

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050423\
 Data File : BP014959.D
 Acq On : 04 May 2023 18:54
 Operator : CG/JU
 Sample : 02447-08 10X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW913

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-BP050323.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014959.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.017	103	107	112	rVB	25158	33622	0.54%	0.057%
2	4.246	142	146	152	rVB	25652	33787	0.55%	0.057%
3	5.622	374	380	386	rBV	98774	147464	2.39%	0.249%
4	6.111	458	463	472	rVB	104157	156030	2.52%	0.263%
5	7.311	658	667	676	rBV2	29534	65418	1.06%	0.110%
6	7.481	691	696	705	rBV	31377	51572	0.83%	0.087%
7	7.687	726	731	741	rVB	34762	58358	0.94%	0.098%
8	8.163	802	812	821	rBV	1193264	1907711	30.87%	3.217%
9	8.875	927	933	940	rBV2	28953	52398	0.85%	0.088%
10	8.999	948	954	961	rBV2	38218	65800	1.06%	0.111%
11	9.340	1005	1012	1018	rBV	28700	51201	0.83%	0.086%
12	10.075	1130	1137	1144	rBV2	17603	34250	0.55%	0.058%
13	10.616	1221	1229	1236	rBV2	27824	53990	0.87%	0.091%
14	10.999	1282	1294	1301	rBV	1501815	2554788	41.34%	4.308%
15	11.052	1301	1303	1308	rVV	31805	41797	0.68%	0.070%
16	12.651	1569	1575	1580	rVB	21751	35583	0.58%	0.060%
17	12.863	1605	1611	1617	rBV	28089	43405	0.70%	0.073%
18	14.128	1817	1826	1831	rBV	36463	64151	1.04%	0.108%
19	14.204	1831	1839	1845	rBV2	70495	126123	2.04%	0.213%
20	14.504	1883	1890	1892	rBV	85814	132027	2.14%	0.223%
21	14.534	1892	1895	1902	rVV	85635	133012	2.15%	0.224%
22	14.810	1931	1942	1948	rVV	1973414	2860111	46.28%	4.823%
23	14.869	1948	1952	1959	rVV	144686	224898	3.64%	0.379%
24	15.204	2003	2009	2016	rVB2	76288	117296	1.90%	0.198%
25	15.422	2038	2046	2051	rBV	21562	42338	0.69%	0.071%
26	15.592	2068	2075	2078	rBV4	17844	37510	0.61%	0.063%
27	15.798	2106	2110	2114	rVV	103590	148185	2.40%	0.250%
28	15.857	2114	2120	2127	rVV	141650	231050	3.74%	0.390%
29	15.981	2138	2141	2144	rVV3	26122	37742	0.61%	0.064%
30	16.022	2144	2148	2151	rVV2	25882	40803	0.66%	0.069%
31	16.063	2151	2155	2159	rVV6	20959	35566	0.58%	0.060%
32	16.151	2165	2170	2176	rVB2	46713	80932	1.31%	0.136%
33	16.298	2190	2195	2203	rVV	42528	90857	1.47%	0.153%
34	16.528	2232	2234	2240	rVB4	29746	37615	0.61%	0.063%
35	16.863	2281	2291	2295	rVV3	41452	96249	1.56%	0.162%
36	16.910	2295	2299	2304	rVB3	28182	46924	0.76%	0.079%

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Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
 Title : SVOA CALIBRATION

37	17.039	2318	2321	2325	rVV	80537	107367	1.74%	0.181%
38	17.086	2326	2329	2333	rVV2	33809	50190	0.81%	0.085%
39	17.134	2333	2337	2338	rVV2	35633	46986	0.76%	0.079%
40	17.157	2338	2341	2346	rVB	59711	78602	1.27%	0.133%
41	17.216	2346	2351	2357	rVB	52820	81359	1.32%	0.137%
42	17.298	2359	2365	2371	rBV5	36272	88701	1.44%	0.150%
43	17.381	2374	2379	2385	rVB	86708	125244	2.03%	0.211%
44	17.563	2403	2410	2414	rBV	1835119	2664531	43.12%	4.493%
45	17.604	2414	2417	2423	rVV	1951956	2692146	43.56%	4.539%
46	17.663	2423	2427	2429	rVV2	110609	163852	2.65%	0.276%
47	17.698	2429	2433	2437	rVV	464851	656702	10.63%	1.107%
48	17.739	2437	2440	2448	rVB5	61648	118054	1.91%	0.199%
49	17.957	2471	2477	2482	rBV2	120420	208746	3.38%	0.352%
50	18.075	2493	2497	2500	rBV	41686	52811	0.85%	0.089%
51	18.139	2503	2508	2516	rVB4	152646	323102	5.23%	0.545%
52	18.269	2526	2530	2539	rVB5	44233	76911	1.24%	0.130%
53	18.416	2547	2555	2559	rBV	246568	401589	6.50%	0.677%
54	18.463	2559	2563	2571	rVB	310263	444359	7.19%	0.749%
55	18.539	2572	2576	2579	rBV	109704	165802	2.68%	0.280%
56	18.610	2583	2588	2591	rVV3	447243	804545	13.02%	1.357%
57	18.639	2591	2593	2600	rVV3	173764	347901	5.63%	0.587%
58	18.704	2600	2604	2609	rVB2	100095	213081	3.45%	0.359%
59	18.898	2629	2637	2640	rBV	185775	332001	5.37%	0.560%
60	18.939	2640	2644	2647	rVB2	80644	114120	1.85%	0.192%
61	19.016	2653	2657	2659	rBV2	36165	45880	0.74%	0.077%
62	19.039	2659	2661	2667	rVB2	43516	49569	0.80%	0.084%
63	19.133	2674	2677	2682	rVB3	65642	79913	1.29%	0.135%
64	19.181	2682	2685	2688	rBV	50587	52189	0.84%	0.088%
65	19.222	2689	2692	2698	rVB2	62016	90806	1.47%	0.153%
66	19.298	2700	2705	2709	rBV	146847	247909	4.01%	0.418%
67	19.363	2709	2716	2718	rVV6	101796	216747	3.51%	0.365%
68	19.386	2718	2720	2723	rVV	115932	137441	2.22%	0.232%
69	19.433	2723	2728	2736	rVB3	161617	326363	5.28%	0.550%
70	19.557	2743	2749	2755	rBV	4004378	5119675	82.84%	8.633%
71	19.610	2756	2758	2761	rVV	52905	71307	1.15%	0.120%
72	19.645	2761	2764	2768	rVV	93658	122745	1.99%	0.207%
73	19.686	2768	2771	2776	rVV3	74501	108660	1.76%	0.183%
74	19.792	2785	2789	2798	rVB	121719	233379	3.78%	0.394%
75	19.880	2798	2804	2805	rBV2	659408	953992	15.44%	1.609%
76	19.910	2805	2809	2816	rVV	3627419	4644115	75.15%	7.831%

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 Title : SVOA CALIBRATION

77	19.975	2816	2820	2825	rVB2	139653	208309	3.37%	0.351%
78	20.080	2834	2838	2841	rVB	137276	171026	2.77%	0.288%
79	20.139	2842	2848	2854	rVB2	128474	272065	4.40%	0.459%
80	20.222	2856	2862	2868	rBV3	194569	482787	7.81%	0.814%
81	20.386	2881	2890	2894	rVB2	429269	817298	13.23%	1.378%
82	20.480	2903	2906	2910	rBV	241586	283777	4.59%	0.479%
83	20.539	2910	2916	2920	rVV2	271796	429491	6.95%	0.724%
84	20.575	2920	2922	2926	rVB	98834	104052	1.68%	0.175%
85	20.675	2936	2939	2943	rBV	175667	248430	4.02%	0.419%
86	20.722	2944	2947	2950	rVB	194538	219522	3.55%	0.370%
87	21.116	3010	3014	3020	rBV	310091	386775	6.26%	0.652%
88	21.280	3039	3042	3045	rVB	243073	257900	4.17%	0.435%
89	21.316	3045	3048	3052	rVB	395139	431140	6.98%	0.727%
90	21.357	3052	3055	3060	rBV2	305059	463059	7.49%	0.781%
91	21.404	3060	3063	3072	rVB2	306633	455264	7.37%	0.768%
92	21.527	3081	3084	3088	rVB	211097	247681	4.01%	0.418%
93	21.633	3096	3102	3108	rBV2	2629643	6179924	100.00%	10.421%
94	21.680	3108	3110	3121	rVB	1804079	2352370	38.06%	3.967%
95	21.816	3129	3133	3138	rBV2	508187	751324	12.16%	1.267%
96	21.874	3140	3143	3147	rVV2	344270	449761	7.28%	0.758%
97	23.445	3403	3410	3415	rBV	1423968	3643544	58.96%	6.144%
98	24.010	3501	3506	3513	rVB	850972	1674339	27.09%	2.823%
99	24.127	3520	3526	3535	rVB	1365559	2816357	45.57%	4.749%
100	24.245	3539	3546	3552	rBV2	1438557	3129001	50.63%	5.276%

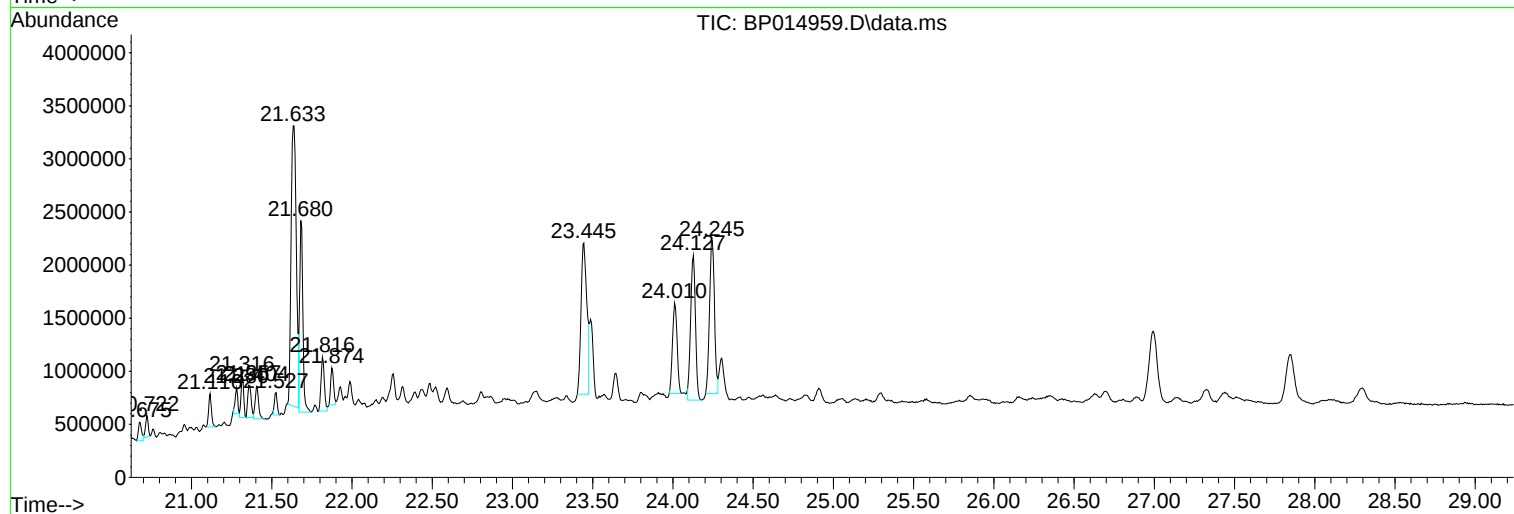
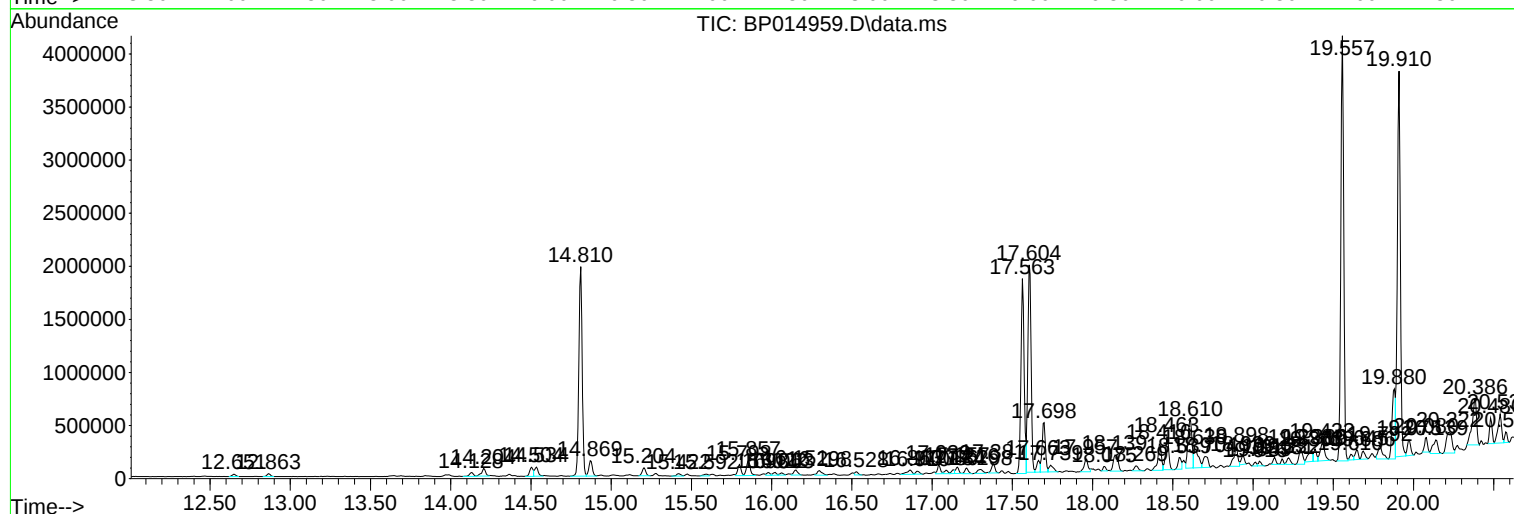
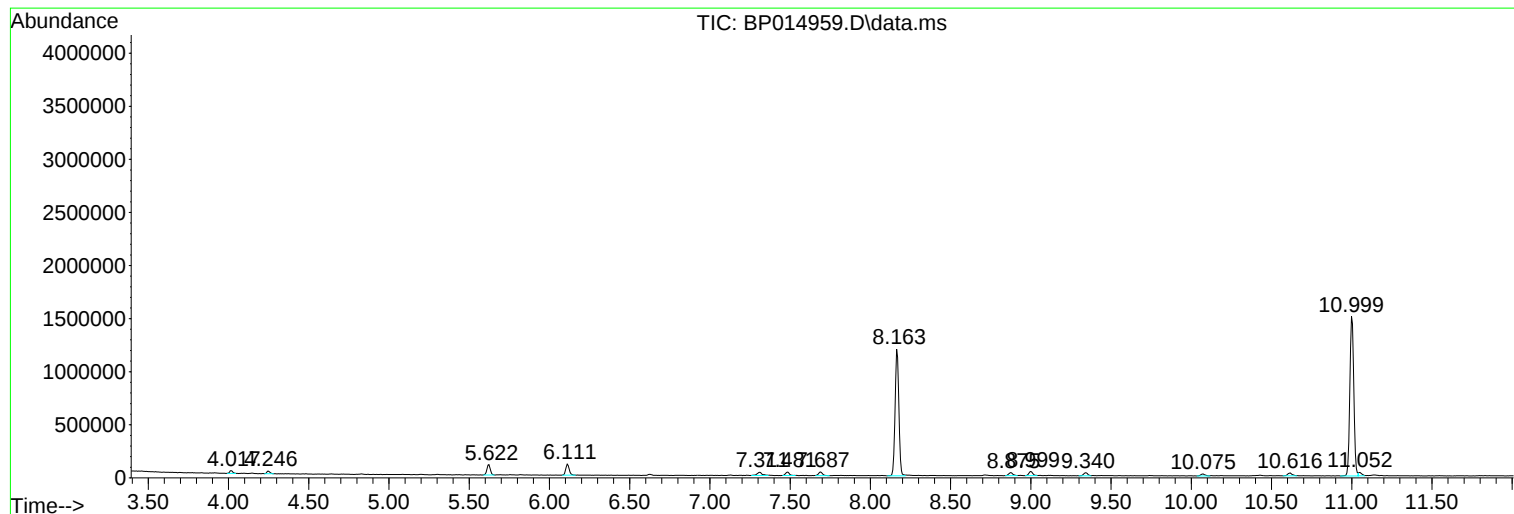
Sum of corrected areas: 59305151

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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



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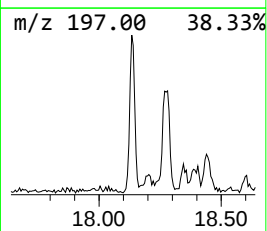
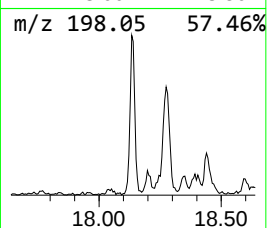
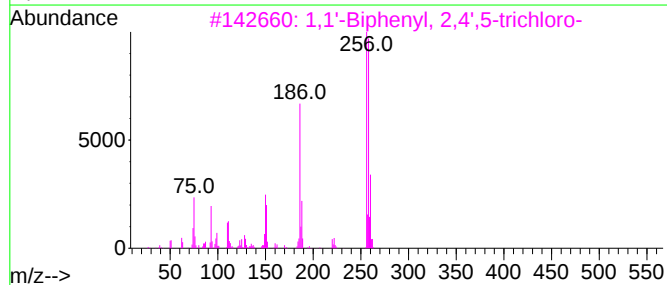
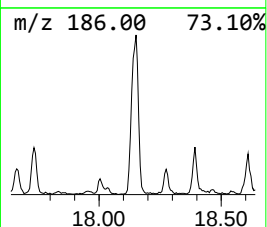
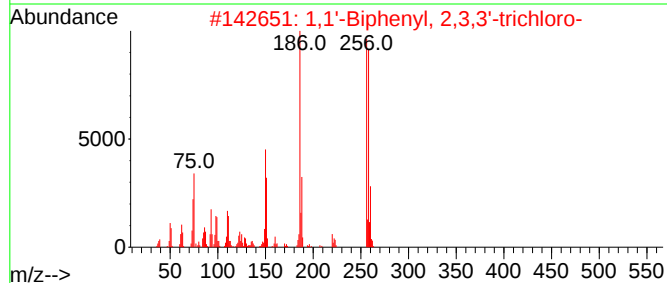
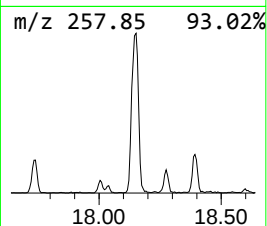
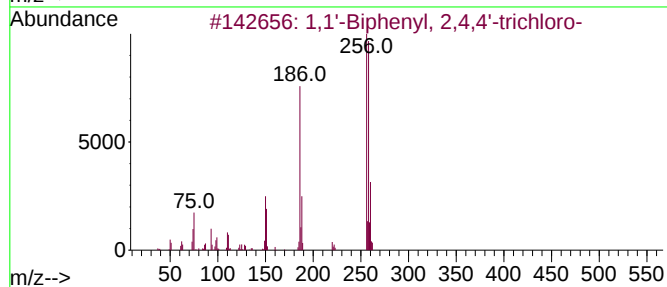
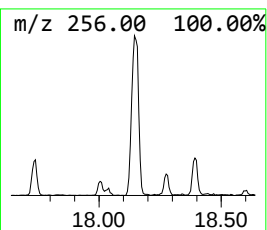
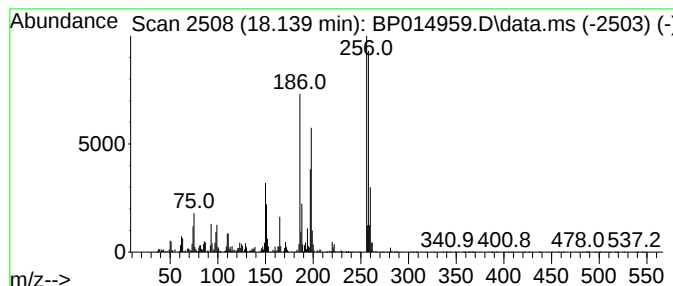
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1,1'-Biphenyl, 2,4,4'-trich... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.139	2.43 ng/ul	323102	Phenanthrene-d10	17.563

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99	
2	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99	
3	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	99	
4	1,1'-Biphenyl, 2,3,4'-Trichloro-	256	C12H7Cl3	038444-85-8	98	
5	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	98	



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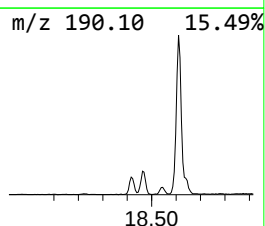
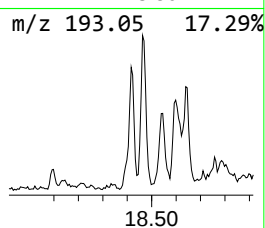
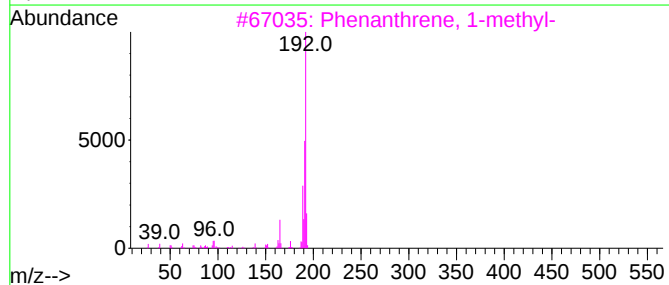
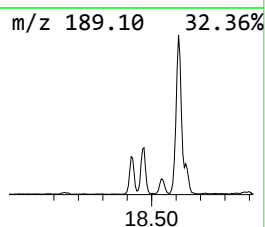
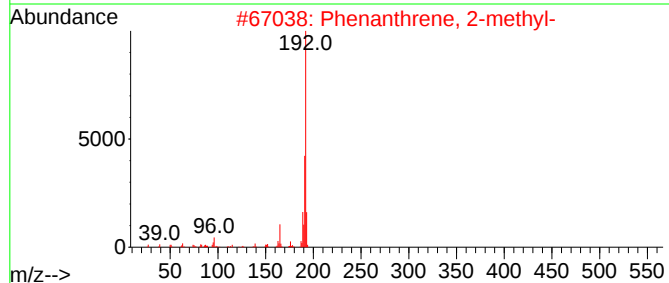
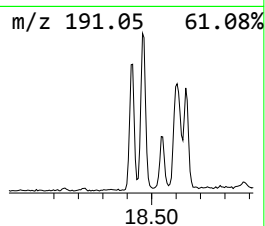
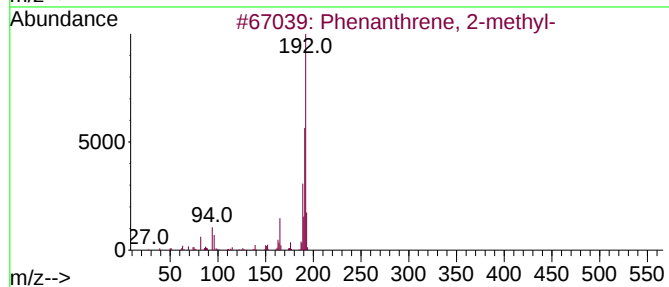
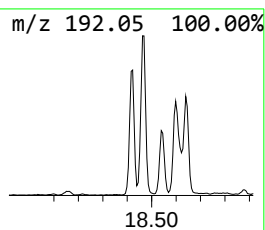
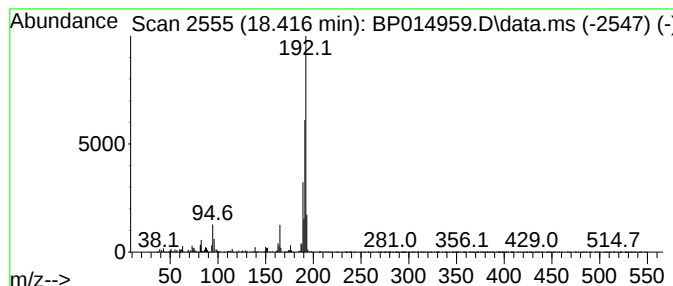
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.416	3.01 ng/ul	401589	Phenanthrene-d10	17.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
4			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
5			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	94



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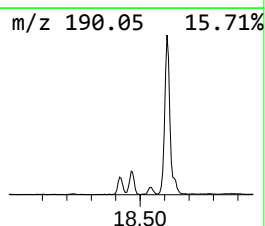
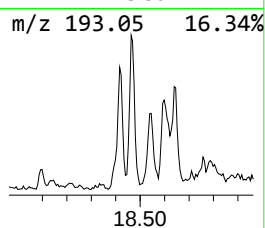
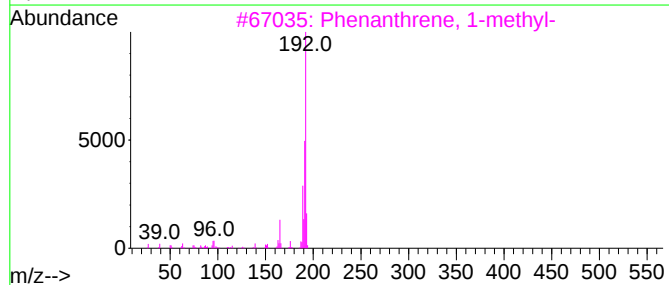
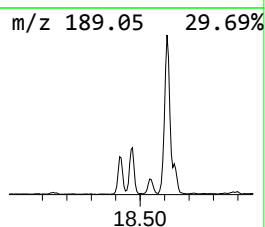
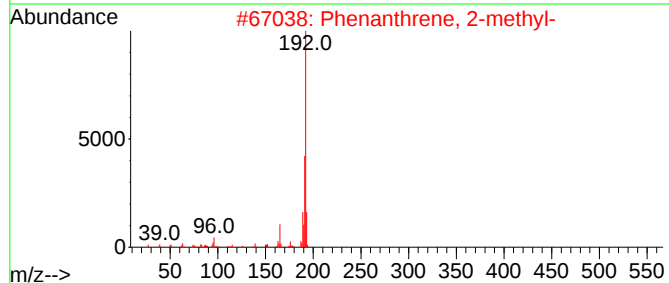
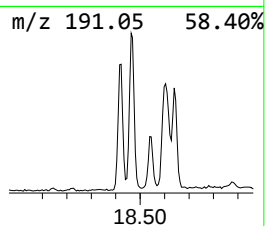
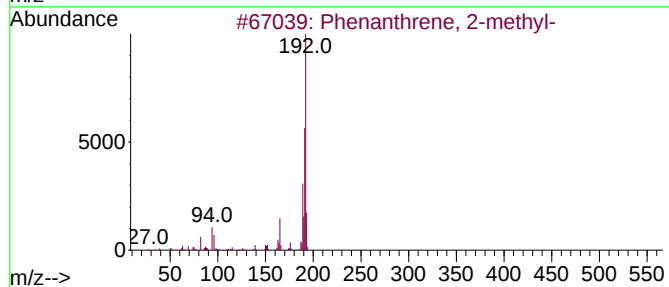
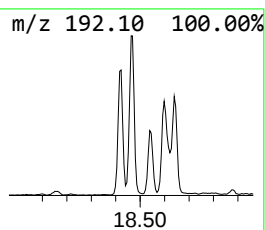
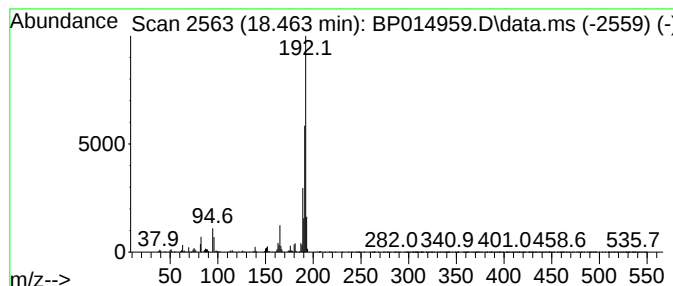
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Phenanthrene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.463	3.34 ng/ul	444359	Phenanthrene-d10	17.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
4			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
5			5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050423\
Data File : BP014959.D
Acq On : 04 May 2023 18:54
Operator : CG/JU
Sample : 02447-08 10X
Misc :
ALS Vial : 18 Sample Multiplier: 1

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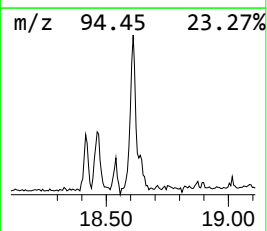
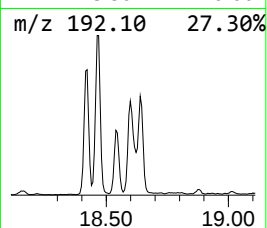
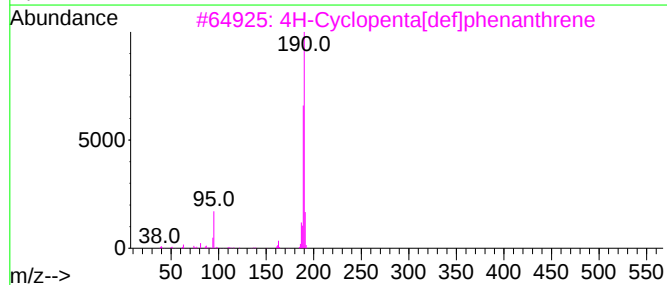
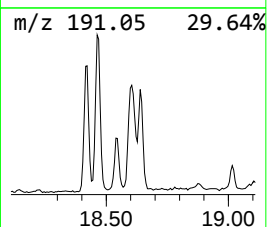
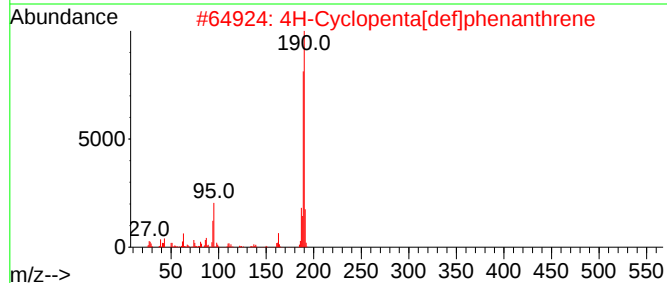
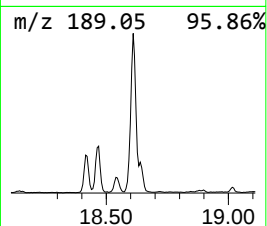
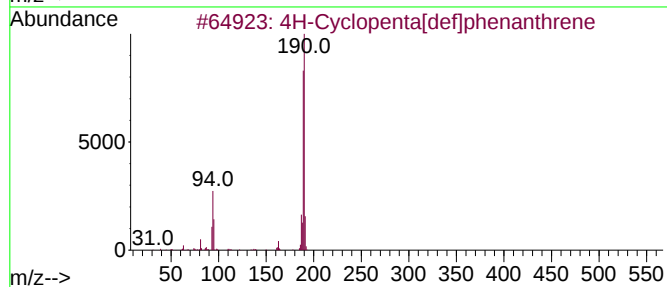
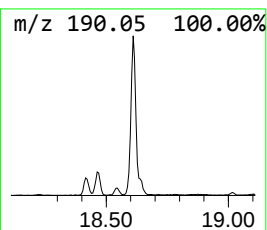
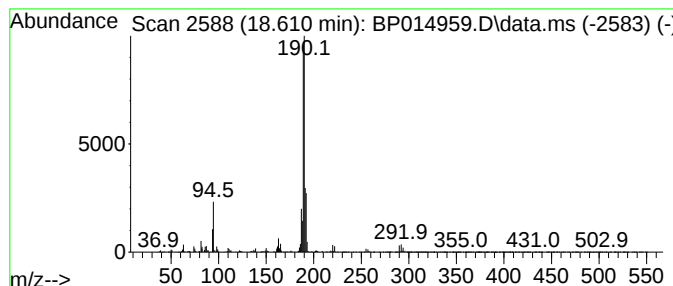
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.610	6.04 ng/ul	804545	Phenanthrene-d10	17.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
4			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	50
5			2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050423\
Data File : BP014959.D
Acq On : 04 May 2023 18:54
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ClientSampleId :
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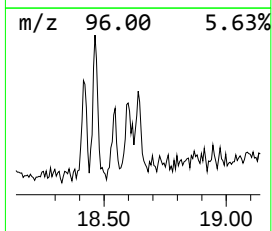
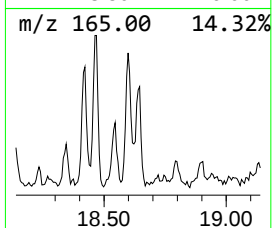
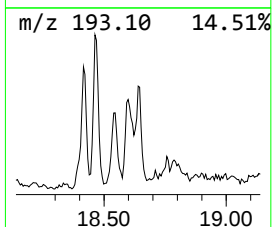
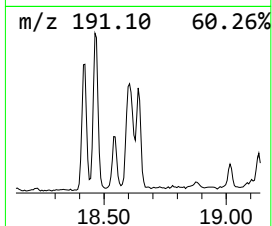
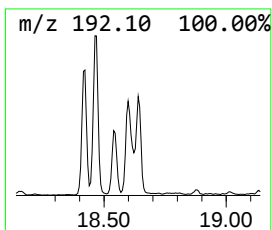
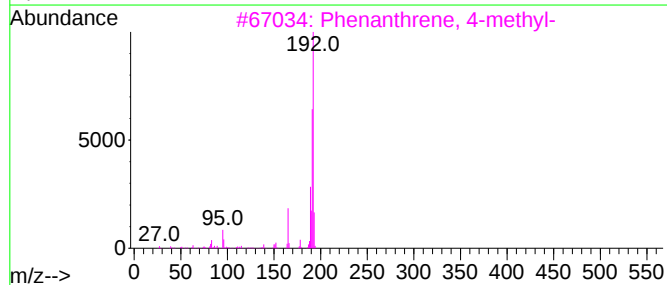
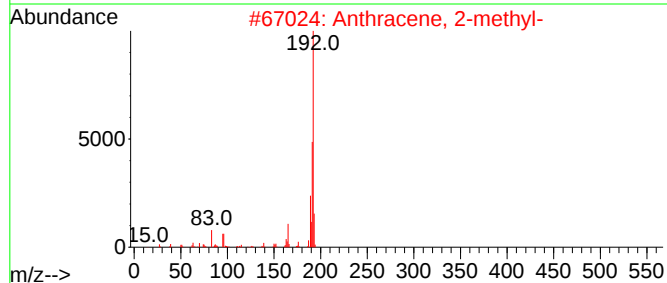
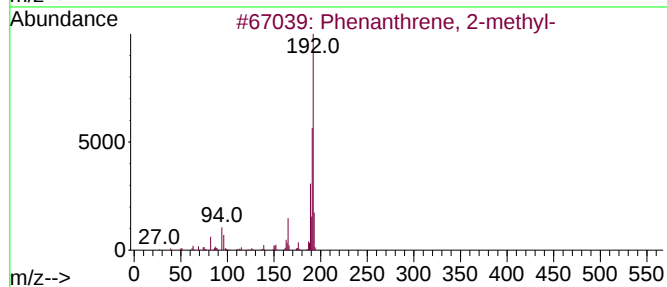
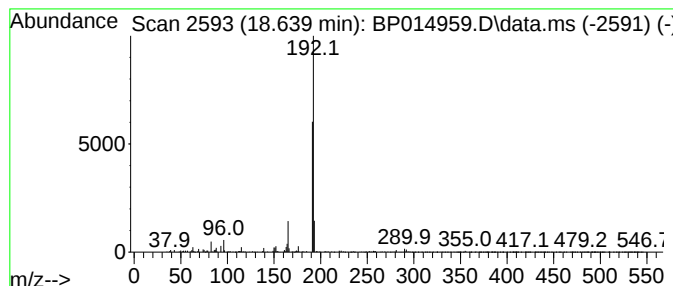
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Anthracene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.639	2.61 ng/ul	347901	Phenanthrene-d10	17.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
3			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	90
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	87
5			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050423\
Data File : BP014959.D
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Misc :
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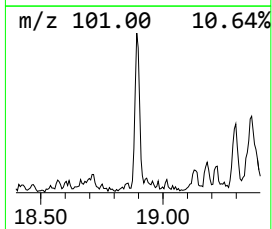
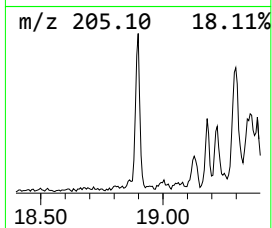
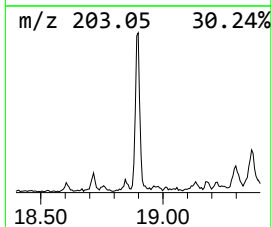
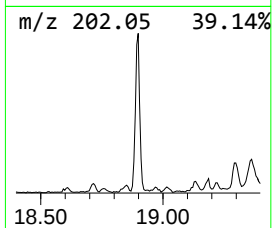
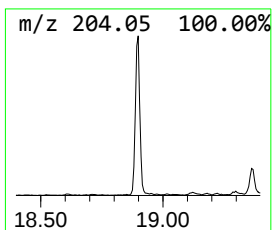
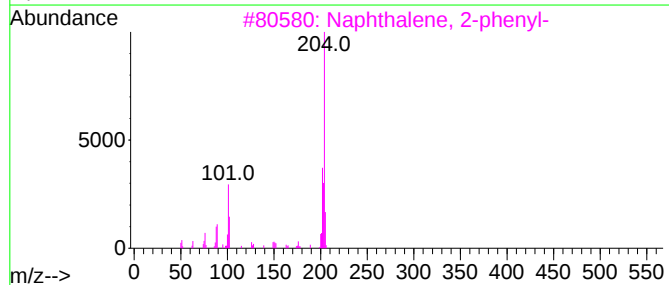
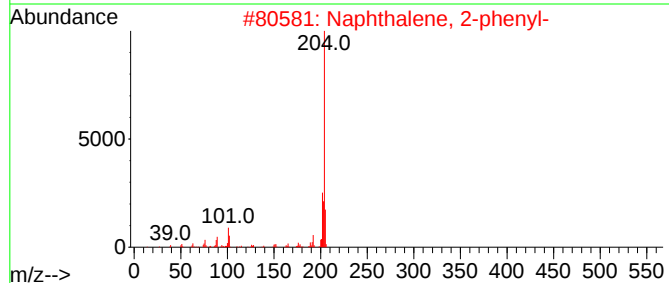
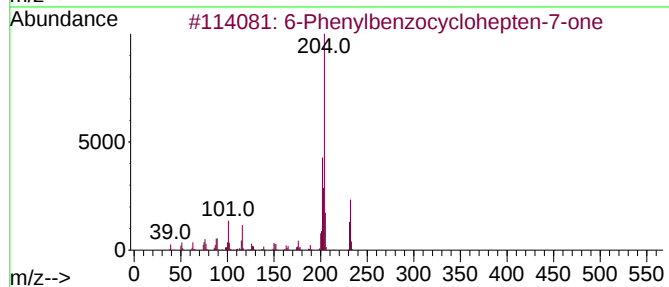
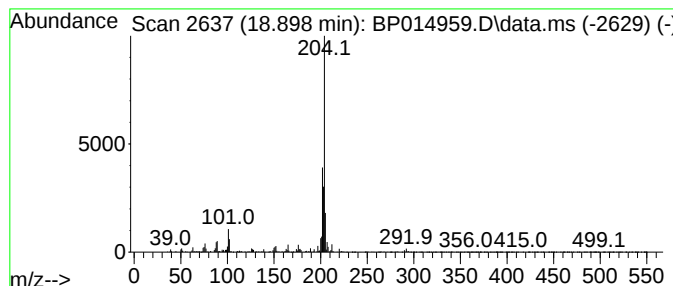
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 6-Phenylbenzocyclohepten-7-one Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.898	2.49 ng/ul	332001	Phenanthrene-d10	17.563

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	87
2		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	83
3		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	83
4		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	78
5		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	64



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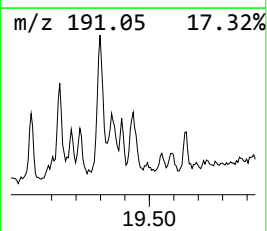
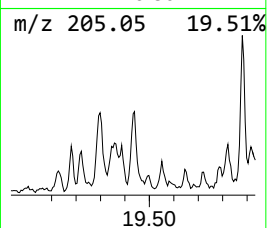
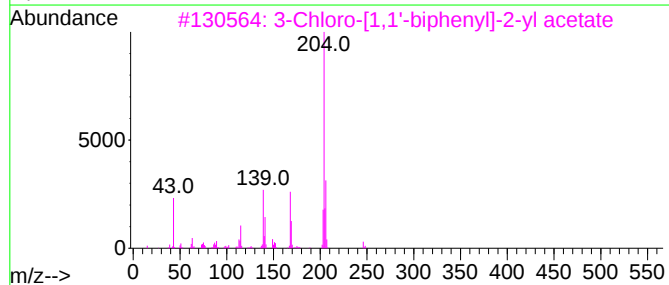
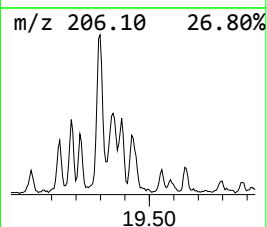
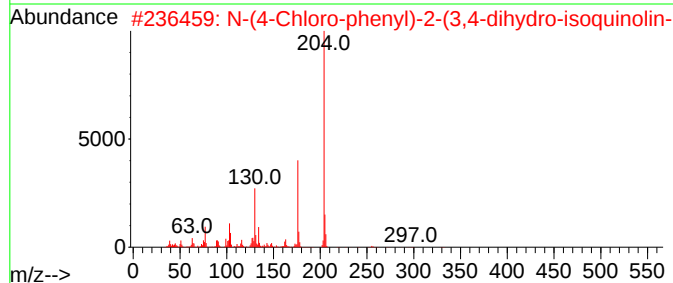
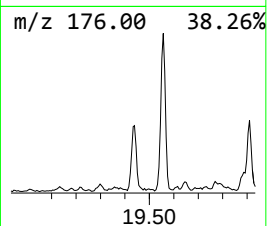
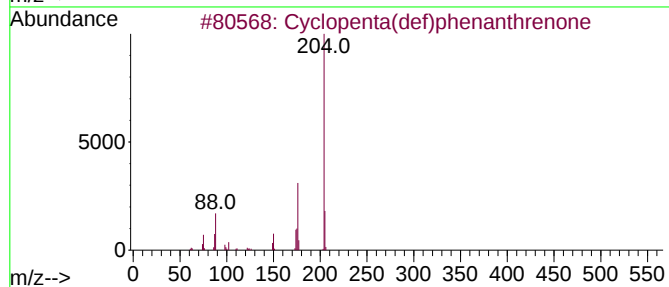
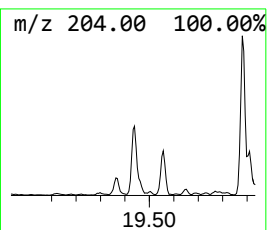
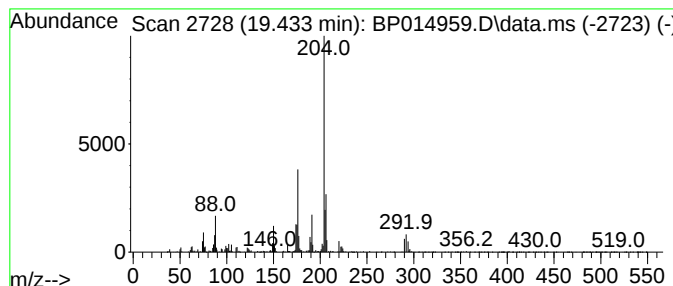
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Cyclopenta(def)phenanthrenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.433	2.45 ng/ul	326363	Phenanthrene-d10	17.563	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	96
2	N-(4-Chloro-phenyl)-2-(3,4-dihyd...	330	C17H15ClN2OS	1010296-34-3	59
3	3-Chloro-[1,1'-biphenyl]-2-yl ac...	246	C14H11ClO2	007460-70-0	47
4	2-Hydroxy-5-chlorobiphenyl, acetate	246	C14H11ClO2	1000447-68-0	47
5	L-Rhamnose, (R,R,S,S)-, 4TMS der...	452	C18H44O5Si4	108392-01-4	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050423\
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ClientSampleId :
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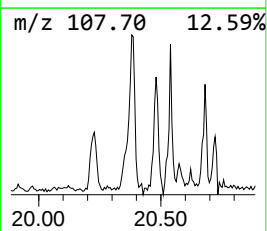
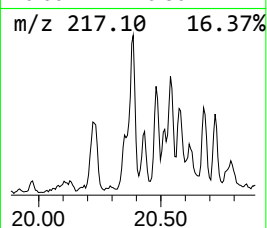
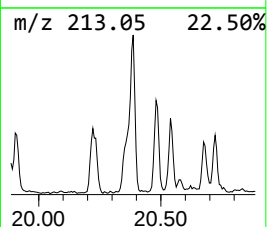
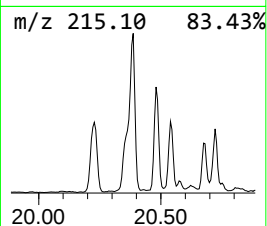
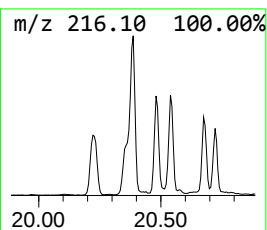
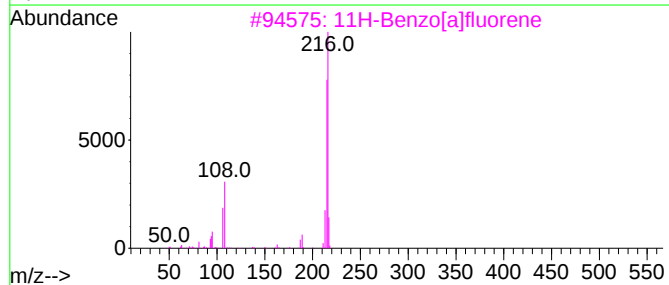
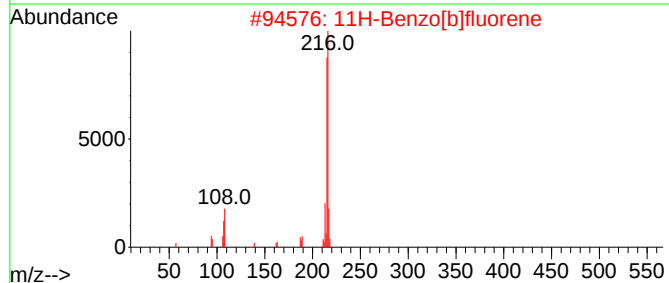
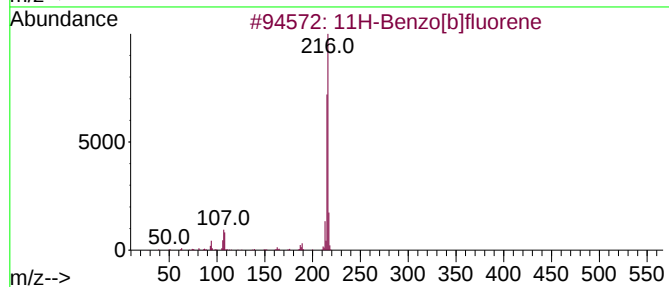
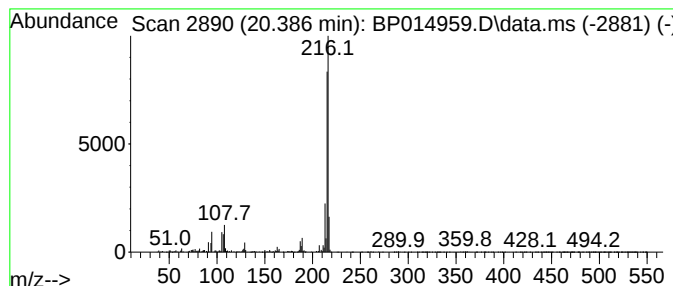
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 11H-Benzo[b]fluorene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.386	2.65 ng/ul	817298	Chrysene-d12	21.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
5			Pyrene, 1-methyl-	216	C17H12	002381-21-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050423\
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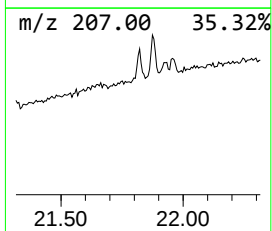
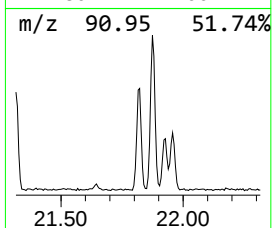
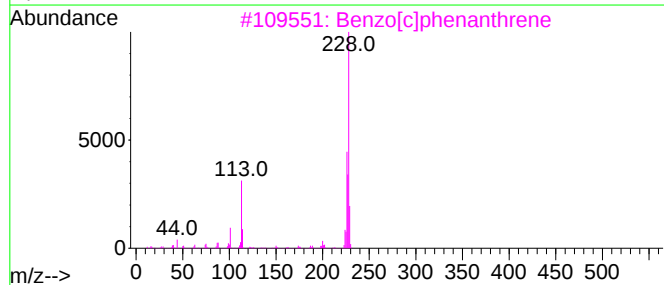
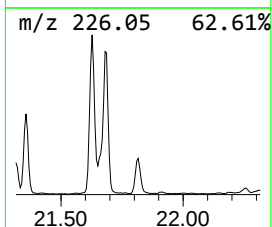
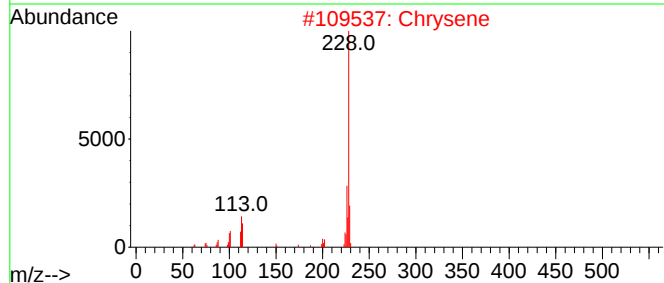
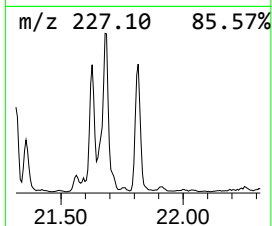
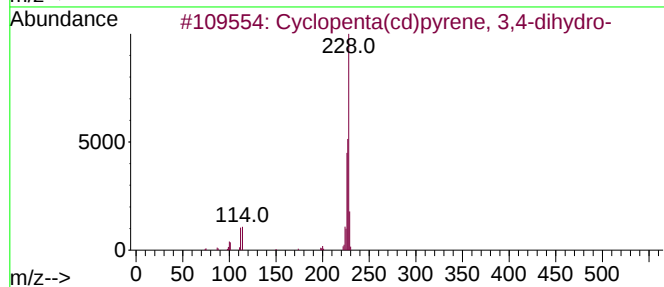
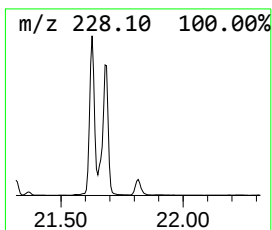
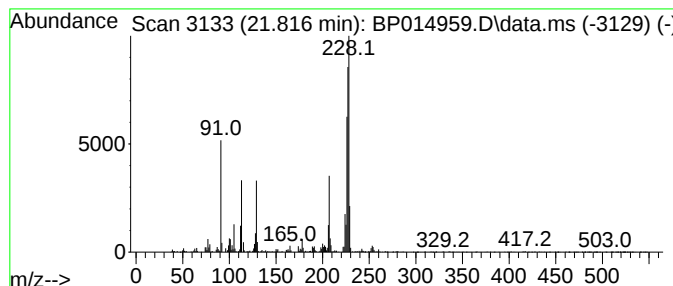
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Cyclopenta(cd)pyrene, 3,4-d... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.816	2.43 ng/ul	751324	Chrysene-d12	21.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(cd)pyrene, 3,4-dihydro-	228	C18H12	025732-74-5	89
2			Chrysene	228	C18H12	000218-01-9	80
3			Benzo[c]phenanthrene	228	C18H12	000195-19-7	60
4			Benzo[c]phenanthrene	228	C18H12	000195-19-7	55
5			5-Phenylthieno[2,3-d]pyrimidin-4-ol	228	C12H8N2OS	035978-39-3	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050423\
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ClientSampleId :
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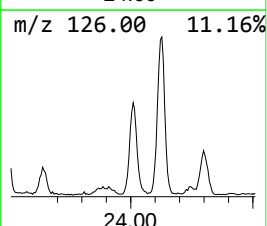
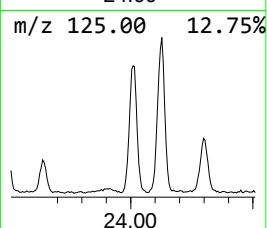
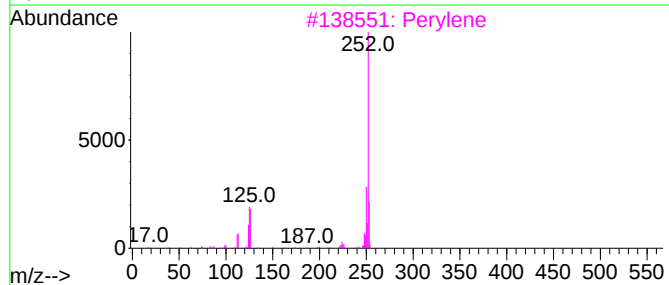
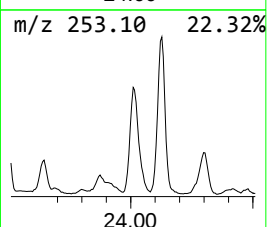
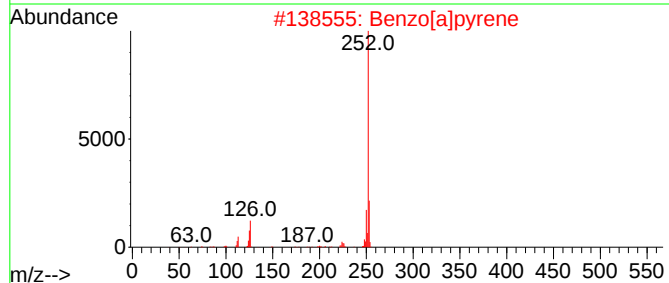
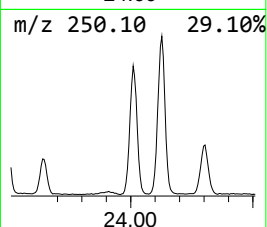
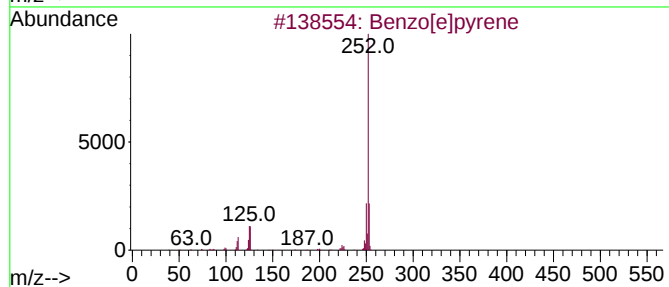
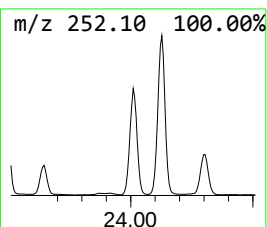
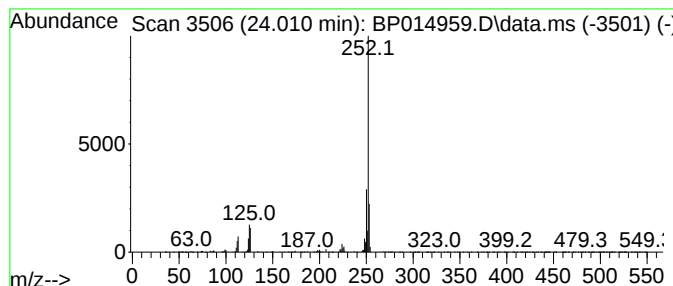
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.010	10.70 ng/ul	1674340	Perylene-d12	24.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	99
2			Benzo[a]pyrene	252	C20H12	000050-32-8	97
3			Perylene	252	C20H12	000198-55-0	97
4			Benzo[k]fluoranthene	252	C20H12	000207-08-9	96
5			Perylene	252	C20H12	000198-55-0	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050423\
Data File : BP014959.D
Acq On : 04 May 2023 18:54
Operator : CG/JU
Sample : 02447-08 10X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW913

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1,1'-Biphenyl, ...	18.139	2.4	ng/u1	323102	4	17.563	2664530	20.0
Phenanthrene, 2...	18.416	3.0	ng/u1	401589	4	17.563	2664530	20.0
Phenanthrene, 1...	18.463	3.3	ng/u1	444359	4	17.563	2664530	20.0
4H-Cyclopenta[d...	18.610	6.0	ng/u1	804545	4	17.563	2664530	20.0
Anthracene, 2-m...	18.639	2.6	ng/u1	347901	4	17.563	2664530	20.0
6-Phenylbenzocy...	18.898	2.5	ng/u1	332001	4	17.563	2664530	20.0
Cyclopenta(def)...	19.433	2.5	ng/u1	326363	4	17.563	2664530	20.0
11H-Benzo[b]flu...	20.386	2.6	ng/u1	817298	5	21.645	6179920	20.0
Cyclopenta(cd)p...	21.816	2.4	ng/u1	751324	5	21.645	6179920	20.0
Benzo[e]pyrene	24.010	10.7	ng/u1	1674340	6	24.245	3129000	20.0