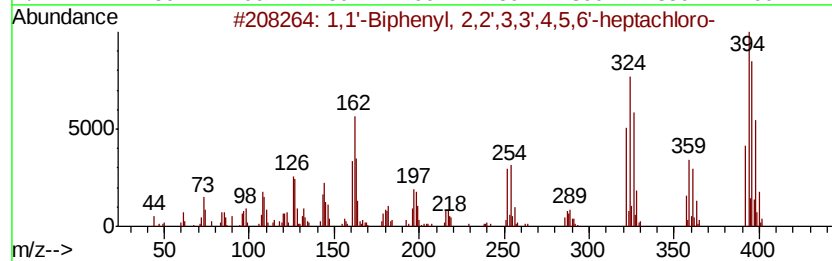
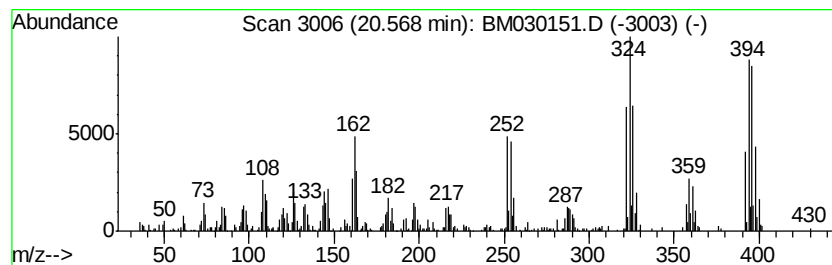
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1,1'-Biphenyl, 2,2',3,3',4,... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.57	2.34 ng/ul	273181	Chrysene-d12	21.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',3,3',4,5,6'-...	392	C12H3Cl7	038411-25-5	99
2		1,1'-Biphenyl, 2,2',3,3',4,4',6-...	392	C12H3Cl7	052663-71-5	99
3		1,1'-Biphenyl, 2,2',3,3',4,4',5-...	392	C12H3Cl7	035065-30-6	99
4		2,2',3,4',5,6,6'-Heptachloro-1,1...	392	C12H3Cl7	074487-85-7	99
5		2,2',3,4,4',6,6'-Heptachloro-1,1...	392	C12H3Cl7	074472-48-3	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052021\
Data File : BM030151.D
Acq On : 22 May 2021 15:39
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Sample : M2455-01
Misc :
ALS Vial : 67 Sample Multiplier: 1

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ClientSampled :
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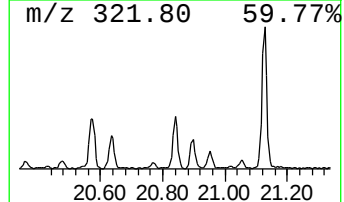
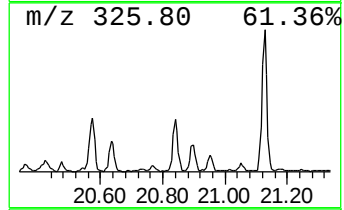
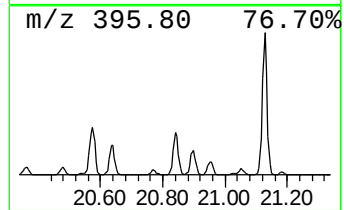
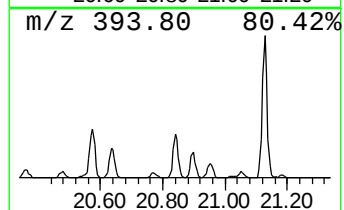
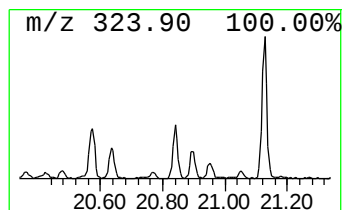
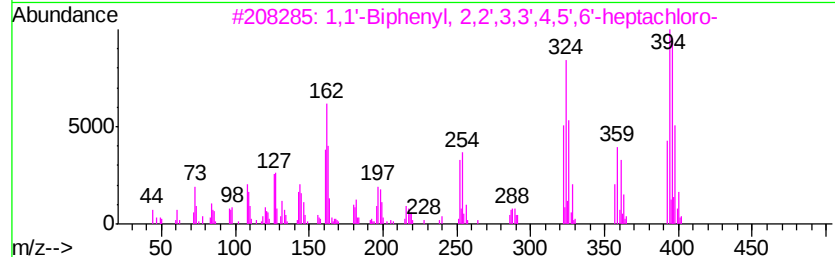
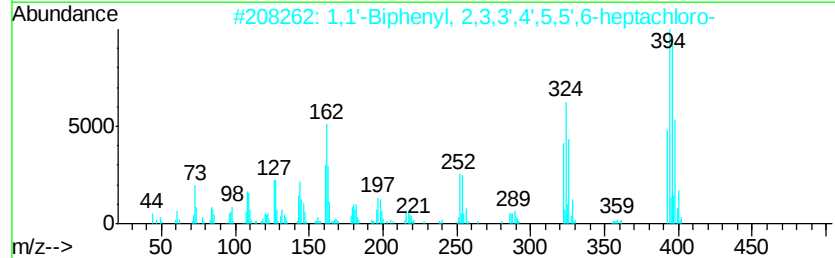
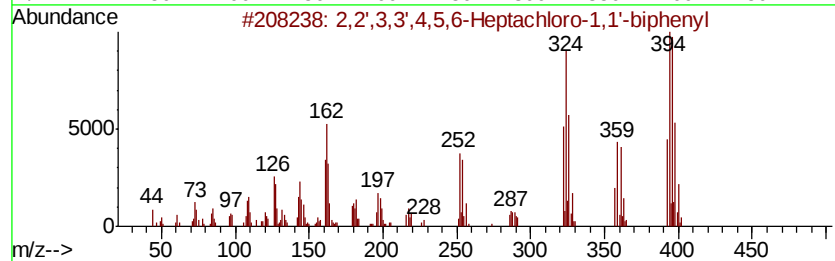
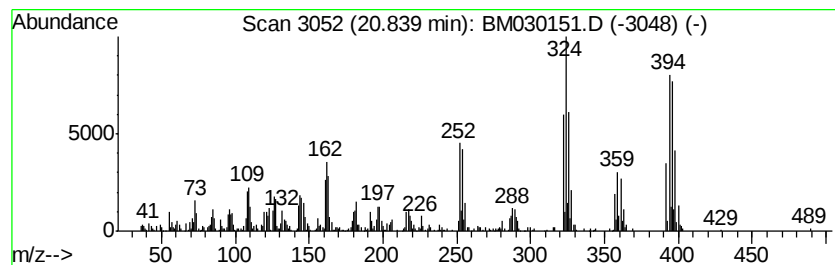
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 2,2',3,3',4,5,6-Heptachloro... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.84	2.67 ng/ul	312159	Chrysene-d12	21.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,3',4,5,6-Heptachloro-1,1'...	392	C12H3Cl7	068194-16-1	99		
2	1,1'-Biphenyl, 2,3,3',4',5,5',6-...	392	C12H3Cl7	069782-91-8	99		
3	1,1'-Biphenyl, 2,2',3,3',4,5',6'...	392	C12H3Cl7	052663-70-4	99		
4	2,3,3',4,4',5',6-Heptachloro-1,1...	392	C12H3Cl7	074472-50-7	99		
5	1,1'-Biphenyl, 2,2',3,4,5,6,6'-H...	392	C12H3Cl7	074472-49-4	99		



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052021\
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Acq On : 22 May 2021 15:39
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Sample : M2455-01
Misc :
ALS Vial : 67 Sample Multiplier: 1

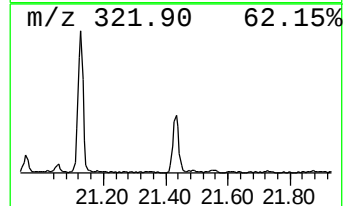
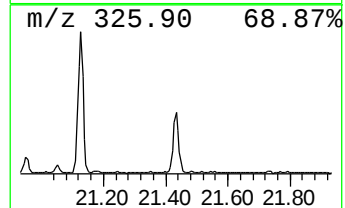
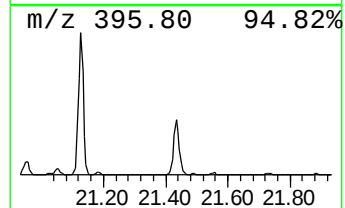
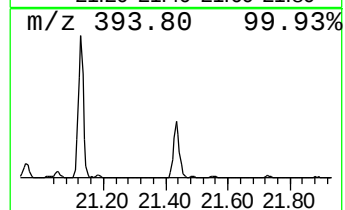
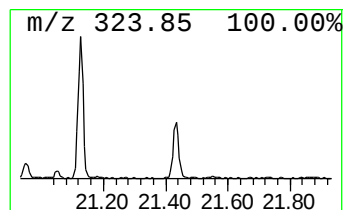
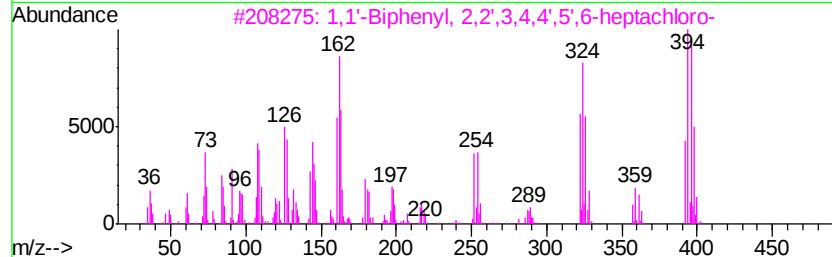
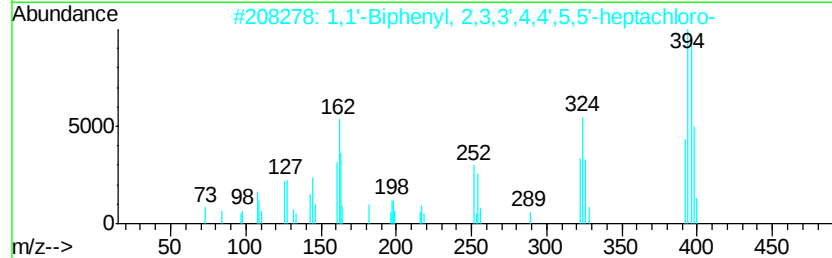
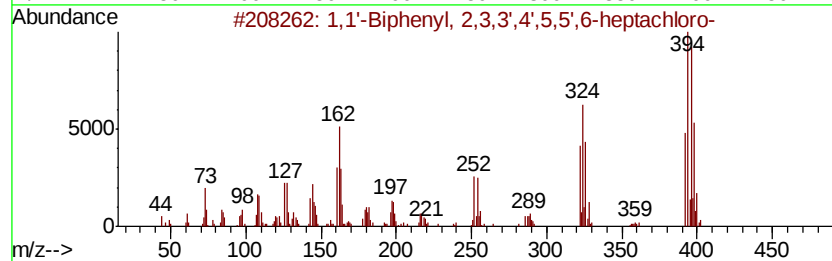
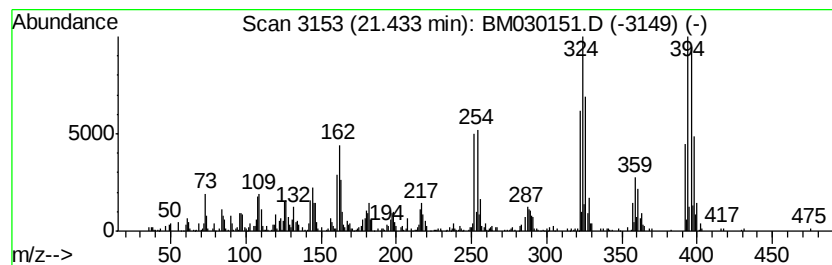
Instrument :
BNA_M
ClientSampled :
C0BA1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050721.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 1,1'-Biphenyl, 2,3,3',4',5,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
21.43	2.99 ng/ul	348637	Chrysene-d12	21.10		
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	1,1'-Biphenyl,	2,3,3',4,4',5,5'-...	392	C12H3Cl7	039635-31-9	99
3	1,1'-Biphenyl,	2,2',3,4,4',5',6-...	392	C12H3Cl7	052663-69-1	99
4	1,1'-Biphenyl,	2,3,3',4,5,5',6-h...	392	C12H3Cl7	074472-51-8	99
5	1,1'-Biphenyl,	2,2',3,4,5,5',6-h...	392	C12H3Cl7	052712-05-7	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052021\
 Data File : BM030151.D
 Acq On : 22 May 2021 15:39
 Operator : CG/JU
 Sample : M2455-01
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0BA1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050721.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Str...	2.94	73.0	ng/ul	3289720	1	7.50	901886	20.0
(DEL) Alkane: Cyc...	3.26	2.2	ng/ul	97958	1	7.50	901886	20.0
1,1'-Biphenyl, 2,...	18.83	3.5	ng/ul	346340	4	16.91	1971920	20.0
2,3,3',5',6-Pent...	19.12	2.9	ng/ul	334522	5	21.10	2334600	20.0
1,1'-Biphenyl, 2,...	19.83	5.3	ng/ul	624362	5	21.10	2334600	20.0
2,3',4,4',5',6-He...	20.11	6.8	ng/ul	794304	5	21.10	2334600	20.0
2,2',3,4',5,6'-He...	20.26	2.4	ng/ul	274123	5	21.10	2334600	20.0
1,1'-Biphenyl, 2,...	20.43	7.3	ng/ul	850315	5	21.10	2334600	20.0
1,1'-Biphenyl, 2,...	20.57	2.3	ng/ul	273181	5	21.10	2334600	20.0
2,2',3,3',4,5,6-H...	20.84	2.7	ng/ul	312159	5	21.10	2334600	20.0
1,1'-Biphenyl, 2,...	21.43	3.0	ng/ul	348637	5	21.10	2334600	20.0