

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
 Data File : BP015703.D
 Acq On : 24 Jun 2023 11:46
 Operator : MA/JU
 Sample : 03085-01
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFN7

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_P\Methods\SFAM-EPA-BP060723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015703.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.840	108	111	124	rBV	232859	370255	8.08%	0.494%
2	3.946	124	129	135	rVV	110241	154282	3.37%	0.206%
3	6.705	587	598	607	rVB4	20556	53472	1.17%	0.071%
4	7.246	684	690	704	rVB	355030	719350	15.70%	0.960%
5	7.399	710	716	730	rBV	357762	596229	13.01%	0.796%
6	7.605	744	751	760	rVV	485358	828383	18.08%	1.106%
7	8.075	825	831	844	rVB	578902	985572	21.51%	1.315%
8	8.804	948	955	968	rBV	365945	738805	16.13%	0.986%
9	9.257	1026	1032	1050	rVV	416748	842813	18.40%	1.125%
10	9.987	1149	1156	1166	rBV	361404	700958	15.30%	0.936%
11	10.540	1243	1250	1270	rBV	364267	932970	20.36%	1.245%
12	10.904	1305	1312	1318	rVV	760275	1302277	28.42%	1.738%
13	10.951	1318	1320	1328	rVV2	43997	74186	1.62%	0.099%
14	11.069	1330	1340	1360	rVB2	109917	363715	7.94%	0.485%
15	11.898	1476	1481	1487	rBV	54548	78899	1.72%	0.105%
16	12.293	1543	1548	1555	rBV	32053	56126	1.23%	0.075%
17	12.575	1590	1596	1601	rBV	27993	57243	1.25%	0.076%
18	13.234	1703	1708	1717	rVB	45465	75527	1.65%	0.101%
19	13.522	1753	1757	1762	rBV	46745	67745	1.48%	0.090%
20	14.045	1840	1846	1851	rBV2	35280	70162	1.53%	0.094%
21	14.134	1851	1861	1866	rVV	954793	1465446	31.98%	1.956%
22	14.175	1866	1868	1873	rVV2	88609	121716	2.66%	0.162%
23	14.422	1904	1910	1925	rVV2	1103112	1887191	41.19%	2.519%
24	14.592	1934	1939	1948	rVB	60971	94164	2.06%	0.126%
25	14.728	1956	1962	1969	rVV	1037845	1566407	34.19%	2.091%
26	14.787	1969	1972	1980	rVB6	28583	61773	1.35%	0.082%
27	14.975	1999	2004	2022	rVV2	126897	442241	9.65%	0.590%
28	15.128	2026	2030	2039	rVV3	66875	144260	3.15%	0.193%
29	15.204	2039	2043	2051	rVB5	26424	61947	1.35%	0.083%
30	15.504	2087	2094	2098	rBV6	40661	83481	1.82%	0.111%
31	15.545	2098	2101	2106	rVB	79839	93409	2.04%	0.125%
32	15.722	2124	2131	2137	rBV	1249046	1908623	41.66%	2.548%
33	15.869	2151	2156	2165	rBV	305679	533584	11.65%	0.712%
34	16.086	2187	2193	2204	rVB2	199323	335912	7.33%	0.448%
35	16.433	2244	2252	2267	rVB	166668	422932	9.23%	0.565%
36	17.151	2368	2374	2380	rBV3	49654	100276	2.19%	0.134%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
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37	17.210	2380	2384	2390	rBV	129117	174872	3.82%	0.233%
38	17.363	2406	2410	2416	rVB2	43339	66302	1.45%	0.088%
39	17.486	2425	2431	2435	rBV	1250025	1768551	38.60%	2.361%
40	17.528	2435	2438	2443	rVV	426048	630676	13.77%	0.842%

41	17.586	2443	2448	2453	rVV2	1289377	1976567	43.14%	2.638%
42	17.628	2453	2455	2463	rVB	172053	225035	4.91%	0.300%
43	17.904	2497	2502	2506	rBV2	73967	127025	2.77%	0.170%
44	17.951	2506	2510	2515	rVB	127597	163818	3.58%	0.219%
45	18.086	2528	2533	2537	rBV5	31959	52110	1.14%	0.070%

46	18.286	2563	2567	2572	rBV	108146	160015	3.49%	0.214%
47	18.345	2572	2577	2582	rVV2	1052590	1348400	29.43%	1.800%
48	18.398	2583	2586	2592	rVV2	187106	318955	6.96%	0.426%
49	18.475	2596	2599	2604	rVV	61229	96620	2.11%	0.129%
50	18.533	2604	2609	2613	rVV3	157611	293044	6.40%	0.391%

51	18.575	2613	2616	2620	rVB2	77440	104391	2.28%	0.139%
52	18.616	2620	2623	2633	rBV	142181	206953	4.52%	0.276%
53	18.745	2641	2645	2650	rVB	344954	373295	8.15%	0.498%
54	18.827	2655	2659	2664	rVV2	100107	158193	3.45%	0.211%
55	18.886	2665	2669	2674	rVB3	64766	104631	2.28%	0.140%

56	19.063	2696	2699	2705	rVB3	51973	59061	1.29%	0.079%
57	19.116	2705	2708	2711	rVB	53656	55605	1.21%	0.074%
58	19.175	2712	2718	2724	rBV	1238345	2035561	44.43%	2.717%
59	19.227	2724	2727	2732	rVV2	323387	419426	9.15%	0.560%
60	19.286	2733	2737	2740	rVV3	80052	131126	2.86%	0.175%

61	19.322	2740	2743	2747	rVV2	98009	148091	3.23%	0.198%
62	19.386	2750	2754	2759	rVB3	276597	395494	8.63%	0.528%
63	19.433	2759	2762	2764	rBV3	70823	87121	1.90%	0.116%
64	19.492	2764	2772	2778	rVV	1696526	2424260	52.91%	3.236%
65	19.545	2778	2781	2782	rVV	50249	52961	1.16%	0.071%

66	19.569	2782	2785	2793	rVV3	317871	502225	10.96%	0.670%
67	19.639	2794	2797	2800	rVB	76286	95217	2.08%	0.127%
68	19.786	2818	2822	2823	rBV	127699	145843	3.18%	0.195%
69	19.816	2823	2827	2830	rVV	1961443	2876636	62.79%	3.840%
70	19.845	2830	2832	2839	rVV	1413270	1617774	35.31%	2.159%

71	19.910	2839	2843	2849	rVB	245837	317071	6.92%	0.423%
72	20.022	2859	2862	2866	rVB	79370	105713	2.31%	0.141%
73	20.174	2880	2888	2892	rVB6	165507	407042	8.88%	0.543%
74	20.298	2903	2909	2912	rBV3	374158	625487	13.65%	0.835%
75	20.422	2927	2930	2932	rBV2	127173	161419	3.52%	0.215%

76	20.445	2932	2934	2937	rVV	114243	132304	2.89%	0.177%
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77	20.622	2960	2964	2968	rBV	207523	310893	6.79%	0.415%
78	20.692	2974	2976	2981	rVB2	106376	114800	2.51%	0.153%
79	20.751	2983	2986	2988	rBV3	94078	108958	2.38%	0.145%
80	20.786	2989	2992	2996	rVB	193762	209378	4.57%	0.279%
81	21.063	3035	3039	3049	rBV	372225	618035	13.49%	0.825%
82	21.221	3062	3066	3067	rBV2	358992	426193	9.30%	0.569%
83	21.251	3067	3071	3076	rVB3	731902	1269309	27.70%	1.694%
84	21.357	3086	3089	3096	rVB	311415	422223	9.22%	0.564%
85	21.451	3101	3105	3112	rVB	2369146	2639381	57.61%	3.523%
86	21.580	3120	3127	3131	rBV3	1851788	3943282	86.07%	5.263%
87	21.621	3131	3134	3142	rVB	1024978	1343916	29.33%	1.794%
88	21.692	3144	3146	3151	rVB3	185351	200308	4.37%	0.267%
89	21.927	3183	3186	3191	rVB3	190750	270050	5.89%	0.360%
90	22.174	3224	3228	3233	rBV2	508875	778769	17.00%	1.039%
91	23.292	3413	3418	3422	rVB	649604	1054338	23.01%	1.407%
92	23.363	3423	3430	3443	rVB	1460054	4581681	100.00%	6.115%
93	23.915	3519	3524	3528	rBV	616075	1160313	25.33%	1.549%
94	23.974	3528	3534	3539	rVV2	1228610	2608201	56.93%	3.481%
95	24.027	3539	3543	3554	rVB	831167	1628791	35.55%	2.174%
96	24.139	3556	3562	3568	rBV	949615	2021136	44.11%	2.698%
97	24.739	3659	3664	3679	rVB	1098290	2534977	55.33%	3.384%
98	24.880	3682	3688	3702	rVV3	515506	1453831	31.73%	1.941%
99	26.727	3995	4002	4015	rBV	759902	2550615	55.67%	3.404%
100	27.821	4182	4188	4205	rVB4	819615	3133520	68.39%	4.182%

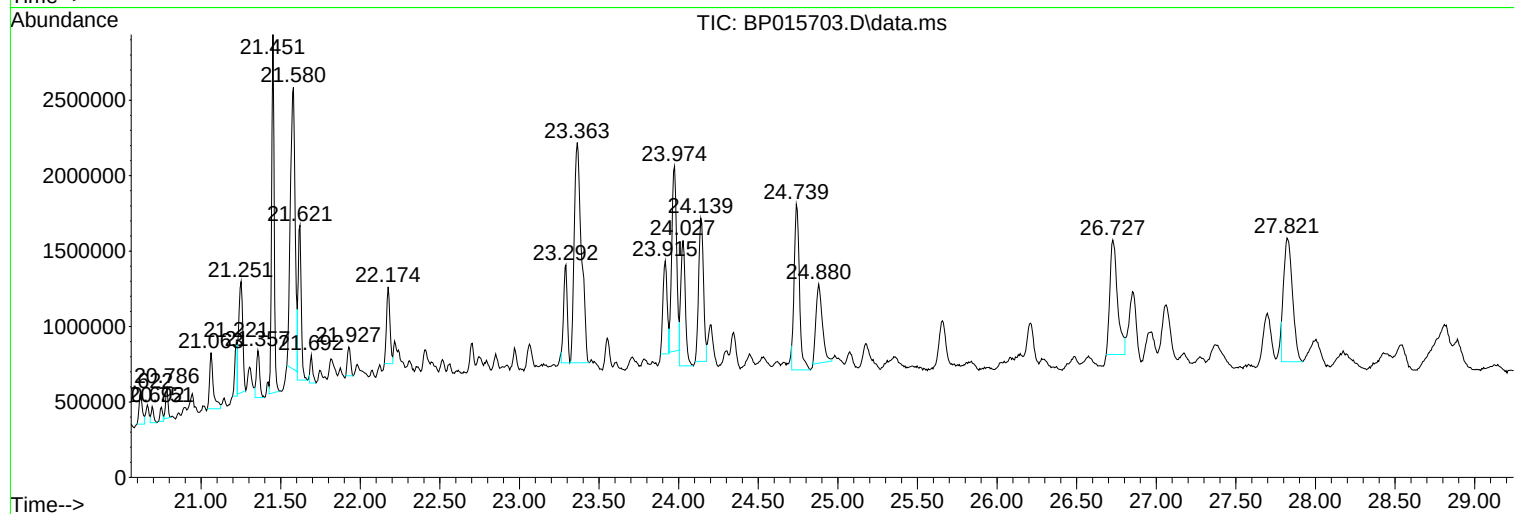
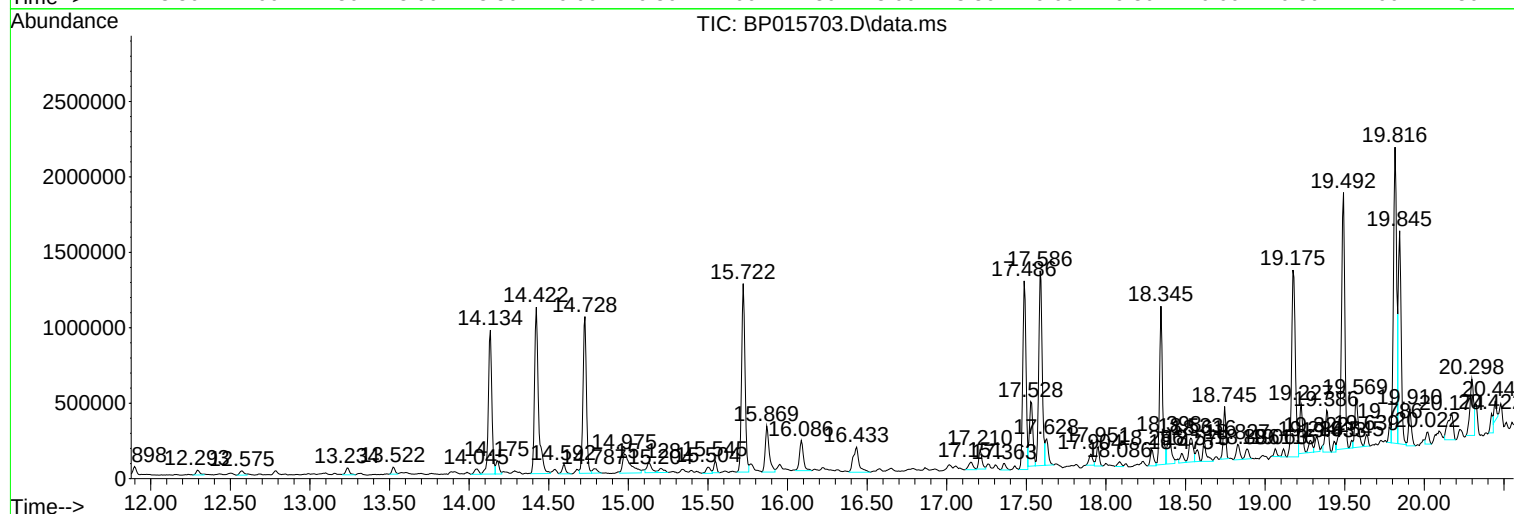
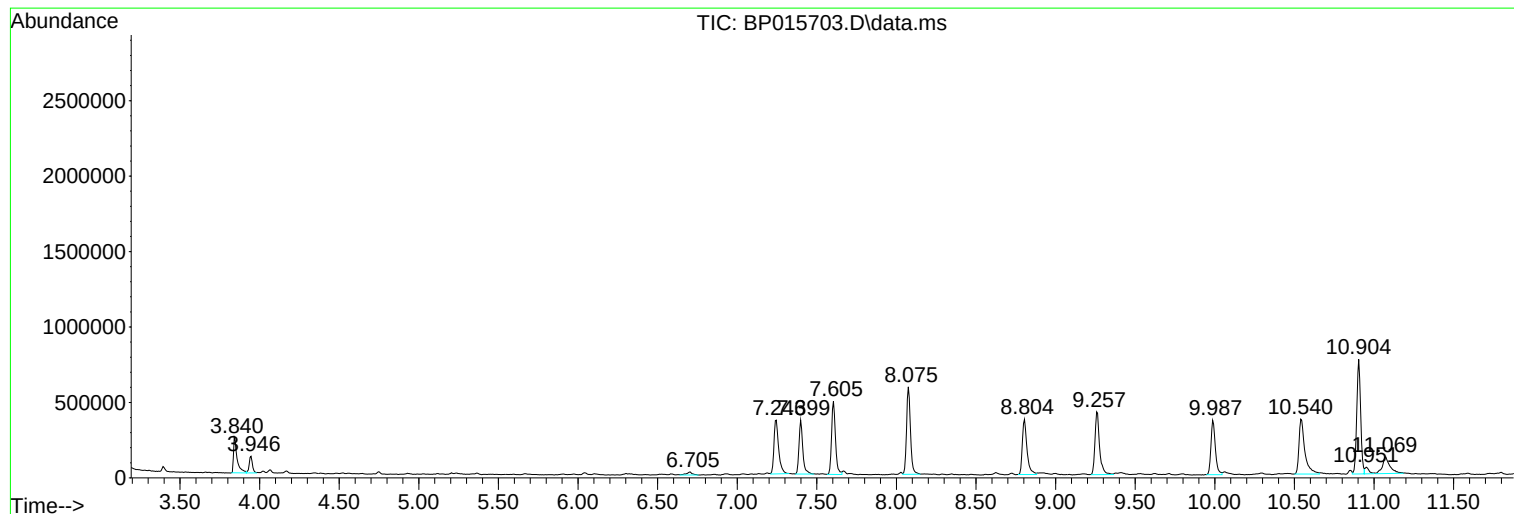
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ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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



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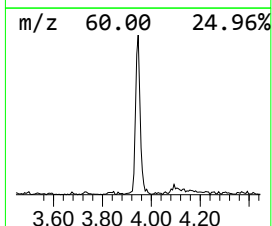
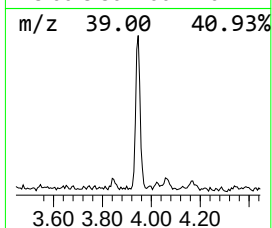
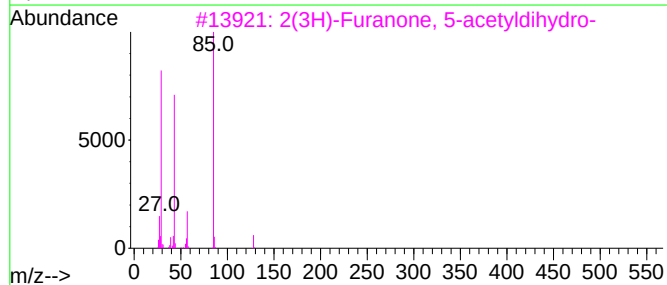
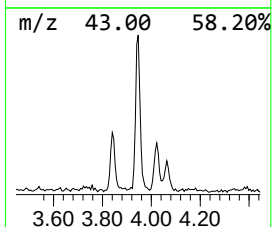
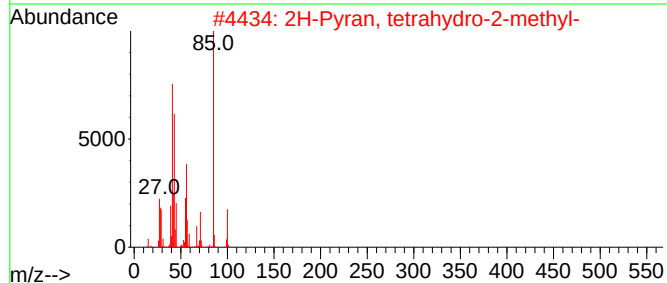
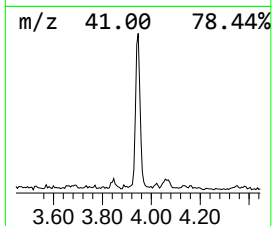
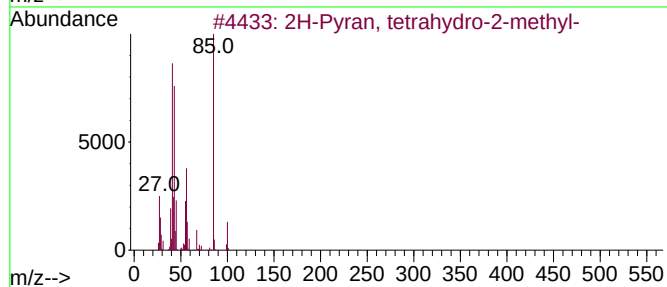
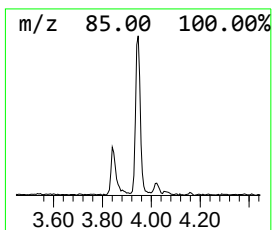
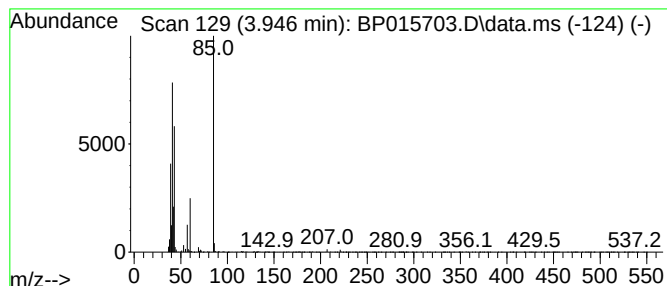
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2H-Pyran, tetrahydro-2-methyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.946	3.13 ng/ul	154282	1,4-Dichlorobenzene-d4	8.075

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	50		
2	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	33		
3	2(3H)-Furanone, 5-acetyldihydro-	128	C6H8O3	029393-32-6	23		
4	Hexane, 1,1'-[ethylidenebis(oxy)...]	230	C14H30O2	005405-58-3	23		
5	2,2'-Bi-2H-pyran, octahydro-	170	C10H18O2	016282-29-4	17		



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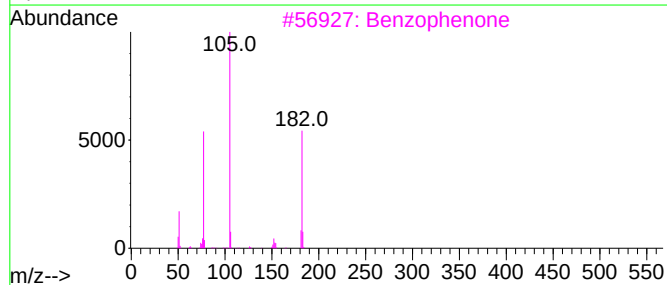
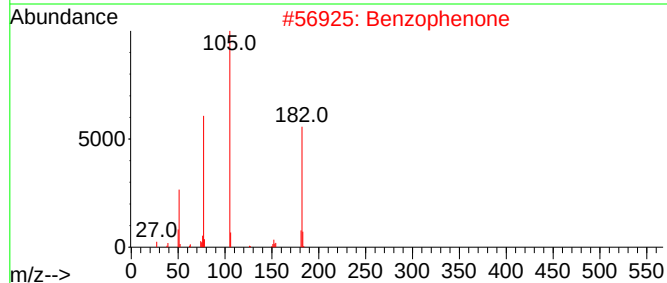
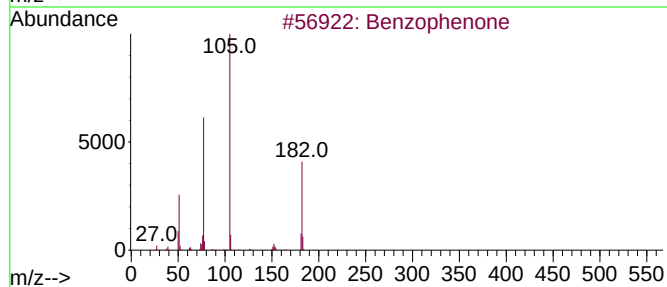
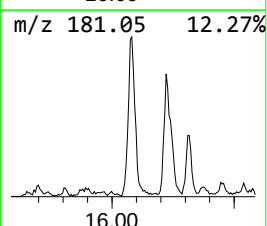
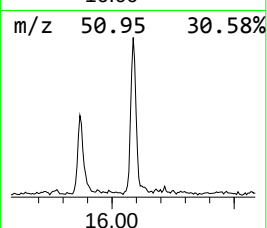
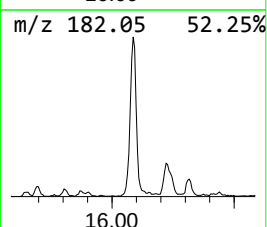
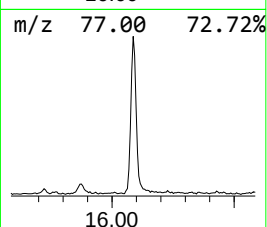
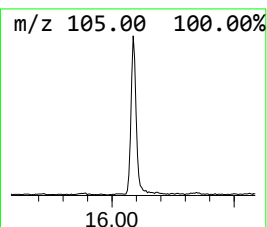
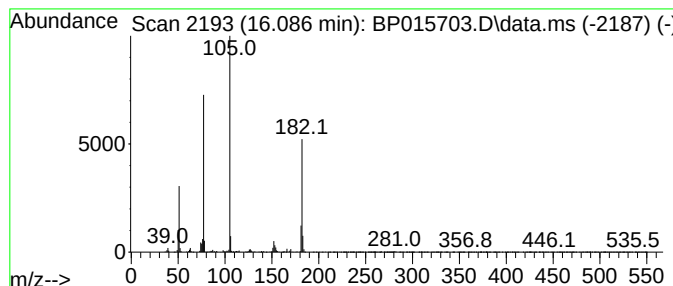
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.086	4.29 ng/ul	335912	Acenaphthene-d10	14.728

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	87



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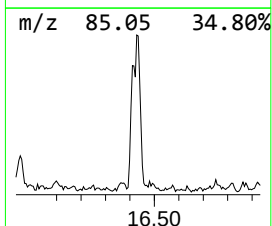
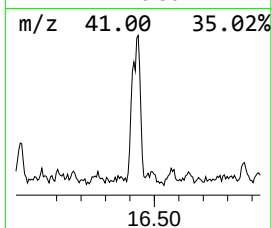
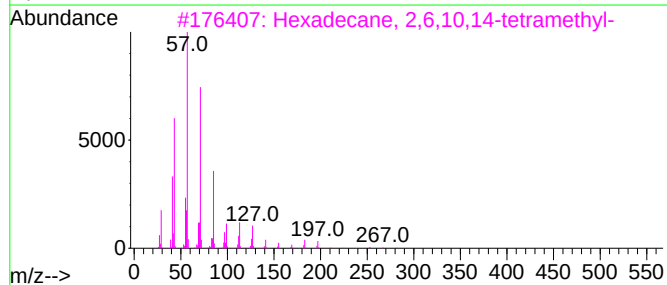
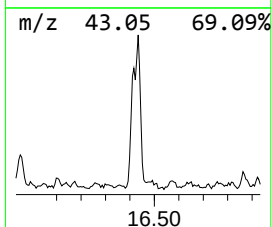
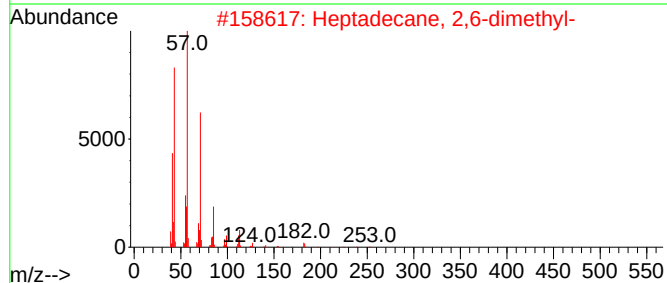
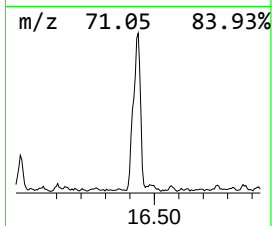
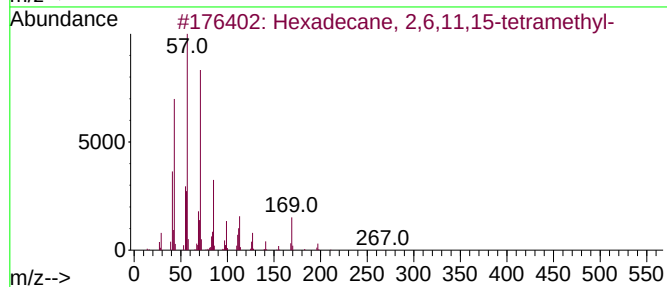
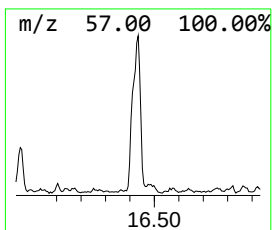
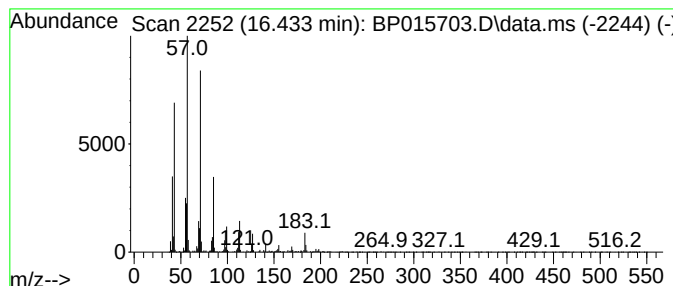
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.433	4.78 ng/ul	422932	Phenanthrene-d10	17.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	91
2			Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	87
3			Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	86
4			Dodecane	170	C12H26	000112-40-3	83
5			Undecane, 2-methyl-	170	C12H26	007045-71-8	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
Acq On : 24 Jun 2023 11:46
Operator : MA/JU
Sample : 03085-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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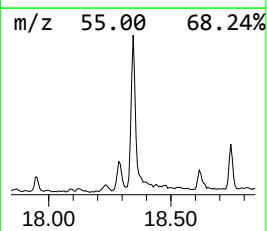
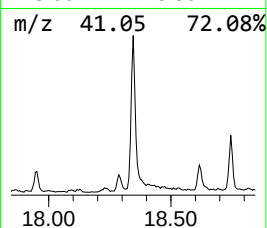
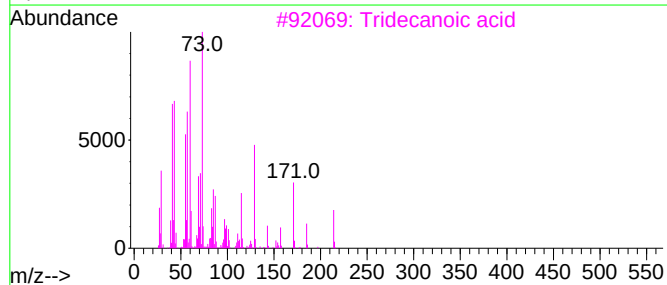
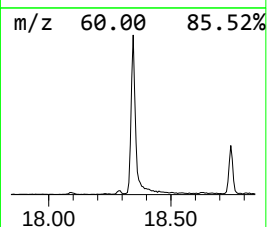
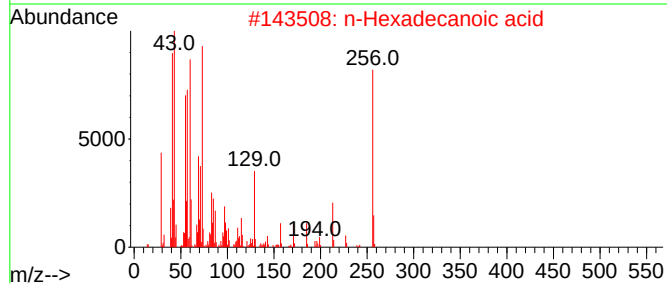
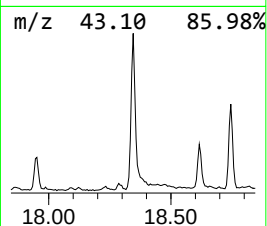
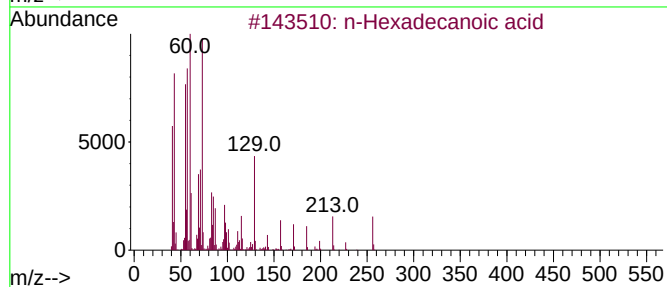
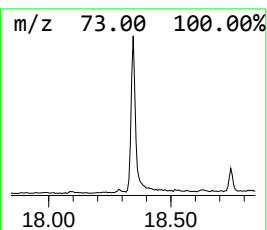
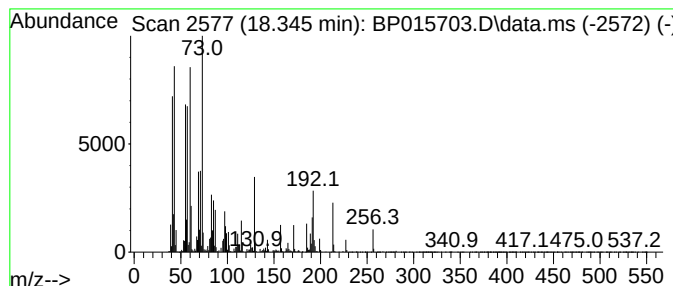
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.345	15.25 ng/ul	1348400	Phenanthrene-d10	17.486

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			Tridecanoic acid	214	C13H26O2	000638-53-9	95
4			Tridecanoic acid	214	C13H26O2	000638-53-9	93
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
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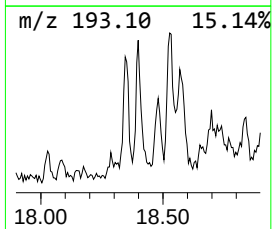
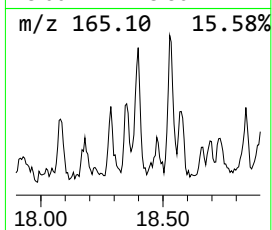
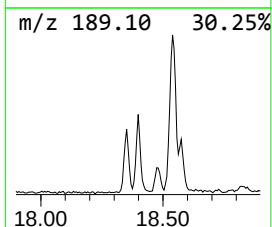
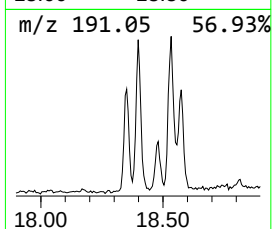
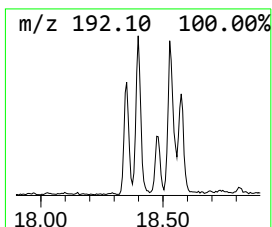
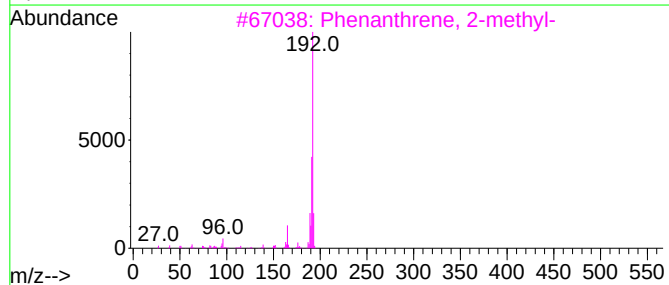
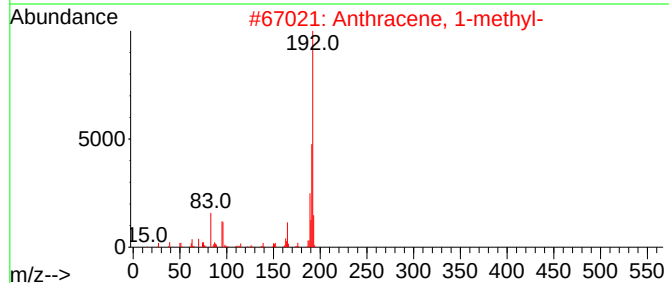
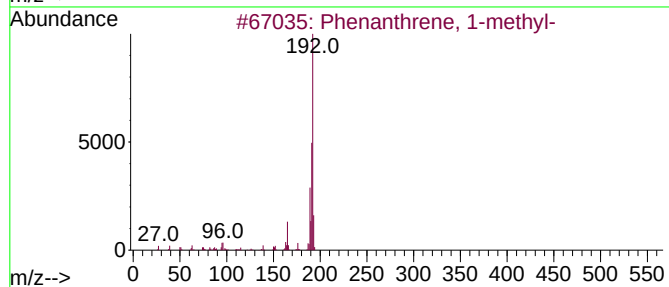
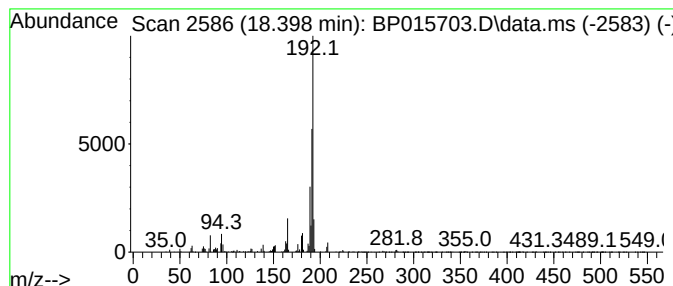
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Phenanthrene, 1-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.398	3.61 ng/ul	318955	Phenanthrene-d10	17.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94
4			Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
5			Anthracene, 2-methyl-	192	C15H12	000613-12-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
Acq On : 24 Jun 2023 11:46
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Sample : 03085-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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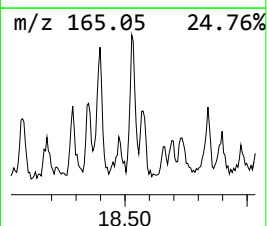
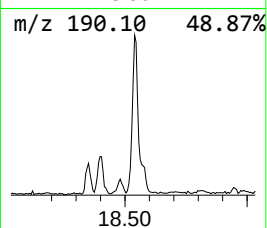
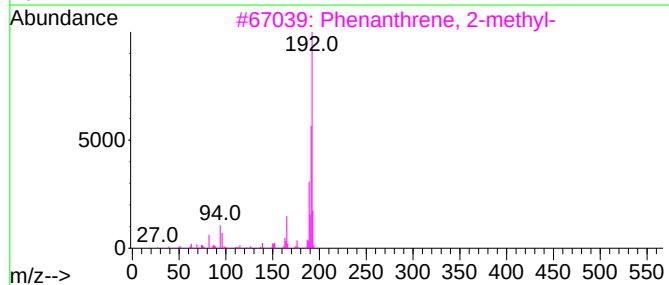
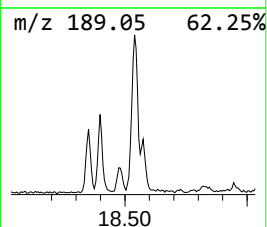
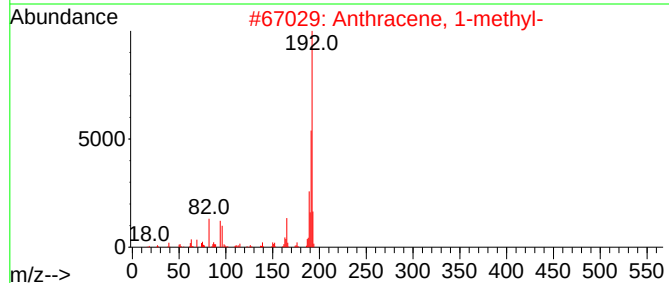
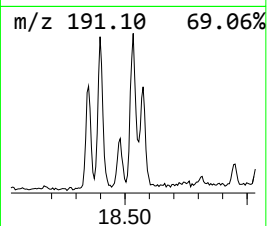
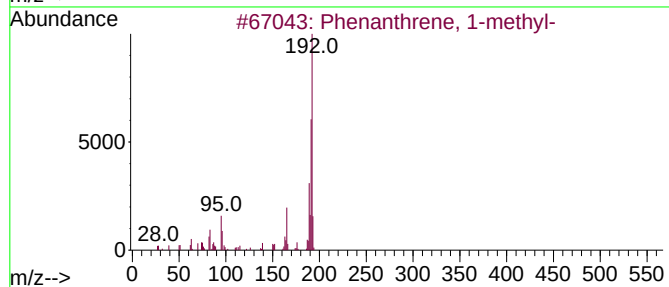
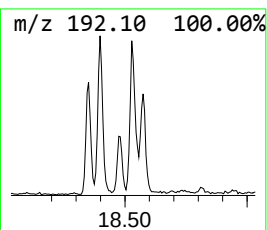
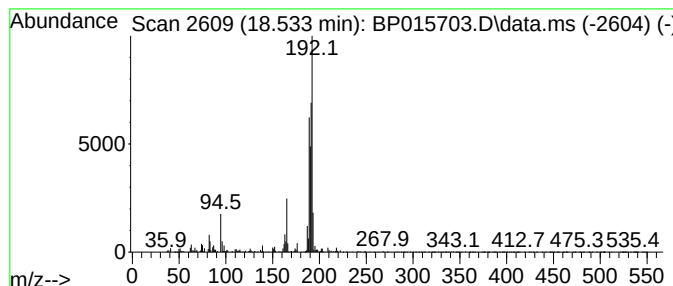
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Anthracene, 1-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.533	3.31 ng/ul	293044	Phenanthrene-d10	17.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	78
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	64
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	64
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	60
5			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
Acq On : 24 Jun 2023 11:46
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Sample : 03085-01
Misc :
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Instrument :
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ClientSampleId :
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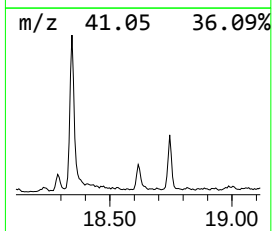
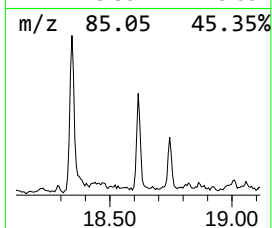
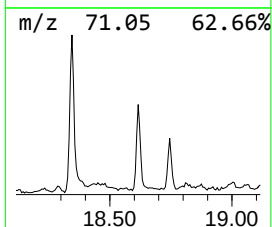
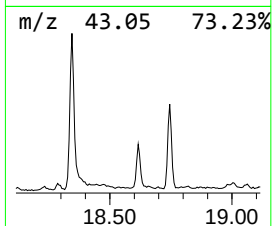
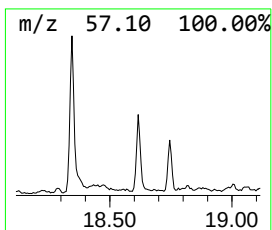
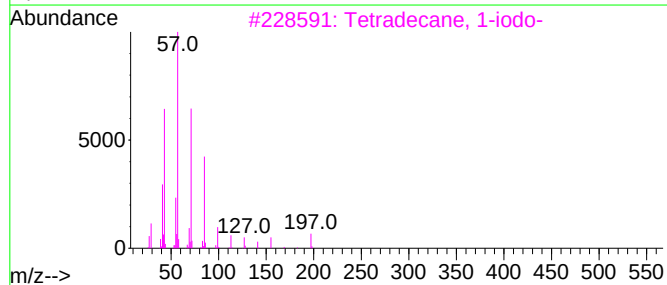
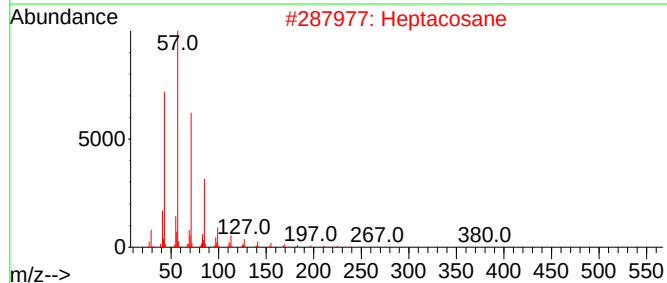
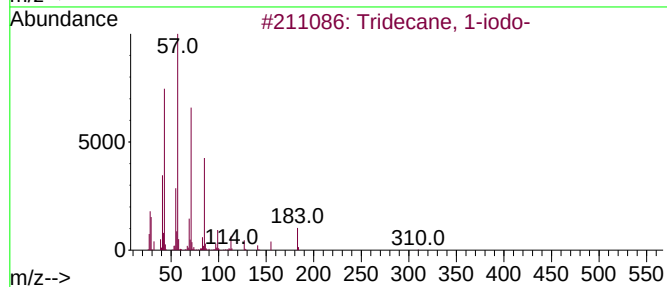
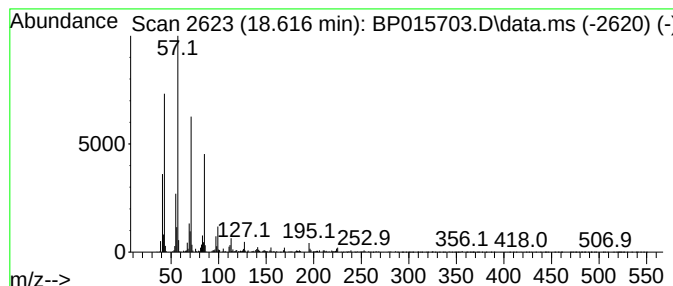
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Tridecane, 1-iodo- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.616	2.34 ng/ul	206953	Phenanthrene-d10	17.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	87
2			Heptacosane	380	C27H56	000593-49-7	87
3			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	86
4			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	86
5			Heneicosane	296	C21H44	000629-94-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
Acq On : 24 Jun 2023 11:46
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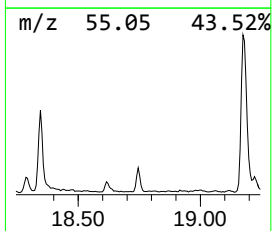
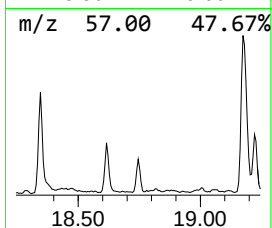
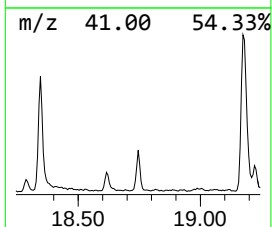
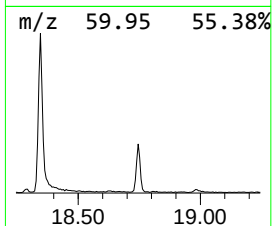
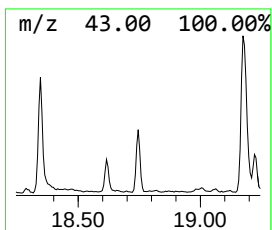
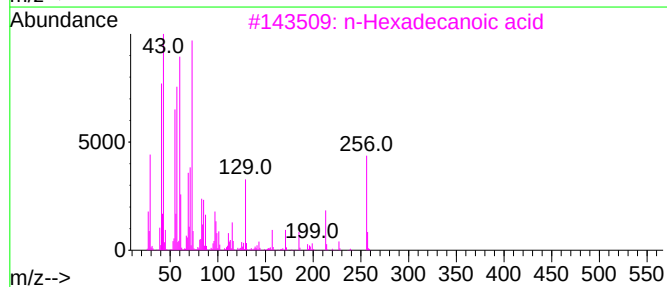
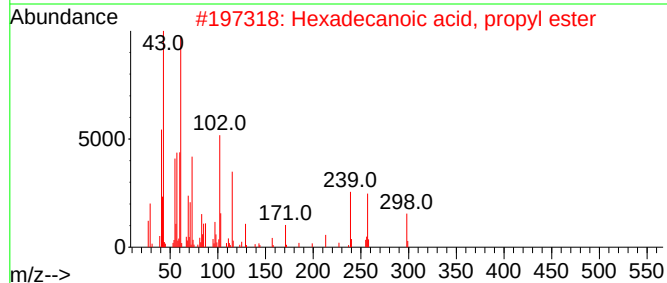
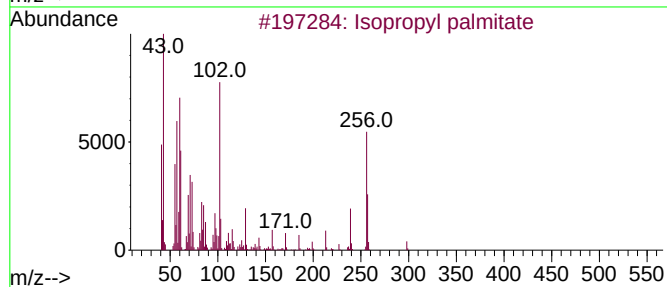
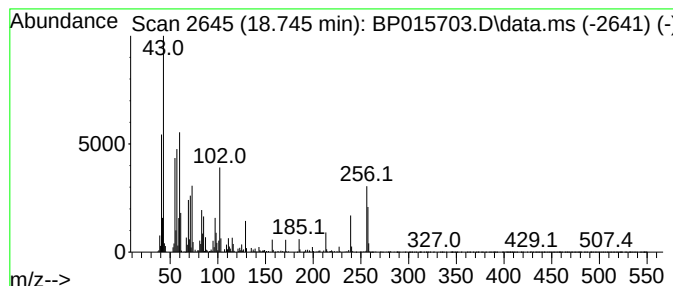
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Isopropyl palmitate Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.745	4.22 ng/ul	373295	Phenanthrene-d10	17.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl palmitate	298	C19H38O2	000142-91-6	64
2			Hexadecanoic acid, propyl ester	298	C19H38O2	002239-78-3	59
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	52
4			Isopropyl palmitate	298	C19H38O2	000142-91-6	50
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
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Acq On : 24 Jun 2023 11:46
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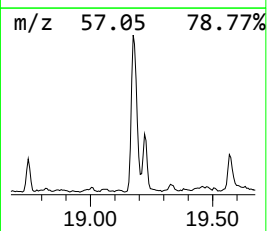
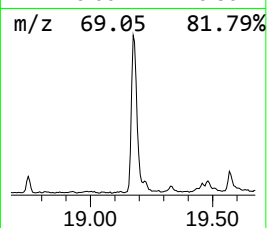
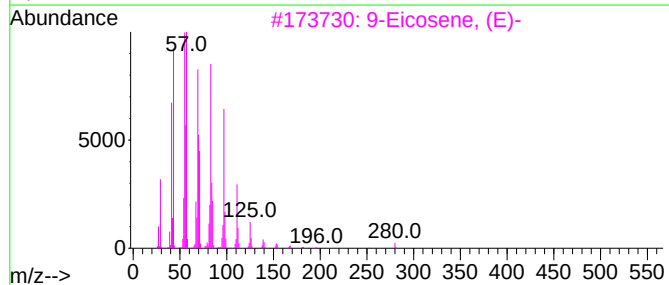
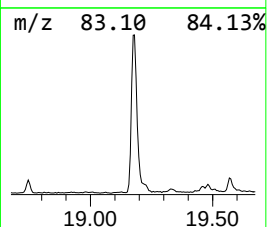
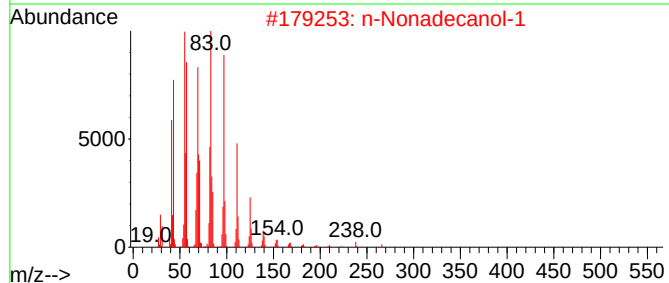
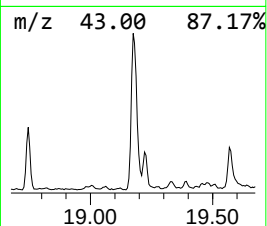
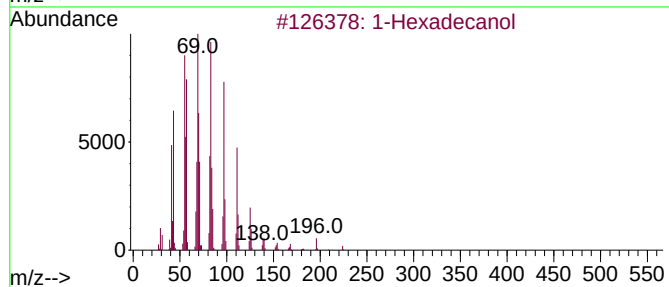
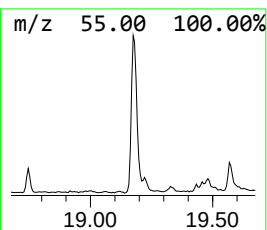
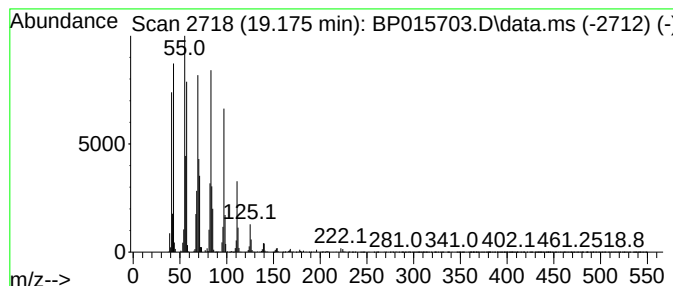
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1-Hexadecanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.174	23.02 ng/ul	2035560	Phenanthrene-d10	17.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexadecanol	242	C16H34O	036653-82-4	94
2			n-Nonadecanol-1	284	C19H40O	001454-84-8	94
3			9-Eicosene, (E)-	280	C20H40	074685-29-3	94
4			Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	94
5			1-Hexadecanol	242	C16H34O	036653-82-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
Acq On : 24 Jun 2023 11:46
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Sample : 03085-01
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ALS Vial : 41 Sample Multiplier: 1

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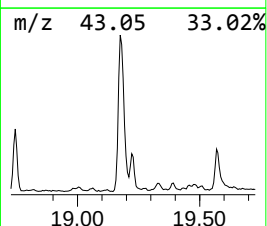
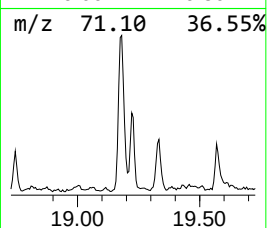
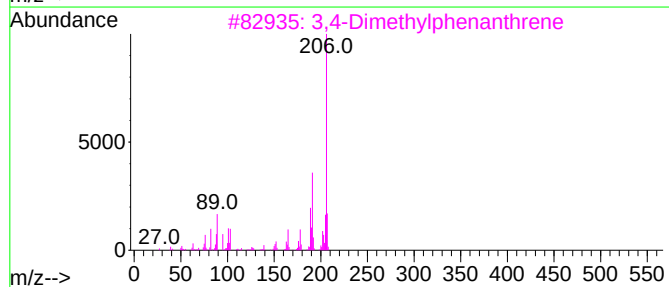
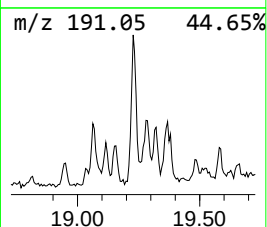
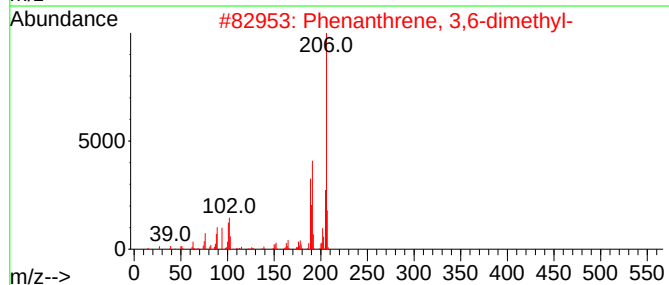
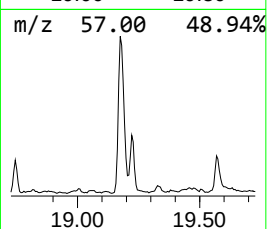
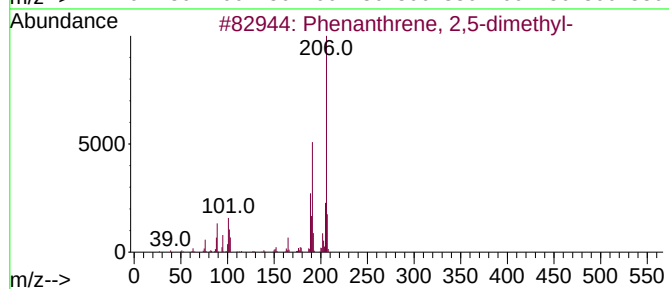
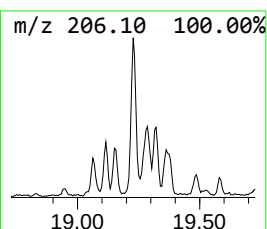
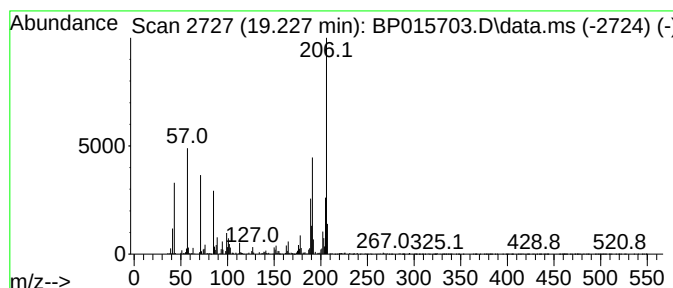
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Phenanthrene, 2,5-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.227	4.74 ng/ul	419426	Phenanthrene-d10	17.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	89
2			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	86
3			3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	86
4			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	83
5			Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
Acq On : 24 Jun 2023 11:46
Operator : MA/JU
Sample : 03085-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN7

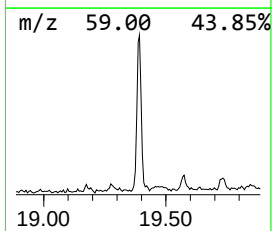
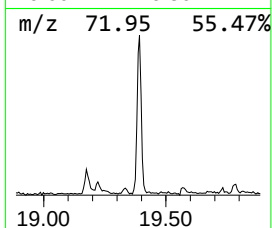
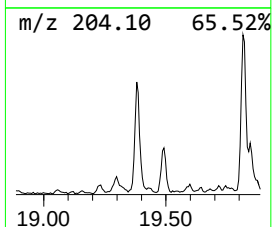
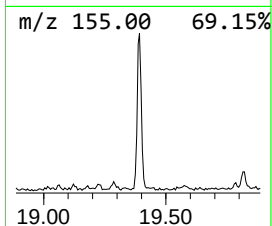
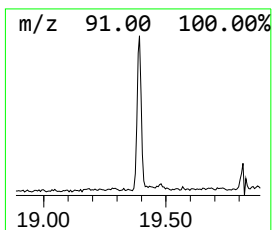
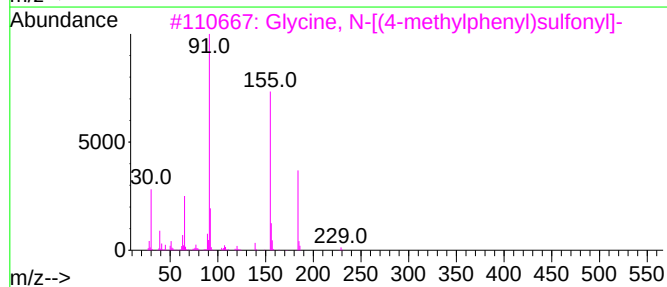
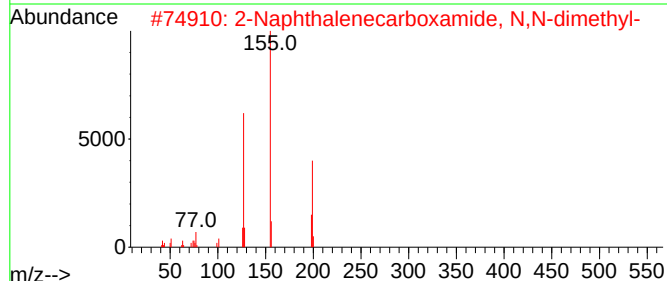
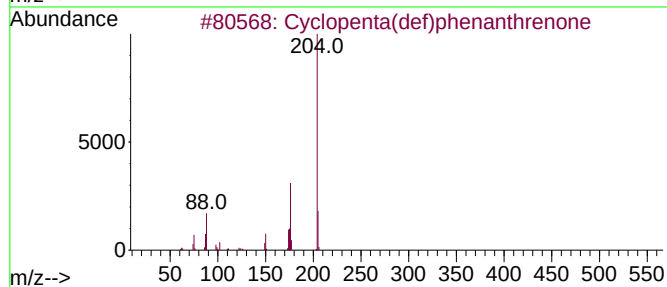
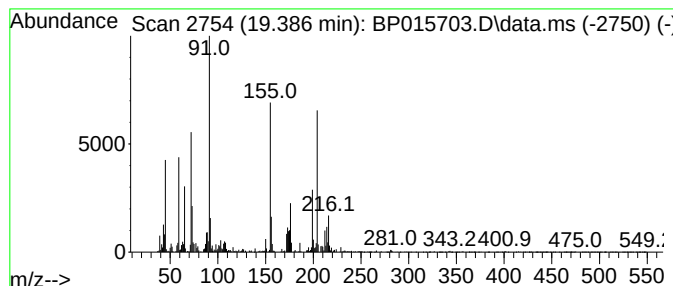
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-01 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.386	4.47 ng/ul	395494	Phenanthrene-d10	17.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	35
2			2-Naphthalenecarboxamide, N,N-di...	199	C13H13NO	013577-85-0	18
3			Glycine, N-[(4-methylphenyl)sulf...	229	C9H11NO4S	001080-44-0	15
4			2-(Toluene-4-sulfonyloxy)propion...	244	C10H12O5S	1000210-27-2	11
5			Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
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Operator : MA/JU
Sample : 03085-01
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ALS Vial : 41 Sample Multiplier: 1

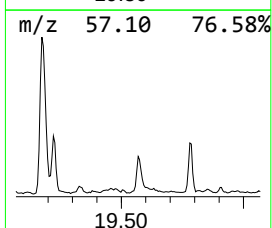
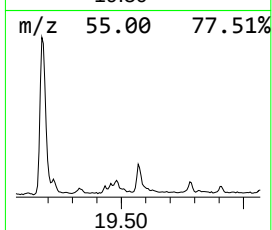
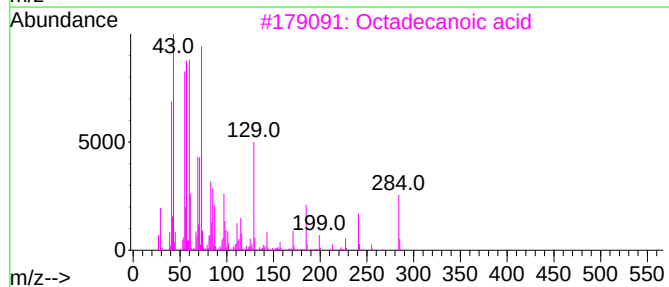
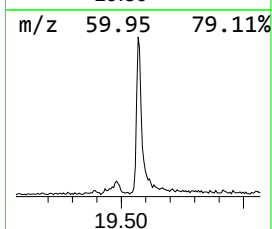
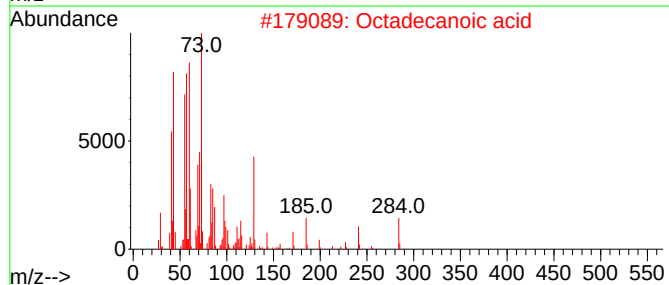
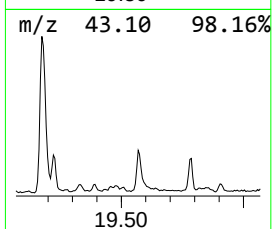
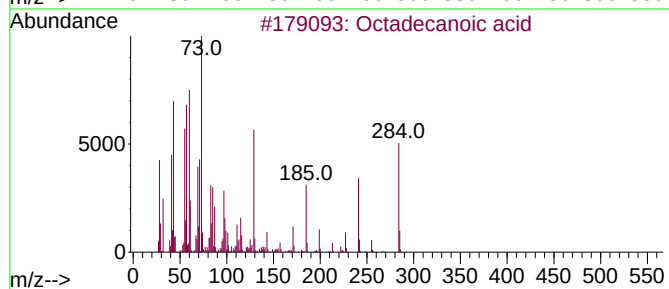
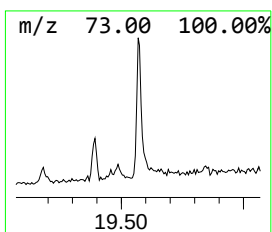
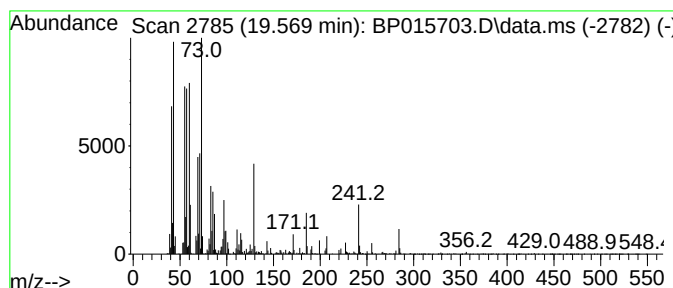
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 Octadecanoic acid Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.569	2.55 ng/ul	502225	Chrysene-d12	21.580	
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	99
3	Octadecanoic acid	284	C18H36O2	000057-11-4	99
4	Octadecanoic acid	284	C18H36O2	000057-11-4	96
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
Acq On : 24 Jun 2023 11:46
Operator : MA/JU
Sample : 03085-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
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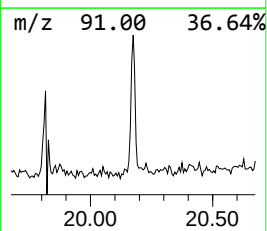
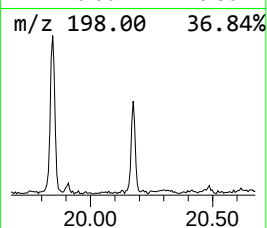
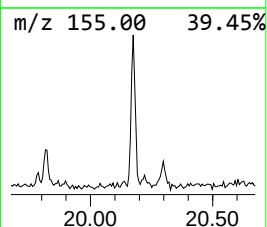
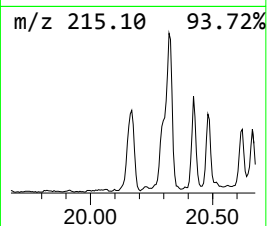
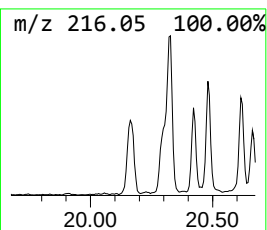
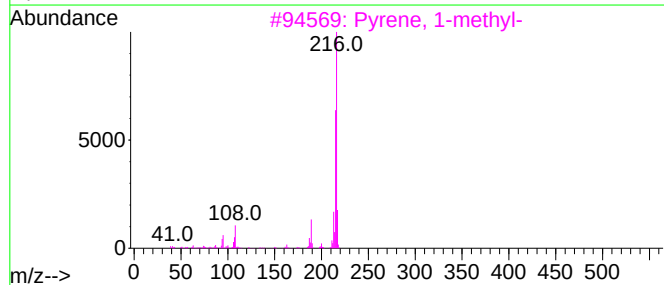
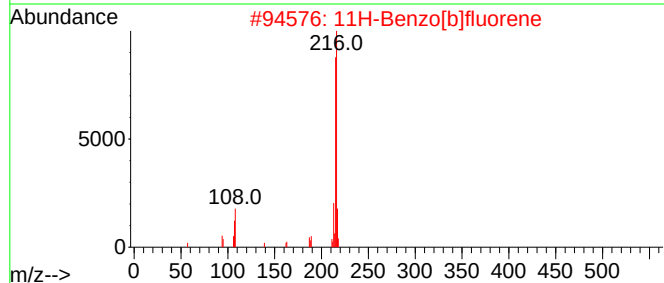
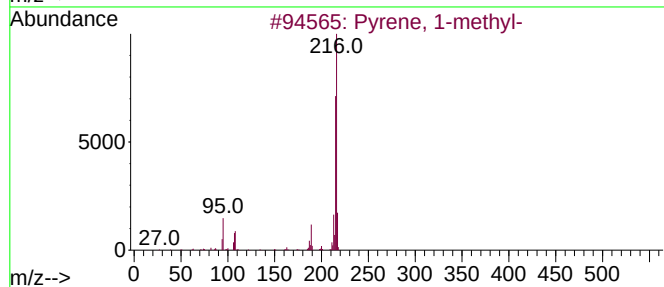
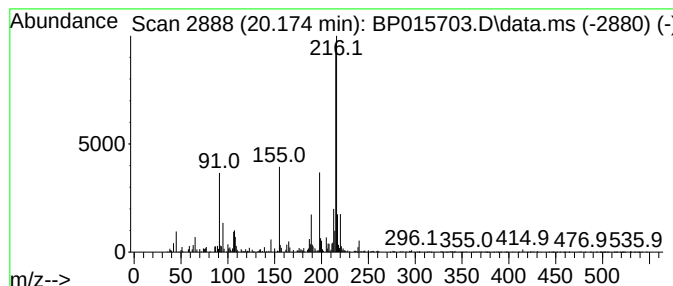
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Pyrene, 1-methyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.174	2.06 ng/ul	407042	Chrysene-d12	21.580

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pyrene, 1-methyl-	216	C17H12	002381-21-7	86
2		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	83
3		Pyrene, 1-methyl-	216	C17H12	002381-21-7	83
4		Pyrene, 2-methyl-	216	C17H12	003442-78-2	70
5		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
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Acq On : 24 Jun 2023 11:46
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Sample : 03085-01
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ALS Vial : 41 Sample Multiplier: 1

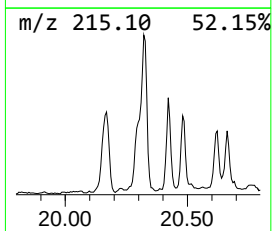
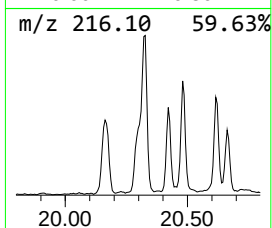
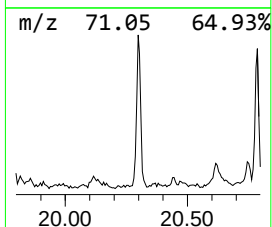
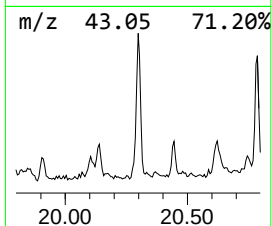
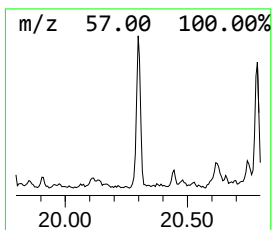
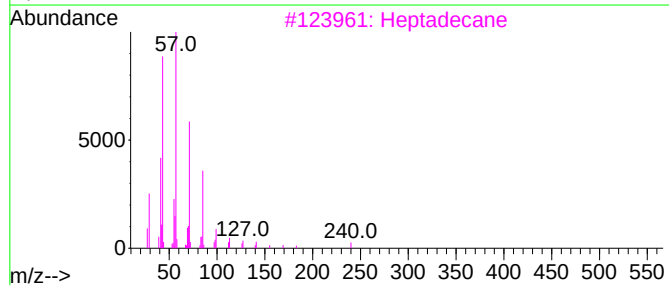
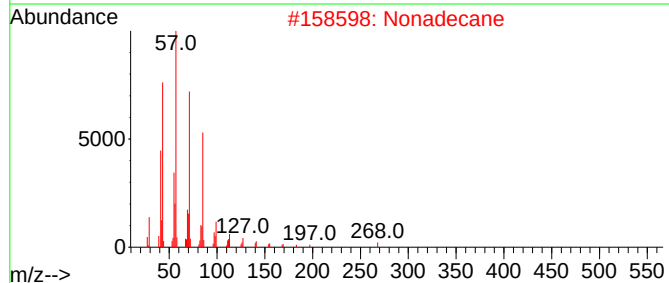
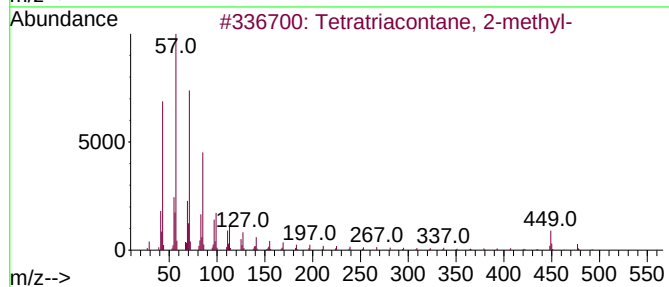
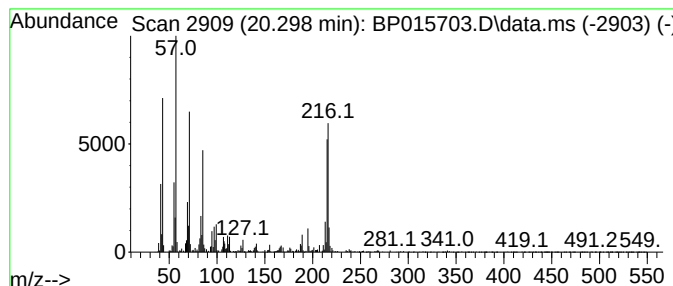
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.298	3.17 ng/ul	625487	Chrysene-d12		21.580	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tetratriacontane, 2-methyl-		493	C35H72	014167-65-8	93
2	Nonadecane		268	C19H40	000629-92-5	86
3	Heptadecane		240	C17H36	000629-78-7	83
4	Pyrene, 2-methyl-		216	C17H12	003442-78-2	60
5	Tetradecane, 1-iodo-		324	C14H29I	019218-94-1	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
Acq On : 24 Jun 2023 11:46
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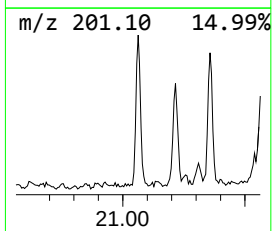
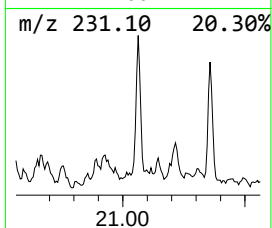
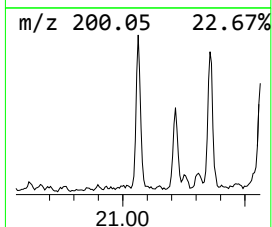
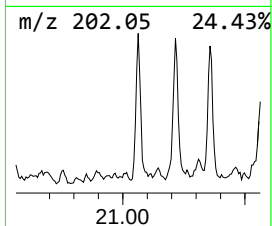
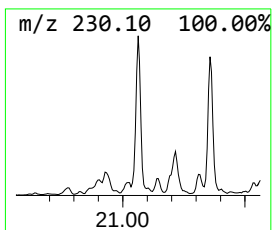
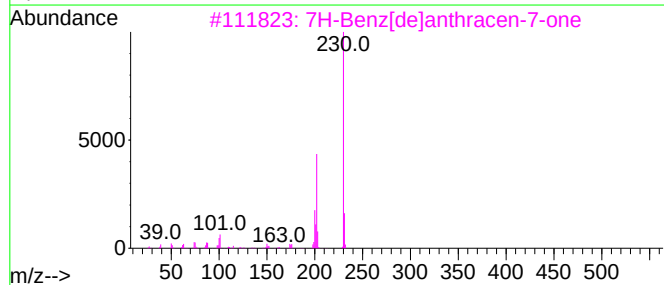
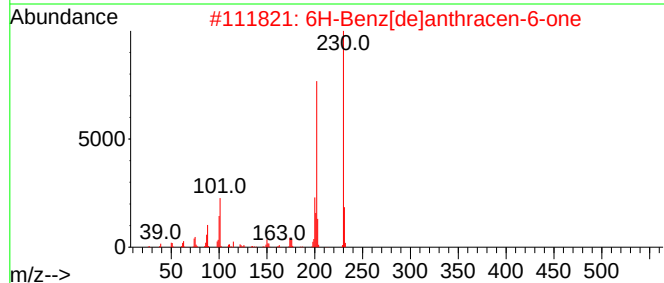
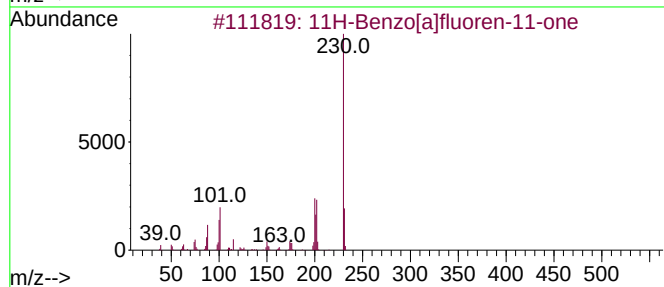
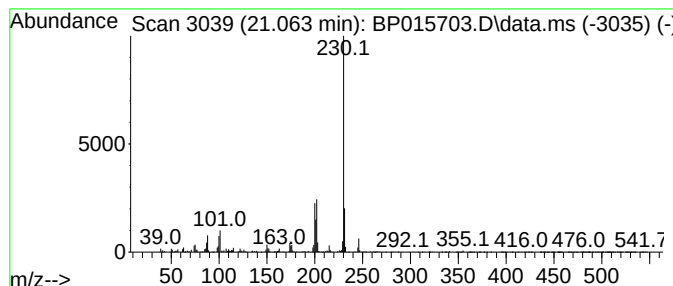
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 11H-Benzo[a]fluoren-11-one Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.063	3.13 ng/ul	618035	Chrysene-d12	21.580

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	97
2			6H-Benz[de]anthracen-6-one	230	C17H10O	080252-14-8	93
3			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	90
4			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	87
5			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
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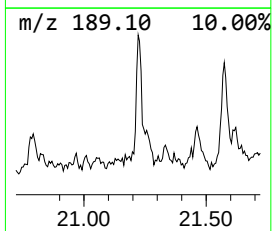
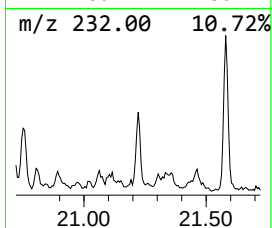
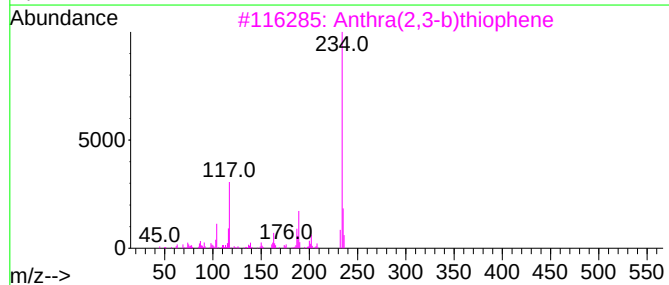
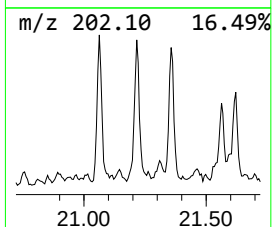
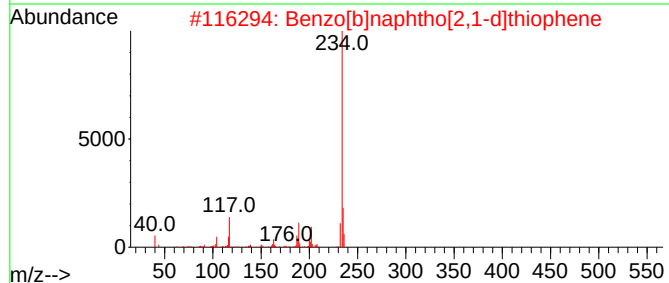
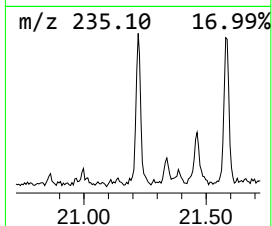
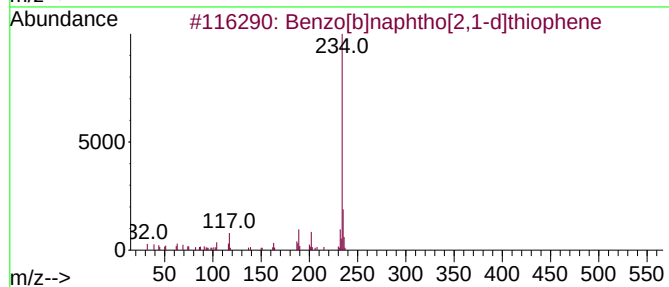
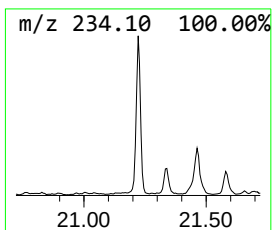
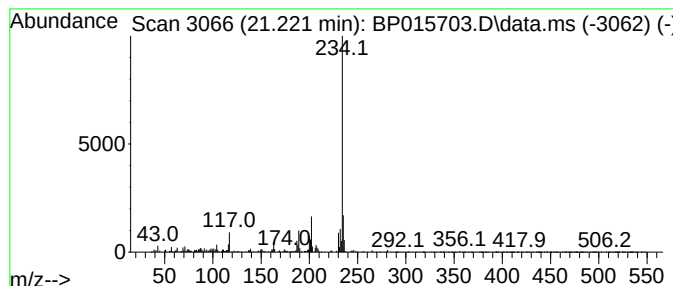
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.221	2.16 ng/ul	426193	Chrysene-d12	21.580

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	95
2			Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	94
3			Anthra(2,3-b)thiophene	234	C16H10S	022108-55-0	93
4			Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	90
5			Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
Acq On : 24 Jun 2023 11:46
Operator : MA/JU
Sample : 03085-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

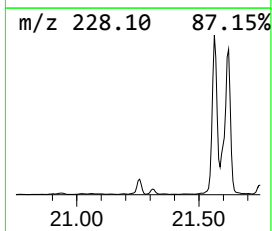
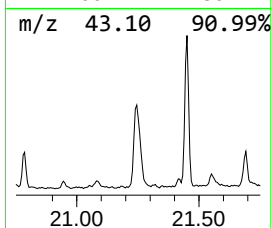
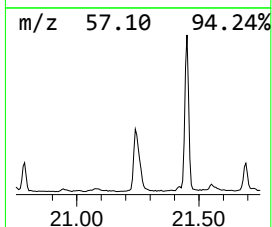
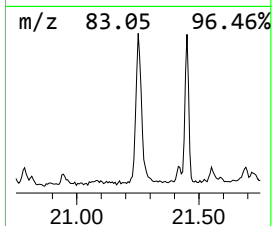
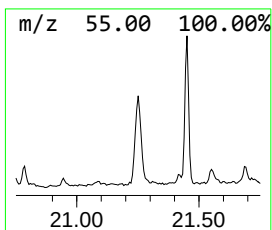
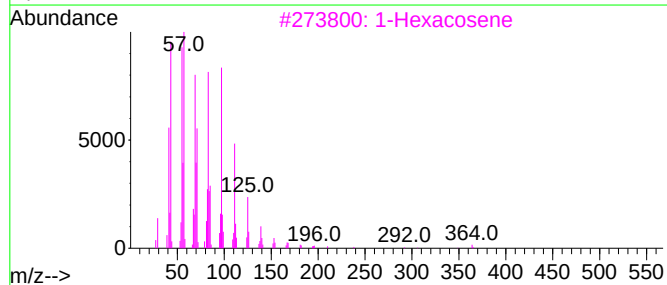
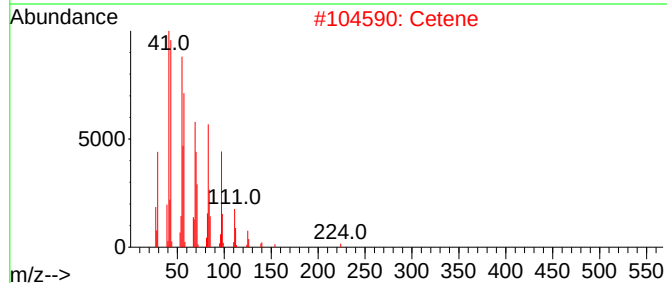
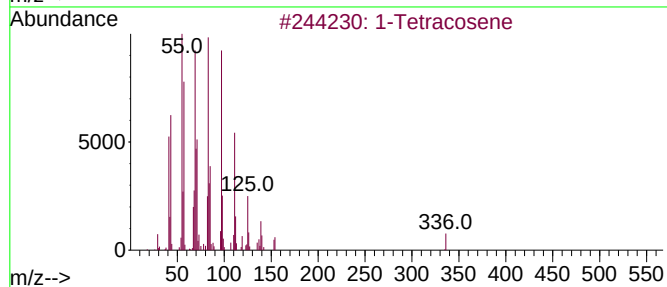
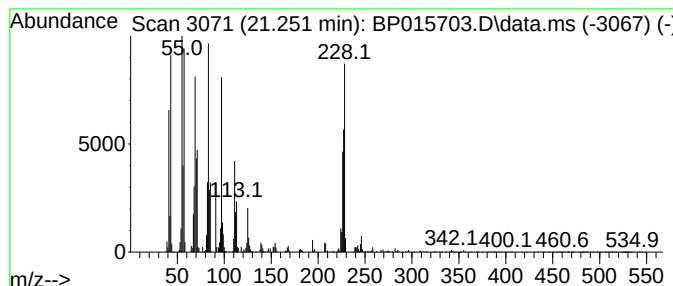
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.251	6.44 ng/ul	1269310	Chrysene-d12	21.580	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene	336	C24H48	010192-32-2	90
2	Cetene	224	C16H32	000629-73-2	78
3	1-Hexacosene	364	C26H52	018835-33-1	78
4	Cyclopentadecane	210	C15H30	000295-48-7	60
5	Dodecane, 1-cyclopentyl-4-(3-cyc...	348	C25H48	007225-68-5	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
Acq On : 24 Jun 2023 11:46
Operator : MA/JU
Sample : 03085-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

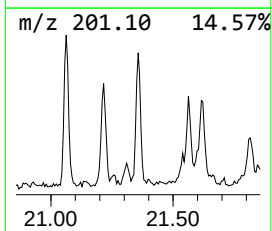
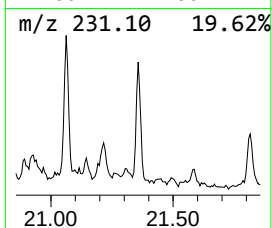
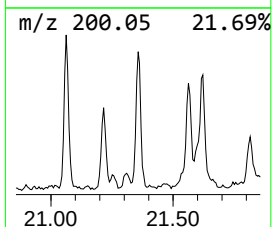
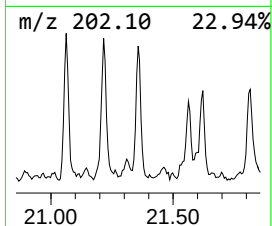
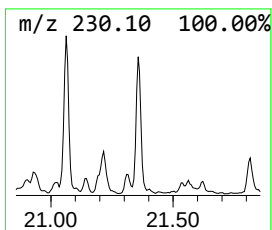
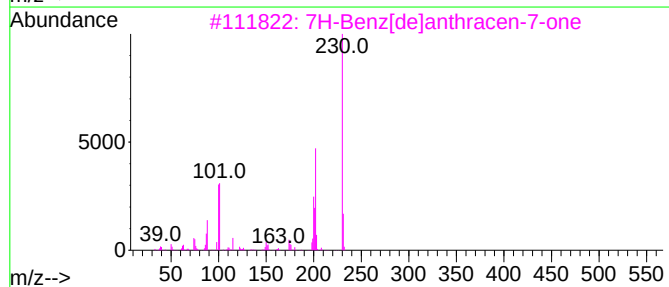
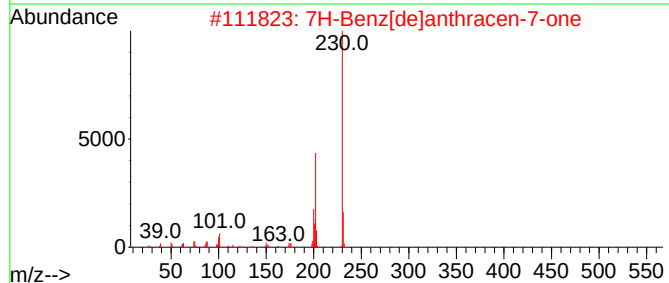
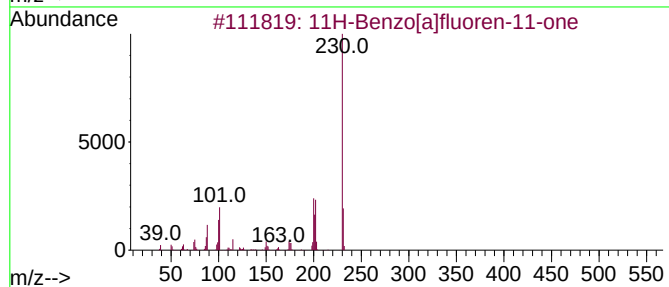
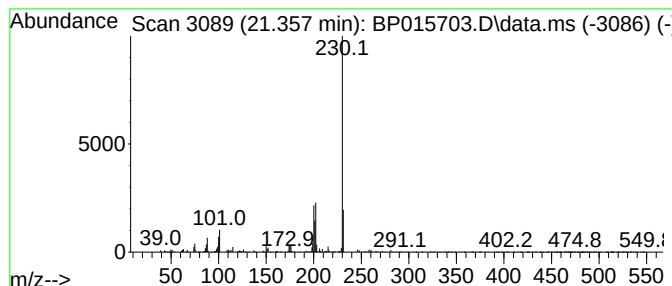
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 18 7H-Benz[de]anthracen-7-one Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.357	2.14 ng/ul	422223	Chrysene-d12		21.580	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluoren-11-one		230	C17H10O	000479-79-8	96
2	7H-Benz[de]anthracen-7-one		230	C17H10O	000082-05-3	90
3	7H-Benz[de]anthracen-7-one		230	C17H10O	000082-05-3	87
4	1-Pyrenecarboxaldehyde		230	C17H10O	003029-19-4	72
5	7H-Benz[de]anthracen-7-one		230	C17H10O	000082-05-3	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
Acq On : 24 Jun 2023 11:46
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Sample : 03085-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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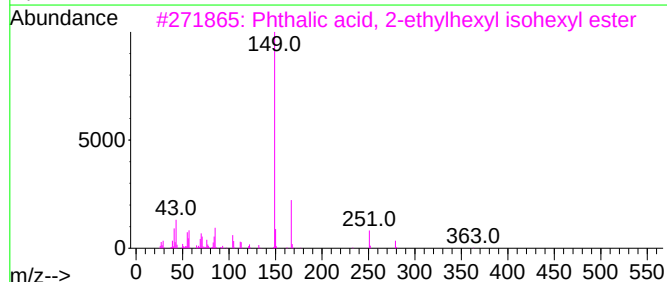
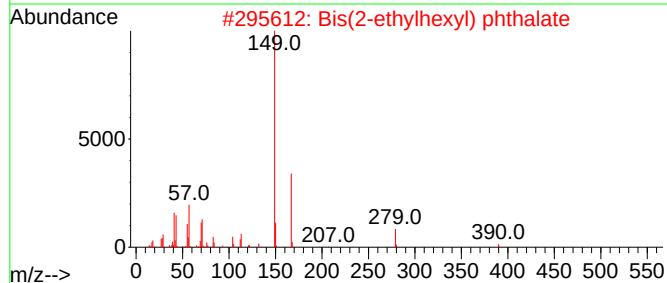
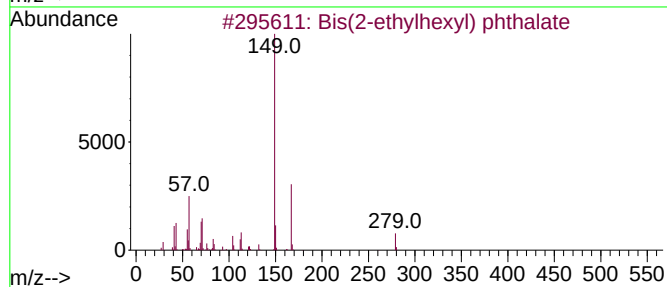
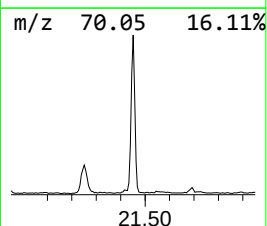
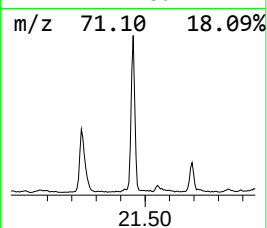
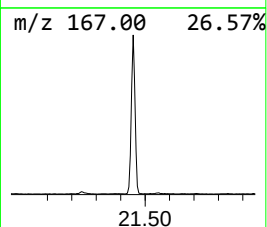
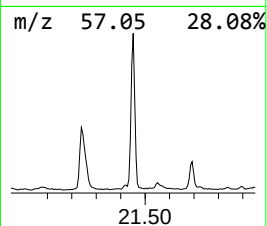
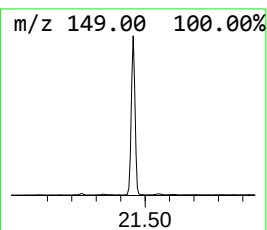
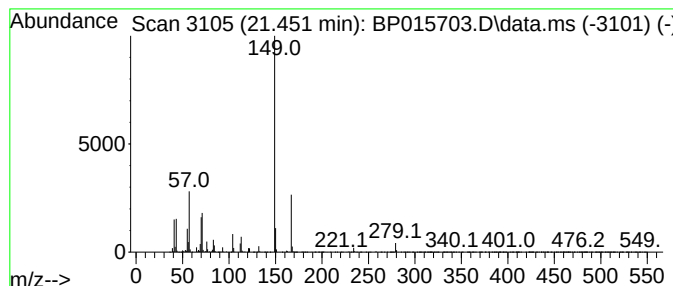
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 Bis(2-ethylhexyl) phthalate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.451	13.39 ng/ul	2639380	Chrysene-d12	21.580

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	91
2			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	91
3			Phthalic acid, 2-ethylhexyl isoh...	362	C22H34O4	1000308-98-5	86
4			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	86
5			Phthalic acid, isohexyl 2-methyl...	334	C20H30O4	1000356-91-0	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
Acq On : 24 Jun 2023 11:46
Operator : MA/JU
Sample : 03085-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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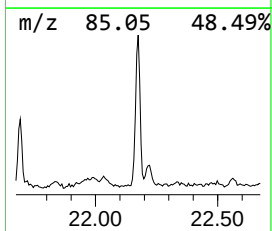
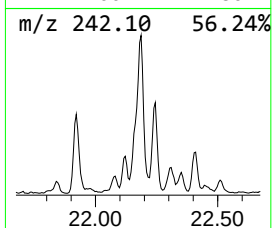
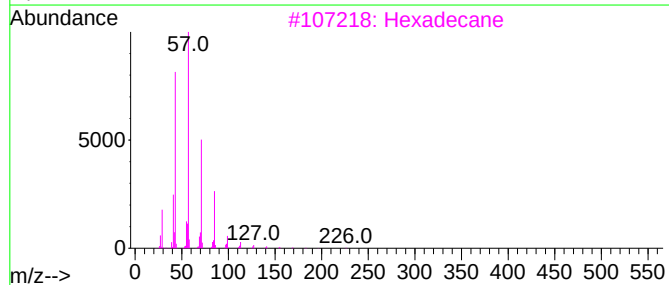
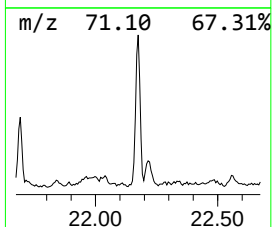
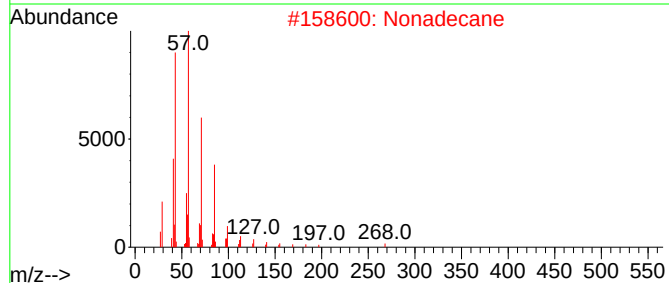
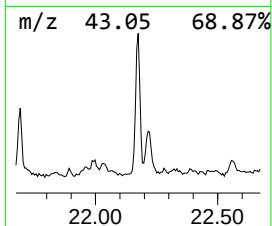
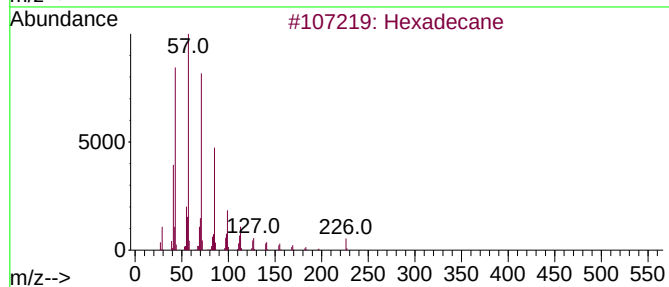
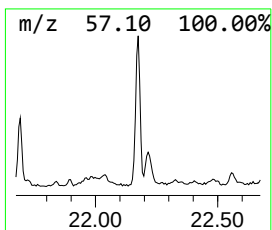
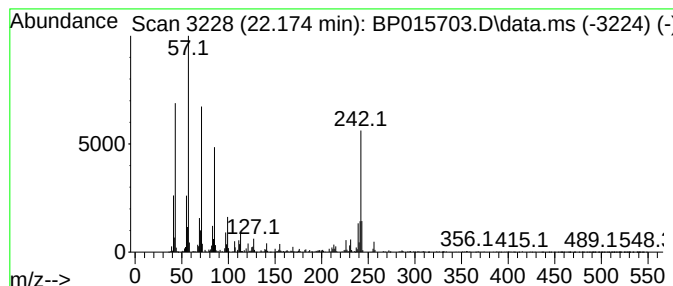
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.174	3.95 ng/ul	778769	Chrysene-d12	21.580

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	78
2			Nonadecane	268	C19H40	000629-92-5	78
3			Hexadecane	226	C16H34	000544-76-3	70
4			Heptadecane	240	C17H36	000629-78-7	70
5			Heptadecane	240	C17H36	000629-78-7	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
Acq On : 24 Jun 2023 11:46
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Sample : 03085-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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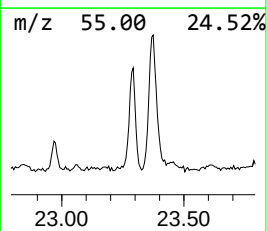
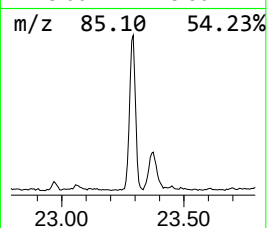
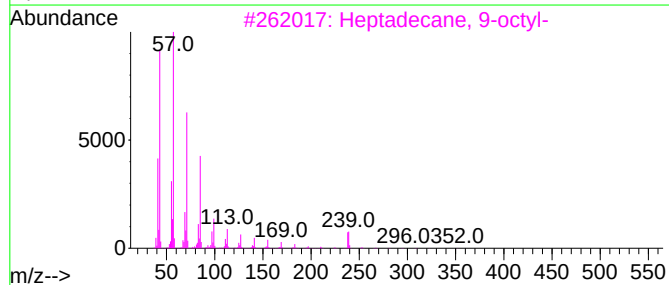
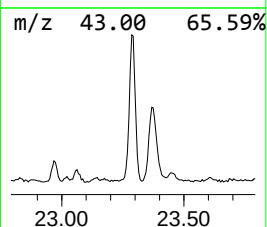
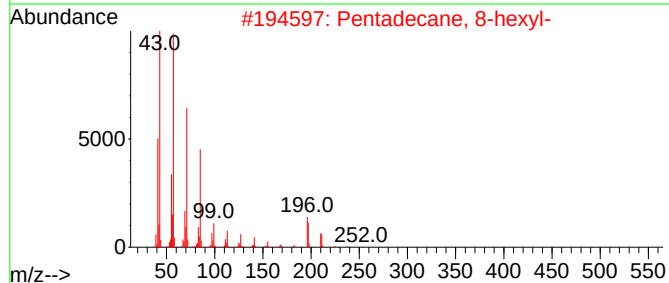
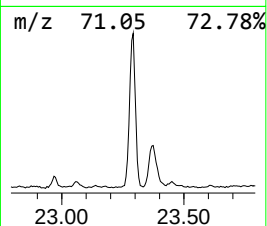
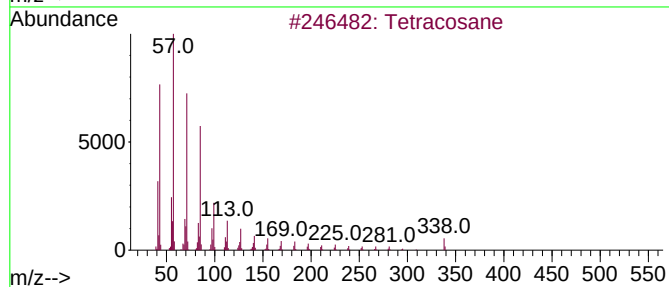
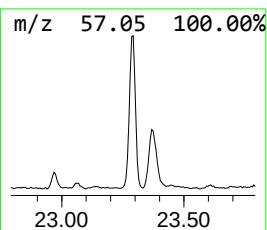
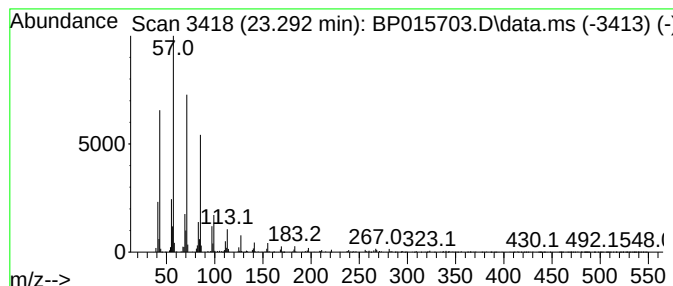
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.292	10.43 ng/ul	1054340	Perylene-d12	24.139

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	94
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
Acq On : 24 Jun 2023 11:46
Operator : MA/JU
Sample : 03085-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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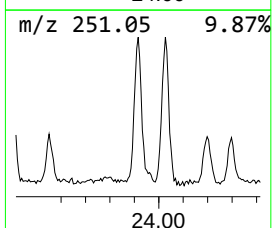
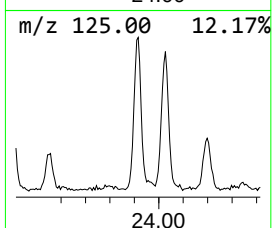
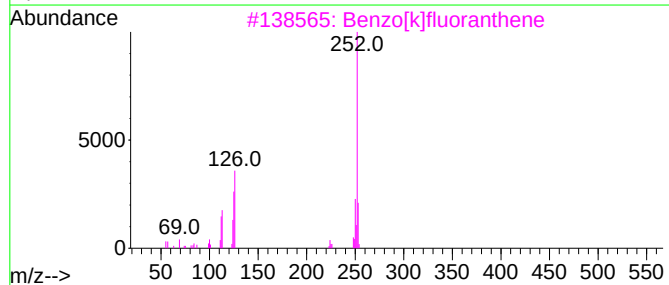
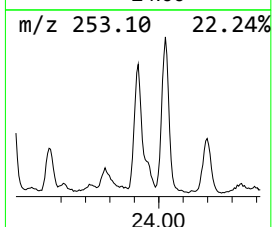
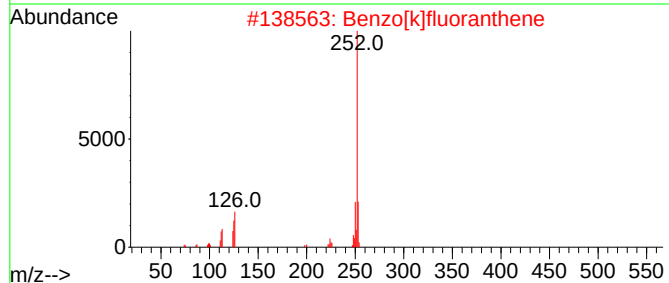
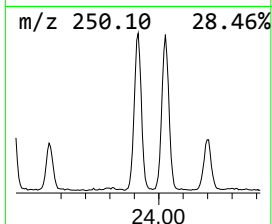
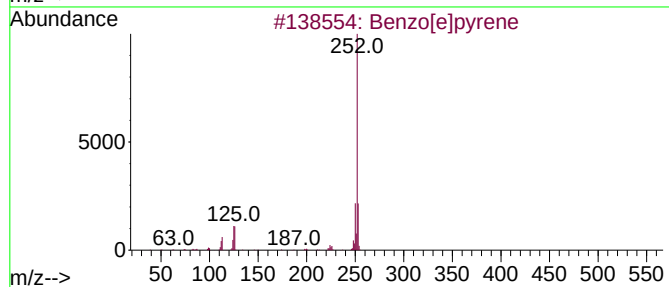
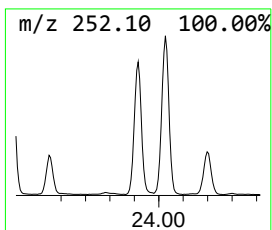
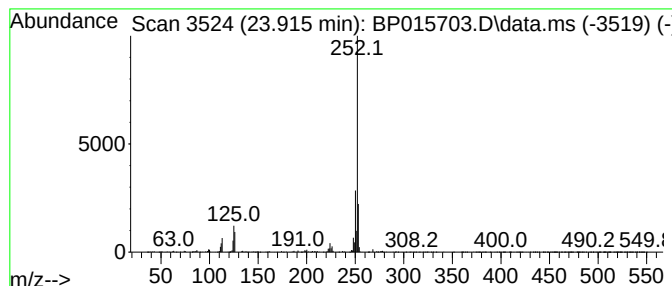
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 Benzo[e]pyrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.916	11.48 ng/ul	1160310	Perylene-d12	24.139

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	99
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
4			Benzo[b]fluoranthene	252	C20H12	000205-99-2	94
5			Benzo[a]pyrene	252	C20H12	000050-32-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
Acq On : 24 Jun 2023 11:46
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Sample : 03085-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
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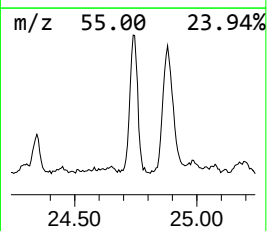
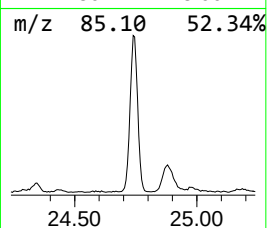
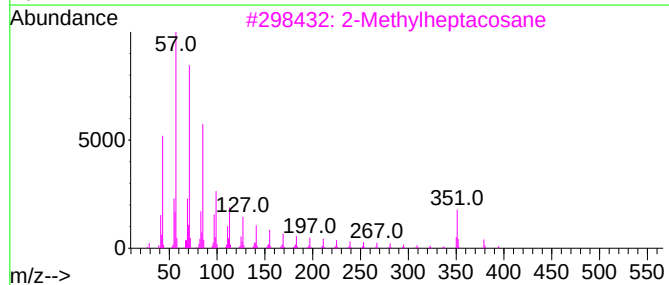
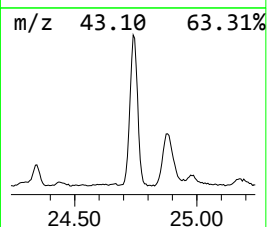
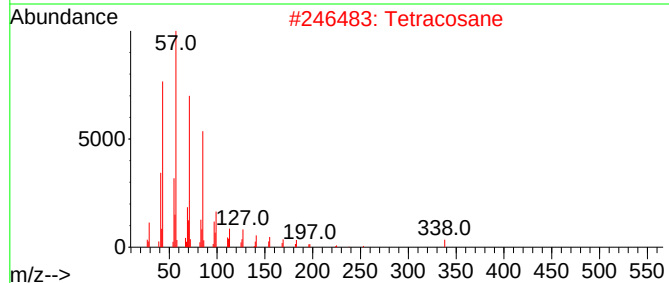
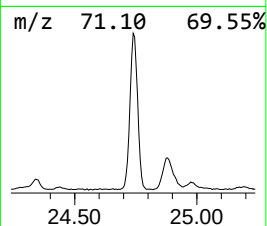
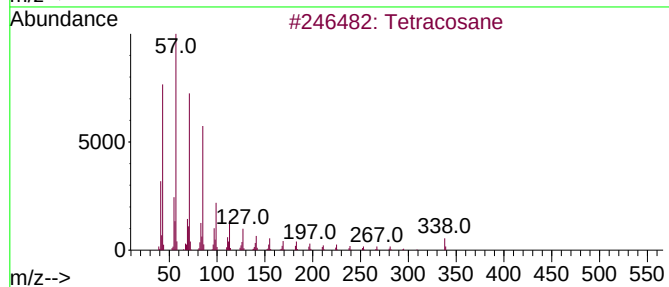
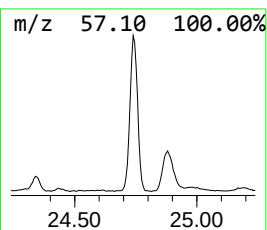
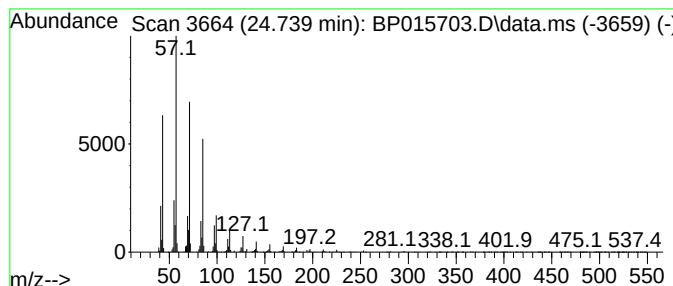
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.739	25.08 ng/ul	2534980	Perylene-d12	24.139

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	96
2			Tetracosane	338	C24H50	000646-31-1	94
3			2-Methylheptacosane	394	C28H58	001561-00-8	94
4			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
Acq On : 24 Jun 2023 11:46
Operator : MA/JU
Sample : 03085-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN7

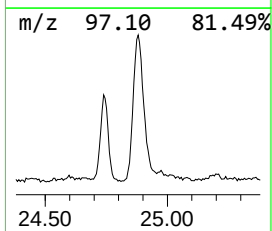
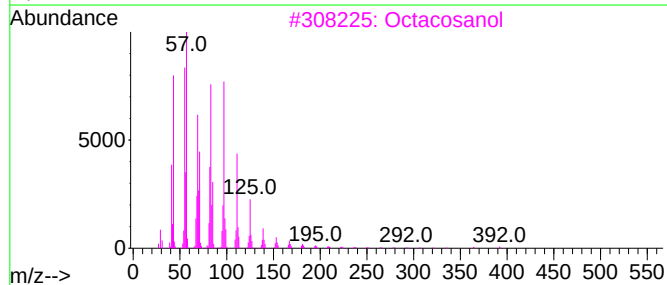
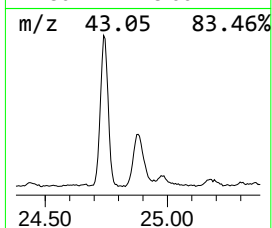
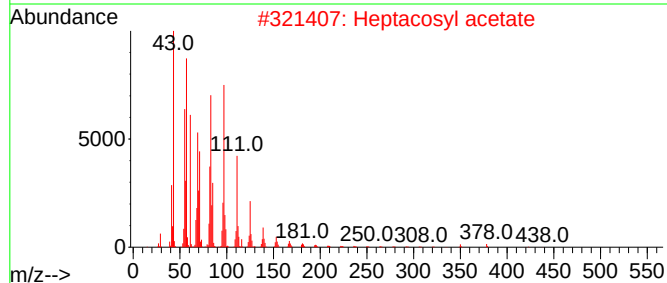
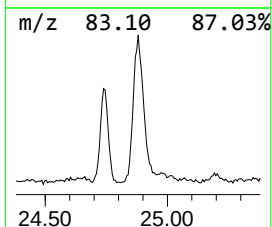
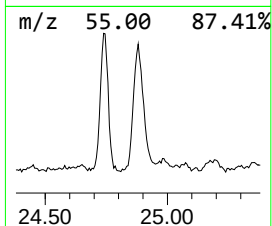
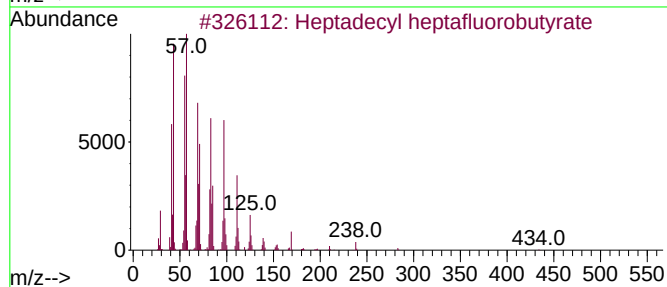
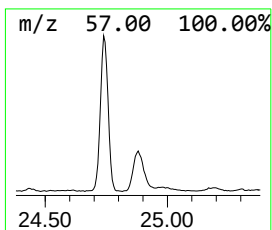
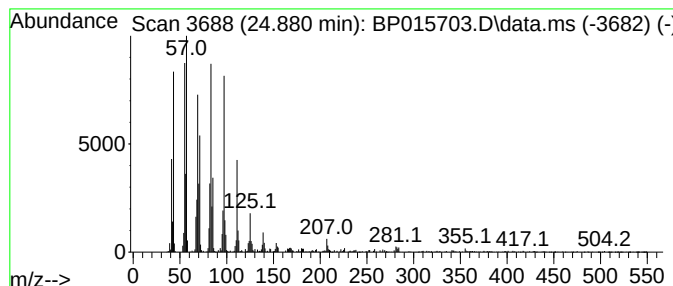
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 Heptadecyl heptafluorobutyrate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.880	14.39 ng/ul	1453830	Perylene-d12	24.139

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
2			Heptacosyl acetate	438	C29H58O2	1000351-78-2	94
3			Octacosanol	410	C28H58O	000557-61-9	93
4			1-Hexacosene	364	C26H52	018835-33-1	93
5			1-Tetracosene	336	C24H48	010192-32-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
Acq On : 24 Jun 2023 11:46
Operator : MA/JU
Sample : 03085-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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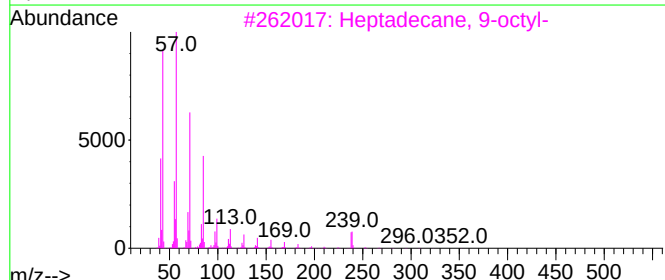
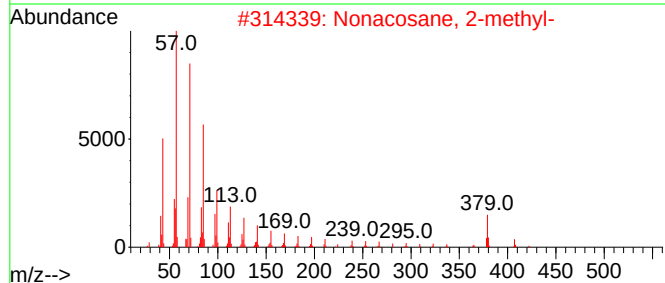
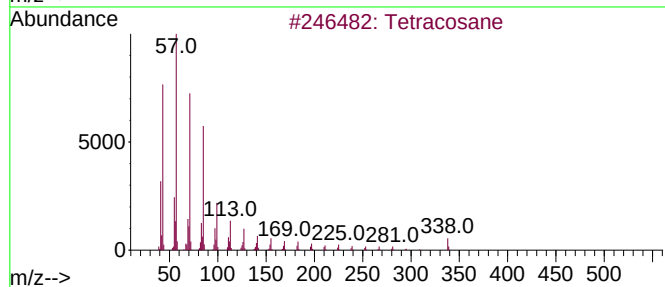
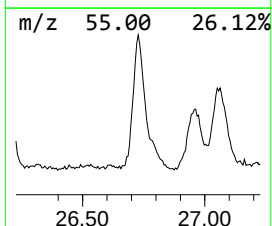
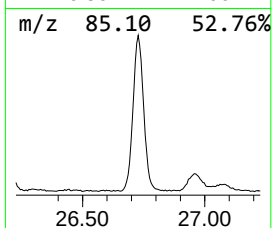
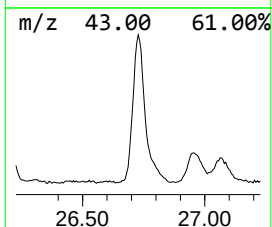
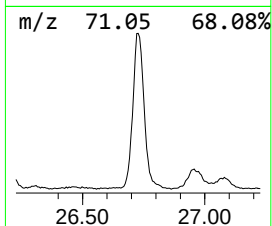
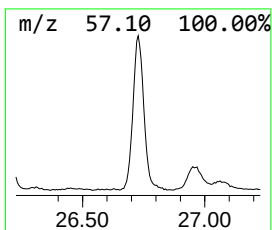
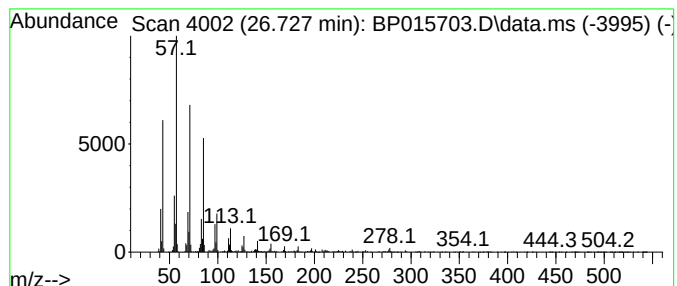
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.727	25.24 ng/ul	2550620	Perylene-d12	24.139

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	97
2			Nonacosane, 2-methyl-	422	C30H62	001560-75-4	94
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
4			Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	93
5			Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015703.D
Acq On : 24 Jun 2023 11:46
Operator : MA/JU
Sample : 03085-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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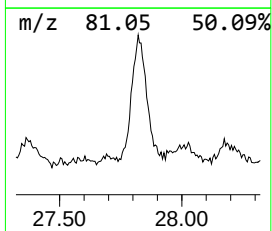
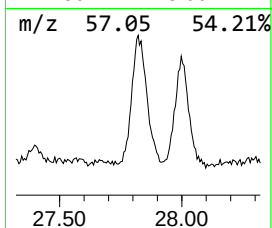
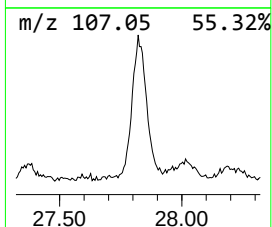
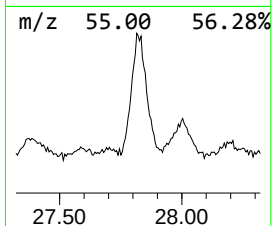
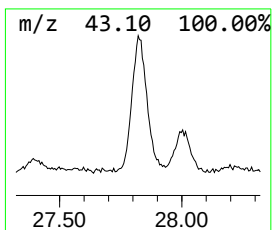
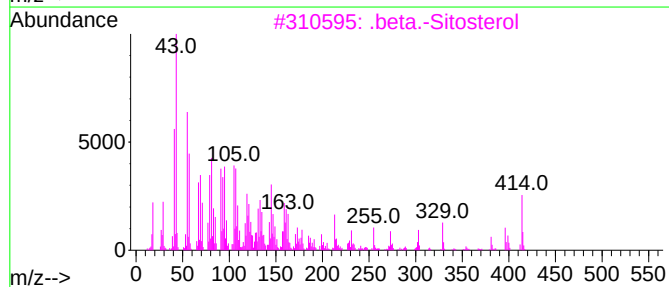
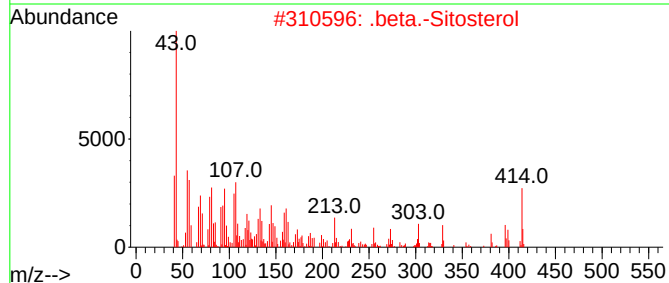
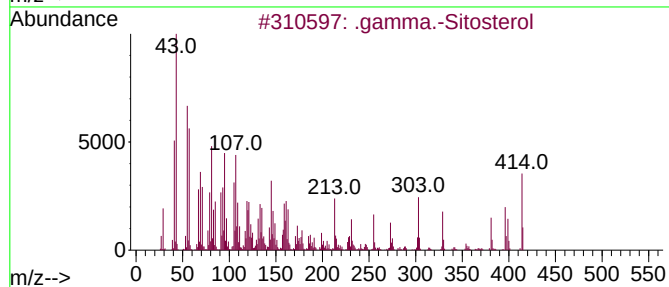
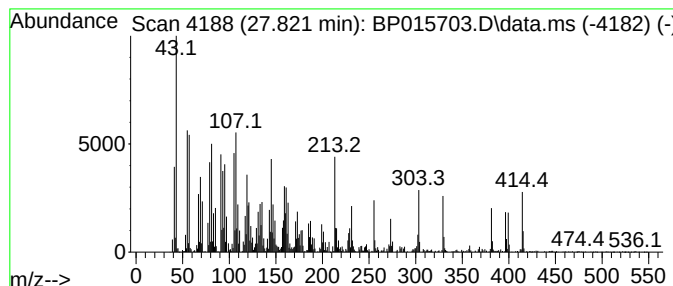
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 .gamma.-Sitosterol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.821	31.01 ng/ul	3133520	Perylene-d12	24.139

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	gamma.-Sitosterol	414	C29H50O	000083-47-6	99	
2	.	beta.-Sitosterol	414	C29H50O	000083-46-5	95	
3	.	beta.-Sitosterol	414	C29H50O	000083-46-5	91	
4	.	gamma.-Sitosterol	414	C29H50O	000083-47-6	46	
5	(3S,8S,9S,10R,13R,14S,17R)-17-((...	414	C29H50O	029365-29-5	45		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
 Data File : BP015703.D
 Acq On : 24 Jun 2023 11:46
 Operator : MA/JU
 Sample : 03085-01
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFN7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.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
2H-Pyran, tetra...	3.946	3.1	ng/ul	154282	1	8.075	985572	20.0
Benzophenone	16.086	4.3	ng/ul	335912	3	14.728	1566410	20.0
(DEL) Alkane: S...	16.433	4.8	ng/ul	422932	4	17.486	1768550	20.0
n-Hexadecanoic ...	18.345	15.3	ng/ul	1348400	4	17.486	1768550	20.0
Phenanthrene, 1...	18.398	3.6	ng/ul	318955	4	17.486	1768550	20.0
Anthracene, 1-m...	18.533	3.3	ng/ul	293044	4	17.486	1768550	20.0
Tridecane, 1-iodo-	18.616	2.3	ng/ul	206953	4	17.486	1768550	20.0
Isopropyl palmi...	18.745	4.2	ng/ul	373295	4	17.486	1768550	20.0
1-Hexadecanol	19.174	23.0	ng/ul	2035560	4	17.486	1768550	20.0
Phenanthrene, 2...	19.227	4.7	ng/ul	419426	4	17.486	1768550	20.0
unknown-01	19.386	4.5	ng/ul	395494	4	17.486	1768550	20.0
Octadecanoic acid	19.569	2.5	ng/ul	502225	5	21.580	3943280	20.0
Pyrene, 1-methyl-	20.174	2.1	ng/ul	407042	5	21.580	3943280	20.0
(DEL) Alkane: S...	20.298	3.2	ng/ul	625487	5	21.580	3943280	20.0
11H-Benzo[a]flu...	21.063	3.1	ng/ul	618035	5	21.580	3943280	20.0
Benzo[b]naphtho...	21.221	2.2	ng/ul	426193	5	21.580	3943280	20.0
(DEL) Alkane: S...	21.251	6.4	ng/ul	1269310	5	21.580	3943280	20.0
7H-Benz[de]anth...	21.357	2.1	ng/ul	422223	5	21.580	3943280	20.0
Bis(2-ethylhexy...	21.451	13.4	ng/ul	2639380	5	21.580	3943280	20.0
(DEL) Alkane: S...	22.174	4.0	ng/ul	778769	5	21.580	3943280	20.0
(DEL) Alkane: S...	23.292	10.4	ng/ul	1054340	6	24.139	2021140	20.0
Benzo[e]pyrene	23.916	11.5	ng/ul	1160310	6	24.139	2021140	20.0
(DEL) Alkane: S...	24.739	25.1	ng/ul	2534980	6	24.139	2021140	20.0
Heptadecyl hept...	24.880	14.4	ng/ul	1453830	6	24.139	2021140	20.0
(DEL) Alkane: S...	26.727	25.2	ng/ul	2550620	6	24.139	2021140	20.0
.gamma.-Sitosterol	27.821	31.0	ng/ul	3133520	6	24.139	2021140	20.0