

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
 Data File : BM029855.D
 Acq On : 06 May 2021 08:55
 Operator : CG/JU
 Sample : M2240-01
 Misc :
 ALS Vial : 74 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B64

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-BM050421.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	2.987	14	17	20	rBV	207942	214480	7.06%	0.579%
2	7.563	791	795	812	rVB	354217	696703	22.95%	1.880%
3	10.339	1261	1267	1280	rVB	436053	730782	24.07%	1.972%
4	12.074	1560	1562	1566	rVB	2624	3584	0.12%	0.010%
5	12.280	1595	1597	1601	rVB3	3197	3125	0.10%	0.008%
6	13.939	1875	1879	1882	rBV6	2002	3453	0.11%	0.009%
7	14.074	1898	1902	1903	rBV3	2104	3045	0.10%	0.008%
8	14.210	1918	1925	1941	rBV	786020	1104085	36.37%	2.980%
9	14.586	1985	1989	1993	rBV7	2172	3607	0.12%	0.010%
10	15.004	2055	2060	2061	rVB5	3377	3822	0.13%	0.010%
11	15.092	2070	2075	2077	rBV5	4673	7379	0.24%	0.020%
12	15.562	2152	2155	2156	rBV3	4654	5247	0.17%	0.014%
13	16.498	2310	2314	2317	rBV5	15178	22462	0.74%	0.061%
14	16.957	2386	2392	2408	rBV	833236	1378660	45.41%	3.721%
15	17.139	2419	2423	2428	rBV2	58526	81867	2.70%	0.221%
16	17.562	2490	2495	2502	rVB	367453	509364	16.78%	1.375%
17	17.674	2510	2514	2521	rVB4	70305	108090	3.56%	0.292%
18	17.809	2533	2537	2541	rBV	108602	147264	4.85%	0.397%
19	17.856	2542	2545	2550	rVB3	73759	99139	3.27%	0.268%
20	18.027	2569	2574	2580	rBV	433898	568384	18.72%	1.534%
21	18.086	2580	2584	2588	rVV	245822	320044	10.54%	0.864%
22	18.127	2588	2591	2596	rVB2	106795	140496	4.63%	0.379%
23	18.315	2618	2623	2627	rBV	293304	406292	13.38%	1.097%
24	18.362	2628	2631	2638	rVB3	92540	131060	4.32%	0.354%
25	18.451	2644	2646	2649	rBV	52484	62778	2.07%	0.169%
26	18.492	2649	2653	2657	rVB	202364	258206	8.50%	0.697%
27	18.592	2667	2670	2677	rVB3	52484	73303	2.41%	0.198%
28	18.803	2702	2706	2710	rBV	166487	206320	6.80%	0.557%
29	18.880	2711	2719	2727	rVB3	818076	1446708	47.65%	3.904%
30	19.015	2740	2742	2750	rVB	56671	81796	2.69%	0.221%
31	19.103	2751	2757	2763	rVV2	182380	459051	15.12%	1.239%
32	19.168	2763	2768	2773	rVV	1040092	1347746	44.39%	3.637%
33	19.227	2775	2778	2783	rVB	108967	145107	4.78%	0.392%
34	19.303	2788	2791	2796	rVB2	65554	79974	2.63%	0.216%

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
 Data File : BM029855.D
 Acq On : 06 May 2021 08:55
 Operator : CG/JU
 Sample : M2240-01
 Misc :
 ALS Vial : 74 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B64

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-BM050421.M
 Title : SVOA CALIBRATION

35	19.427	2808	2812	2818	rVB2	120091	153895	5.07%	0.415%
36	19.503	2821	2825	2830	rVB	221394	274508	9.04%	0.741%
37	19.556	2831	2834	2836	rVV2	48707	61216	2.02%	0.165%
38	19.609	2836	2843	2849	rVB2	620106	1101178	36.27%	2.972%
39	19.733	2860	2864	2868	rBV	543388	685521	22.58%	1.850%
40	19.780	2868	2872	2879	rVV	234760	417313	13.75%	1.126%
41	19.874	2883	2888	2893	rBV	1849039	2219261	73.10%	5.989%
42	19.927	2893	2897	2905	rVB	375133	484561	15.96%	1.308%
43	20.003	2906	2910	2914	rVB3	93273	111782	3.68%	0.302%
44	20.074	2918	2922	2928	rVB	238081	312312	10.29%	0.843%
45	20.156	2931	2936	2940	rBV	2527381	2982269	98.23%	8.049%
46	20.203	2940	2944	2955	rVB2	630085	929732	30.62%	2.509%
47	20.309	2957	2962	2967	rBV2	690028	987148	32.51%	2.664%
48	20.403	2974	2978	2982	rBV3	175747	226151	7.45%	0.610%
49	20.468	2982	2989	2996	rVV	1711533	3036075	100.00%	8.194%
50	20.521	2996	2998	3001	rVB	111519	103878	3.42%	0.280%
51	20.556	3001	3004	3007	rBV4	60045	60774	2.00%	0.164%
52	20.615	3010	3014	3020	rVB	874944	1008395	33.21%	2.722%
53	20.680	3021	3025	3033	rVB	503470	597867	19.69%	1.614%
54	20.774	3037	3041	3045	rBV	232044	265367	8.74%	0.716%
55	20.815	3045	3048	3054	rVB6	84244	100578	3.31%	0.271%
56	20.886	3055	3060	3065	rBV2	795545	995552	32.79%	2.687%
57	20.939	3065	3069	3074	rBV	458275	551948	18.18%	1.490%
58	20.992	3075	3078	3081	rBV2	238584	306517	10.10%	0.827%
59	21.021	3081	3083	3089	rVB2	179474	213166	7.02%	0.575%
60	21.092	3093	3095	3099	rVB3	101602	110455	3.64%	0.298%
61	21.139	3099	3103	3105	rBV	920652	1143301	37.66%	3.086%
62	21.174	3105	3109	3114	rVB	2190430	2768736	91.19%	7.472%
63	21.309	3129	3132	3136	rVB2	94986	102504	3.38%	0.277%
64	21.474	3156	3160	3166	rBV2	993385	1285162	42.33%	3.468%
65	21.533	3166	3170	3175	rVV3	259469	320002	10.54%	0.864%
66	21.597	3177	3181	3187	rVV	352746	413012	13.60%	1.115%
67	21.927	3235	3237	3241	rBV2	90964	120717	3.98%	0.326%
68	22.144	3270	3274	3277	rBV2	266167	375906	12.38%	1.015%
69	23.297	3465	3470	3482	rVB2	721982	1372528	45.21%	3.704%

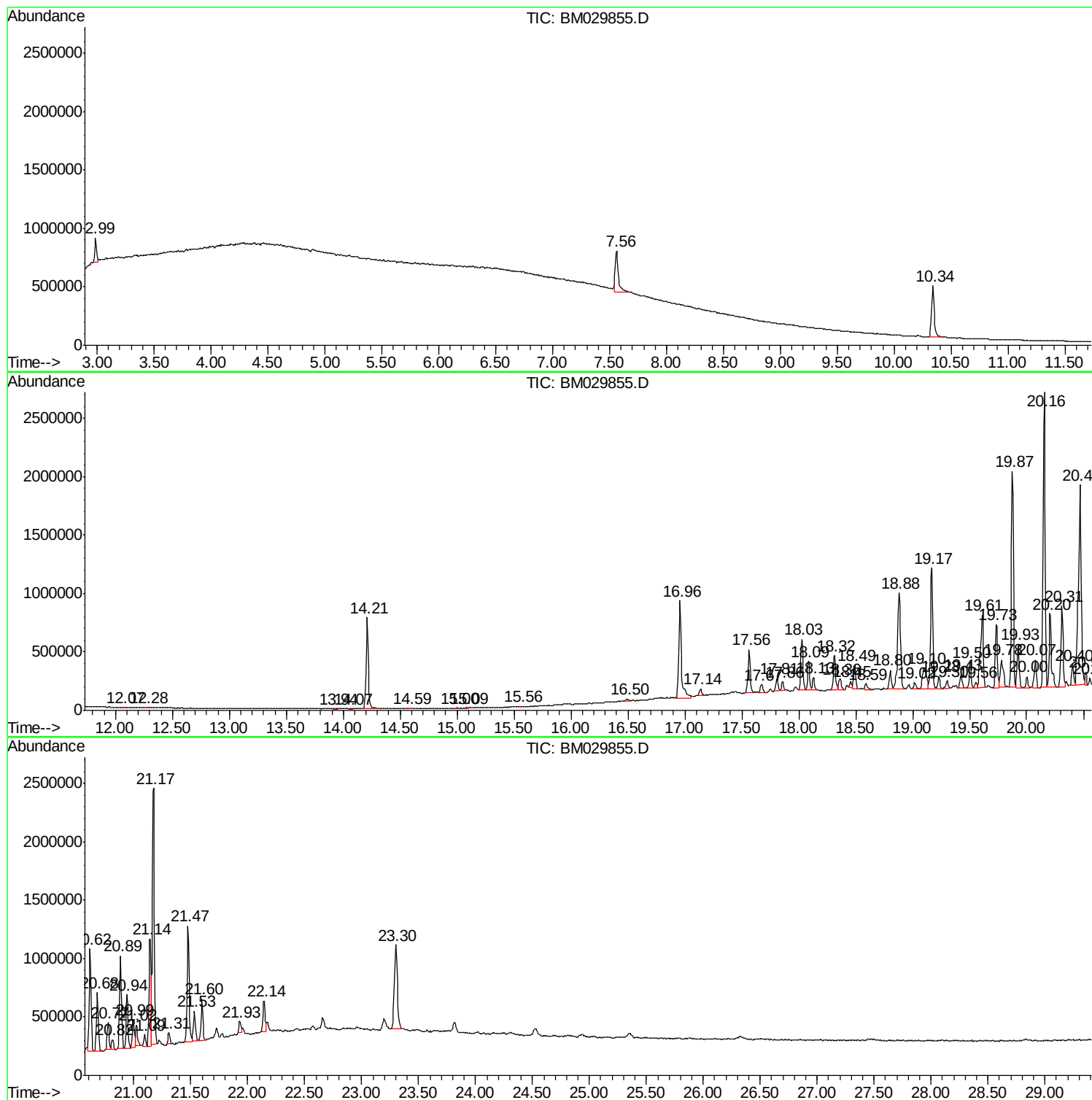
Sum of corrected areas: 37052710

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

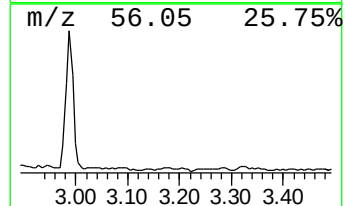
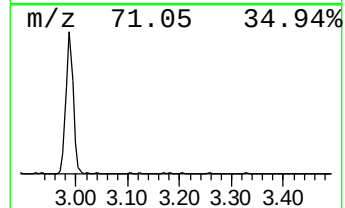
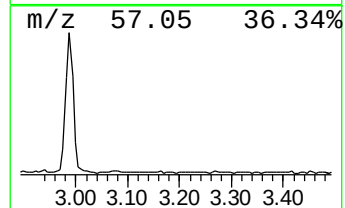
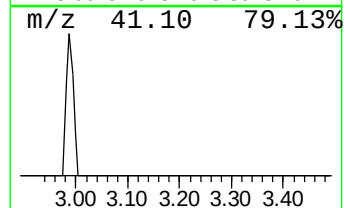
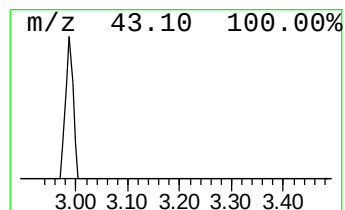
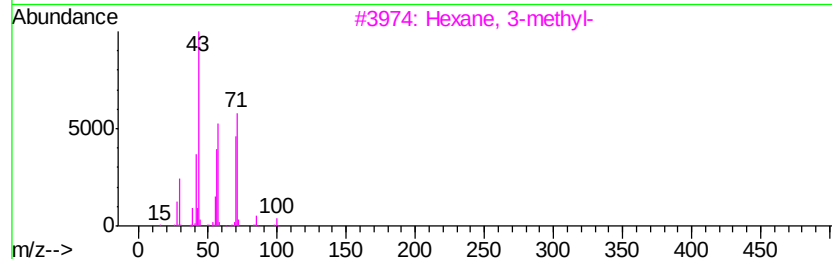
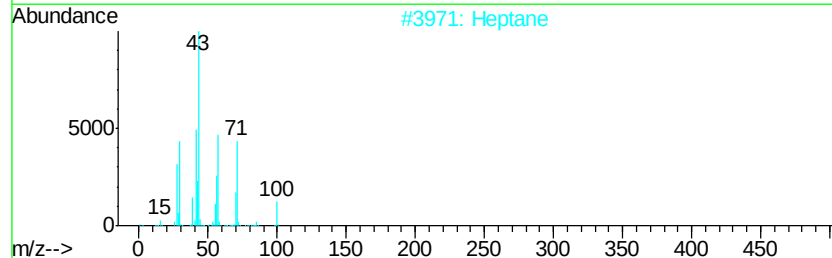
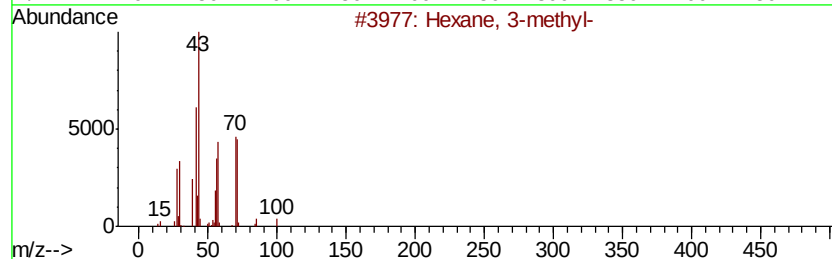
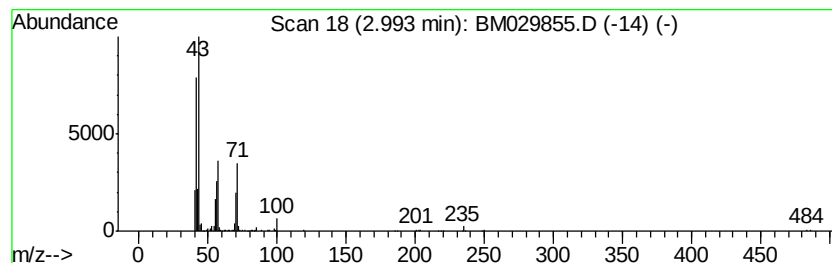
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.99	6.16 ng/ul	214480	1,4-Dichlorobenzene-d4	7.56

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 3-methyl-			100	C7H16	000589-34-4	52
2	Heptane			100	C7H16	000142-82-5	50
3	Hexane, 3-methyl-			100	C7H16	000589-34-4	50
4	Heptane			100	C7H16	000142-82-5	50
5	Heptane			100	C7H16	000142-82-5	40



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

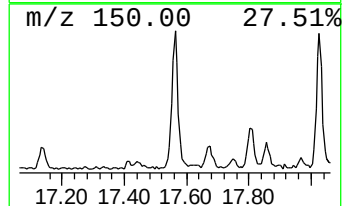
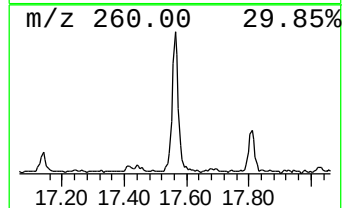
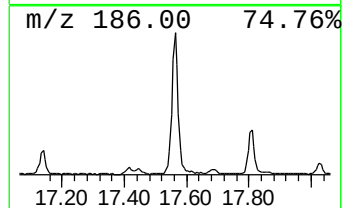
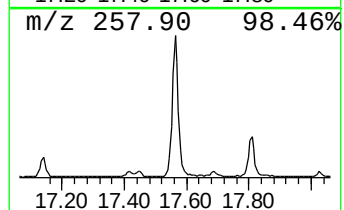
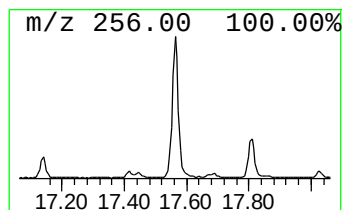
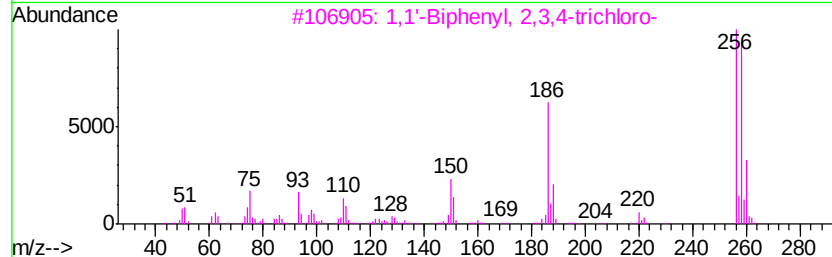
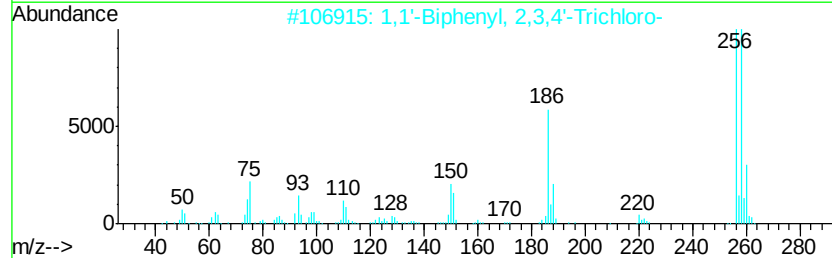
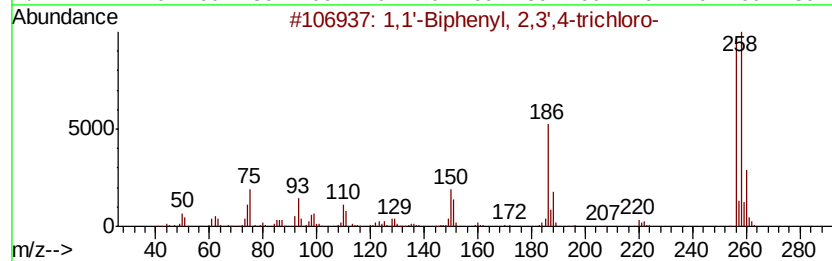
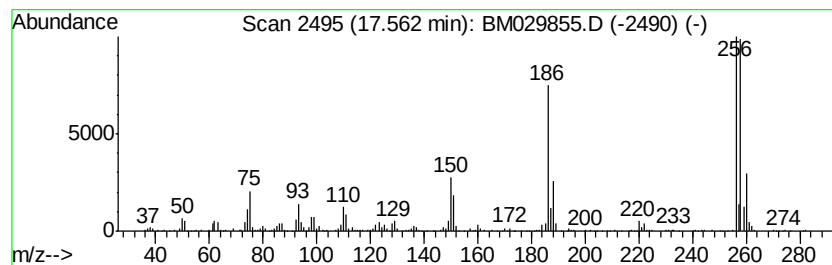
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1,1'-Biphenyl, 2,3',4-trich... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.56	7.39 ng/ul	509364	Phenanthrene-d10	16.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3',4-trichloro-	256	C12H7Cl3	055712-37-3	99
2		1,1'-Biphenyl, 2,3,4'-Trichloro-	256	C12H7Cl3	038444-85-8	99
3		1,1'-Biphenyl, 2,3,4-trichloro-	256	C12H7Cl3	055702-46-0	99
4		1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99
5		1,1'-Biphenyl, 2,4,6-trichloro-	256	C12H7Cl3	035693-92-6	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

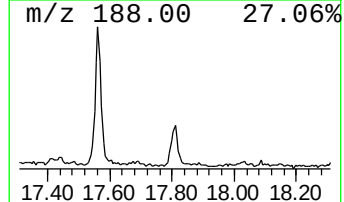
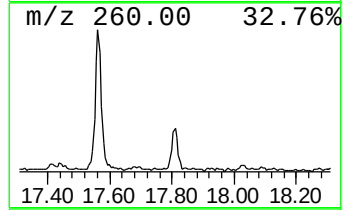
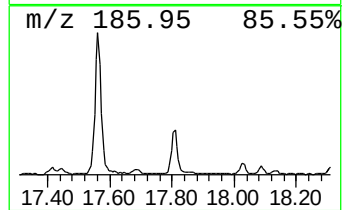
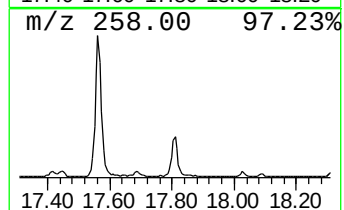
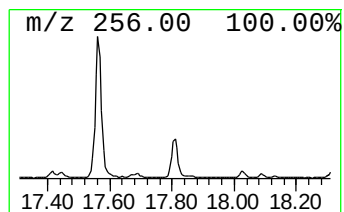
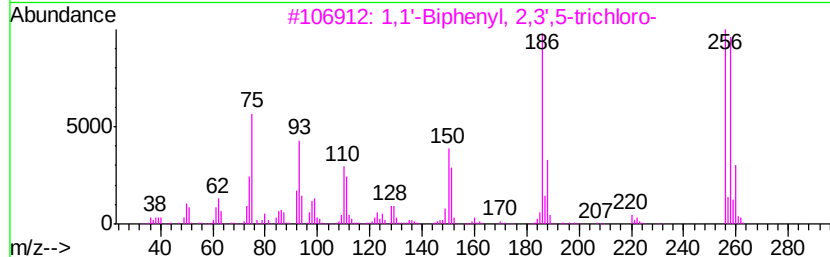
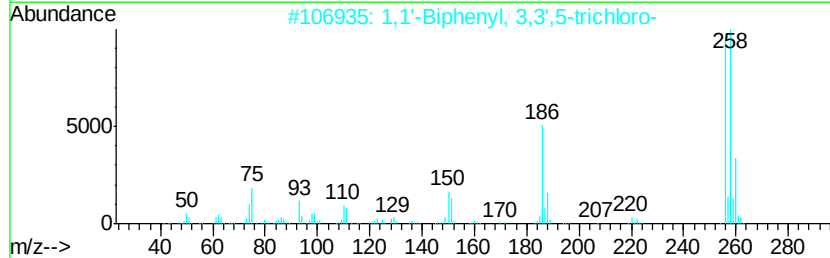
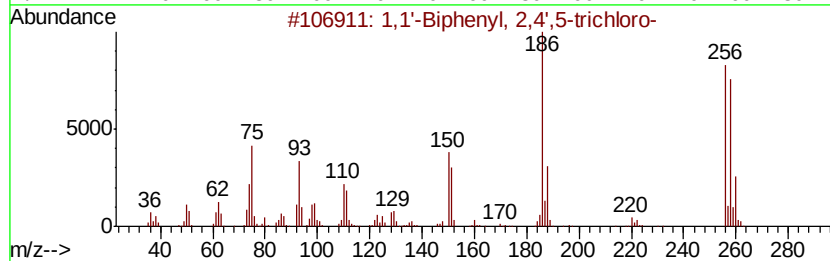
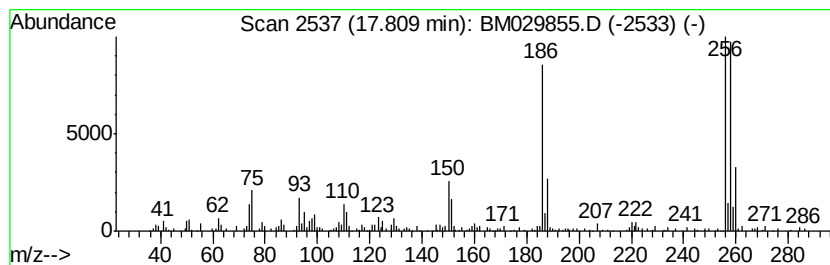
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1,1'-Biphenyl, 2,4',5-trich... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.81	2.14 ng/ul	147264	Phenanthrene-d10	16.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	96
2		1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	96
3		1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	95
4		1,1'-Biphenyl, 2,4,6-trichloro-	256	C12H7Cl3	035693-92-6	95
5		2,4',6-Trichloro-1,1'-biphenyl	256	C12H7Cl3	038444-77-8	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

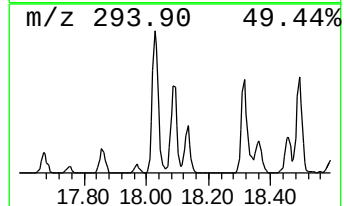
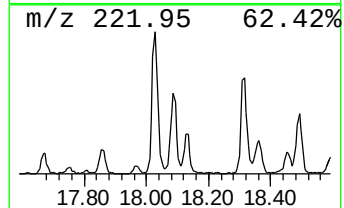
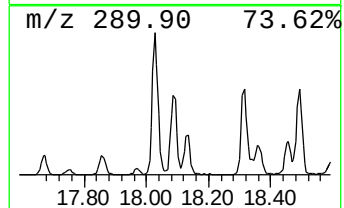
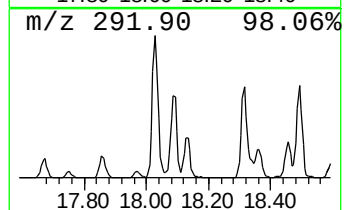
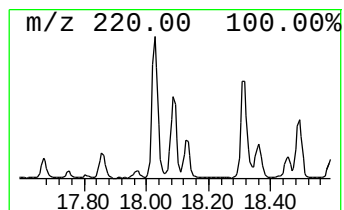
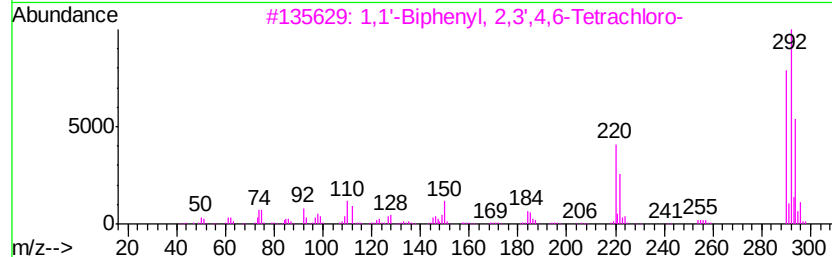
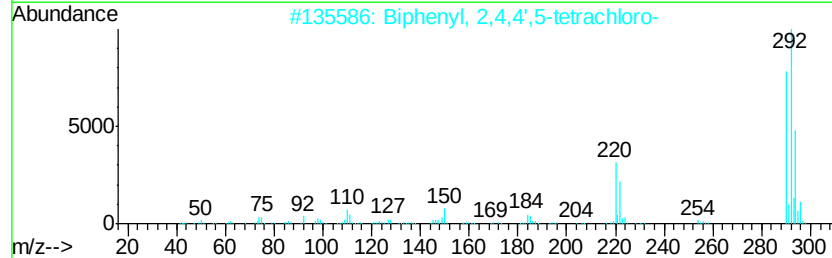
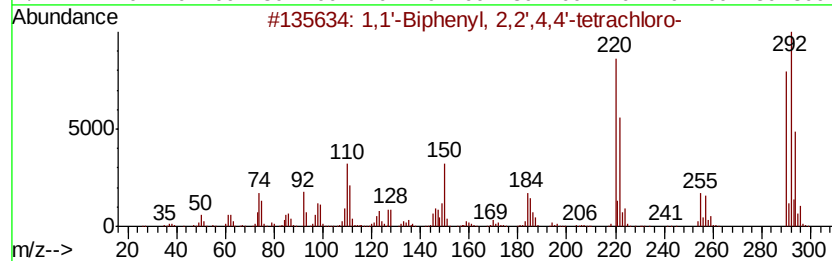
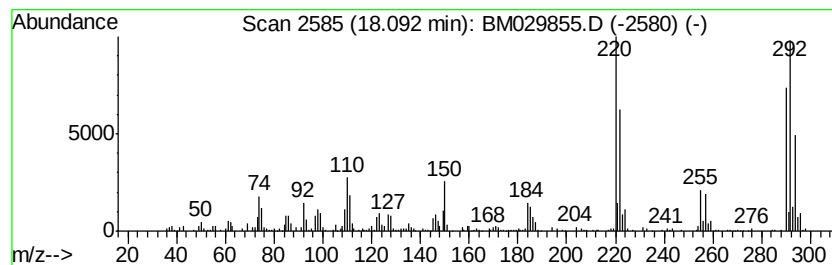
Instrument :
BNA_M
ClientSampleId :
C0B64

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,1'-Biphenyl, 2,2',4,4'-te... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.09	4.64 ng/ul	320044	Phenanthrene-d10	16.96	
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	Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	99
3	1,1'-Biphenyl, 2,3',4,6-Tetrachl...	290	C12H6Cl4	060233-24-1	99
4	2,3,3',6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074472-33-6	99
5	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

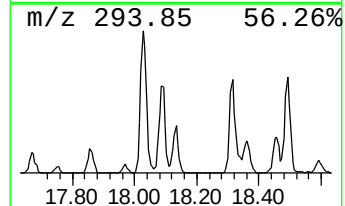
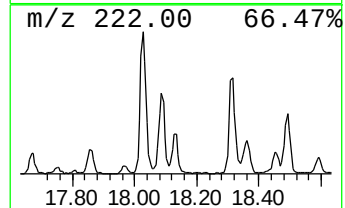
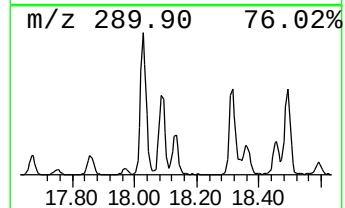
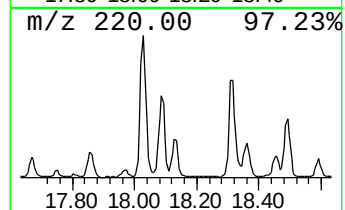
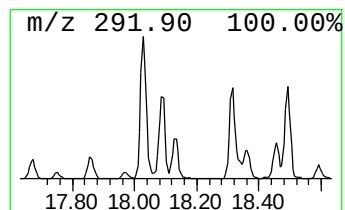
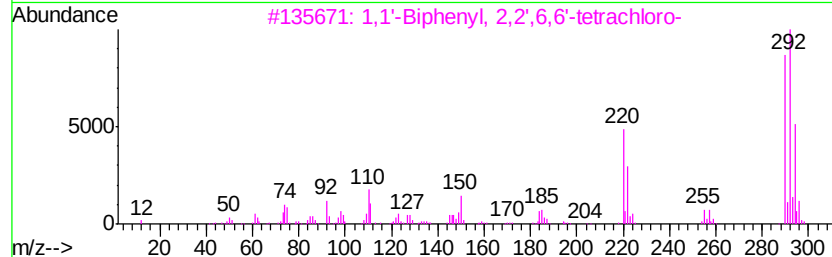
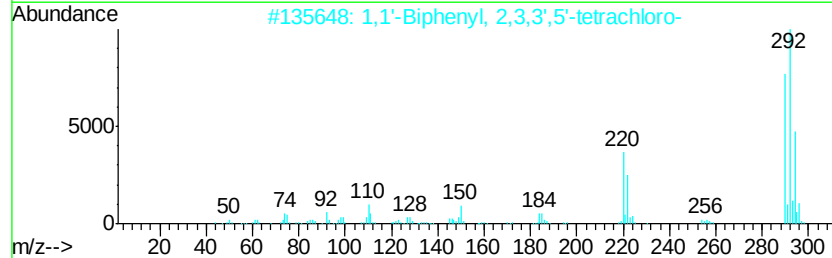
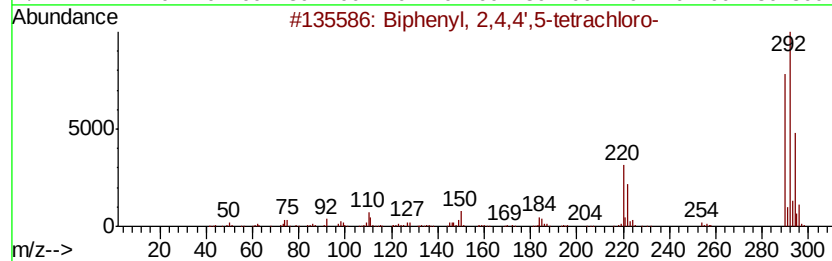
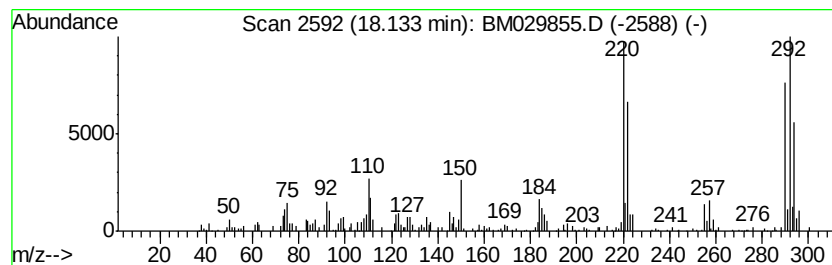
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Biphenyl, 2,4,4',5-tetrachl... Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.13	2.04 ng/ul	140496	Phenanthrene-d10	16.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	99
2			1,1'-Biphenyl, 2,3,3',5'-tetrach...	290	C12H6Cl4	041464-49-7	99
3			1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	98
4			1,1'-Biphenyl, 2,4,4',6-tetrachl...	290	C12H6Cl4	032598-12-2	98
5			1,1'-Biphenyl, 2,3',5,5'-tetrach...	290	C12H6Cl4	041464-42-0	98



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

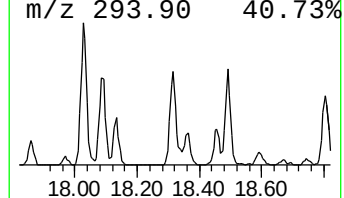
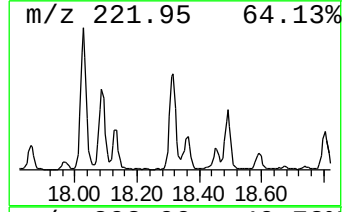
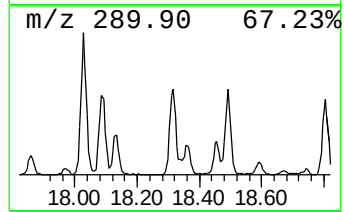
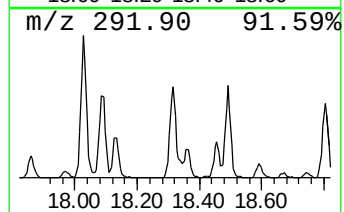
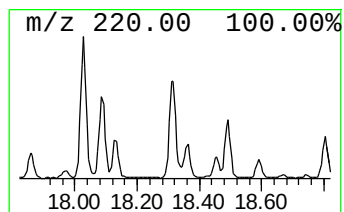
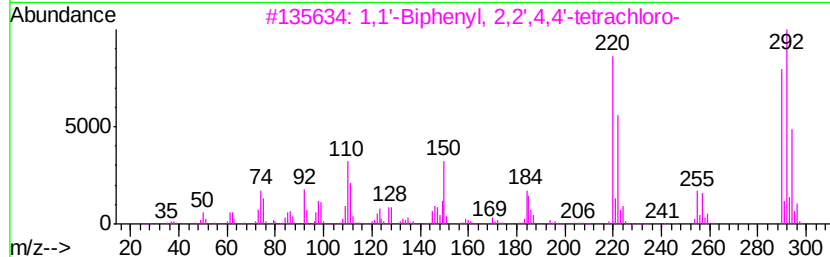
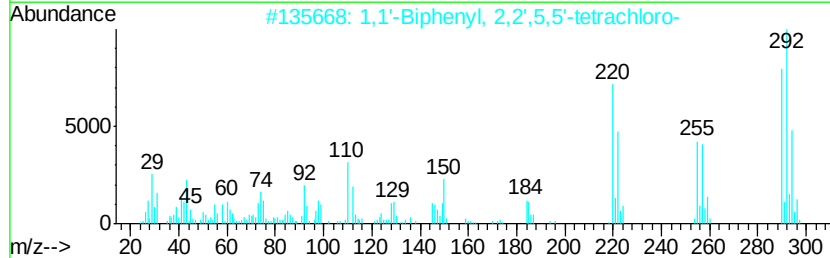
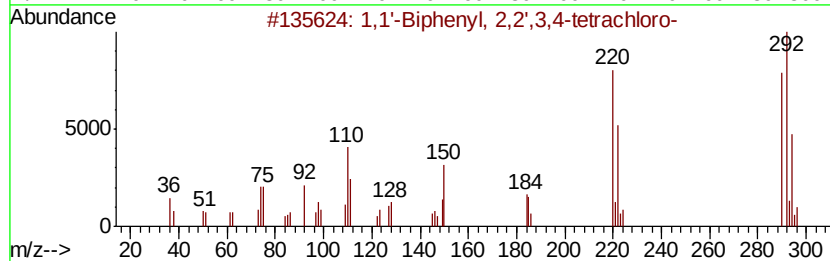
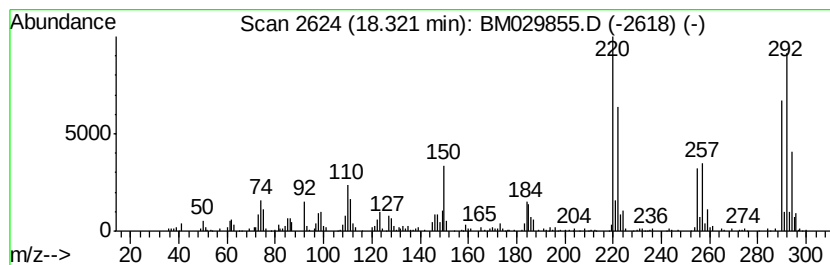
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 1,1'-Biphenyl, 2,2',3,4-tet... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.32	5.89 ng/ul	406292	Phenanthrene-d10	16.96

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',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	99
3		1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
4		2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99
5		2,2',3,6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-45-7	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

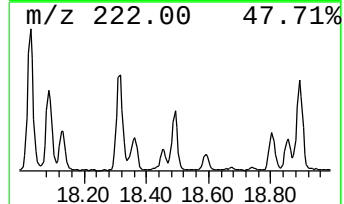
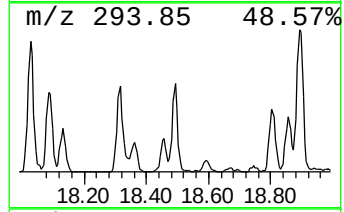
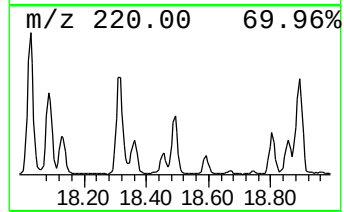
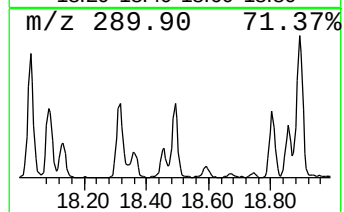
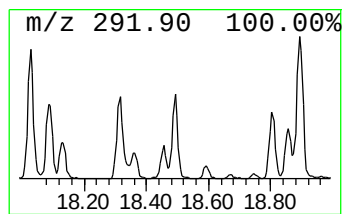
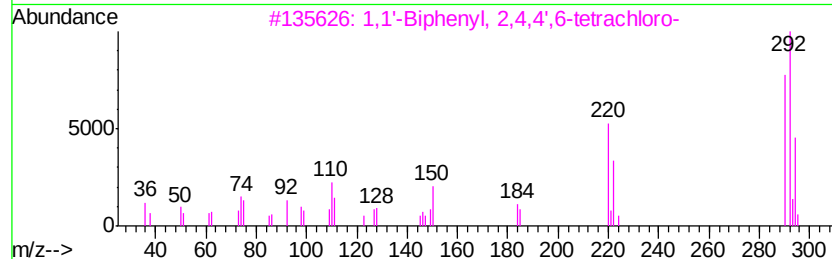
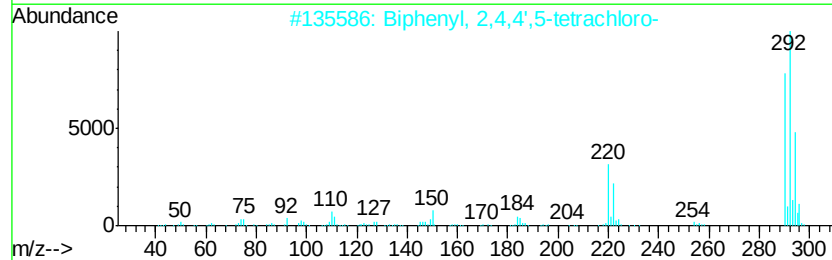
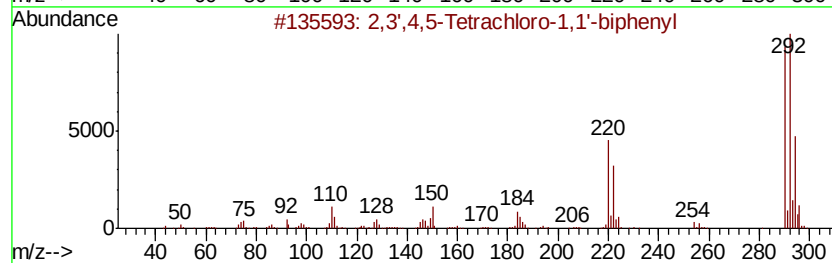
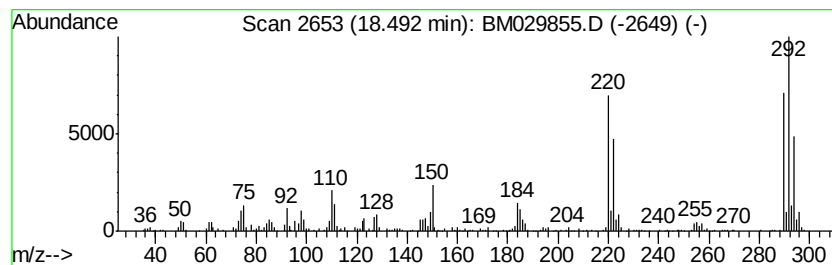
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 2,3',4,5-Tetrachloro-1,1'-b... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.49	3.75 ng/ul	258206	Phenanthrene-d10	16.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-53-8	99
2		Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	99
3		1,1'-Biphenyl, 2,4,4',6-tetrachl...	290	C12H6Cl4	032598-12-2	99
4		1,1'-Biphenyl, 2,3',5,5'-tetrach...	290	C12H6Cl4	041464-42-0	99
5		1,1'-Biphenyl, 3,3',5,5'-tetrach...	290	C12H6Cl4	033284-52-5	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

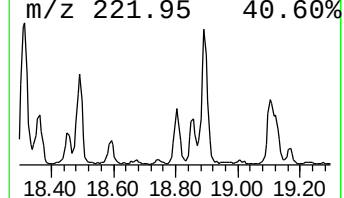
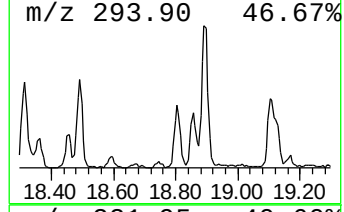
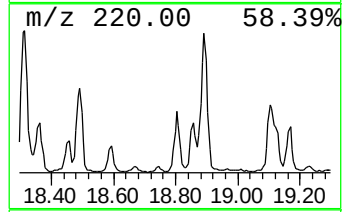
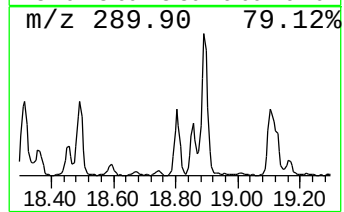
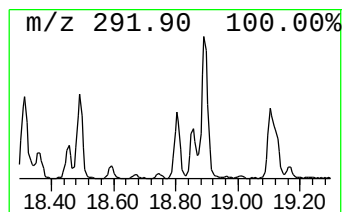
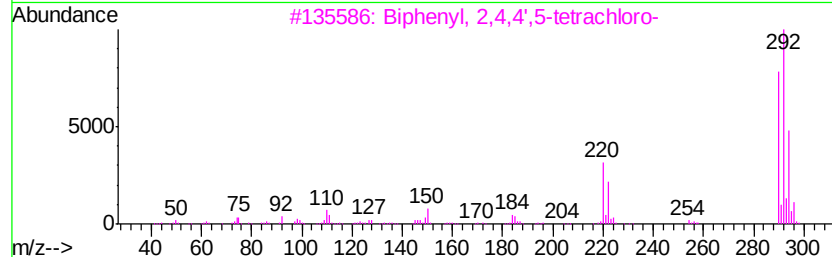
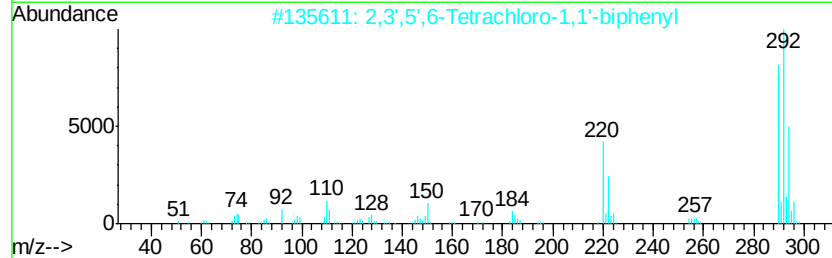
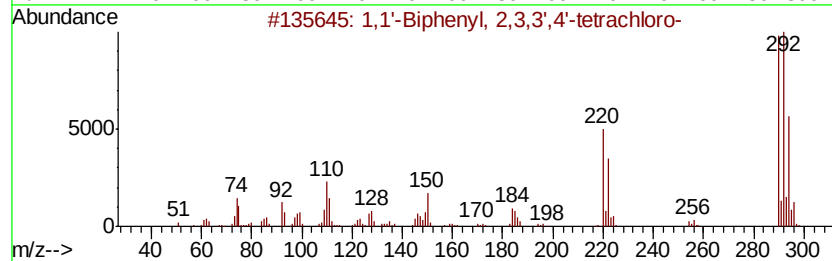
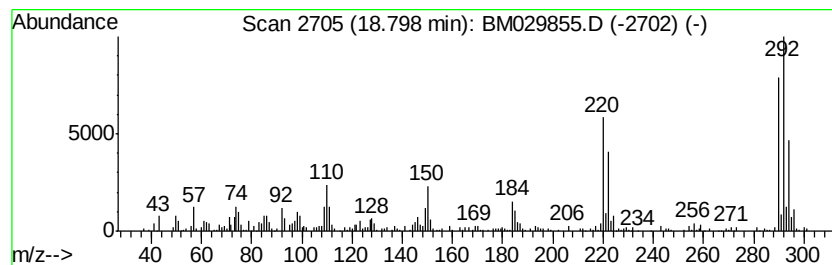
Instrument :
BNA_M
ClientSampleId :
C0B64

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 1,1'-Biphenyl, 2,3,3',4'-te... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.80	2.99 ng/ul	206320	Phenanthrene-d10	16.96		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,3',4'-tetrach...	290	C12H6Cl4	041464-43-1	99	
2	2,3',5',6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074338-23-1	99	
3	Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	99	
4	1,1'-Biphenyl, 3,3',5,5'-tetrach...	290	C12H6Cl4	033284-52-5	99	
5	1,1'-Biphenyl, 2,2',3,4-tetrachl...	290	C12H6Cl4	052663-59-9	99	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

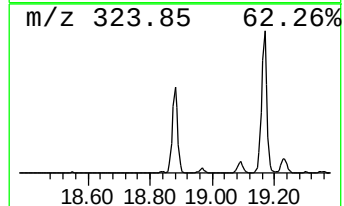
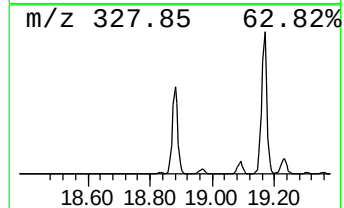
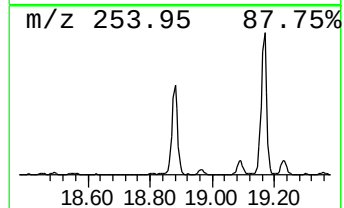
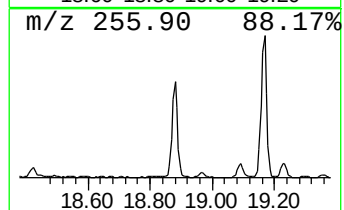
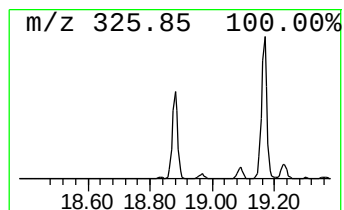
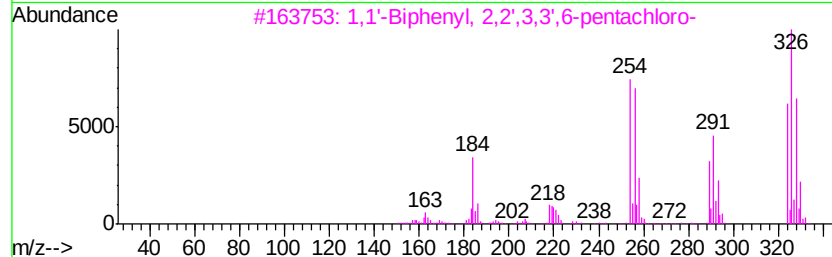
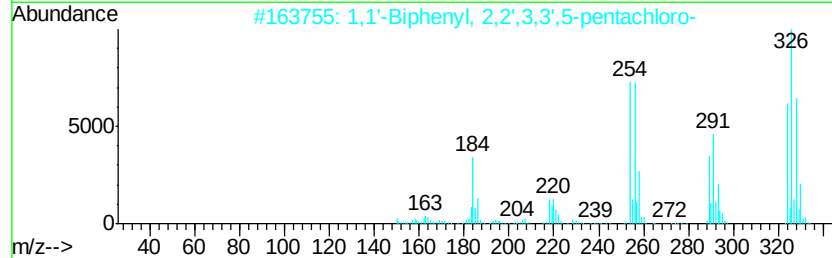
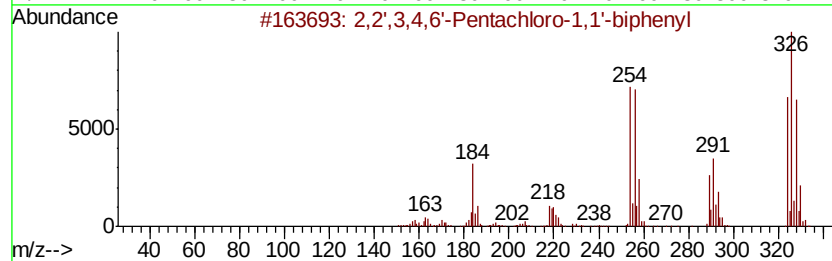
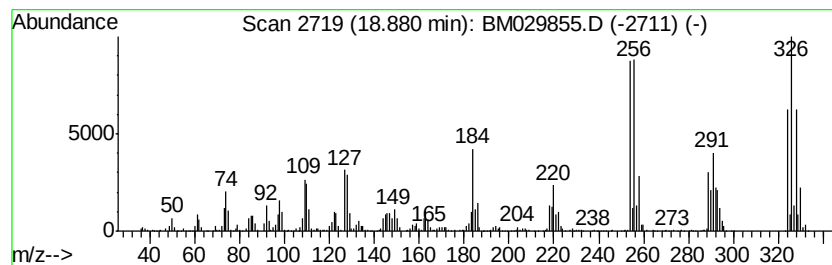
Instrument :
BNA_M
ClientSampleId :
C0B64

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 1,1'-Biphenyl, 2,2',3,3',5-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.88	20.99 ng/ul	1446710	Phenanthrene-d10	16.96		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,4,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-57-2	99	
2	1,1'-Biphenyl, 2,2',3,3',5-penta...	324	C12H5Cl5	060145-20-2	99	
3	1,1'-Biphenyl, 2,2',3,3',6-penta...	324	C12H5Cl5	052663-60-2	99	
4	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99	
5	1,1'-Biphenyl, 2,2',3,4,6-Pentac...	324	C12H5Cl5	055215-17-3	99	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

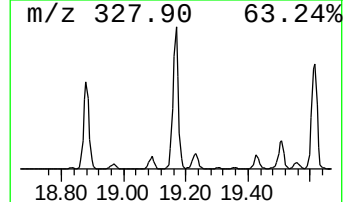
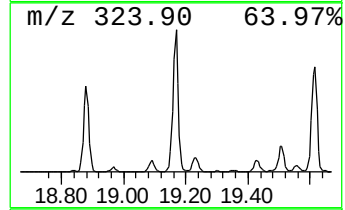
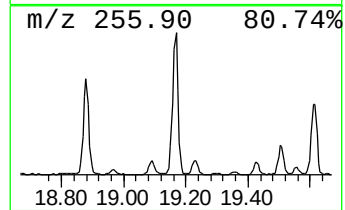
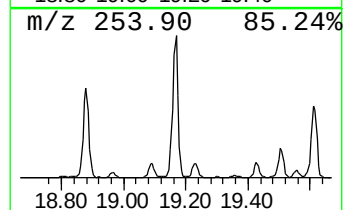
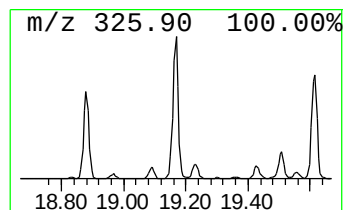
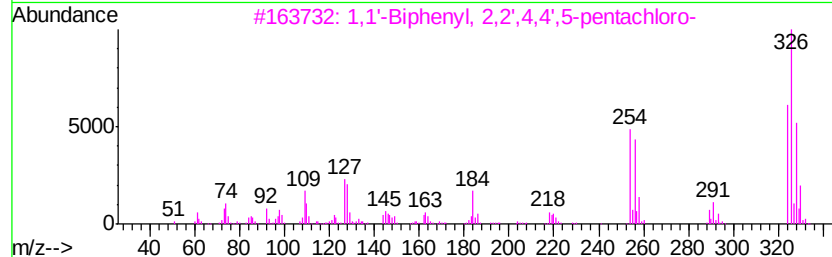
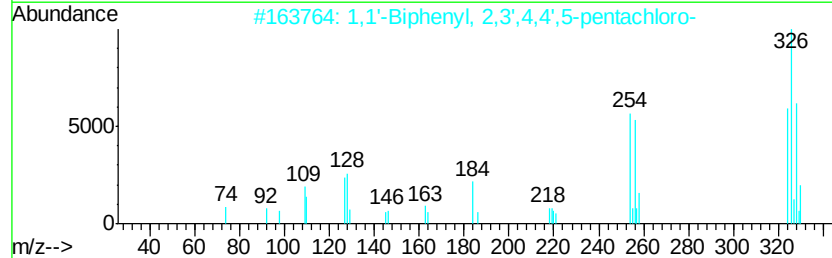
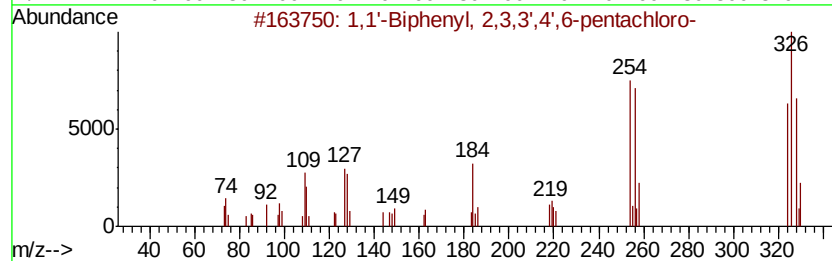
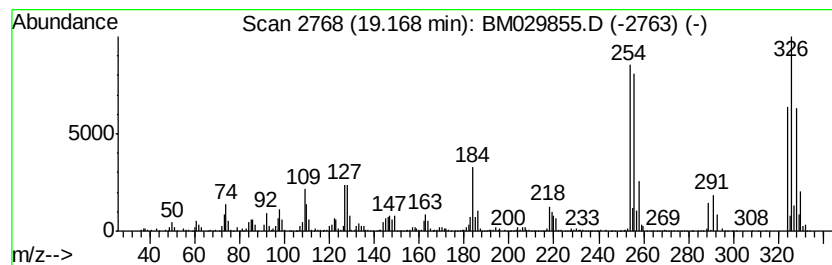
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 1,1'-Biphenyl, 2,3,3',4',6-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.17	23.58 ng/ul	1347750	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99
2		1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	99
3		1,1'-Biphenyl, 2,2',4,4',5-penta...	324	C12H5Cl5	038380-01-7	99
4		2,3,3',5',6-Pentachloro-1,1'-bi...	324	C12H5Cl5	068194-10-5	99
5		2,3',4',5',6-Pentachloro-1,1'-bi...	324	C12H5Cl5	074472-39-2	98



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

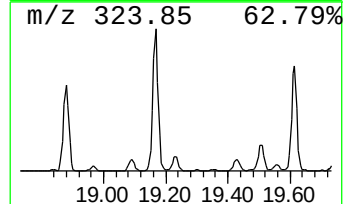
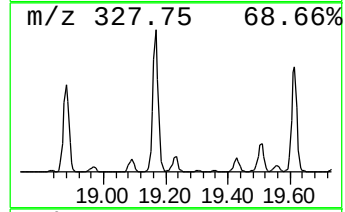
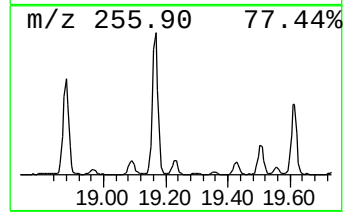
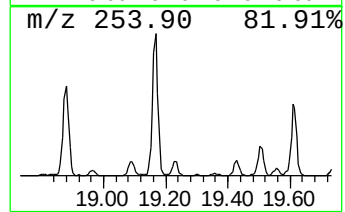
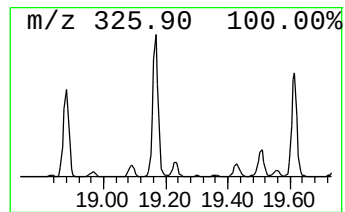
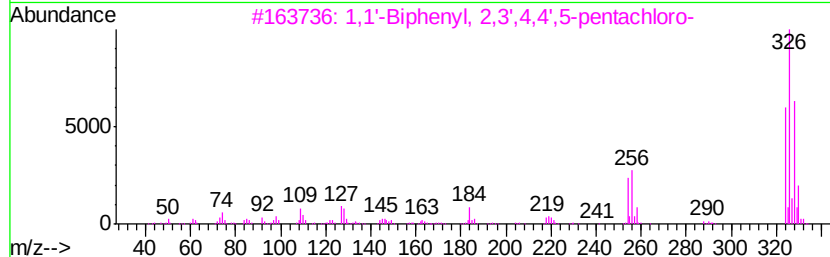
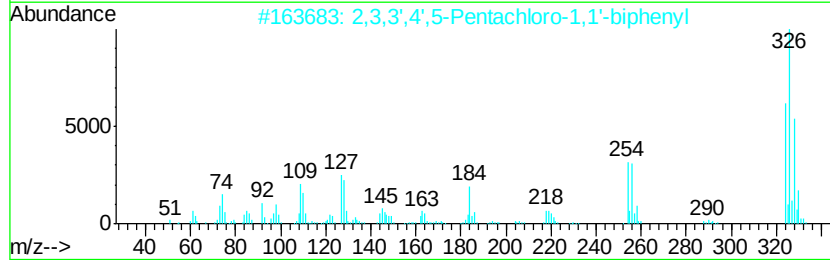
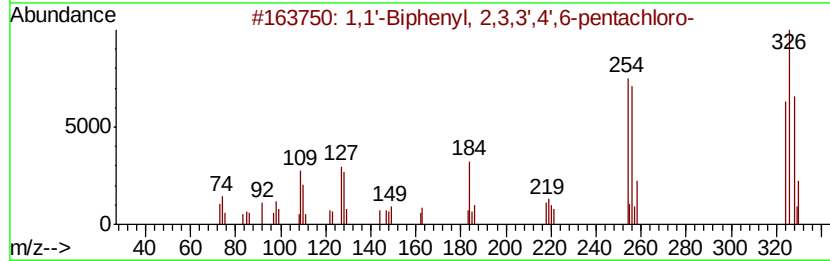
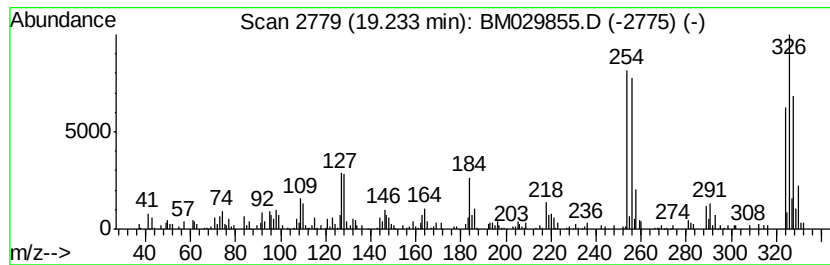
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 13 1,1'-Biphenyl, 2,3',4,4',5-... Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.23	2.54 ng/ul	145107	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99
2		2,3,3',4',5-Pentachloro-1,1'-bip...	324	C12H5Cl5	070424-68-9	99
3		1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	99
4		1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99
5		3,3',4,5,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	039635-33-1	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

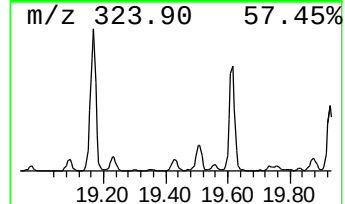
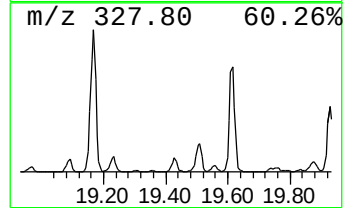
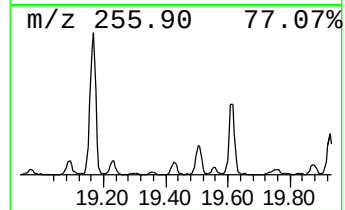
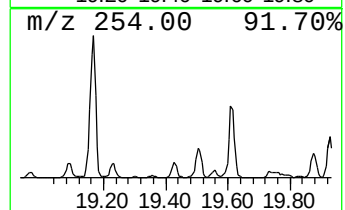
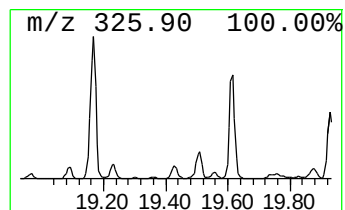
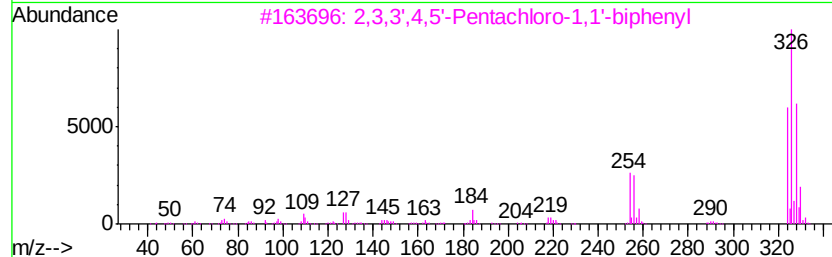
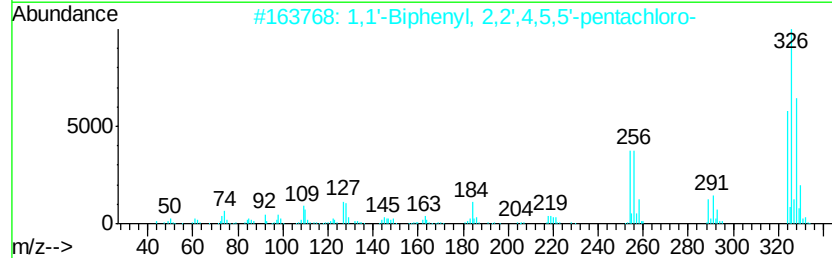
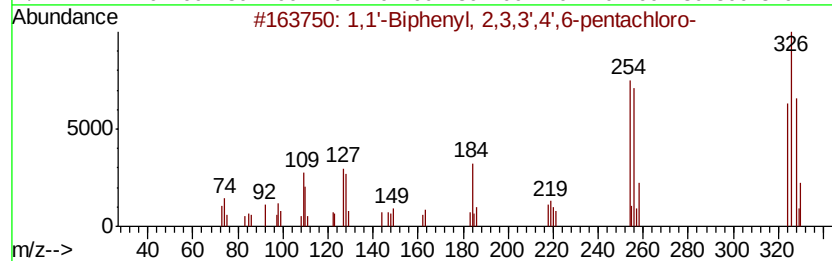
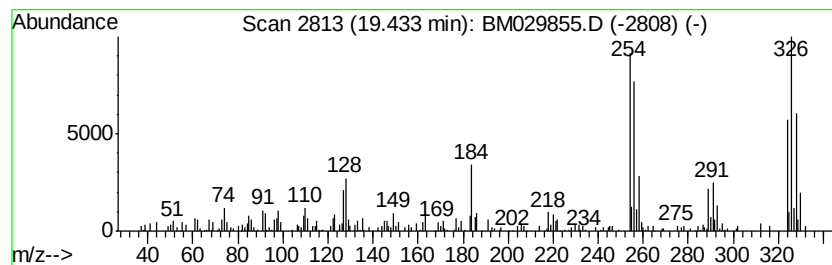
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 14 1,1'-Biphenyl, 2,2',4,5,5'-... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.43	2.69 ng/ul	153895	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99
2		1,1'-Biphenyl, 2,2',4,5,5'-penta...	324	C12H5Cl5	037680-73-2	99
3		2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	070362-41-3	99
4		2,3,3',5',6-Pentachloro-1,1'-bi...	324	C12H5Cl5	068194-10-5	99
5		2,2',3,4',5-Pentachloro-1,1'-bip...	324	C12H5Cl5	068194-07-0	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

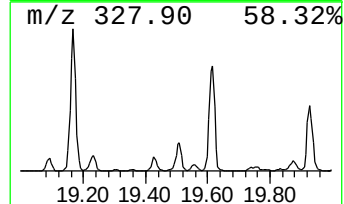
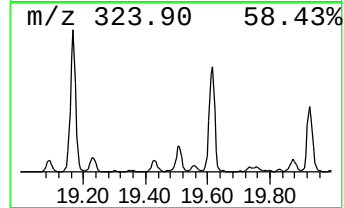
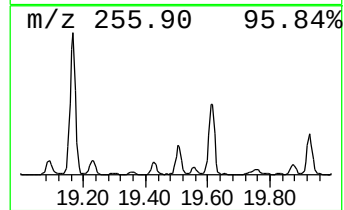
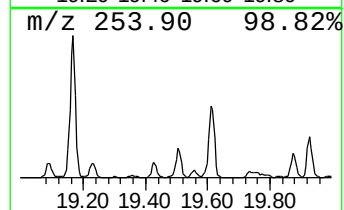
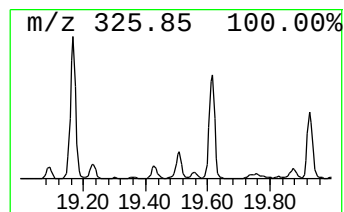
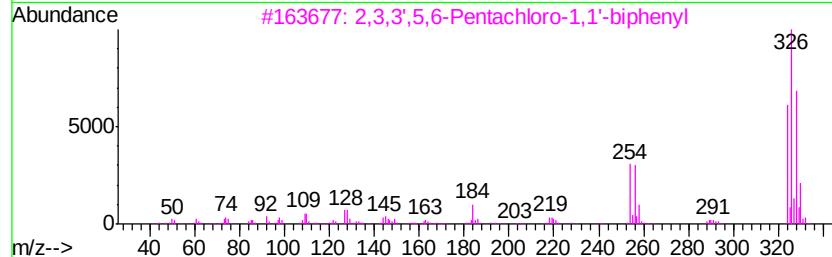
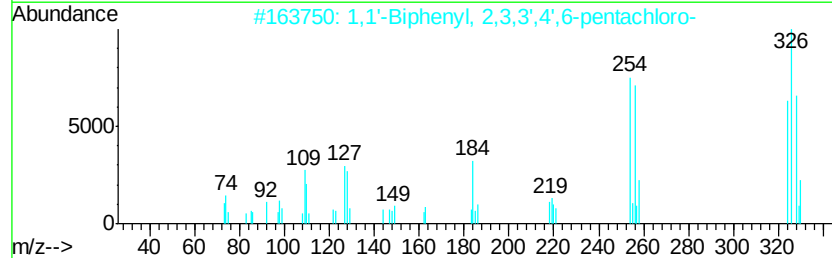
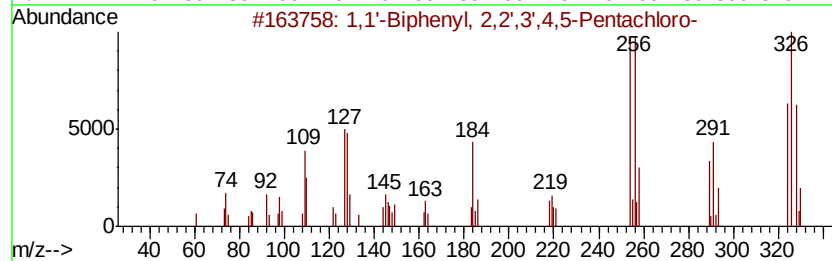
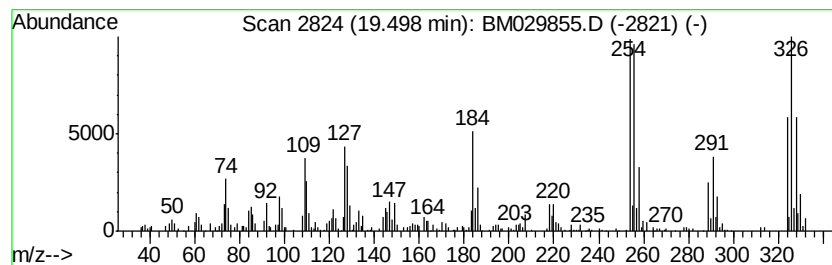
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 15 1,1'-Biphenyl, 2,2',3',4,5-... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.50	4.80 ng/ul	274508	Chrysene-d12	21.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'	-Biphenyl, 2,2',3',4,5-Penta...	324	C12H5Cl5	041464-51-1	99	
2	1,1'	-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99	
3	2,3,3',5,6-Pentachloro-1,1'	-biph...	324	C12H5Cl5	074472-36-9	99	
4	1,1'	-Biphenyl, 2,2',3,3',4-penta...	324	C12H5Cl5	052663-62-4	99	
5	1,1'	-Biphenyl, 2,2',4,5,6'-penta...	324	C12H5Cl5	068194-06-9	99	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
C0B64

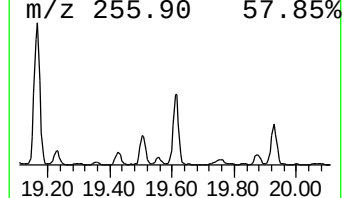
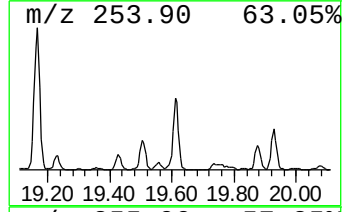
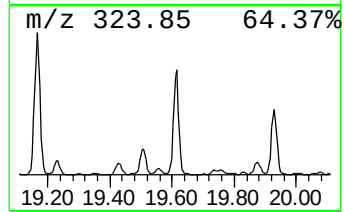
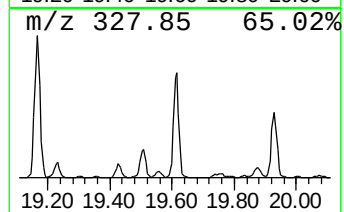
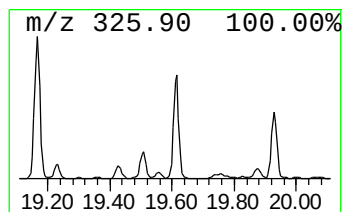
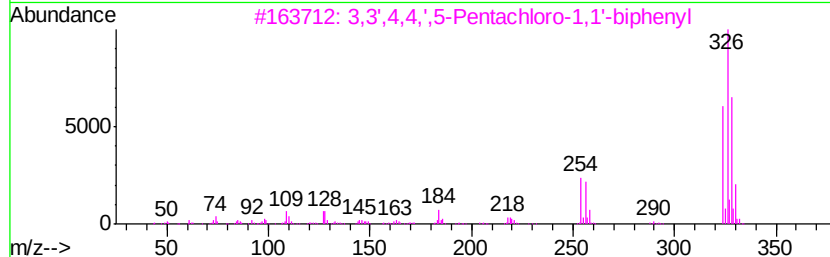
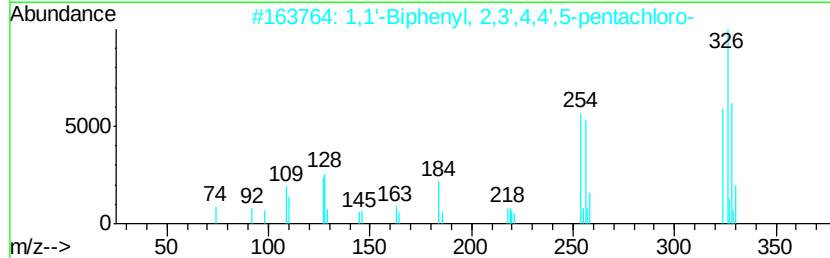
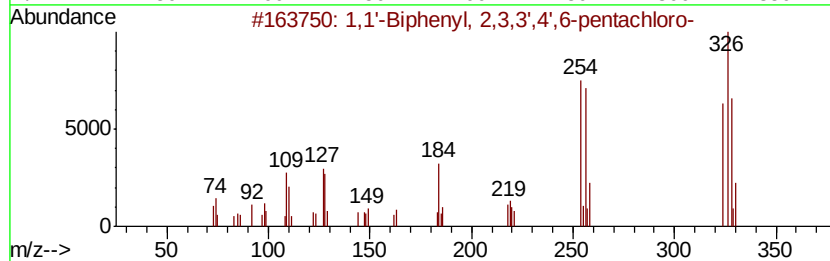
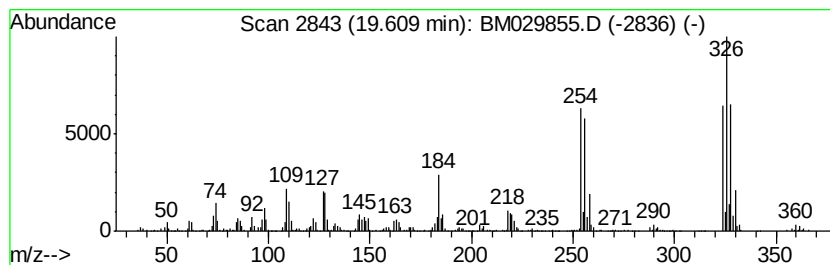
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 16 3,3',4,4',5-Pentachloro-1,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.61	19.26 ng/ul	1101180	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99
2		1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	99
3		3,3',4,4',5-Pentachloro-1,1'-bi...	324	C12H5Cl5	057465-28-8	99
4		1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	99
5		3,3',4,5,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	039635-33-1	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

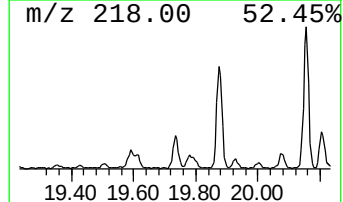
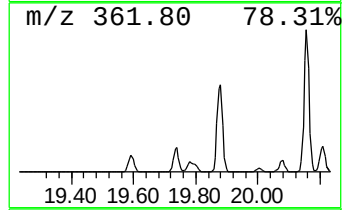
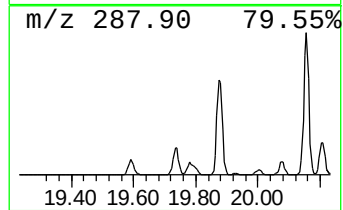
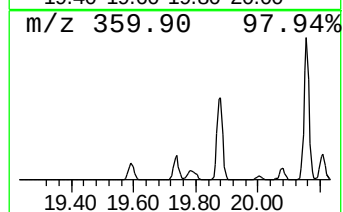
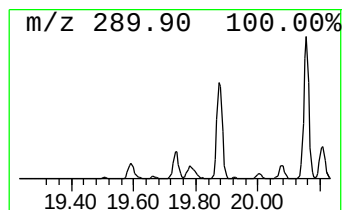
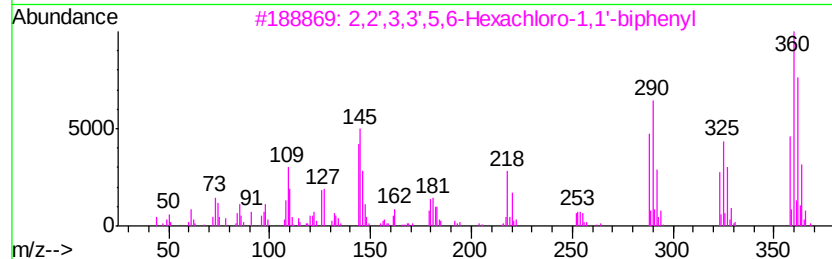
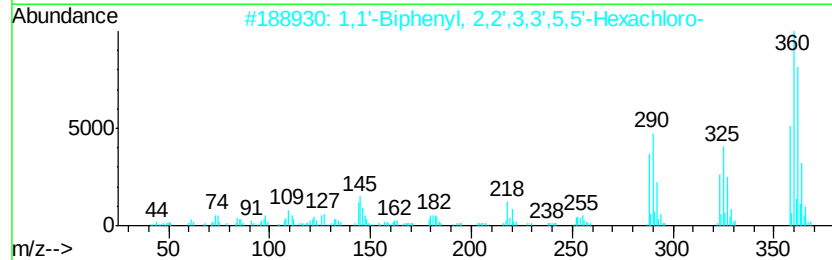
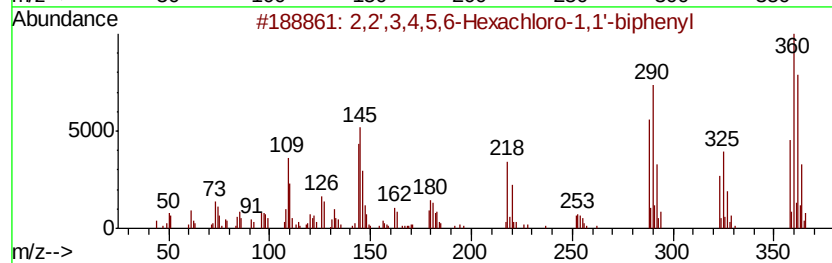
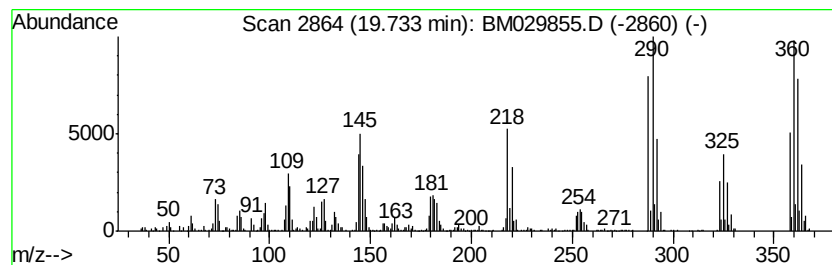
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 17 2,2',3,4,5,6-Hexachloro-1,1... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.73	11.99 ng/ul	685521	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2',3,4,5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-61-4	99
2		1,1'-Biphenyl, 2,2',3,3',5,5'-He...	358	C12H4Cl6	035694-04-3	99
3		2,2',3,3',5,6-Hexachloro-1,1'-bi...	358	C12H4Cl6	052704-70-8	99
4		1,1'-Biphenyl, 2,2',3,3',4,5-hex...	358	C12H4Cl6	055215-18-4	99
5		2,2',3,4',5,6-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-13-8	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

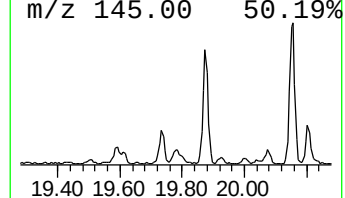
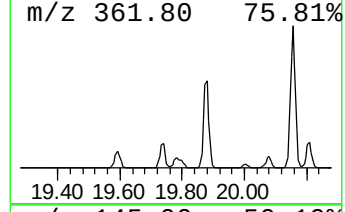
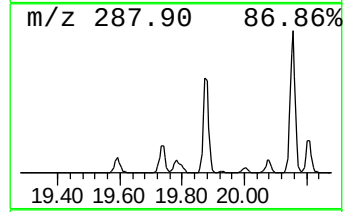
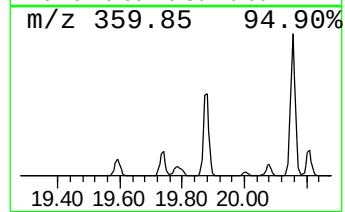
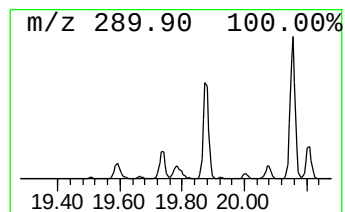
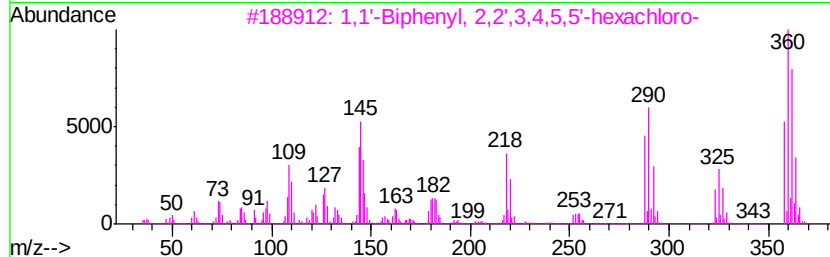
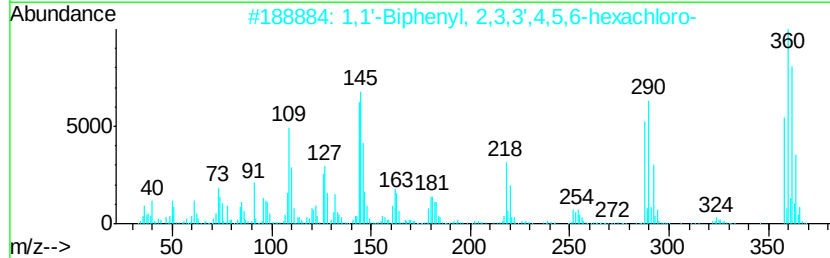
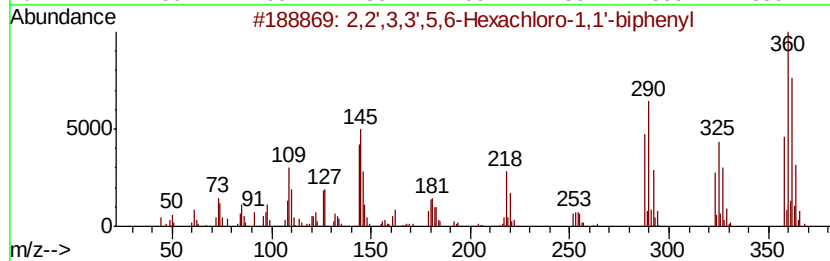
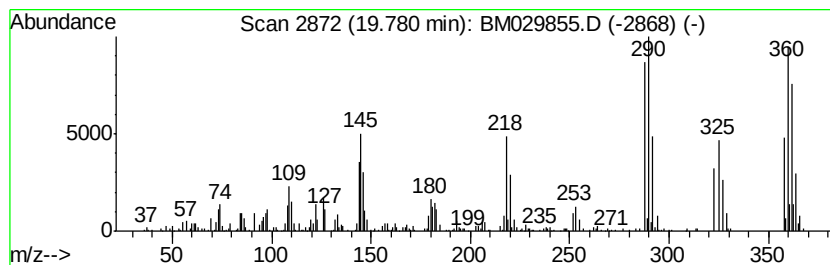
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 18 2,2',3,3',5,6-Hexachloro-1,... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.78	7.30 ng/ul	417313	Chrysene-d12	21.14

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		1,1'-Biphenyl, 2,3,3',4,5,6-hexa...	358	C12H4Cl6	041411-62-5	99
3		1,1'-Biphenyl, 2,2',3,4,5,5'-hex...	358	C12H4Cl6	052712-04-6	99
4		2,2',3,4,5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-61-4	96
5		2,2',3,4',5,6'-Hexachloro-1,1'-b...	358	C12H4Cl6	074472-41-6	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

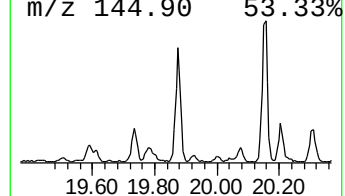
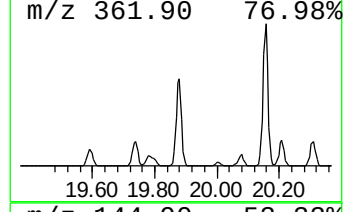
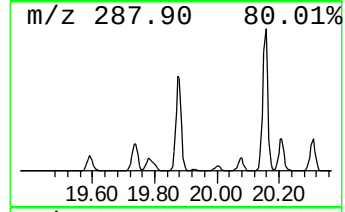
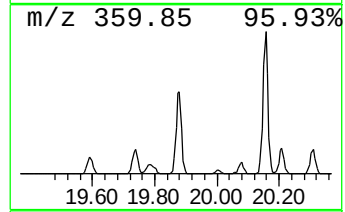
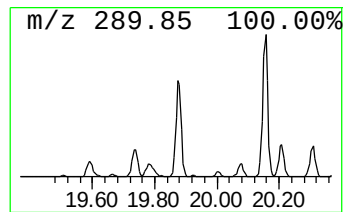
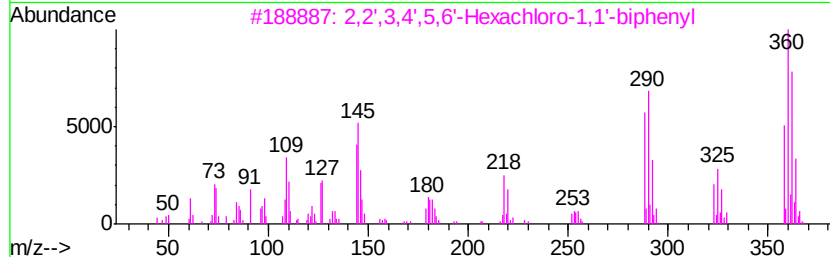
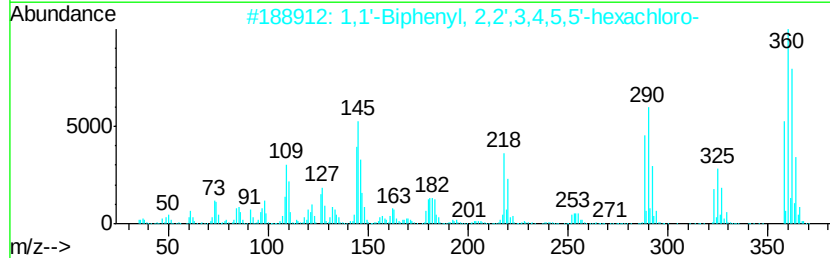
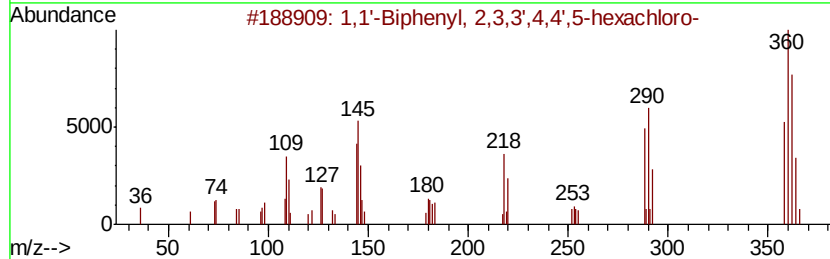
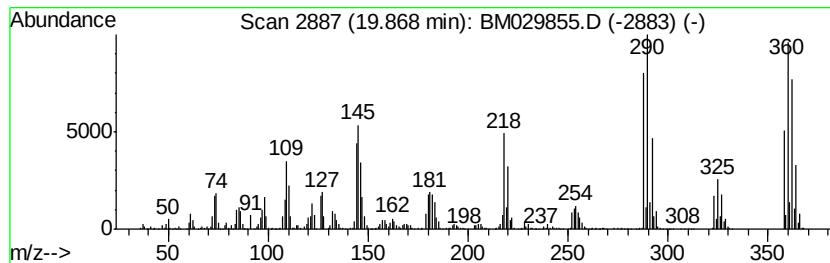
Instrument :
BNA_M
ClientSampleId :
C0B64

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
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 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
19.87	38.82 ng/ul	2219260	Chrysene-d12	21.14		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'	-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	99
2	1,1'	-Biphenyl, 2,2',3,4,5,5'-hex...	358	C12H4Cl6	052712-04-6	99
3	2,2',3,4',5,6'	-Hexachloro-1,1'-b...	358	C12H4Cl6	074472-41-6	99
4	1,1'	-Biphenyl, 2,2',3,4',5',6-he...	358	C12H4Cl6	038380-04-0	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\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

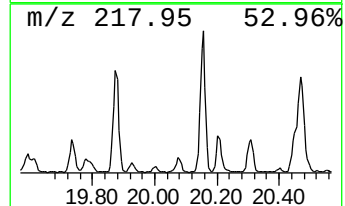
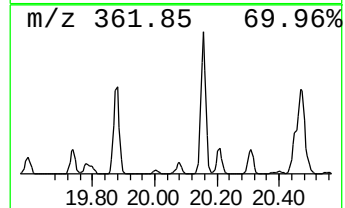
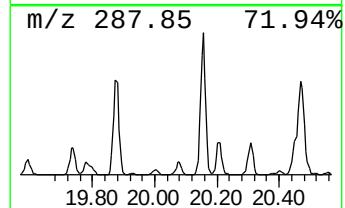
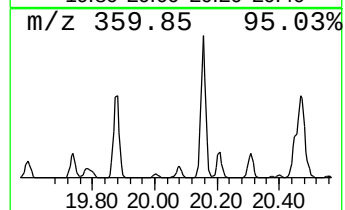
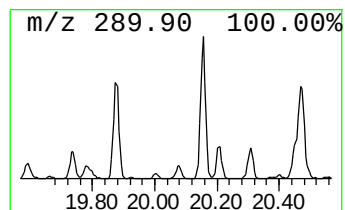
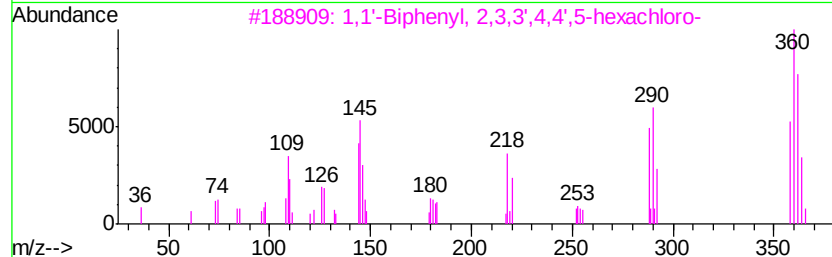
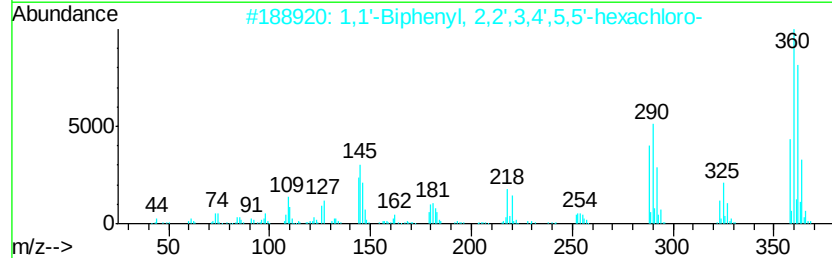
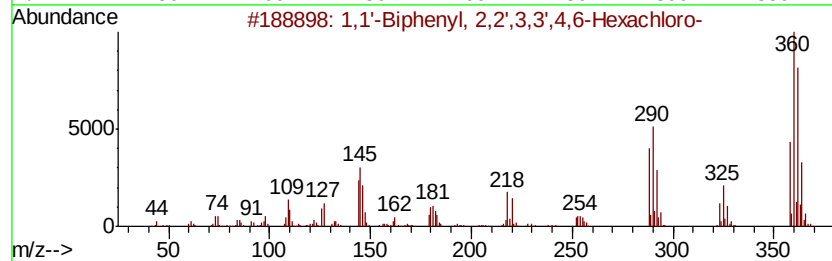
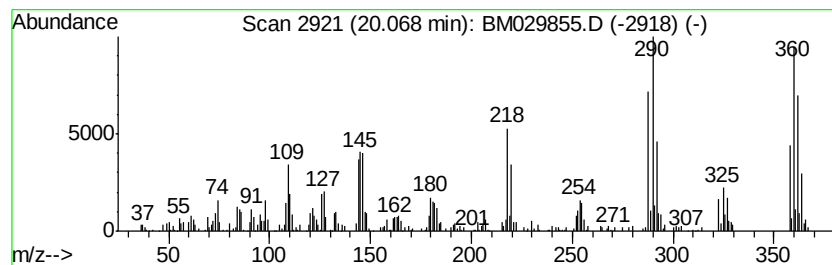
Instrument :
BNA_M
ClientSampleId :
C0B64

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 21 1,1'-Biphenyl, 2,2',3,3',4,... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.		
20.07	5.46 ng/ul	312312	Chrysene-d12	21.14		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,3',4,6-Hex...	358	C12H4Cl6	061798-70-7	99	
2	1,1'-Biphenyl, 2,2',3,4',5,5'-he...	358	C12H4Cl6	051908-16-8	99	
3	1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	99	
4	1,1'-Biphenyl, 2,3',4,4',5,5'-he...	358	C12H4Cl6	052663-72-6	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\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

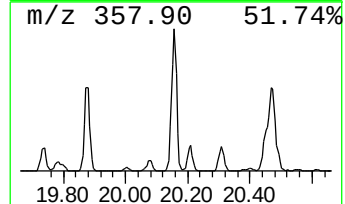
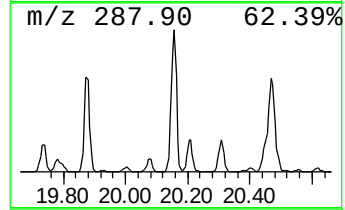
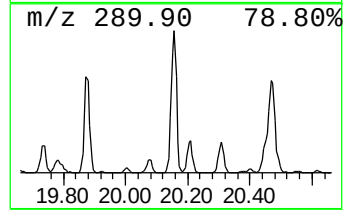
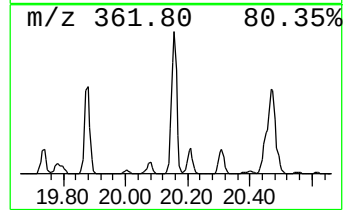
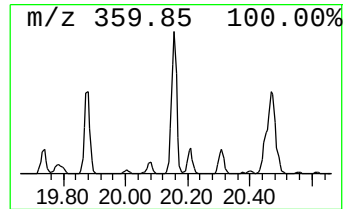
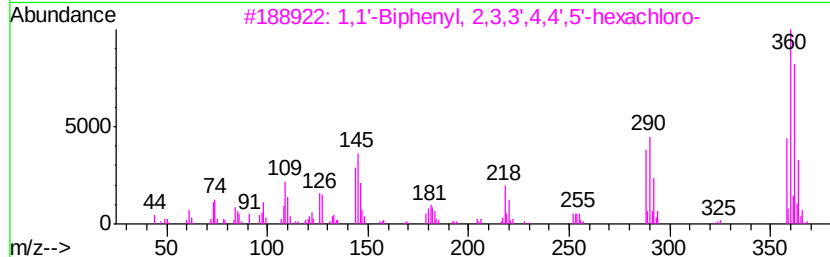
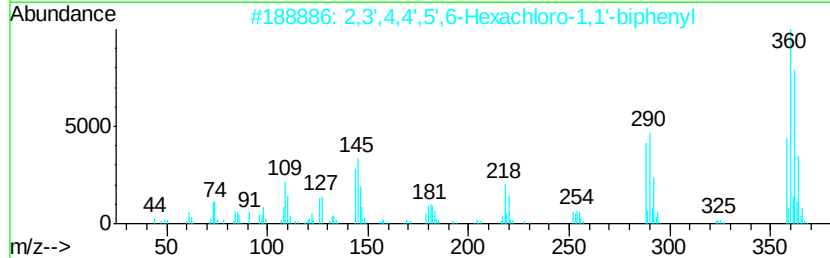
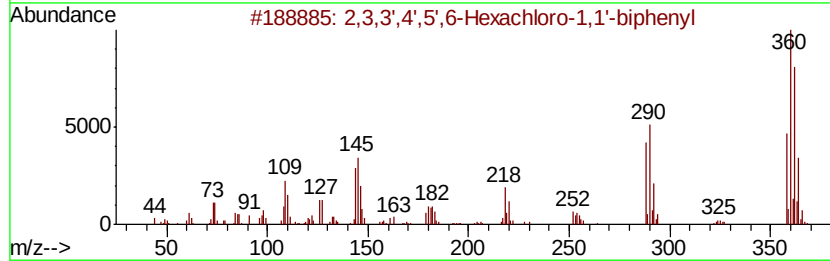
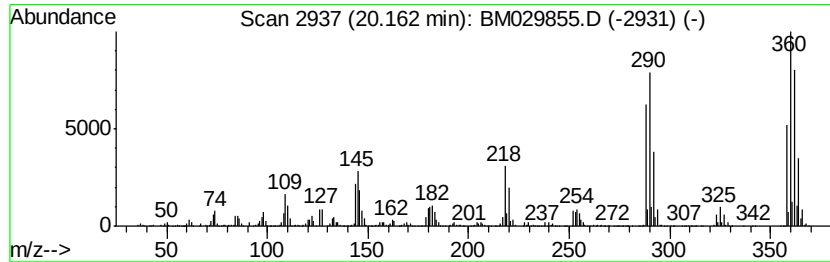
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 22 2,3,3',4',5',6-Hexachloro-1... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.16	52.17 ng/ul	2982270	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,3,3',4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	074472-45-0	99
2		2,3',4,4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-65-5	99
3		1,1'-Biphenyl, 2,3,3',4,4',5'-he...	358	C12H4Cl6	069782-90-7	99
4		1,1'-Biphenyl, 2,3',4,4',5,5'-he...	358	C12H4Cl6	052663-72-6	99
5		2,2',3,4',6,6'-Hexachloro-1,1'-b...	358	C12H4Cl6	068194-08-1	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

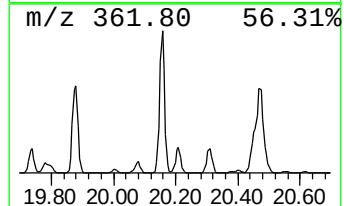
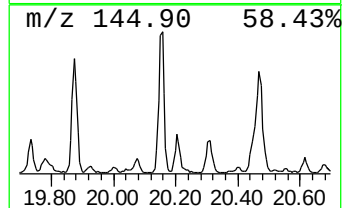
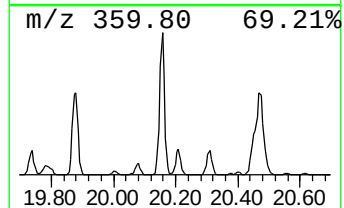
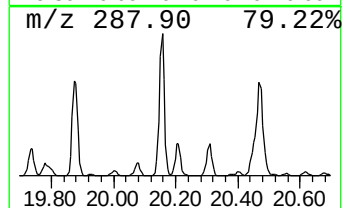
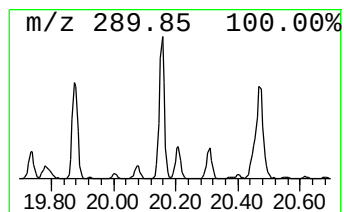
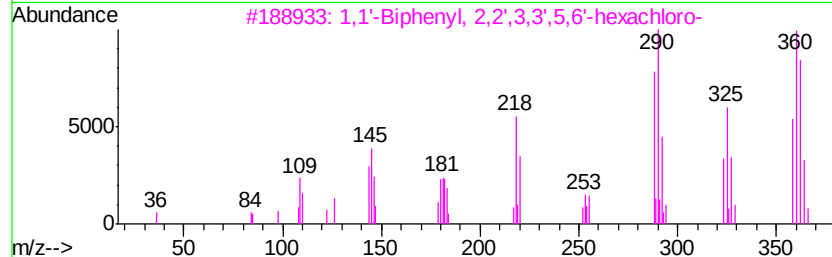
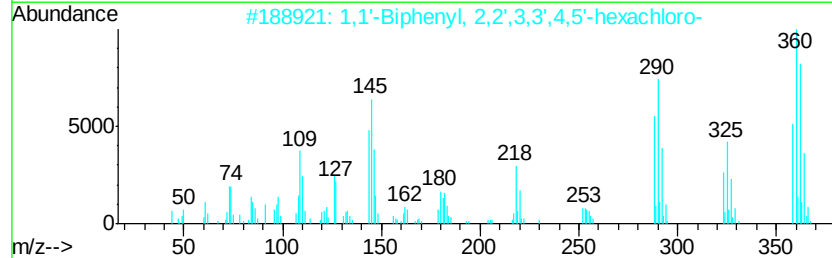
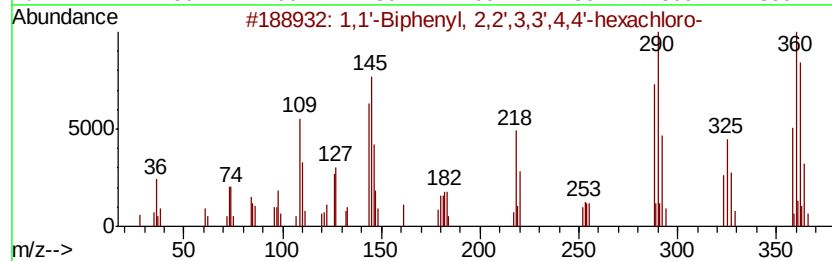
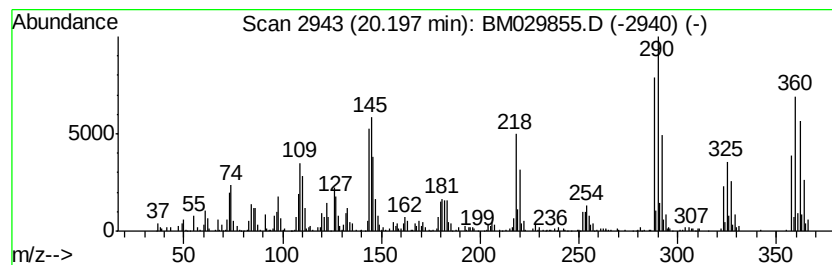
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 23 1,1'-Biphenyl, 2,2',4,4',6,... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.20	16.26 ng/ul	929732	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',3,3',4,4'-he...	358	C12H4Cl6	038380-07-3	99
2		1,1'-Biphenyl, 2,2',3,3',4,5'-he...	358	C12H4Cl6	052663-66-8	99
3		1,1'-Biphenyl, 2,2',3,3',5,6'-he...	358	C12H4Cl6	052744-13-5	99
4		1,1'-Biphenyl, 2,2',3,3',4,6'-he...	358	C12H4Cl6	038380-05-1	99
5		1,1'-Biphenyl, 2,2',4,4',6,6'-he...	358	C12H4Cl6	033979-03-2	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

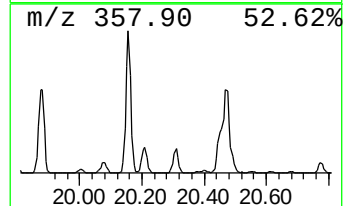
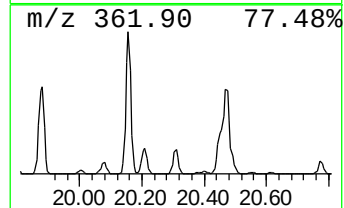
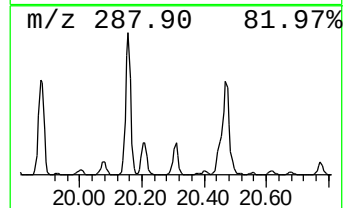
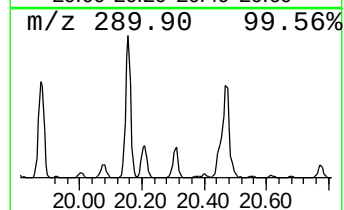
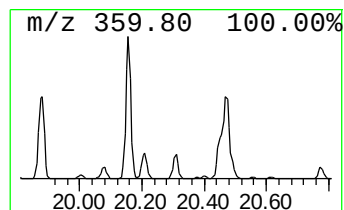
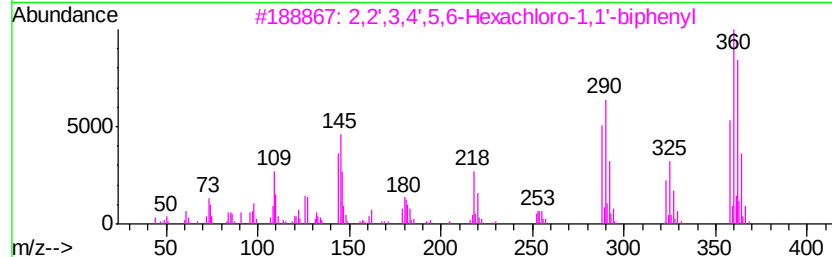
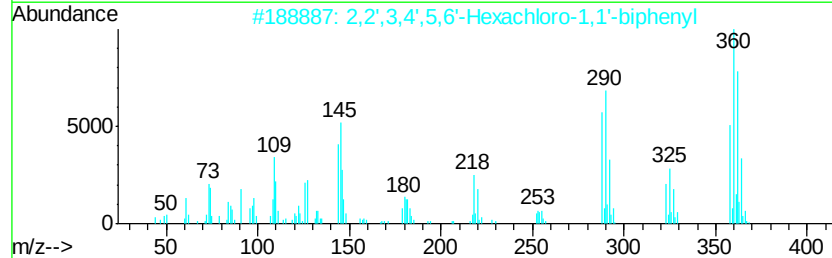
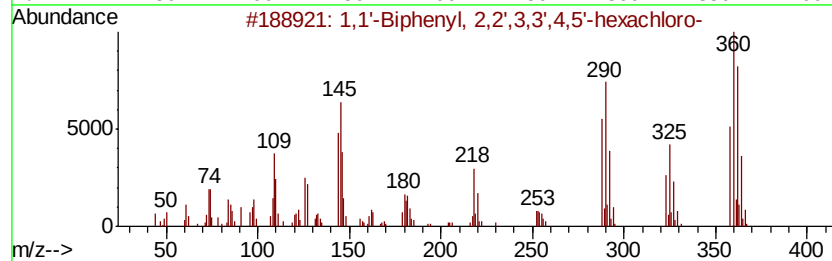
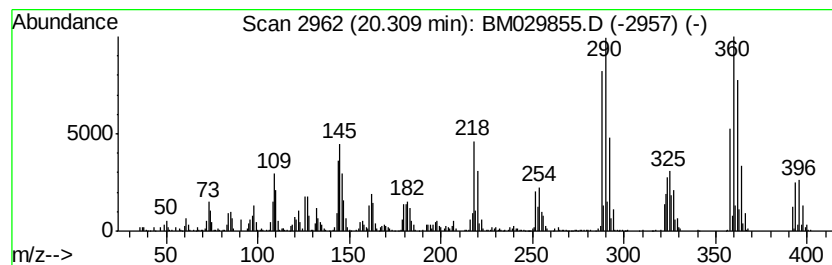
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 24 1,1'-Biphenyl, 2,2',3,3',4,... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.31	17.27 ng/ul	987148	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',3,3',4,5'-he...	358	C12H4Cl6	052663-66-8	99
2		2,2',3,4',5,6'-Hexachloro-1,1'-b...	358	C12H4Cl6	074472-41-6	99
3		2,2',3,4',5,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-13-8	99
4		1,1'-Biphenyl, 2,2',3,4,4',5'-he...	358	C12H4Cl6	035065-28-2	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\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

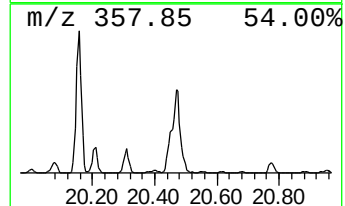
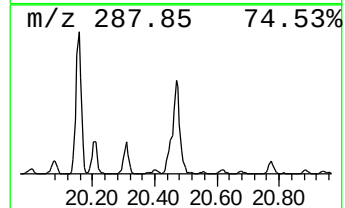
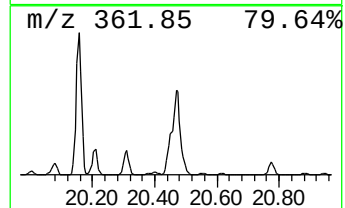
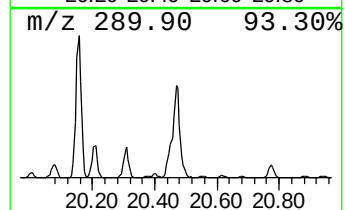
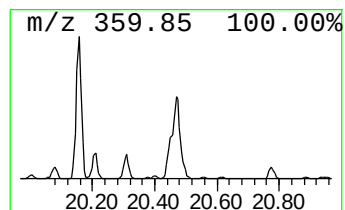
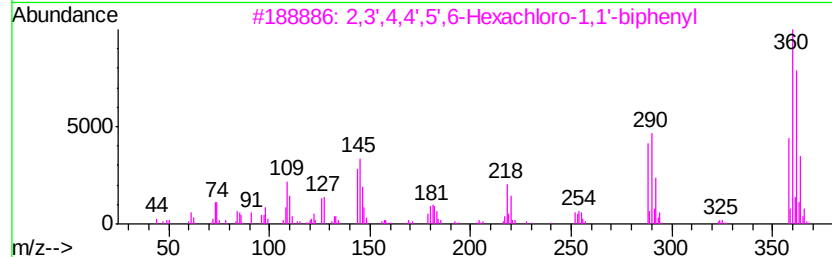
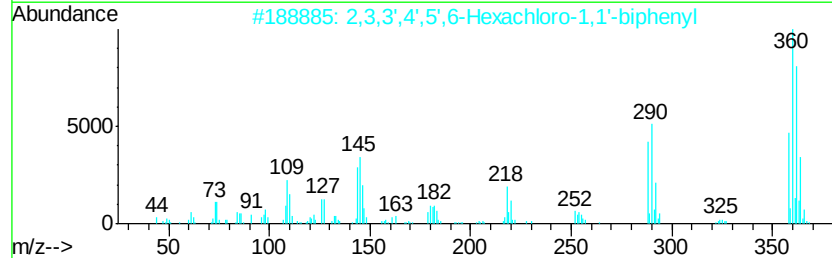
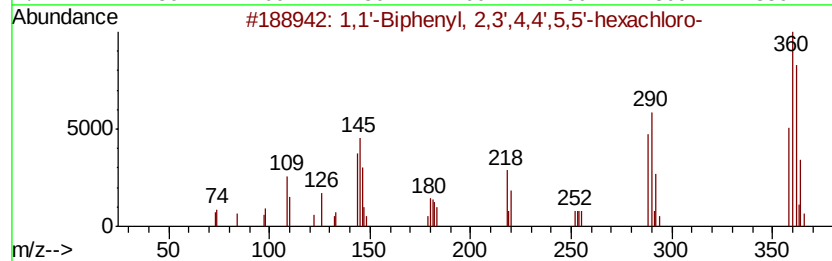
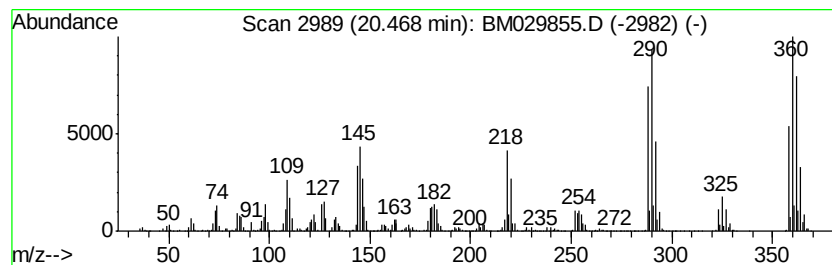
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.47	53.11 ng/ul	3036080	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3',4,4',5,5'-he...	358	C12H4Cl6	052663-72-6	99
2		2,3,3',4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	074472-45-0	99
3		2,3',4,4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-65-5	99
4		1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	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\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

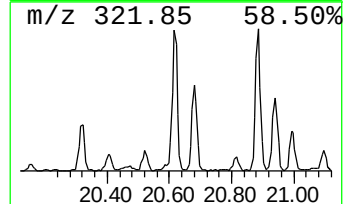
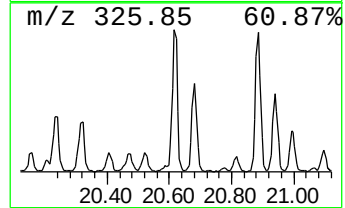
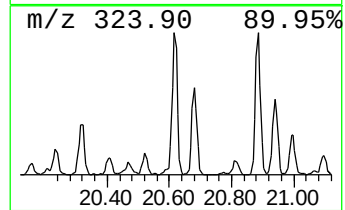
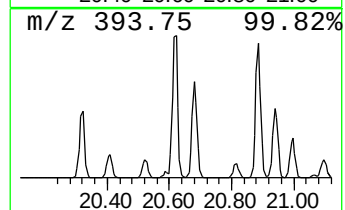
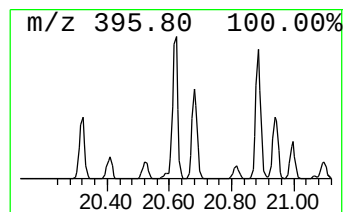
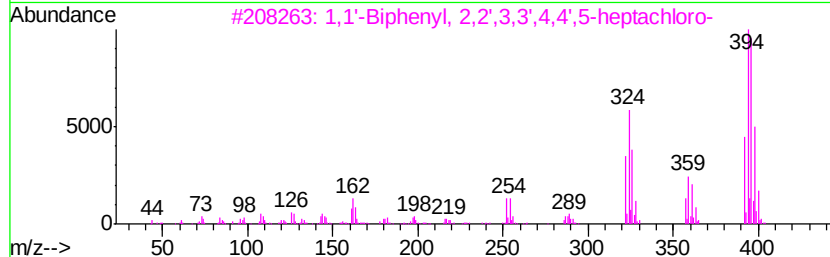
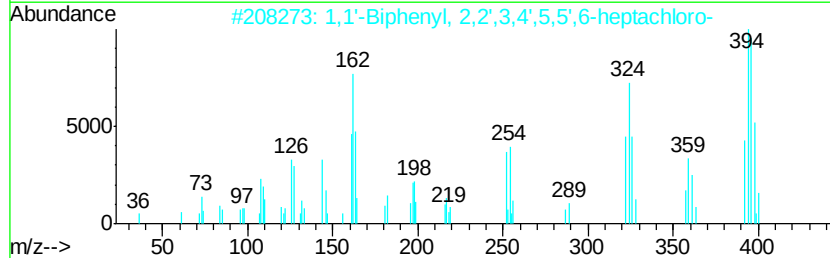
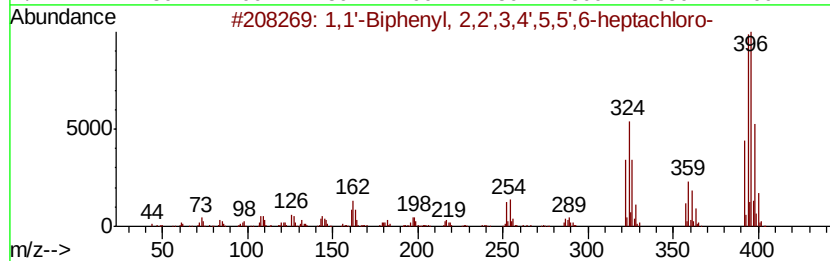
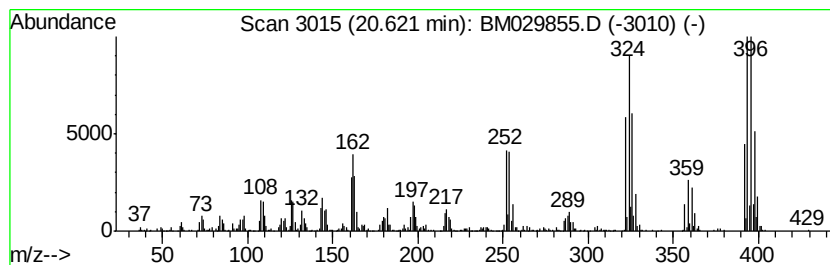
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 27 1,1'-Biphenyl, 2,2',3,4',5,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.62	17.64 ng/ul	1008400	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',3,4',5,5',6-...	392	C12H3Cl7	052663-68-0	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		1,1'-Biphenyl, 2,2',3,3',4,4',6-...	392	C12H3Cl7	052663-71-5	99
5		1,1'-Biphenyl, 2,2',3,4,4',5,6'-...	392	C12H3Cl7	060145-23-5	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

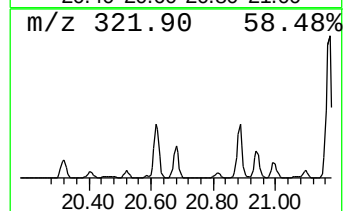
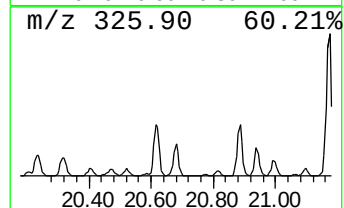
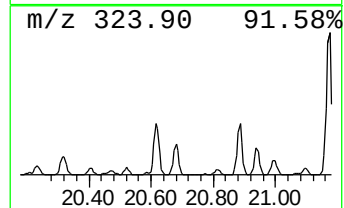
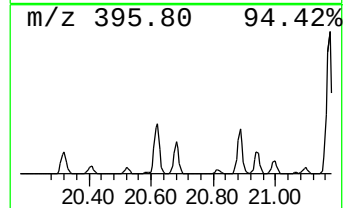
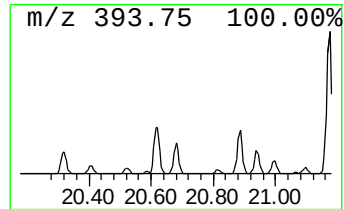
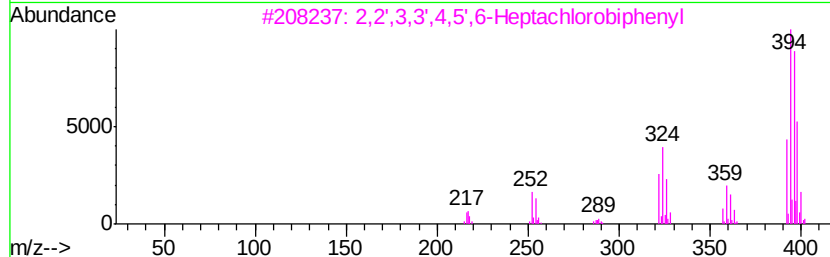
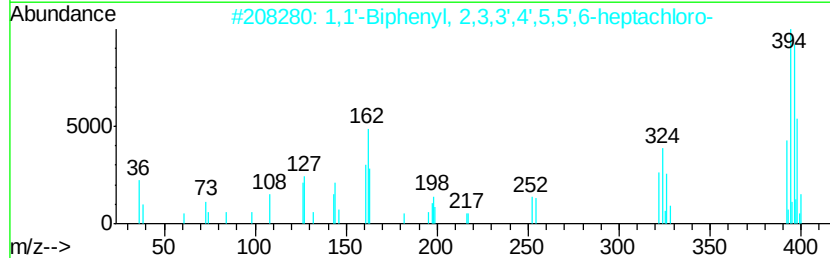
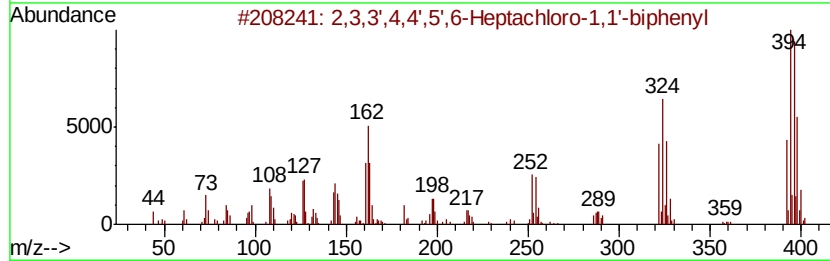
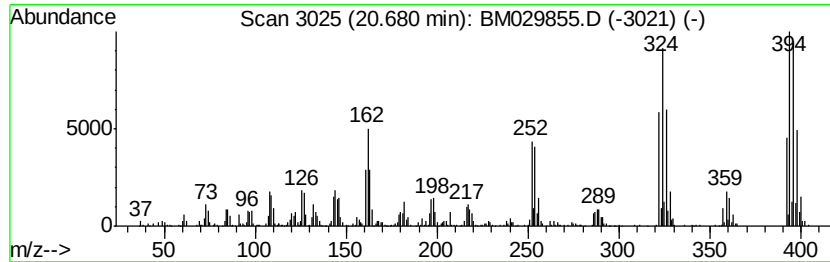
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 28 2,3,3',4,4',5',6-Heptachlor... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.68	10.46 ng/ul	597867	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,3,3',4,4',5',6-Heptachloro-1,1...	392	C12H3Cl7	074472-50-7	99
2		1,1'-Biphenyl, 2,3,3',4,4',5,5',6-...	392	C12H3Cl7	069782-91-8	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\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

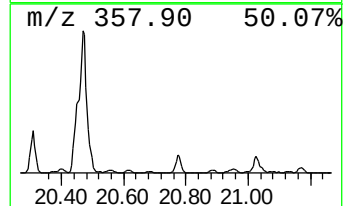
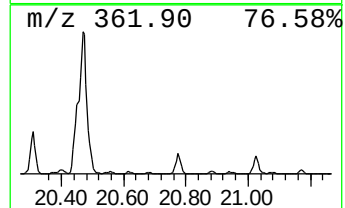
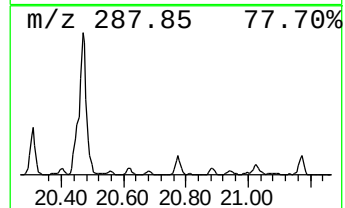
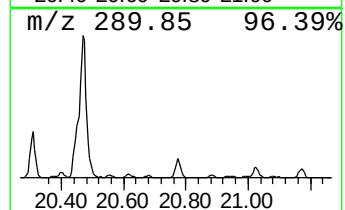
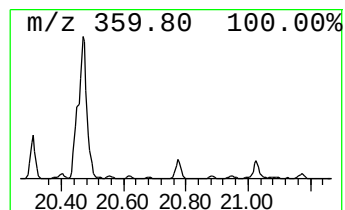
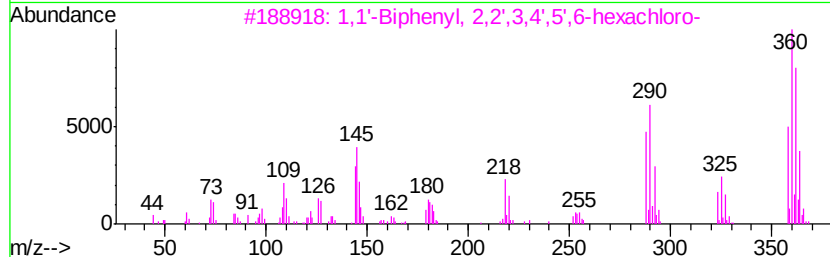
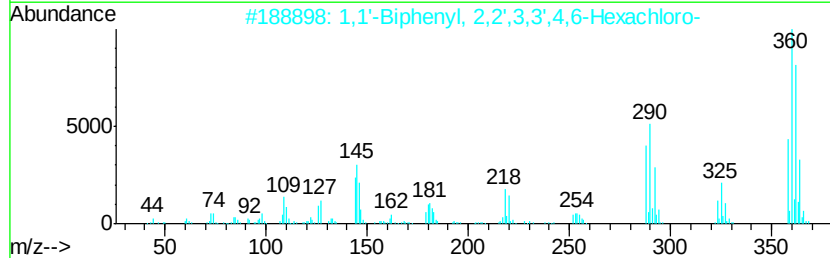
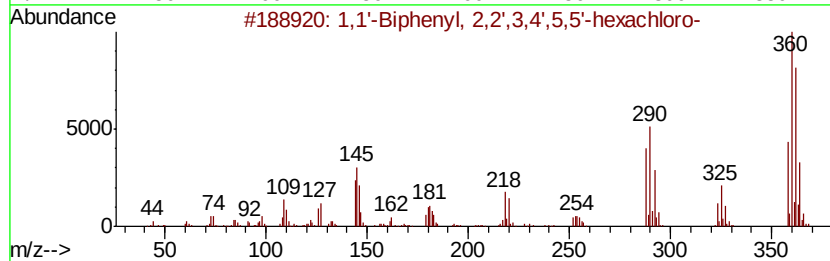
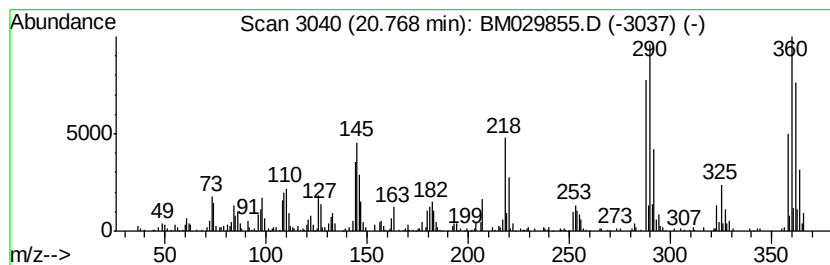
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 29 1,1'-Biphenyl, 2,2',3,4',5,... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.77	4.64 ng/ul	265367	Chrysene-d12	21.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'	-Biphenyl, 2,2',3,4',5,5'-he...	358	C12H4Cl6	051908-16-8	99	
2	1,1'	-Biphenyl, 2,2',3,3',4,6-Hex...	358	C12H4Cl6	061798-70-7	99	
3	1,1'	-Biphenyl, 2,2',3,4',5',6-he...	358	C12H4Cl6	038380-04-0	99	
4	2,2',3,4,4',6'	-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-64-4	99	
5	1,1'	-Biphenyl, 2,2',3,4,4',6-Hex...	358	C12H4Cl6	056030-56-9	99	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

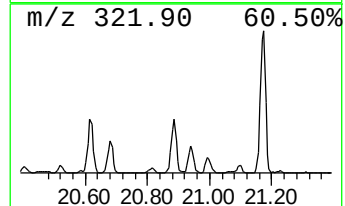
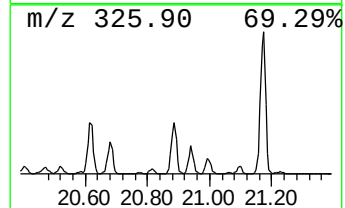
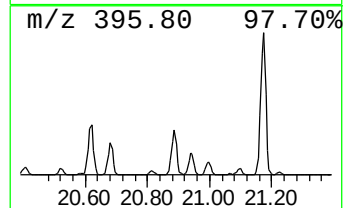
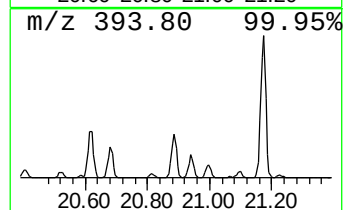
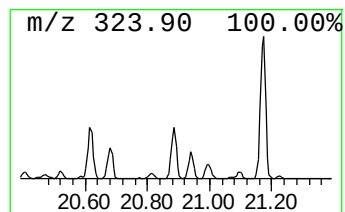
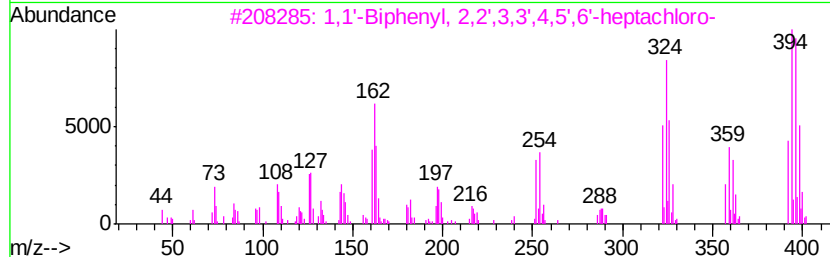
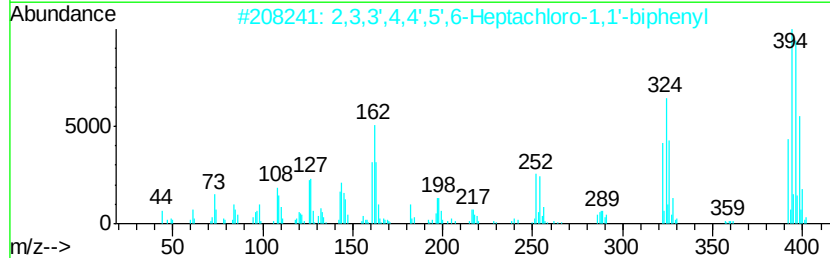
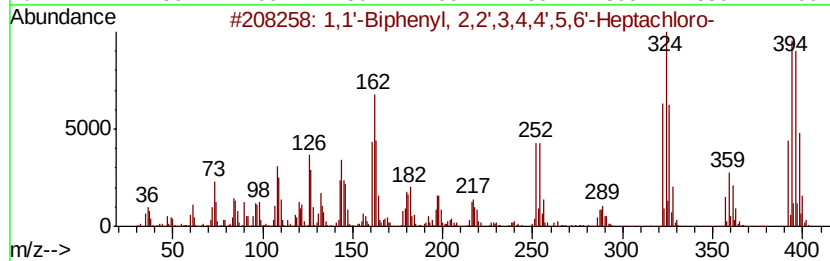
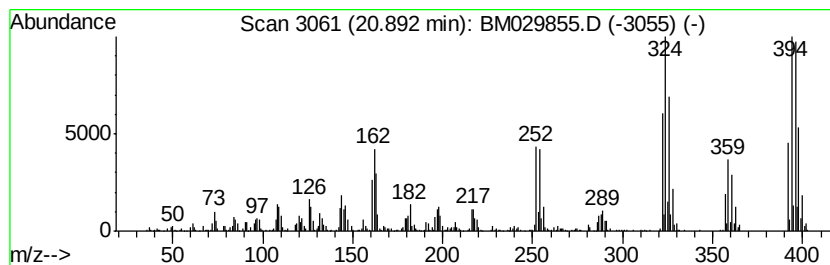
Instrument :
BNA_M
ClientSampleId :
C0B64

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 30 1,1'-Biphenyl, 2,2',3,4,4',... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
20.89	17.42 ng/ul	995552	Chrysene-d12	21.14		
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	2,3,3',4,4',5',6-Heptachloro-1,1...	392	C12H3Cl7	074472-50-7	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',5,5'-...	392	C12H3Cl7	039635-31-9	99	
5	2,2',3,3',4,5,6-Heptachloro-1,1'...	392	C12H3Cl7	068194-16-1	99	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

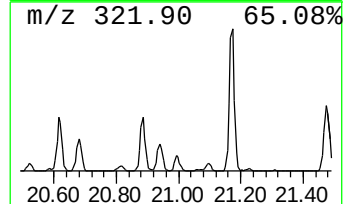
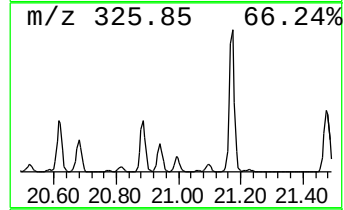
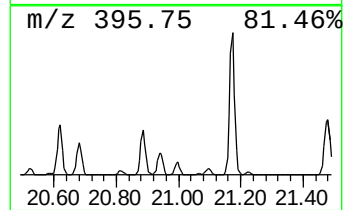
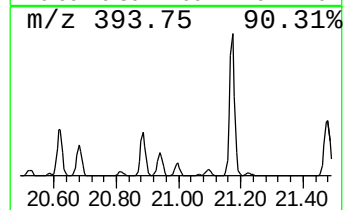
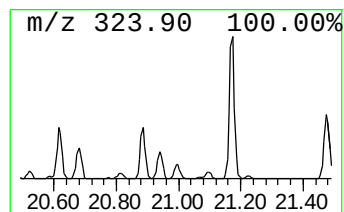
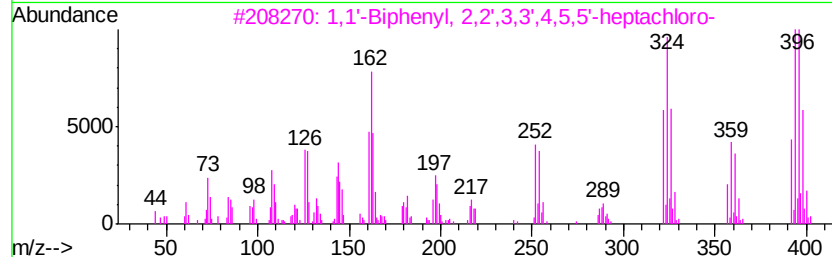
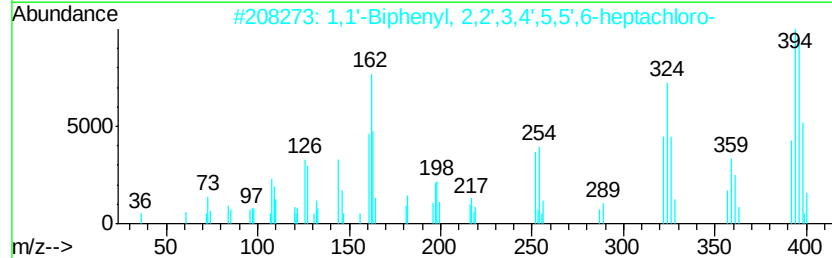
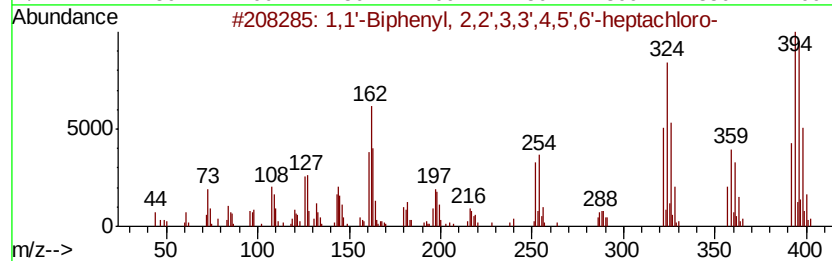
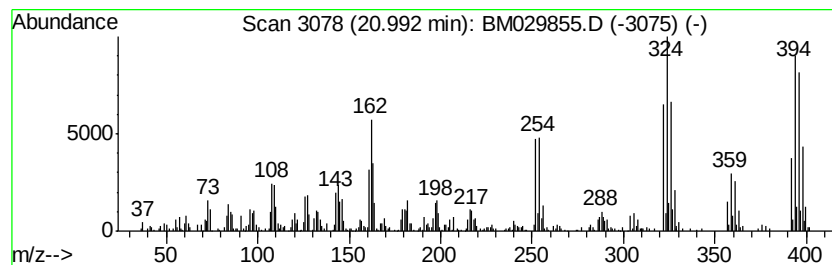
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 32 1,1'-Biphenyl, 2,2',3,3',4,... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.99	5.36 ng/ul	306517	Chrysene-d12	21.14

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',4,5,5'-...	392	C12H3Cl7	052663-74-8	99	
4	1,1'	-Biphenyl, 2,2',3,4,4',5,6'-...	392	C12H3Cl7	060145-23-5	99	
5	2,2',3,4',5,6,6'	-Heptachloro-1,1...	392	C12H3Cl7	074487-85-7	99	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029855.D
Acq On : 06 May 2021 08:55
Operator : CG/JU
Sample : M2240-01
Misc :
ALS Vial : 74 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B64

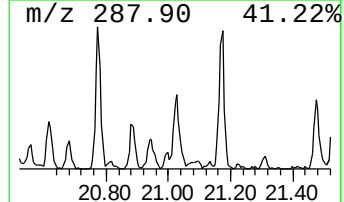
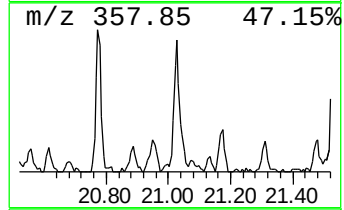
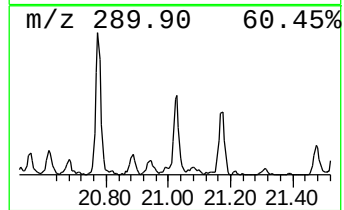
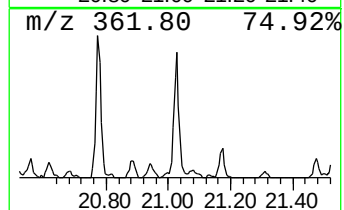
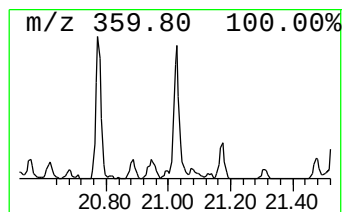
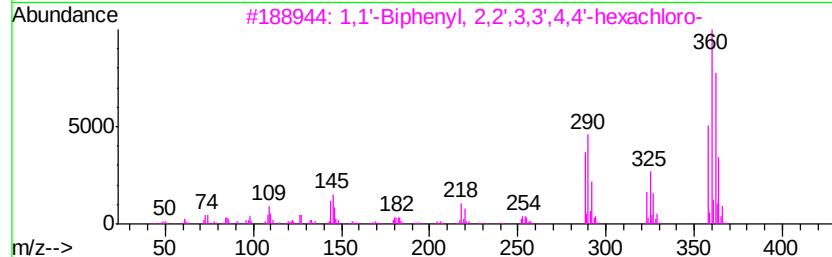
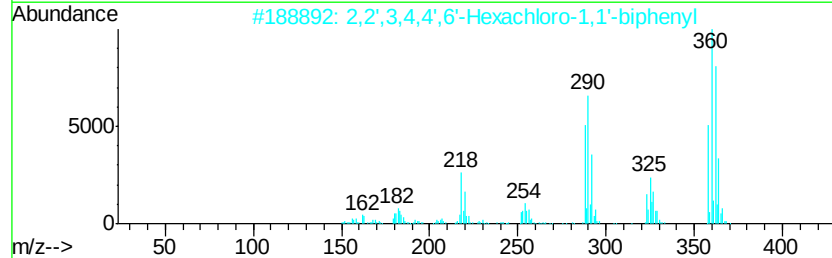
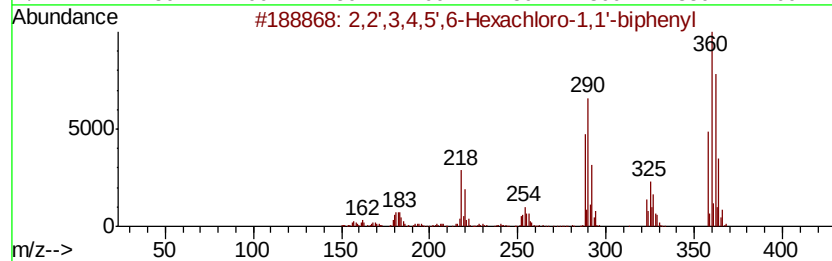
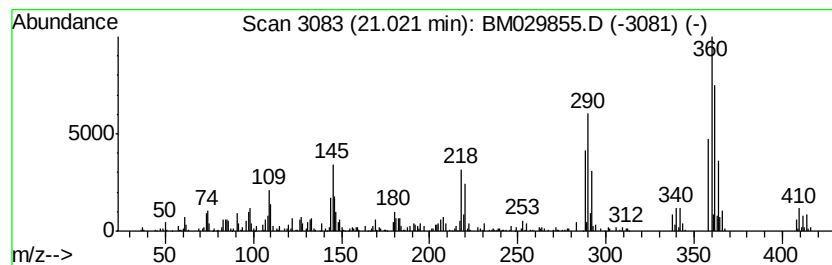
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 33 2,2',3,4,5',6-Hexachloro-1,... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.02	3.73 ng/ul	213166	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2',3,4,5',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-14-9	98
2		2,2',3,4,4',6'-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-64-4	98
3		1,1'-Biphenyl, 2,2',3,3',4,4'-he...	358	C12H4Cl6	038380-07-3	97
4		2,2',3,4',5,6-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-13-8	96
5		1,1'-Biphenyl, 2,2',3,3',4,5'-he...	358	C12H4Cl6	052663-66-8	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM050421\
 Data File : BM029855.D
 Acq On : 06 May 2021 08:55
 Operator : CG/JU
 Sample : M2240-01
 Misc :
 ALS Vial : 74 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B64

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.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...	2.99	6.2	ng/ul	214480	1	7.56	696703	20.0
1,1'-Biphenyl, 2,...	17.56	7.4	ng/ul	509364	4	16.96	1378660	20.0
1,1'-Biphenyl, 2,...	17.81	2.1	ng/ul	147264	4	16.96	1378660	20.0
1,1'-Biphenyl, 2,...	18.09	4.6	ng/ul	320044	4	16.96	1378660	20.0
Biphenyl, 2,4,4',...	18.13	2.0	ng/ul	140496	4	16.96	1378660	20.0
1,1'-Biphenyl, 2,...	18.32	5.9	ng/ul	406292	4	16.96	1378660	20.0
2,3',4,5-Tetrachl...	18.49	3.8	ng/ul	258206	4	16.96	1378660	20.0
1,1'-Biphenyl, 2,...	18.80	3.0	ng/ul	206320	4	16.96	1378660	20.0
1,1'-Biphenyl, 2,...	18.88	21.0	ng/ul	1446710	4	16.96	1378660	20.0
1,1'-Biphenyl, 2,...	19.17	23.6	ng/ul	1347750	5	21.14	1143300	20.0
1,1'-Biphenyl, 2,...	19.23	2.5	ng/ul	145107	5	21.14	1143300	20.0
1,1'-Biphenyl, 2,...	19.43	2.7	ng/ul	153895	5	21.14	1143300	20.0
1,1'-Biphenyl, 2,...	19.50	4.8	ng/ul	274508	5	21.14	1143300	20.0
3,3',4,4',5-Pent...	19.61	19.3	ng/ul	1101180	5	21.14	1143300	20.0
2,2',3,4,5,6-Hexa...	19.73	12.0	ng/ul	685521	5	21.14	1143300	20.0
2,2',3,3',5,6-Hex...	19.78	7.3	ng/ul	417313	5	21.14	1143300	20.0
1,1'-Biphenyl, 2,...	19.87	38.8	ng/ul	2219260	5	21.14	1143300	20.0
1,1'-Biphenyl, 2,...	20.07	5.5	ng/ul	312312	5	21.14	1143300	20.0
2,3,3',4',5',6-He...	20.16	52.2	ng/ul	2982270	5	21.14	1143300	20.0
1,1'-Biphenyl, 2,...	20.20	16.3	ng/ul	929732	5	21.14	1143300	20.0
1,1'-Biphenyl, 2,...	20.31	17.3	ng/ul	987148	5	21.14	1143300	20.0
1,1'-Biphenyl, 2,...	20.47	53.1	ng/ul	3036080	5	21.14	1143300	20.0
1,1'-Biphenyl, 2,...	20.62	17.6	ng/ul	1008400	5	21.14	1143300	20.0
2,3,3',4,4',5',6-...	20.68	10.5	ng/ul	597867	5	21.14	1143300	20.0
1,1'-Biphenyl, 2,...	20.77	4.6	ng/ul	265367	5	21.14	1143300	20.0
1,1'-Biphenyl, 2,...	20.89	17.4	ng/ul	995552	5	21.14	1143300	20.0
1,1'-Biphenyl, 2,...	20.99	5.4	ng/ul	306517	5	21.14	1143300	20.0
2,2',3,4,5',6-Hex...	21.02	3.7	ng/ul	213166	5	21.14	1143300	20.0