

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
 Data File : BM026347.D
 Acq On : 11 Jun 2020 01:35
 Operator : JU/CG
 Sample : L2850-06
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ETTX9

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.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	2.922	3	6	15	rVB	568705	536699	4.75%	0.516%
2	3.240	56	60	64	rVB2	30190	35946	0.32%	0.035%
3	3.693	134	137	141	rVB	10045	13206	0.12%	0.013%
4	5.934	514	518	525	rVB6	6673	12902	0.11%	0.012%
5	7.416	761	770	779	rBV	610091	920686	8.15%	0.885%
6	7.616	800	804	812	rVV7	6596	12309	0.11%	0.012%
7	10.175	1232	1239	1249	rBV	711750	1166202	10.32%	1.121%
8	14.069	1894	1901	1914	rBV2	903769	1291542	11.43%	1.241%
9	14.204	1918	1924	1934	rVV	55032	81993	0.73%	0.079%
10	15.028	2059	2064	2070	rVB2	13362	20346	0.18%	0.020%
11	15.339	2112	2117	2125	rBV	845942	1116282	9.88%	1.073%
12	15.786	2188	2193	2199	rBV	124549	156138	1.38%	0.150%
13	15.951	2215	2221	2227	rBV	331000	425670	3.77%	0.409%
14	16.069	2235	2241	2255	rBV	2082821	2626372	23.24%	2.524%
15	16.363	2286	2291	2297	rBV	397010	520513	4.61%	0.500%
16	16.716	2345	2351	2355	rBV	4953730	6627968	58.65%	6.371%
17	16.751	2355	2357	2363	rVV	1659187	1803563	15.96%	1.734%
18	16.822	2363	2369	2373	rVV	1451784	1952919	17.28%	1.877%
19	16.863	2373	2376	2384	rVB	240722	346332	3.06%	0.333%
20	17.010	2394	2401	2413	rVB2	2347747	3598703	31.85%	3.459%
21	17.127	2417	2421	2428	rBV4	12275	20016	0.18%	0.019%
22	17.204	2430	2434	2441	rVB3	33793	45785	0.41%	0.044%
23	17.286	2443	2448	2452	rBV	640552	916589	8.11%	0.881%
24	17.322	2452	2454	2461	rVV	245080	244359	2.16%	0.235%
25	17.380	2461	2464	2466	rVV3	11005	11656	0.10%	0.011%
26	17.433	2466	2473	2485	rVV	5882163	11300295	100.00%	10.862%
27	17.563	2487	2495	2501	rVV2	3303262	4857360	42.98%	4.669%
28	17.627	2501	2506	2510	rVV	232995	298412	2.64%	0.287%
29	17.680	2510	2515	2520	rVV	1592793	2089448	18.49%	2.008%
30	17.733	2520	2524	2530	rVB	818031	1068094	9.45%	1.027%
31	17.845	2537	2543	2548	rBV	346211	444826	3.94%	0.428%
32	17.910	2548	2554	2560	rVV	4199014	5539154	49.02%	5.324%
33	17.969	2560	2564	2567	rVV	2720472	3356555	29.70%	3.226%
34	18.010	2567	2571	2577	rVB	2213529	2671753	23.64%	2.568%

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
 Data File : BM026347.D
 Acq On : 11 Jun 2020 01:35
 Operator : JU/CG
 Sample : L2850-06
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ETTX9

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
 Title : SVOA CALIBRATION

35	18.192	2596	2602	2607	rBV	3819691	5123597	45.34%	4.925%
36	18.239	2607	2610	2615	rVV	1163552	1514843	13.41%	1.456%
37	18.298	2615	2620	2623	rVV	971252	1315453	11.64%	1.264%
38	18.333	2623	2626	2629	rVV	1091038	1411052	12.49%	1.356%
39	18.369	2629	2632	2638	rVV	2499334	3158613	27.95%	3.036%
40	18.433	2639	2643	2645	rVV2	46565	59188	0.52%	0.057%
41	18.469	2645	2649	2656	rVB	770810	1009572	8.93%	0.970%
42	18.551	2656	2663	2671	rBV	120788	164249	1.45%	0.158%
43	18.627	2671	2676	2680	rBV	175277	227201	2.01%	0.218%
44	18.686	2680	2686	2690	rBV	2170183	2625561	23.23%	2.524%
45	18.739	2690	2695	2698	rVV	4027488	5824719	51.54%	5.599%
46	18.780	2698	2702	2707	rVV2	4046155	5506270	48.73%	5.293%
47	18.851	2709	2714	2718	rVV	328764	405017	3.58%	0.389%
48	18.886	2718	2720	2725	rVB	61858	73566	0.65%	0.071%
49	18.992	2731	2738	2744	rBV	2064964	4226411	37.40%	4.062%
50	19.051	2744	2748	2755	rVV	1540648	2253393	19.94%	2.166%
51	19.116	2755	2759	2765	rVB	762825	902575	7.99%	0.868%
52	19.192	2768	2772	2776	rBV5	32749	48361	0.43%	0.046%
53	19.245	2777	2781	2785	rVB	124574	146775	1.30%	0.141%
54	19.316	2788	2793	2799	rBV	636434	788066	6.97%	0.757%
55	19.392	2799	2806	2811	rVV	814031	1074794	9.51%	1.033%
56	19.445	2811	2815	2819	rVV	484382	573912	5.08%	0.552%
57	19.504	2819	2825	2829	rVV	1451345	1749787	15.48%	1.682%
58	19.545	2829	2832	2840	rVB	328785	417200	3.69%	0.401%
59	19.645	2842	2849	2859	rBV	401091	576796	5.10%	0.554%
60	19.721	2859	2862	2865	rVV	58967	68685	0.61%	0.066%
61	19.763	2865	2869	2874	rVV3	205236	352495	3.12%	0.339%
62	19.821	2874	2879	2884	rVB	1230934	1421144	12.58%	1.366%
63	19.898	2889	2892	2895	rVB4	20552	20227	0.18%	0.019%
64	19.951	2897	2901	2908	rBV3	85458	167001	1.48%	0.161%
65	20.051	2913	2918	2922	rBV	173483	203889	1.80%	0.196%
66	20.098	2922	2926	2927	rBV2	113615	112085	0.99%	0.108%
67	20.127	2927	2931	2936	rVB	836867	983849	8.71%	0.946%
68	20.204	2940	2944	2948	rBV3	57237	71497	0.63%	0.069%
69	20.274	2952	2956	2958	rBV5	23840	32988	0.29%	0.032%
70	20.298	2958	2960	2963	rVB4	22447	20915	0.19%	0.020%
71	20.363	2964	2971	2979	rBV	220255	377498	3.34%	0.363%

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETTX9

Integration Parameters: LSCINT.P

Integrator: RTE
Smoothing : OFF
Sampling : 1
Start Thrs: 0.2
Stop Thrs : 0
Filtering: 5
Min Area: 0.1 % of largest Peak
Max Peaks: 100
Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Title : SVOA CALIBRATION

72	20.451	2983	2986	2989	rVV4	25199	28292	0.25%	0.027%
73	20.515	2994	2997	3002	rVB6	29256	37434	0.33%	0.036%
74	20.668	3020	3023	3028	rBV2	74588	87296	0.77%	0.084%
75	20.786	3040	3043	3047	rVB5	35640	38173	0.34%	0.037%
76	20.921	3064	3066	3070	rVB3	42450	38877	0.34%	0.037%
77	21.033	3080	3085	3089	rBV	1011634	1279029	11.32%	1.229%
78	21.068	3089	3091	3097	rVB5	77378	94959	0.84%	0.091%
79	23.127	3436	3441	3449	rVB	765591	1293570	11.45%	1.243%

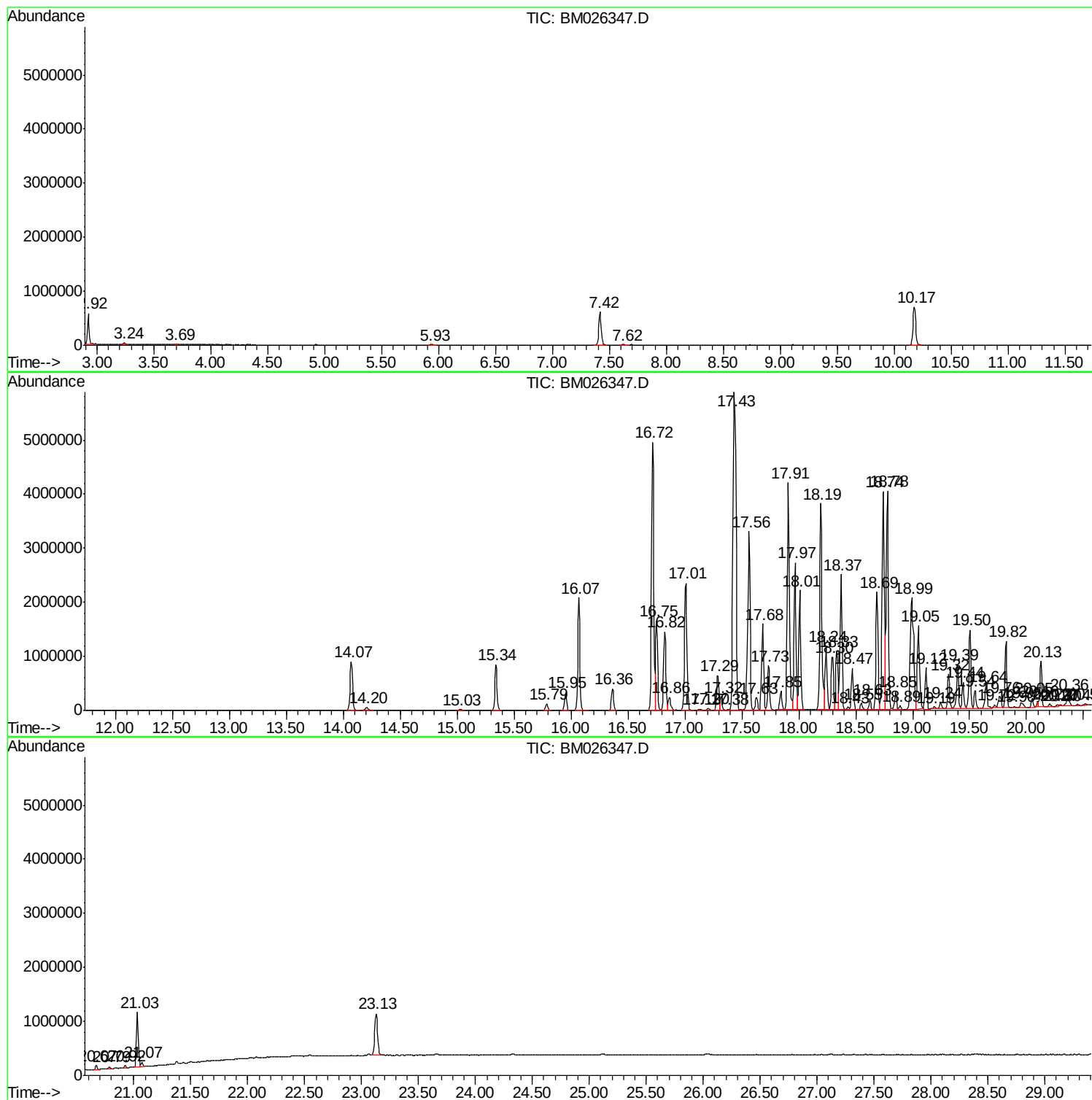
Sum of corrected areas: 104037467

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETTX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETTX9

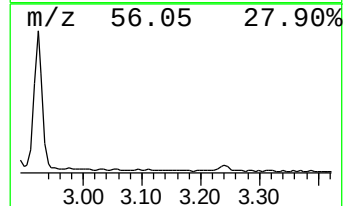
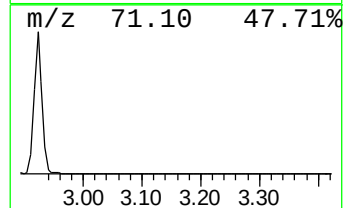
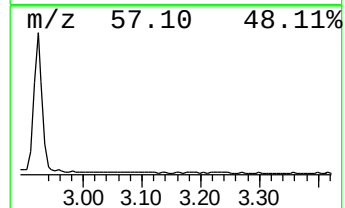
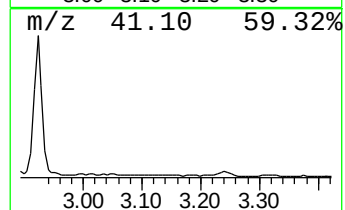
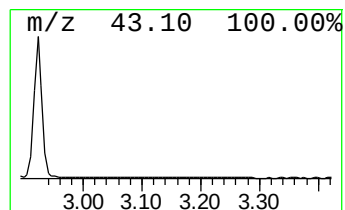
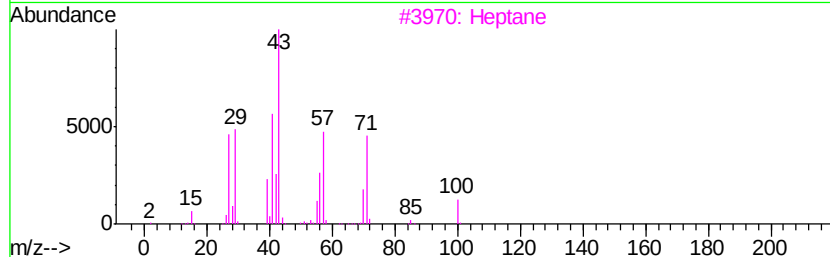
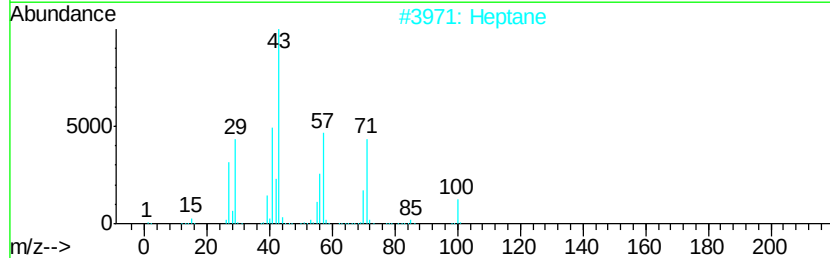
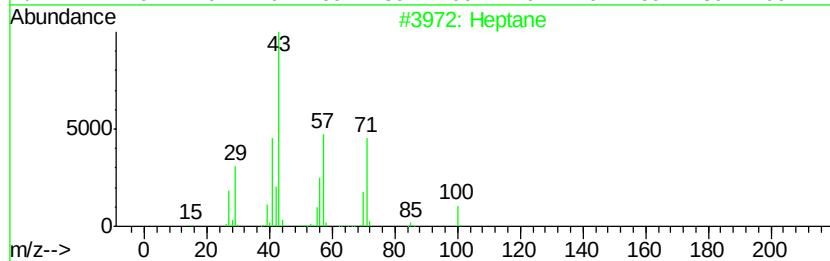
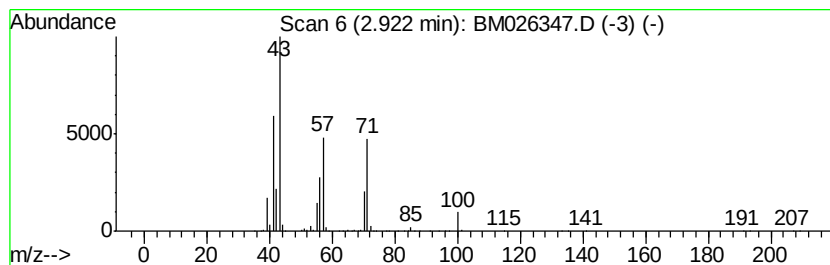
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.92	11.66 ng/ul	536699	1,4-Dichlorobenzene-d4	7.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane	100	C7H16	000142-82-5	94
2		Heptane	100	C7H16	000142-82-5	94
3		Heptane	100	C7H16	000142-82-5	91
4		Heptane	100	C7H16	000142-82-5	90
5		Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETTX9

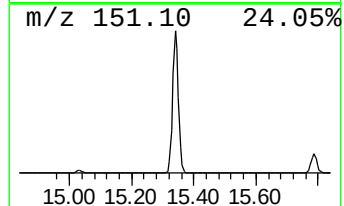
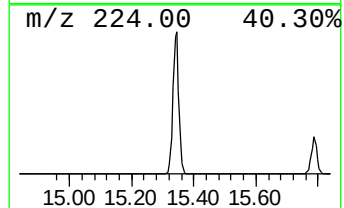
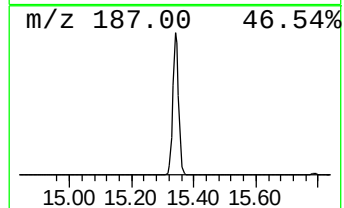
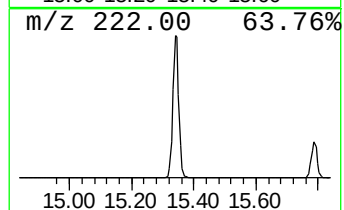
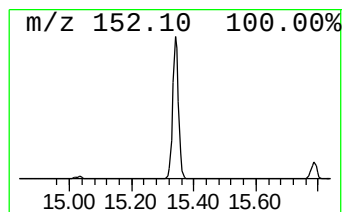
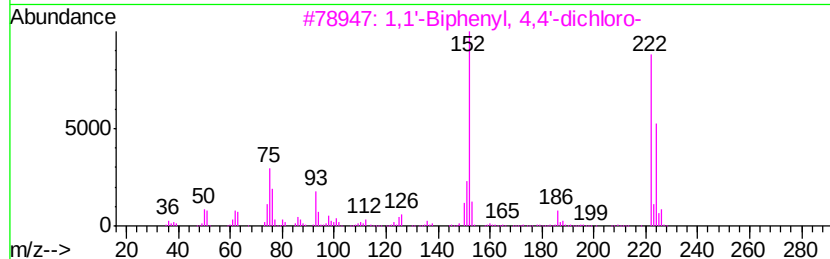
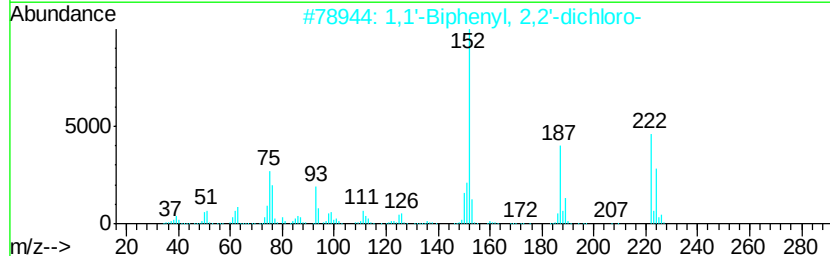
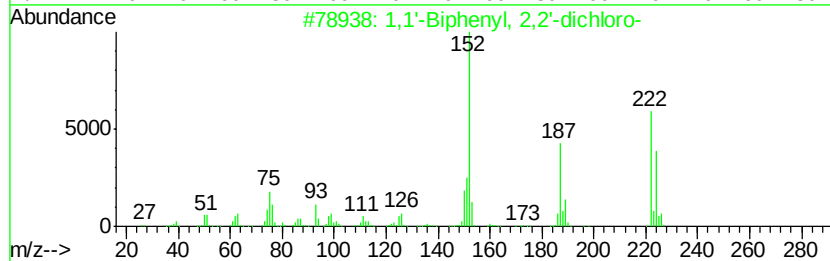
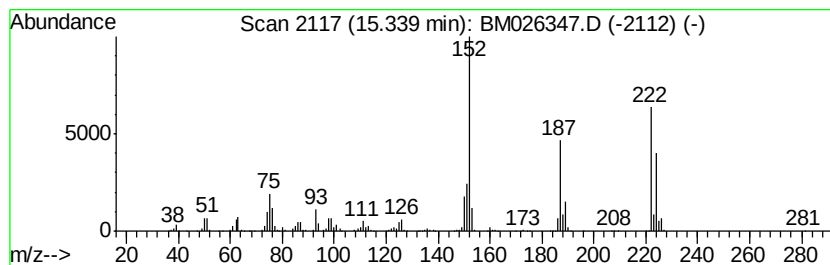
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1,1'-Biphenyl, 2,2'-dichloro- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.34	17.29 ng/ul	1116280	Acenaphthene-d10	14.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	99
2		1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	99
3		1,1'-Biphenyl, 4,4'-dichloro-	222	C12H8Cl2	002050-68-2	98
4		1,1'-Biphenyl, 2,4'-dichloro-	222	C12H8Cl2	034883-43-7	98
5		1,1'-Biphenyl, 2,5-dichloro-	222	C12H8Cl2	034883-39-1	97



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETTX9

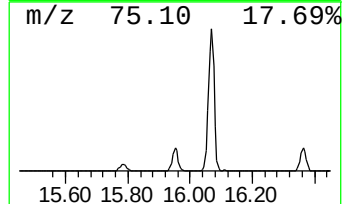
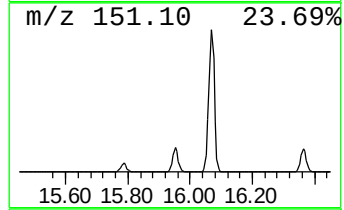
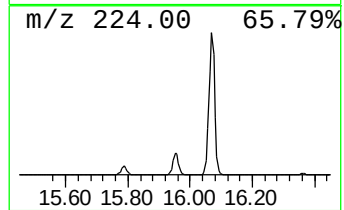
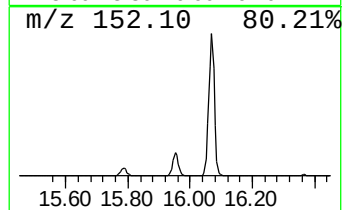
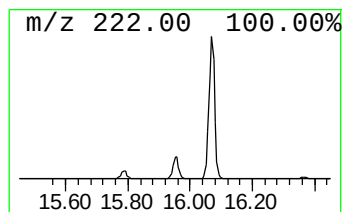
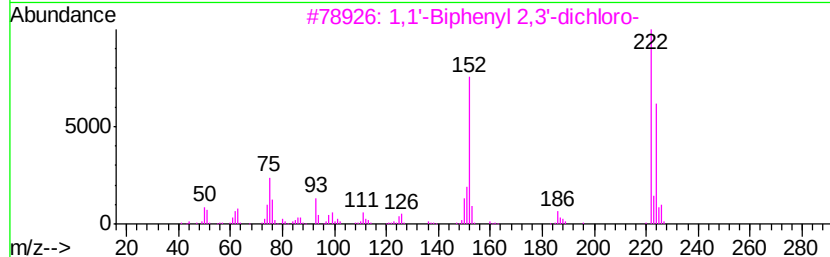
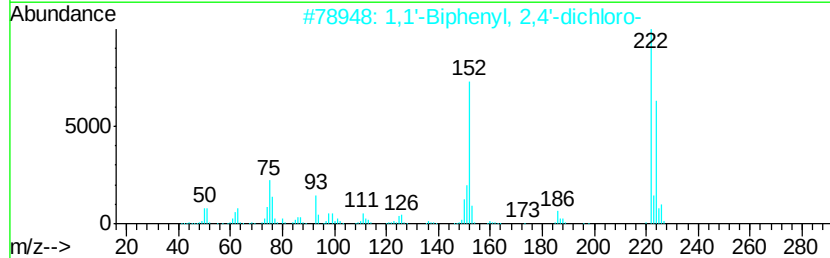
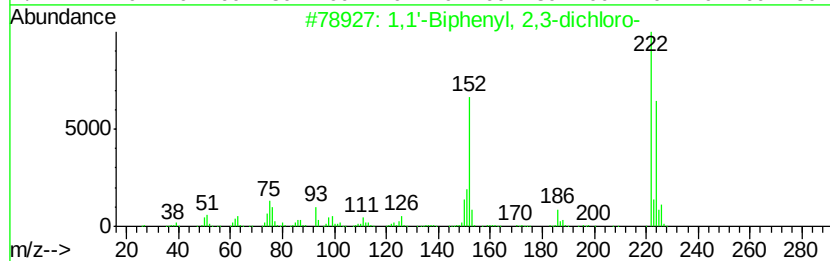
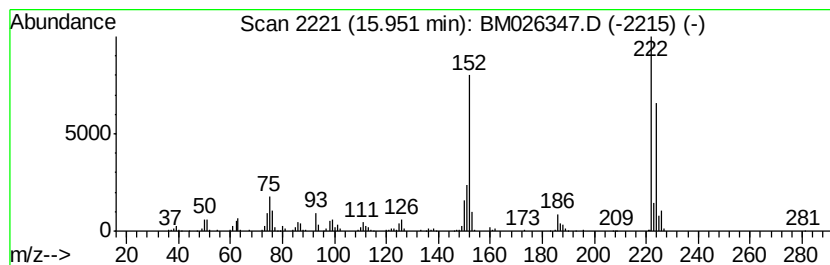
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1,1'-Biphenyl, 2,3-dichloro- Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.95	4.36 ng/ul	425670	Phenanthrene-d10	16.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3-dichloro-	222	C12H8Cl2	016605-91-7	99
2		1,1'-Biphenyl, 2,4'-dichloro-	222	C12H8Cl2	034883-43-7	99
3		1,1'-Biphenyl 2,3'-dichloro-	222	C12H8Cl2	025569-80-6	99
4		1,1'-Biphenyl, 3,5-dichloro-	222	C12H8Cl2	034883-41-5	98
5		1,1'-Biphenyl, 2,5-dichloro-	222	C12H8Cl2	034883-39-1	98



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETTX9

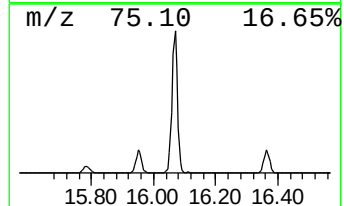
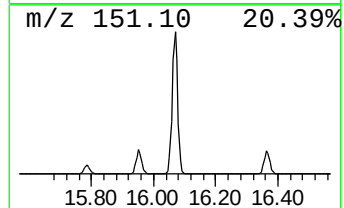
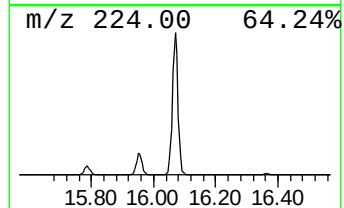
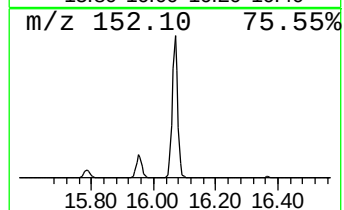
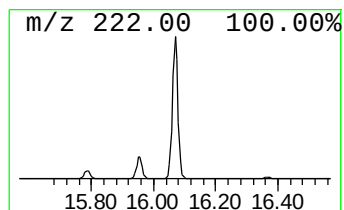
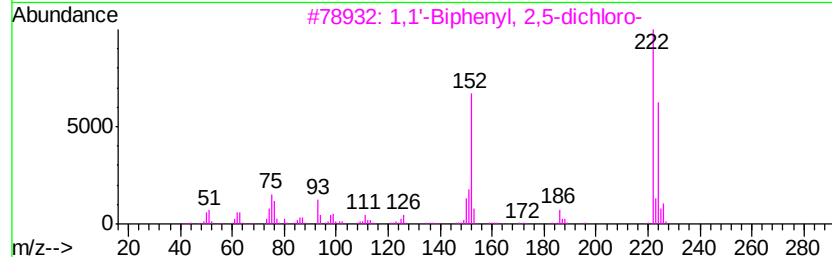
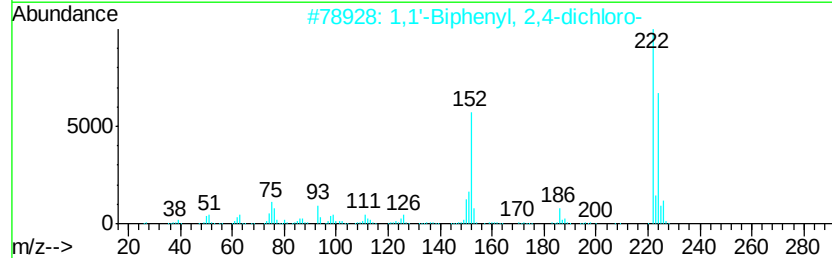
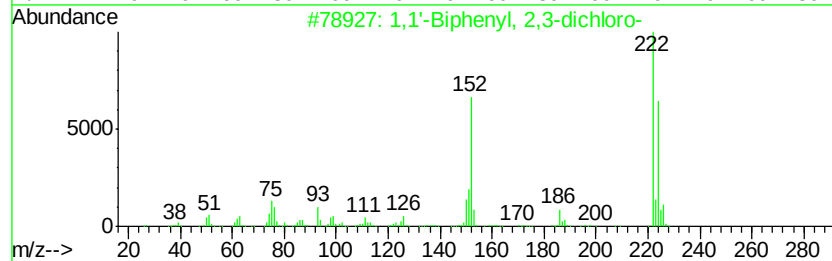
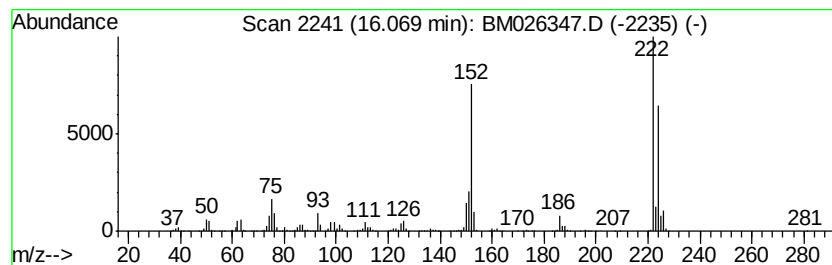
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 1,1'-Biphenyl, 2,4-dichloro- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.07	26.90 ng/ul	2626370	Phenanthrene-d10	16.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3-dichloro-	222	C12H8Cl2	016605-91-7	99
2		1,1'-Biphenyl, 2,4-dichloro-	222	C12H8Cl2	033284-50-3	99
3		1,1'-Biphenyl, 2,5-dichloro-	222	C12H8Cl2	034883-39-1	99
4		1,1'-Biphenyl, 3,3'-dichloro-	222	C12H8Cl2	002050-67-1	99
5		1,1'-Biphenyl, 2,4'-dichloro-	222	C12H8Cl2	034883-43-7	98



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

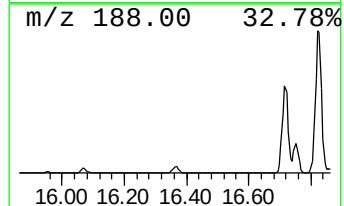
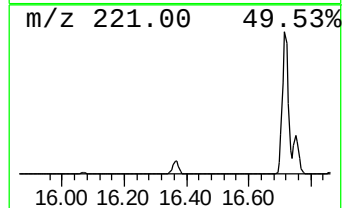
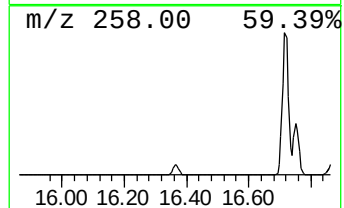
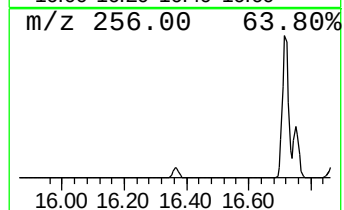
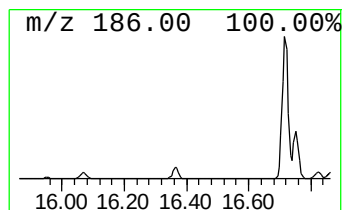
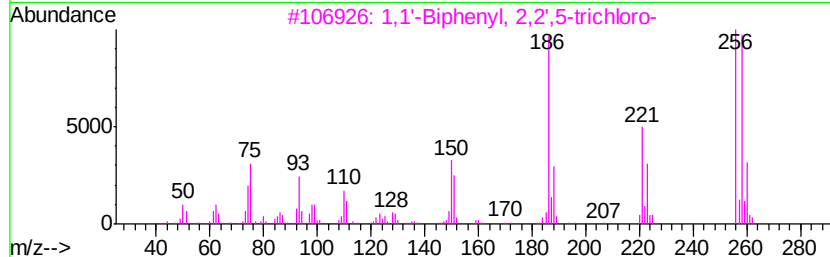
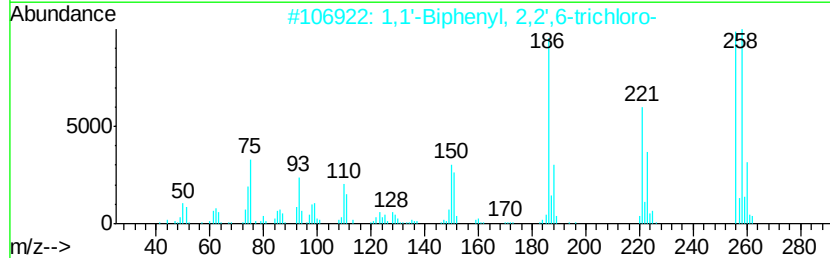
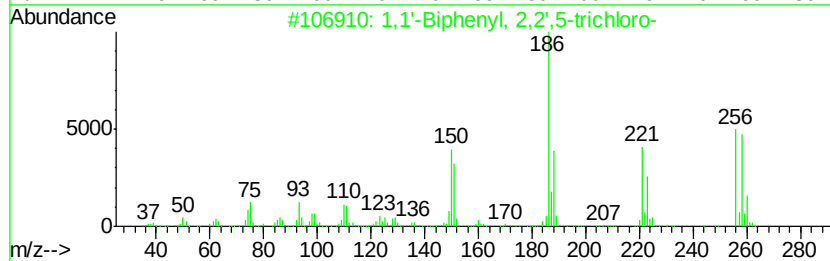
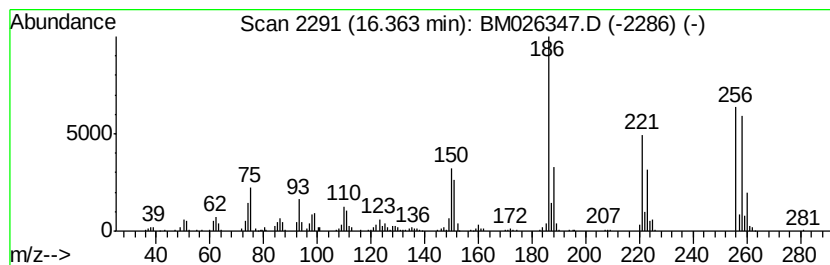
Instrument :
BNA_M
ClientSampleId :
ETTX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,1'-Biphenyl, 2,2',5-trich... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.36	5.33 ng/ul	520513	Phenanthrene-d10	16.82	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99
2	1,1'-Biphenyl, 2,2',6-trichloro-	256	C12H7Cl3	038444-73-4	99
3	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99
4	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	98
5	2,2',4-Trichloro-1,1'-biphenyl	256	C12H7Cl3	037680-66-3	97



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

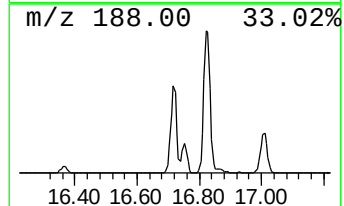
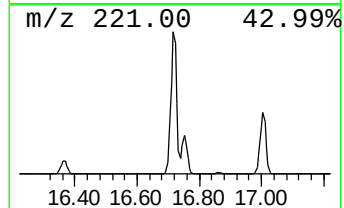
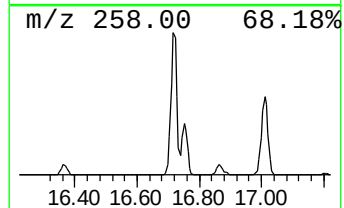
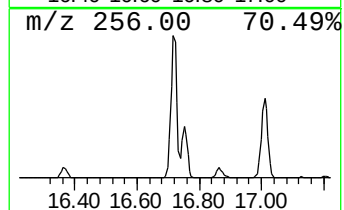
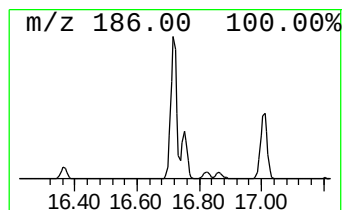
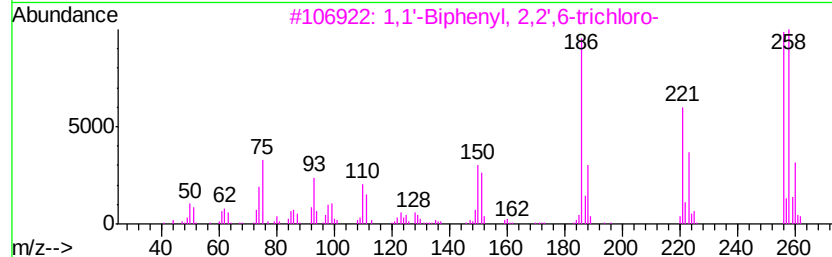
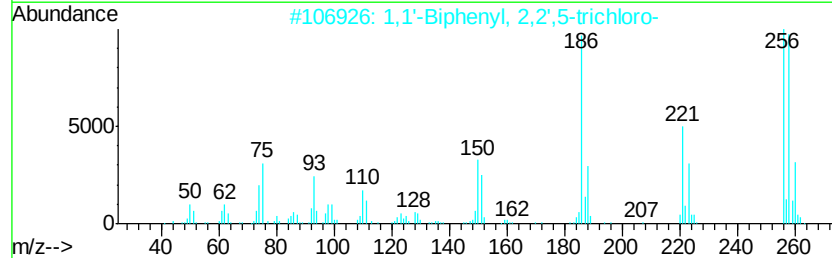
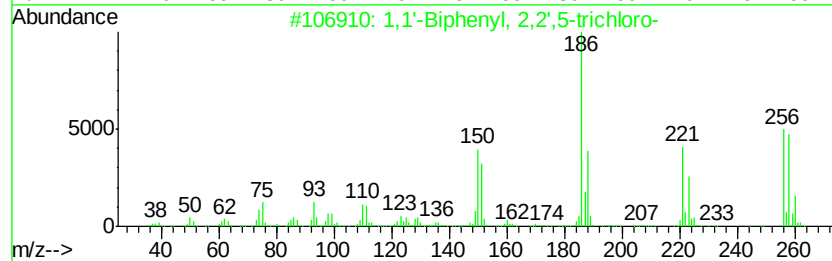
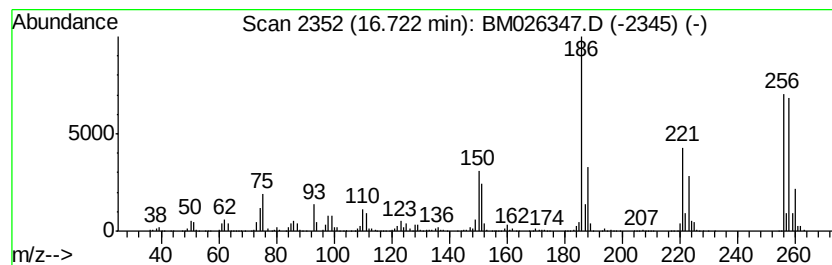
Instrument :
BNA_M
ClientSampleId :
ETTX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD		R.T.		
16.72	67.88 ng/ul	6627970	Phenanthrene-d10		16.82		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'	-Biphenyl,	2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99
2	1,1'	-Biphenyl,	2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99
3	1,1'	-Biphenyl,	2,2',6-trichloro-	256	C12H7Cl3	038444-73-4	99
4	1,1'	-Biphenyl,	2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99
5	2,2',3-	Trichloro-1,1'-biphenyl		256	C12H7Cl3	038444-78-9	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETTX9

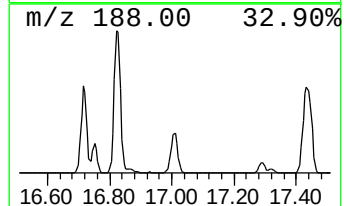
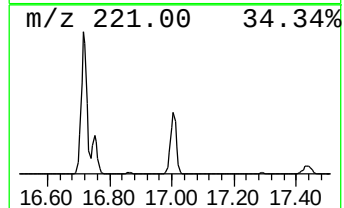
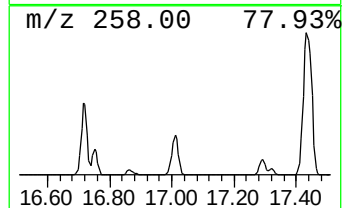
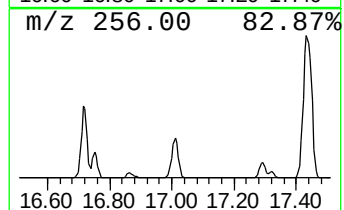
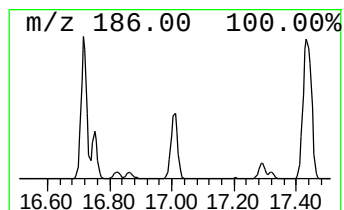
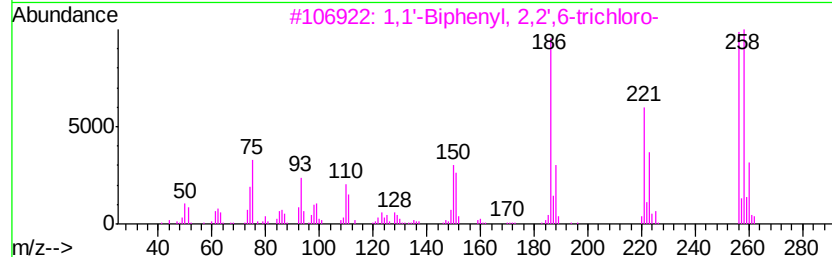
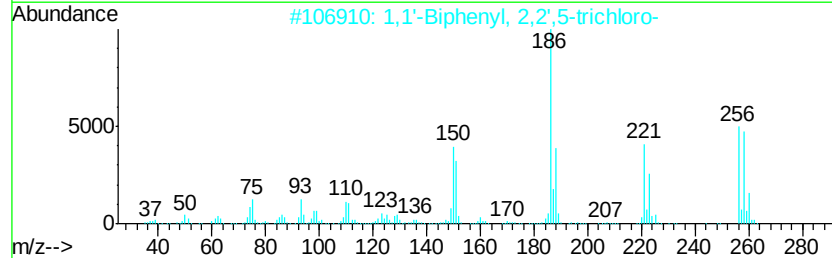
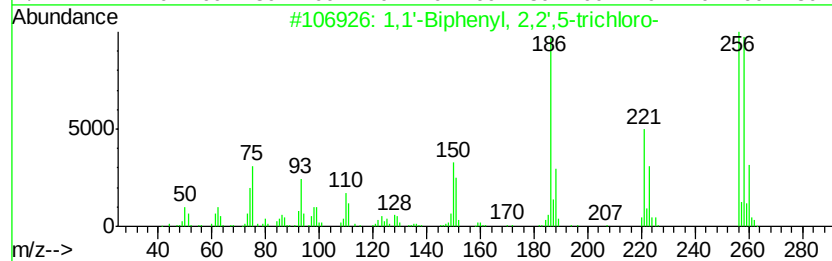
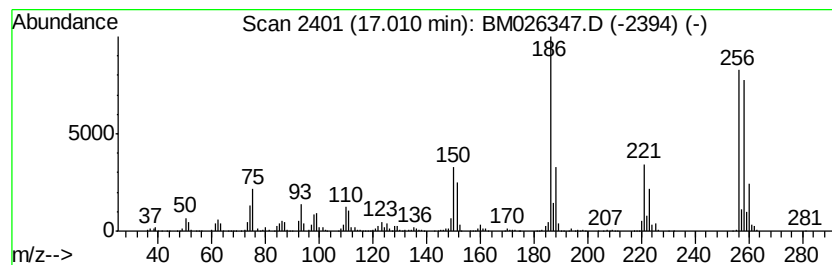
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 2,2',4-Trichloro-1,1'-biphenyl Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.01	36.85 ng/ul	3598700	Phenanthrene-d10	16.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99
2		1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99
3		1,1'-Biphenyl, 2,2',6-trichloro-	256	C12H7Cl3	038444-73-4	99
4		2,2',4-Trichloro-1,1'-biphenyl	256	C12H7Cl3	037680-66-3	99
5		1,1'-Biphenyl, 2,4,5-trichloro-	256	C12H7Cl3	015862-07-4	98



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

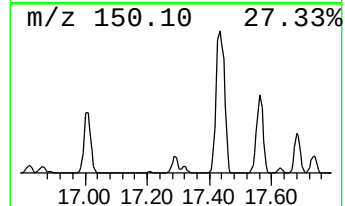
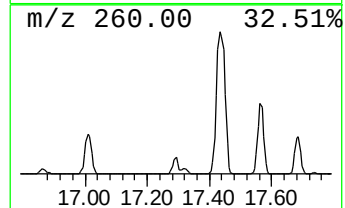
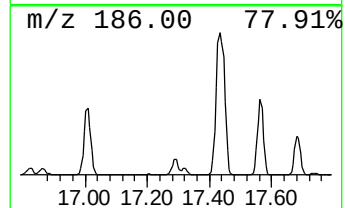
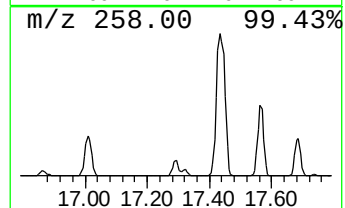
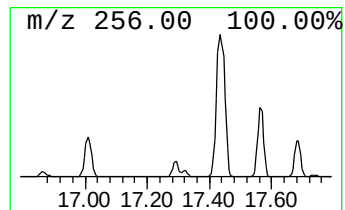
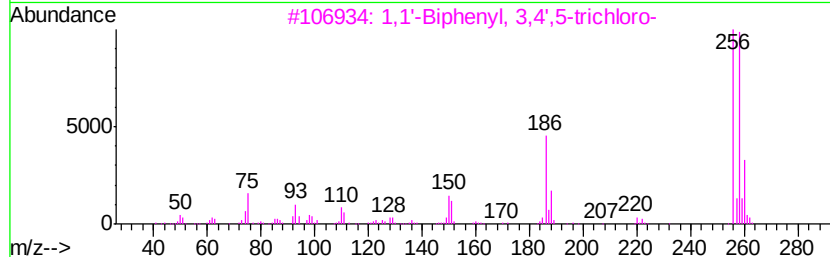
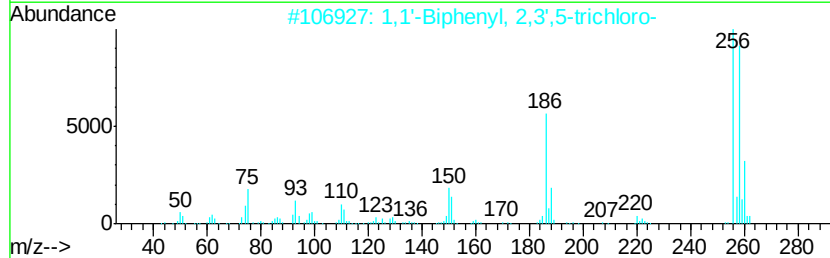
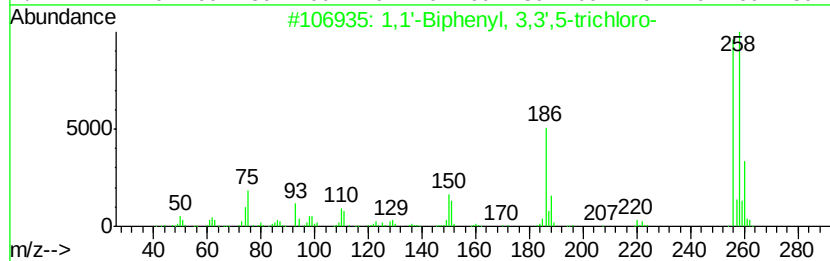
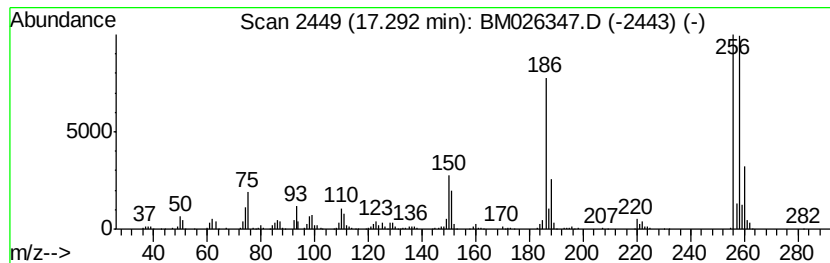
Instrument :
BNA_M
ClientSampleId :
ETTX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 1,1'-Biphenyl, 3,4',5-trich... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.29	9.39 ng/ul	916589	Phenanthrene-d10	16.82	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	99
2	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99
3	1,1'-Biphenyl, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	99
4	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	99
5	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETTX9

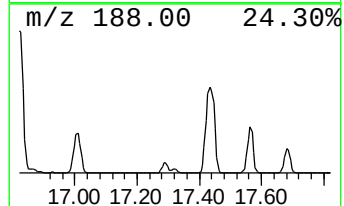
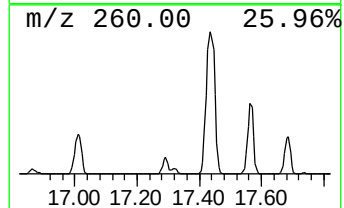
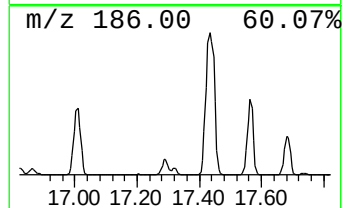
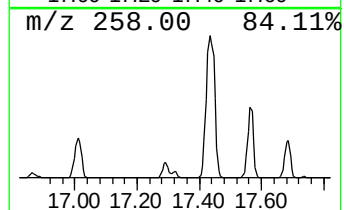
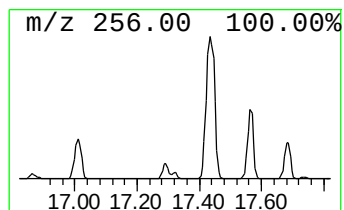
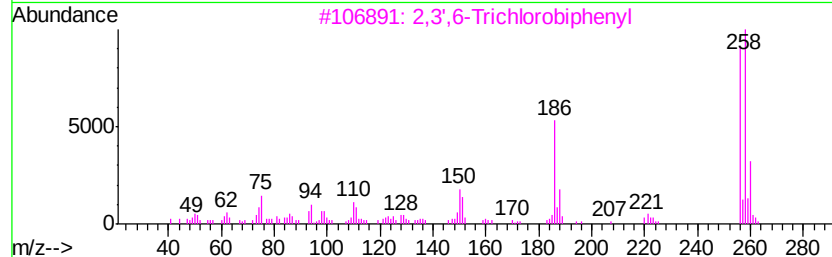
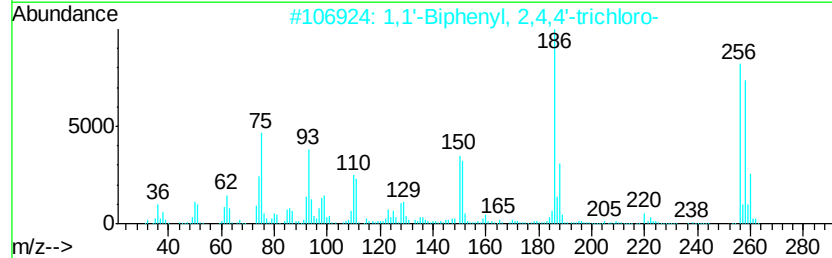
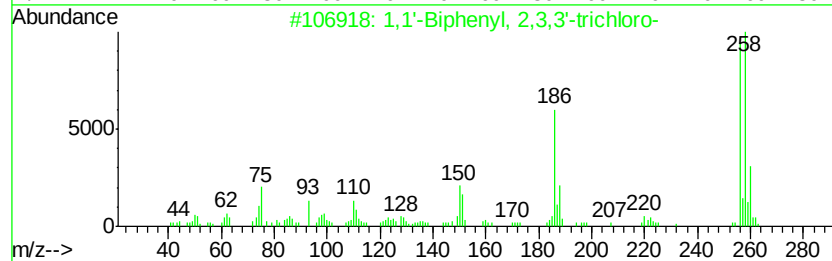
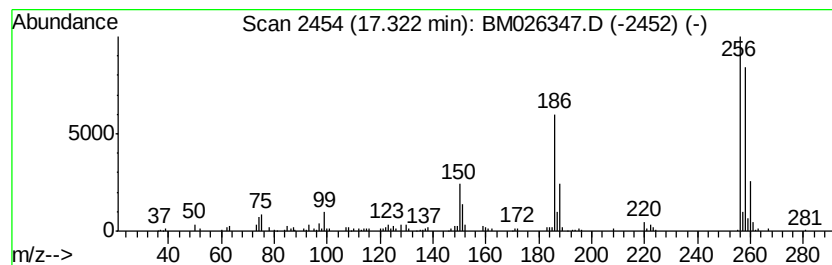
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 1,1'-Biphenyl, 2,3,3'-trich... Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.32	2.50 ng/ul	244359	Phenanthrene-d10	16.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	94
2			1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	94
3			2,3',6-Trichlorobiphenyl	256	C12H7Cl3	038444-76-7	94
4			3,3',4-Trichloro-1,1'-biphenyl	256	C12H7Cl3	037680-69-6	94
5			2,4',6-Trichloro-1,1'-biphenyl	256	C12H7Cl3	038444-77-8	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

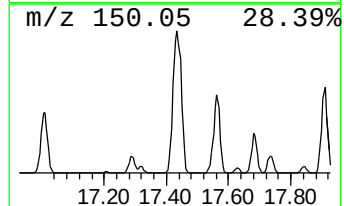
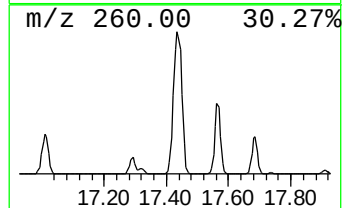
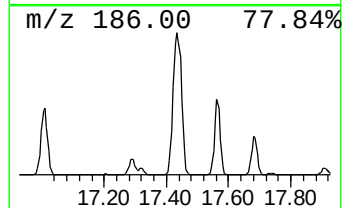
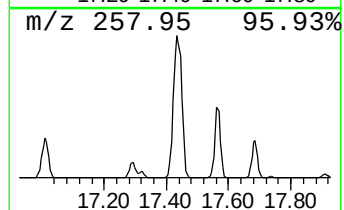
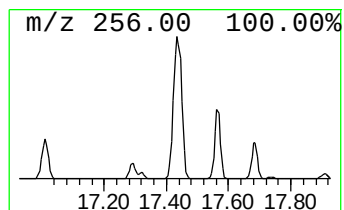
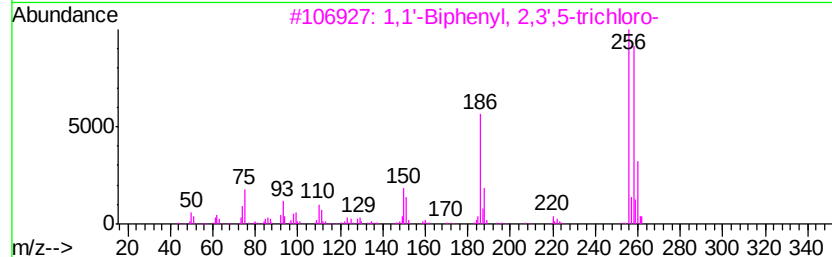
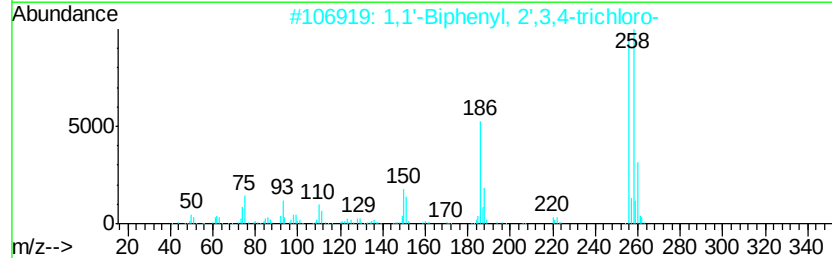
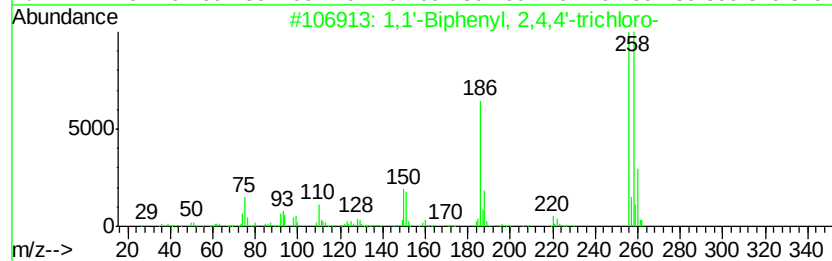
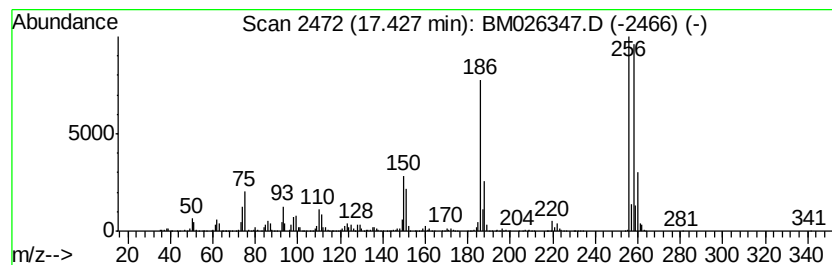
Instrument :
BNA_M
ClientSampleId :
ETTX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
17.43	115.73 ng/ul	11300300	Phenanthrene-d10	16.82		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl,	2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99
2	1,1'-Biphenyl,	2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99
3	1,1'-Biphenyl,	2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99
4	1,1'-Biphenyl,	3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	99
5	1,1'-Biphenyl,	2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETTX9

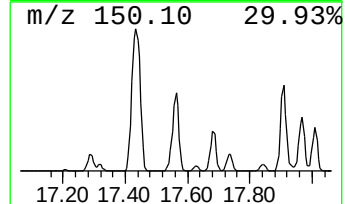
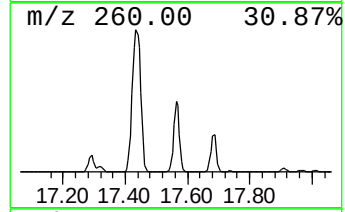
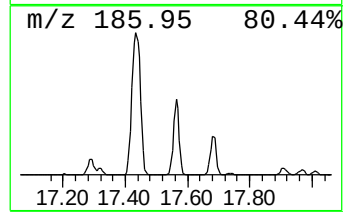
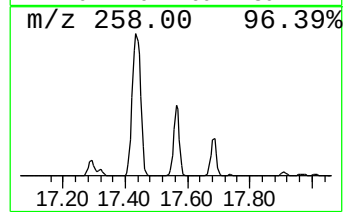
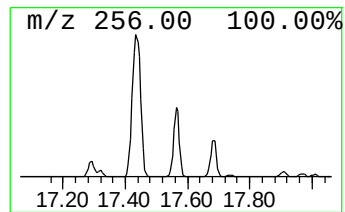
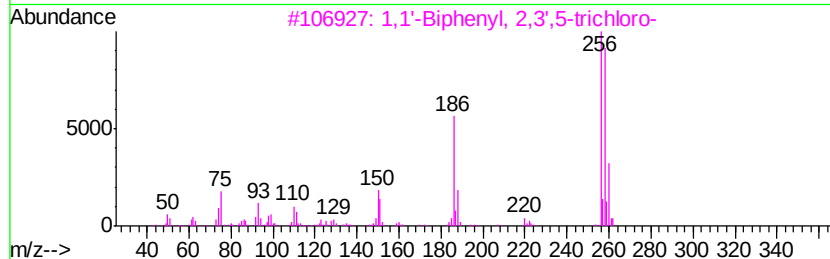
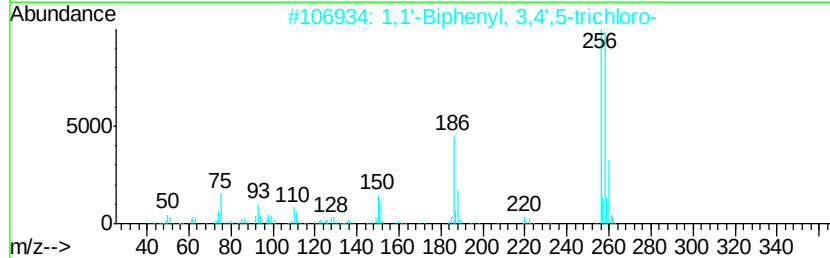
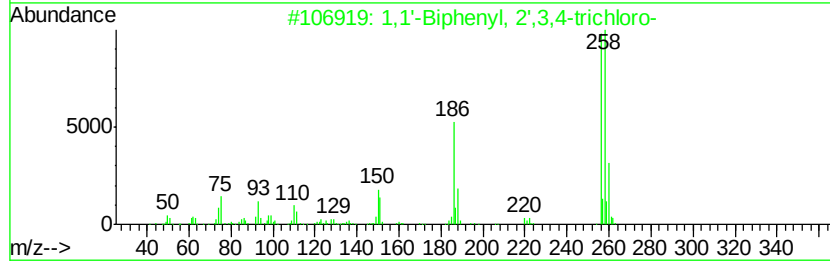
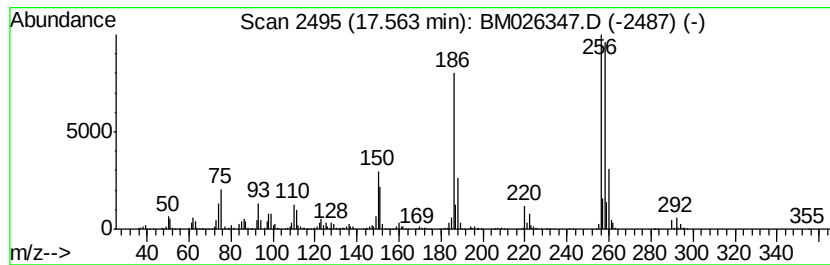
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 1,1'-Biphenyl, 3,3',5-trich... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.56	49.74 ng/ul	4857360	Phenanthrene-d10	16.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99
2		1,1'-Biphenyl, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	99
3		1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99
4		1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	99
5		1,1'-Biphenyl, 2,3,6-trichloro-	256	C12H7Cl3	055702-45-9	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETTX9

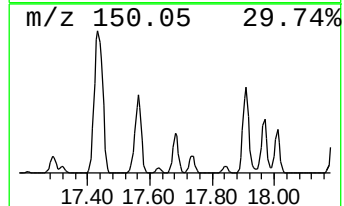
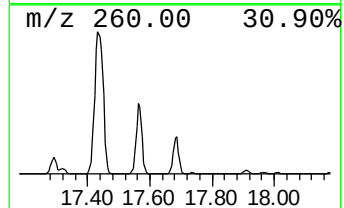
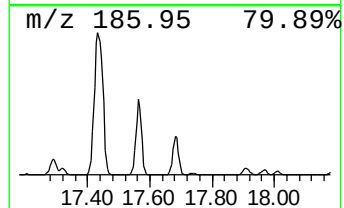
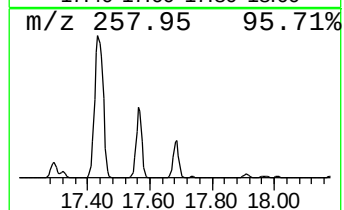
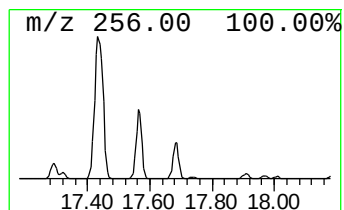
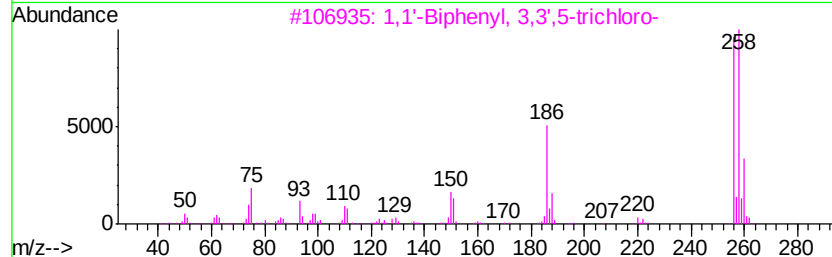
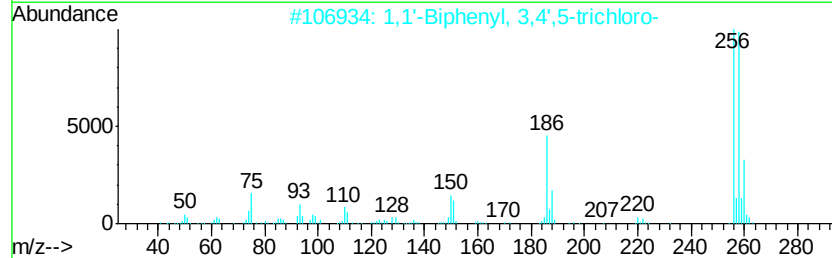
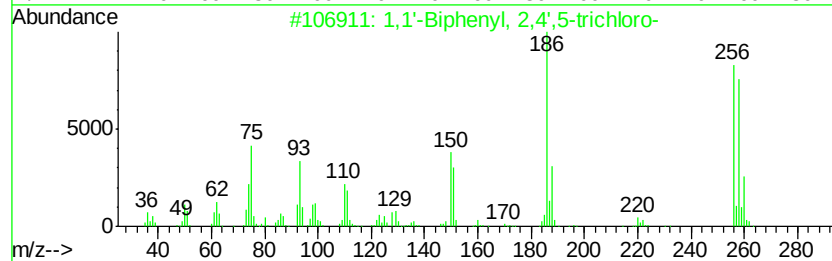
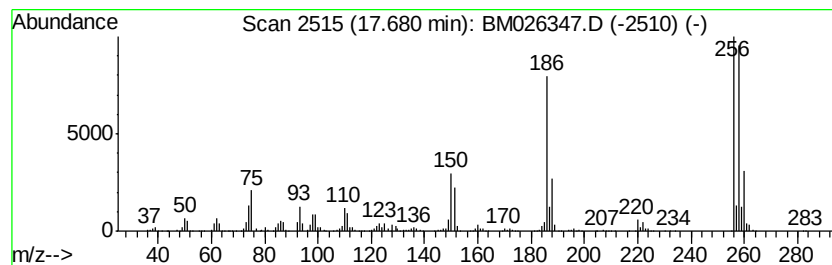
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 14 1,1'-Biphenyl, 2,4',5-trich... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.68	21.40 ng/ul	2089450	Phenanthrene-d10	16.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	99
2		1,1'-Biphenyl, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	99
3		1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	99
4		1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99
5		1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

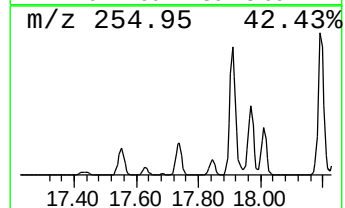
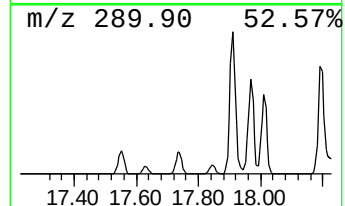
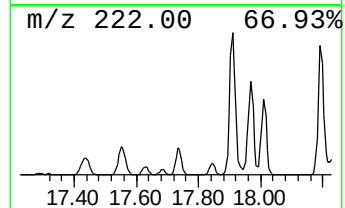
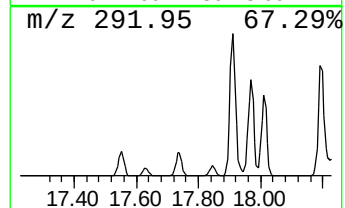
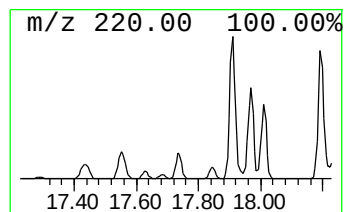
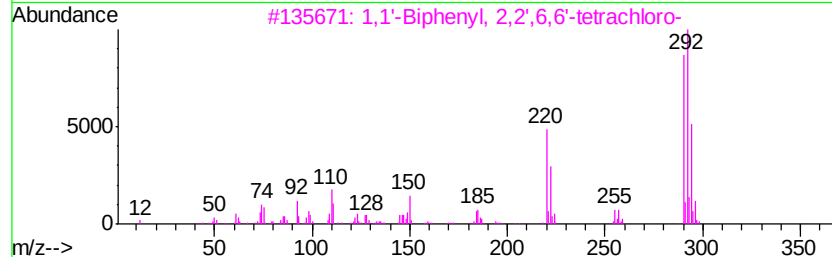
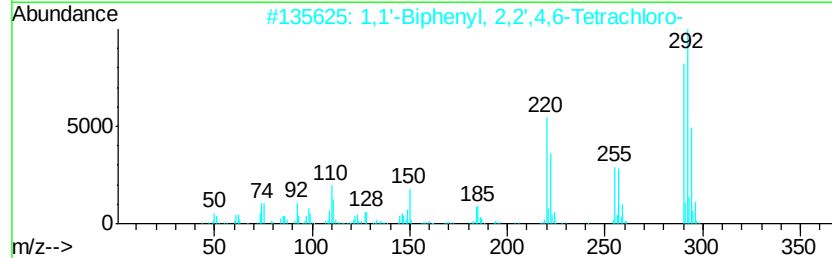
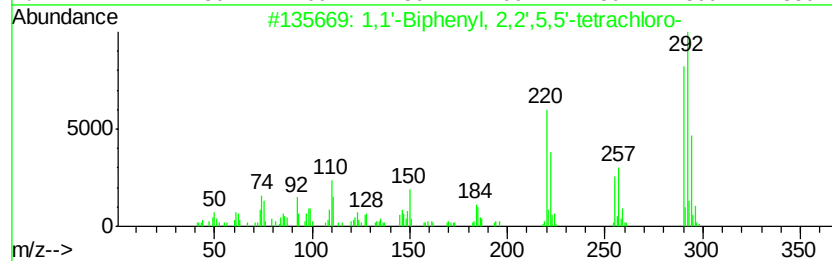
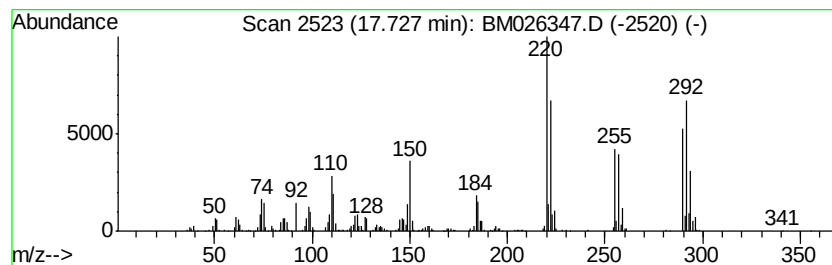
Instrument :
BNA_M
ClientSampleId :
ETTX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 15 1,1'-Biphenyl, 2,2',5,5'-te... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.		
17.73	10.94 ng/ul	1068090	Phenanthrene-d10	16.82		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	99	
2	1,1'-Biphenyl, 2,2',4,6-Tetrachl...	290	C12H6Cl4	062796-65-0	99	
3	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99	
4	1,1'-Biphenyl, 2,2',5,6'-Tetrach...	290	C12H6Cl4	041464-41-9	99	
5	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

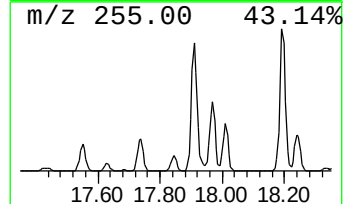
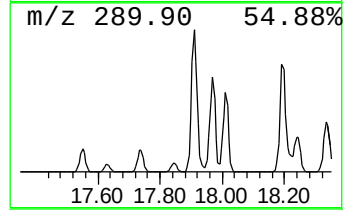
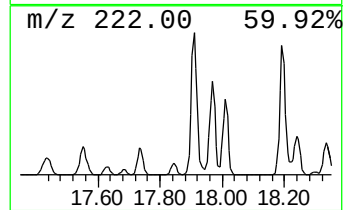
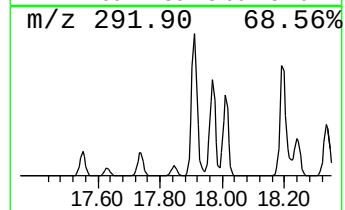
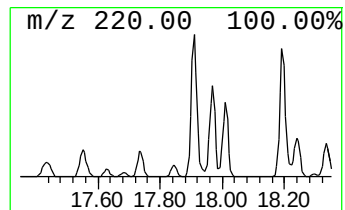
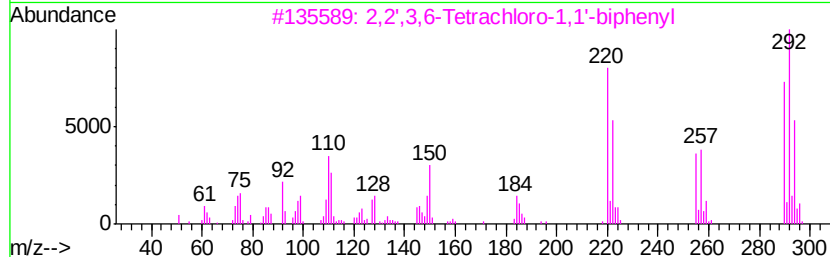
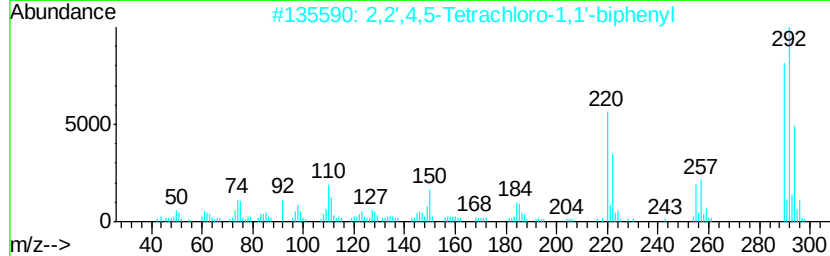
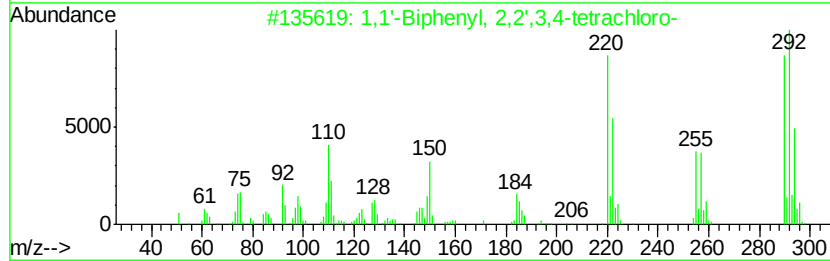
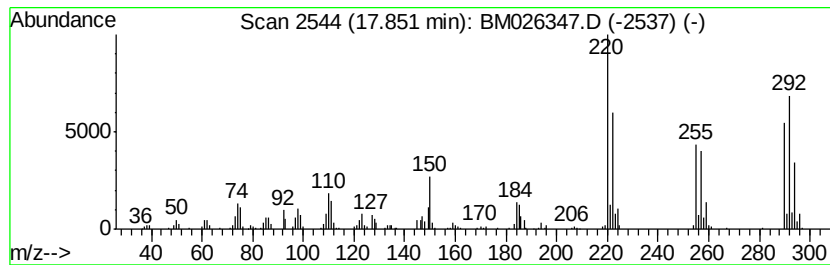
Instrument :
BNA_M
ClientSampleId :
ETTX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 16 1,1'-Biphenyl, 2,2',3,4-tet... Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.85	4.56 ng/ul	444826	Phenanthrene-d10	16.82	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,4-tetrachl...	290	C12H6Cl4	052663-59-9	99
2	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99
3	2,2',3,6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-45-7	99
4	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99
5	1,1'-Biphenyl, 2,2',3,6'-tetrach...	290	C12H6Cl4	041464-47-5	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETTX9

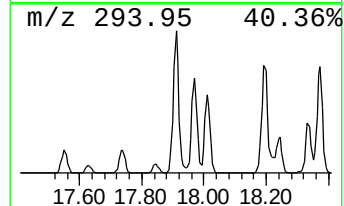
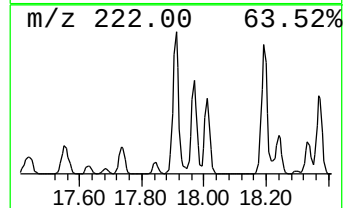
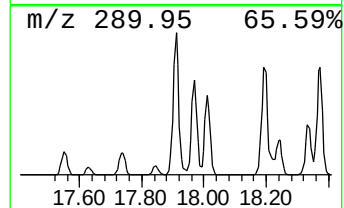
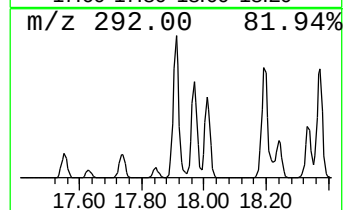
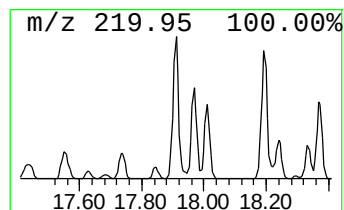
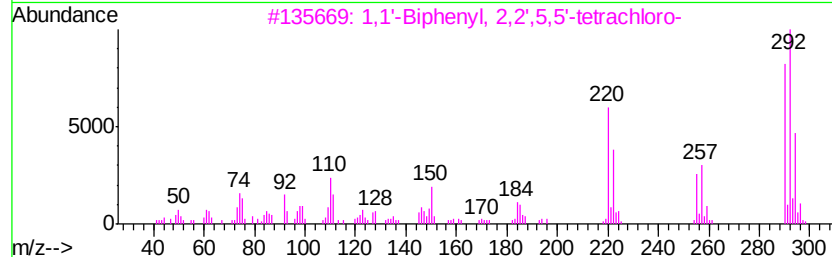
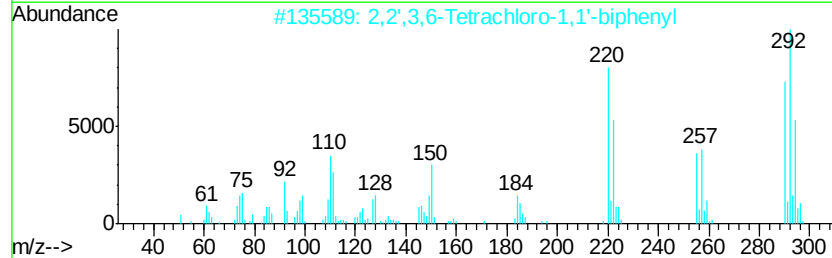
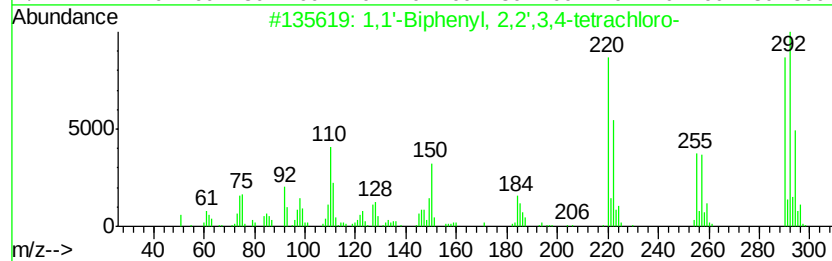
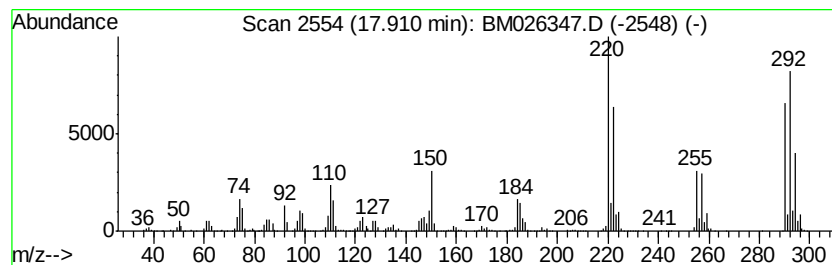
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 17 2,2',3,6-Tetrachloro-1,1'-b... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.91	56.73 ng/ul	5539150	Phenanthrene-d10	16.82

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		2,2',3,6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-45-7	99
3		1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	99
4		1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
5		1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

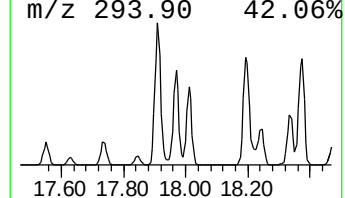
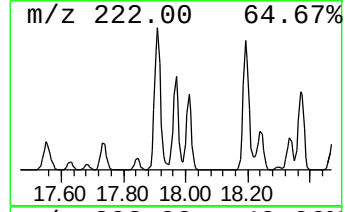
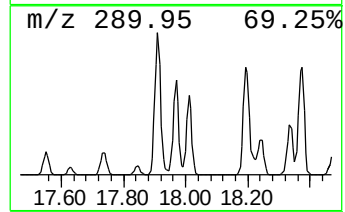
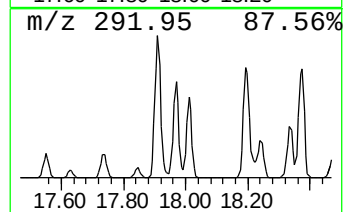
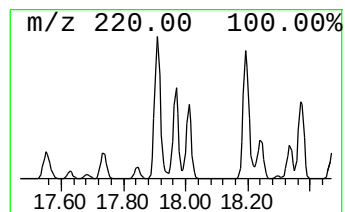
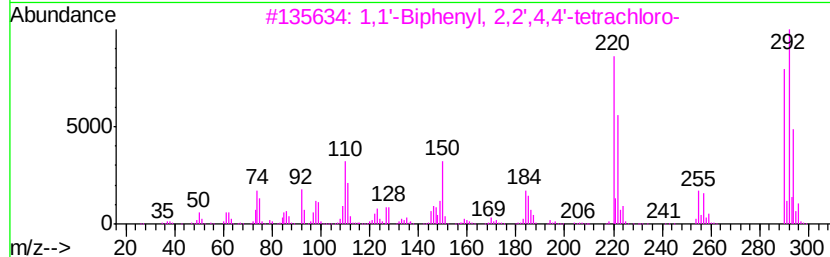
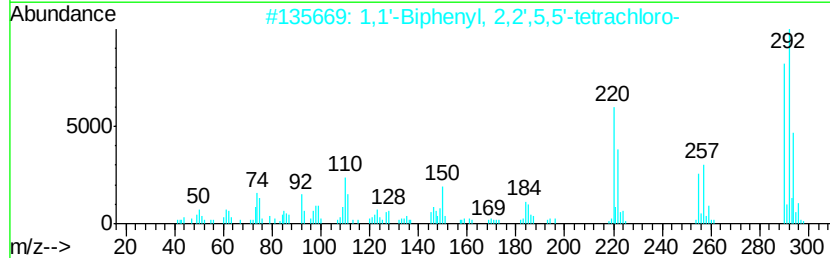
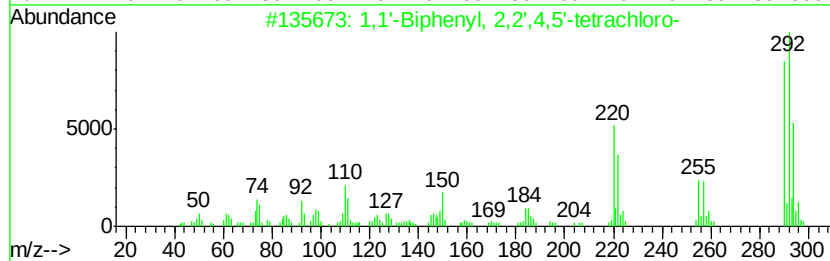
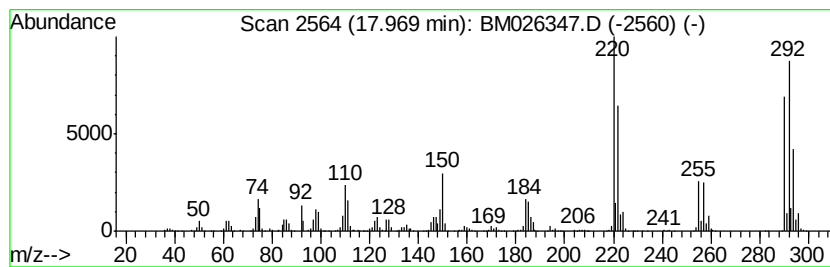
Instrument :
BNA_M
ClientSampleId :
ETTX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 18 1,1'-Biphenyl, 2,2',4,5'-te... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.		
17.97	34.37 ng/ul	3356560	Phenanthrene-d10	16.82		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'	-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	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	1,1'	-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99
5	2,2',4,5-Tetrachloro-1,1'-biphenyl		290	C12H6Cl4	070362-47-9	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

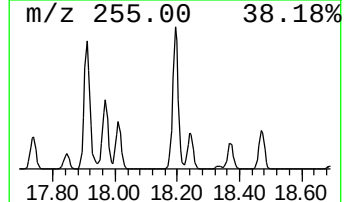
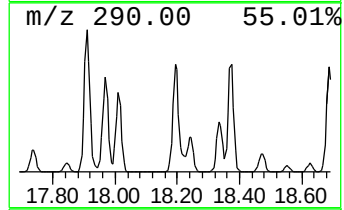
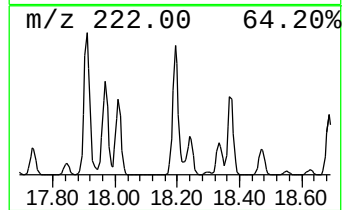
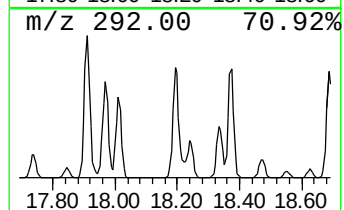
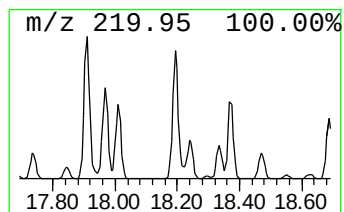
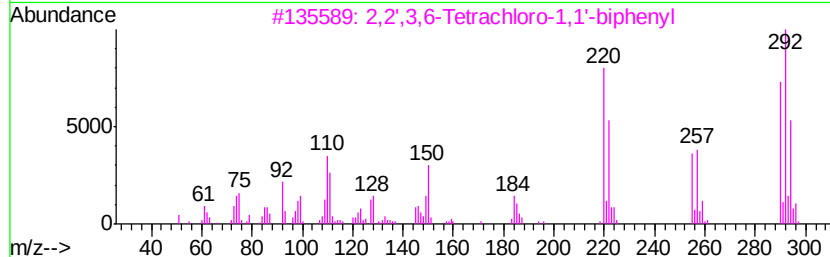
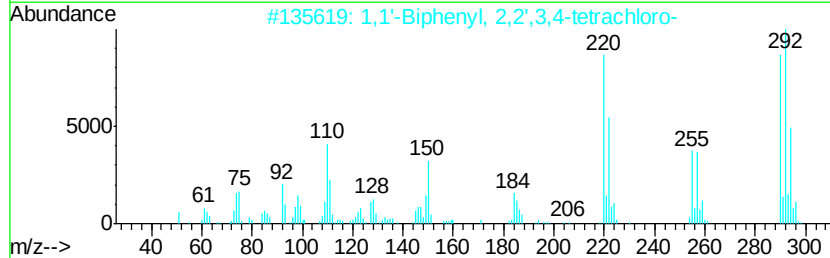
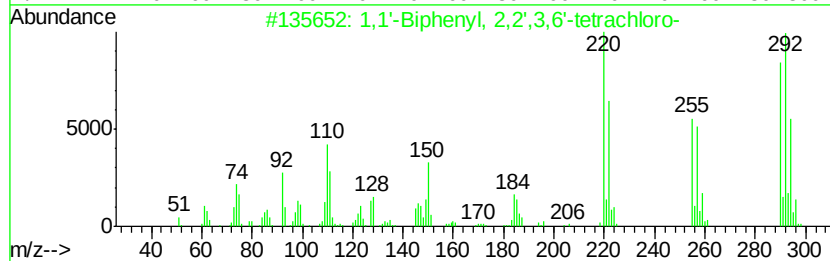
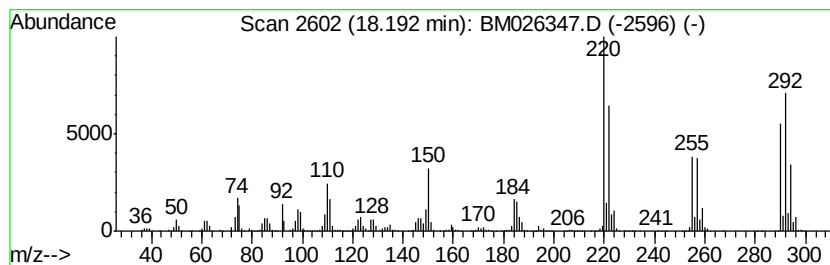
Instrument :
BNA_M
ClientSampleId :
ETTX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 20 1,1'-Biphenyl, 2,2',3,6'-te... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.19	52.47 ng/ul	5123600	Phenanthrene-d10	16.82		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,6'-tetrachl...	290	C12H6Cl4		041464-47-5	99
2	1,1'-Biphenyl, 2,2',3,4'-tetrachl...	290	C12H6Cl4		052663-59-9	99
3	2,2',3,6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4		070362-45-7	99
4	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4		002437-79-8	99
5	1,1'-Biphenyl, 2,2',4,6-Tetrachl...	290	C12H6Cl4		062796-65-0	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

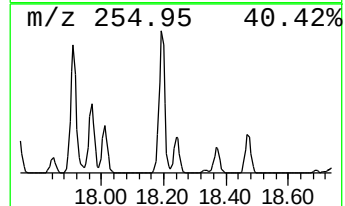
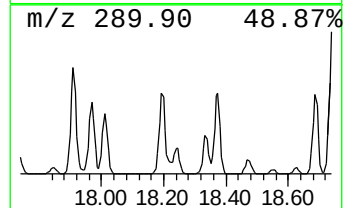
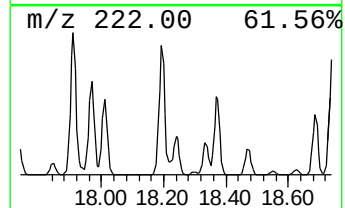
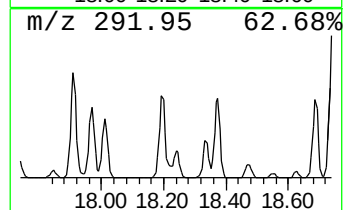
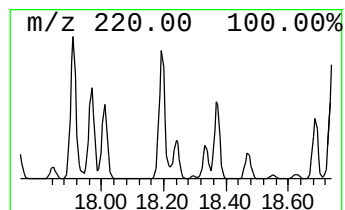
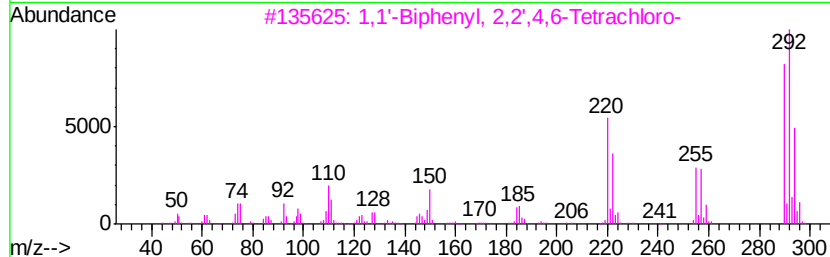
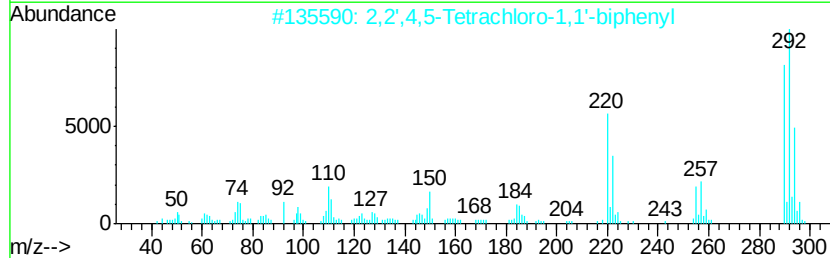
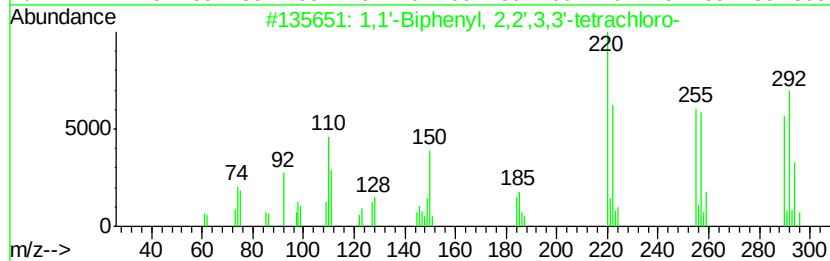
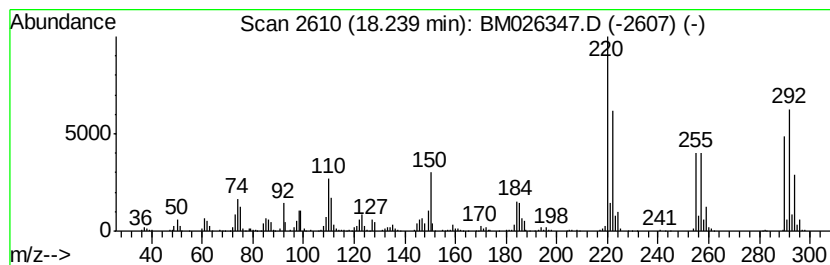
Instrument :
BNA_M
ClientSampleId :
ETTX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.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'-te... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.			
18.24	15.51 ng/ul	1514840	Phenanthrene-d10	16.82			
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'	-Biphenyl,	2,2',3,3'-tetrach...	290	C12H6Cl4	038444-93-8	99
2	2,2',	4,5-Tetrachloro-	1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99
3	1,1'	-Biphenyl,	2,2',4,6-Tetrachl...	290	C12H6Cl4	062796-65-0	99
4	1,1'	-Biphenyl,	2,2',3,5'-tetrach...	290	C12H6Cl4	041464-39-5	99
5	2,2',	4,5-Tetrachloro-	1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETTX9

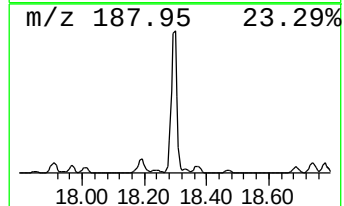
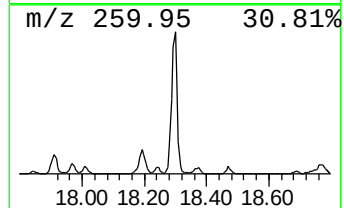
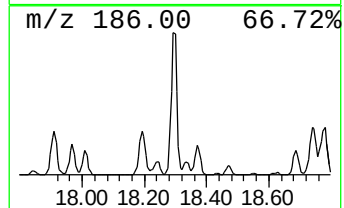
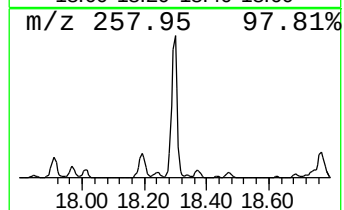
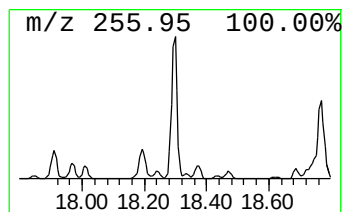
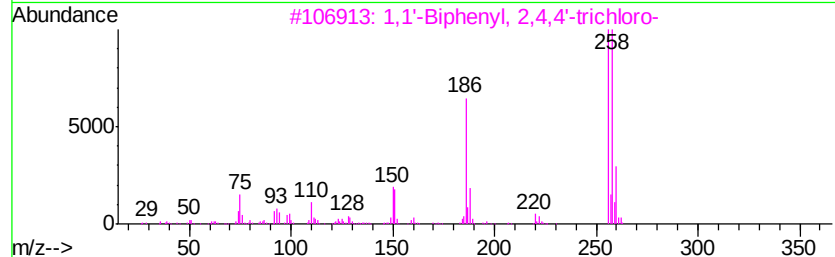
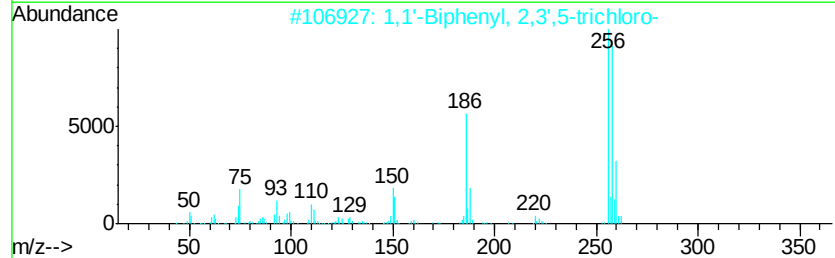
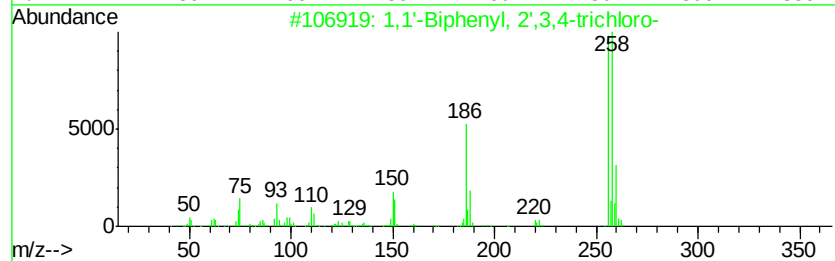
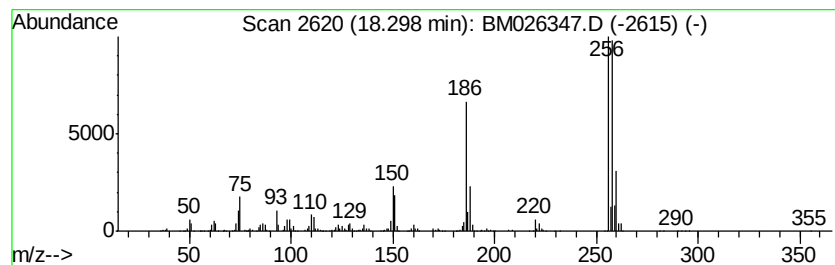
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 22 1,1'-Biphenyl, 2',3,4-trich... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.30	13.47 ng/ul	1315450	Phenanthrene-d10	16.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99
2		1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99
3		1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99
4		1,1'-Biphenyl, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	99
5		1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETTX9

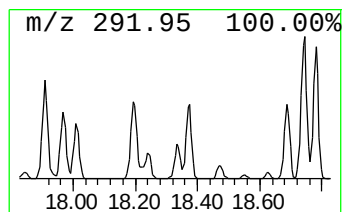
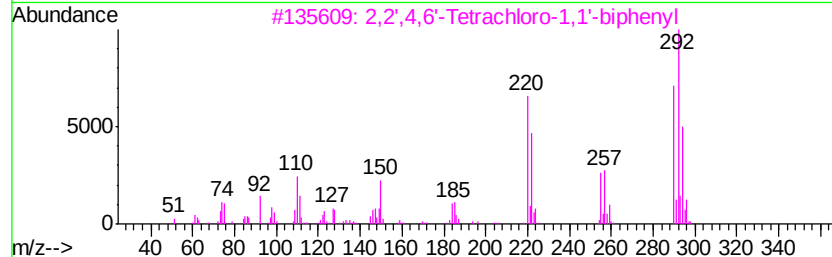
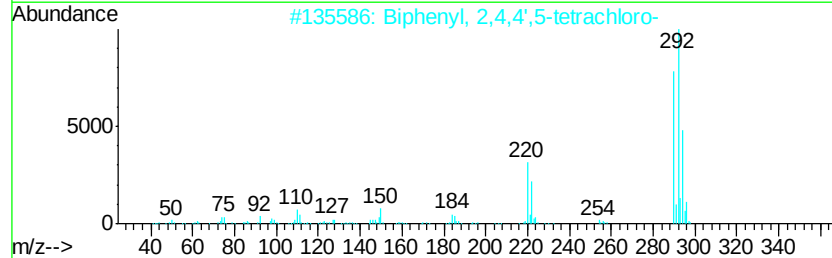
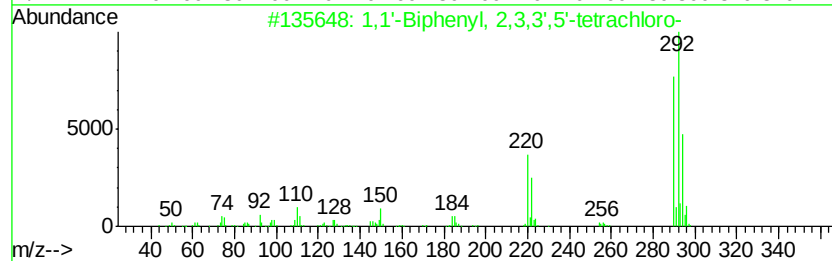
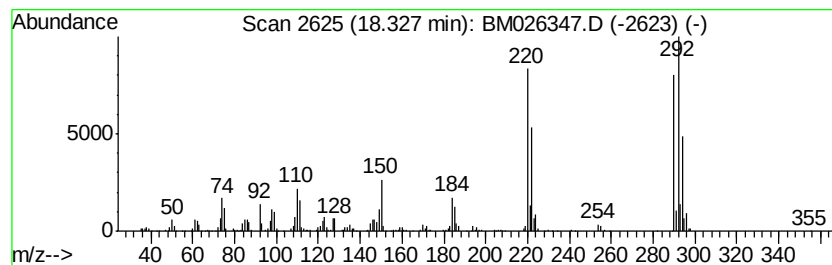
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 23 1,1'-Biphenyl, 2,3,3',5'-te... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.33	14.45 ng/ul	1411050	Phenanthrene-d10	16.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3,3',5'-tetrach...	290	C12H6Cl4	041464-49-7	99
2		Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	99
3		2,2',4,6'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	068194-04-7	99
4		1,1'-Biphenyl, 3,3',5,5'-tetrach...	290	C12H6Cl4	033284-52-5	99
5		1,1'-Biphenyl, 2,2',5,6'-Tetrach...	290	C12H6Cl4	041464-41-9	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETTX9

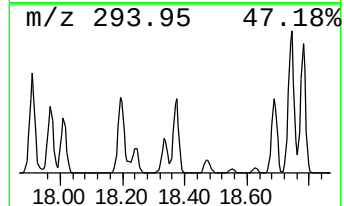
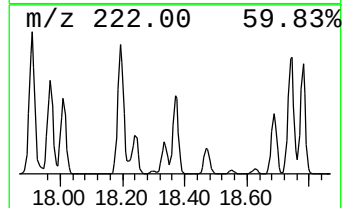
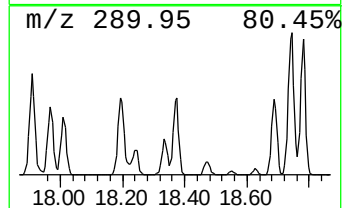
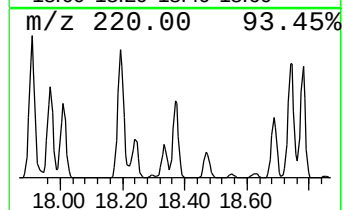
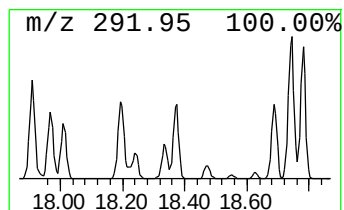
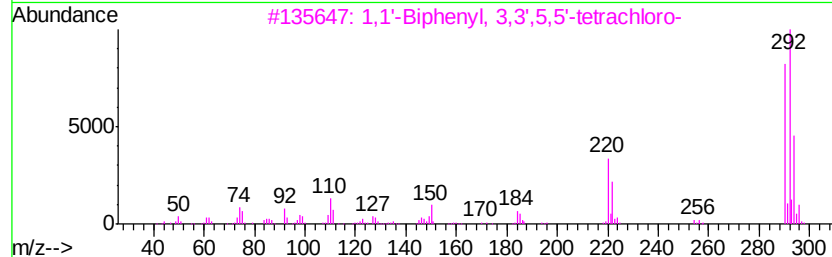
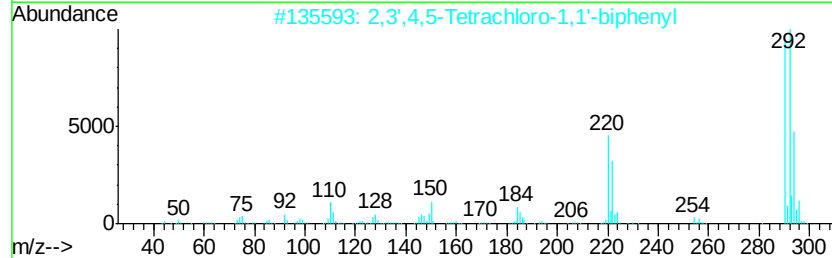
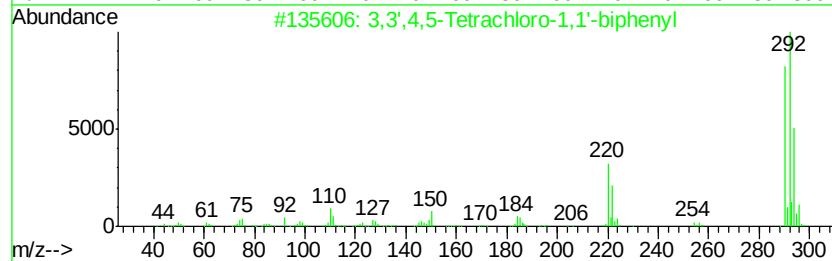
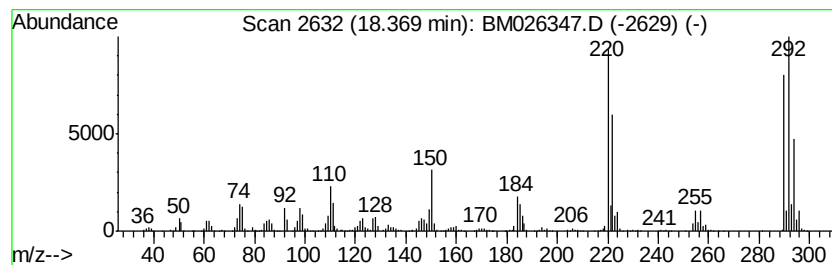
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 24 3,3',4,5-Tetrachloro-1,1'-b... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.37	32.35 ng/ul	3158610	Phenanthrene-d10	16.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-49-1	99
2		2,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-53-8	99
3		1,1'-Biphenyl, 3,3',5,5'-tetrach...	290	C12H6Cl4	033284-52-5	99
4		2,3,3',6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074472-33-6	99
5		2,3',4,5'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-52-7	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETTX9

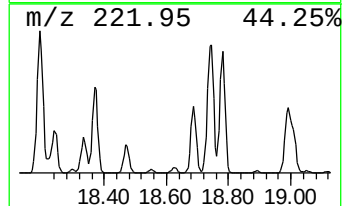
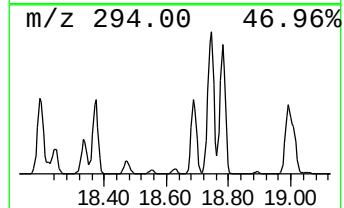
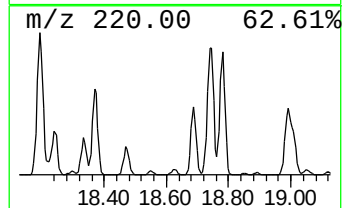
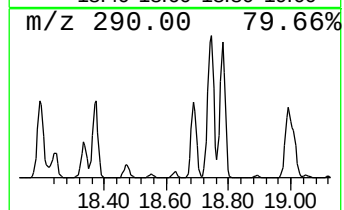
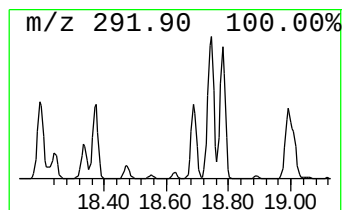
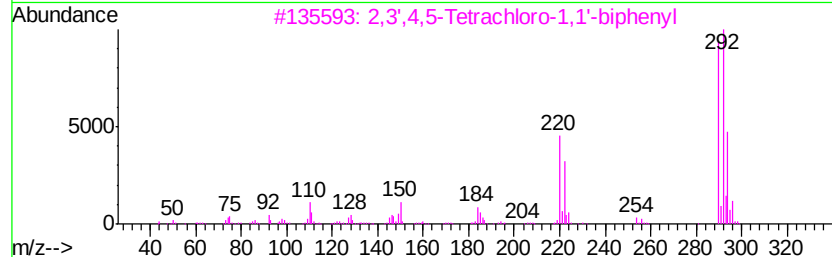
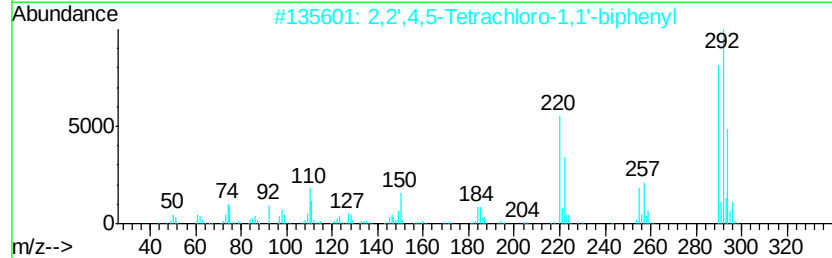
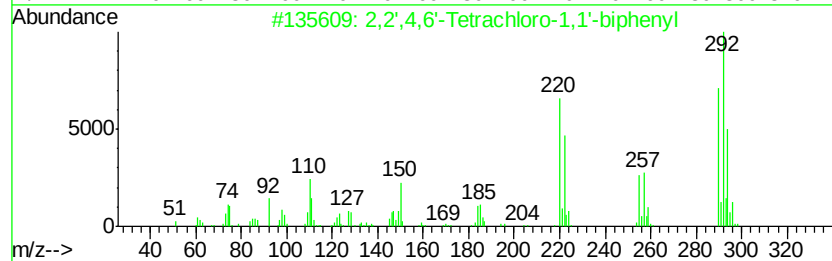
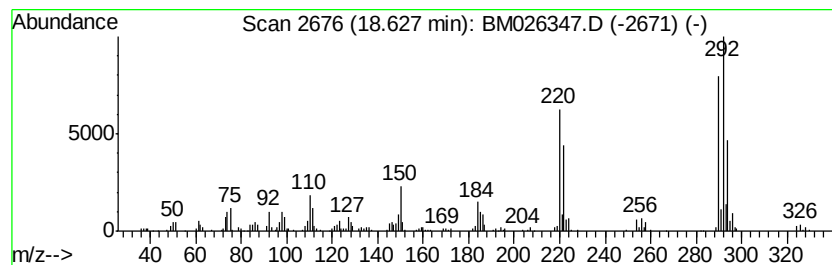
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 26 2,2',4,6'-Tetrachloro-1,1'-... Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.63	2.33 ng/ul	227201	Phenanthrene-d10	16.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2',4,6'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	068194-04-7	99
2		2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99
3		2,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-53-8	99
4		1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	99
5		1,1'-Biphenyl, 2,2',4,6-Tetrachl...	290	C12H6Cl4	062796-65-0	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

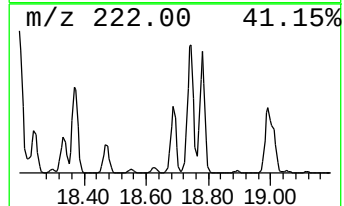
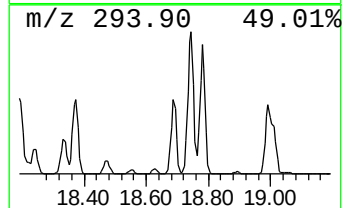
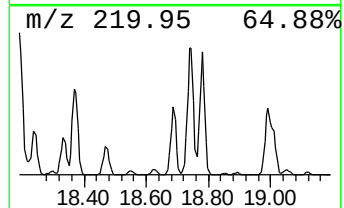
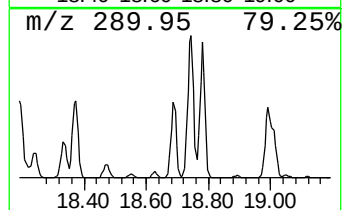
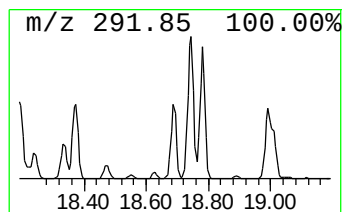
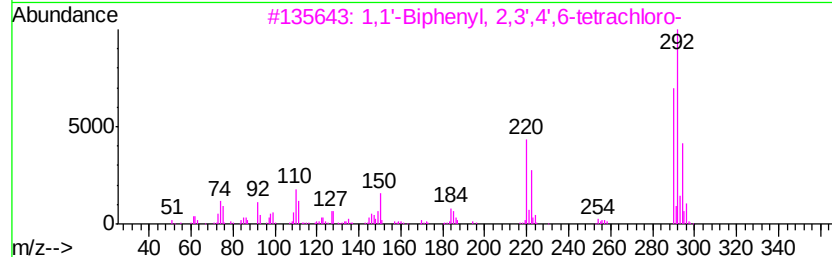
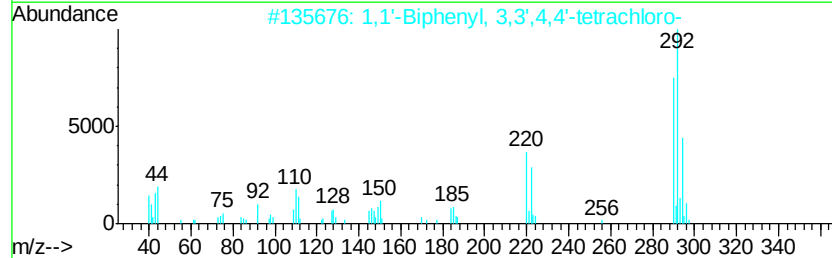
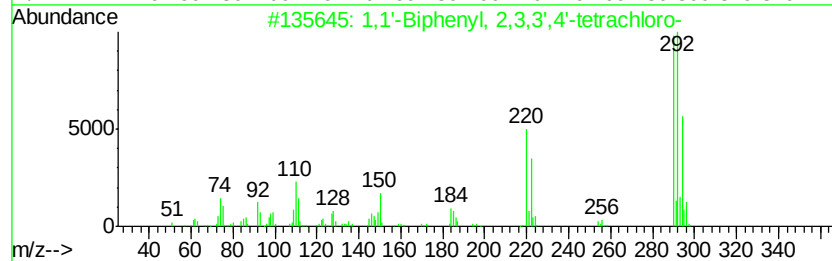
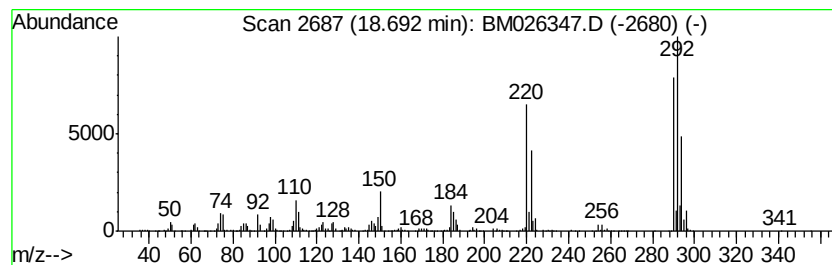
Instrument :
BNA_M
ClientSampleId :
ETTX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 27 1,1'-Biphenyl, 2,3,3',4'-te... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.69	26.89 ng/ul	2625560	Phenanthrene-d10	16.82		
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	1,1'-Biphenyl, 3,3',4,4'-tetrach...	290	C12H6Cl4	032598-13-3	99	
3	1,1'-Biphenyl, 2,3',4',6-tetrach...	290	C12H6Cl4	041464-46-4	99	
4	2,3',5',6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074338-23-1	99	
5	2,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-53-8	99	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

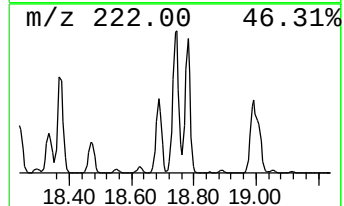
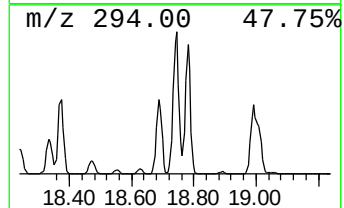
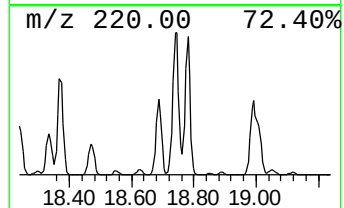
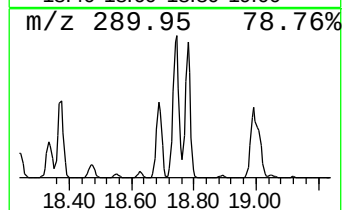
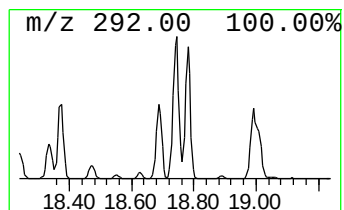
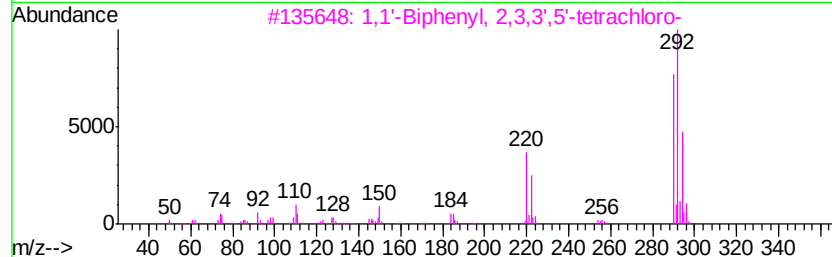
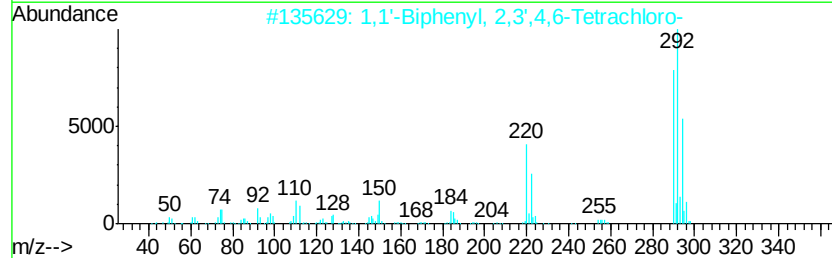
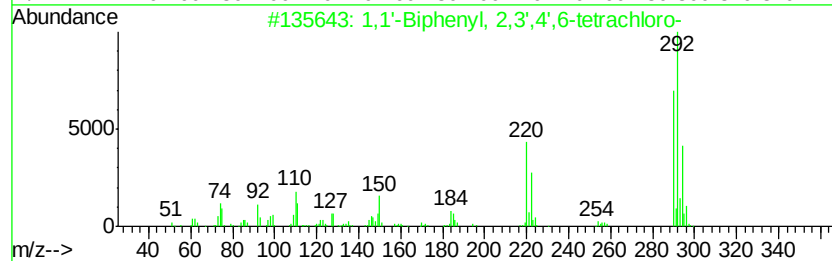
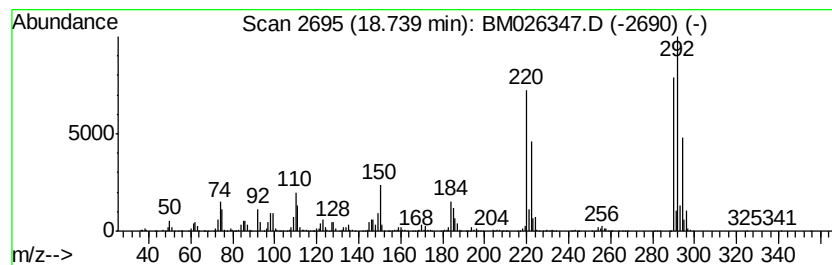
Instrument :
BNA_M
ClientSampleId :
ETTX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 28 1,1'-Biphenyl, 2,3',4',6-te... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.74	59.65 ng/ul	5824720	Phenanthrene-d10	16.82	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3',4',6-tetrach...	290	C12H6Cl4	041464-46-4	99
2	1,1'-Biphenyl, 2,3',4,6-Tetrachl...	290	C12H6Cl4	060233-24-1	99
3	1,1'-Biphenyl, 2,3,3',5'-tetrach...	290	C12H6Cl4	041464-49-7	99
4	1,1'-Biphenyl, 3,3',4,5'-tetrach...	290	C12H6Cl4	041464-48-6	99
5	1,1'-Biphenyl, 3,3',4,4'-tetrach...	290	C12H6Cl4	032598-13-3	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

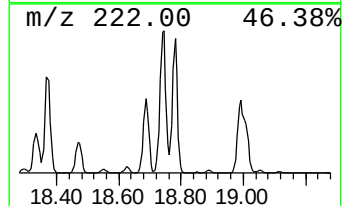
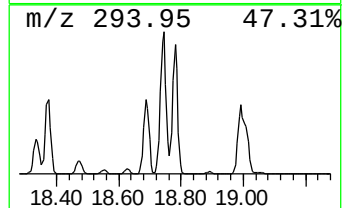
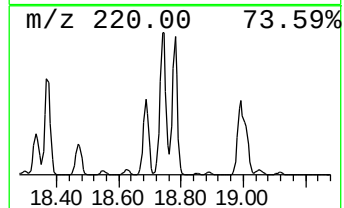
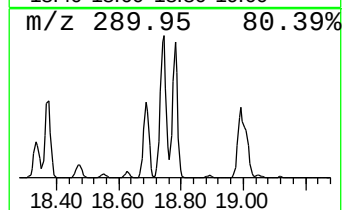
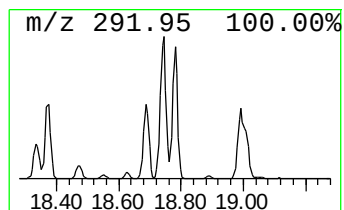
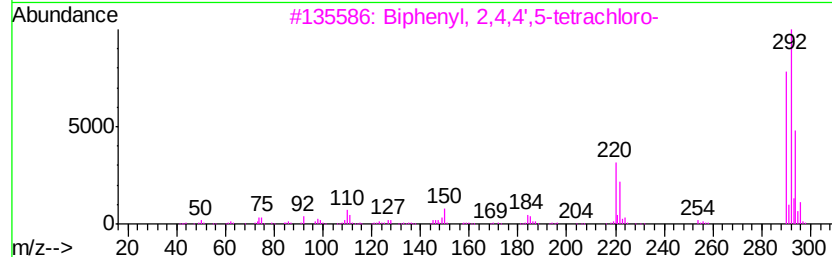
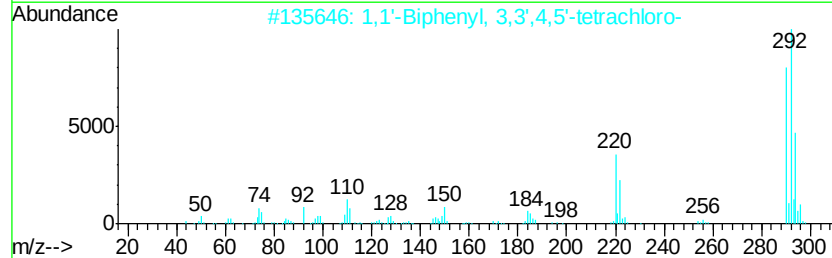
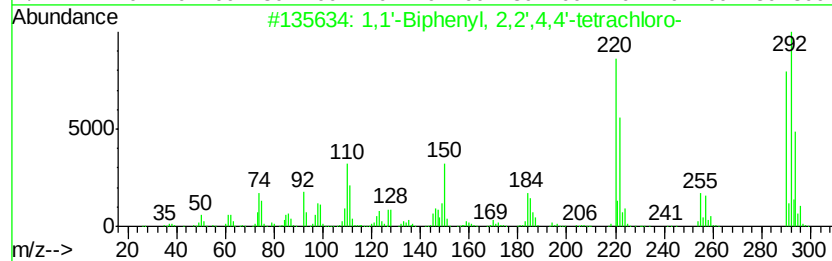
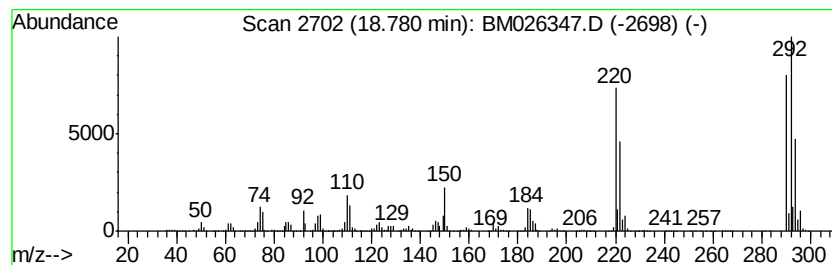
Instrument :
BNA_M
ClientSampleId :
ETTX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 29 1,1'-Biphenyl, 2,2',4,4'-te... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.78	56.39 ng/ul	5506270	Phenanthrene-d10	16.82	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,4'-tetrachloro-	290	C12H6Cl4	002437-79-8	99
2	1,1'-Biphenyl, 3,3',4,5'-tetrachloro-	290	C12H6Cl4	041464-48-6	99
3	Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	99
4	2,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-53-8	99
5	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026347.D
Acq On : 11 Jun 2020 01:35
Operator : JU/CG
Sample : L2850-06
Misc :
ALS Vial : 28 Sample Multiplier: 1

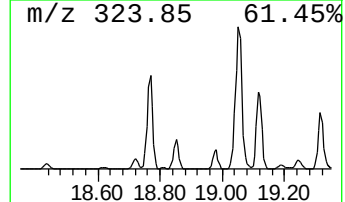
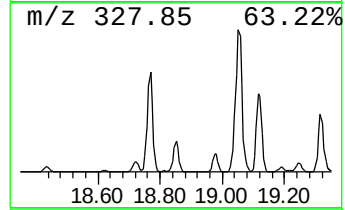
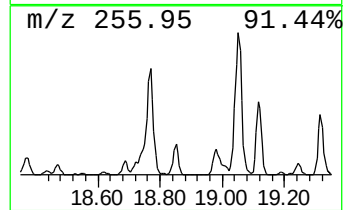
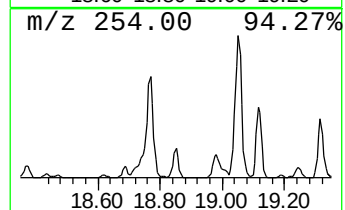
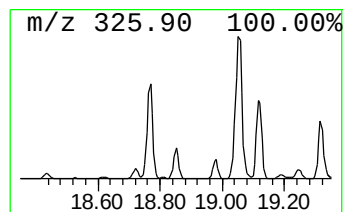
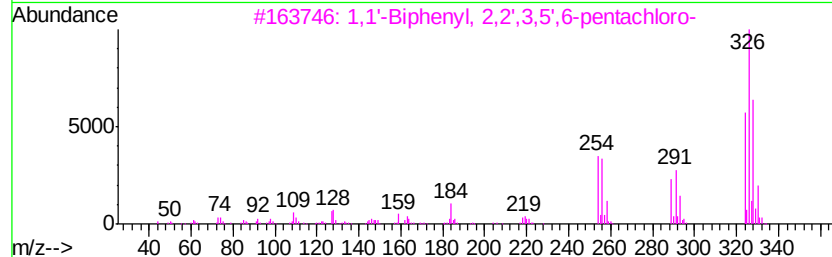
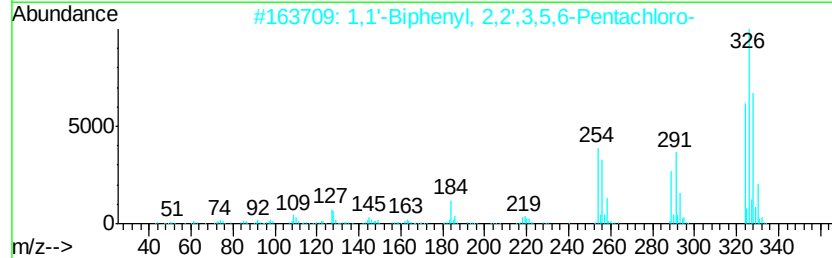
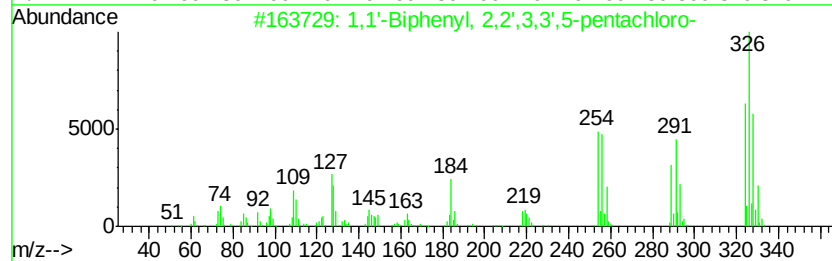
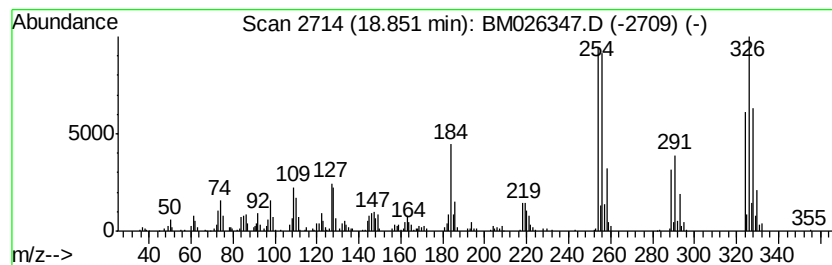
Instrument :
BNA_M
ClientSampleId :
ETTX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 30 1,1'-Biphenyl, 2,2',3,3',5-... Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.85	4.15 ng/ul	405017	Phenanthrene-d10	16.82	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,3',5-penta...	324	C12H5Cl5	060145-20-2	99
2	1,1'-Biphenyl, 2,2',3,5,6-Pentac...	324	C12H5Cl5	073575-56-1	99
3	1,1'-Biphenyl, 2,2',3,5',6-penta...	324	C12H5Cl5	038379-99-6	99
4	2,3,3',5,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	039635-32-0	99
5	3,3',4,4',5-Pentachloro-1,1'-bi...	324	C12H5Cl5	057465-28-8	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM061020\
 Data File : BM026347.D
 Acq On : 11 Jun 2020 01:35
 Operator : JU/CG
 Sample : L2850-06
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ETTX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.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...	2.92	11.7	ng/ul	536699	1	7.42	920686	20.0
1,1'-Biphenyl, 2,...	15.34	17.3	ng/ul	1116280	3	14.07	1291540	20.0
1,1'-Biphenyl, 2,...	15.95	4.4	ng/ul	425670	4	16.82	1952920	20.0
1,1'-Biphenyl, 2,...	16.07	26.9	ng/ul	2626370	4	16.82	1952920	20.0
1,1'-Biphenyl, 2,...	16.36	5.3	ng/ul	520513	4	16.82	1952920	20.0
1,1'-Biphenyl, 2,...	16.72	67.9	ng/ul	6627970	4	16.82	1952920	20.0
2,2',4-Trichloro-...	17.01	36.9	ng/ul	3598700	4	16.82	1952920	20.0
1,1'-Biphenyl, 3,...	17.29	9.4	ng/ul	916589	4	16.82	1952920	20.0
1,1'-Biphenyl, 2,...	17.32	2.5	ng/ul	244359	4	16.82	1952920	20.0
1,1'-Biphenyl, 2,...	17.43	115.7	ng/ul	11300300	4	16.82	1952920	20.0
1,1'-Biphenyl, 3,...	17.56	49.7	ng/ul	4857360	4	16.82	1952920	20.0
1,1'-Biphenyl, 2,...	17.68	21.4	ng/ul	2089450	4	16.82	1952920	20.0
1,1'-Biphenyl, 2,...	17.73	10.9	ng/ul	1068090	4	16.82	1952920	20.0
1,1'-Biphenyl, 2,...	17.85	4.6	ng/ul	444826	4	16.82	1952920	20.0
2,2',3,6-Tetrachl...	17.91	56.7	ng/ul	5539150	4	16.82	1952920	20.0
1,1'-Biphenyl, 2,...	17.97	34.4	ng/ul	3356560	4	16.82	1952920	20.0
1,1'-Biphenyl, 2,...	18.19	52.5	ng/ul	5123600	4	16.82	1952920	20.0
1,1'-Biphenyl, 2,...	18.24	15.5	ng/ul	1514840	4	16.82	1952920	20.0
1,1'-Biphenyl, 2'...	18.30	13.5	ng/ul	1315450	4	16.82	1952920	20.0
1,1'-Biphenyl, 2,...	18.33	14.4	ng/ul	1411050	4	16.82	1952920	20.0
3,3',4,5-Tetrachl...	18.37	32.4	ng/ul	3158610	4	16.82	1952920	20.0
2,2',4,6'-Tetrach...	18.63	2.3	ng/ul	227201	4	16.82	1952920	20.0
1,1'-Biphenyl, 2,...	18.69	26.9	ng/ul	2625560	4	16.82	1952920	20.0
1,1'-Biphenyl, 2,...	18.74	59.6	ng/ul	5824720	4	16.82	1952920	20.0
1,1'-Biphenyl, 2,...	18.78	56.4	ng/ul	5506270	4	16.82	1952920	20.0
1,1'-Biphenyl, 2,...	18.85	4.2	ng/ul	405017	4	16.82	1952920	20.0