

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121823\
 Data File : BP018905.D
 Acq On : 18 Dec 2023 14:10
 Operator : MA/JU
 Sample : 05794-03 5X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGRF9

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-BP120123.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP018905.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.258	26	29	37	rVB3	10505	20416	0.33%	0.039%
2	3.687	99	102	108	rVB3	16722	22279	0.36%	0.042%
3	3.781	114	118	123	rVB	38352	46259	0.74%	0.087%
4	3.999	151	155	159	rBV	17996	22261	0.36%	0.042%
5	4.558	248	250	257	rVB5	11399	17100	0.27%	0.032%
6	4.928	308	313	323	rVB4	9917	20688	0.33%	0.039%
7	5.022	327	329	336	rVB8	6338	9768	0.16%	0.018%
8	5.134	344	348	353	rBV4	6125	9490	0.15%	0.018%
9	5.458	399	403	408	rBV4	7627	12147	0.19%	0.023%
10	5.816	459	464	465	rBV5	7905	11522	0.18%	0.022%
11	5.834	465	467	471	rVB5	7230	8128	0.13%	0.015%
12	6.999	655	665	682	rVB2	92648	238219	3.80%	0.450%
13	7.163	684	693	700	rBV	79231	122914	1.96%	0.232%
14	7.357	718	726	733	rVB	95921	153536	2.45%	0.290%
15	7.446	737	741	744	rBV5	5509	9114	0.15%	0.017%
16	7.822	799	805	821	rVB	838592	1357048	21.65%	2.561%
17	8.040	839	842	848	rBV7	3809	6303	0.10%	0.012%
18	8.369	892	898	906	rVB3	8315	16340	0.26%	0.031%
19	8.540	920	927	935	rBV	85915	146201	2.33%	0.276%
20	8.652	940	946	954	rVB	256623	399897	6.38%	0.755%
21	8.775	963	967	973	rBV2	13787	23037	0.37%	0.043%
22	8.987	996	1003	1010	rBV	79716	132087	2.11%	0.249%
23	9.710	1118	1126	1131	rBV	59177	102975	1.64%	0.194%
24	9.869	1146	1153	1160	rBV	44426	84339	1.35%	0.159%
25	10.069	1182	1187	1192	rVB2	23885	36694	0.59%	0.069%
26	10.246	1210	1217	1226	rBV	102338	166318	2.65%	0.314%
27	10.493	1251	1259	1262	rBV5	10941	22081	0.35%	0.042%
28	10.622	1272	1281	1286	rBV	1017141	1654678	26.40%	3.123%
29	10.669	1286	1289	1297	rVV	115791	185206	2.96%	0.350%
30	10.769	1300	1306	1322	rVB3	30463	82049	1.31%	0.155%
31	11.169	1368	1374	1381	rBV3	18241	39222	0.63%	0.074%
32	11.416	1409	1416	1421	rBV	17672	33792	0.54%	0.064%
33	11.657	1454	1457	1463	rVB7	4783	6718	0.11%	0.013%
34	11.816	1479	1484	1487	rBV7	3972	7704	0.12%	0.015%
35	11.975	1504	1511	1519	rBV2	23898	66215	1.06%	0.125%
36	12.134	1533	1538	1543	rVB8	6057	11583	0.18%	0.022%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
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37	12.245	1543	1557	1558	rBV10	13527	41567	0.66%	0.078%
38	12.281	1558	1563	1570	rVB	91244	159294	2.54%	0.301%
39	12.410	1580	1585	1586	rBV5	6621	8557	0.14%	0.016%
40	12.504	1596	1601	1608	rVB	57628	99144	1.58%	0.187%
41	12.616	1616	1620	1623	rBV6	4892	8506	0.14%	0.016%
42	13.122	1702	1706	1711	rBV2	13010	20577	0.33%	0.039%
43	13.292	1724	1735	1741	rBV	39885	87599	1.40%	0.165%
44	13.492	1765	1769	1772	rVB	11552	14125	0.23%	0.027%
45	13.628	1784	1792	1801	rVB3	29371	81512	1.30%	0.154%
46	13.781	1810	1818	1823	rBV3	51290	108232	1.73%	0.204%
47	13.834	1823	1827	1830	rVV	31985	51861	0.83%	0.098%
48	13.875	1830	1834	1839	rVV	170614	235785	3.76%	0.445%
49	13.916	1839	1841	1845	rVB5	7167	8269	0.13%	0.016%
50	14.016	1851	1858	1862	rBV2	26753	50502	0.81%	0.095%
51	14.151	1875	1881	1885	rBV	210858	369780	5.90%	0.698%
52	14.322	1908	1910	1913	rVB4	7729	7475	0.12%	0.014%
53	14.463	1925	1934	1939	rBV	1349858	2077538	33.15%	3.921%
54	14.522	1939	1944	1953	rVB	323203	576150	9.19%	1.087%
55	14.681	1966	1971	1977	rVB4	41454	76587	1.22%	0.145%
56	14.857	1992	2001	2010	rBV	295811	559304	8.92%	1.056%
57	14.934	2011	2014	2018	rVB2	24442	33615	0.54%	0.063%
58	15.081	2035	2039	2043	rVB2	20549	28537	0.46%	0.054%
59	15.134	2044	2048	2054	rVB	21992	30759	0.49%	0.058%
60	15.457	2097	2103	2106	rBV	272476	389048	6.21%	0.734%
61	15.504	2106	2111	2122	rVB	518243	883799	14.10%	1.668%
62	15.634	2130	2133	2136	rBV	25653	28365	0.45%	0.054%
63	15.669	2136	2139	2144	rVB5	45393	68010	1.09%	0.128%
64	15.757	2152	2154	2158	rVB	21489	22183	0.35%	0.042%
65	15.804	2158	2162	2171	rVB	81836	129902	2.07%	0.245%
66	15.951	2183	2187	2196	rVB	82546	170844	2.73%	0.322%
67	16.186	2223	2227	2234	rVB6	42790	78343	1.25%	0.148%
68	16.516	2277	2283	2288	rVB2	71194	129356	2.06%	0.244%
69	16.563	2288	2291	2297	rVB3	39785	51843	0.83%	0.098%
70	16.745	2320	2322	2325	rBV3	20416	21704	0.35%	0.041%
71	16.786	2326	2329	2333	rVB4	37095	44324	0.71%	0.084%
72	16.869	2339	2343	2350	rVB3	59125	88702	1.42%	0.167%
73	16.945	2352	2356	2357	rBV	36676	48771	0.78%	0.092%
74	17.028	2366	2370	2379	rVB	166258	254048	4.05%	0.479%
75	17.216	2396	2402	2405	rBV	1665790	2572937	41.05%	4.856%
76	17.257	2405	2409	2414	rVV	2978181	4278789	68.27%	8.076%

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77	17.310	2414	2418	2421	rVV2	366383	650157	10.37%	1.227%
78	17.345	2421	2424	2430	rVV	859787	1257087	20.06%	2.373%
79	17.404	2431	2434	2440	rVB	142125	238399	3.80%	0.450%
80	17.622	2467	2471	2480	rVB	312139	436701	6.97%	0.824%
81	18.080	2545	2549	2554	rBV	285280	537408	8.58%	1.014%
82	18.128	2554	2557	2563	rVB	400168	561738	8.96%	1.060%
83	18.210	2567	2571	2575	rVV	143296	194938	3.11%	0.368%
84	18.275	2576	2582	2585	rVV2	502280	871048	13.90%	1.644%
85	18.304	2585	2587	2593	rVB	205660	278538	4.44%	0.526%
86	18.569	2628	2632	2636	rBV	227868	281748	4.50%	0.532%
87	18.610	2636	2639	2645	rVB2	126053	150810	2.41%	0.285%
88	18.663	2645	2648	2651	rBV	80581	97149	1.55%	0.183%
89	18.975	2697	2701	2705	rBV	184314	284184	4.53%	0.536%
90	19.227	2739	2744	2751	rBV	3314721	4227746	67.46%	7.979%
91	19.557	2795	2800	2801	rBV3	869585	1266305	20.21%	2.390%
92	19.580	2801	2804	2812	rVB	2904205	3709423	59.19%	7.001%
93	20.063	2883	2886	2892	rVB2	442572	650656	10.38%	1.228%
94	20.163	2900	2903	2906	rBV	308015	399741	6.38%	0.754%
95	21.304	3091	3097	3101	rBV2	3104630	6267143	100.00%	11.828%
96	21.345	3101	3104	3114	rVB2	1692654	3881724	61.94%	7.326%
97	22.957	3374	3378	3383	rBV	937244	2010100	32.07%	3.794%
98	23.574	3479	3483	3492	rVB2	1307651	2679199	42.75%	5.057%
99	23.680	3496	3501	3506	rVB	1561424	2751362	43.90%	5.193%

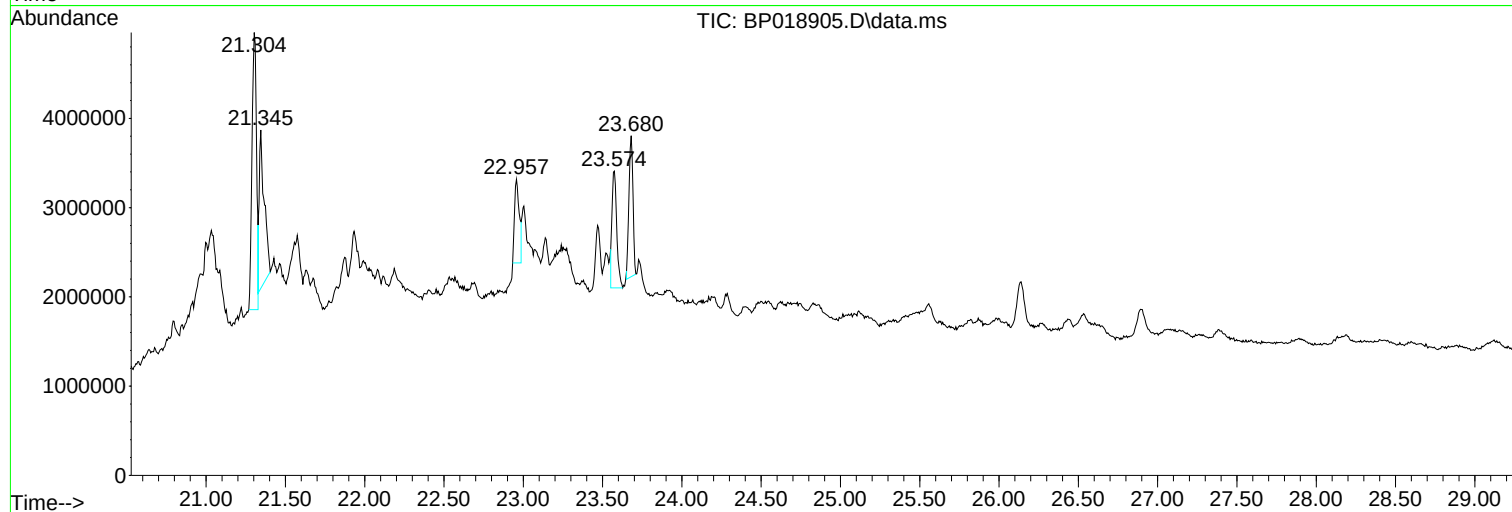
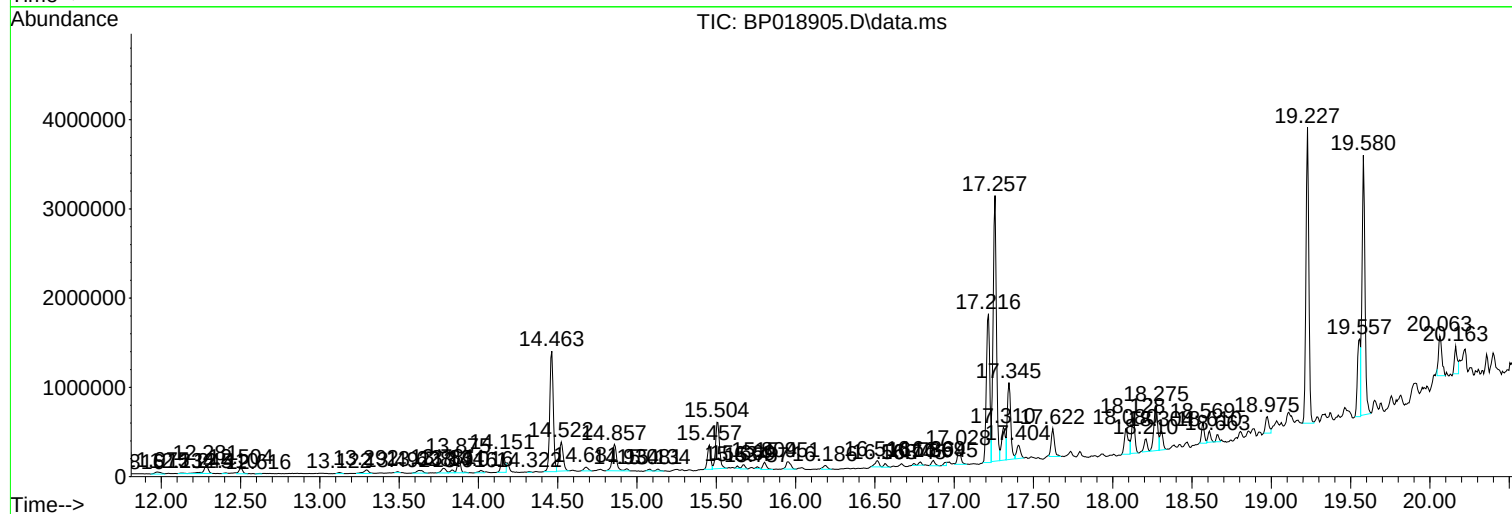
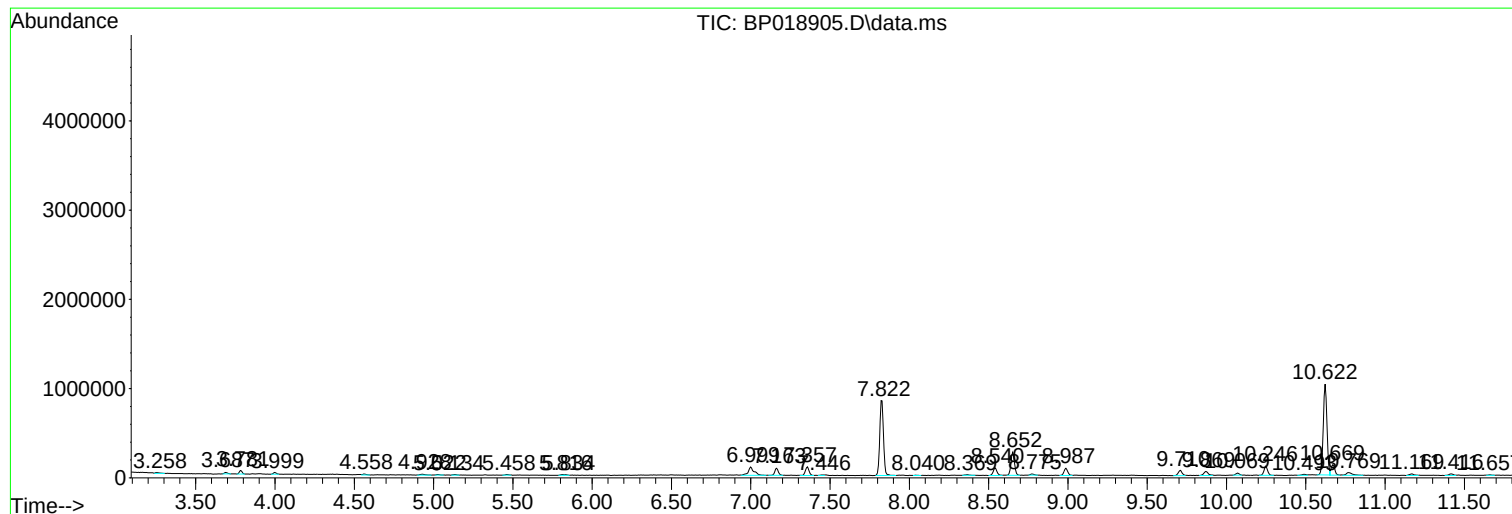
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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



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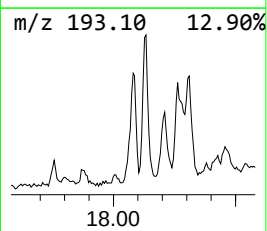
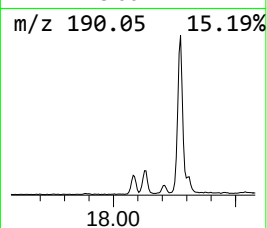
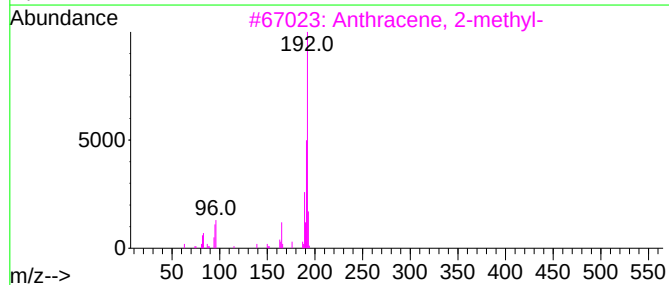
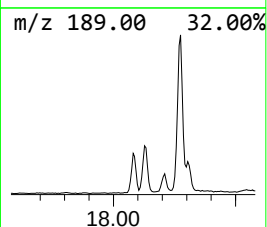
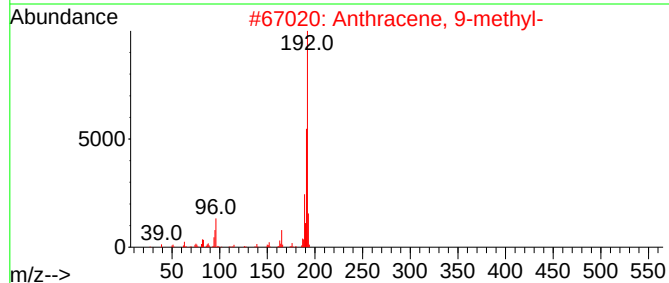
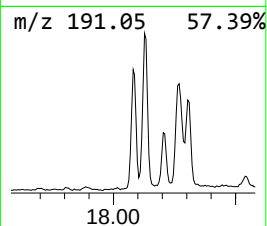
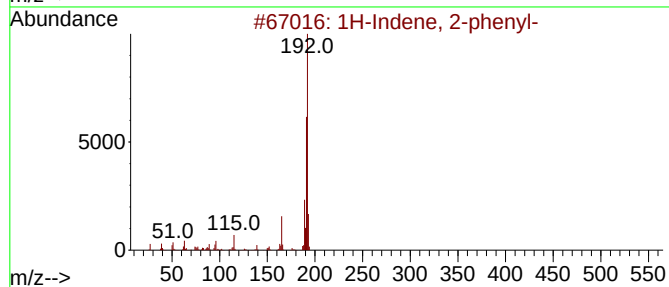
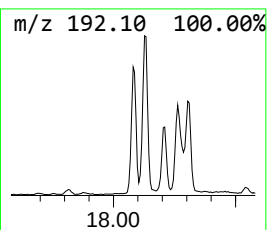
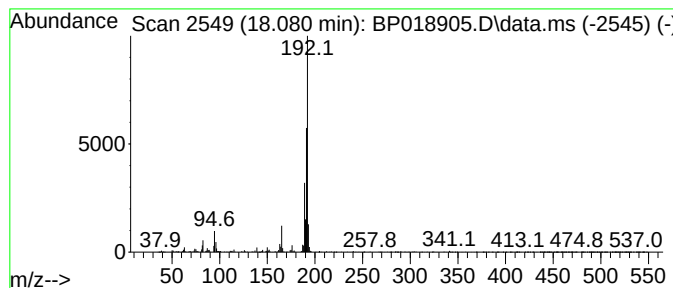
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1H-Indene, 2-phenyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.081	4.18 ng/ul	537408	Phenanthrene-d10	17.216

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	91
2		Anthracene, 9-methyl-	192	C15H12	000779-02-2	90
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
4		Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	89



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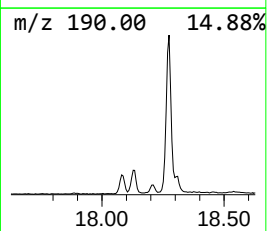
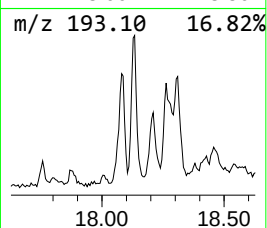
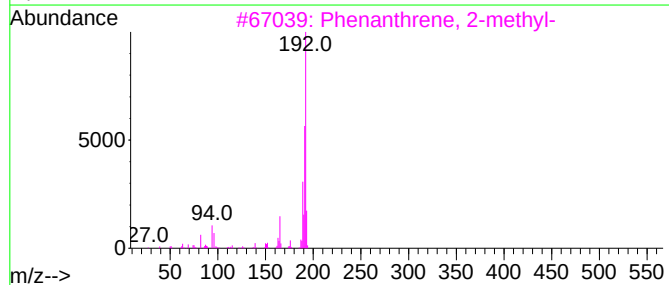
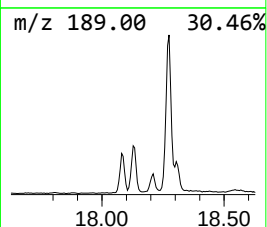
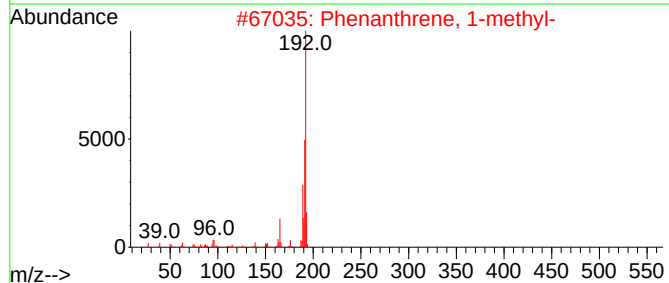
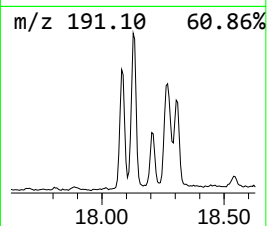
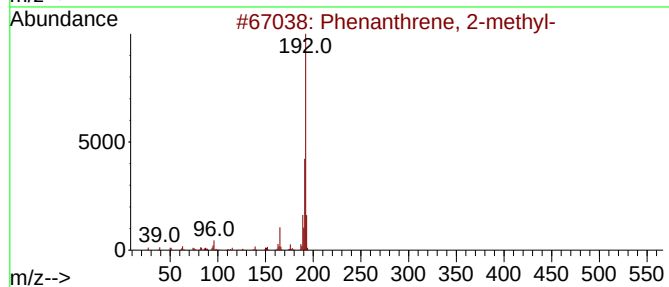
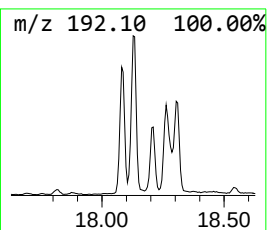
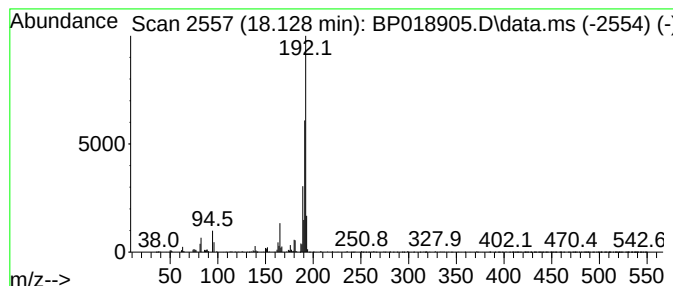
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.128	4.37 ng/ul	561738	Phenanthrene-d10	17.216

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
4		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	95
5		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95



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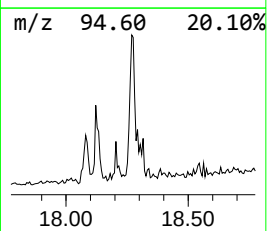
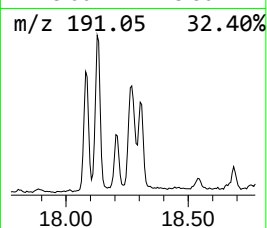
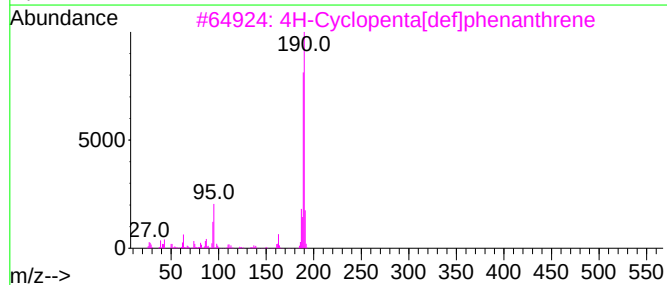
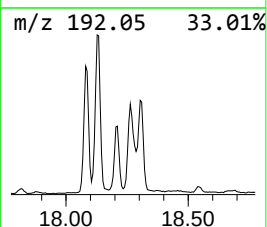
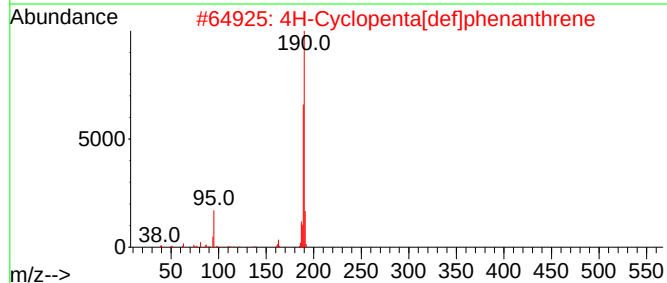
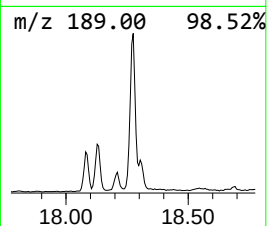
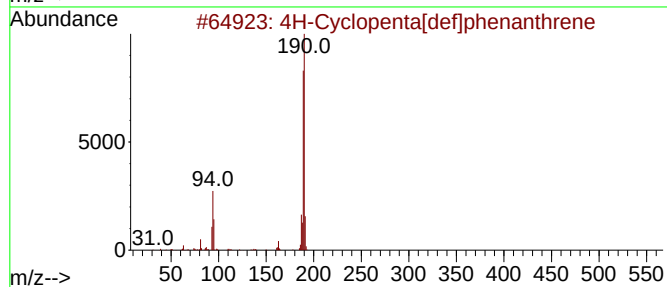
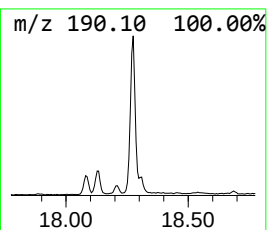
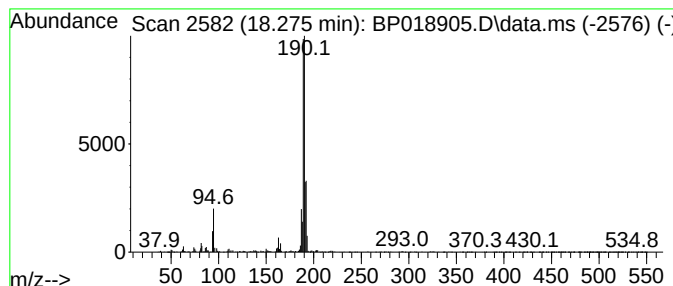
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.275	6.77 ng/ul	871048	Phenanthrene-d10	17.216

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	89
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	68
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
4			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	59
5			Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121823\
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Acq On : 18 Dec 2023 14:10
Operator : MA/JU
Sample : 05794-03 5X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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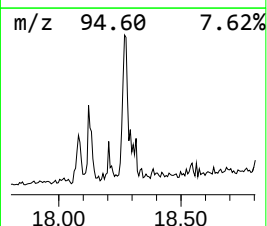
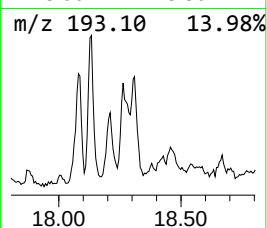
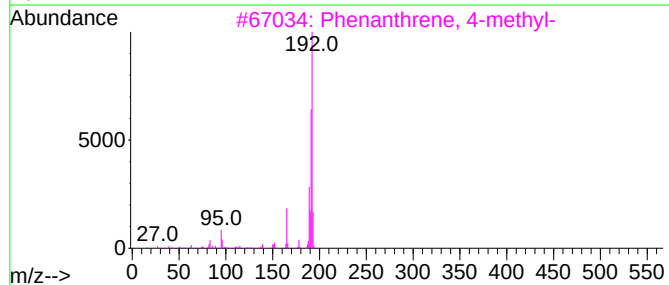
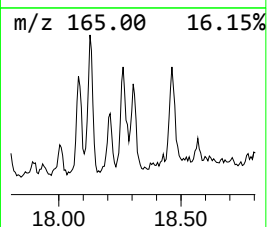
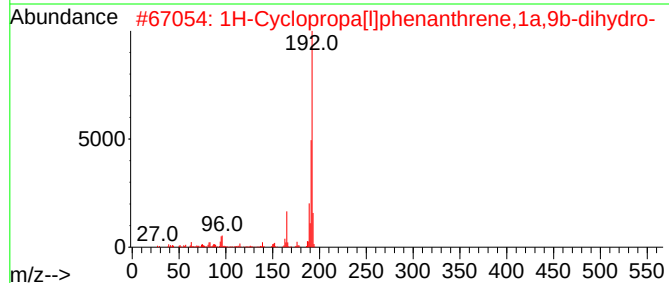
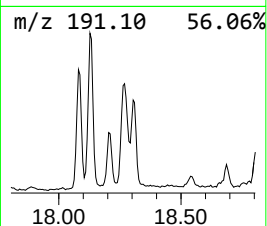
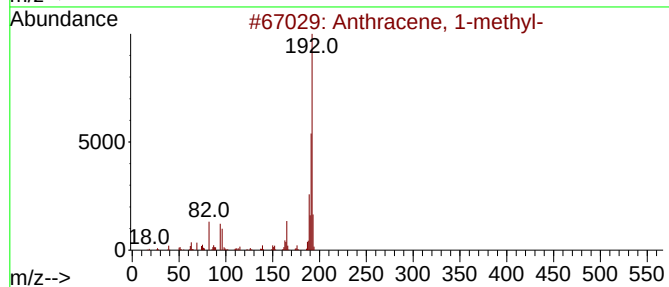
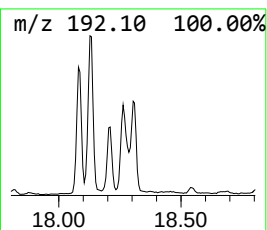
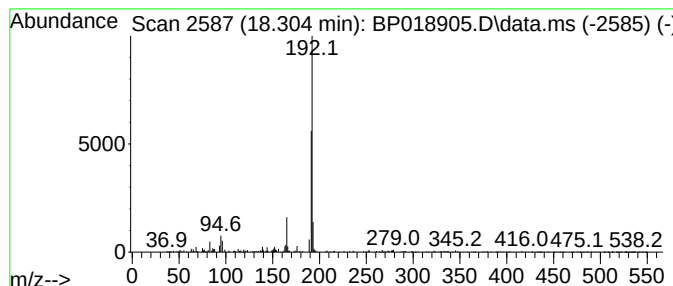
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Anthracene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.304	2.17 ng/ul	278538	Phenanthrene-d10	17.216

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
2			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	90
3			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	90
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
5			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121823\
Data File : BP018905.D
Acq On : 18 Dec 2023 14:10
Operator : MA/JU
Sample : 05794-03 5X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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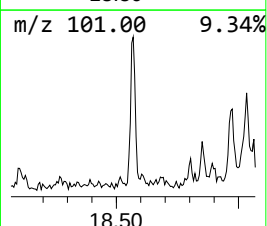
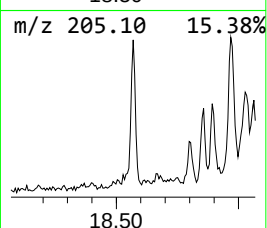
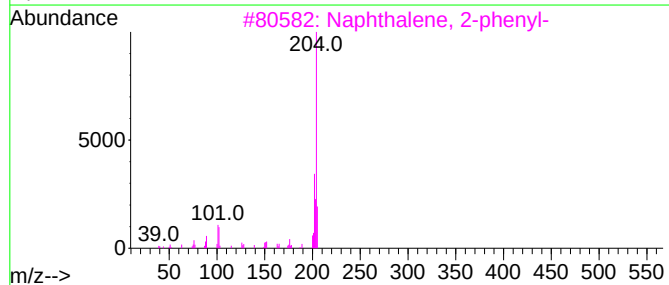
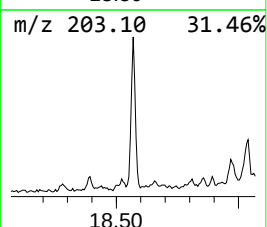
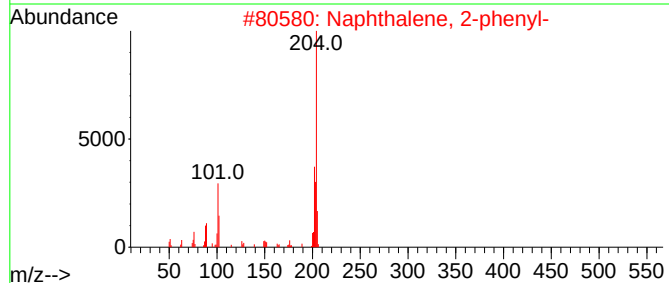
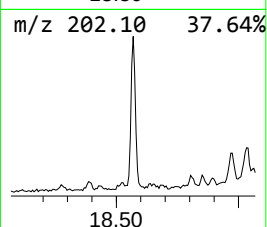
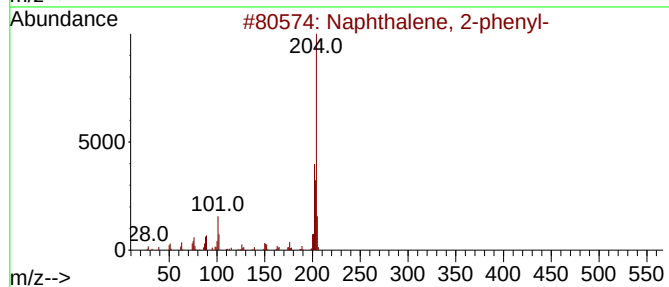
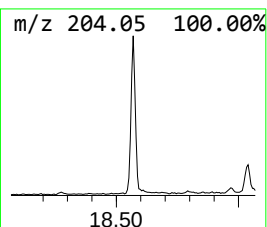
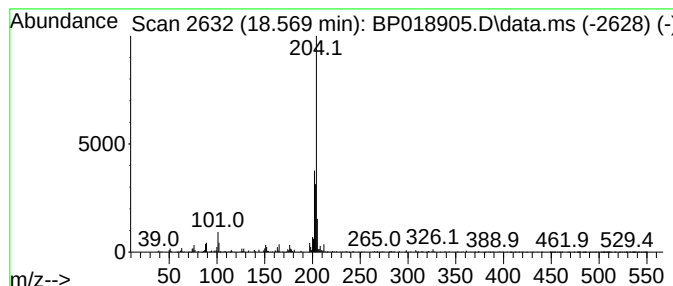
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Naphthalene, 2-phenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.569	2.19 ng/ul	281748	Phenanthrene-d10	17.216

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	95
2			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	90
3			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	86
4			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	76
5			Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121823\
Data File : BP018905.D
Acq On : 18 Dec 2023 14:10
Operator : MA/JU
Sample : 05794-03 5X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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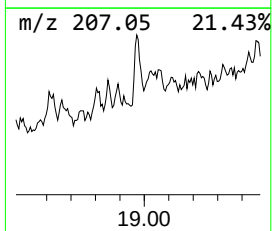
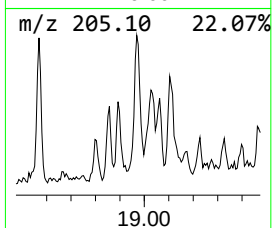
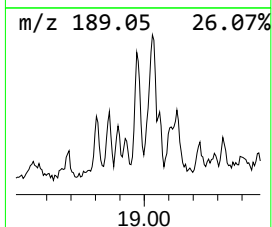
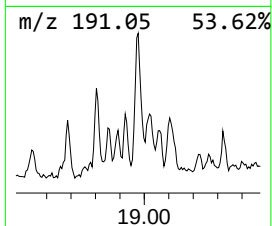
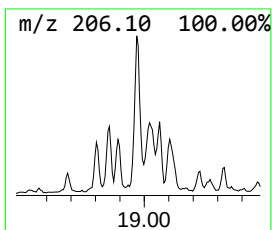
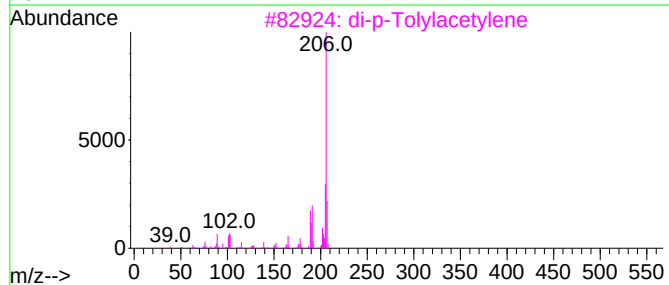
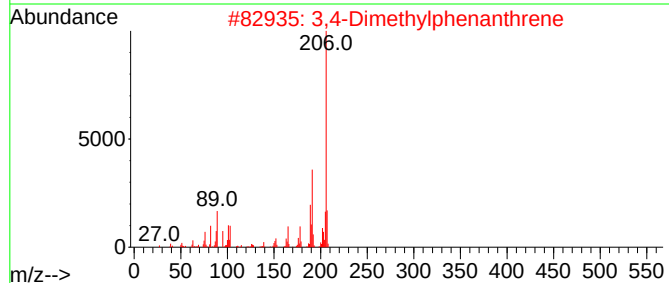
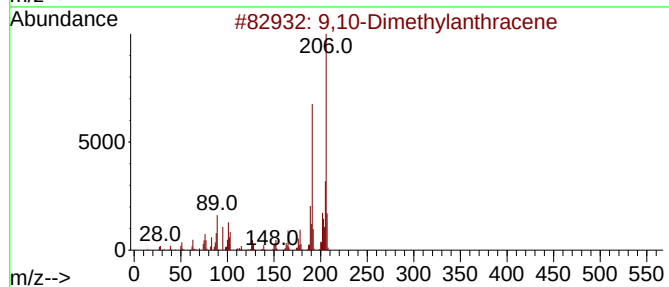
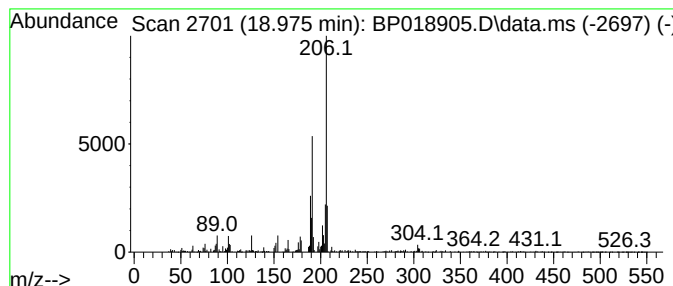
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 9,10-Dimethylantracene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.974	2.21 ng/ul	284184	Phenanthrene-d10	17.216

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Dimethylantracene	206	C16H14	000781-43-1	90
2			3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	90
3			di-p-Tolylacetylene	206	C16H14	002789-88-0	90
4			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	87
5			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121823\
Data File : BP018905.D
Acq On : 18 Dec 2023 14:10
Operator : MA/JU
Sample : 05794-03 5X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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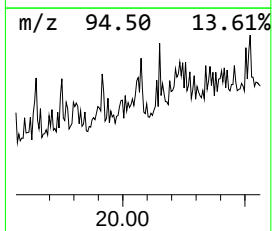
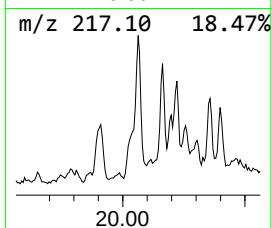
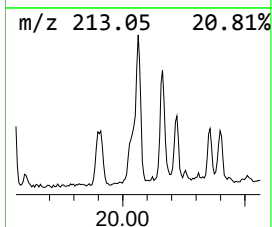
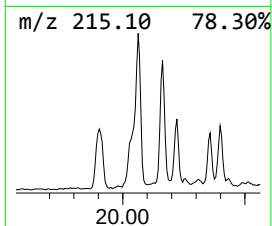
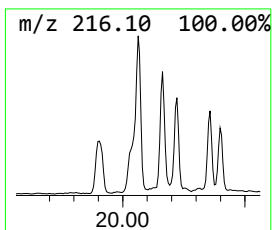
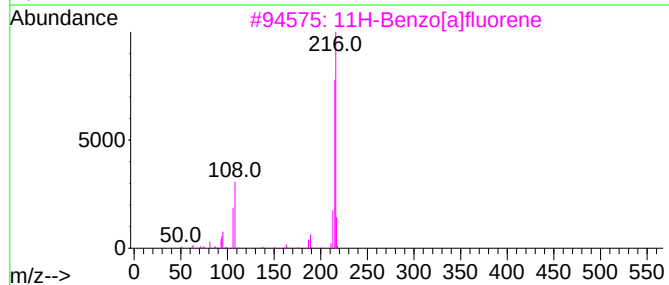
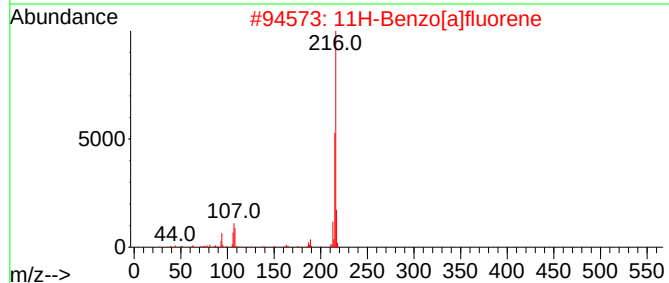
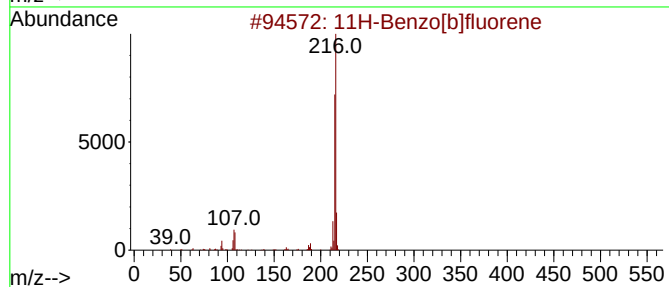
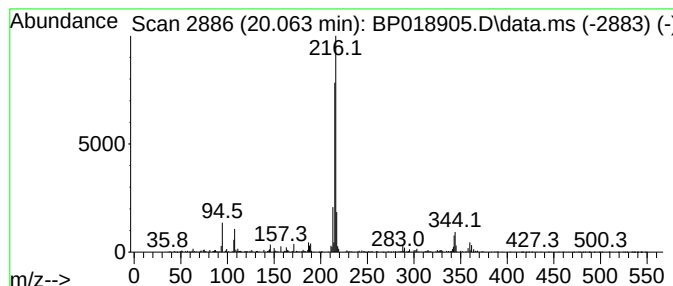
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 11H-Benzo[b]fluorene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.063	2.08 ng/ul	650656	Chrysene-d12	21.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
2			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
4			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	81
5			Pyrene, 2-methyl-	216	C17H12	003442-78-2	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121823\
Data File : BP018905.D
Acq On : 18 Dec 2023 14:10
Operator : MA/JU
Sample : 05794-03 5X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.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
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Phenanthrene, 2...	18.128	4.4	ng/ul	561738	4	17.216	2572940	20.0
4H-Cyclopenta[d...	18.275	6.8	ng/ul	871048	4	17.216	2572940	20.0
Anthracene, 1-m...	18.304	2.2	ng/ul	278538	4	17.216	2572940	20.0
Naphthalene, 2-...	18.569	2.2	ng/ul	281748	4	17.216	2572940	20.0
9,10-Dimethylan...	18.974	2.2	ng/ul	284184	4	17.216	2572940	20.0
11H-Benzo[b]flu...	20.063	2.1	ng/ul	650656	5	21.310	6267140	20.0