

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052824\
 Data File : BG061303.D
 Acq On : 28 May 2024 17:19
 Operator : MA/JU
 Sample : P2568-22
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BH643

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: BG061303.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	48	rVB	1292823	1992013	100.00%	13.351%
2	3.340	54	60	65	rVB4	4712	8107	0.41%	0.054%
3	3.745	125	129	138	rBV	60657	98798	4.96%	0.662%
4	3.863	145	149	154	rVV4	5676	7863	0.39%	0.053%
5	4.309	219	225	229	rBV2	11326	15377	0.77%	0.103%
6	4.721	289	295	301	rVB2	5951	8590	0.43%	0.058%
7	7.065	688	694	709	rVB2	92989	161764	8.12%	1.084%
8	7.212	712	719	727	rBV	60645	95344	4.79%	0.639%
9	7.418	748	754	763	rBV	84566	146023	7.33%	0.979%
10	7.882	826	833	840	rBV	88078	146949	7.38%	0.985%
11	8.599	948	955	965	rBV	109008	181325	9.10%	1.215%
12	9.039	1023	1030	1037	rBV	90266	157352	7.90%	1.055%
13	9.762	1146	1153	1160	rBV	84070	144654	7.26%	0.969%
14	10.308	1240	1246	1255	rBV	117490	207559	10.42%	1.391%
15	10.449	1265	1270	1278	rBV	4900	8022	0.40%	0.054%
16	10.679	1302	1309	1316	rBV	135100	230379	11.57%	1.544%
17	10.814	1325	1332	1342	rBV	104932	186522	9.36%	1.250%
18	11.413	1429	1434	1440	rBV3	9389	16173	0.81%	0.108%
19	13.922	1852	1861	1869	rVV	246254	347018	17.42%	2.326%
20	14.210	1903	1910	1924	rVB2	257696	397717	19.97%	2.666%
21	14.515	1955	1962	1969	rBV	240471	361403	18.14%	2.422%
22	14.738	1989	2000	2017	rBV3	157953	374260	18.79%	2.508%
23	14.921	2027	2031	2037	rVV5	4610	8884	0.45%	0.060%
24	15.514	2125	2132	2138	rBV	353223	518081	26.01%	3.472%
25	15.643	2149	2154	2162	rBV	106381	160273	8.05%	1.074%
26	16.918	2366	2371	2378	rVB4	8634	13201	0.66%	0.088%
27	17.265	2423	2430	2434	rBV	321345	474619	23.83%	3.181%
28	17.306	2434	2437	2442	rVV	107921	149034	7.48%	0.999%
29	17.365	2442	2447	2458	rVB	365255	564261	28.33%	3.782%
30	17.664	2489	2498	2503	rBV	10942	18454	0.93%	0.124%
31	17.847	2522	2529	2537	rVB5	3312	7638	0.38%	0.051%
32	18.140	2572	2579	2582	rVV2	20990	30062	1.51%	0.201%
33	18.187	2582	2587	2592	rVV	31693	52689	2.65%	0.353%
34	18.228	2592	2594	2597	rVV2	7457	8227	0.41%	0.055%
35	18.270	2597	2601	2606	rVV3	5681	8618	0.43%	0.058%
36	18.334	2606	2612	2616	rVV4	26188	48776	2.45%	0.327%

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Integrator: RTE

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Start Thrs: 0.2

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Stop Thrs : 0

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

37	18.375	2616	2619	2624	rVB2	17526	22780	1.14%	0.153%
38	18.452	2627	2632	2635	rBV6	6365	8132	0.41%	0.055%
39	18.640	2659	2664	2668	rBV2	20004	28122	1.41%	0.188%
40	18.687	2668	2672	2678	rVV2	26873	41673	2.09%	0.279%
41	18.887	2703	2706	2711	rVB4	7395	9939	0.50%	0.067%
42	19.057	2731	2735	2741	rBV4	19035	37473	1.88%	0.251%
43	19.110	2741	2744	2748	rVB4	8513	11326	0.57%	0.076%
44	19.151	2748	2751	2755	rVB3	8015	10132	0.51%	0.068%
45	19.204	2755	2760	2770	rBV4	20278	44674	2.24%	0.299%
46	19.321	2770	2780	2786	rVB	323070	451541	22.67%	3.026%
47	19.415	2793	2796	2802	rVB6	6877	8412	0.42%	0.056%
48	19.468	2802	2805	2809	rBV	6922	8773	0.44%	0.059%
49	19.521	2809	2814	2817	rBV4	8908	14638	0.73%	0.098%
50	19.568	2817	2822	2830	rVB3	10428	21778	1.09%	0.146%
51	19.656	2831	2837	2840	rBV	535670	843028	42.32%	5.650%
52	19.686	2840	2842	2849	rVB	270040	312440	15.68%	2.094%
53	19.756	2849	2854	2863	rVB2	17696	30263	1.52%	0.203%
54	19.862	2866	2872	2877	rBV	18129	27539	1.38%	0.185%
55	19.921	2877	2882	2888	rVB7	7461	13621	0.68%	0.091%
56	20.003	2891	2896	2898	rBV4	24927	42336	2.13%	0.284%
57	20.097	2908	2912	2915	rBV6	5571	9784	0.49%	0.066%
58	20.144	2915	2920	2922	rBV3	19485	32848	1.65%	0.220%
59	20.179	2922	2926	2934	rVB2	32799	63300	3.18%	0.424%
60	20.279	2939	2943	2947	rBV2	14629	22759	1.14%	0.153%
61	20.338	2947	2953	2956	rVV2	34529	55797	2.80%	0.374%
62	20.373	2956	2959	2963	rVB4	10016	14260	0.72%	0.096%
63	20.420	2963	2967	2971	rBV6	7382	11609	0.58%	0.078%
64	20.479	2973	2977	2980	rBV	21049	30312	1.52%	0.203%
65	20.520	2980	2984	2988	rVB	22201	32708	1.64%	0.219%
66	20.632	2999	3003	3009	rVB8	10438	16679	0.84%	0.112%
67	20.690	3009	3013	3015	rBV3	12167	13055	0.66%	0.087%
68	20.737	3017	3021	3024	rVV6	13285	17361	0.87%	0.116%
69	20.931	3048	3054	3059	rVB2	52031	88290	4.43%	0.592%
70	21.102	3077	3083	3088	rBV3	35219	77985	3.91%	0.523%
71	21.149	3088	3091	3093	rVV	31375	45811	2.30%	0.307%
72	21.184	3093	3097	3105	rVV4	61588	158339	7.95%	1.061%
73	21.254	3105	3109	3117	rVB	48301	89177	4.48%	0.598%
74	21.378	3127	3130	3131	rVV2	17896	19824	1.00%	0.133%
75	21.413	3132	3136	3140	rVV	42529	65676	3.30%	0.440%
76	21.513	3144	3153	3158	rVV3	416965	930180	46.70%	6.234%

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77	21.560	3158	3161	3173	rVB	175791	287314	14.42%	1.926%
78	21.707	3182	3186	3193	rBV4	25982	46019	2.31%	0.308%
79	21.771	3194	3197	3201	rBV4	12879	19140	0.96%	0.128%
80	21.907	3215	3220	3228	rBV4	21468	49625	2.49%	0.333%
81	21.965	3228	3230	3236	rVB8	10198	14648	0.74%	0.098%
82	22.142	3258	3260	3264	rBV4	6689	10350	0.52%	0.069%
83	22.224	3269	3274	3278	rVB2	28714	52649	2.64%	0.353%
84	22.294	3280	3286	3292	rBV6	41841	77056	3.87%	0.516%
85	22.488	3314	3319	3322	rBV4	19476	34803	1.75%	0.233%
86	22.617	3335	3341	3347	rBV7	12342	31500	1.58%	0.211%
87	22.870	3377	3384	3386	rVV7	10154	21428	1.08%	0.144%
88	22.952	3388	3398	3405	rVB9	25713	85581	4.30%	0.574%
89	23.281	3445	3454	3462	rVB9	17882	53118	2.67%	0.356%
90	23.622	3504	3512	3520	rBV2	134376	434632	21.82%	2.913%
91	23.869	3549	3554	3561	rVB2	22841	48842	2.45%	0.327%
92	24.315	3623	3630	3635	rVV2	74380	195114	9.79%	1.308%
93	24.392	3635	3643	3650	rVV2	283216	778579	39.09%	5.218%
94	24.456	3650	3654	3664	rVB2	103723	238364	11.97%	1.598%
95	24.603	3669	3679	3687	rBV	218315	615770	30.91%	4.127%
96	25.690	3858	3864	3877	rVB3	44962	127111	6.38%	0.852%
97	26.977	4076	4083	4090	rVV7	14535	47758	2.40%	0.320%
98	28.093	4261	4273	4292	rVB3	51354	237997	11.95%	1.595%
99	28.528	4336	4347	4359	rBV7	28213	117638	5.91%	0.788%
100	29.186	4446	4459	4470	rVB2	36166	155154	7.79%	1.040%

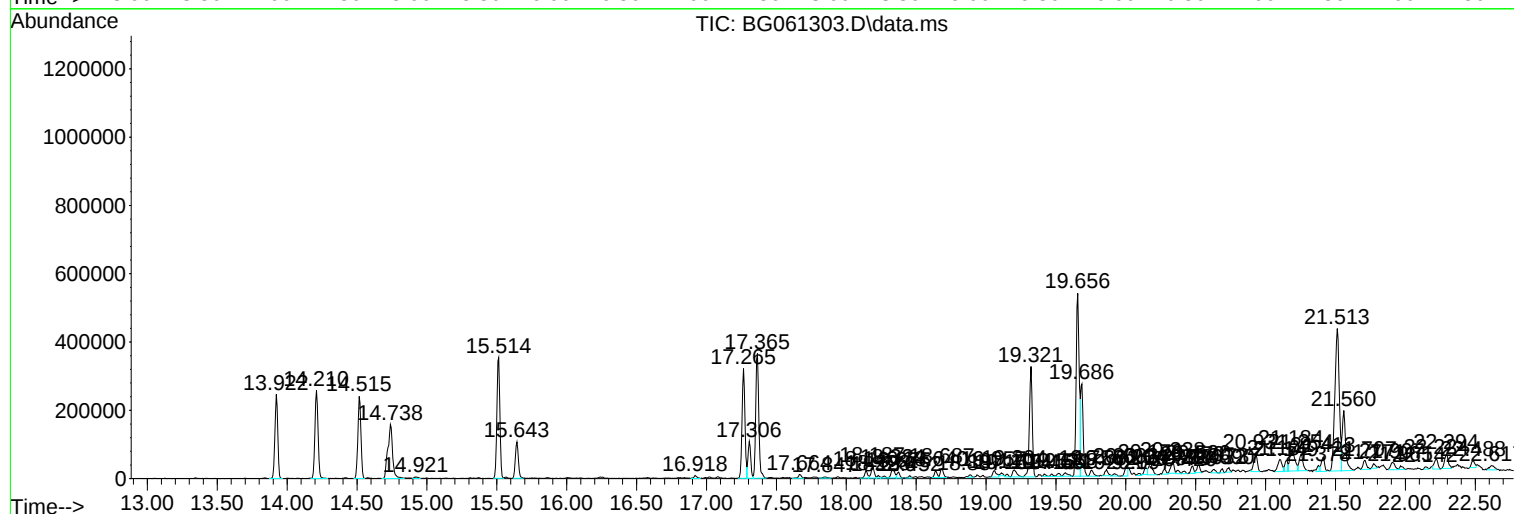
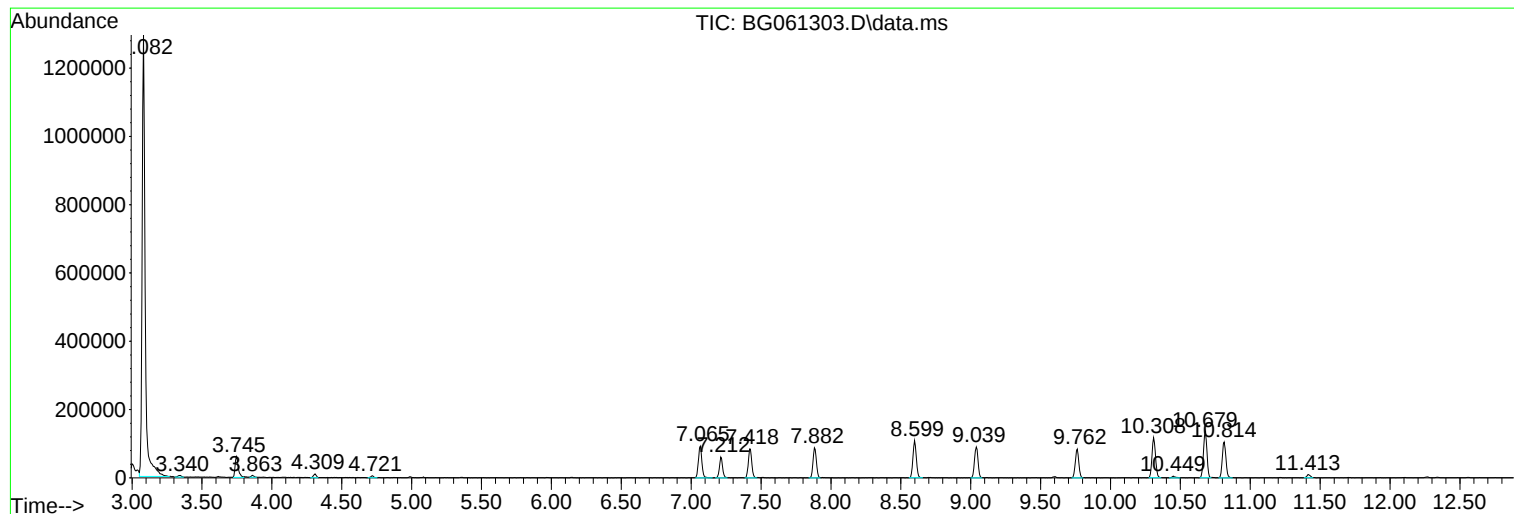
Sum of corrected areas: 14920593

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ALS Vial : 11 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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Acq On : 28 May 2024 17:19
Operator : MA/JU
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Misc :
ALS Vial : 11 Sample Multiplier: 1

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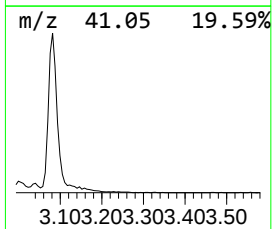
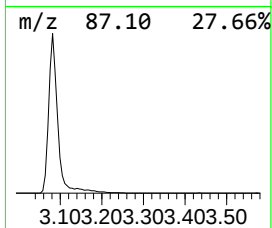
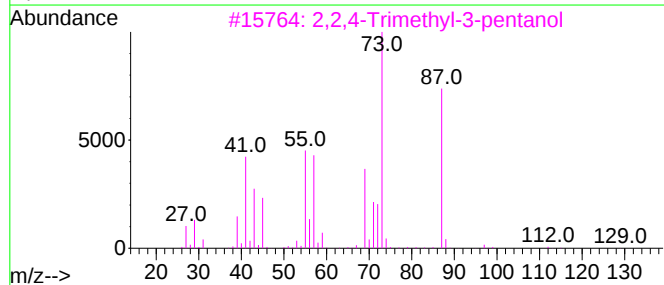
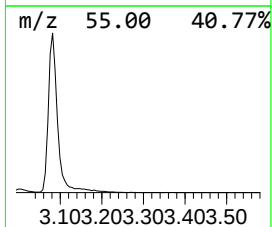
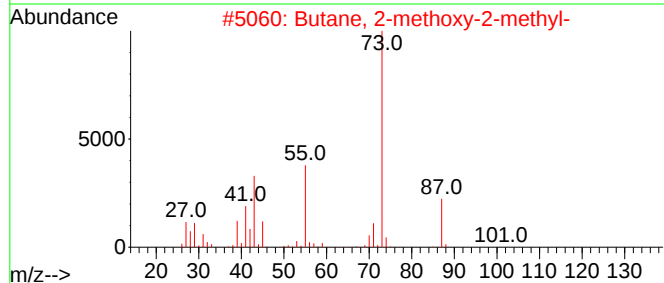
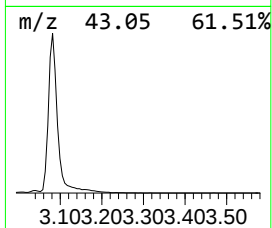
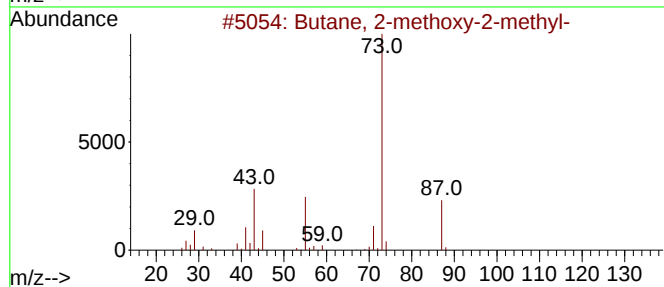
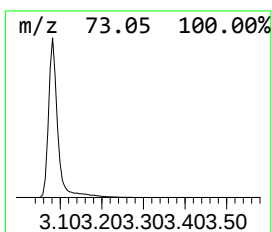
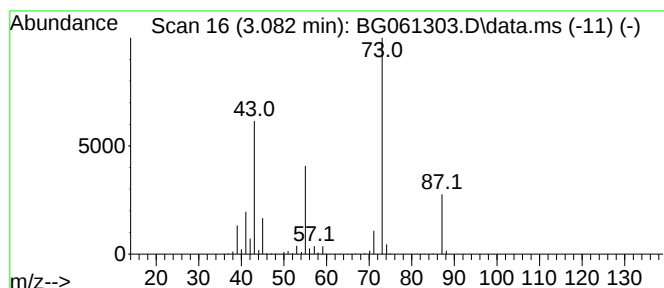
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.082	271.12 ng/ul	1992010	1,4-Dichlorobenzene-d4	7.888

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			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	38
4			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	37
5			3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	33



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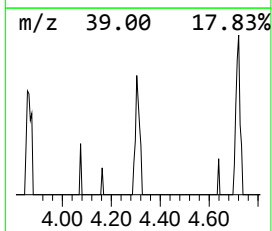
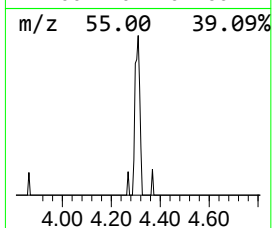
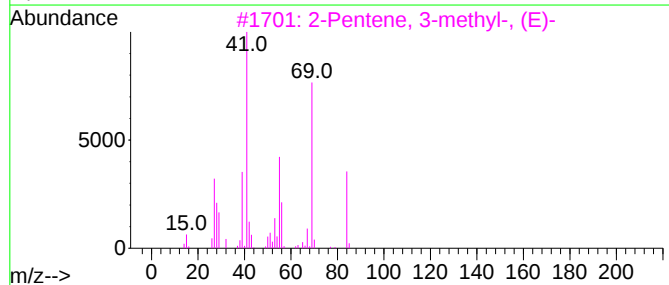
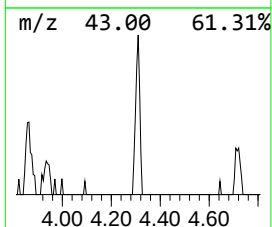
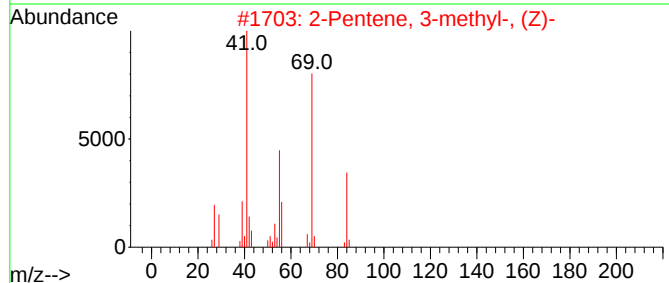
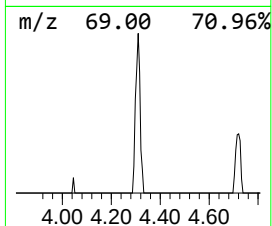
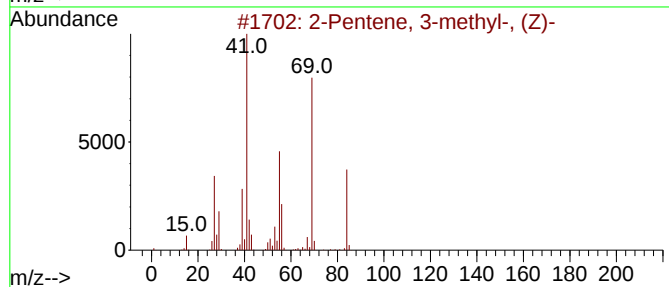
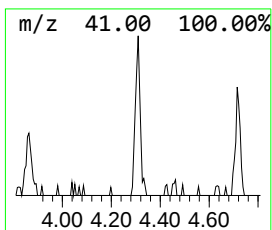
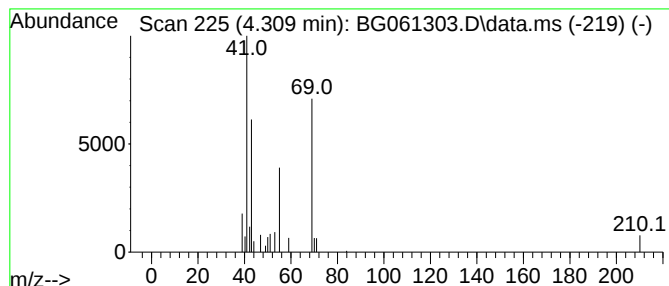
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.309	2.09 ng/ul	15377	1,4-Dichlorobenzene-d4	7.888

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentene, 3-methyl-, (Z)-	84	C6H12	000922-62-3	53
2		2-Pentene, 3-methyl-, (Z)-	84	C6H12	000922-62-3	53
3		2-Pentene, 3-methyl-, (E)-	84	C6H12	000616-12-6	53
4		1-Pentene, 2,3-dimethyl-	98	C7H14	003404-72-6	45
5		2-Pentene, 3-methyl-, (E)-	84	C6H12	000616-12-6	42



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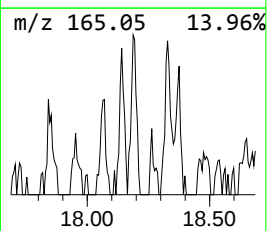
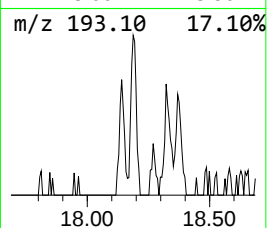
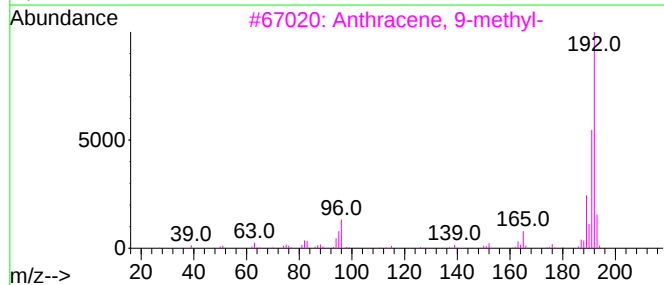
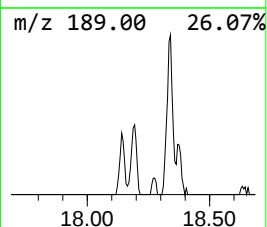
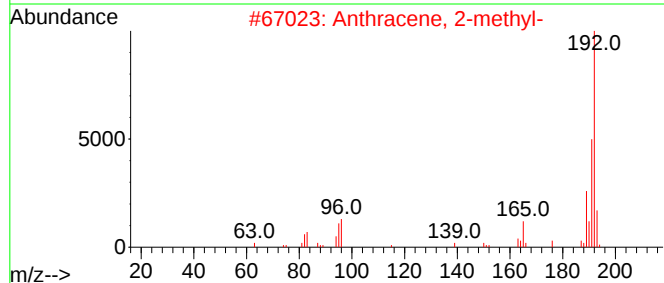
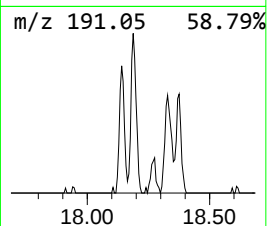
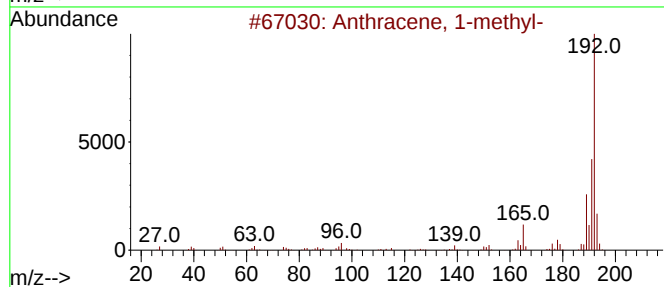
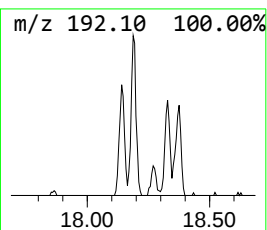
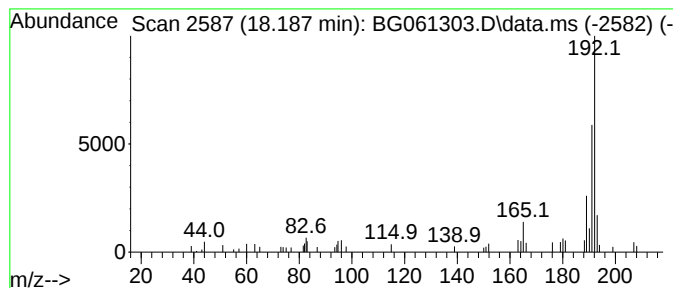
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.187	2.22 ng/ul	52689	Phenanthrene-d10	17.265

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 1-methyl-	192	C15H12	000610-48-0	95
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	90
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052824\
Data File : BG061303.D
Acq On : 28 May 2024 17:19
Operator : MA/JU
Sample : P2568-22
Misc :
ALS Vial : 11 Sample Multiplier: 1

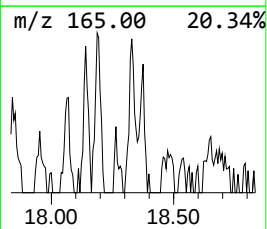
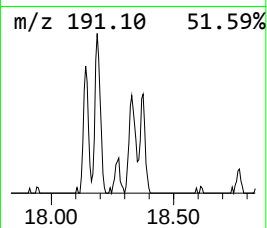
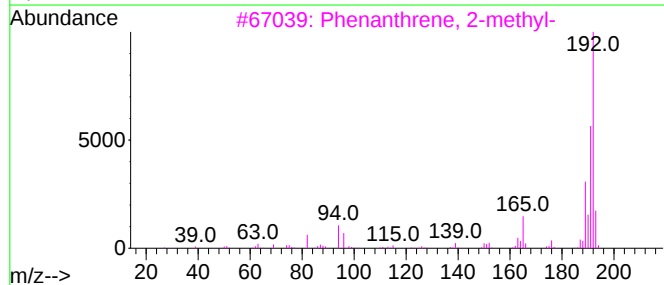
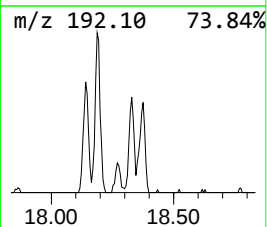
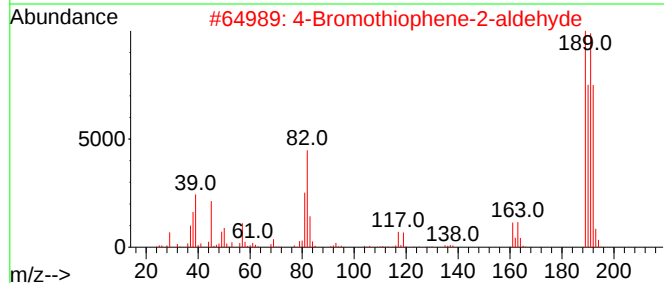
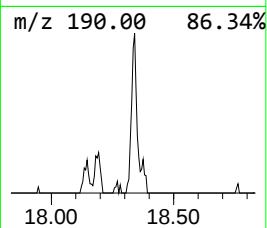
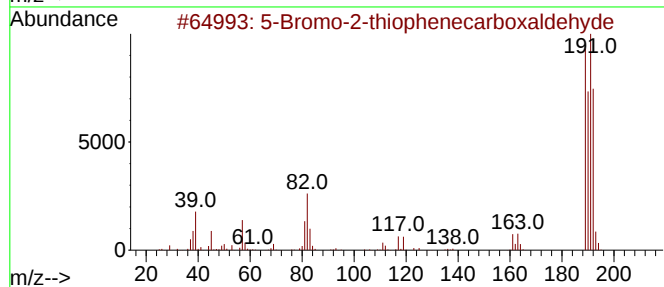
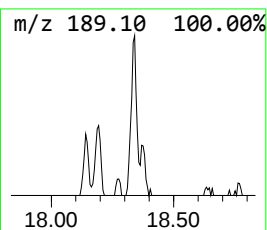
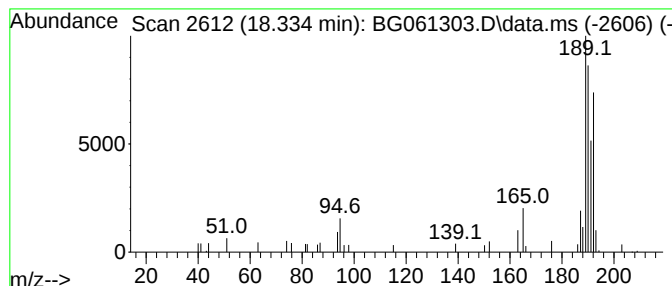
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 5-Bromo-2-thiophenecarboxal... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.334	2.06 ng/ul	48776	Phenanthrene-d10	17.265	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Bromo-2-thiophenecarboxaldehyde	190	C5H3BrOS	004701-17-1	53
2	4-Bromothiophene-2-aldehyde	190	C5H3BrOS	018791-75-8	53
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	50
4	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052824\
Data File : BG061303.D
Acq On : 28 May 2024 17:19
Operator : MA/JU
Sample : P2568-22
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ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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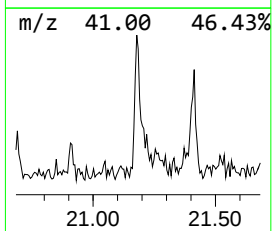
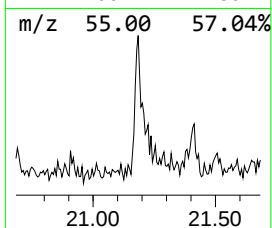
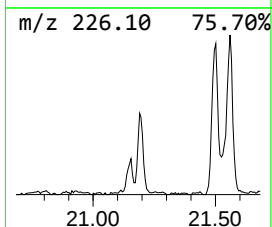
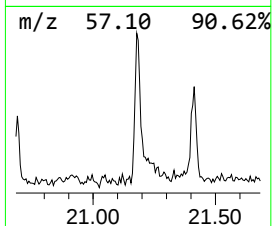
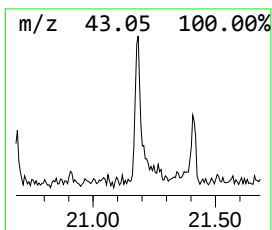
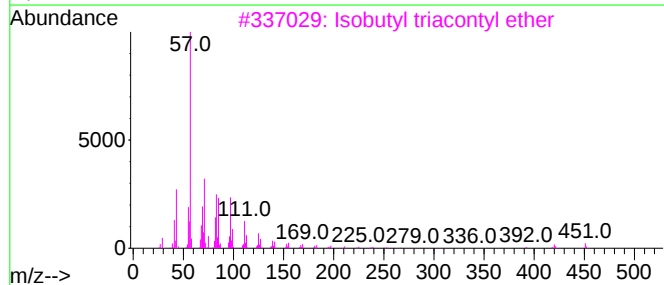
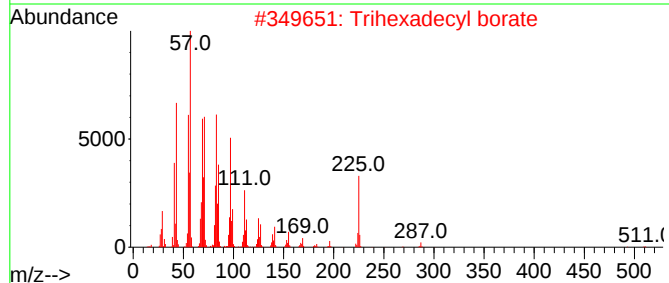
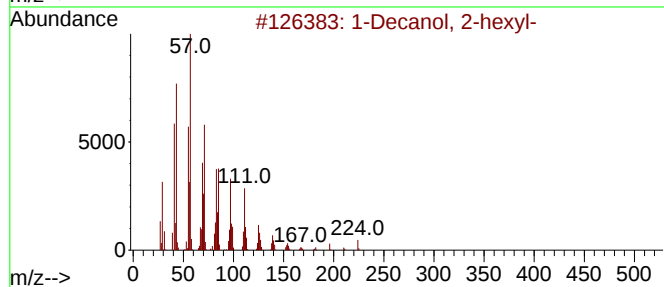
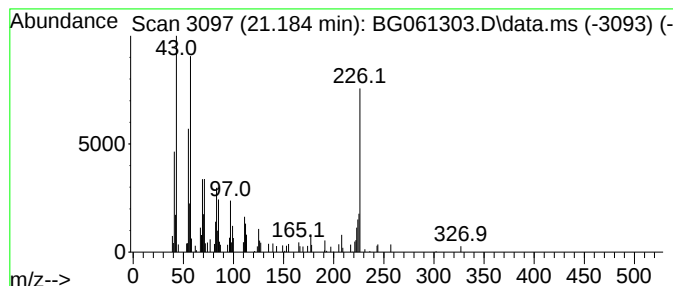
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.184	3.40 ng/ul	158339	Chrysene-d12	21.513

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Decanol, 2-hexyl-			242	C16H34O	002425-77-6	20
2	Trihexadecyl borate			735	C48H99BO3	002665-11-4	18
3	Isobutyl triacontyl ether			495	C34H70O	1000406-33-5	15
4	Octatriacontyl pentafluoropropio...			697	C41H77F5O2	1000351-89-1	15
5	Isobutyl tetratriacontyl ether			551	C38H78O	1000406-33-7	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052824\
Data File : BG061303.D
Acq On : 28 May 2024 17:19
Operator : MA/JU
Sample : P2568-22
Misc :
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Instrument :
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ClientSampleId :
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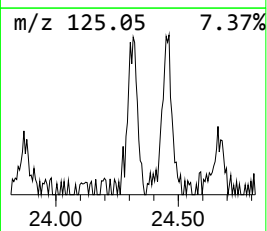
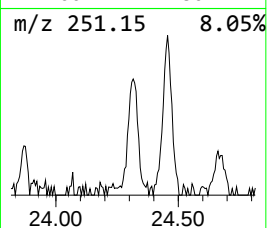
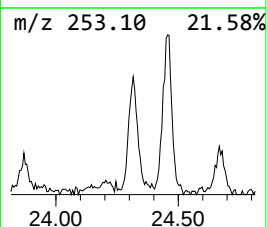
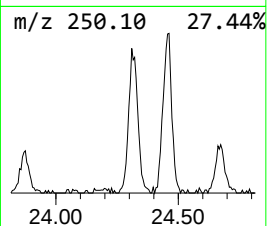
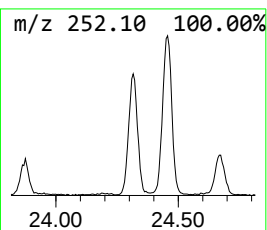
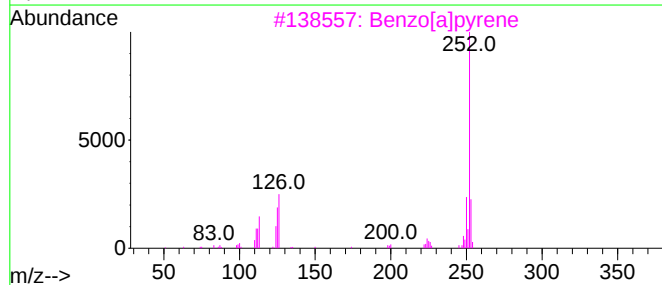
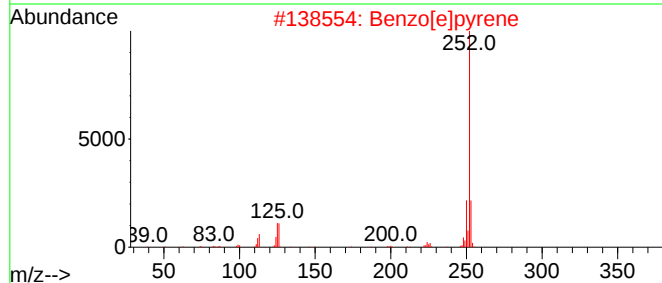
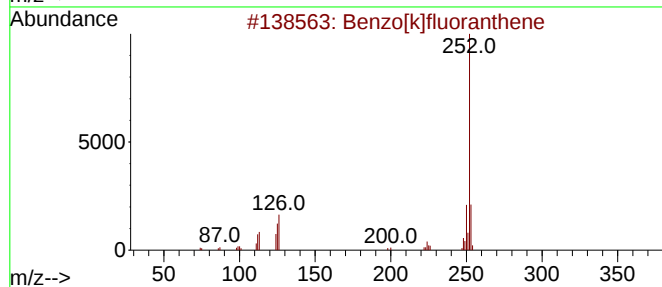
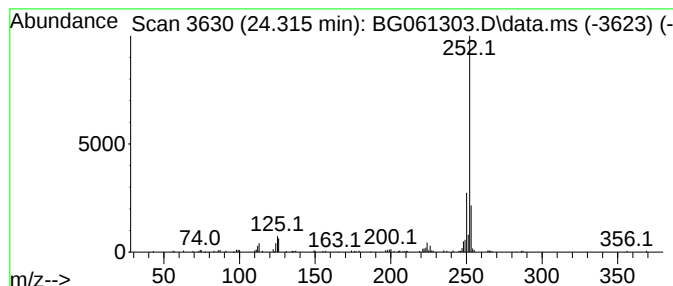
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.316	6.34 ng/ul	195114	Perylene-d12	24.598

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
2			Benzo[e]pyrene	252	C20H12	000192-97-2	93
3			Benzo[a]pyrene	252	C20H12	000050-32-8	90
4			Benzo[b]fluoranthene	252	C20H12	000205-99-2	90
5			Benzo[j]fluoranthene	252	C20H12	000205-82-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052824\
Data File : BG061303.D
Acq On : 28 May 2024 17:19
Operator : MA/JU
Sample : P2568-22
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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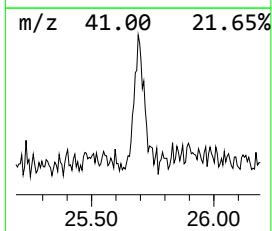
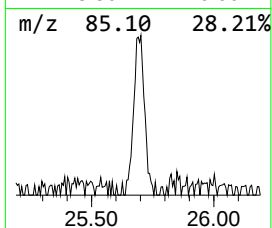
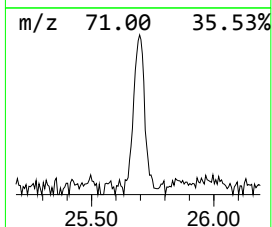
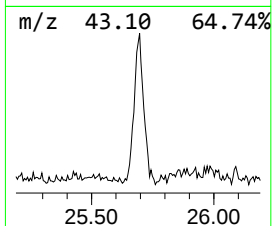
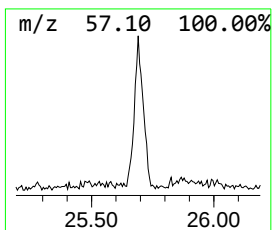
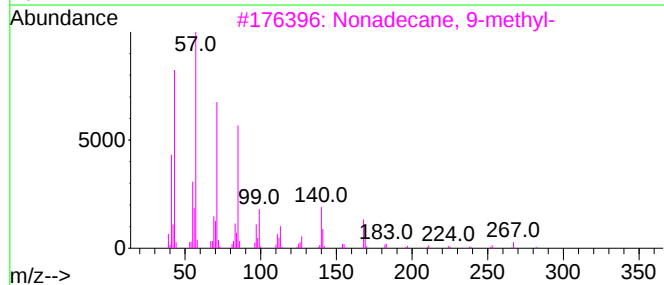
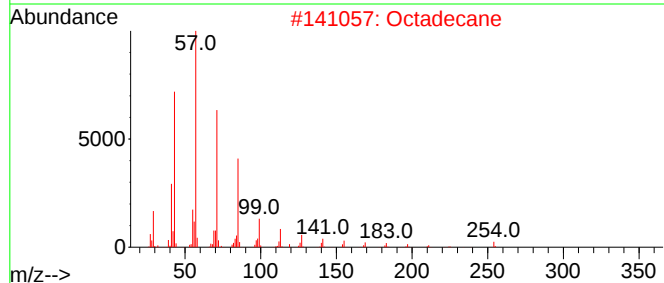
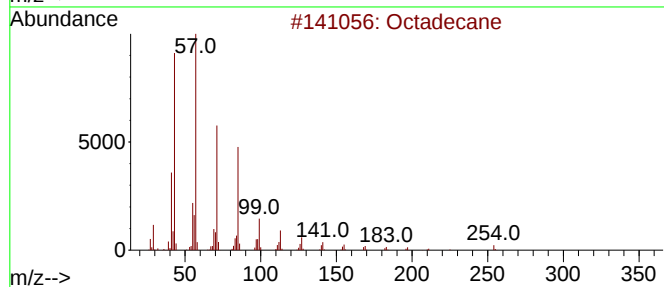
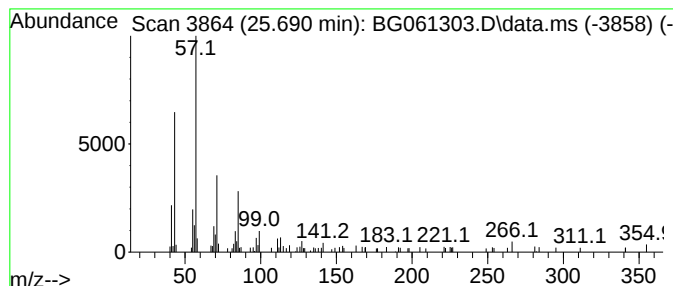
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.690	4.13 ng/ul	127111	Perylene-d12	24.598

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	93
2			Octadecane	254	C18H38	000593-45-3	91
3			Nonadecane, 9-methyl-	282	C20H42	013287-24-6	83
4			Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	68
5			Octadecane	254	C18H38	000593-45-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052824\
Data File : BG061303.D
Acq On : 28 May 2024 17:19
Operator : MA/JU
Sample : P2568-22
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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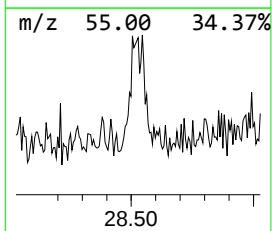
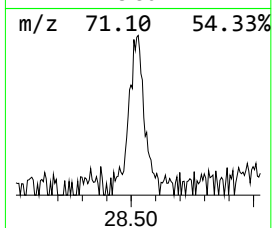
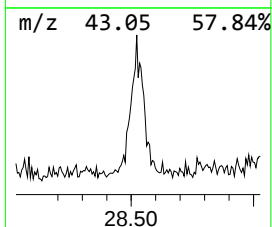
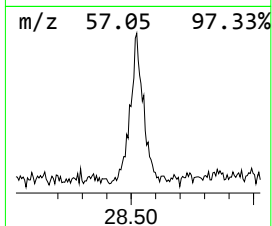
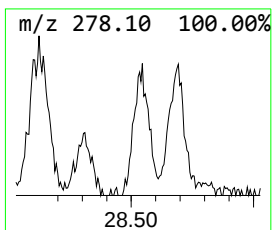
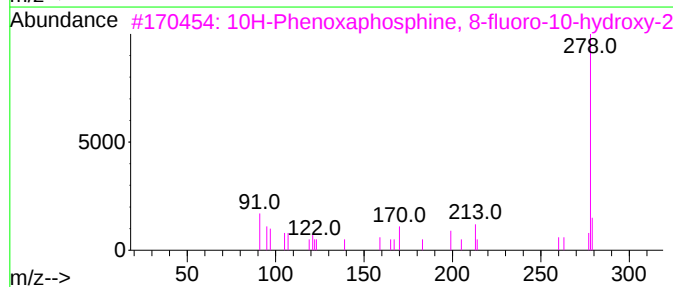
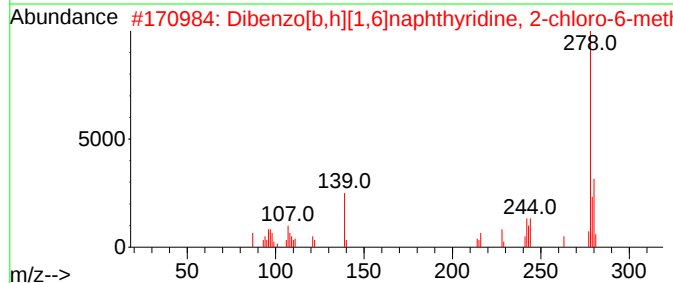
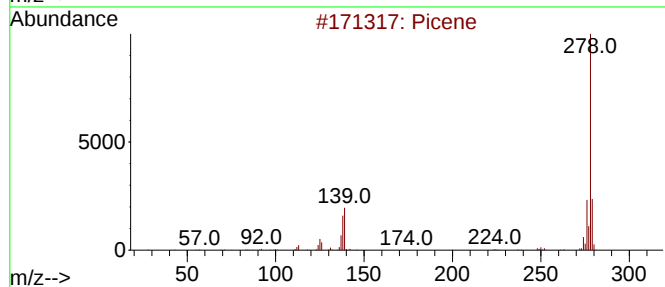
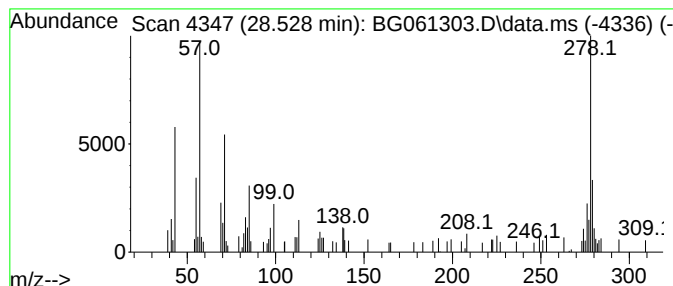
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Picene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.528	3.82 ng/ul	117638	Perylene-d12	24.598

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Picene	278	C22H14	000213-46-7	58
2			Dibenzo[b,h][1,6]naphthyridine, ...	278	C17H11ClN2	004240-91-9	46
3			10H-Phenoxaphosphine, 8-fluoro-1...	278	C14H12FO3P	037041-13-7	43
4			Curan, 16,17,19,20-tetradecydro-	278	C19H22N2	056053-17-9	43
5			Benzo[b]triphenylene	278	C22H14	000215-58-7	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052824\
Data File : BG061303.D
Acq On : 28 May 2024 17:19
Operator : MA/JU
Sample : P2568-22
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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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\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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(DEL) Alkane: S...	4.309	2.1	ng/u1	15377	1	7.888	146949	20.0
Anthracene, 1-m...	18.187	2.2	ng/u1	52689	4	17.265	474619	20.0
5-Bromo-2-thiop...	18.334	2.1	ng/u1	48776	4	17.265	474619	20.0
unknown-01	21.184	3.4	ng/u1	158339	5	21.513	930180	20.0
Benzo[e]pyrene	24.316	6.3	ng/u1	195114	6	24.598	615770	20.0
(DEL) Alkane: S...	25.690	4.1	ng/u1	127111	6	24.598	615770	20.0
Picene	28.528	3.8	ng/u1	117638	6	24.598	615770	20.0