

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061719\
 Data File : BM021000.D
 Acq On : 18 Jun 2019 04:26
 Operator : HP/JU
 Sample : AR1248 50PPM
 Misc : GCMS CONFIRMATION
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 AR1248 50PPM

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-BM061419MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.046	8	10	18	rVB2	89842	106255	2.50%	0.380%
2	3.210	34	38	42	rVB	259748	253893	5.98%	0.908%
3	7.787	807	816	826	rBV	977894	1542269	36.34%	5.515%
4	7.998	849	852	859	rBV6	4389	7985	0.19%	0.029%
5	8.439	924	927	931	rBV3	4133	6566	0.15%	0.023%
6	10.580	1283	1291	1303	rBV	1209398	2103988	49.57%	7.524%
7	14.427	1938	1945	1956	rBV2	2040933	3086337	72.72%	11.037%
8	14.557	1961	1967	1972	rBV3	6258	8665	0.20%	0.031%
9	15.427	2109	2115	2118	rBV5	6629	10772	0.25%	0.039%
10	15.686	2154	2159	2163	rBV	26664	33436	0.79%	0.120%
11	16.127	2229	2234	2241	rBV5	7372	11323	0.27%	0.040%
12	16.292	2257	2262	2267	rBV3	11019	13164	0.31%	0.047%
13	16.410	2277	2282	2288	rVB	68699	89850	2.12%	0.321%
14	16.704	2328	2332	2336	rVB4	14016	17708	0.42%	0.063%
15	16.927	2364	2370	2373	rBV3	6520	10027	0.24%	0.036%
16	16.962	2373	2376	2381	rVB	9292	12898	0.30%	0.046%
17	17.051	2386	2391	2395	rBV	289170	379252	8.94%	1.356%
18	17.086	2395	2397	2406	rVV	75927	84762	2.00%	0.303%
19	17.174	2406	2412	2423	rVV2	3061847	4244374	100.00%	15.178%
20	17.268	2427	2428	2434	rVV4	7276	10539	0.25%	0.038%
21	17.345	2436	2441	2447	rVB2	119900	179240	4.22%	0.641%
22	17.627	2481	2489	2492	rBV3	26021	40200	0.95%	0.144%
23	17.656	2492	2494	2500	rVB6	8345	9924	0.23%	0.035%
24	17.762	2506	2512	2523	rVB	376941	729412	17.19%	2.608%
25	17.898	2529	2535	2541	rVB2	176668	293204	6.91%	1.048%
26	17.962	2543	2546	2552	rBV5	16378	22546	0.53%	0.081%
27	18.021	2552	2556	2561	rVV	84850	110311	2.60%	0.394%
28	18.074	2561	2565	2569	rVB	66527	83726	1.97%	0.299%
29	18.186	2580	2584	2589	rBV5	30609	37686	0.89%	0.135%
30	18.239	2589	2593	2599	rVV	421382	575769	13.57%	2.059%
31	18.298	2599	2603	2607	rVV	247502	332375	7.83%	1.189%
32	18.345	2607	2611	2618	rVB2	188996	258756	6.10%	0.925%
33	18.527	2637	2642	2647	rBV	351827	497699	11.73%	1.780%
34	18.574	2647	2650	2654	rVV	106564	128155	3.02%	0.458%

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35	18.633	2656	2660	2662	rVV	37993	50415	1.19%	0.180%
36	18.662	2662	2665	2668	rVV	97495	127887	3.01%	0.457%
37	18.703	2668	2672	2677	rVB	236532	302701	7.13%	1.082%
38	18.803	2685	2689	2695	rBV3	69348	87953	2.07%	0.315%
39	18.951	2711	2714	2718	rVB3	12868	13602	0.32%	0.049%
40	19.009	2719	2724	2728	rVV	193926	238945	5.63%	0.854%
41	19.062	2728	2733	2736	rVV	435261	603381	14.22%	2.158%
42	19.098	2736	2739	2745	rVV2	389716	522426	12.31%	1.868%
43	19.174	2748	2752	2756	rVV2	36112	40526	0.95%	0.145%
44	19.315	2769	2776	2782	rBV	175358	380344	8.96%	1.360%
45	19.368	2782	2785	2792	rVV	195124	259465	6.11%	0.928%
46	19.433	2792	2796	2801	rVB	84381	102685	2.42%	0.367%
47	19.492	2802	2806	2811	rBV2	49567	59798	1.41%	0.214%
48	19.568	2815	2819	2822	rBV	49020	55821	1.32%	0.200%
49	19.609	2822	2826	2827	rVV	38206	44715	1.05%	0.160%
50	19.633	2827	2830	2835	rVB2	68598	83383	1.96%	0.298%
51	19.709	2839	2843	2847	rVV2	85685	103127	2.43%	0.369%
52	19.762	2848	2852	2856	rVV2	45063	55547	1.31%	0.199%
53	19.815	2857	2861	2866	rVV	157064	190896	4.50%	0.683%
54	19.862	2867	2869	2873	rVB5	16775	20701	0.49%	0.074%
55	19.962	2883	2886	2890	rVB4	43146	49275	1.16%	0.176%
56	20.133	2911	2915	2920	rVB	112219	140165	3.30%	0.501%
57	20.439	2963	2967	2968	rBV2	74936	79122	1.86%	0.283%
58	20.462	2968	2971	2976	rVB	824823	902871	21.27%	3.229%
59	20.539	2981	2984	2988	rVB5	39063	41212	0.97%	0.147%
60	20.597	2991	2994	2997	rBV	81917	88076	2.08%	0.315%
61	21.062	3070	3073	3077	rVB2	76846	82357	1.94%	0.295%
62	21.350	3117	3122	3131	rBV	2919247	3755080	88.47%	13.428%
63	21.497	3143	3147	3151	rBV	196105	222302	5.24%	0.795%
64	21.933	3218	3221	3227	rBV	174644	217247	5.12%	0.777%
65	22.403	3298	3301	3307	rVB	212264	267031	6.29%	0.955%
66	22.915	3385	3388	3396	rVB	172306	250163	5.89%	0.895%
67	23.627	3503	3509	3518	rVB	1737539	3191878	75.20%	11.414%

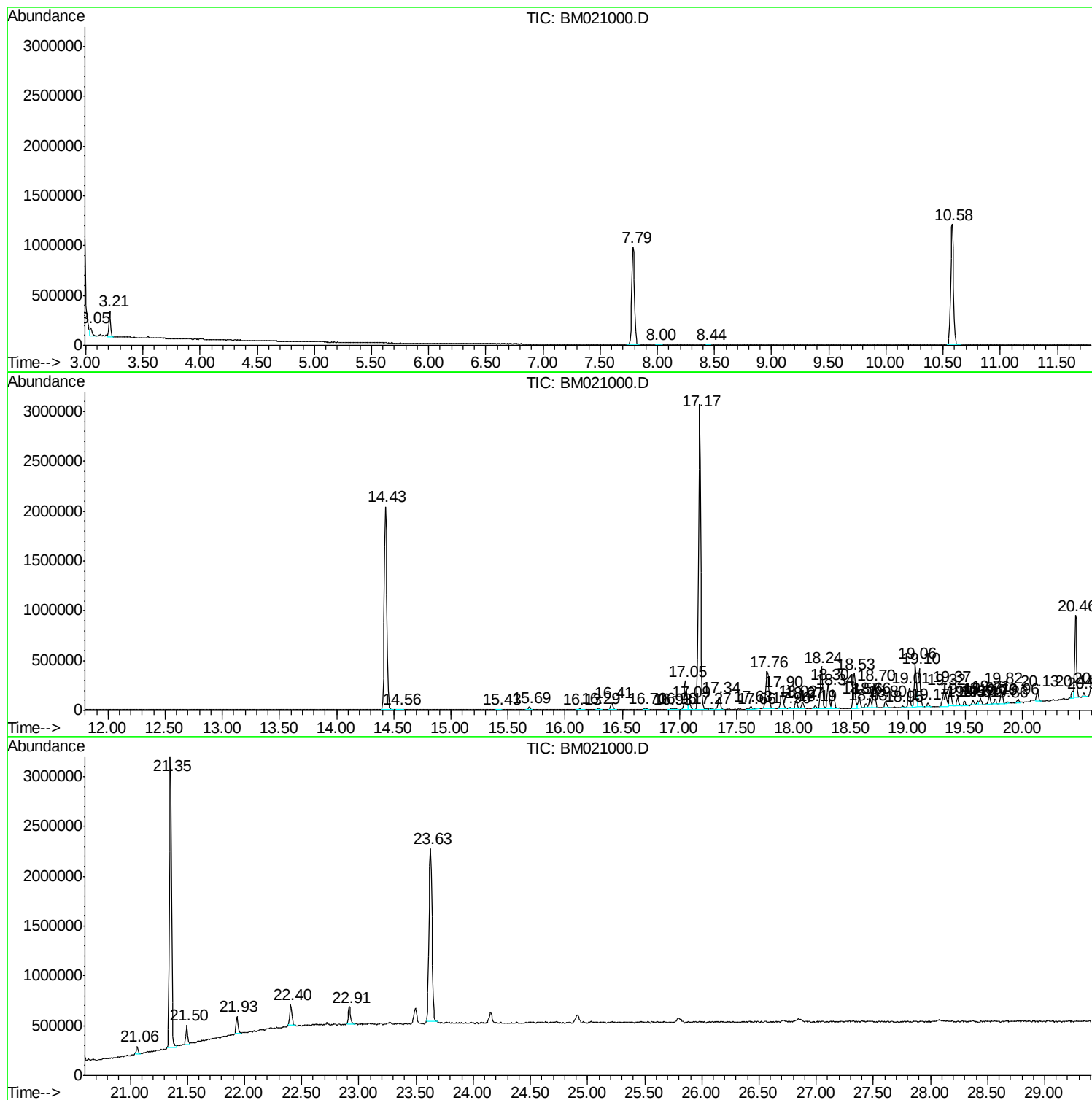
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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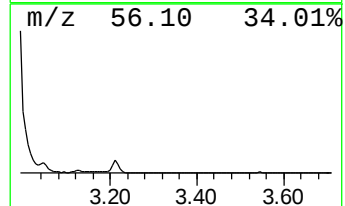
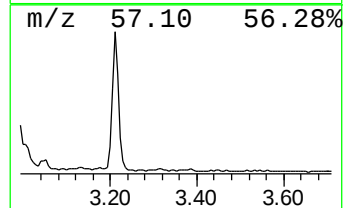
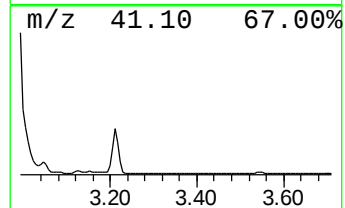
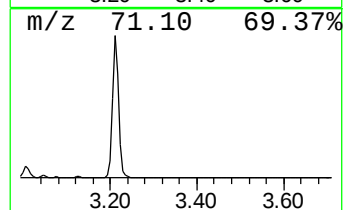
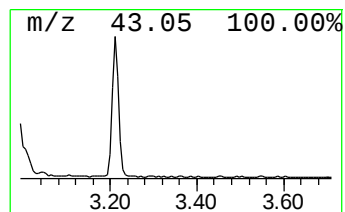
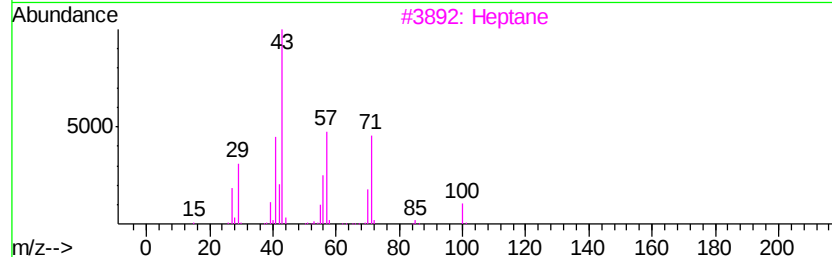
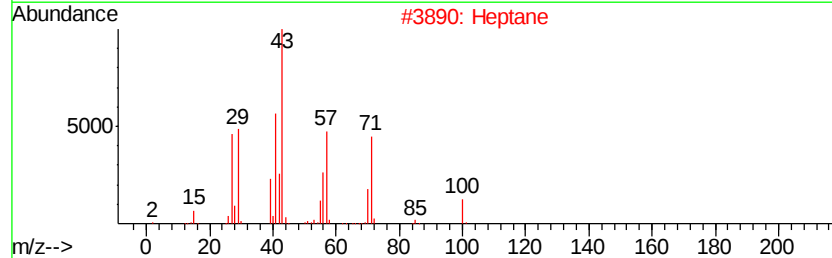
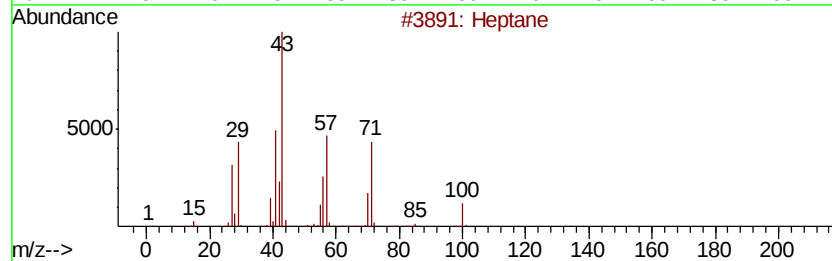
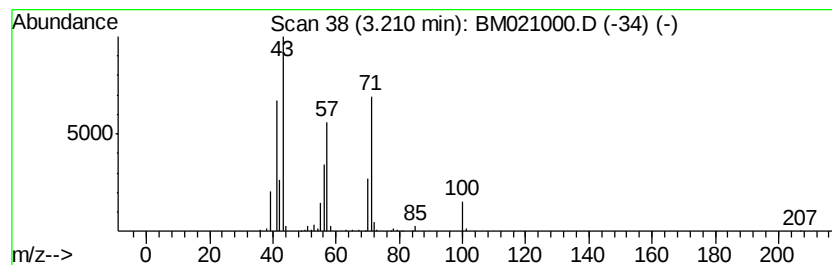
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.21	3.29 ng/ul	253893	1,4-Dichlorobenzene-d4	7.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane	100	C7H16	000142-82-5	91
2		Heptane	100	C7H16	000142-82-5	90
3		Heptane	100	C7H16	000142-82-5	87
4		Heptane	100	C7H16	000142-82-5	86
5		Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	47



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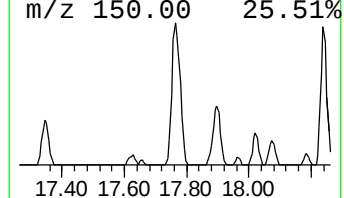
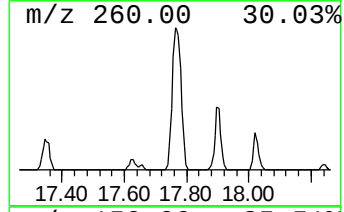
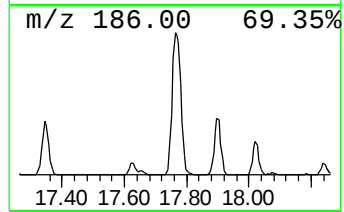
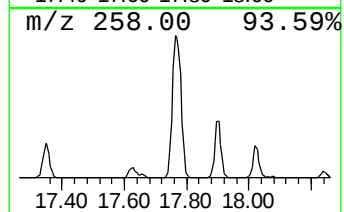
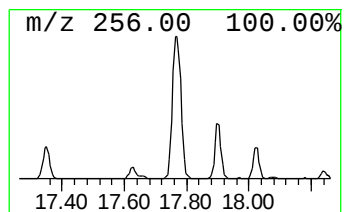
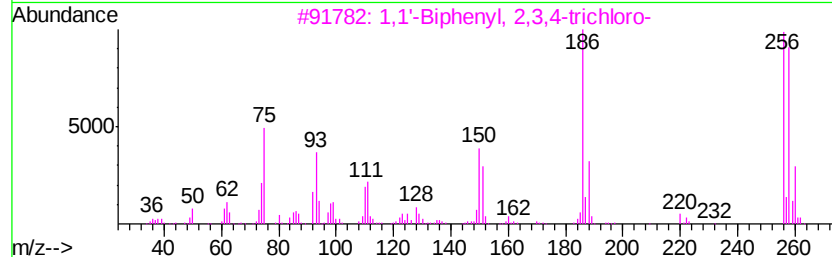
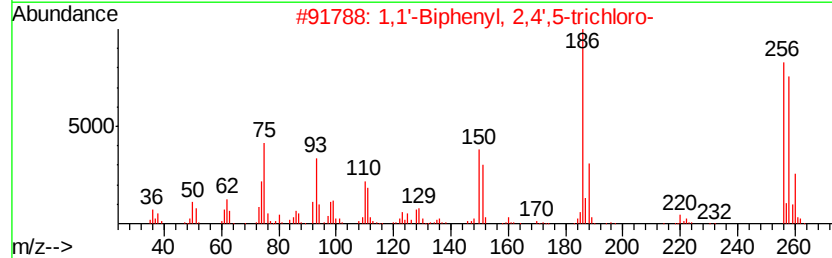
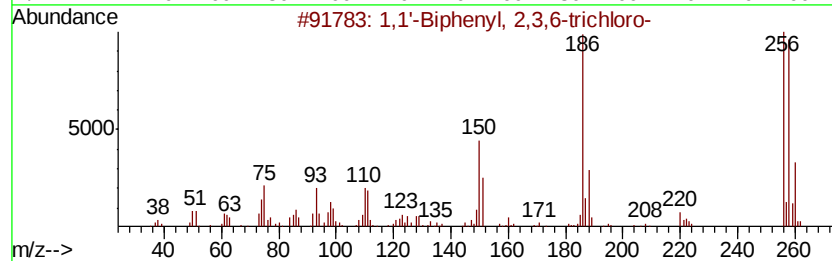
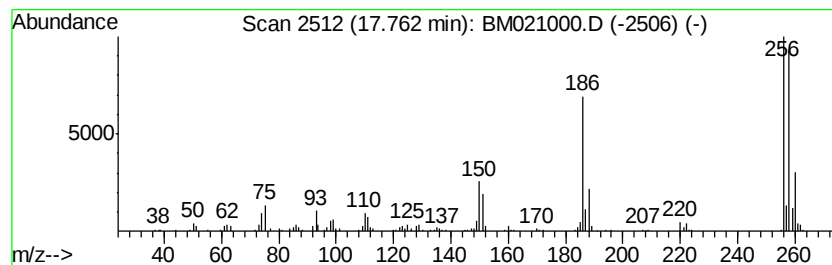
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.76	3.44 ng/ul	729412	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3,6-trichloro-	256	C12H7Cl3	055702-45-9	98
2		1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	98
3		1,1'-Biphenyl, 2,3,4-trichloro-	256	C12H7Cl3	055702-46-0	98
4		1,1'-Biphenyl, 2,3,4'-Trichloro-	256	C12H7Cl3	038444-85-8	98
5		1,1'-Biphenyl, 2,4,5-trichloro-	256	C12H7Cl3	015862-07-4	98



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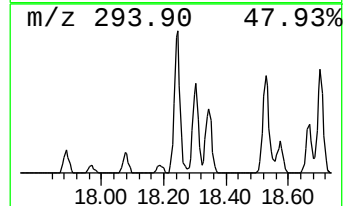
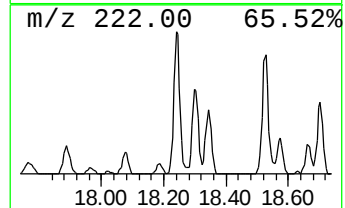
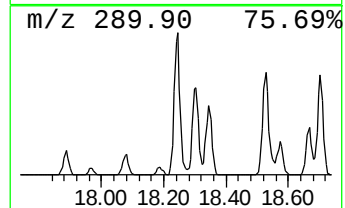
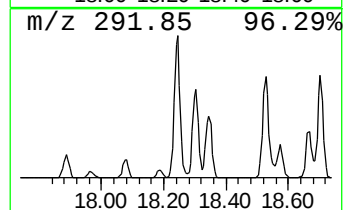
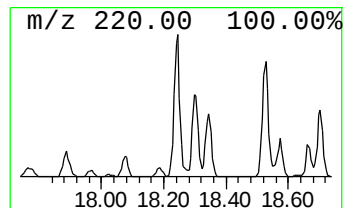
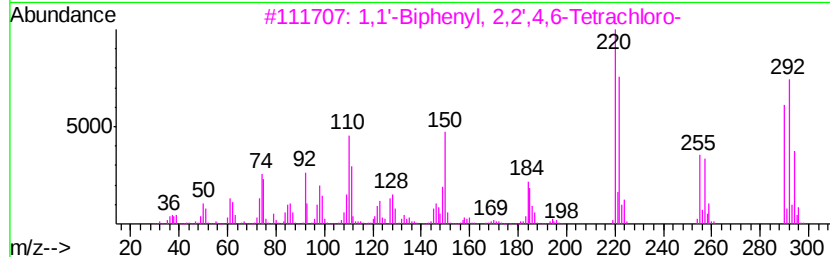
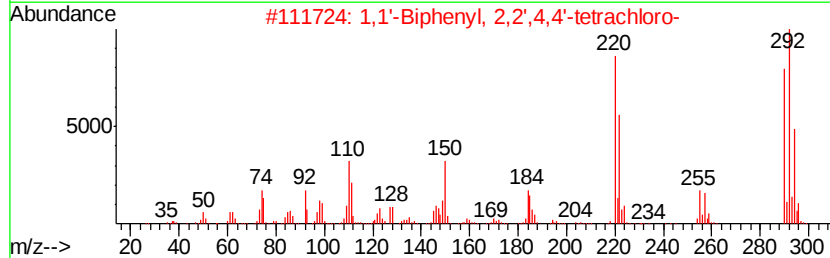
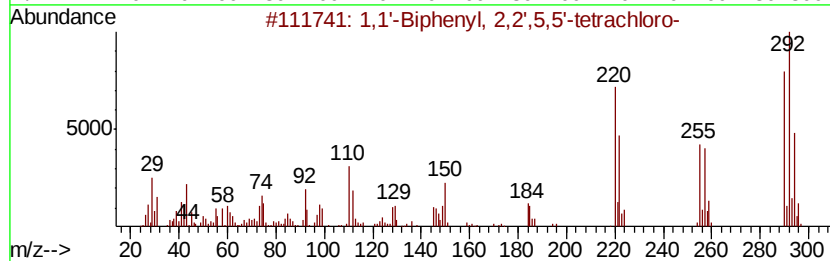
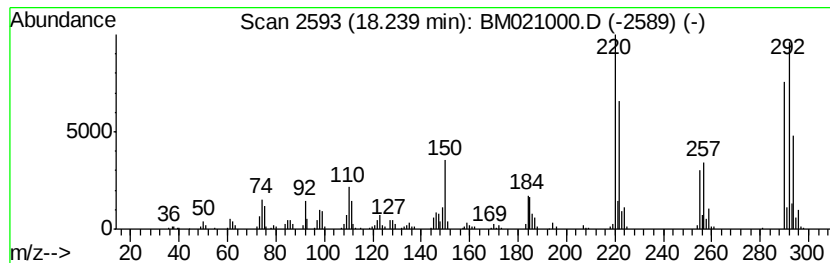
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Peak Number 3 1,1'-Biphenyl, 2,2',5,5'-te... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.24	2.71 ng/ul	575769	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	99
2		1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
3		1,1'-Biphenyl, 2,2',4,6-Tetrachl...	290	C12H6Cl4	062796-65-0	99
4		1,1'-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	99
5		1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99



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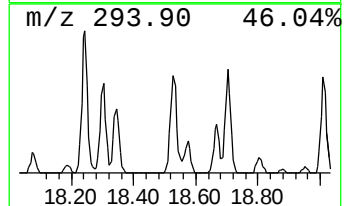
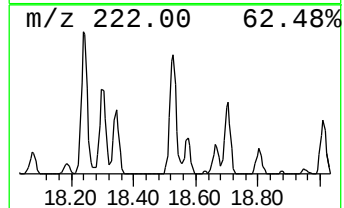
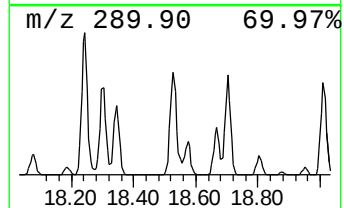
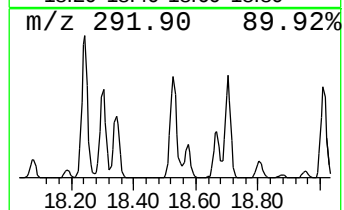
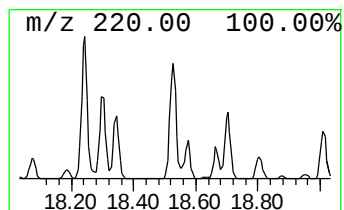
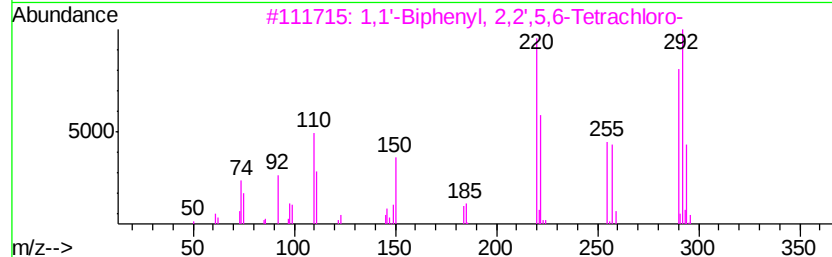
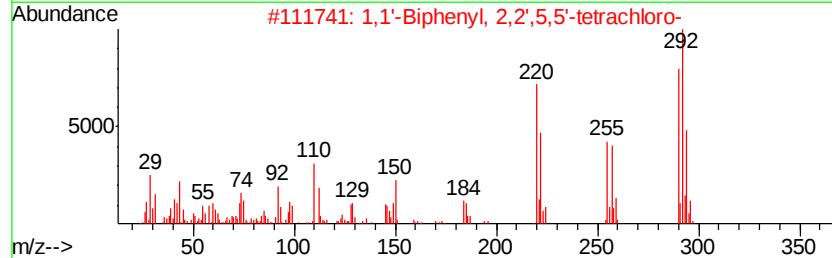
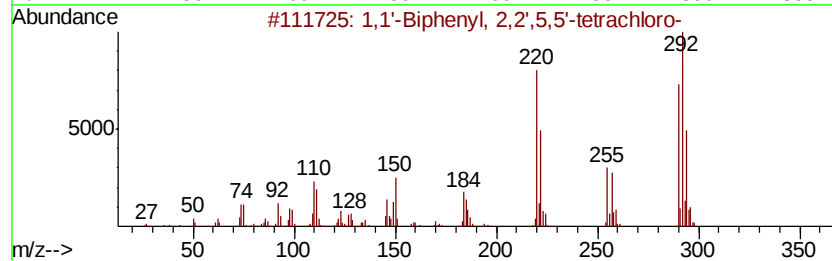
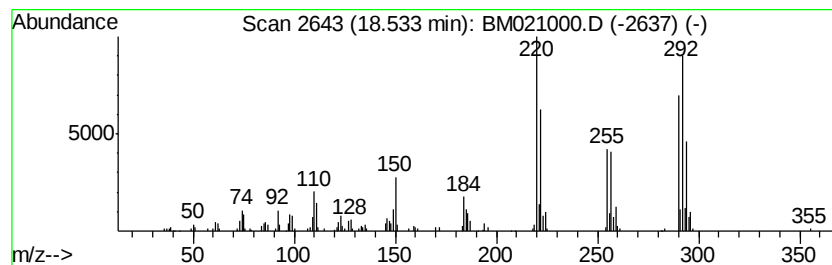
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Peak Number 4 1,1'-Biphenyl, 2,2',5,6-Tet... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.53	2.35 ng/ul	497699	Phenanthrene-d10	17.17		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5,5'-tetrachloro-	290	C12H6Cl4		035693-99-3	99
2	1,1'-Biphenyl, 2,2',5,5'-tetrachloro-	290	C12H6Cl4		035693-99-3	99
3	1,1'-Biphenyl, 2,2',5,6-Tetrachloro-	290	C12H6Cl4		041464-41-9	99
4	1,1'-Biphenyl, 2,2',4,4'-tetrachloro-	290	C12H6Cl4		002437-79-8	99
5	1,1'-Biphenyl, 2,2',5,6-Tetrachloro-	290	C12H6Cl4		041464-41-9	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061719\
Data File : BM021000.D
Acq On : 18 Jun 2019 04:26
Operator : HP/JU
Sample : AR1248 50PPM
Misc : GCMS CONFIRMATION
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
AR1248 50PPM

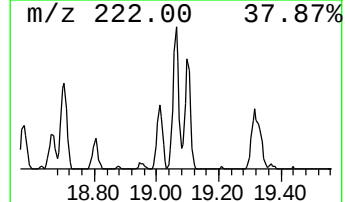
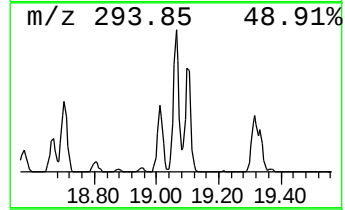
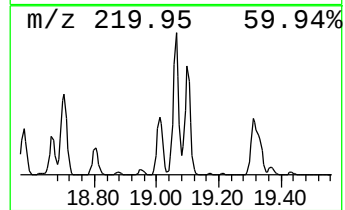
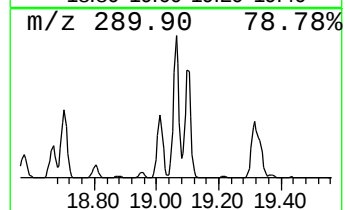
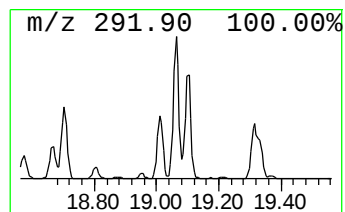
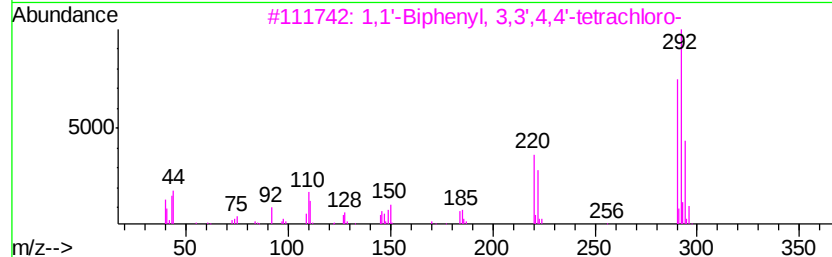
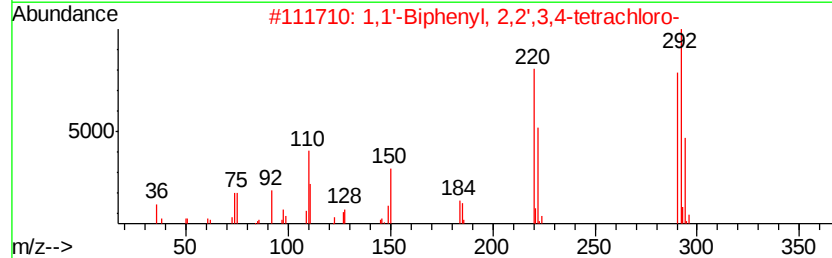
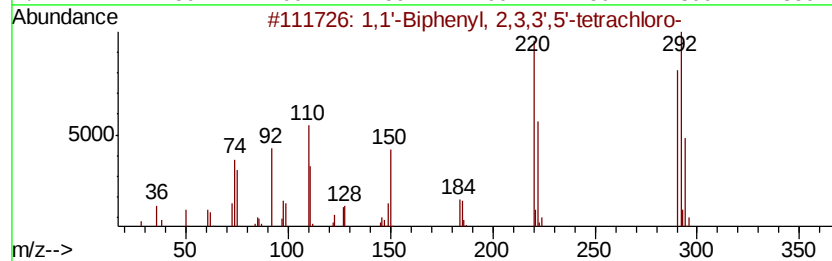
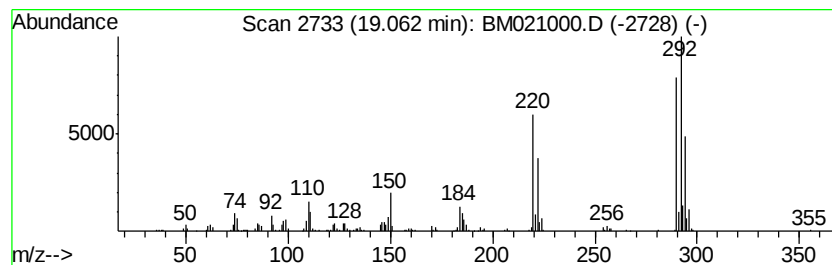
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.06	2.84 ng/ul	603381	Phenanthrene-d10	17.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'	-Biphenyl	2,3,3',5'-tetrach...	290	C12H6Cl4	041464-49-7	99
2	1,1'	-Biphenyl	2,2',3,4-tetrachl...	290	C12H6Cl4	052663-59-9	99
3	1,1'	-Biphenyl	3,3',4,4'-tetrach...	290	C12H6Cl4	032598-13-3	99
4	1,1'	-Biphenyl	2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99
5	1,1'	-Biphenyl	2,3',4',5-tetrach...	290	C12H6Cl4	032598-11-1	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061719\
Data File : BM021000.D
Acq On : 18 Jun 2019 04:26
Operator : HP/JU
Sample : AR1248 50PPM
Misc : GCMS CONFIRMATION
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
AR1248 50PPM

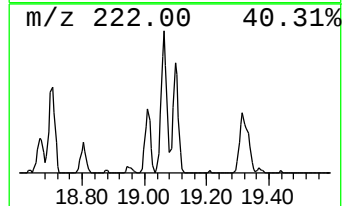
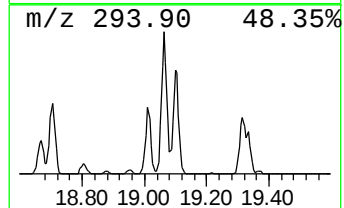
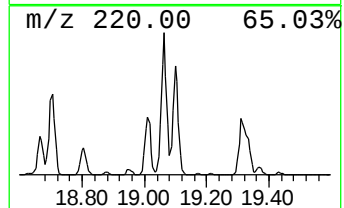
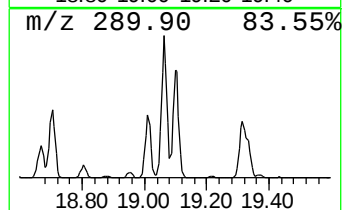
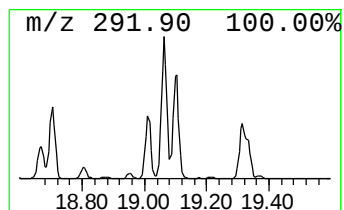
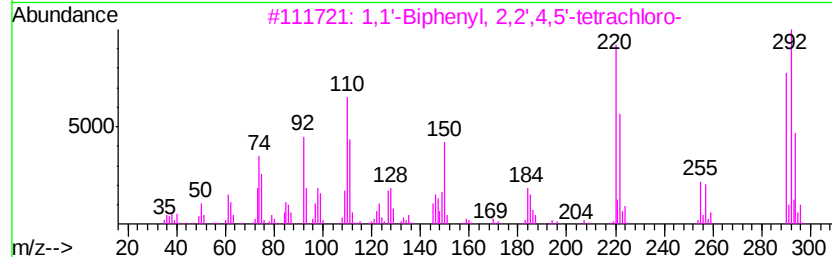
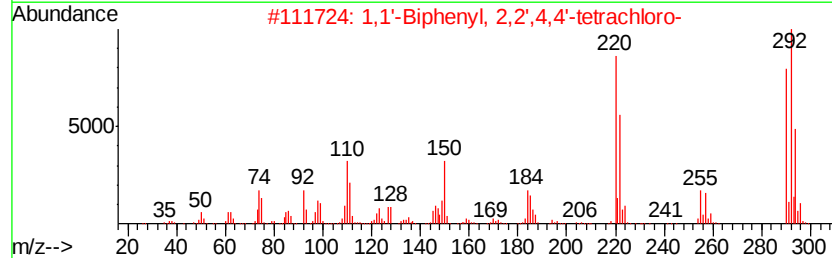
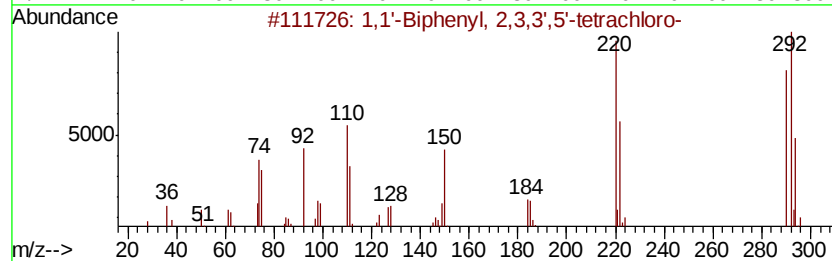
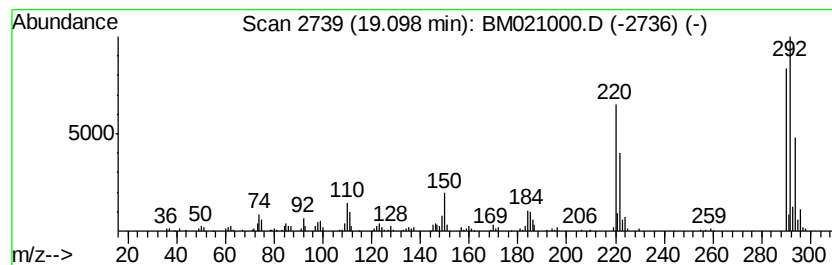
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.10	2.46 ng/ul	522426	Phenanthrene-d10	17.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1'-Biphenyl, 2,3,3',5'-tetrach...	290	C12H6Cl4	041464-49-7	99
2			1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
3			1,1'-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	99
4			1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99
5			1,1'-Biphenyl, 2,3',5,5'-tetrach...	290	C12H6Cl4	041464-42-0	98



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061719\
Data File : BM021000.D
Acq On : 18 Jun 2019 04:26
Operator : HP/JU
Sample : AR1248 50PPM
Misc : GCMS CONFIRMATION
ALS Vial : 30 Sample Multiplier: 1

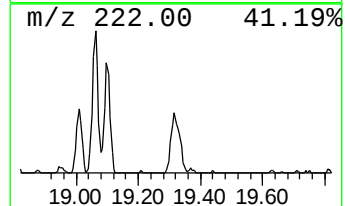
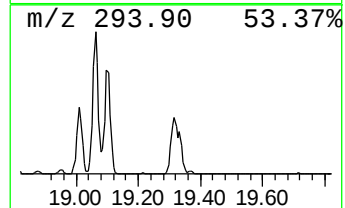
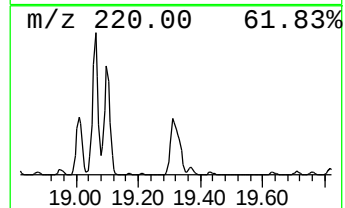
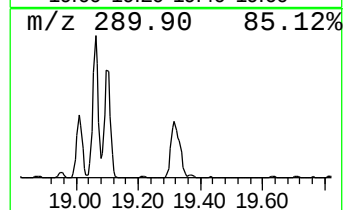
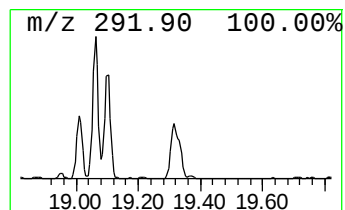
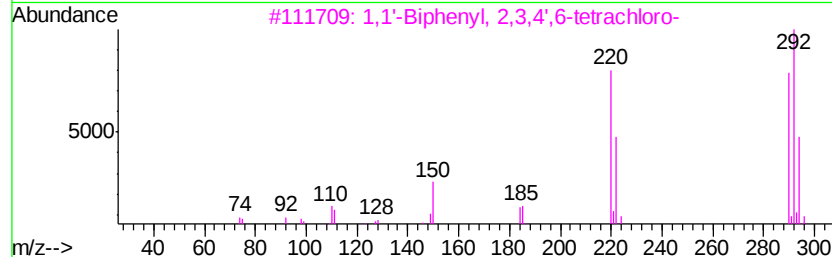
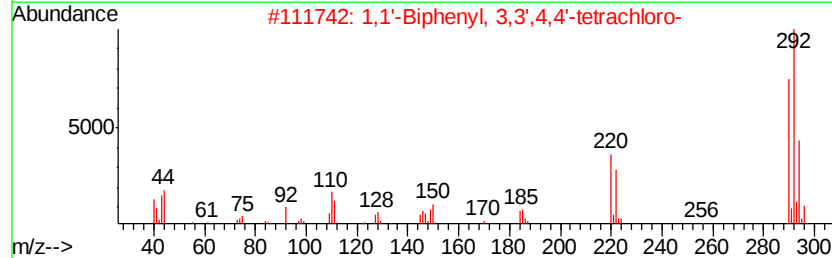
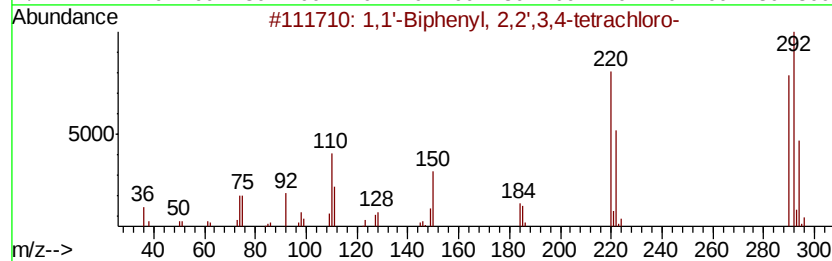
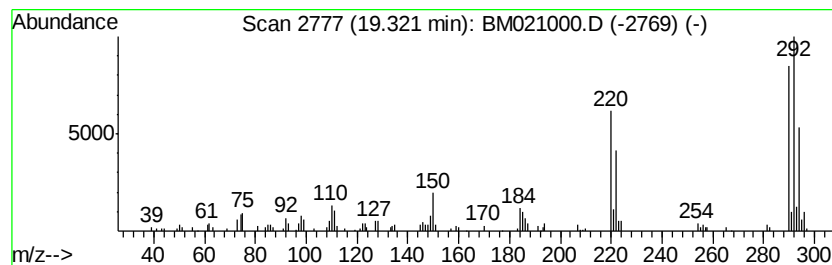
Instrument :
BNA_M
ClientSampleId :
AR1248 50PPM

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 1,1'-Biphenyl, 2,2',3,4-tet... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.			
19.32	2.03 ng/ul	380344	Chrysene-d12	21.35			
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	1,1'	-Biphenyl	3,3',4,4'-tetrach...	290	C12H6Cl4	032598-13-3	99
3	1,1'	-Biphenyl	2,3,4',6-tetrachl...	290	C12H6Cl4	052663-58-8	99
4	1,1'	-Biphenyl	2,3',4',6-tetrach...	290	C12H6Cl4	041464-46-4	99
5	1,1'	-Biphenyl	3,3',5,5'-tetrach...	290	C12H6Cl4	033284-52-5	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061719\
Data File : BM021000.D
Acq On : 18 Jun 2019 04:26
Operator : HP/JU
Sample : AR1248 50PPM
Misc : GCMS CONFIRMATION
ALS Vial : 30 Sample Multiplier: 1

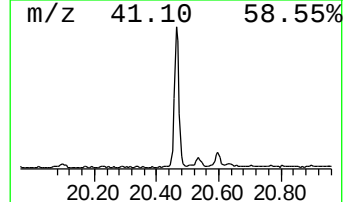
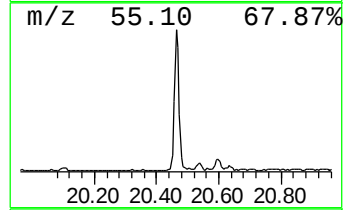
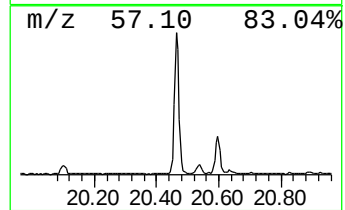
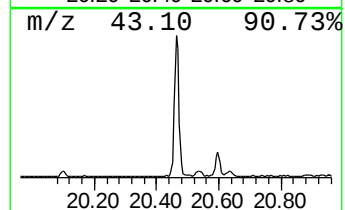
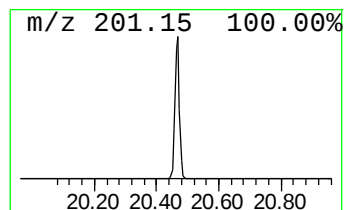
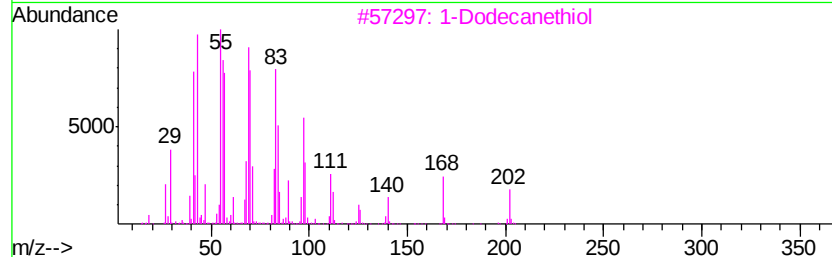
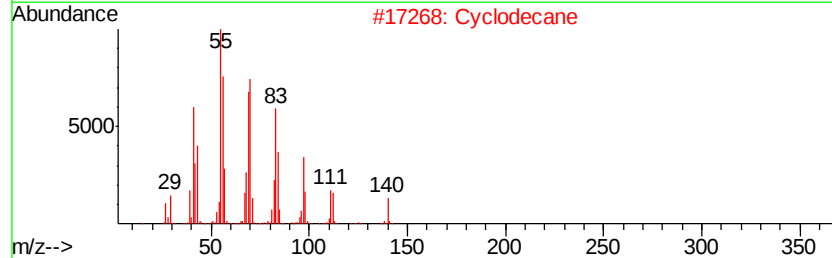
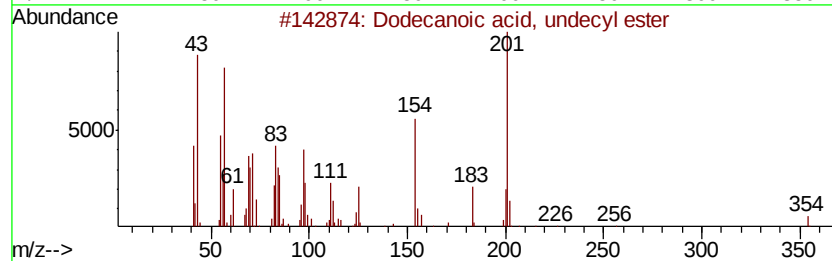
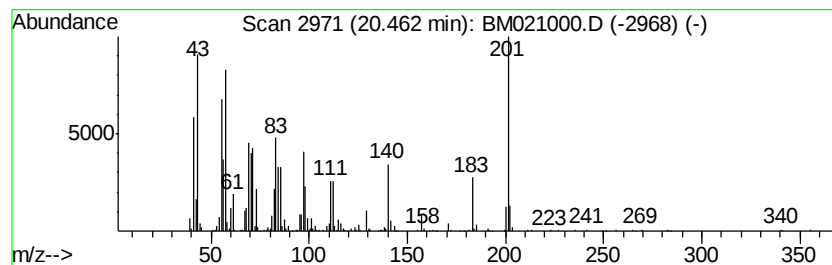
Instrument :
BNA_M
ClientSampleId :
AR1248 50PPM

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Dodecanoic acid, undecyl ester Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.46	4.81 ng/ul	902871	Chrysene-d12	21.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Dodecanoic acid, undecyl ester	354 C23H46O2	003658-44-4	78
2	Cyclodecane	140 C10H20	000293-96-9	70
3	1-Dodecanethiol	202 C12H26S	000112-55-0	53
4	Cyclodecane	140 C10H20	000293-96-9	45
5	3-Chloropropionic acid, dodecyl ...	276 C15H29ClO2	074316-16-8	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM061719\
Data File : BM021000.D
Acq On : 18 Jun 2019 04:26
Operator : HP/JU
Sample : AR1248 50PPM
Misc : GCMS CONFIRMATION
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
AR1248 50PPM

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Str...	3.21	3.3	ng/ul	253893	1	7.79	1542270	20.0
1,1'-Biphenyl, 2,...	17.76	3.4	ng/ul	729412	4	17.17	4244370	20.0
1,1'-Biphenyl, 2,...	18.24	2.7	ng/ul	575769	4	17.17	4244370	20.0
1,1'-Biphenyl, 2,...	18.53	2.4	ng/ul	497699	4	17.17	4244370	20.0
1,1'-Biphenyl, 2,...	19.06	2.8	ng/ul	603381	4	17.17	4244370	20.0
1,1'-Biphenyl, 2,...	19.10	2.5	ng/ul	522426	4	17.17	4244370	20.0
1,1'-Biphenyl, 2,...	19.32	2.0	ng/ul	380344	5	21.35	3755080	20.0
Dodecanoic acid, ...	20.46	4.8	ng/ul	902871	5	21.35	3755080	20.0