

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111524\
 Data File : BM048728.D
 Acq On : 16 Nov 2024 04:23
 Operator : RC/JU
 Sample : P4751-22
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BH9G3

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_M\Methods\SFAM-EPA-BM110724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM048728.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.087	31	34	40	rVB3	15621	22166	0.66%	0.064%
2	3.605	118	122	128	rBV	28394	36741	1.09%	0.106%
3	3.816	155	158	166	rVB	25106	35127	1.04%	0.101%
4	4.375	249	253	259	rBV3	12000	18194	0.54%	0.052%
5	6.810	657	667	685	rBV	153656	339283	10.08%	0.976%
6	6.963	687	693	704	rVV	153272	262477	7.80%	0.755%
7	7.157	720	726	737	rBV	198404	343620	10.21%	0.989%
8	7.622	798	805	815	rBV	433292	715154	21.26%	2.058%
9	8.345	922	928	941	rBV2	161547	328767	9.77%	0.946%
10	8.451	941	946	961	rVB	94776	182741	5.43%	0.526%
11	8.781	996	1002	1012	rBV	161853	301681	8.97%	0.868%
12	9.504	1119	1125	1139	rBV	123645	241363	7.17%	0.694%
13	9.716	1153	1161	1169	rBV5	7368	24341	0.72%	0.070%
14	9.875	1183	1188	1193	rBV3	7134	14513	0.43%	0.042%
15	10.051	1210	1218	1239	rBV	157702	384184	11.42%	1.105%
16	10.404	1269	1278	1295	rBV	569773	1008297	29.97%	2.901%
17	10.575	1300	1307	1321	rBV2	55400	172571	5.13%	0.497%
18	12.086	1559	1564	1571	rBV2	6543	12905	0.38%	0.037%
19	12.316	1598	1603	1613	rVB3	6811	14797	0.44%	0.043%
20	13.592	1816	1820	1824	rBV2	9300	13051	0.39%	0.038%
21	13.692	1831	1837	1854	rVB	375807	580481	17.25%	1.670%
22	13.969	1875	1884	1894	rBV	449889	730828	21.72%	2.103%
23	14.274	1930	1936	1942	rVV	865194	1217740	36.19%	3.504%
24	14.339	1942	1947	1956	rVV	71503	107589	3.20%	0.310%
25	14.533	1974	1980	1987	rBV5	18578	60076	1.79%	0.173%
26	14.680	2000	2005	2010	rBV	32827	53054	1.58%	0.153%
27	15.074	2065	2072	2076	rBV9	6468	14999	0.45%	0.043%
28	15.268	2098	2105	2111	rBV	588345	906689	26.95%	2.609%
29	15.327	2111	2115	2124	rVB	69207	111710	3.32%	0.321%
30	15.416	2124	2130	2134	rBV	69016	113087	3.36%	0.325%
31	15.492	2140	2143	2148	rVB5	9617	13830	0.41%	0.040%
32	15.645	2161	2169	2177	rBV2	105085	172939	5.14%	0.498%
33	15.774	2184	2191	2197	rBV	19854	40764	1.21%	0.117%
34	15.863	2202	2206	2213	rVB5	7620	12993	0.39%	0.037%
35	15.998	2224	2229	2237	rVB8	14178	31549	0.94%	0.091%
36	16.333	2283	2286	2291	rVB4	14970	20280	0.60%	0.058%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110724.MA.M
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37	16.380	2291	2294	2300	rVB4	9336	10920	0.32%	0.031%
38	16.563	2322	2325	2328	rVV3	11513	13717	0.41%	0.039%
39	16.604	2328	2332	2335	rVB3	11588	15917	0.47%	0.046%
40	16.686	2341	2346	2353	rVB	32369	55962	1.66%	0.161%

41	16.757	2353	2358	2360	rBV3	16360	24655	0.73%	0.071%
42	16.845	2368	2373	2382	rVB	45377	69632	2.07%	0.200%
43	17.021	2397	2403	2407	rBV	997912	1460265	43.40%	4.202%
44	17.062	2407	2410	2415	rVV	992464	1305006	38.79%	3.755%
45	17.121	2415	2420	2424	rVV	664749	990522	29.44%	2.850%

46	17.157	2424	2426	2432	rVV	196525	240810	7.16%	0.693%
47	17.227	2432	2438	2442	rVB2	18217	32759	0.97%	0.094%
48	17.433	2466	2473	2486	rVV2	77833	182163	5.41%	0.524%
49	17.545	2486	2492	2499	rVV4	21173	55608	1.65%	0.160%
50	17.604	2499	2502	2510	rVB6	13609	31316	0.93%	0.090%

51	17.745	2523	2526	2530	rVB4	9980	14376	0.43%	0.041%
52	17.898	2543	2552	2554	rBV	103210	162695	4.84%	0.468%
53	17.927	2554	2557	2558	rVV	149771	181957	5.41%	0.524%
54	17.945	2558	2560	2567	rVV	175295	255965	7.61%	0.736%
55	18.027	2570	2574	2579	rVV	48684	81501	2.42%	0.235%

56	18.092	2579	2585	2589	rVV2	207466	344846	10.25%	0.992%
57	18.127	2589	2591	2600	rVB	76073	105816	3.15%	0.304%
58	18.221	2602	2607	2609	rBV4	23697	36828	1.09%	0.106%
59	18.257	2609	2613	2617	rVV2	18557	22861	0.68%	0.066%
60	18.304	2617	2621	2625	rBV7	8529	12551	0.37%	0.036%

61	18.404	2634	2638	2642	rBV	86025	118811	3.53%	0.342%
62	18.451	2642	2646	2651	rVV	91899	126329	3.75%	0.363%
63	18.498	2651	2654	2658	rVV	25395	31957	0.95%	0.092%
64	18.651	2676	2680	2684	rBV3	29443	40212	1.20%	0.116%
65	18.704	2685	2689	2691	rBV	23619	31510	0.94%	0.091%

66	18.733	2691	2694	2699	rVB3	31526	48834	1.45%	0.141%
67	18.774	2699	2701	2704	rVB4	11543	10110	0.30%	0.029%
68	18.821	2704	2709	2713	rBV	67203	108078	3.21%	0.311%
69	18.862	2713	2716	2717	rBV3	20918	24205	0.72%	0.070%
70	18.968	2729	2734	2741	rBV2	84306	181293	5.39%	0.522%

71	19.086	2748	2754	2762	rVB	2120813	2826638	84.01%	8.133%
72	19.151	2762	2765	2768	rBV	21269	23158	0.69%	0.067%
73	19.186	2768	2771	2772	rBV2	35500	36280	1.08%	0.104%
74	19.421	2805	2811	2813	rBV2	1069727	1607245	47.77%	4.625%
75	19.445	2813	2815	2824	rVV	1262465	1594068	47.38%	4.587%

76	19.521	2824	2828	2834	rVB3	81100	129264	3.84%	0.372%
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77	19.633	2842	2847	2853	rBV	102011	156081	4.64%	0.449%
78	19.692	2854	2857	2862	rVB	43100	61049	1.81%	0.176%
79	19.774	2866	2871	2879	rBV3	110407	277742	8.26%	0.799%
80	19.915	2890	2895	2897	rBV4	86705	163278	4.85%	0.470%
81	19.945	2897	2900	2909	rVB2	254633	354061	10.52%	1.019%
82	20.051	2913	2918	2921	rBV	173424	232854	6.92%	0.670%
83	20.103	2921	2927	2931	rVV2	157162	289615	8.61%	0.833%
84	20.139	2931	2933	2938	rVB2	62513	80549	2.39%	0.232%
85	20.186	2938	2941	2945	rVB3	40701	52541	1.56%	0.151%
86	20.245	2948	2951	2954	rBV	65626	74881	2.23%	0.215%
87	20.292	2955	2959	2963	rVB2	83285	119176	3.54%	0.343%
88	20.698	3024	3028	3035	rVB	163915	215154	6.39%	0.619%
89	20.862	3052	3056	3060	rBV3	149398	206687	6.14%	0.595%
90	20.903	3060	3063	3066	rVV2	155025	204072	6.07%	0.587%
91	20.945	3066	3070	3076	rVV3	205228	425569	12.65%	1.224%
92	21.003	3077	3080	3087	rVB	163225	250344	7.44%	0.720%
93	21.245	3114	3121	3125	rBV3	1522613	3364471	100.00%	9.681%
94	21.292	3125	3129	3140	rVB	922796	1465071	43.55%	4.215%
95	21.427	3149	3152	3156	rVB2	104546	130133	3.87%	0.374%
96	21.909	3231	3234	3239	rVB	144563	219591	6.53%	0.632%
97	23.227	3451	3458	3464	rBV	731922	2002810	59.53%	5.763%
98	23.933	3573	3578	3584	rBV	222038	436540	12.97%	1.256%
99	23.991	3584	3588	3600	rVB	268677	674209	20.04%	1.940%
100	24.127	3604	3611	3621	rBV	808649	1990886	59.17%	5.728%

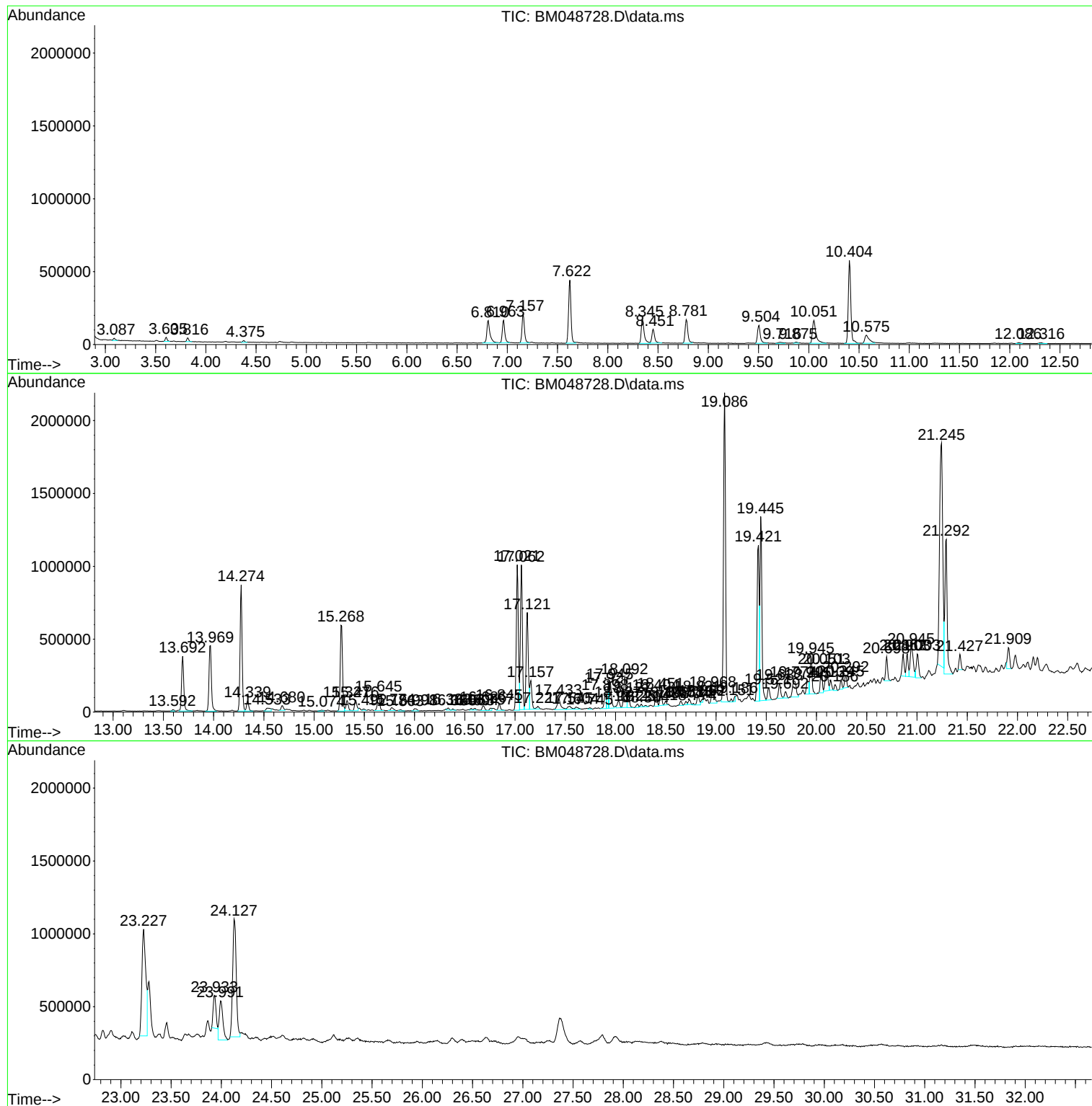
Sum of corrected areas: 34754585

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110724.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111524\
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Instrument :
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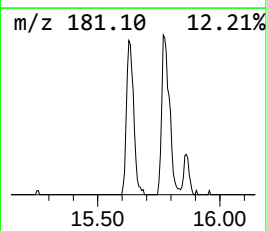
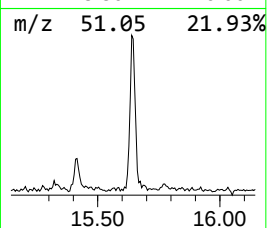
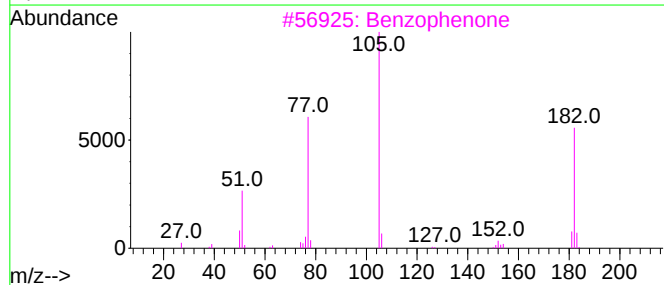
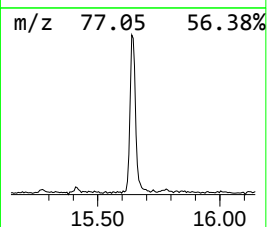
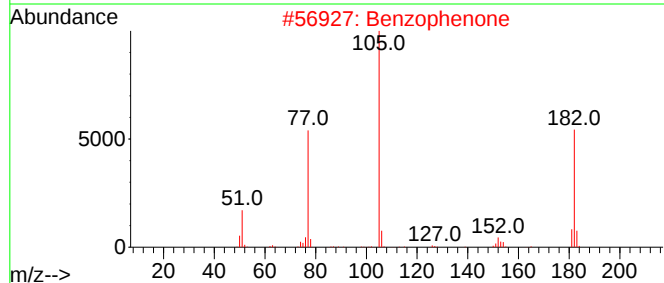
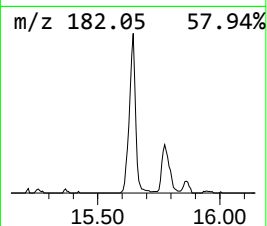
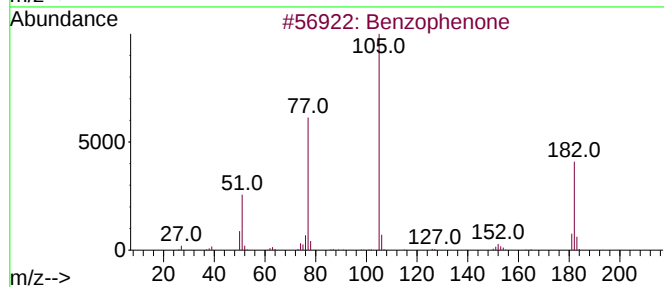
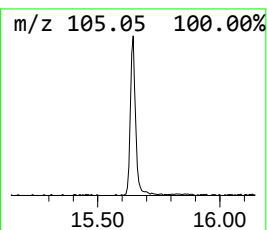
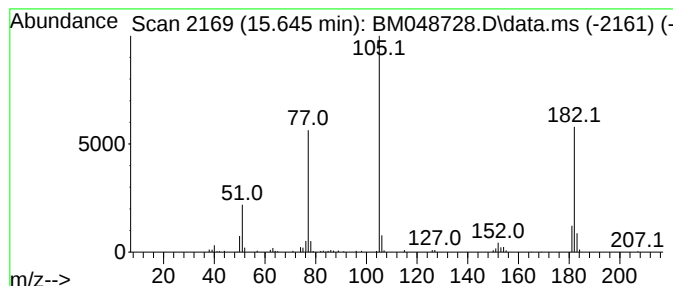
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.645	2.84 ng/ul	172939	Acenaphthene-d10	14.274

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	96
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	91



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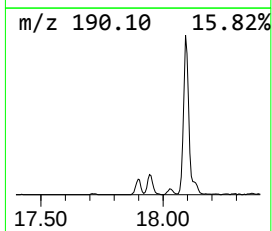
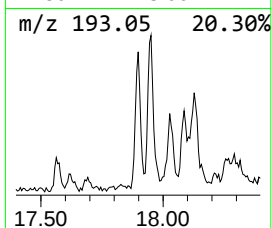
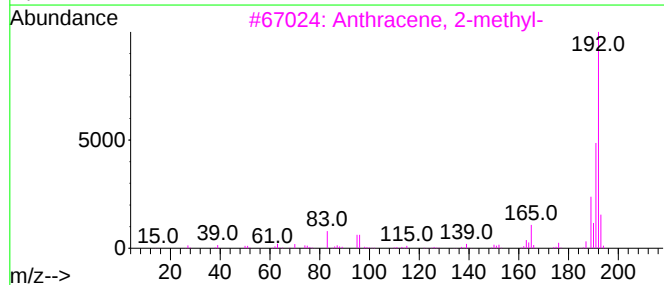
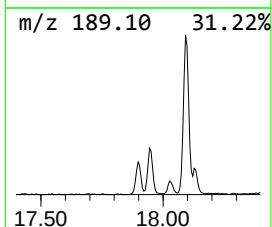
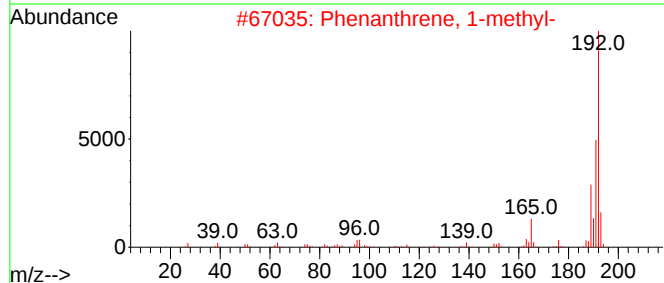
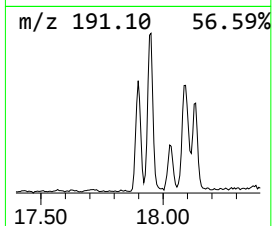
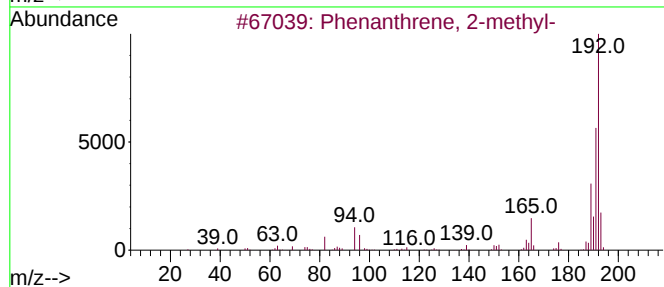
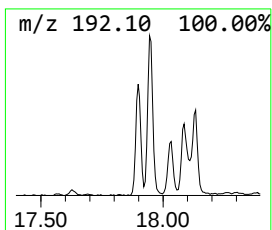
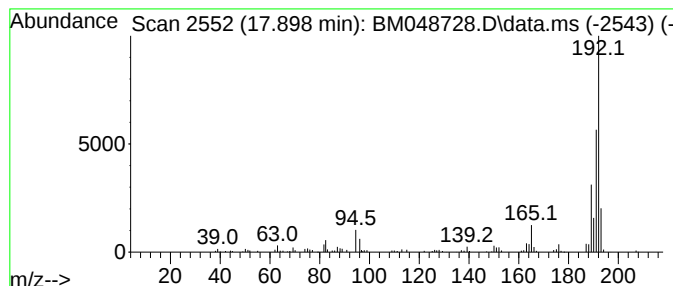
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Phenanthrene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.898	2.23 ng/ul	162695	Phenanthrene-d10	17.021

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94



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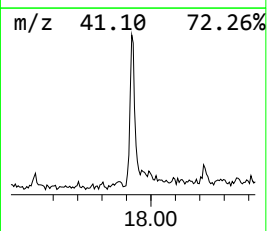
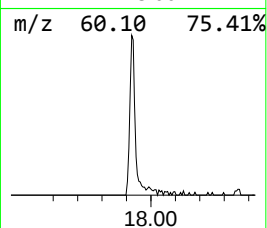
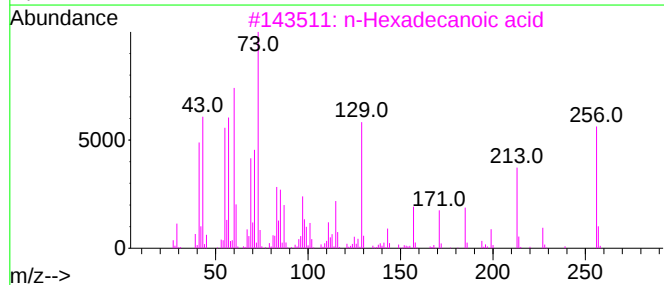
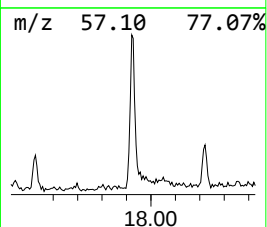
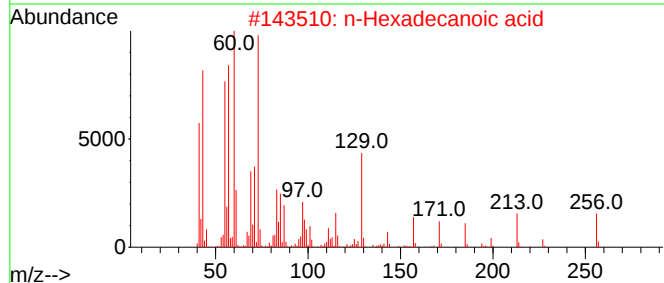
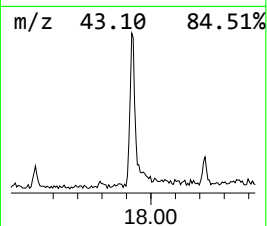
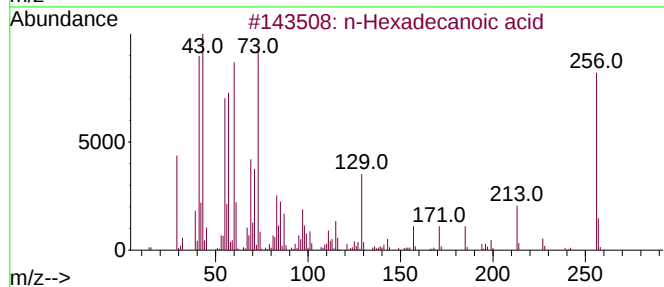
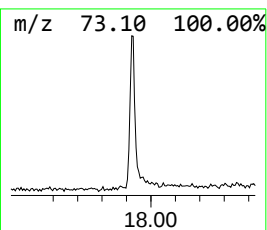
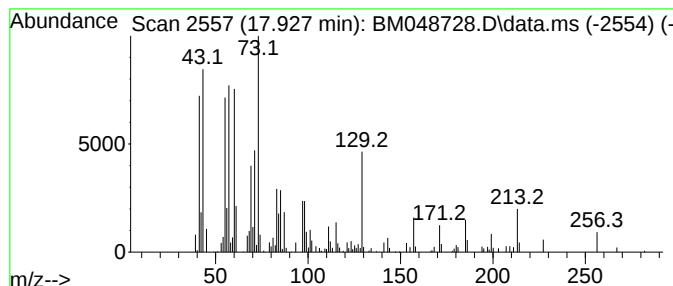
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Quant Title : SVOA CALIBRATION

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TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.927	2.49 ng/ul	181957	Phenanthrene-d10	17.021

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	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5			Tridecanoic acid	214	C13H26O2	000638-53-9	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111524\
Data File : BM048728.D
Acq On : 16 Nov 2024 04:23
Operator : RC/JU
Sample : P4751-22
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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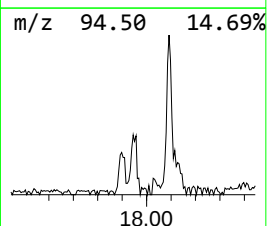
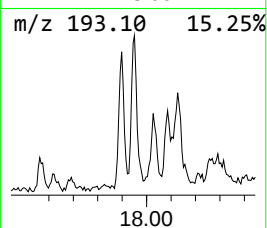
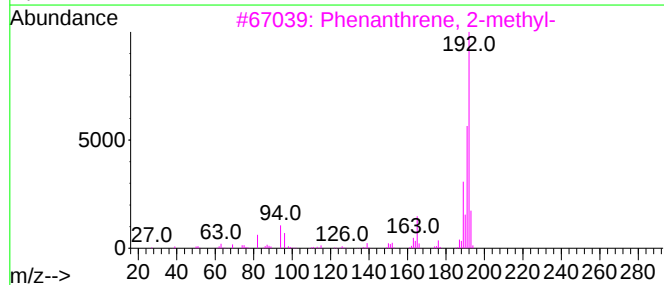
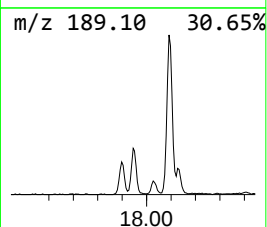
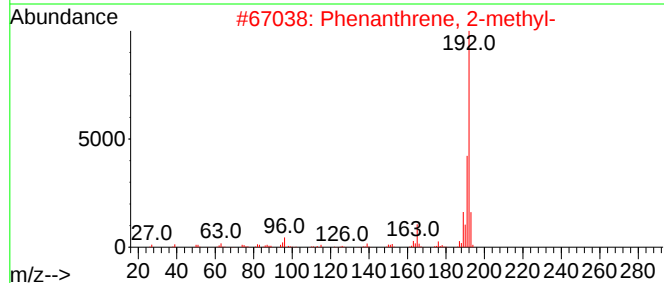
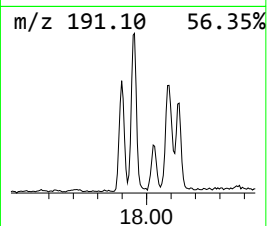
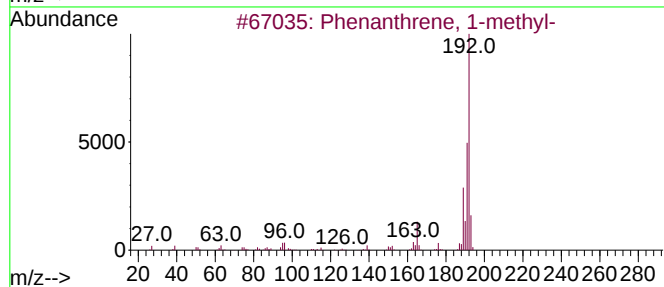
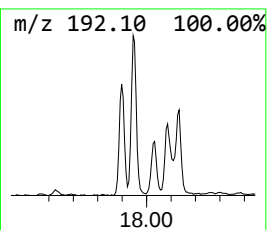
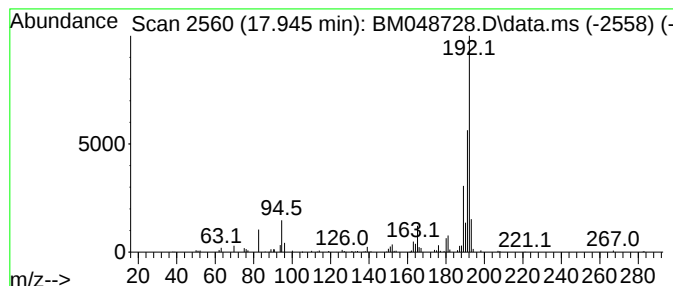
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenanthrene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.945	3.51 ng/ul	255965	Phenanthrene-d10	17.021

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111524\
Data File : BM048728.D
Acq On : 16 Nov 2024 04:23
Operator : RC/JU
Sample : P4751-22
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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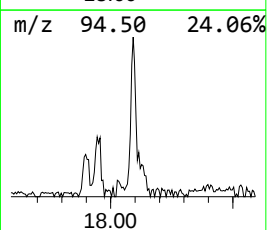
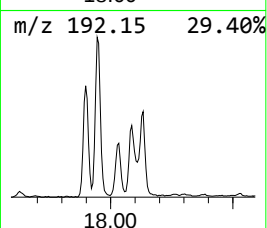
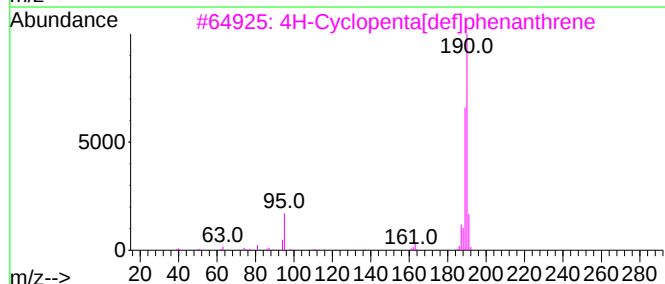
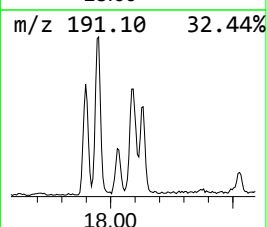
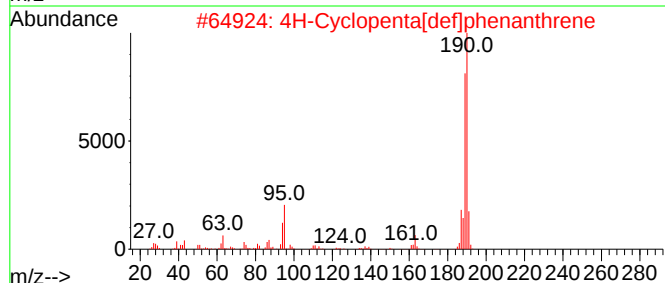
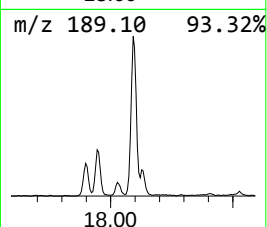
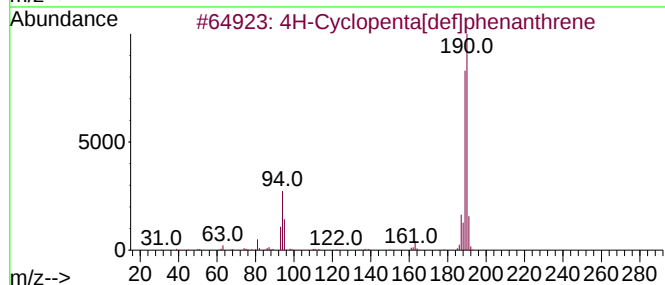
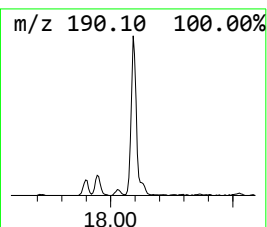
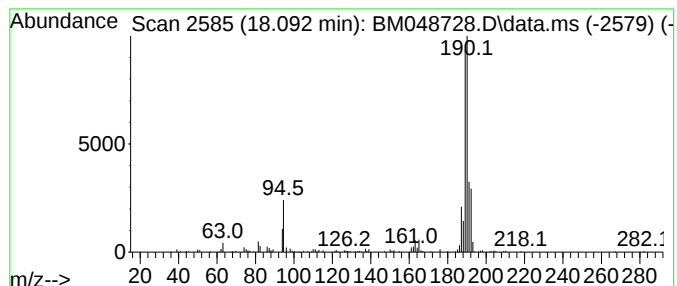
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.092	4.72 ng/ul	344846	Phenanthrene-d10	17.021

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	94
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	68
4			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	50
5			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111524\
Data File : BM048728.D
Acq On : 16 Nov 2024 04:23
Operator : RC/JU
Sample : P4751-22
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BH9G3

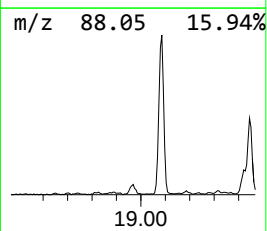
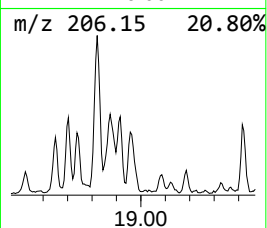
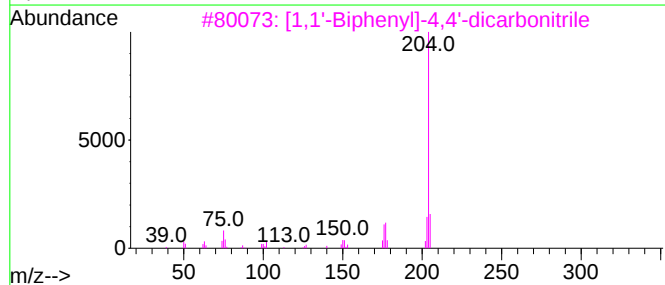
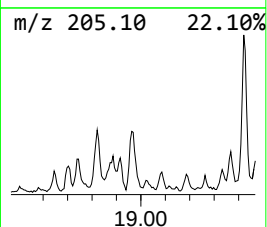
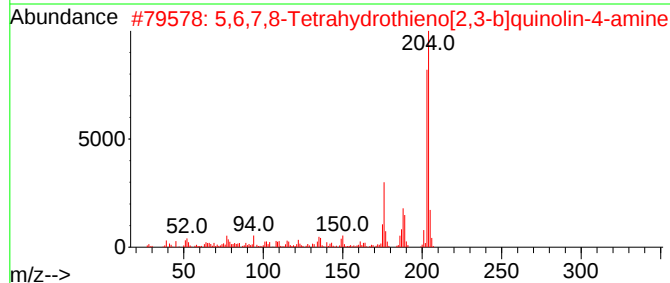
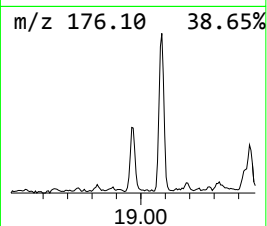
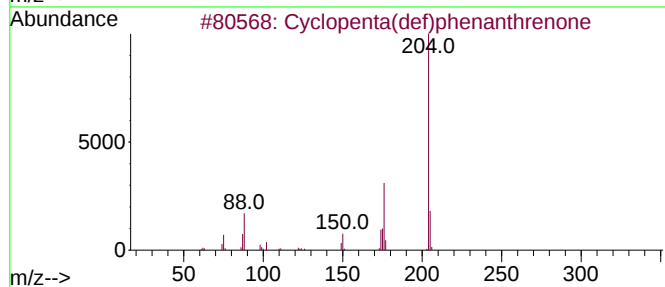
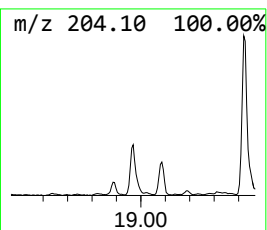
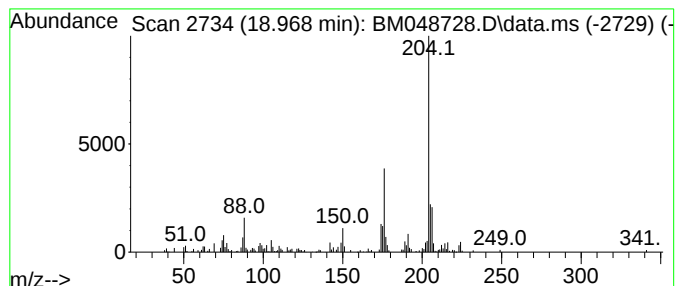
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Cyclopenta(def)phenanthrenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.968	2.48 ng/ul	181293	Phenanthrene-d10	17.021

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	96
2			5,6,7,8-Tetrahydrothieno[2,3-b]q...	204	C11H12N2S	122914-50-5	72
3			[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	50
4			Tricyclo[5.1.0.0(2,4)]octane, 5-...	247	C17H29N	1000063-59-3	47
5			[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111524\
Data File : BM048728.D
Acq On : 16 Nov 2024 04:23
Operator : RC/JU
Sample : P4751-22
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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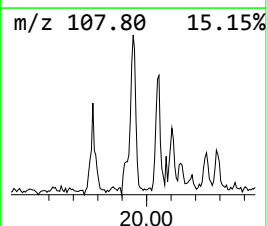
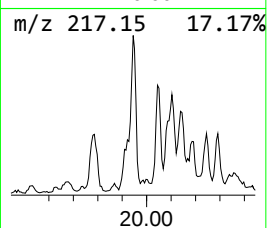
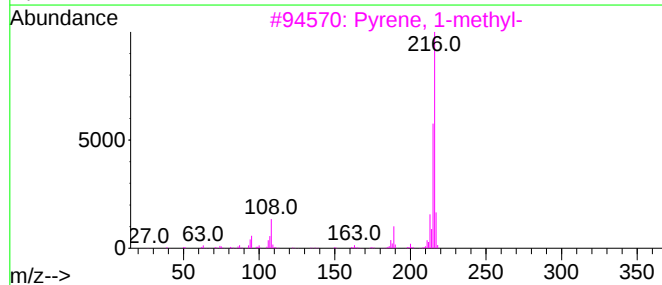
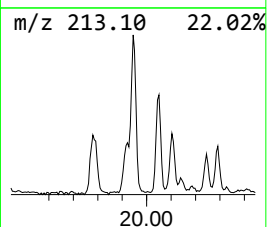
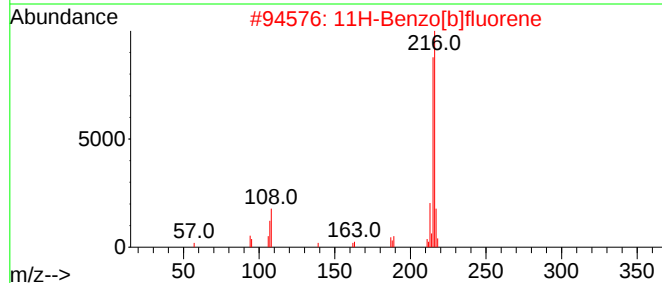
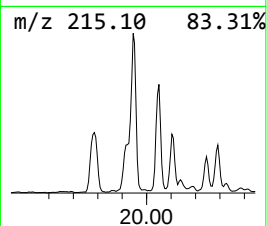
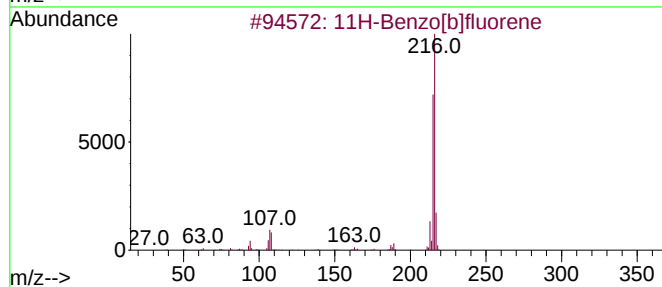
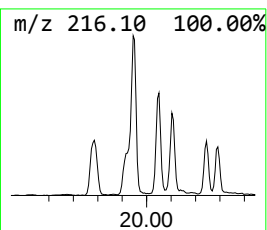
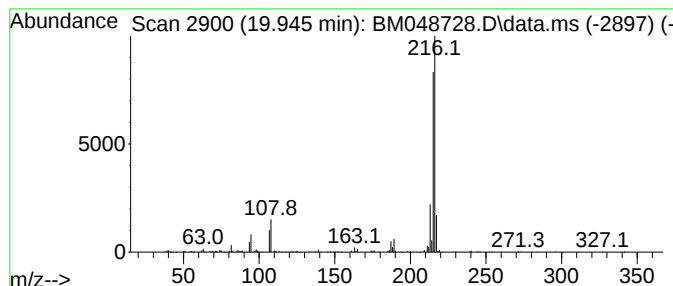
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 11H-Benzo[b]fluorene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.945	2.10 ng/ul	354061	Chrysene-d12	21.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	97
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
3			Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
4			Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
5			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111524\
Data File : BM048728.D
Acq On : 16 Nov 2024 04:23
Operator : RC/JU
Sample : P4751-22
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110724.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...	17.898	2.2	ng/u1	162695	4	17.021	1460270	20.0
n-Hexadecanoic ...	17.927	2.5	ng/u1	181957	4	17.021	1460270	20.0
Phenanthrene, 1...	17.945	3.5	ng/u1	255965	4	17.021	1460270	20.0
4H-Cyclopenta[d...	18.092	4.7	ng/u1	344846	4	17.021	1460270	20.0
Cyclopenta(def)...	18.968	2.5	ng/u1	181293	4	17.021	1460270	20.0
11H-Benzo[b]flu...	19.945	2.1	ng/u1	354061	5	21.245	3364470	20.0