

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052924\
 Data File : BG061355.D
 Acq On : 30 May 2024 9:45
 Operator : MA/JU
 Sample : P2571-20
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BH6A4

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_G\Methods\SFAM-EPA-BG052124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG061355.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.082	11	16	37	rVV	274402	462806	58.00%	4.633%
2	3.335	56	59	66	rVB4	5078	7380	0.92%	0.074%
3	3.617	101	107	112	rBV4	3430	6366	0.80%	0.064%
4	3.746	125	129	138	rBV	37117	64935	8.14%	0.650%
5	5.967	502	507	512	rVB2	3495	5091	0.64%	0.051%
6	7.066	686	694	702	rBV	117044	200170	25.09%	2.004%
7	7.213	713	719	728	rVB	71938	121006	15.16%	1.211%
8	7.418	748	754	761	rBV	111591	188288	23.60%	1.885%
9	7.882	826	833	840	rBV	91251	148732	18.64%	1.489%
10	7.941	840	843	847	rVB3	3210	4683	0.59%	0.047%
11	8.599	948	955	966	rVB2	126918	217718	27.28%	2.180%
12	9.034	1023	1029	1039	rVB	113780	204097	25.58%	2.043%
13	9.592	1119	1124	1129	rBV2	11514	17329	2.17%	0.173%
14	9.757	1146	1152	1162	rVB	103983	181020	22.69%	1.812%
15	10.309	1236	1246	1256	rBV	141205	256992	32.21%	2.573%
16	10.456	1267	1271	1276	rBV3	1937	3617	0.45%	0.036%
17	10.673	1301	1308	1315	rBV	120678	215493	27.01%	2.157%
18	10.814	1324	1332	1339	rBV	110793	201440	25.24%	2.017%
19	11.208	1394	1399	1404	rBV3	2906	4335	0.54%	0.043%
20	11.413	1428	1434	1442	rBV4	10251	18932	2.37%	0.190%
21	13.341	1758	1762	1769	rVB	3426	4936	0.62%	0.049%
22	13.834	1842	1846	1850	rBV3	2413	4080	0.51%	0.041%
23	13.922	1854	1861	1867	rVB	258376	365833	45.85%	3.662%
24	14.210	1899	1910	1921	rBV	274461	440391	55.19%	4.409%
25	14.416	1940	1945	1950	rVB2	2574	3988	0.50%	0.040%
26	14.516	1953	1962	1969	rVV	207962	310795	38.95%	3.111%
27	14.745	1992	2001	2011	rBV5	73029	155610	19.50%	1.558%
28	14.915	2027	2030	2034	rVB4	3242	4270	0.54%	0.043%
29	15.362	2104	2106	2113	rVB3	2484	4049	0.51%	0.041%
30	15.509	2124	2131	2137	rBV	378396	541144	67.82%	5.417%
31	15.562	2137	2140	2144	rVB3	3169	3756	0.47%	0.038%
32	15.644	2149	2154	2162	rBV	72384	104304	13.07%	1.044%
33	15.779	2173	2177	2181	rVB3	2707	3793	0.48%	0.038%
34	15.855	2186	2190	2196	rVB4	2811	4791	0.60%	0.048%
35	16.243	2247	2256	2259	rBV6	5017	11542	1.45%	0.116%
36	16.919	2367	2371	2377	rVV3	6815	9278	1.16%	0.093%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
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37	17.025	2385	2389	2392	rVV3	4161	4906	0.61%	0.049%
38	17.083	2395	2399	2403	rVV4	3350	5048	0.63%	0.051%
39	17.265	2423	2430	2434	rVV	257246	384906	48.24%	3.853%
40	17.307	2434	2437	2441	rVV	69397	95243	11.94%	0.953%

41	17.365	2441	2447	2458	rVB	381407	578853	72.54%	5.795%
42	17.665	2496	2498	2502	rVB2	3338	4072	0.51%	0.041%
43	17.841	2523	2528	2533	rVB3	5924	8640	1.08%	0.086%
44	17.994	2552	2554	2559	rVB2	2715	3544	0.44%	0.035%
45	18.065	2562	2566	2570	rBV5	2714	5119	0.64%	0.051%

46	18.141	2574	2579	2581	rBV	22406	32340	4.05%	0.324%
47	18.170	2581	2584	2585	rVV3	28287	36774	4.61%	0.368%
48	18.188	2585	2587	2598	rVV3	33532	49445	6.20%	0.495%
49	18.270	2598	2601	2607	rVV6	4946	8441	1.06%	0.085%
50	18.329	2607	2611	2615	rVV4	26383	44845	5.62%	0.449%

51	18.370	2615	2618	2625	rVB2	17504	26262	3.29%	0.263%
52	18.452	2629	2632	2633	rBV3	6035	5238	0.66%	0.052%
53	18.499	2636	2640	2643	rVV6	6534	12140	1.52%	0.122%
54	18.529	2643	2645	2646	rVV2	4799	4599	0.58%	0.046%
55	18.540	2646	2647	2652	rVV4	6160	5256	0.66%	0.053%

56	18.640	2660	2664	2667	rVV	15290	20758	2.60%	0.208%
57	18.681	2667	2671	2678	rVB4	21255	32638	4.09%	0.327%
58	18.887	2703	2706	2710	rVB4	7130	7487	0.94%	0.075%
59	18.940	2711	2715	2718	rVB2	4995	6422	0.80%	0.064%
60	18.981	2718	2722	2725	rVB3	6710	9825	1.23%	0.098%

61	19.057	2730	2735	2740	rBV3	22101	46349	5.81%	0.464%
62	19.116	2742	2745	2748	rVV5	9186	12237	1.53%	0.123%
63	19.146	2748	2750	2754	rVB3	7995	8371	1.05%	0.084%
64	19.198	2754	2759	2765	rBV5	18185	38194	4.79%	0.382%
65	19.322	2771	2780	2787	rBV	144110	224546	28.14%	2.248%

66	19.381	2787	2790	2793	rBV5	6625	9547	1.20%	0.096%
67	19.522	2810	2814	2818	rBV6	10734	14250	1.79%	0.143%
68	19.569	2820	2822	2826	rVB4	4902	6155	0.77%	0.062%
69	19.657	2831	2837	2840	rBV	534551	797950	100.00%	7.988%
70	19.751	2851	2853	2857	rBV5	12211	16268	2.04%	0.163%

71	19.845	2867	2869	2870	rBV2	8079	6025	0.76%	0.060%
72	19.857	2870	2871	2875	rVB2	9118	8186	1.03%	0.082%
73	19.921	2879	2882	2886	rVB5	12336	18392	2.30%	0.184%
74	19.951	2886	2887	2891	rBV3	3845	6043	0.76%	0.060%
75	20.003	2891	2896	2897	rBV5	18393	25711	3.22%	0.257%

76	20.133	2916	2918	2919	rBV2	14545	11396	1.43%	0.114%
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77	20.280	2938	2943	2946	rBV4	16162	24872	3.12%	0.249%
78	20.338	2948	2953	2956	rVB3	25466	37817	4.74%	0.379%
79	20.421	2963	2967	2970	rBV5	7378	12476	1.56%	0.125%
80	20.473	2973	2976	2980	rVV	19913	27265	3.42%	0.273%
81	20.520	2981	2984	2987	rVB3	16720	20352	2.55%	0.204%
82	20.620	2999	3001	3006	rVB6	11891	18560	2.33%	0.186%
83	20.691	3009	3013	3016	rVB5	19195	24405	3.06%	0.244%
84	20.932	3050	3054	3059	rVB	32131	46903	5.88%	0.470%
85	21.149	3088	3091	3093	rBV3	14840	21116	2.65%	0.211%
86	21.179	3093	3096	3105	rVV4	65042	135196	16.94%	1.353%
87	21.408	3132	3135	3141	rVB	55259	72200	9.05%	0.723%
88	21.513	3145	3153	3158	rVV3	322056	634813	79.56%	6.355%
89	21.560	3158	3161	3169	rVB	85413	127911	16.03%	1.281%
90	21.707	3182	3186	3191	rBV4	29928	42177	5.29%	0.422%
91	21.972	3227	3231	3237	rBV9	11038	29348	3.68%	0.294%
92	22.289	3281	3285	3288	rBV	33763	46125	5.78%	0.462%
93	22.929	3389	3394	3402	rVB5	45852	107543	13.48%	1.077%
94	23.623	3507	3512	3521	rBV2	54927	167099	20.94%	1.673%
95	24.322	3625	3631	3634	rBV2	22236	50933	6.38%	0.510%
96	24.392	3635	3643	3650	rBV2	274405	684309	85.76%	6.851%
97	24.604	3670	3679	3691	rBV	173445	497953	62.40%	4.985%
98	28.076	4268	4270	4271	rBV2	8382	4023	0.50%	0.040%
99	29.974	4591	4593	4594	rBV2	5827	4514	0.57%	0.045%
100	32.230	4975	4977	4980	rBV3	6239	5765	0.72%	0.058%

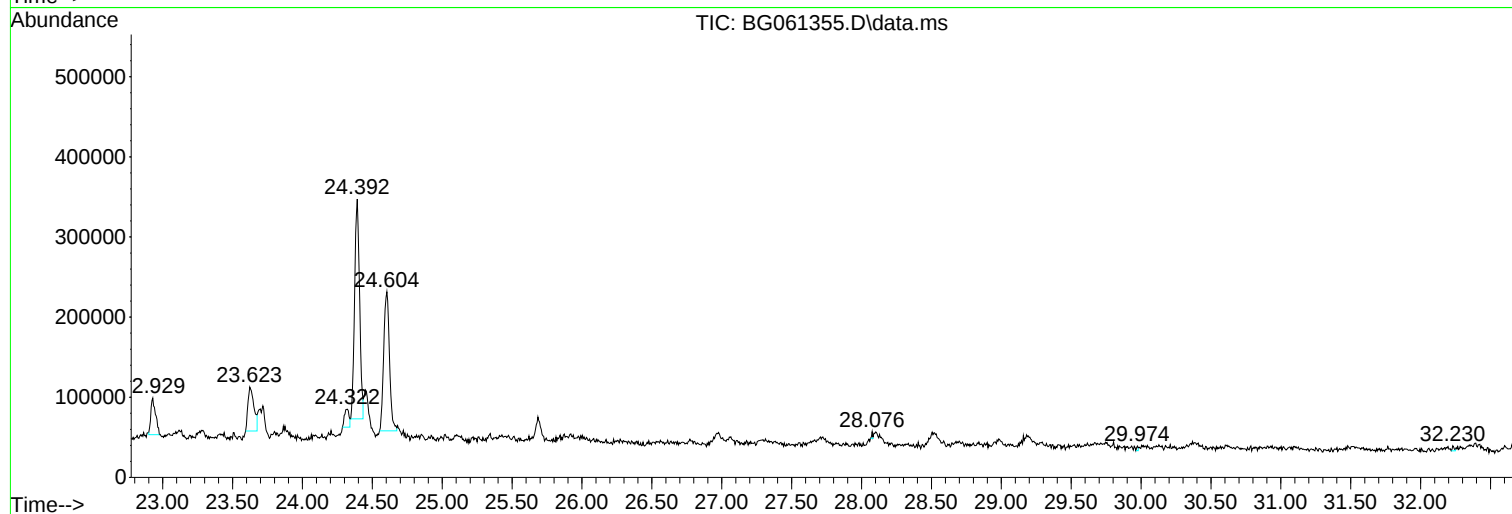
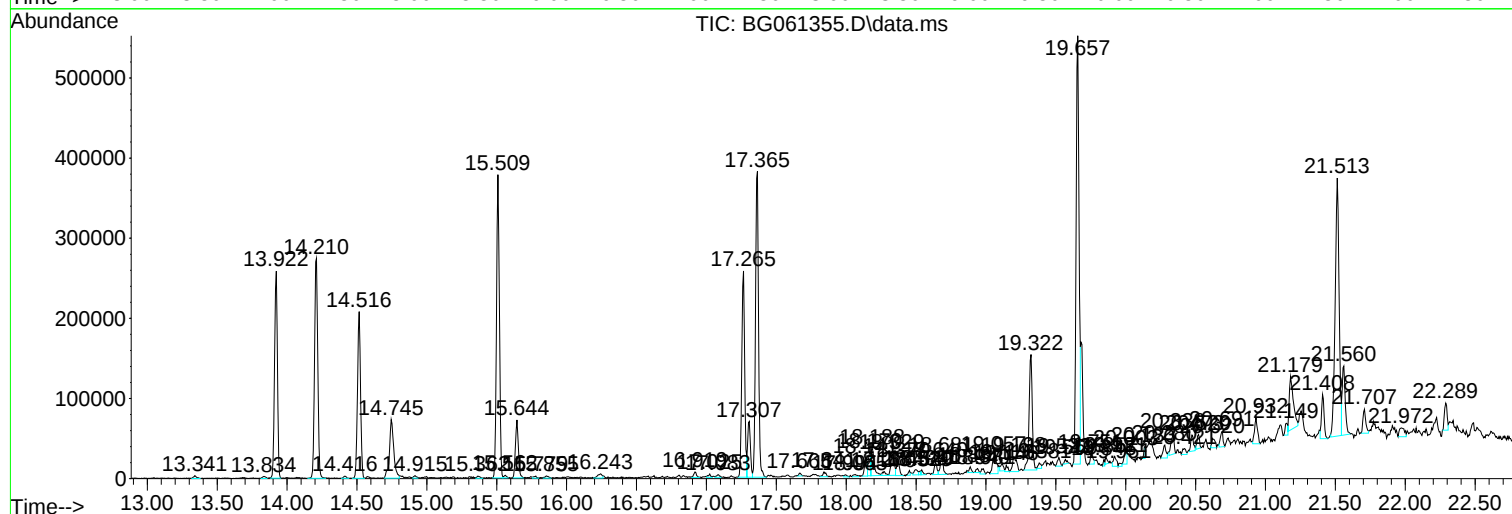
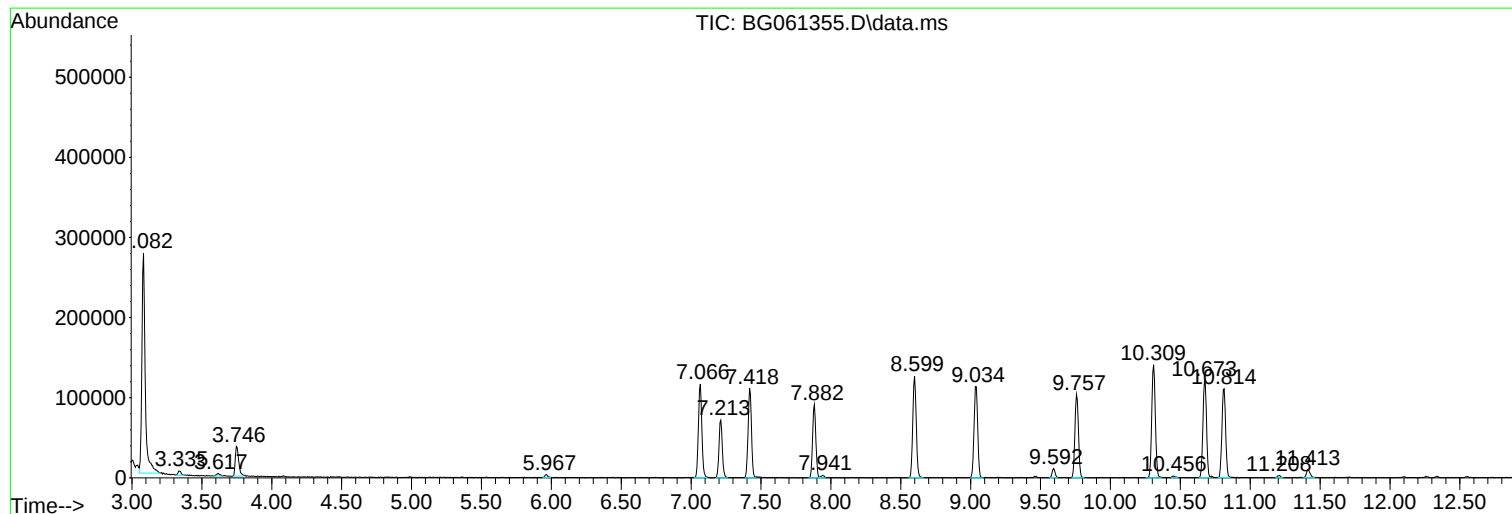
Sum of corrected areas: 9989086

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ALS Vial : 28 Sample Multiplier: 1

Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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



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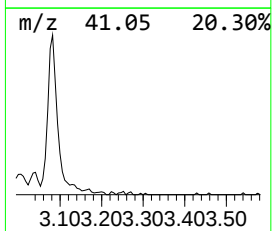
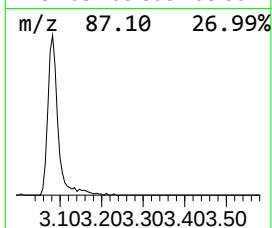
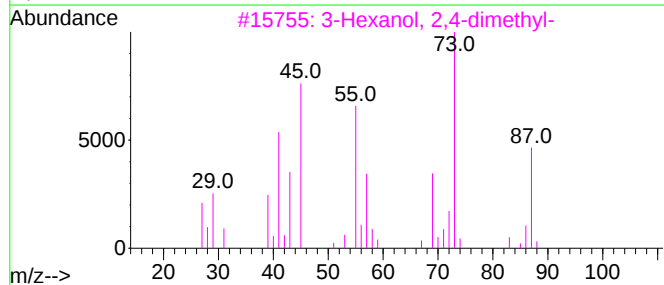
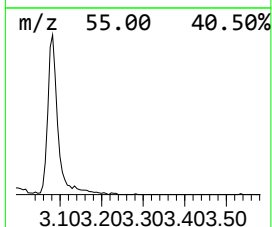
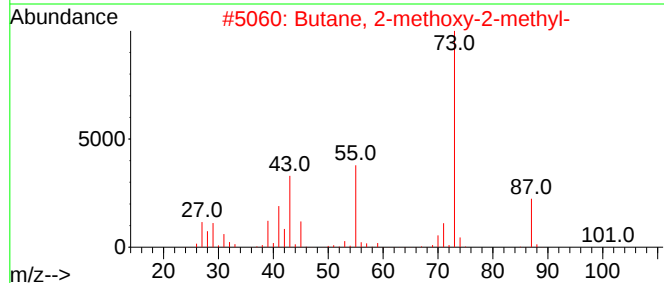
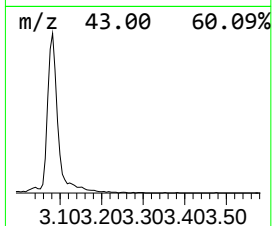
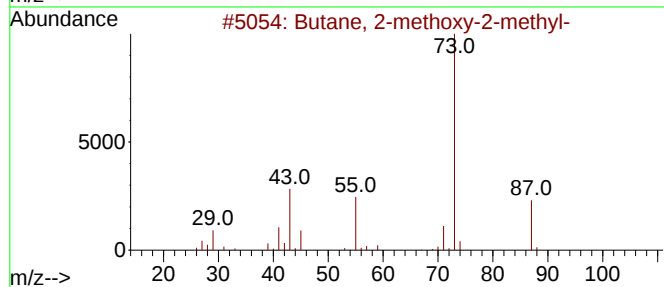
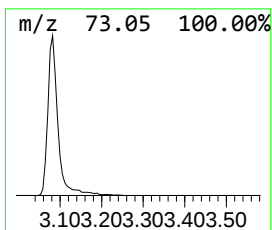
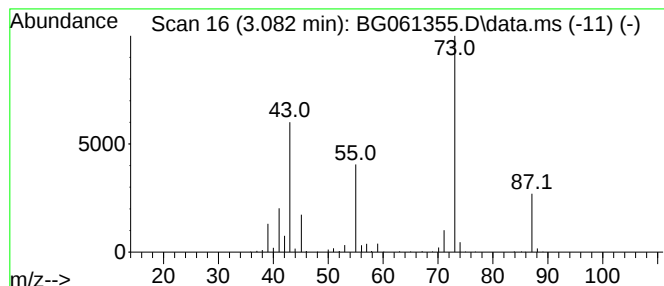
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.082	62.23 ng/ul	462806	1,4-Dichlorobenzene-d4	7.882

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3			3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	50
4			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
5			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	36



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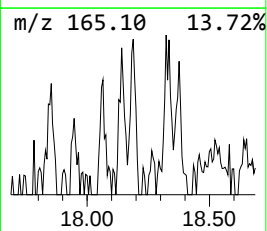
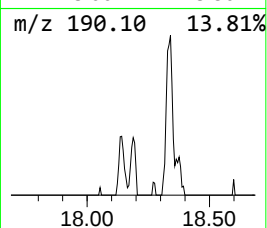
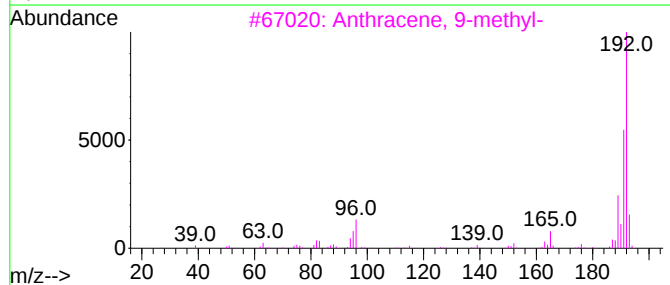
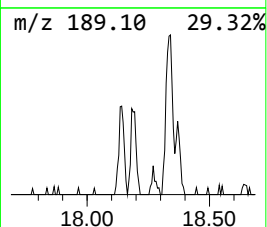
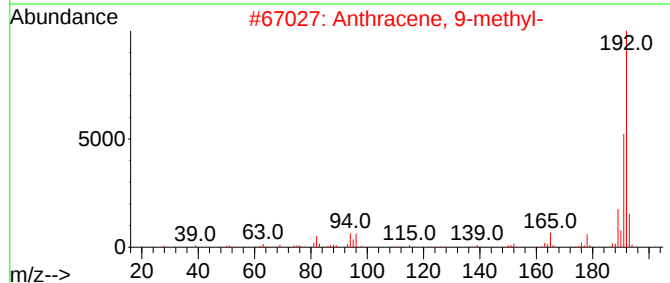
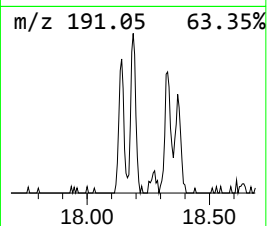
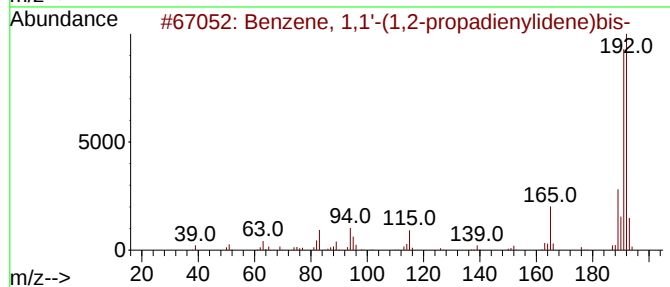
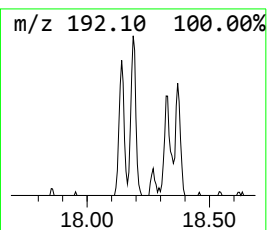
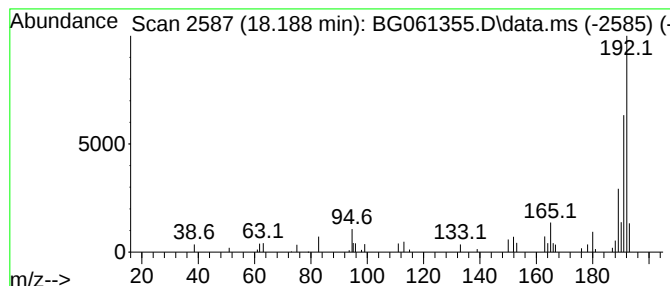
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,1'-(1,2-propadienylidene) Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.188	2.57 ng/ul	49445	Phenanthrene-d10	17.265

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,1'-(1,2-propadienylidene)bis-	192	C15H12	014251-57-1	90
2			Anthracene, 9-methyl-	192	C15H12	000779-02-2	81
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	76
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	76
5			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	76



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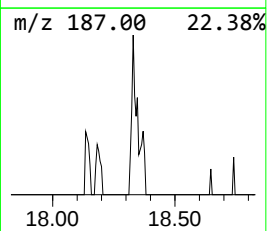
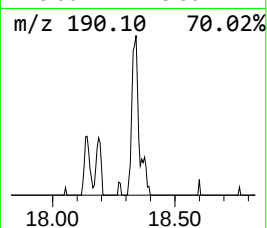
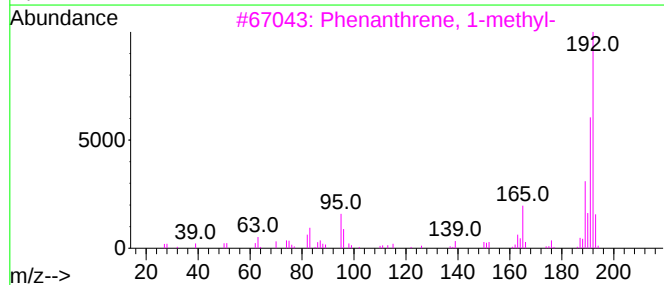
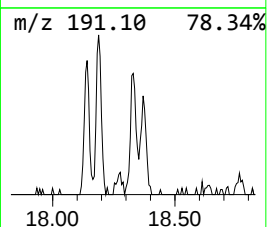
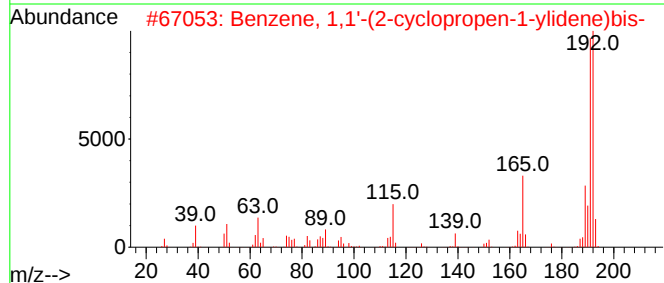
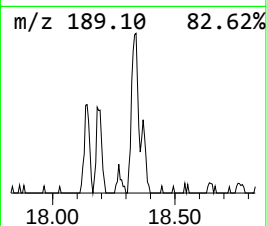
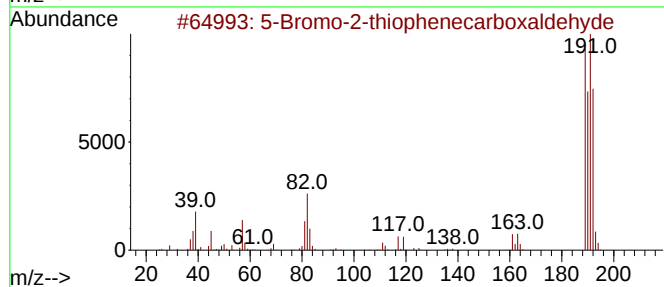
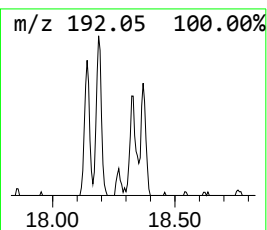
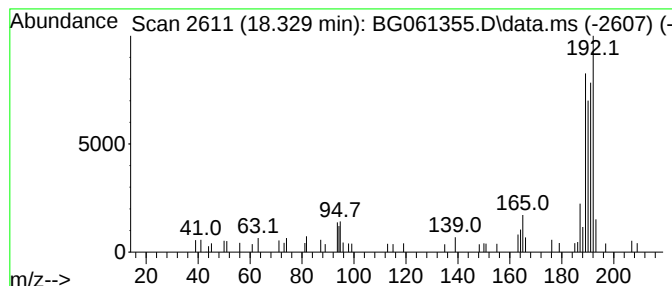
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 5-Bromo-2-thiophenecarboxal... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.329	2.33 ng/ul	44845	Phenanthrene-d10	17.265

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Bromo-2-thiophenecarboxaldehyde	190	C5H3BrOS	004701-17-1	58
2			Benzene, 1,1'-(2-cyclopropen-1-y...	192	C15H12	022825-21-4	49
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	49
4			Benzene, 1,1'-(1,2-propadienylid...	192	C15H12	014251-57-1	46
5			Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052924\
Data File : BG061355.D
Acq On : 30 May 2024 9:45
Operator : MA/JU
Sample : P2571-20
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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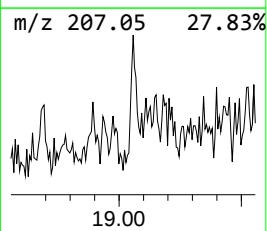
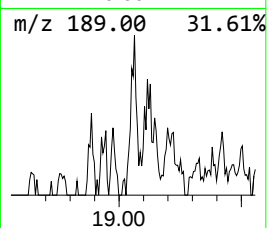
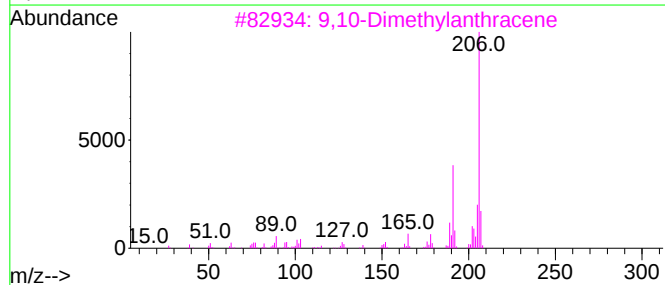
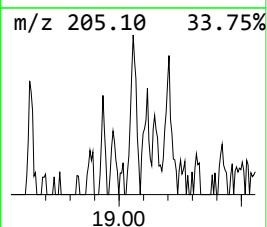
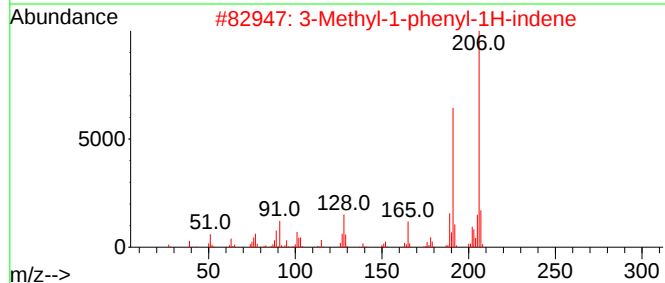
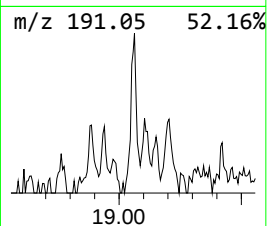
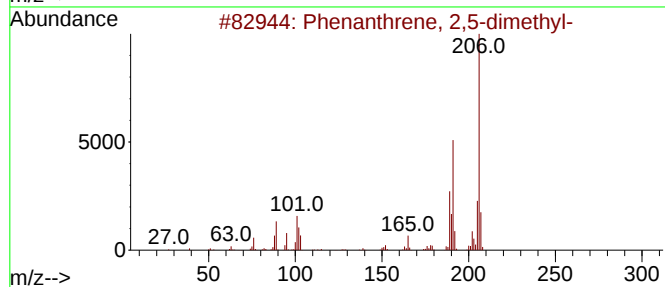
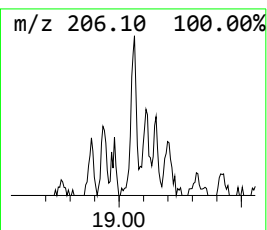
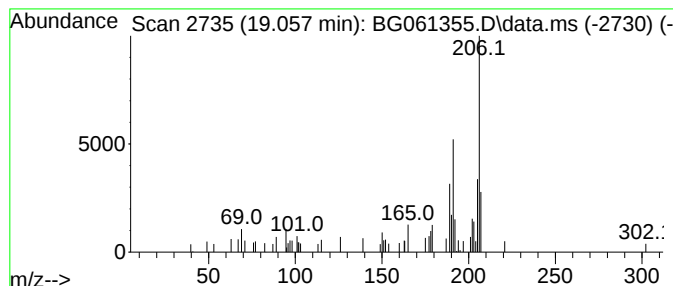
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.057	2.41 ng/ul	46349	Phenanthrene-d10	17.265

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	83
2		3-Methyl-1-phenyl-1H-indene	206	C16H14	022360-63-0	76
3		9,10-Dimethylantracene	206	C16H14	000781-43-1	76
4		Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	74
5		Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052924\
Data File : BG061355.D
Acq On : 30 May 2024 9:45
Operator : MA/JU
Sample : P2571-20
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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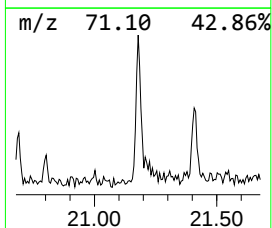
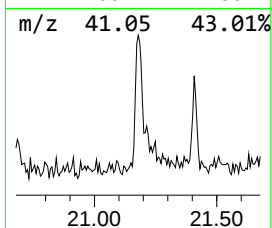
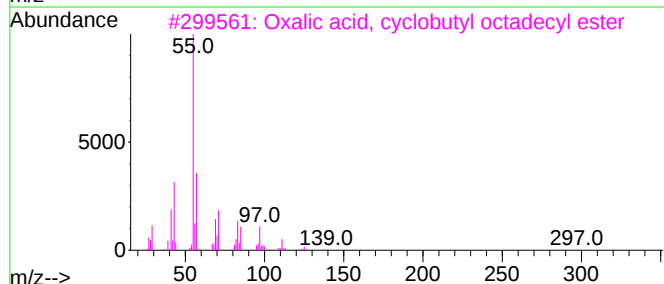
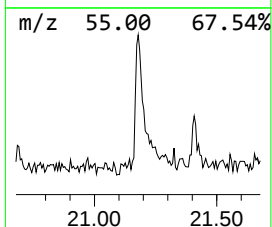
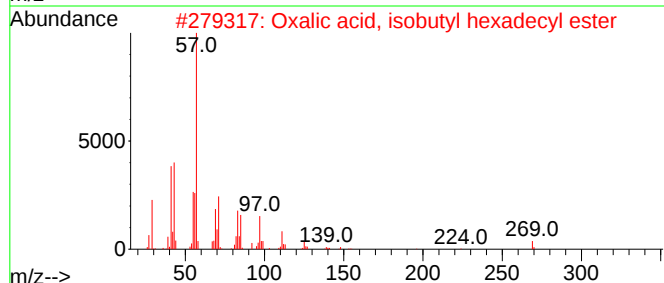
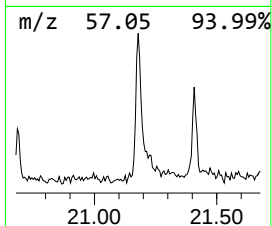
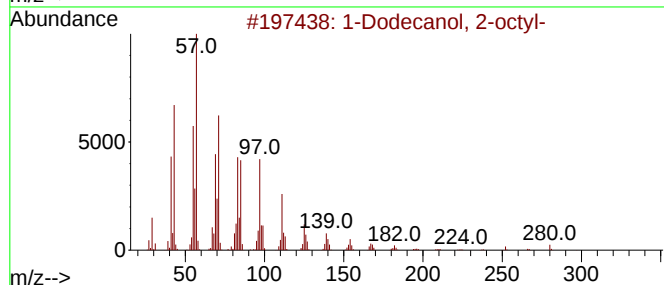
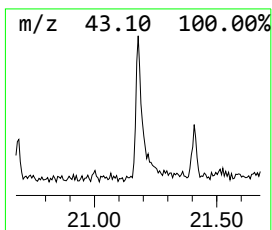
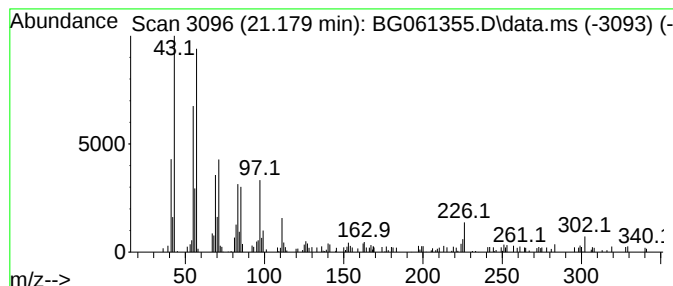
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1-Dodecanol, 2-octyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.179	4.26 ng/ul	135196	Chrysene-d12	21.513

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Dodecanol, 2-octyl-			298	C20H42O	005333-42-6	74
2	Oxalic acid, isobutyl hexadecyl ...			370	C22H42O4	1000309-38-1	74
3	Oxalic acid, cyclobutyl octadecyl...			396	C24H44O4	1000309-70-8	74
4	Oxalic acid, isobutyl heptadecyl...			384	C23H44O4	1000309-38-2	72
5	1-Chloroeicosane			316	C20H41Cl	042217-02-7	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052924\
Data File : BG061355.D
Acq On : 30 May 2024 9:45
Operator : MA/JU
Sample : P2571-20
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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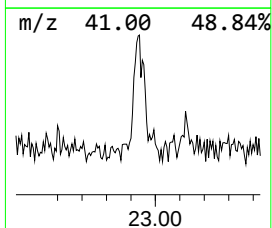
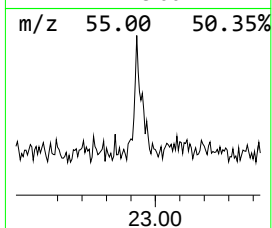
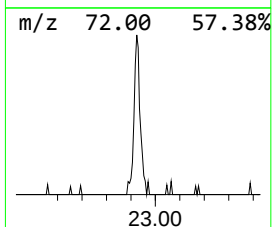
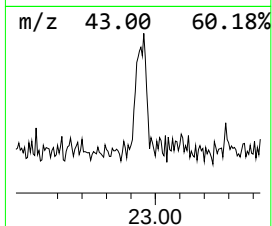
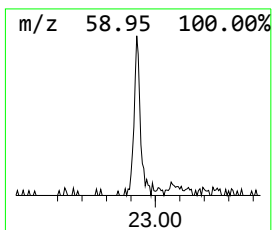
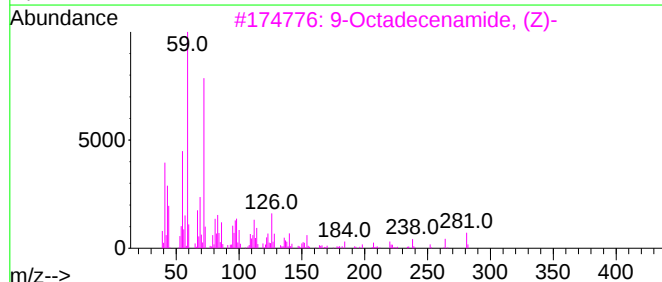
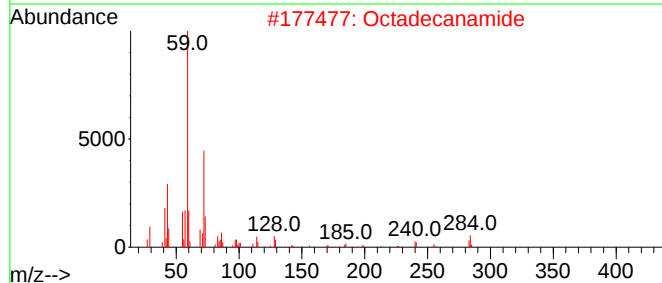
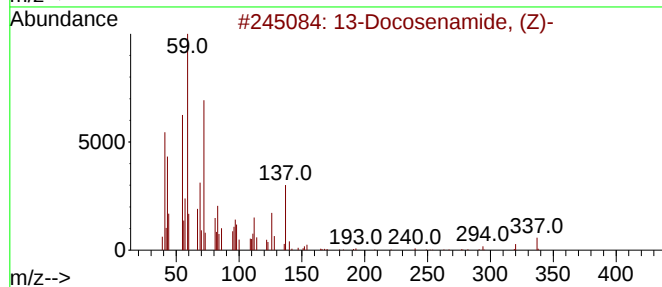
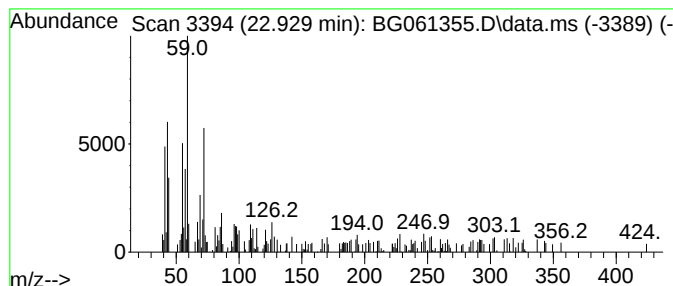
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 13-Docosenamide, (Z)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.929	3.39 ng/ul	107543	Chrysene-d12	21.513

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	93
2			Octadecanamide	283	C18H37NO	000124-26-5	64
3			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	64
4			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	64
5			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052924\
Data File : BG061355.D
Acq On : 30 May 2024 9:45
Operator : MA/JU
Sample : P2571-20
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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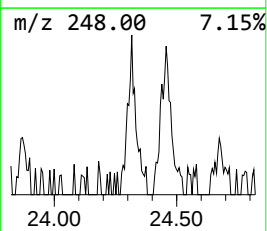
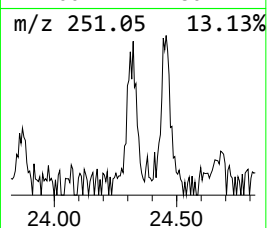
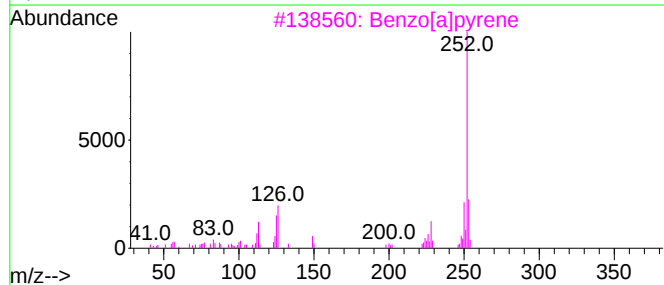
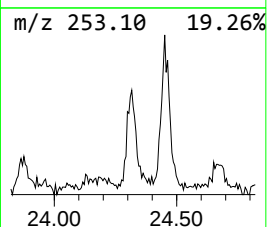
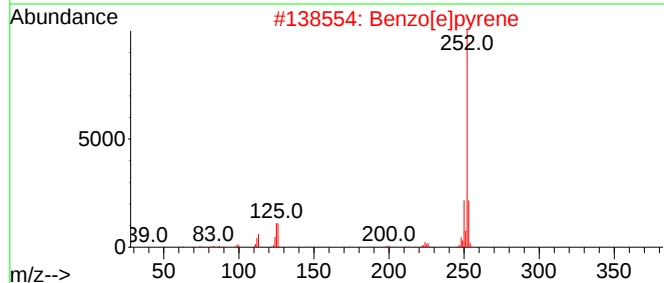
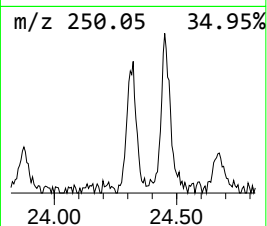
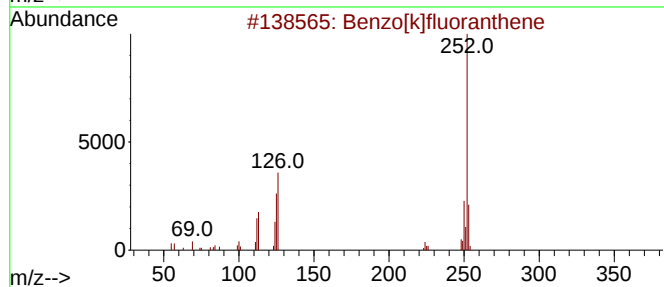
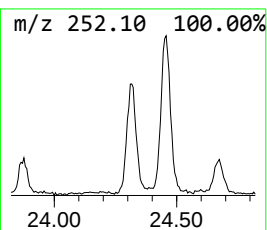
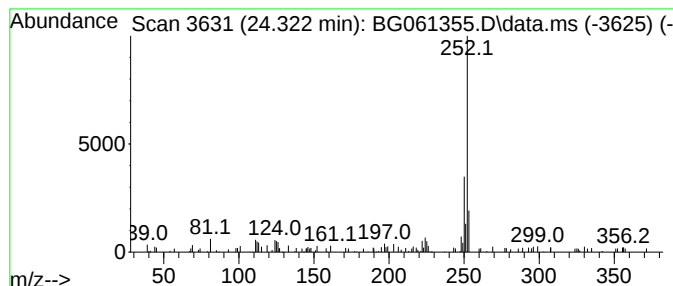
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzo[e]pyrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.322	2.05 ng/ul	50933	Perylene-d12	24.604

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[k]fluoranthene	252	C20H12	000207-08-9	81
2			Benzo[e]pyrene	252	C20H12	000192-97-2	74
3			Benzo[a]pyrene	252	C20H12	000050-32-8	74
4			(1-Cyclopentenyl)ferrocene	252	C15H16Fe	012260-67-2	72
5			Benzo[k]fluoranthene	252	C20H12	000207-08-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052924\
Data File : BG061355.D
Acq On : 30 May 2024 9:45
Operator : MA/JU
Sample : P2571-20
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH6A4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.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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Benzene, 1,1'-(...	18.188	2.6	ng/ul	49445	4	17.265	384906	20.0
5-Bromo-2-thiop...	18.329	2.3	ng/ul	44845	4	17.265	384906	20.0
Phenanthrene, 2...	19.057	2.4	ng/ul	46349	4	17.265	384906	20.0
1-Dodecanol, 2-...	21.179	4.3	ng/ul	135196	5	21.513	634813	20.0
13-Docosenamide...	22.929	3.4	ng/ul	107543	5	21.513	634813	20.0
Benzo[e]pyrene	24.322	2.0	ng/ul	50933	6	24.604	497953	20.0