

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102524\
 Data File : BP022621.D
 Acq On : 25 Oct 2024 22:01
 Operator : RC/JU
 Sample : P4436-05
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4F47

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-BP100724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP022621.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.605	101	105	118	rBV	112551	184494	3.99%	0.340%
2	6.840	650	655	669	rVV	194002	348482	7.55%	0.642%
3	6.993	673	681	692	rVV	128716	208580	4.52%	0.384%
4	7.193	707	715	724	rBV	181080	309147	6.69%	0.569%
5	7.646	785	792	801	rBV	330855	550473	11.92%	1.013%
6	8.357	905	913	931	rBV	241232	447337	9.69%	0.824%
7	8.793	979	987	996	rBV	172180	301220	6.52%	0.555%
8	9.351	1073	1082	1089	rBV2	24913	45221	0.98%	0.083%
9	9.504	1100	1108	1122	rBV	151381	284191	6.15%	0.523%
10	10.040	1191	1199	1213	rBV	253320	486170	10.53%	0.895%
11	10.404	1249	1261	1266	rBV	440874	767529	16.62%	1.413%
12	10.451	1266	1269	1277	rVB	63226	106991	2.32%	0.197%
13	10.546	1277	1285	1301	rBV	198182	374960	8.12%	0.690%
14	12.069	1538	1544	1550	rBV	56125	84763	1.84%	0.156%
15	12.293	1575	1582	1591	rVB2	40267	73738	1.60%	0.136%
16	13.087	1712	1717	1724	rBV	22662	38431	0.83%	0.071%
17	13.422	1763	1774	1775	rBV	22349	32565	0.71%	0.060%
18	13.581	1794	1801	1806	rBV	45285	78461	1.70%	0.144%
19	13.669	1811	1816	1825	rVB	538568	808261	17.50%	1.488%
20	13.957	1856	1865	1878	rBV	558727	881024	19.08%	1.622%
21	14.269	1907	1918	1924	rBV	745421	1184664	25.65%	2.181%
22	14.334	1924	1929	1937	rVV	201704	330768	7.16%	0.609%
23	14.504	1948	1958	1966	rBV	171085	342093	7.41%	0.630%
24	14.575	1968	1970	1979	rVV	20779	39747	0.86%	0.073%
25	14.669	1979	1986	1996	rVV	189741	308020	6.67%	0.567%
26	15.269	2081	2088	2094	rVV	782935	1237258	26.79%	2.278%
27	15.328	2094	2098	2108	rVV	277899	459865	9.96%	0.847%
28	15.422	2108	2114	2118	rVV	230015	334720	7.25%	0.616%
29	15.457	2118	2120	2123	rVV	31610	42641	0.92%	0.079%
30	15.492	2123	2126	2132	rVV2	38794	65310	1.41%	0.120%
31	15.645	2145	2152	2159	rVB2	163154	319696	6.92%	0.589%
32	15.786	2171	2176	2184	rVV	77408	152932	3.31%	0.282%
33	15.869	2184	2190	2197	rVB2	19159	38278	0.83%	0.070%
34	16.339	2262	2270	2271	rBV2	31965	46639	1.01%	0.086%
35	16.363	2271	2274	2280	rVV4	44928	79183	1.71%	0.146%
36	16.592	2305	2313	2317	rBV4	26925	50890	1.10%	0.094%

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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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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Title : SVOA CALIBRATION

37	16.722	2330	2335	2344	rVB	59388	102588	2.22%	0.189%
38	16.804	2344	2349	2356	rVV3	33705	79971	1.73%	0.147%
39	16.886	2357	2363	2372	rVB	149441	228539	4.95%	0.421%
40	17.075	2388	2395	2398	rBV	1070405	1545985	33.47%	2.846%
41	17.122	2398	2403	2408	rVV	2730994	4133077	89.49%	7.609%
42	17.180	2408	2413	2416	rVV	995688	1637140	35.45%	3.014%
43	17.210	2416	2418	2424	rVV	781637	928273	20.10%	1.709%
44	17.275	2424	2429	2434	rVB2	30310	59204	1.28%	0.109%
45	17.498	2458	2467	2480	rBV	243702	388581	8.41%	0.715%
46	17.604	2481	2485	2493	rVB2	37331	58109	1.26%	0.107%
47	17.675	2493	2497	2505	rVB3	31292	62440	1.35%	0.115%
48	17.980	2543	2549	2552	rBV	203763	286203	6.20%	0.527%
49	18.028	2552	2557	2565	rVB2	383027	640540	13.87%	1.179%
50	18.122	2567	2573	2577	rVV	115409	154915	3.35%	0.285%
51	18.180	2577	2583	2587	rVV3	347702	601327	13.02%	1.107%
52	18.222	2587	2590	2598	rVB	127160	173922	3.77%	0.320%
53	18.304	2599	2604	2607	rBV4	20629	39877	0.86%	0.073%
54	18.457	2625	2630	2634	rBV6	17581	34441	0.75%	0.063%
55	18.504	2634	2638	2642	rVV	206698	273236	5.92%	0.503%
56	18.545	2642	2645	2650	rVB	107476	146835	3.18%	0.270%
57	18.763	2678	2682	2686	rBV2	44737	59336	1.28%	0.109%
58	18.816	2687	2691	2694	rBV	36291	43349	0.94%	0.080%
59	18.857	2694	2698	2701	rBV2	37675	42273	0.92%	0.078%
60	18.933	2706	2711	2713	rBV2	70215	110670	2.40%	0.204%
61	18.969	2714	2717	2720	rVV4	25192	34727	0.75%	0.064%
62	19.086	2732	2737	2745	rVB4	58639	135141	2.93%	0.249%
63	19.210	2751	2758	2765	rBV	3465410	4618461	100.00%	8.503%
64	19.275	2766	2769	2772	rVV2	40821	55993	1.21%	0.103%
65	19.316	2772	2776	2778	rVV	56418	75984	1.65%	0.140%
66	19.345	2778	2781	2788	rVB6	61576	103008	2.23%	0.190%
67	19.469	2798	2802	2806	rVB	85185	114109	2.47%	0.210%
68	19.563	2812	2818	2820	rBV2	1976407	2784792	60.30%	5.127%
69	19.592	2820	2823	2831	rVV	2697141	3376596	73.11%	6.216%
70	19.663	2831	2835	2842	rVB2	101984	174587	3.78%	0.321%
71	19.780	2851	2855	2860	rVB	138252	192077	4.16%	0.354%
72	19.851	2863	2867	2872	rVB	99202	138919	3.01%	0.256%
73	19.927	2874	2880	2881	rBV2	129895	194853	4.22%	0.359%
74	19.986	2887	2890	2893	rBV	52470	58948	1.28%	0.109%
75	20.069	2899	2904	2907	rBV4	90313	175160	3.79%	0.322%
76	20.116	2907	2912	2916	rVB	321827	469273	10.16%	0.864%

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Integrator: RTE

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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

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77	20.216	2925	2929	2933	rBV	285058	375025	8.12%	0.690%
78	20.280	2933	2940	2943	rVV2	161804	282534	6.12%	0.520%
79	20.357	2950	2953	2958	rBV3	42743	64748	1.40%	0.119%
80	20.416	2960	2963	2967	rVV	77584	106759	2.31%	0.197%
81	20.469	2968	2972	2981	rVB3	79742	180128	3.90%	0.332%
82	20.904	3042	3046	3051	rVB	133831	188558	4.08%	0.347%
83	21.080	3072	3076	3080	rBV2	221461	307326	6.65%	0.566%
84	21.121	3080	3083	3087	rVB	165588	210417	4.56%	0.387%
85	21.186	3087	3094	3100	rBV3	239314	568422	12.31%	1.046%
86	21.333	3114	3119	3123	rBV	170866	284999	6.17%	0.525%
87	21.392	3127	3129	3135	rVV2	182107	250928	5.43%	0.462%
88	21.504	3137	3148	3153	rVV2	1716513	4329397	93.74%	7.970%
89	21.557	3153	3157	3163	rVB	951152	1463828	31.70%	2.695%
90	21.716	3180	3184	3193	rBV	166589	270701	5.86%	0.498%
91	22.263	3274	3277	3282	rVB	136458	213216	4.62%	0.393%
92	22.363	3292	3294	3298	rVB2	74714	89513	1.94%	0.165%
93	23.757	3522	3531	3536	rBV	618119	1855529	40.18%	3.416%
94	23.910	3553	3557	3566	rVB2	255018	544320	11.79%	1.002%
95	24.004	3569	3573	3582	rVB	155060	349269	7.56%	0.643%
96	24.486	3648	3655	3660	rBV	279227	646510	14.00%	1.190%
97	24.562	3660	3668	3673	rVV	793524	1887605	40.87%	3.475%
98	24.621	3673	3678	3694	rVB	604606	1500370	32.49%	2.762%
99	24.774	3696	3704	3713	rBV	800554	2131482	46.15%	3.924%
100	28.480	4329	4334	4335	rBV	154562	238857	5.17%	0.440%

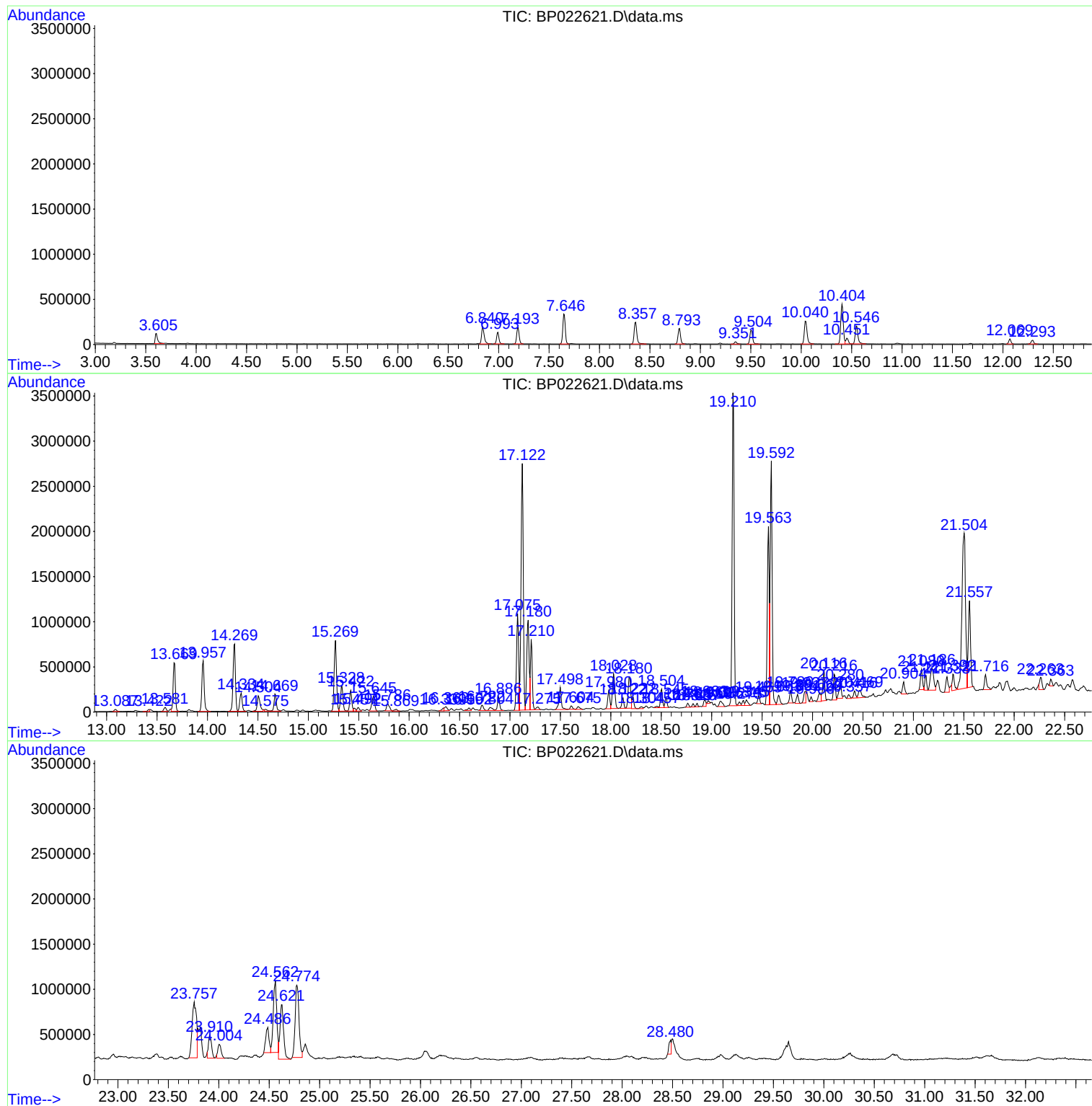
Sum of corrected areas: 54318642

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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



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Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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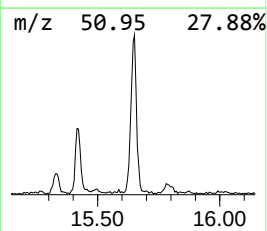
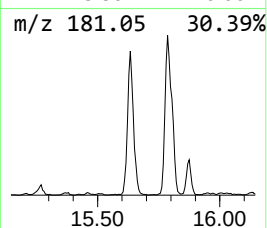
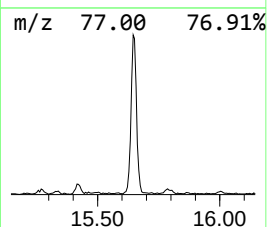
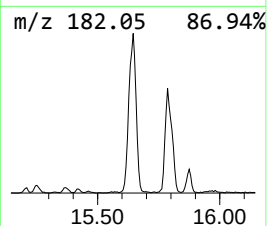
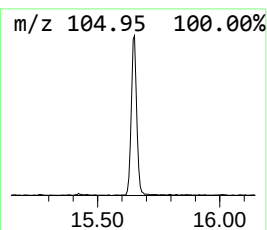
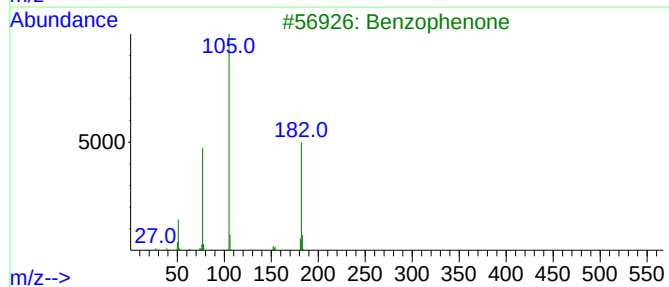
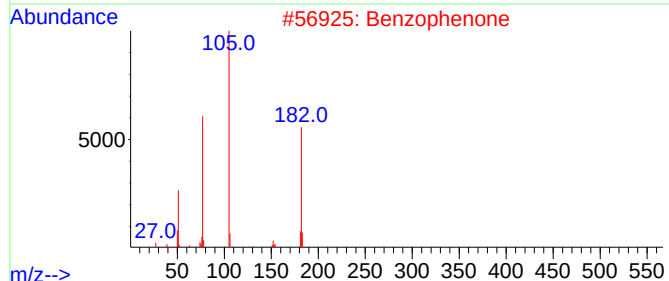
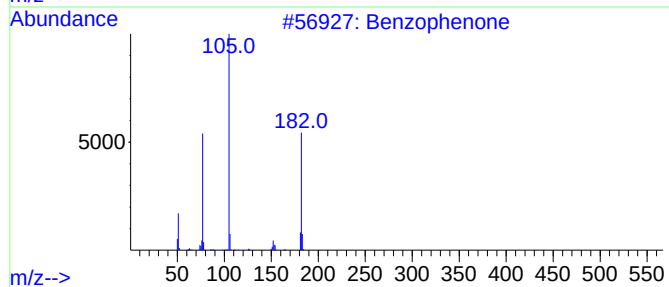
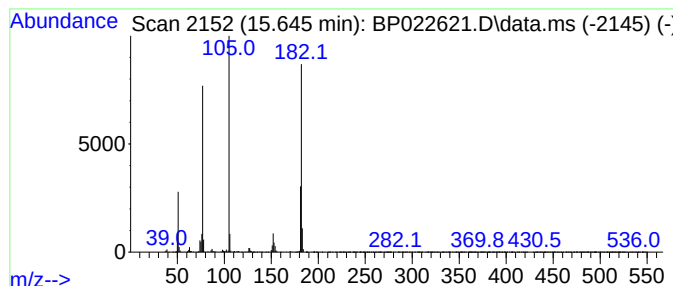
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.645	5.40 ng/ul	319696	Acenaphthene-d10	14.269

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	95
2			Benzophenone	182	C13H10O	000119-61-9	90
3			Benzophenone	182	C13H10O	000119-61-9	87
4			Benzophenone	182	C13H10O	000119-61-9	76
5			Benzophenone	182	C13H10O	000119-61-9	68



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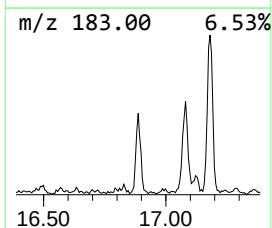
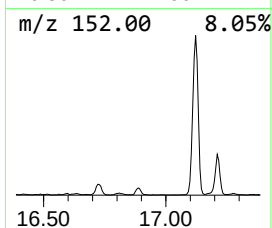
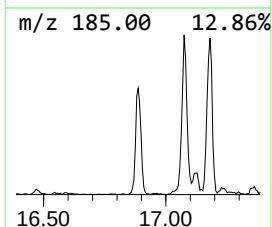
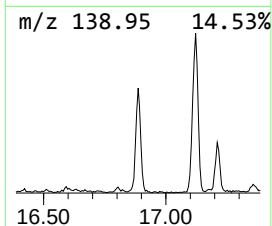
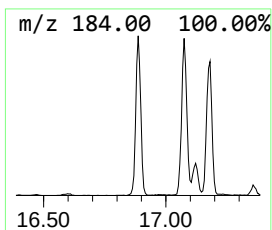
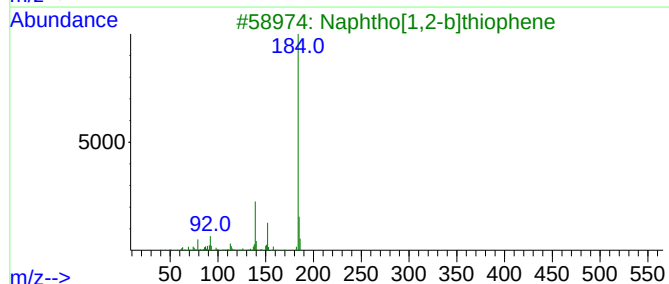
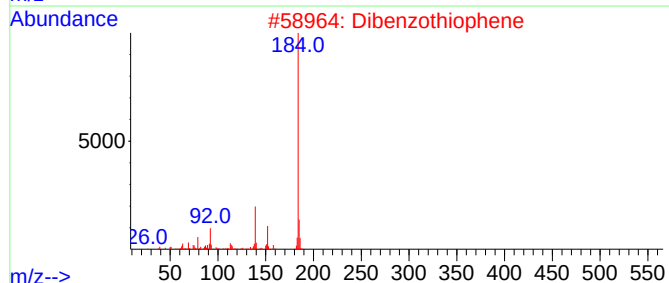
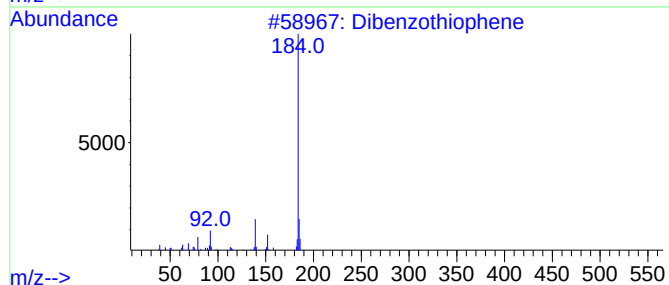
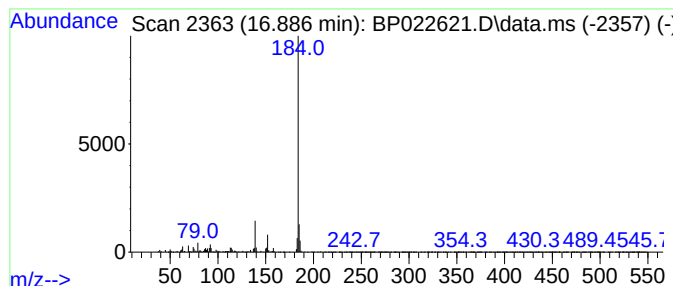
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Dibenzothiophene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.886	2.96 ng/ul	228539	Phenanthrene-d10	17.075

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzothiophene	184	C12H8S	000132-65-0	97
2			Dibenzothiophene	184	C12H8S	000132-65-0	95
3			Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	95
4			Dibenzothiophene	184	C12H8S	000132-65-0	95
5			Dibenzothiophene	184	C12H8S	000132-65-0	95



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Instrument :
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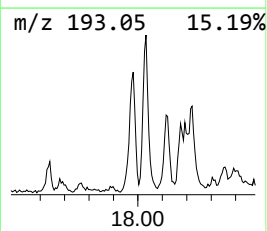
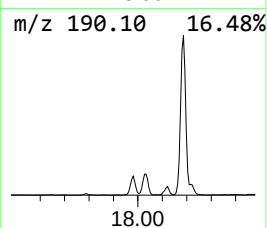
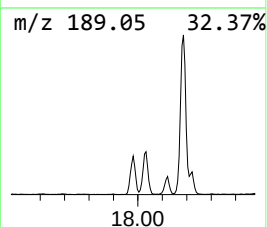
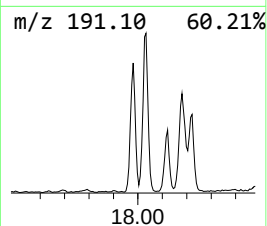
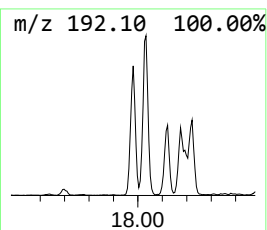
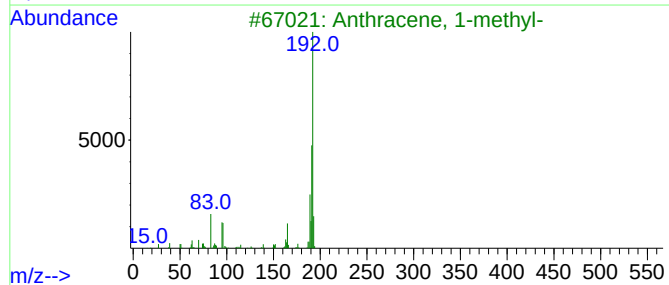
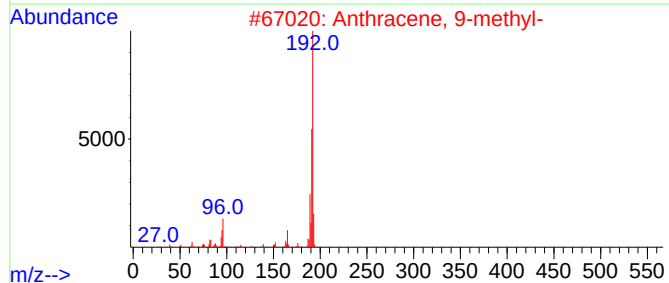
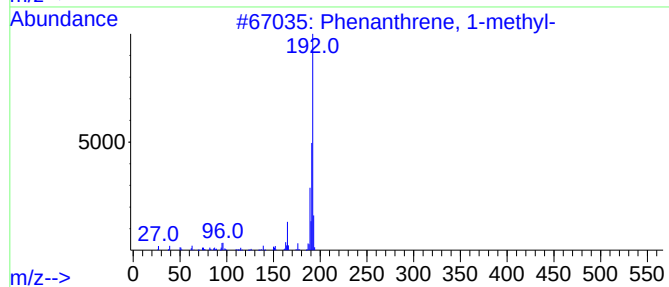
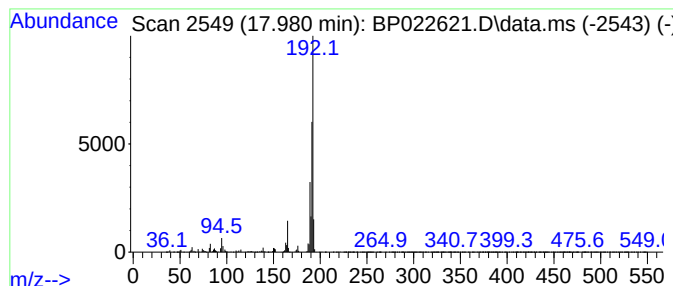
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Phenanthrene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.980	3.70 ng/ul	286203	Phenanthrene-d10	17.075

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
2			Anthracene, 9-methyl-	192	C15H12	000779-02-2	94
3			Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94
5			Anthracene, 9-methyl-	192	C15H12	000779-02-2	93



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Data File : BP022621.D
Acq On : 25 Oct 2024 22:01
Operator : RC/JU
Sample : P4436-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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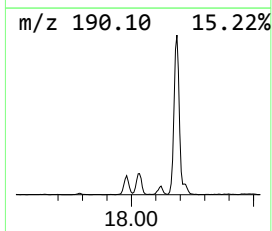
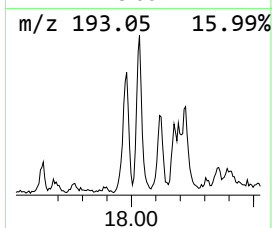
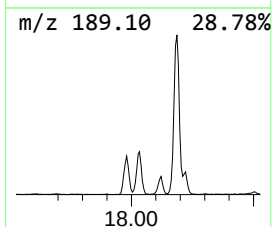
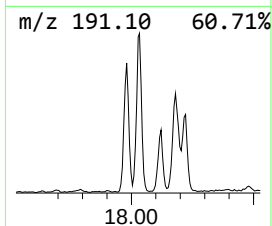
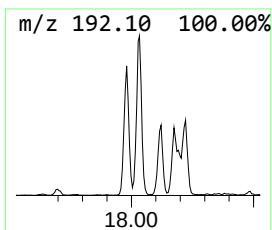
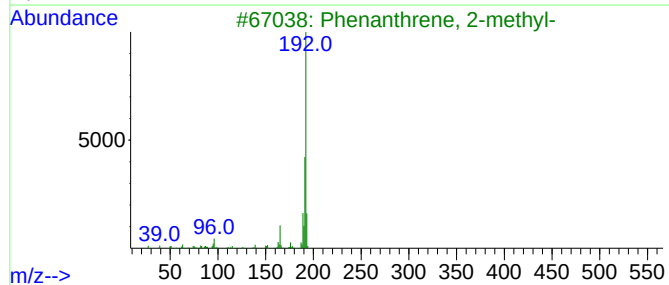
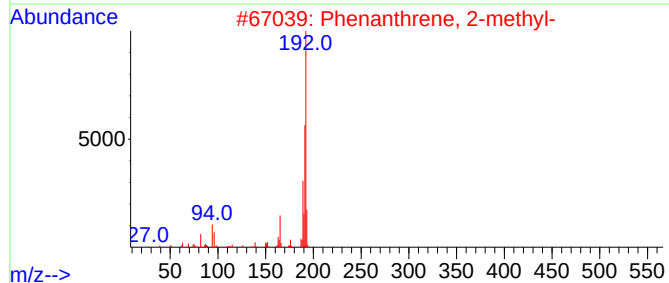
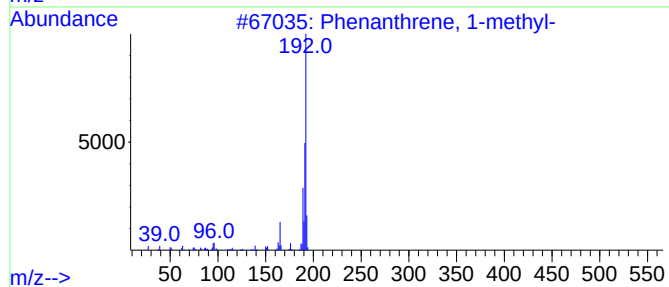
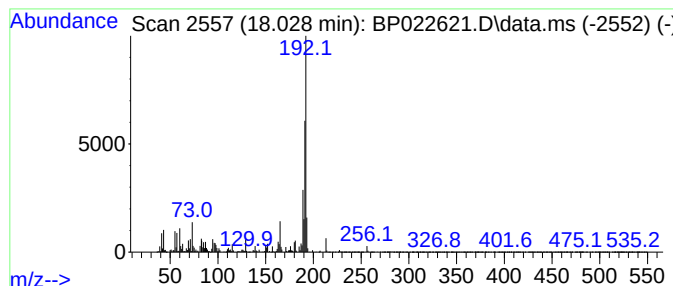
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.028	8.29 ng/ul	640540	Phenanthrene-d10	17.075

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			Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
5			Anthracene, 9-methyl-	192	C15H12	000779-02-2	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102524\
Data File : BP022621.D
Acq On : 25 Oct 2024 22:01
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Sample : P4436-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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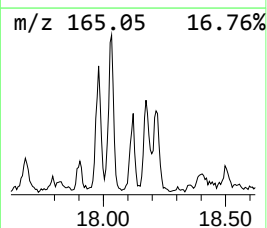
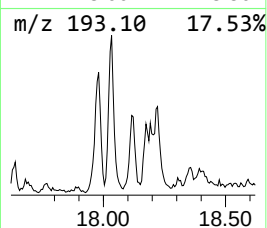
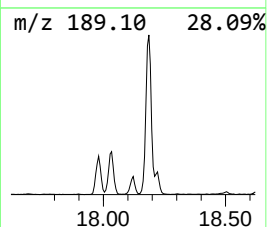
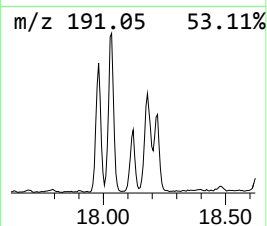
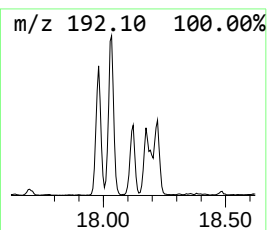
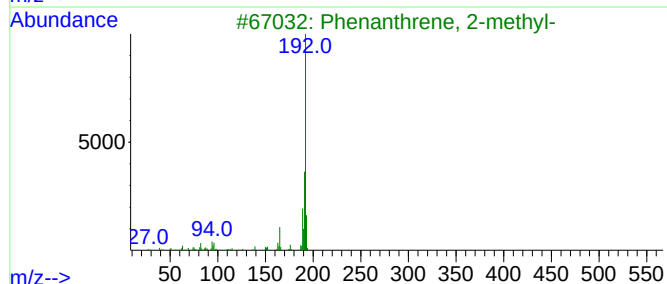
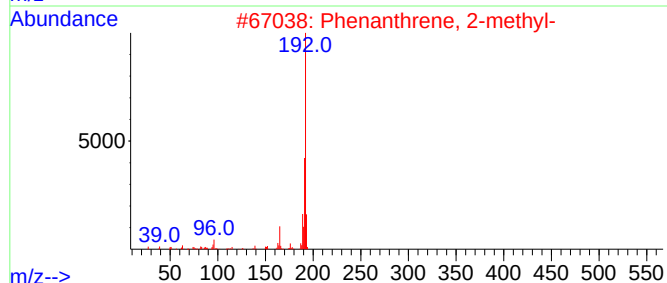
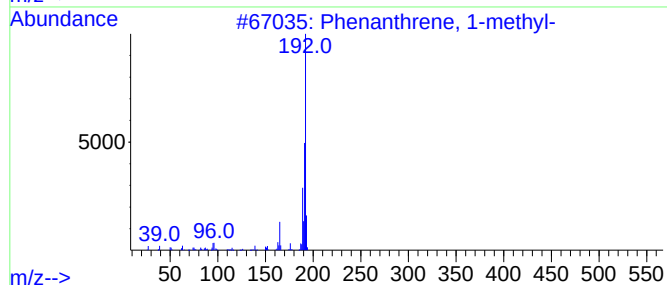
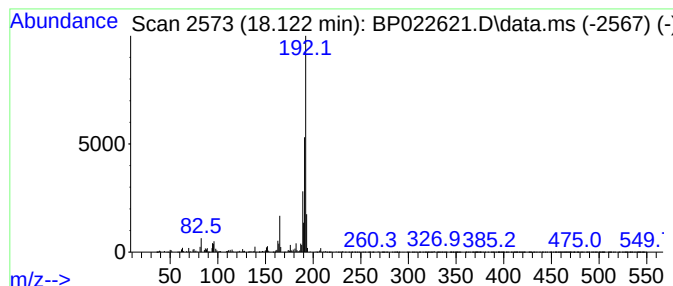
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Anthracene, 9-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.122	2.00 ng/ul	154915	Phenanthrene-d10	17.075

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	96
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
5			Anthracene, 9-methyl-	192	C15H12	000779-02-2	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102524\
Data File : BP022621.D
Acq On : 25 Oct 2024 22:01
Operator : RC/JU
Sample : P4436-05
Misc :
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Instrument :
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ClientSampleId :
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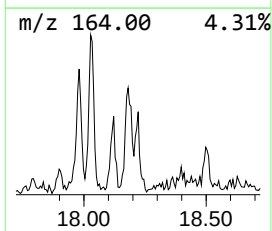
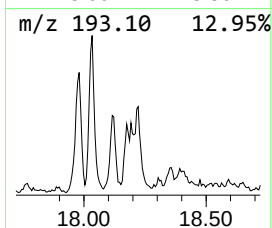
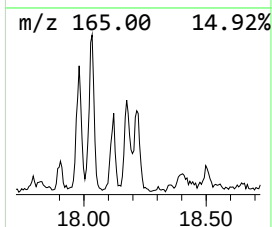
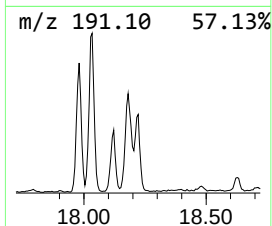
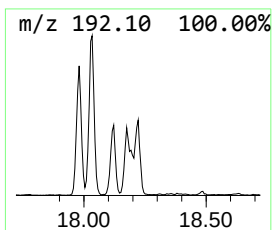
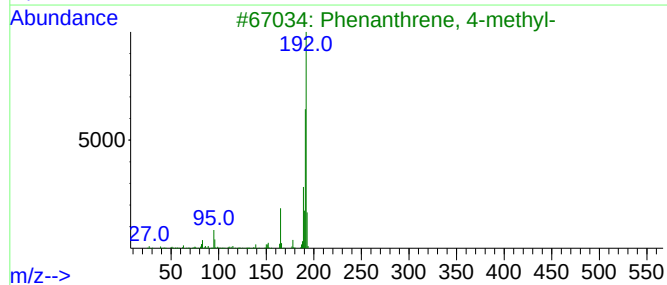
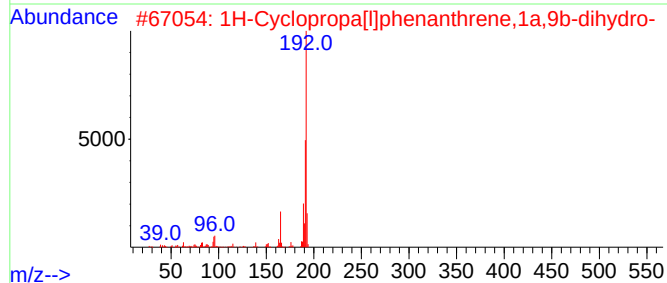
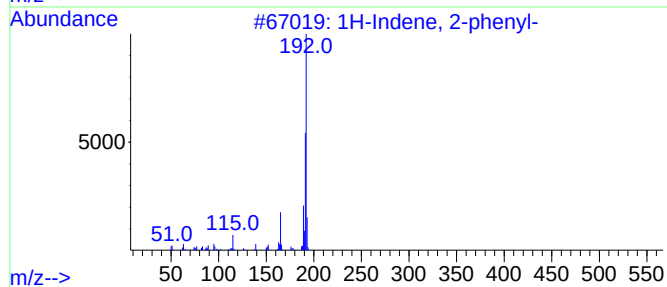
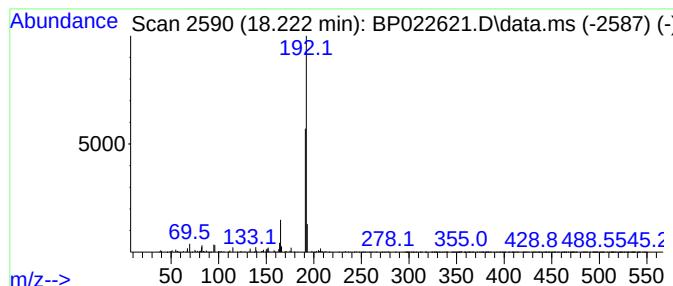
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1H-Indene, 2-phenyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.222	2.25 ng/ul	173922	Phenanthrene-d10	17.075

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	90
2			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	90
3			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	90
4			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\BP102524\
Data File : BP022621.D
Acq On : 25 Oct 2024 22:01
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ClientSampleId :
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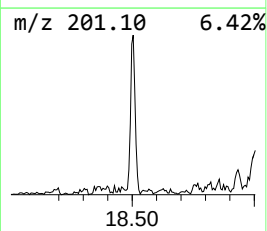
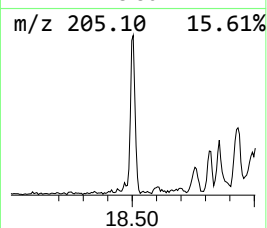
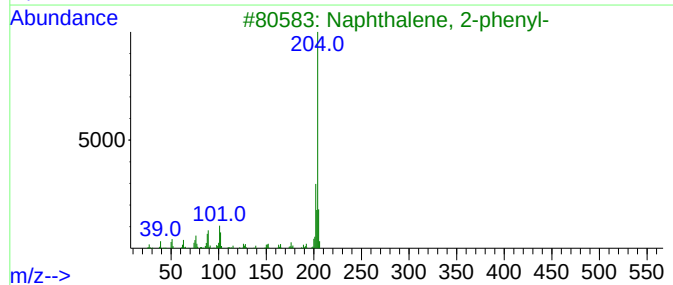
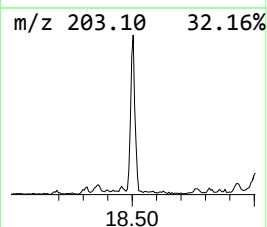
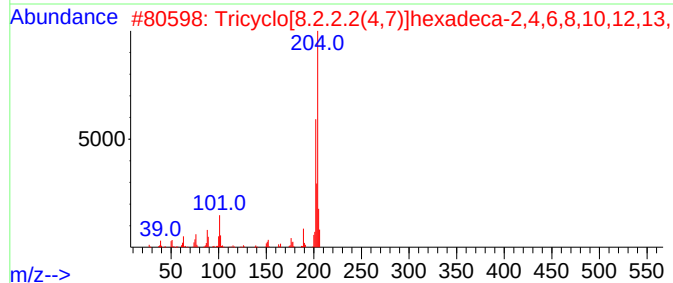
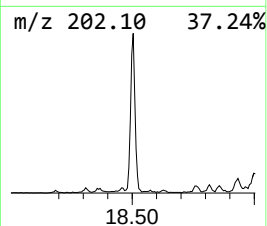
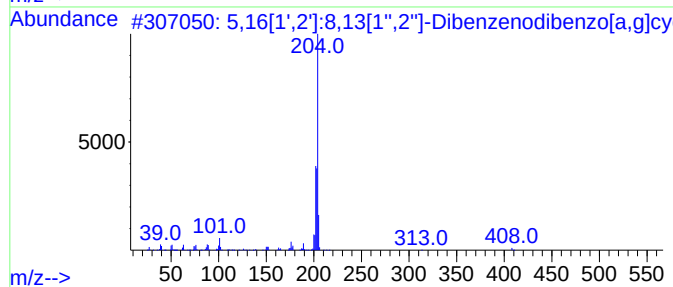
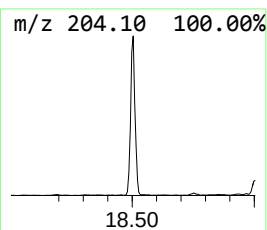
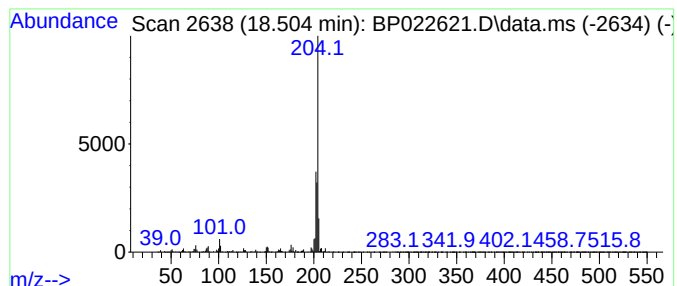
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.504	3.53 ng/ul	273236	Phenanthrene-d10	17.075

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	91		
2	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	81		
3	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	76		
4	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	76		
5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	76		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102524\
Data File : BP022621.D
Acq On : 25 Oct 2024 22:01
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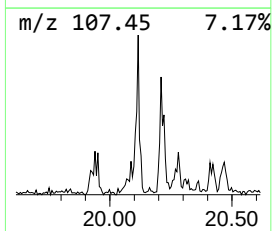
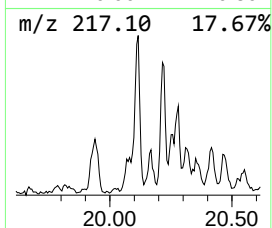
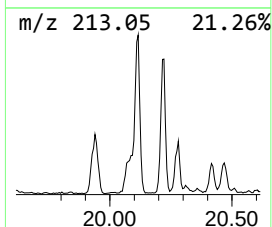
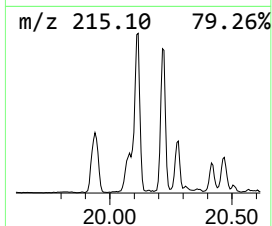
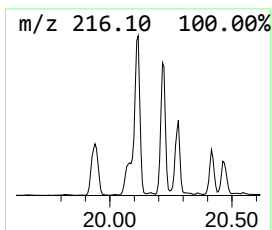
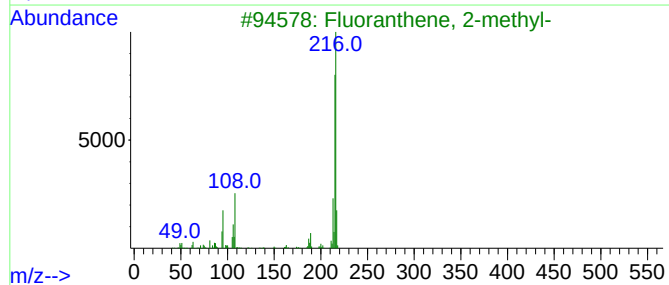
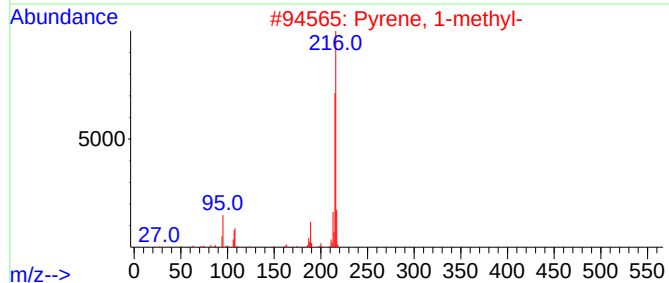
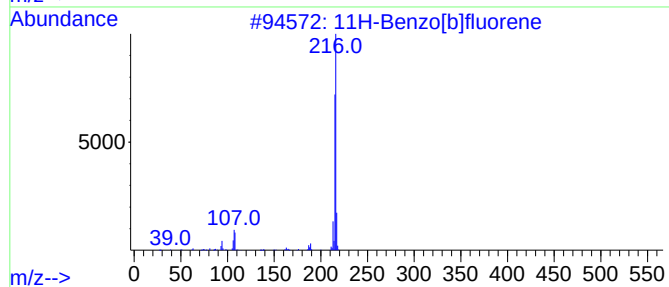
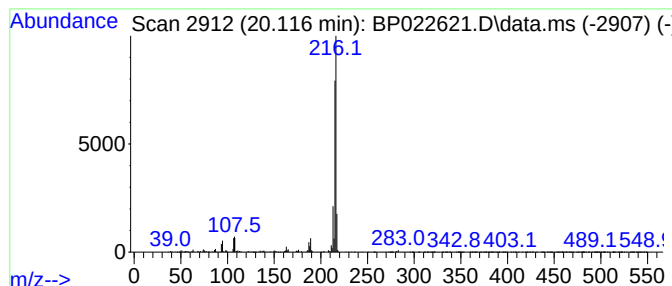
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 11H-Benzo[b]fluorene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.116	2.17 ng/ul	469273	Chrysene-d12	21.510

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102524\
Data File : BP022621.D
Acq On : 25 Oct 2024 22:01
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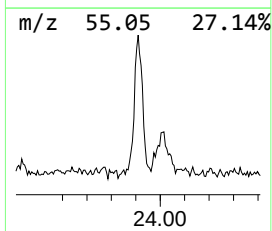
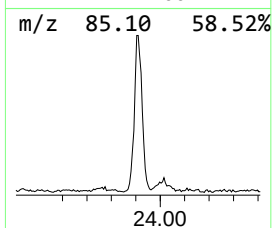
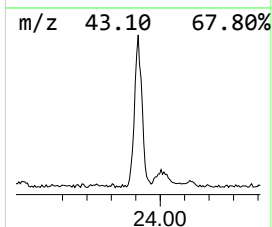
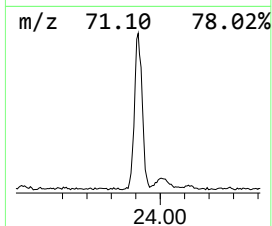
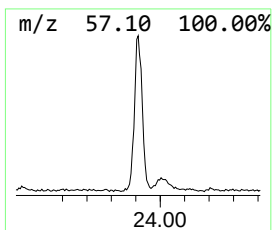
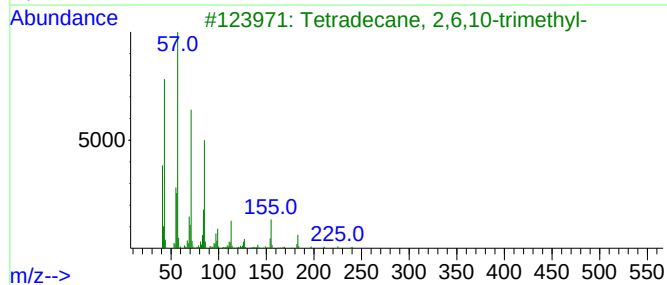
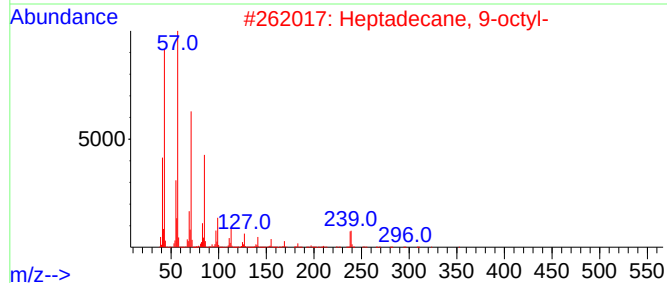
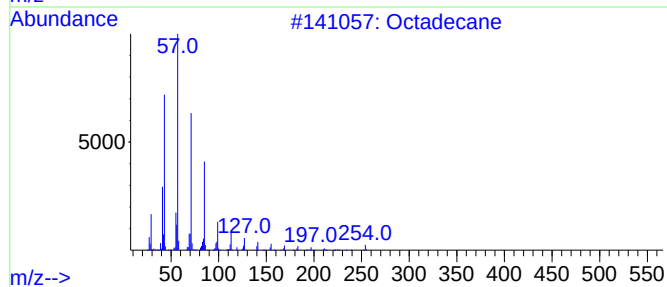
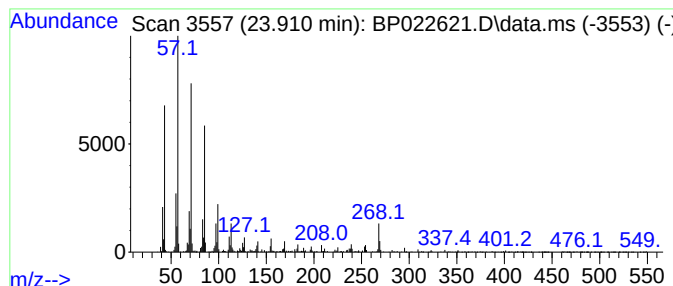
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.910	5.11 ng/ul	544320	Perylene-d12	24.774

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	96
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
3			Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	91
4			Octadecane	254	C18H38	000593-45-3	90
5			Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102524\
Data File : BP022621.D
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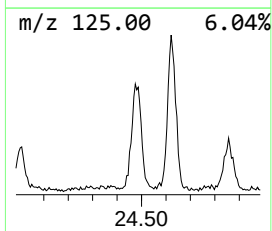
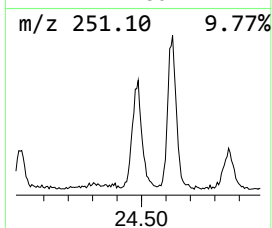
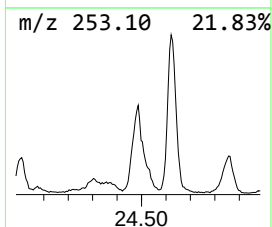
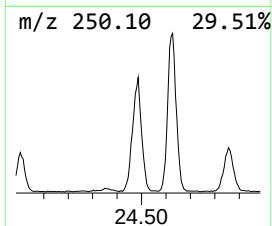
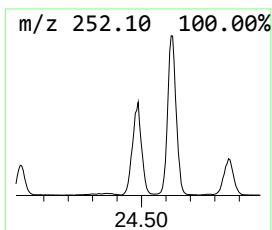
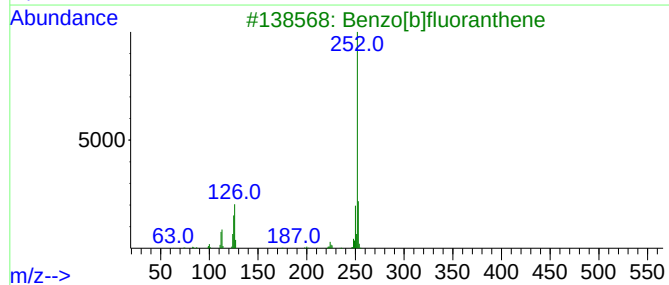
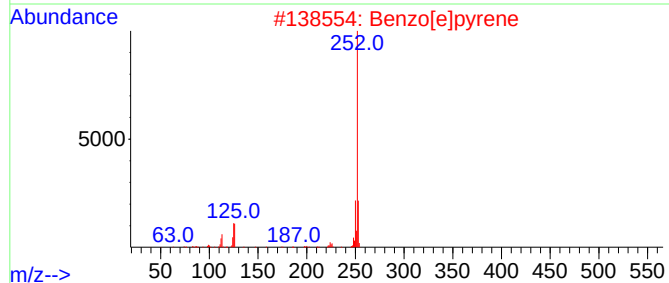
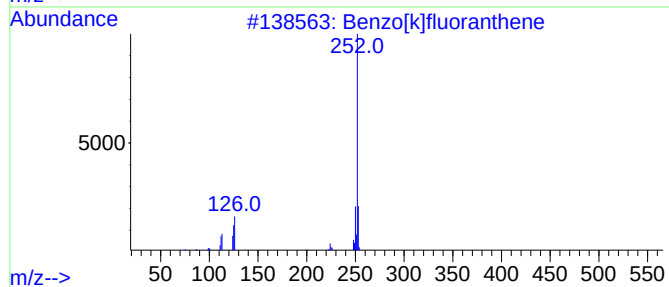
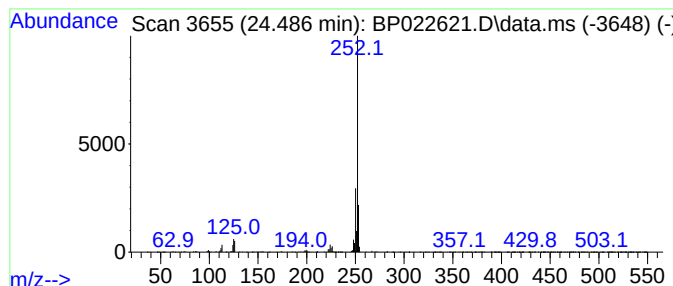
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.486	6.07 ng/ul	646510	Perylene-d12	24.774

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
2			Benzo[e]pyrene	252	C20H12	000192-97-2	95
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	91
4			Benzo[j]fluoranthene	252	C20H12	000205-82-3	90
5			Benzo[j]fluoranthene	252	C20H12	000205-82-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102524\
Data File : BP022621.D
Acq On : 25 Oct 2024 22:01
Operator : RC/JU
Sample : P4436-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4F47

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.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
Benzophenone	15.645	5.4	ng/ul	319696	3	14.269	1184660	20.0
Dibenzothiophene	16.886	3.0	ng/ul	228539	4	17.075	1545990	20.0
Phenanthrene, 1...	17.980	3.7	ng/ul	286203	4	17.075	1545990	20.0
Phenanthrene, 2...	18.028	8.3	ng/ul	640540	4	17.075	1545990	20.0
Anthracene, 9-m...	18.122	2.0	ng/ul	154915	4	17.075	1545990	20.0
1H-Indene, 2-ph...	18.222	2.3	ng/ul	173922	4	17.075	1545990	20.0
5,16[1',2']:8,1...	18.504	3.5	ng/ul	273236	4	17.075	1545990	20.0
11H-Benzo[b]flu...	20.116	2.2	ng/ul	469273	5	21.510	4329400	20.0
(DEL) Alkane: S...	23.910	5.1	ng/ul	544320	6	24.774	2131480	20.0
Benzo[e]pyrene	24.486	6.1	ng/ul	646510	6	24.774	2131480	20.0