

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
 Data File : BP012356.D
 Acq On : 27 Oct 2022 22:57
 Operator : CG/JU
 Sample : N5144-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BK1

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-BP102522.M
 Title : SVOA CALIBRATION

Signal : TIC: BP012356.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.246	41	44	52	rVB	63627	87728	0.88%	0.067%
2	3.693	118	120	129	rVV	40805	74621	0.75%	0.057%
3	3.770	129	133	141	rVV	92294	139256	1.40%	0.106%
4	3.887	148	153	162	rVV	78279	125116	1.26%	0.096%
5	4.317	223	226	231	rBV	34475	50220	0.51%	0.038%
6	4.546	261	265	273	rVB	55004	80594	0.81%	0.062%
7	4.922	321	329	336	rBV2	78718	183931	1.85%	0.141%
8	6.975	667	678	698	rBV	1163018	2056765	20.72%	1.571%
9	7.146	700	707	719	rBV	975854	1624370	16.36%	1.241%
10	7.334	733	739	749	rBV	1163987	1927723	19.42%	1.473%
11	7.805	807	819	828	rBV	1197789	1966841	19.81%	1.503%
12	8.346	904	911	921	rBV	29790	67249	0.68%	0.051%
13	8.522	934	941	956	rBV	1356303	2299318	23.16%	1.757%
14	8.640	956	961	969	rVB	46030	85777	0.86%	0.066%
15	8.975	1010	1018	1034	rBV	1158311	1974319	19.89%	1.508%
16	9.363	1079	1084	1090	rBV	37704	61915	0.62%	0.047%
17	9.693	1132	1140	1165	rBV	951953	1670520	16.83%	1.276%
18	10.234	1224	1232	1253	rBV	1556130	2723236	27.43%	2.080%
19	10.399	1255	1260	1275	rBV	77519	187754	1.89%	0.143%
20	10.610	1286	1296	1302	rBV	1762395	3125502	31.48%	2.388%
21	10.657	1302	1304	1311	rVB	75538	104917	1.06%	0.080%
22	10.757	1314	1321	1340	rBV	782687	1563888	15.75%	1.195%
23	12.199	1562	1566	1572	rVV	30547	50056	0.50%	0.038%
24	12.275	1574	1579	1586	rVB	69084	115391	1.16%	0.088%
25	12.493	1611	1616	1624	rVB	77583	129343	1.30%	0.099%
26	13.269	1742	1748	1756	rBV2	40595	84800	0.85%	0.065%
27	13.634	1800	1810	1818	rBV5	27735	78817	0.79%	0.060%
28	13.775	1828	1834	1839	rBV2	70743	128149	1.29%	0.098%
29	13.828	1839	1843	1845	rVV	44151	61459	0.62%	0.047%
30	13.869	1845	1850	1864	rVB	2794495	4107526	41.37%	3.138%
31	14.151	1884	1898	1915	rBV	3523033	5448280	54.88%	4.162%
32	14.457	1940	1950	1957	rBV	3032615	4441255	44.74%	3.393%
33	14.522	1957	1961	1968	rVB2	45706	82759	0.83%	0.063%
34	14.675	1981	1987	2008	rBV	1136145	2051248	20.66%	1.567%
35	14.857	2009	2018	2021	rBV3	85672	205099	2.07%	0.157%
36	15.075	2049	2055	2060	rBV4	28538	54752	0.55%	0.042%

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Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
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37	15.251	2078	2085	2087	rBV5	33855	68016	0.69%	0.052%
38	15.451	2113	2119	2125	rBV	4431016	6569241	66.17%	5.019%
39	15.510	2125	2129	2137	rVV3	103845	175302	1.77%	0.134%
40	15.587	2137	2142	2154	rVV	1093646	1524922	15.36%	1.165%
41	15.710	2158	2163	2172	rVB7	30627	85988	0.87%	0.066%
42	15.804	2174	2179	2185	rVB	45858	76737	0.77%	0.059%
43	15.951	2201	2204	2212	rVV	49381	91398	0.92%	0.070%
44	16.192	2237	2245	2250	rBV4	79142	166120	1.67%	0.127%
45	16.510	2291	2299	2305	rBV6	42490	111075	1.12%	0.085%
46	16.751	2335	2340	2344	rBV2	40012	74634	0.75%	0.057%
47	16.875	2356	2361	2368	rBV	169674	261672	2.64%	0.200%
48	16.969	2374	2377	2381	rVV3	43004	57388	0.58%	0.044%
49	17.034	2383	2388	2396	rVB	132379	219703	2.21%	0.168%
50	17.116	2397	2402	2409	rBV	109491	177707	1.79%	0.136%
51	17.181	2409	2413	2414	rBV	104980	114822	1.16%	0.088%
52	17.216	2414	2419	2423	rVV	3650291	5399132	54.38%	4.125%
53	17.263	2423	2427	2431	rVV	2433396	3380341	34.05%	2.582%
54	17.316	2431	2436	2447	rVB2	4871150	7392313	74.46%	5.647%
55	17.628	2482	2489	2496	rBV	242426	431559	4.35%	0.330%
56	17.739	2505	2508	2514	rVB2	77969	108017	1.09%	0.083%
57	17.798	2514	2518	2524	rBV2	85402	132612	1.34%	0.101%
58	17.886	2530	2533	2538	rVB2	62476	77169	0.78%	0.059%
59	17.992	2546	2551	2558	rBV3	83470	199114	2.01%	0.152%
60	18.051	2558	2561	2563	rBV2	66835	81385	0.82%	0.062%
61	18.110	2563	2571	2573	rVV2	1088443	2032395	20.47%	1.553%
62	18.134	2573	2575	2580	rVV	697972	870556	8.77%	0.665%
63	18.175	2580	2582	2585	rVV	125874	147943	1.49%	0.113%
64	18.210	2585	2588	2593	rVB	96296	131750	1.33%	0.101%
65	18.275	2593	2599	2602	rBV2	446210	786495	7.92%	0.601%
66	18.310	2602	2605	2610	rVB	330829	447545	4.51%	0.342%
67	18.392	2614	2619	2622	rBV4	127158	218409	2.20%	0.167%
68	18.428	2623	2625	2629	rVB4	47366	55129	0.56%	0.042%
69	18.575	2645	2650	2653	rBV	330319	453998	4.57%	0.347%
70	18.616	2653	2657	2666	rVV	333828	483248	4.87%	0.369%
71	18.810	2687	2690	2693	rVV2	88064	96736	0.97%	0.074%
72	18.857	2695	2698	2701	rVV	94311	108734	1.10%	0.083%
73	18.892	2701	2704	2708	rVB	77613	100615	1.01%	0.077%
74	18.975	2713	2718	2724	rBV2	287764	616143	6.21%	0.471%
75	19.063	2731	2733	2737	rVB	119375	121526	1.22%	0.093%
76	19.116	2737	2742	2753	rBV3	303810	698365	7.03%	0.534%

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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 >
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 Title : SVOA CALIBRATION

77	19.233	2753	2762	2769	rBV	4295821	5901373	59.44%	4.508%
78	19.339	2775	2780	2784	rBV2	651158	924287	9.31%	0.706%
79	19.469	2799	2802	2811	rVB	144936	233686	2.35%	0.179%
80	19.557	2811	2817	2820	rVV	6735925	9927870	100.00%	7.584%
81	19.586	2820	2822	2829	rVV	4074143	4713959	47.48%	3.601%
82	19.657	2830	2834	2839	rVB2	211501	305347	3.08%	0.233%
83	19.757	2848	2851	2856	rVB	152098	216491	2.18%	0.165%
84	19.898	2870	2875	2882	rBV4	290273	690513	6.96%	0.528%
85	20.033	2893	2898	2899	rBV3	238595	329970	3.32%	0.252%
86	20.222	2926	2930	2934	rVB2	397614	549146	5.53%	0.420%
87	20.357	2950	2953	2957	rBV	289317	356076	3.59%	0.272%
88	20.404	2958	2961	2965	rVB	261504	342692	3.45%	0.262%
89	20.557	2985	2987	2992	rVB3	156107	160444	1.62%	0.123%
90	20.804	3025	3029	3035	rBV	458184	625289	6.30%	0.478%
91	21.022	3059	3066	3074	rBV5	835079	2325009	23.42%	1.776%
92	21.092	3075	3078	3083	rVB	439747	588484	5.93%	0.450%
93	21.216	3095	3099	3104	rVB	876002	924719	9.31%	0.706%
94	21.310	3108	3115	3119	rVV2	5436594	9326555	93.94%	7.125%
95	21.345	3119	3121	3131	rVB	2070428	2582929	26.02%	1.973%
96	22.974	3392	3398	3403	rBV2	1149582	2806801	28.27%	2.144%
97	23.498	3482	3487	3491	rBV	827127	1581588	15.93%	1.208%
98	23.551	3491	3496	3501	rBV2	3294985	6224498	62.70%	4.755%
99	23.710	3517	3523	3529	rBV2	2444880	4530962	45.64%	3.461%
100	24.298	3619	3623	3632	rVB2	567021	1162877	11.71%	0.888%

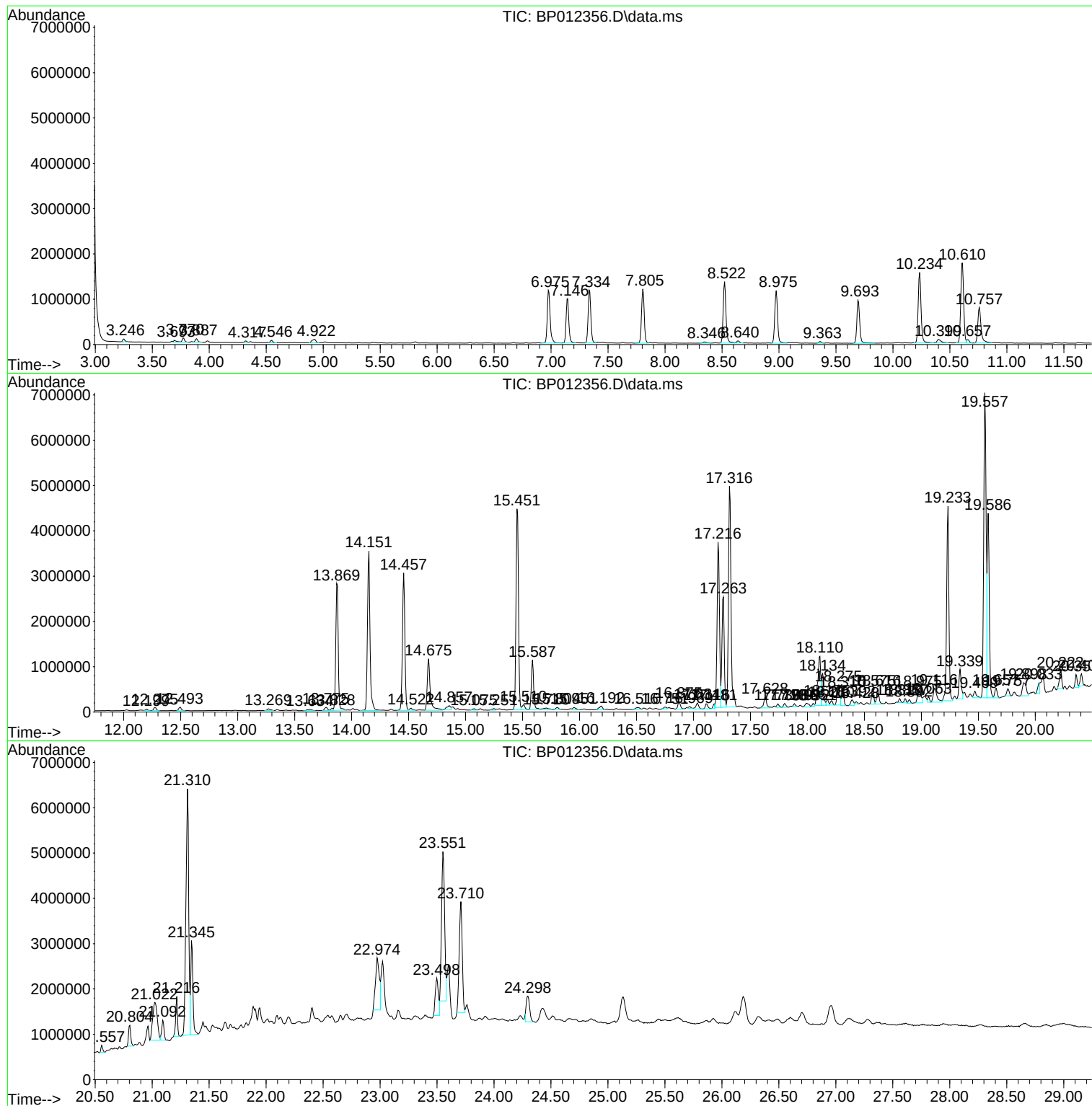
Sum of corrected areas: 130898013

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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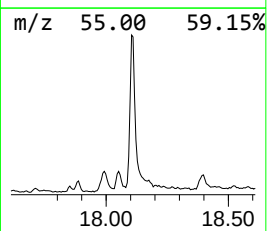
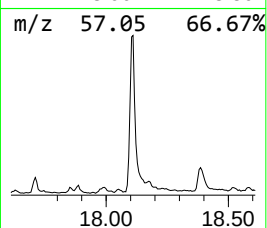
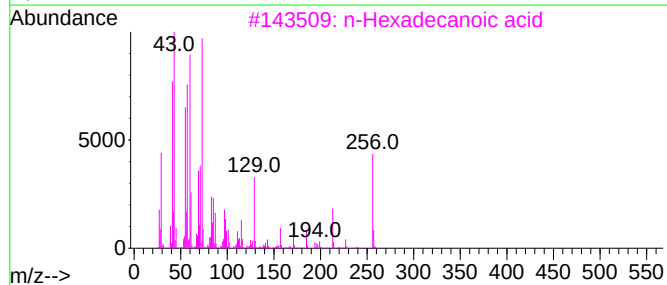
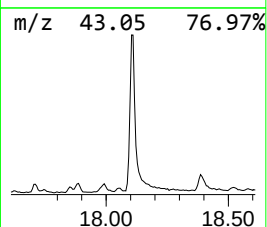
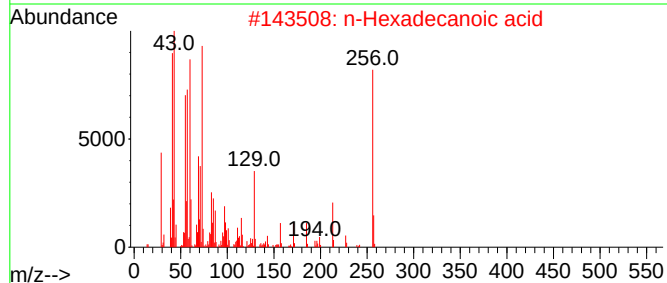
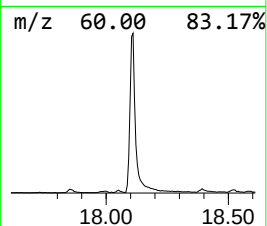
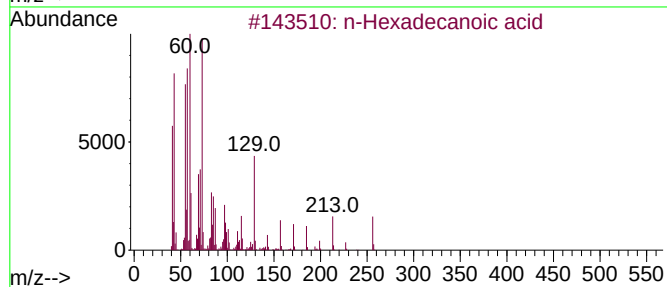
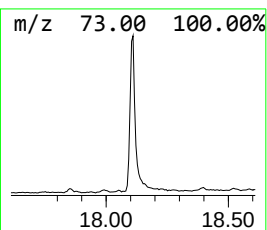
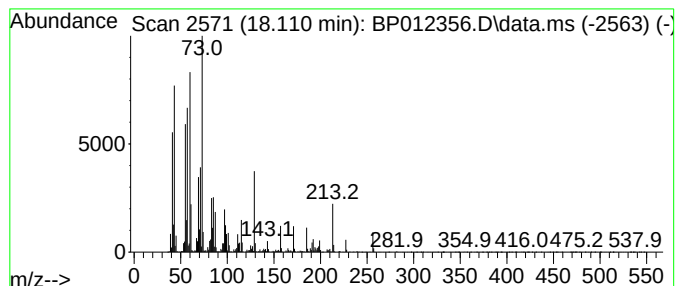
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.110	7.53 ng/ul	2032400	Phenanthrene-d10	17.216

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	98
4	n-Hexadecanoic acid			256	C16H32O2	000057-10-3	98
5	n-Hexadecanoic acid			256	C16H32O2	000057-10-3	95



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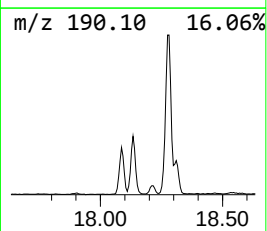
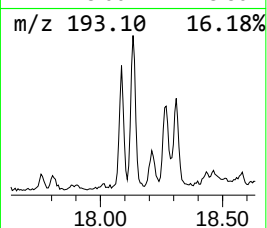
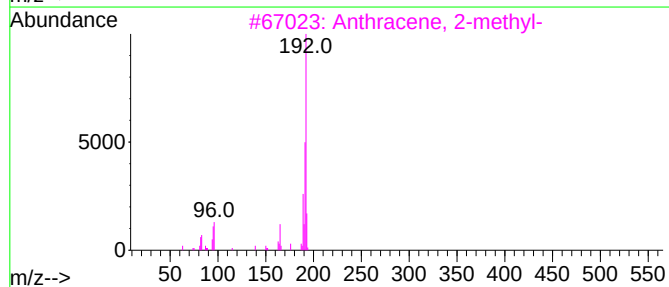
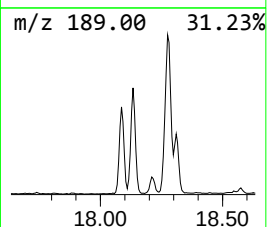
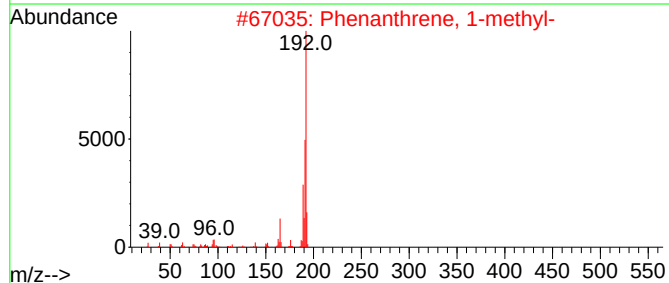
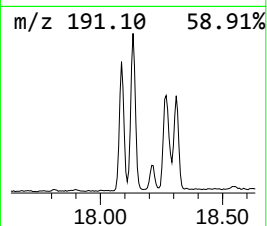
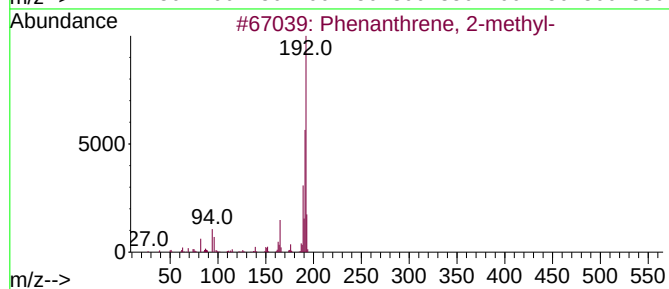
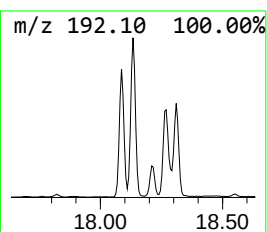
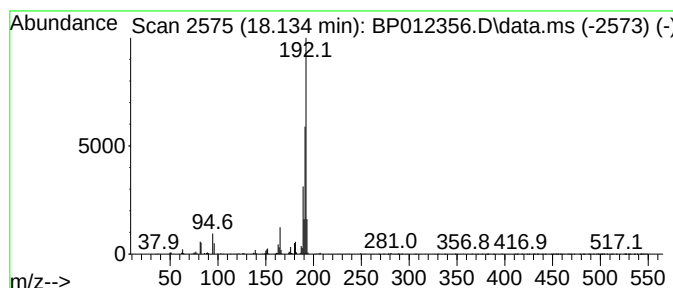
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Phenanthrene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.134	3.22 ng/ul	870556	Phenanthrene-d10	17.216

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	91
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90
5			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	90



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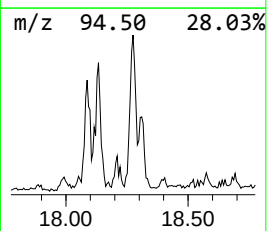
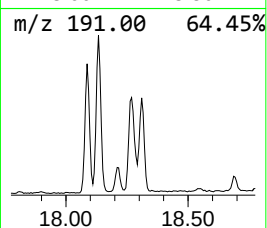
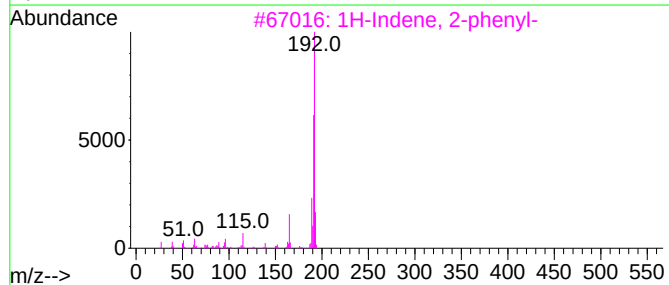
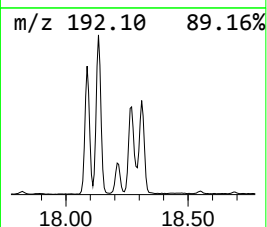
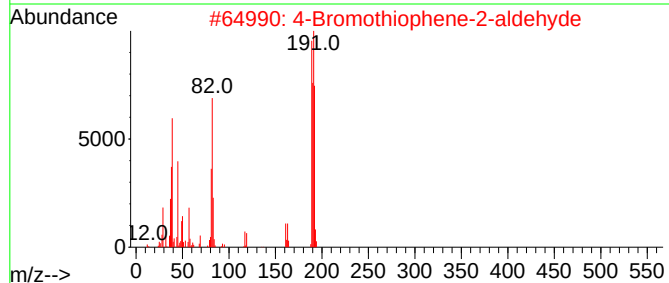
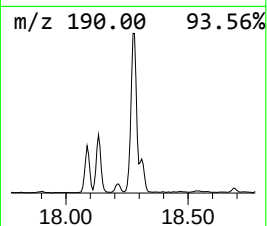
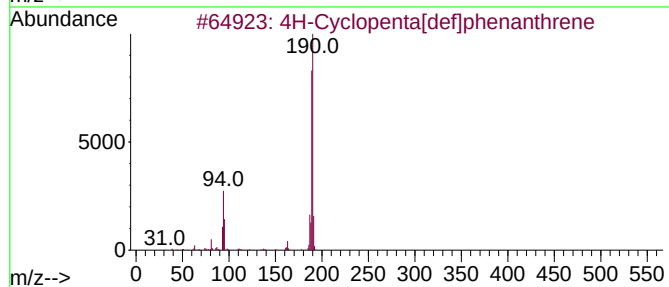
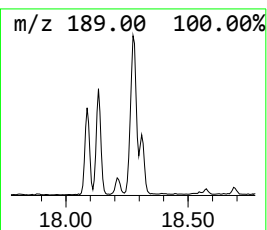
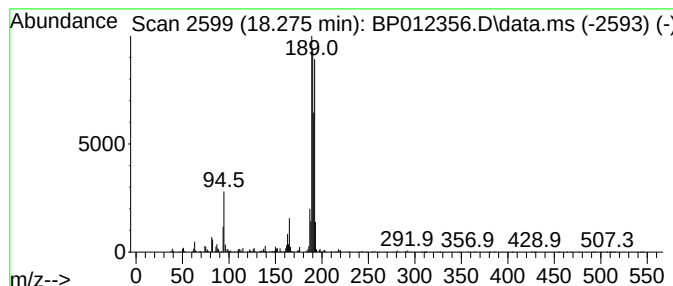
Instrument :
BNA_P
ClientSampleId :
C0BK1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.275	2.91 ng/ul	786495	Phenanthrene-d10	17.216	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	53
2	4-Bromothiophene-2-aldehyde	190	C5H3BrOS	018791-75-8	52
3	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	42
4	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	42
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	42



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Data File : BP012356.D
Acq On : 27 Oct 2022 22:57
Operator : CG/JU
Sample : N5144-01
Misc :
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Instrument :
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ClientSampleId :
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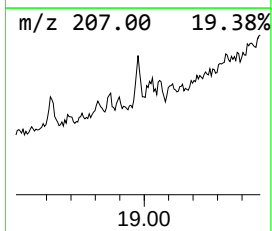
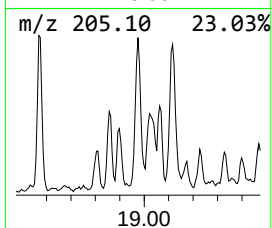
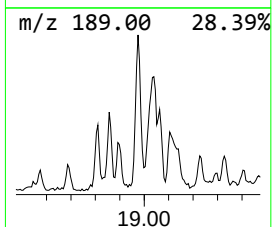
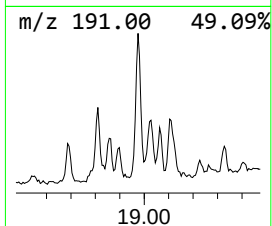
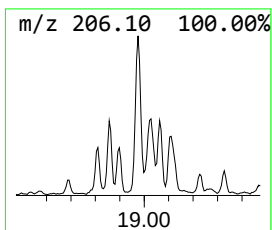
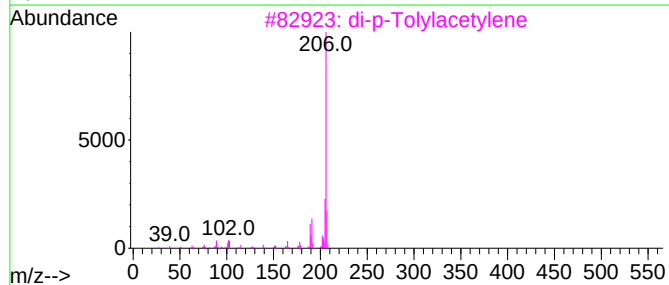
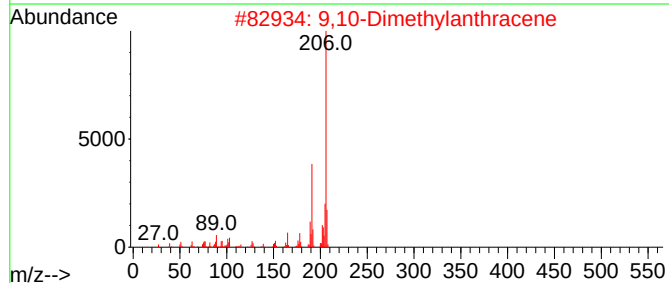
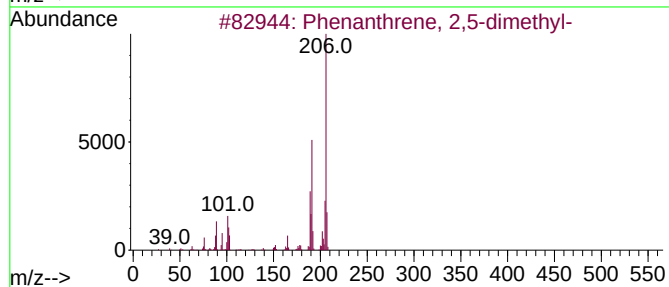
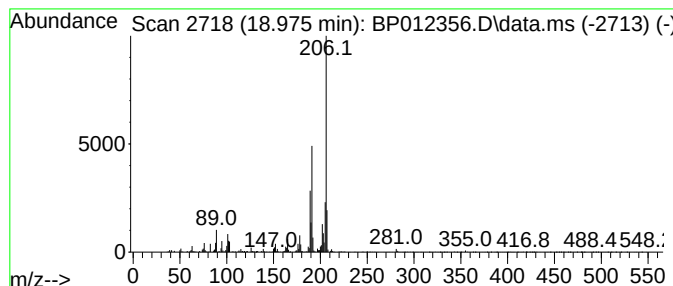
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenanthrene, 2,5-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.975	2.28 ng/ul	616143	Phenanthrene-d10	17.216

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	97
2			9,10-Dimethylantracene	206	C16H14	000781-43-1	94
3			di-p-Tolylacetylene	206	C16H14	002789-88-0	93
4			3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	91
5			9,10-Dimethylantracene	206	C16H14	000781-43-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012356.D
Acq On : 27 Oct 2022 22:57
Operator : CG/JU
Sample : N5144-01
Misc :
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Instrument :
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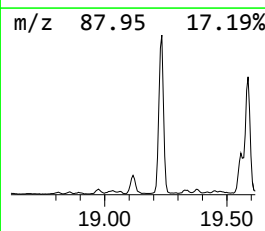
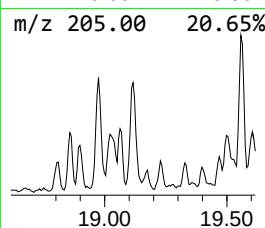
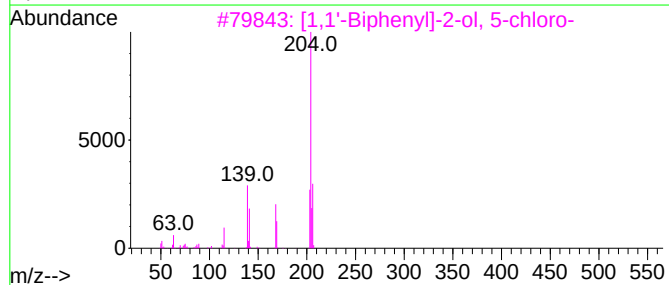
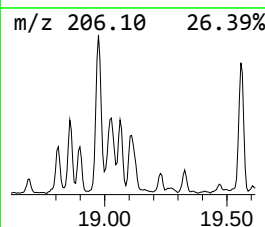
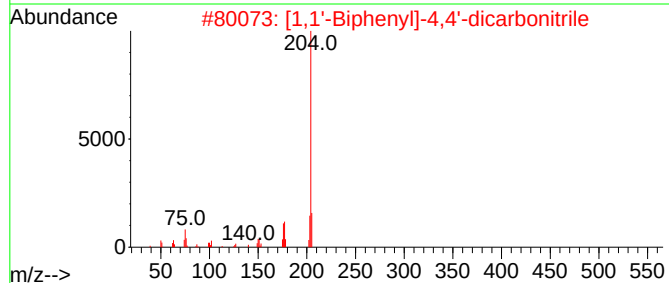
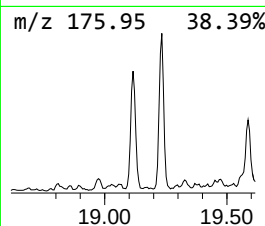
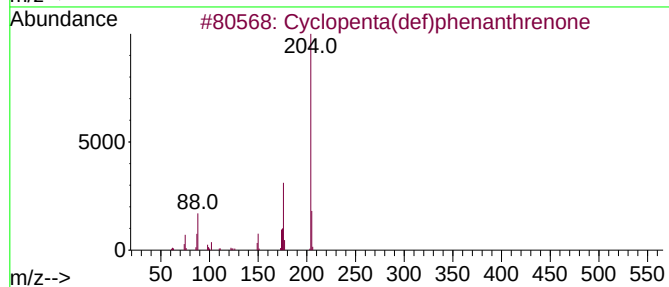
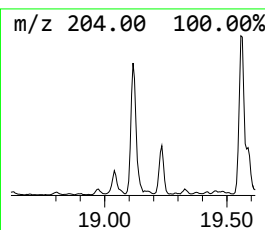
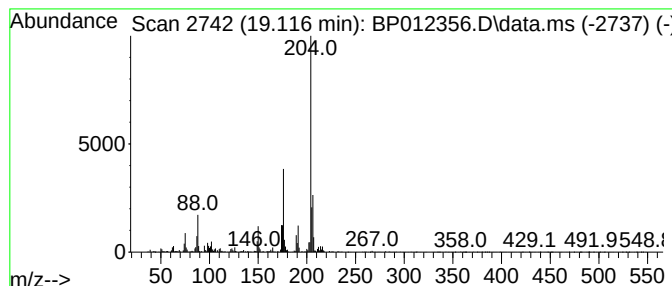
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Cyclopenta(def)phenanthrenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.116	2.59 ng/ul	698365	Phenanthrene-d10	17.216

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	98
2			[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	52
3			[1,1'-Biphenyl]-2-ol, 5-chloro-	204	C12H9ClO	000607-12-5	46
4			[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	43
5			[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012356.D
Acq On : 27 Oct 2022 22:57
Operator : CG/JU
Sample : N5144-01
Misc :
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Instrument :
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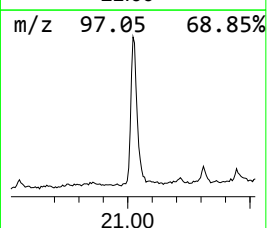
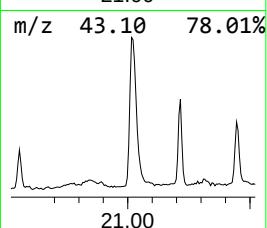
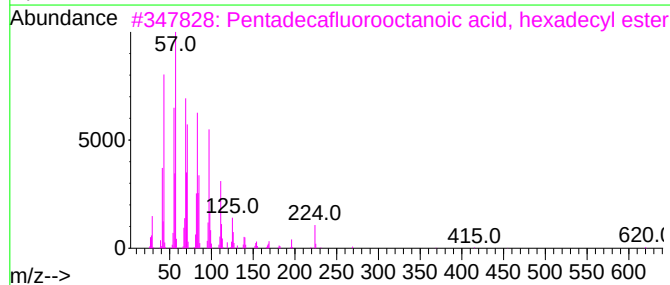
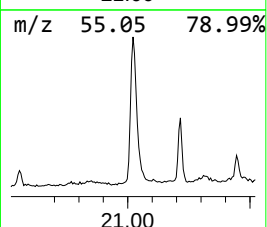
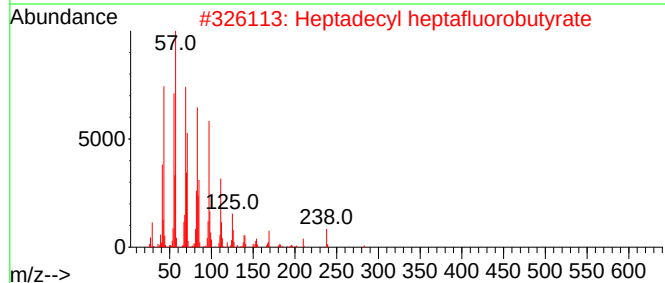
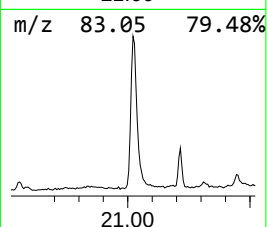
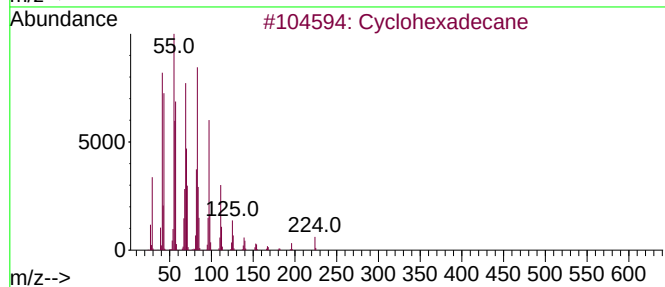
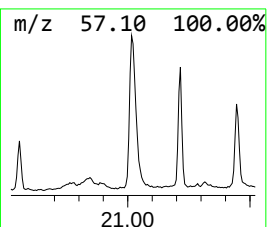
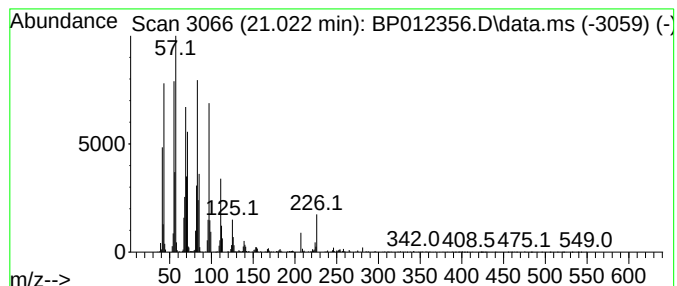
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Cyclic21.022 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.022	4.99 ng/ul	2325010	Chrysene-d12	21.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	94
2			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
3			Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	94
4			1-Hexacosene	364	C26H52	018835-33-1	94
5			Cyclohexadecane	224	C16H32	000295-65-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012356.D
Acq On : 27 Oct 2022 22:57
Operator : CG/JU
Sample : N5144-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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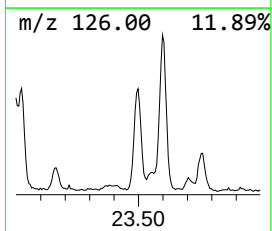
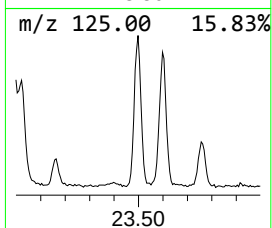
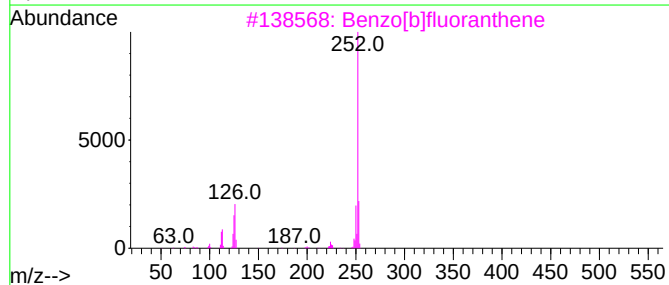
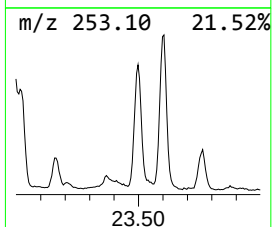
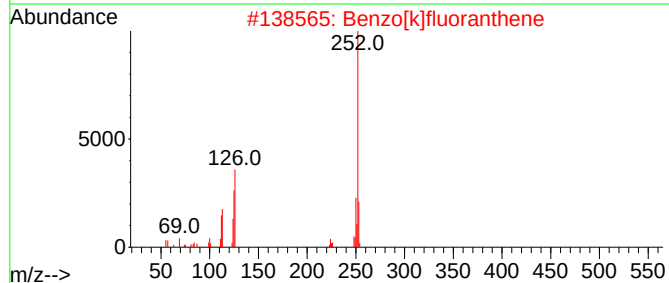
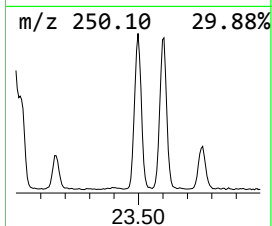
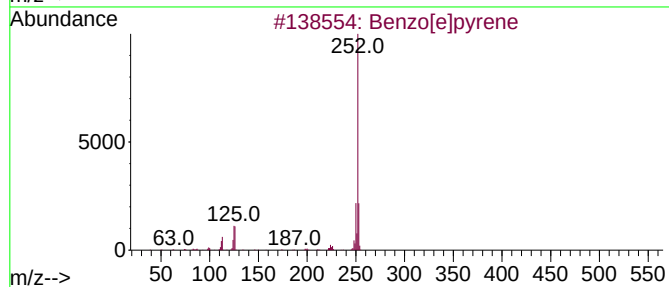
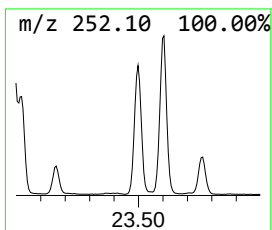
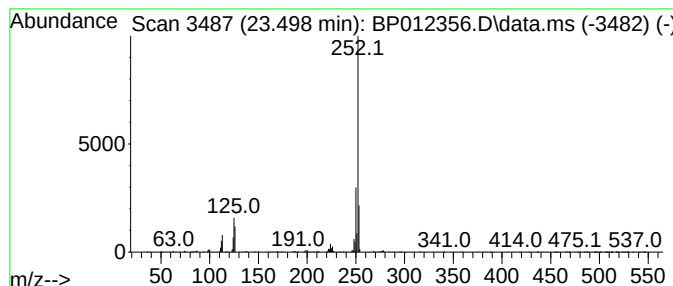
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.498	6.98 ng/ul	1581590	Perylene-d12	23.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	96
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	95
4			Benzo[a]pyrene	252	C20H12	000050-32-8	93
5			Benzo[k]fluoranthene	252	C20H12	000207-08-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012356.D
Acq On : 27 Oct 2022 22:57
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Sample : N5144-01
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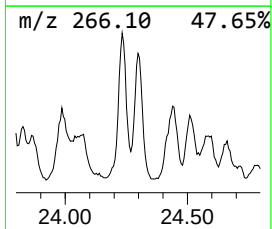
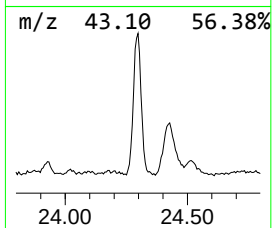
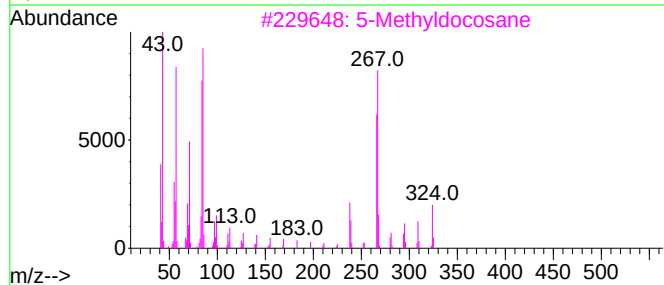
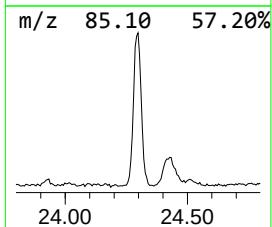
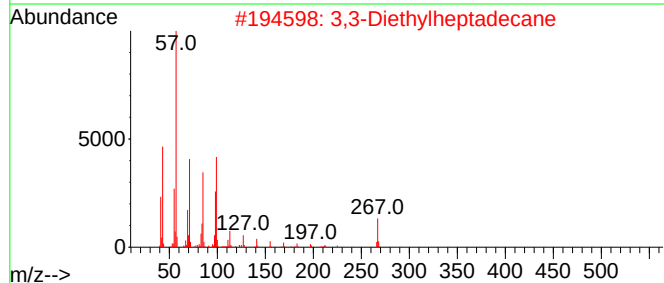
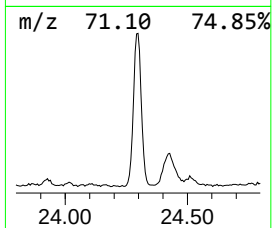
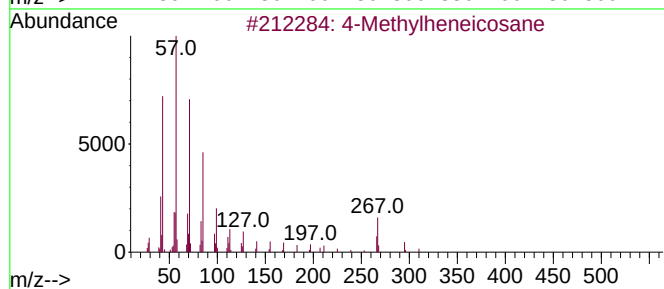
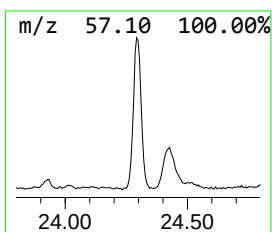
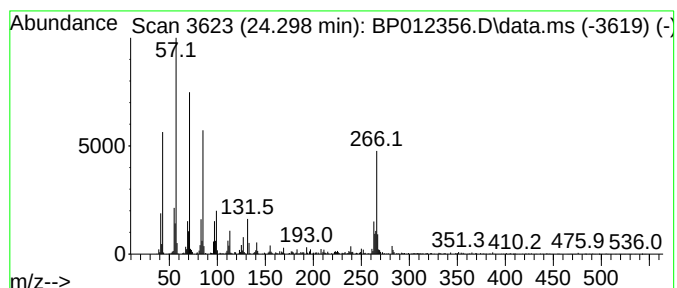
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.298	5.13 ng/ul	1162880	Perylene-d12	23.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Methylheneicosane	310	C22H46	025117-29-7	25
2			3,3-Diethylheptadecane	296	C21H44	1000360-41-6	23
3			5-Methyldocosane	324	C23H48	025163-52-4	17
4			2-Anthracenecarboxamide, 9,10-di...	335	C20H17NO4	1000351-03-6	16
5			9-Octadecenoic acid (Z)-, pentyl...	352	C23H44O2	000142-57-4	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012356.D
Acq On : 27 Oct 2022 22:57
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Sample : N5144-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
n-Hexadecanoic ...	18.110	7.5	ng/u1	2032400	4	17.216	5399130	20.0
Phenanthrene, 2...	18.134	3.2	ng/u1	870556	4	17.216	5399130	20.0
4H-Cyclopenta[d...	18.275	2.9	ng/u1	786495	4	17.216	5399130	20.0
Phenanthrene, 2...	18.975	2.3	ng/u1	616143	4	17.216	5399130	20.0
Cyclopenta(def)...	19.116	2.6	ng/u1	698365	4	17.216	5399130	20.0
(DEL) Alkane: C...	21.022	5.0	ng/u1	2325010	5	21.310	9326560	20.0
Benzo[e]pyrene	23.498	7.0	ng/u1	1581590	6	23.710	4530960	20.0
(DEL) Alkane: S...	24.298	5.1	ng/u1	1162880	6	23.710	4530960	20.0