

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061224\
 Data File : BN031918.D
 Acq On : 11 Jun 2024 20:19
 Operator : MA/JU
 Sample : P2642-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4C18

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN031918.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.175	28	32	42	rVB	66945	99078	0.99%	0.083%
2	3.446	69	78	91	rBV	45990	84817	0.85%	0.071%
3	3.587	98	102	115	rBV	386021	690002	6.92%	0.575%
4	3.687	115	119	126	rVB	39818	60251	0.60%	0.050%
5	5.758	463	471	479	rVV2	39522	66054	0.66%	0.055%
6	6.834	647	654	667	rVB	1318712	2125311	21.32%	1.772%
7	7.005	676	683	696	rVB	1070492	1703577	17.09%	1.420%
8	7.193	707	715	725	rBV	1303780	2075479	20.82%	1.730%
9	7.657	787	794	802	rBV	1015976	1658184	16.63%	1.383%
10	7.734	802	807	814	rVB2	39236	72900	0.73%	0.061%
11	8.369	907	915	923	rBV	1651791	2762001	27.70%	2.303%
12	8.481	930	934	942	rVB	39332	66864	0.67%	0.056%
13	8.816	983	991	999	rBV	1260842	2110344	21.17%	1.759%
14	9.381	1081	1087	1098	rVB	162526	250142	2.51%	0.209%
15	9.534	1105	1113	1126	rBV	1022462	1726586	17.32%	1.440%
16	10.063	1195	1203	1224	rBV	1616618	2870284	28.79%	2.393%
17	10.440	1259	1267	1274	rBV	1557322	2699896	27.08%	2.251%
18	10.493	1274	1276	1283	rVB	42404	54150	0.54%	0.045%
19	10.587	1283	1292	1311	rBV	1084803	2103744	21.10%	1.754%
20	10.993	1354	1361	1369	rBV	46407	83267	0.84%	0.069%
21	11.193	1388	1395	1411	rBV	104326	251209	2.52%	0.209%
22	12.034	1534	1538	1546	rVB	33447	55077	0.55%	0.046%
23	13.122	1712	1723	1734	rBV8	24188	55238	0.55%	0.046%
24	13.722	1818	1825	1839	rVB	3181225	4492484	45.06%	3.746%
25	13.998	1858	1872	1889	rBV	3566925	5499807	55.16%	4.585%
26	14.304	1916	1924	1930	rBV	2652028	3857666	38.69%	3.216%
27	14.369	1930	1935	1942	rVB	120438	190535	1.91%	0.159%
28	14.516	1953	1960	1972	rBV	1390045	2353954	23.61%	1.963%
29	14.675	1982	1987	1989	rBV2	36883	55682	0.56%	0.046%
30	14.704	1989	1992	1995	rVV	47757	64779	0.65%	0.054%
31	15.298	2086	2093	2099	rBV	4639497	6707918	67.28%	5.593%
32	15.357	2099	2103	2108	rVB2	126281	184663	1.85%	0.154%
33	15.428	2108	2115	2122	rBV	1415932	1996920	20.03%	1.665%
34	15.563	2135	2138	2142	rVV3	52695	81476	0.82%	0.068%
35	15.651	2148	2153	2161	rVB	40637	74299	0.75%	0.062%
36	15.798	2174	2178	2186	rVB2	41584	83936	0.84%	0.070%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
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37	16.033	2212	2218	2224	rVB5	36535	73049	0.73%	0.061%
38	16.363	2266	2274	2278	rVV4	31202	76297	0.77%	0.064%
39	16.586	2305	2312	2316	rBV6	40256	98496	0.99%	0.082%
40	16.710	2328	2333	2339	rVB	73639	118675	1.19%	0.099%

41	16.869	2356	2360	2367	rVB	104391	162289	1.63%	0.135%
42	16.933	2367	2371	2374	rBV	63978	87275	0.88%	0.073%
43	16.969	2374	2377	2382	rVB4	51165	67384	0.68%	0.056%
44	17.051	2385	2391	2395	rBV	3227729	4777755	47.92%	3.983%
45	17.098	2395	2399	2403	rVB	1783266	2490918	24.98%	2.077%

46	17.151	2403	2408	2413	rBV	4852031	7078321	71.00%	5.902%
47	17.239	2419	2423	2428	rBV5	30045	71613	0.72%	0.060%
48	17.463	2453	2461	2465	rBV2	170245	310443	3.11%	0.259%
49	17.504	2465	2468	2472	rVB3	38283	56755	0.57%	0.047%
50	17.575	2477	2480	2486	rVB2	43390	66605	0.67%	0.056%

51	17.657	2486	2494	2499	rBV	574282	1018072	10.21%	0.849%
52	17.769	2509	2513	2520	rVB2	78098	132603	1.33%	0.111%
53	17.927	2531	2540	2541	rBV2	241030	398892	4.00%	0.333%
54	17.957	2541	2545	2554	rVB2	943399	1729510	17.35%	1.442%
55	18.057	2559	2562	2567	rVV	113581	172945	1.73%	0.144%

56	18.122	2567	2573	2577	rVV2	621565	993632	9.97%	0.828%
57	18.157	2577	2579	2581	rVV	170154	212614	2.13%	0.177%
58	18.180	2581	2583	2587	rVV	205035	255909	2.57%	0.213%
59	18.227	2587	2591	2597	rVB3	193529	283458	2.84%	0.236%
60	18.280	2597	2600	2605	rBV2	81889	115719	1.16%	0.096%

61	18.410	2618	2622	2624	rBV	194791	271160	2.72%	0.226%
62	18.433	2624	2626	2629	rVV	216521	268030	2.69%	0.223%
63	18.474	2629	2633	2637	rVV2	183023	306843	3.08%	0.256%
64	18.510	2637	2639	2643	rVV	118080	160858	1.61%	0.134%
65	18.551	2643	2646	2649	rVV2	73985	99899	1.00%	0.083%

66	18.586	2649	2652	2657	rVB	166084	202345	2.03%	0.169%
67	18.680	2664	2668	2673	rVB3	118361	175317	1.76%	0.146%
68	18.769	2679	2683	2686	rBV3	59991	69637	0.70%	0.058%
69	18.851	2692	2697	2701	rBV2	126637	247928	2.49%	0.207%
70	18.898	2701	2705	2710	rVV4	249215	408161	4.09%	0.340%

71	18.951	2710	2714	2717	rVV2	365343	513659	5.15%	0.428%
72	18.992	2717	2721	2728	rVB3	426554	730608	7.33%	0.609%
73	19.116	2736	2742	2748	rVB	3276551	4338268	43.51%	3.617%
74	19.198	2749	2756	2757	rBV3	142678	256000	2.57%	0.213%
75	19.245	2760	2764	2772	rVV3	611343	1090545	10.94%	0.909%

76	19.322	2774	2777	2781	rVB3	92794	115242	1.16%	0.096%
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Integration Parameters: LSCINT.P

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

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 Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
 Title : SVOA CALIBRATION

77	19.451	2793	2799	2803	rBV	6172364	9970106	100.00%	8.313%
78	19.480	2803	2804	2813	rVV	2613682	2512225	25.20%	2.095%
79	19.557	2813	2817	2822	rVB3	109333	167745	1.68%	0.140%
80	19.604	2822	2825	2829	rVB4	59327	69183	0.69%	0.058%
81	19.663	2830	2835	2839	rBV	148217	241903	2.43%	0.202%
82	19.710	2839	2843	2850	rBV3	170267	277384	2.78%	0.231%
83	19.804	2855	2859	2866	rBV3	150513	363294	3.64%	0.303%
84	19.945	2878	2883	2884	rBV2	140926	195846	1.96%	0.163%
85	19.980	2885	2889	2892	rVB2	314869	454500	4.56%	0.379%
86	20.080	2902	2906	2910	rBV	231751	315453	3.16%	0.263%
87	20.139	2910	2916	2919	rVV2	226341	390126	3.91%	0.325%
88	20.280	2937	2940	2942	rVV	102197	106994	1.07%	0.089%
89	20.327	2944	2948	2955	rVB2	212429	327282	3.28%	0.273%
90	20.727	3013	3016	3023	rVB	215070	321400	3.22%	0.268%
91	20.939	3049	3052	3055	rBV	210191	244762	2.45%	0.204%
92	20.980	3055	3059	3066	rBV3	752388	1383509	13.88%	1.153%
93	21.198	3093	3096	3101	rVV	653727	800706	8.03%	0.668%
94	21.292	3104	3112	3116	rVV3	3701536	7479236	75.02%	6.236%
95	21.333	3116	3119	3131	rVB	1314777	1921652	19.27%	1.602%
96	21.474	3139	3143	3150	rBV3	206272	385247	3.86%	0.321%
97	22.033	3235	3238	3243	rBV3	227937	314775	3.16%	0.262%
98	23.298	3446	3453	3459	rBV	767718	2177318	21.84%	1.815%
99	24.021	3568	3576	3582	rBV2	2326375	5444017	54.60%	4.539%
100	24.221	3602	3610	3625	rVB2	1722024	4603771	46.18%	3.838%

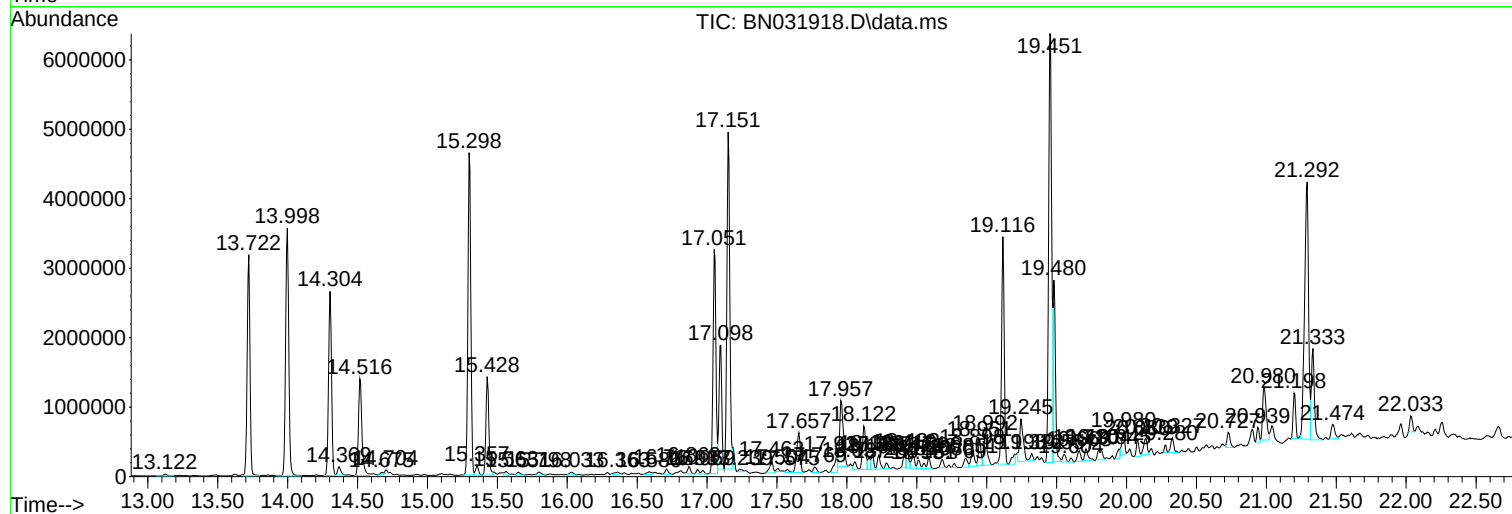
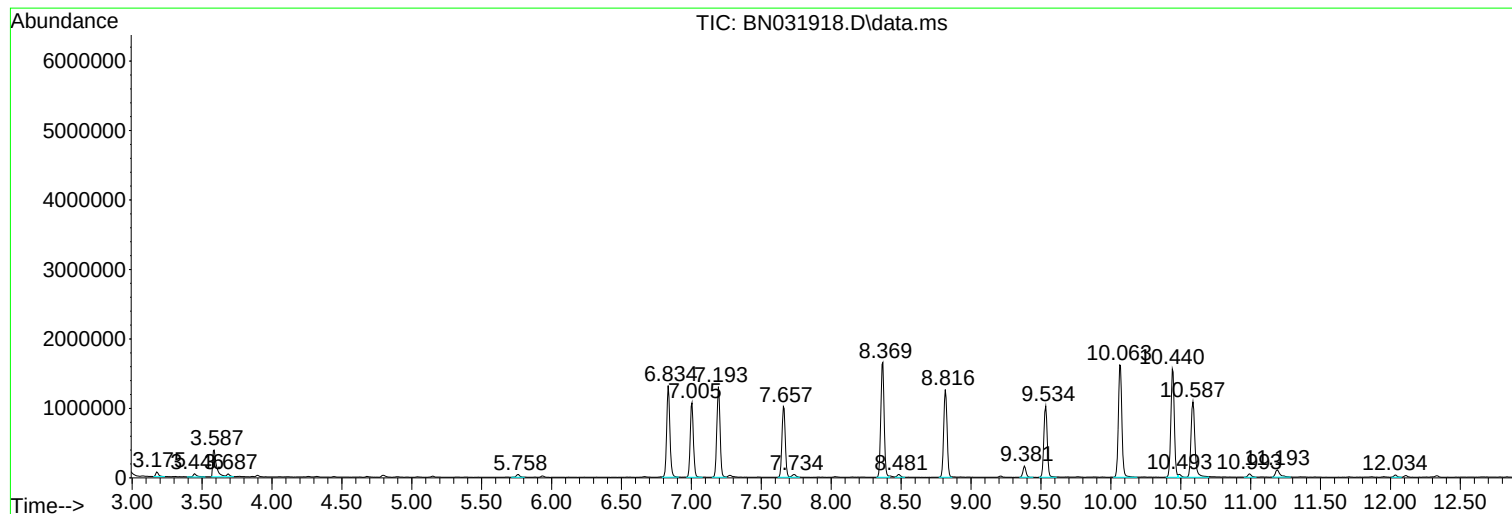
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ALS Vial : 8 Sample Multiplier: 1

Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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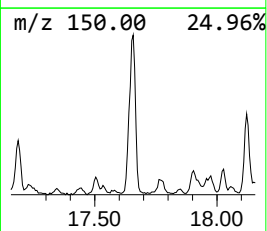
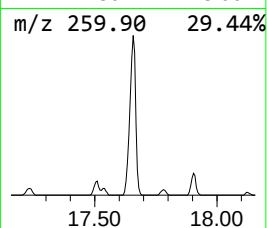
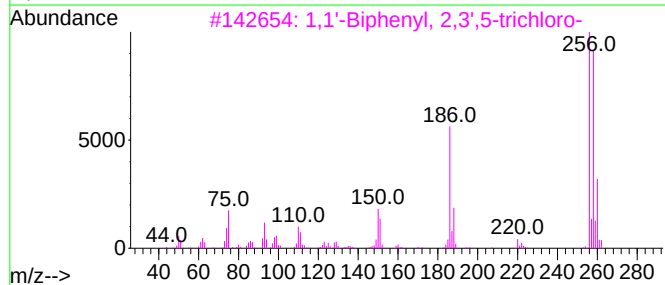
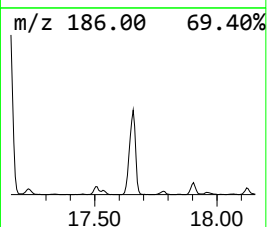
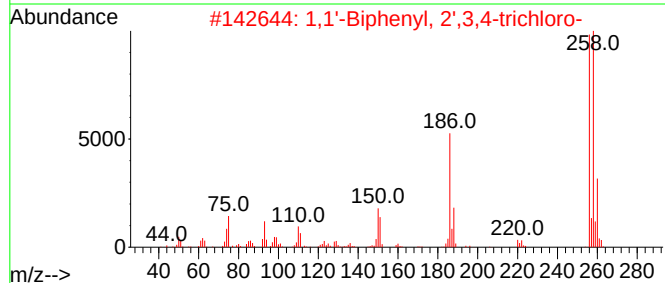
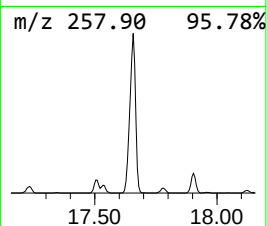
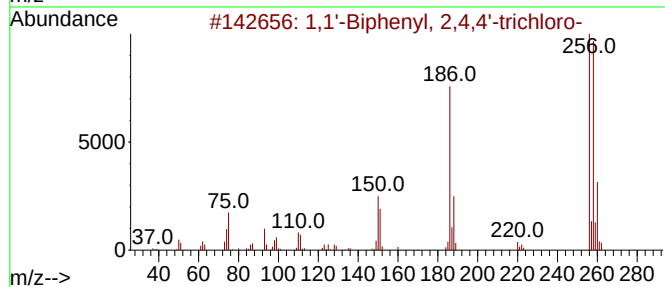
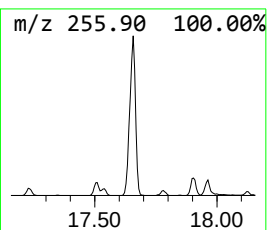
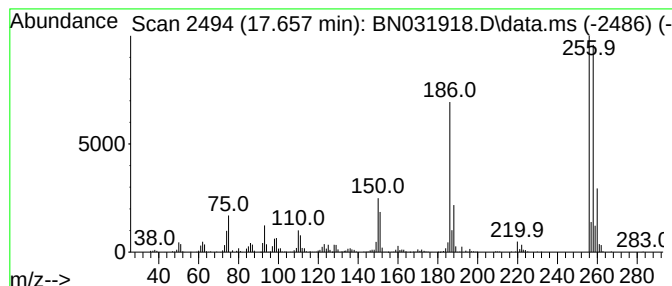
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.657	4.26 ng/ul	1018070	Phenanthrene-d10	17.051

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, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99		
5	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	99		



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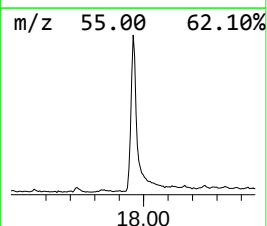
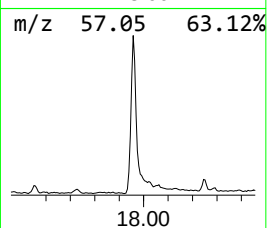
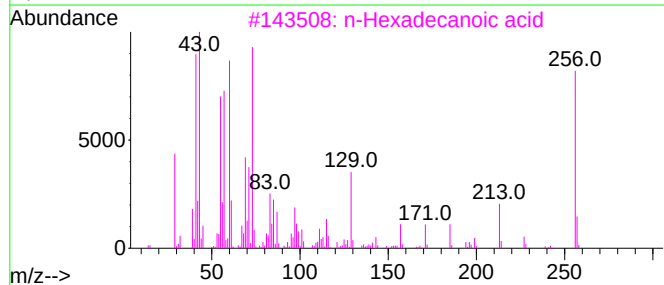
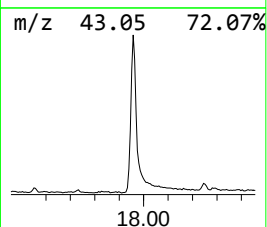
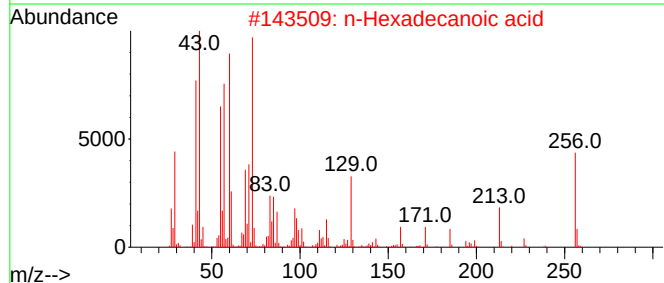
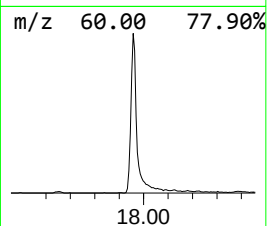
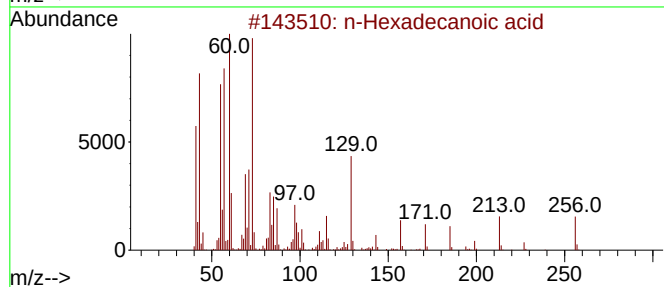
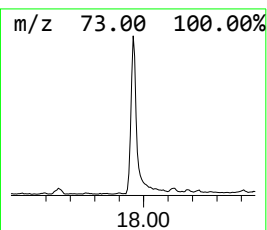
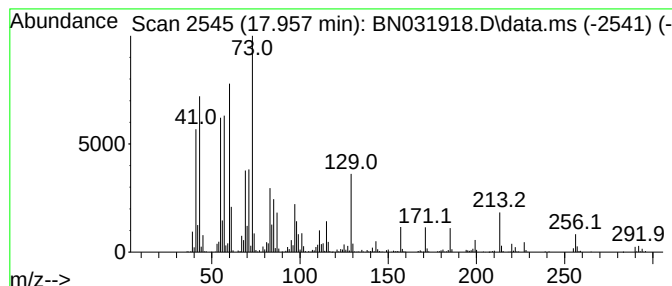
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.957	7.24 ng/ul	1729510	Phenanthrene-d10	17.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
5			Tridecanoic acid	214	C13H26O2	000638-53-9	91



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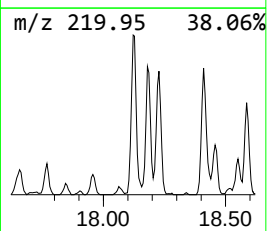
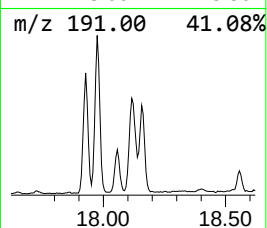
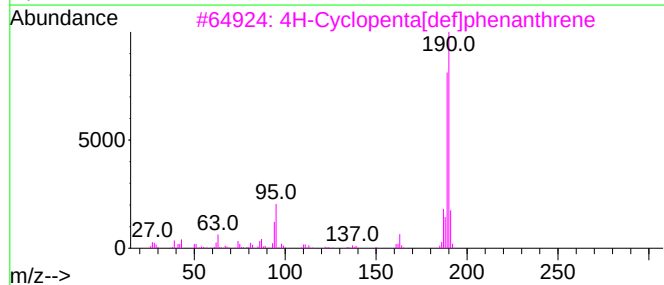
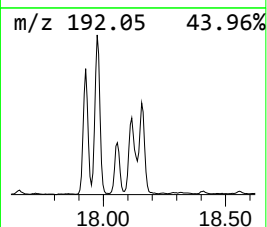
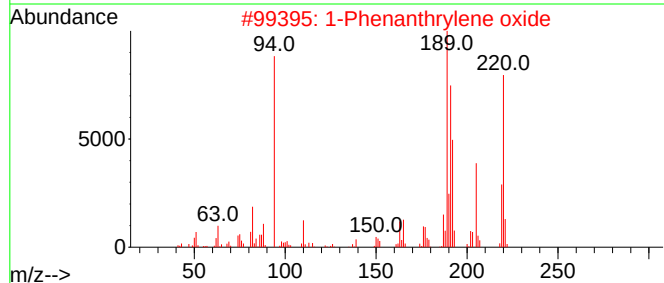
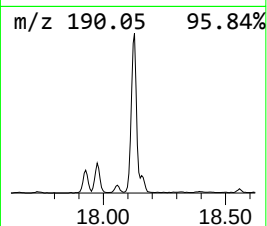
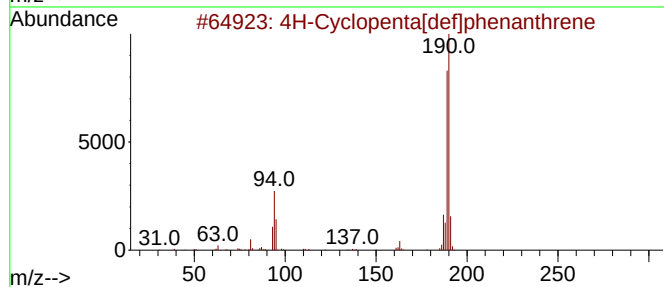
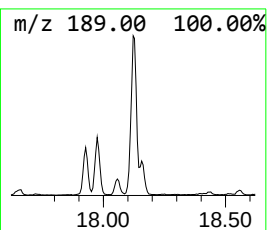
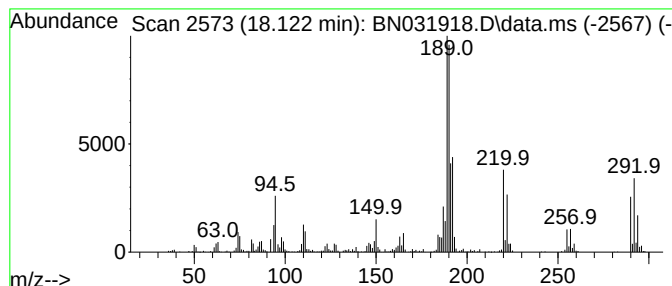
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.122	4.16 ng/ul	993632	Phenanthrene-d10	17.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
2			1-Phenanthrylene oxide	220	C16H12O	026698-45-3	46
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	41
4			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	38
5			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061224\
Data File : BN031918.D
Acq On : 11 Jun 2024 20:19
Operator : MA/JU
Sample : P2642-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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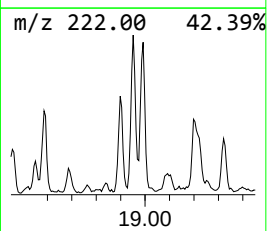
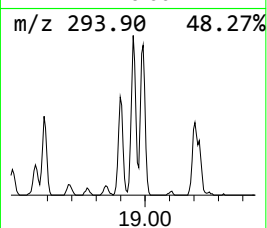
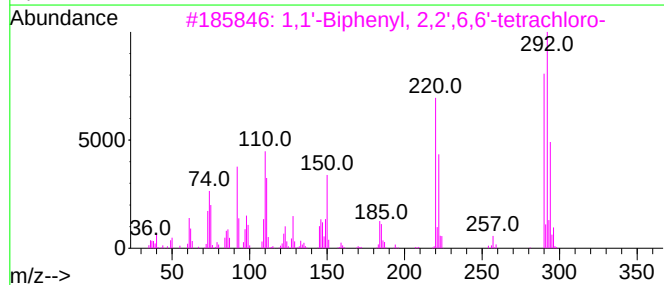
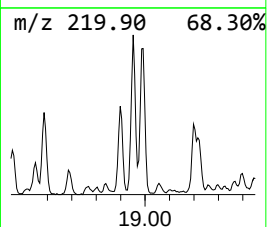
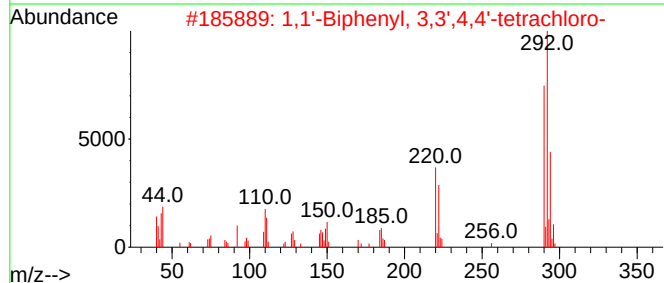
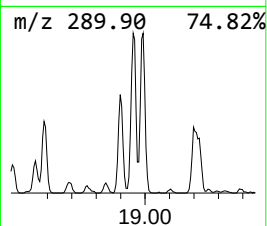
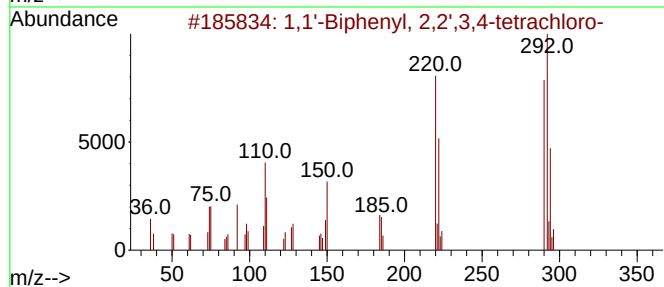
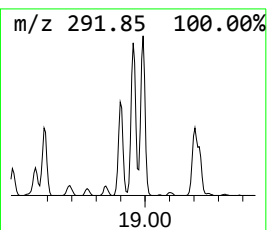
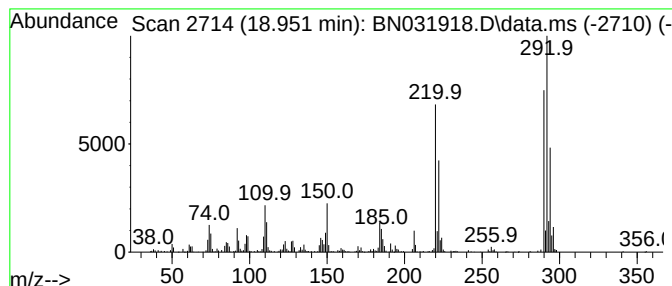
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.951	2.15 ng/ul	513659	Phenanthrene-d10	17.051

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, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99		
4	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	97		
5	1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	97		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061224\
Data File : BN031918.D
Acq On : 11 Jun 2024 20:19
Operator : MA/JU
Sample : P2642-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

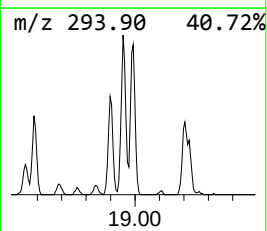
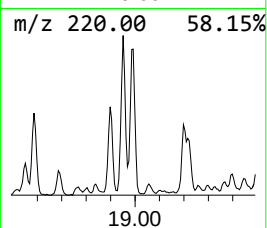
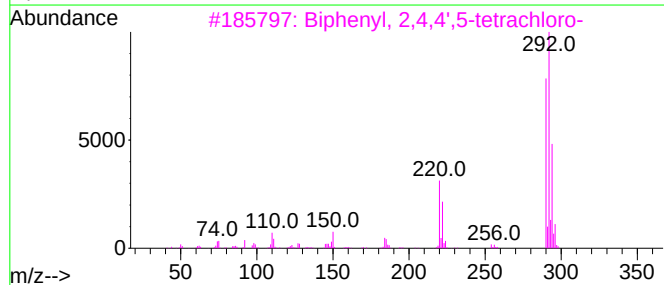
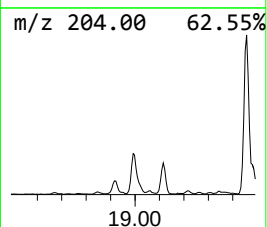
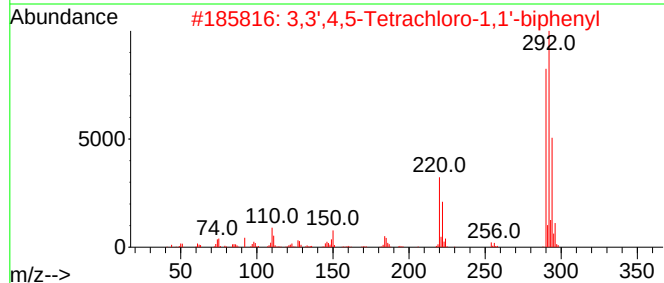
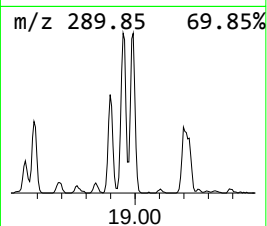
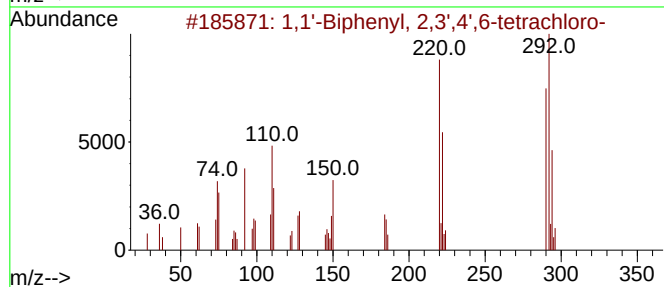
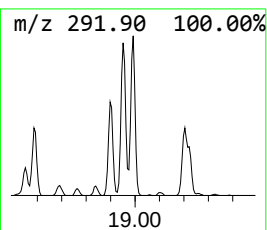
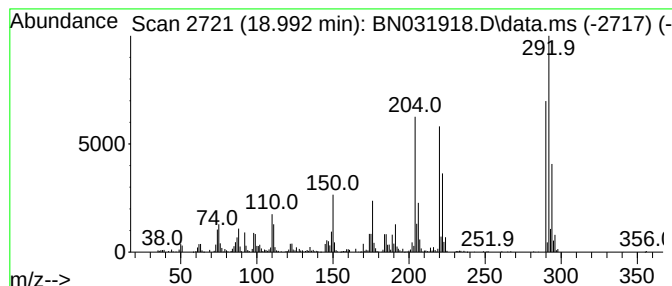
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.992	3.06 ng/ul	730608	Phenanthrene-d10	17.051	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3',4',6-tetrach...	290	C12H6Cl4	041464-46-4	94
2	3,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-49-1	94
3	Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	94
4	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	94
5	1,1'-Biphenyl, 2,2',3,4-tetrachl...	290	C12H6Cl4	052663-59-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061224\
Data File : BN031918.D
Acq On : 11 Jun 2024 20:19
Operator : MA/JU
Sample : P2642-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A4C18

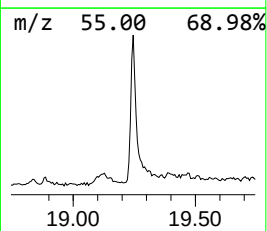
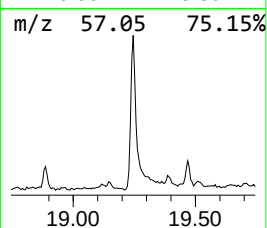
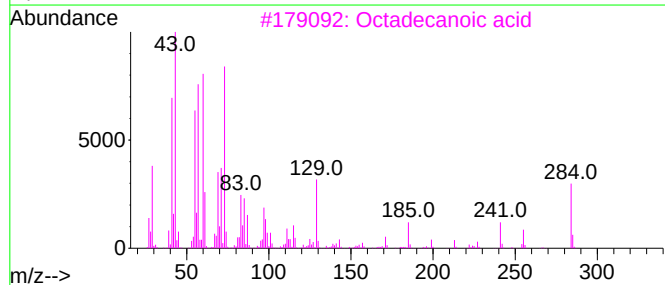
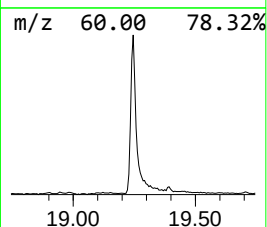
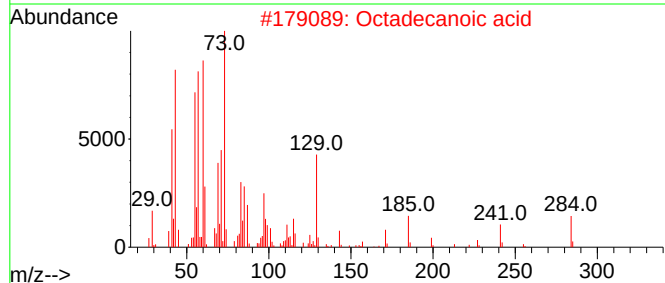
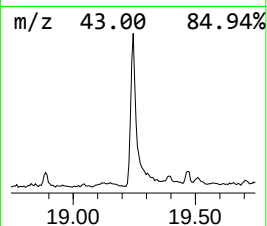
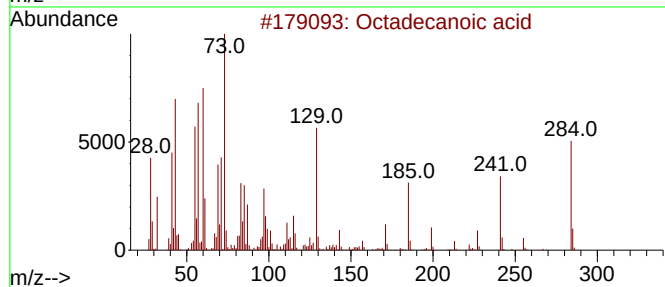
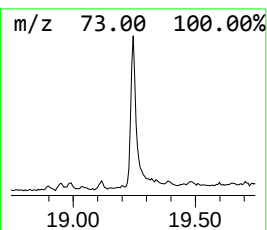
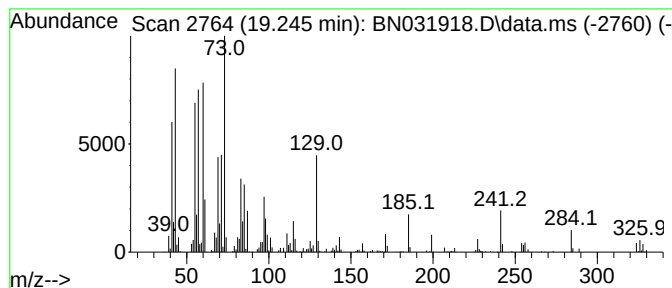
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.245	2.92 ng/ul	1090550	Chrysene-d12	21.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	98
3			Octadecanoic acid	284	C18H36O2	000057-11-4	95
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061224\
Data File : BN031918.D
Acq On : 11 Jun 2024 20:19
Operator : MA/JU
Sample : P2642-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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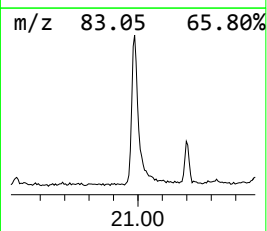
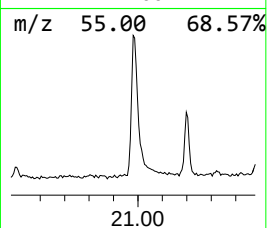
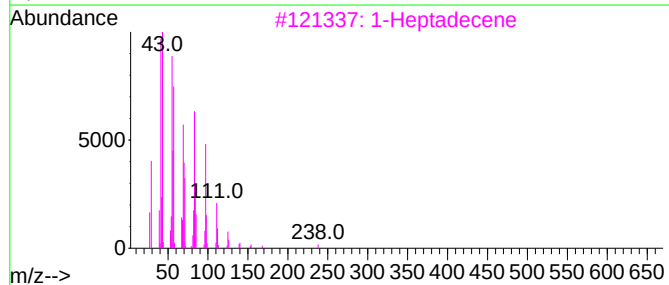
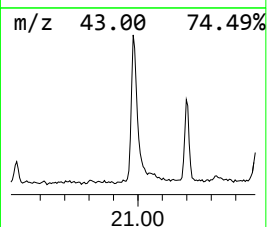
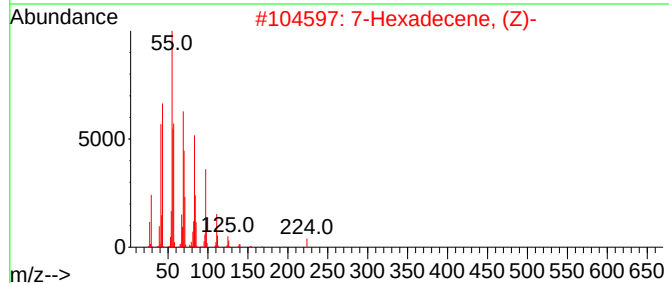
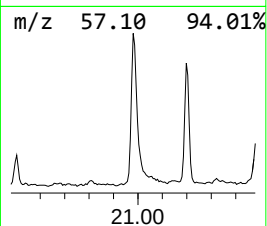
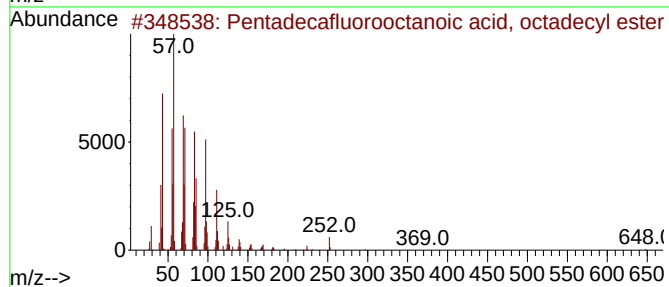
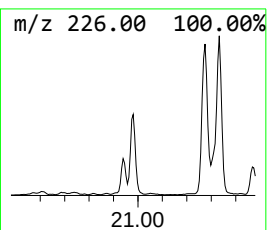
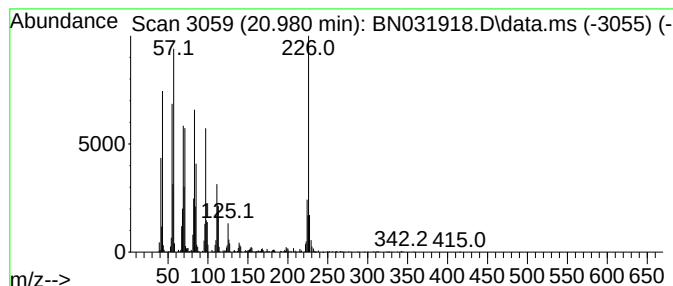
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Pentadecafluorooctanoic aci... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.980	3.70 ng/ul	1383510	Chrysene-d12	21.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	90
2			7-Hexadecene, (Z)-	224	C16H32	035507-09-6	86
3			1-Heptadecene	238	C17H34	006765-39-5	86
4			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	60
5			Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061224\
Data File : BN031918.D
Acq On : 11 Jun 2024 20:19
Operator : MA/JU
Sample : P2642-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1,1'-Biphenyl, ...	17.657	4.3	ng/ul	1018070	4	17.051	4777760	20.0
n-Hexadecanoic ...	17.957	7.2	ng/ul	1729510	4	17.051	4777760	20.0
unknown-01	18.122	4.2	ng/ul	993632	4	17.051	4777760	20.0
1,1'-Biphenyl, ...	18.951	2.1	ng/ul	513659	4	17.051	4777760	20.0
1,1'-Biphenyl, ...	18.992	3.1	ng/ul	730608	4	17.051	4777760	20.0
Octadecanoic acid	19.245	2.9	ng/ul	1090550	5	21.292	7479240	20.0
Pentadecafluoro...	20.980	3.7	ng/ul	1383510	5	21.292	7479240	20.0