

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053024\
 Data File : BG061382.D
 Acq On : 31 May 2024 6:49
 Operator : MA/JU
 Sample : P2570-10
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BH6B9

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: BG061382.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.035	4	8	10	rBV	16396	21686	2.14%	0.149%
2	3.076	10	15	30	rVB	257279	476569	47.03%	3.282%
3	3.335	54	59	66	rVB3	7000	10424	1.03%	0.072%
4	3.605	102	105	114	rVV2	3857	7370	0.73%	0.051%
5	3.746	124	129	136	rBV	22906	39288	3.88%	0.271%
6	4.298	219	223	228	rVB	8853	12291	1.21%	0.085%
7	5.955	501	505	512	rBV4	3537	6398	0.63%	0.044%
8	7.060	686	693	705	rVB	131854	222306	21.94%	1.531%
9	7.207	710	718	725	rBV	86777	140073	13.82%	0.965%
10	7.412	746	753	759	rBV	133264	214280	21.14%	1.476%
11	7.871	822	831	838	rBV	112077	185368	18.29%	1.277%
12	8.593	947	954	963	rBV	147688	247564	24.43%	1.705%
13	9.028	1021	1028	1036	rBV	138651	236767	23.36%	1.630%
14	9.586	1117	1123	1129	rBV2	12998	20218	2.00%	0.139%
15	9.751	1144	1151	1159	rVB2	119789	207232	20.45%	1.427%
16	10.303	1238	1245	1256	rVB	166399	294721	29.08%	2.030%
17	10.438	1260	1268	1279	rBV	32823	55208	5.45%	0.380%
18	10.667	1299	1307	1314	rBV	144789	253643	25.03%	1.747%
19	10.802	1322	1330	1339	rBV2	119548	211001	20.82%	1.453%
20	11.202	1393	1398	1402	rVB2	4453	6538	0.65%	0.045%
21	11.408	1426	1433	1440	rBV3	11478	22948	2.26%	0.158%
22	13.916	1853	1860	1868	rVV	310250	438296	43.25%	3.018%
23	14.204	1902	1909	1918	rVB2	325869	518465	51.16%	3.570%
24	14.510	1952	1961	1967	rBV	247411	360050	35.53%	2.479%
25	14.745	1996	2001	2012	rVB2	63442	125458	12.38%	0.864%
26	15.503	2122	2130	2136	rBV	438534	657757	64.91%	4.530%
27	15.638	2148	2153	2161	rBV	115519	169663	16.74%	1.168%
28	16.913	2365	2370	2376	rVB3	9718	14853	1.47%	0.102%
29	17.077	2394	2398	2402	rBV2	4352	7012	0.69%	0.048%
30	17.260	2422	2429	2433	rBV	295305	446512	44.06%	3.075%
31	17.301	2433	2436	2440	rVV	118084	155607	15.36%	1.072%
32	17.359	2440	2446	2457	rVB	472445	718894	70.94%	4.951%
33	17.659	2489	2497	2504	rBV2	12317	27583	2.72%	0.190%
34	17.835	2524	2527	2534	rVV4	4570	7785	0.77%	0.054%
35	18.135	2569	2578	2580	rBV2	25050	40031	3.95%	0.276%
36	18.158	2580	2582	2584	rVV2	35561	45737	4.51%	0.315%

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

37	18.182	2584	2586	2589	rVV	37337	48195	4.76%	0.332%
38	18.223	2589	2593	2598	rVV	48709	63597	6.28%	0.438%
39	18.270	2598	2601	2604	rVB2	6838	9053	0.89%	0.062%
40	18.329	2604	2611	2615	rBV4	29006	55571	5.48%	0.383%
41	18.364	2615	2617	2623	rVB3	18833	22618	2.23%	0.156%
42	18.487	2635	2638	2641	rVV4	8059	10592	1.05%	0.073%
43	18.529	2641	2645	2649	rVV5	5010	9096	0.90%	0.063%
44	18.634	2658	2663	2667	rBV	22416	30551	3.01%	0.210%
45	18.681	2667	2671	2675	rVB	30878	44419	4.38%	0.306%
46	18.881	2702	2705	2708	rVB3	7259	8706	0.86%	0.060%
47	18.934	2710	2714	2717	rVB2	7966	9391	0.93%	0.065%
48	18.969	2717	2720	2724	rVB6	7587	9056	0.89%	0.062%
49	19.052	2730	2734	2741	rVV7	17976	42453	4.19%	0.292%
50	19.110	2742	2744	2747	rVV4	8157	10401	1.03%	0.072%
51	19.146	2747	2750	2753	rVB4	7806	10288	1.02%	0.071%
52	19.198	2753	2759	2767	rBV3	26591	56482	5.57%	0.389%
53	19.316	2771	2779	2785	rVV	345903	483827	47.74%	3.332%
54	19.375	2787	2789	2792	rVV4	10169	10373	1.02%	0.071%
55	19.428	2792	2798	2802	rVV7	14466	35749	3.53%	0.246%
56	19.463	2802	2804	2809	rVV2	16083	22604	2.23%	0.156%
57	19.516	2809	2813	2816	rVV6	13113	24065	2.37%	0.166%
58	19.557	2818	2820	2824	rVV	12864	19576	1.93%	0.135%
59	19.586	2824	2825	2830	rVV5	8714	7789	0.77%	0.054%
60	19.651	2830	2836	2839	rVV2	672324	1013394	100.00%	6.979%
61	19.680	2839	2841	2848	rVV	311498	370463	36.56%	2.551%
62	19.751	2848	2853	2859	rVB2	20240	32868	3.24%	0.226%
63	19.862	2867	2872	2877	rBV	17231	28083	2.77%	0.193%
64	19.915	2878	2881	2887	rVB3	11420	16962	1.67%	0.117%
65	20.003	2890	2896	2898	rBV3	28394	54451	5.37%	0.375%
66	20.139	2915	2919	2921	rVV4	21972	35250	3.48%	0.243%
67	20.174	2921	2925	2932	rVB2	37162	79527	7.85%	0.548%
68	20.274	2939	2942	2945	rVB	17698	19825	1.96%	0.137%
69	20.332	2945	2952	2955	rBV2	36327	62939	6.21%	0.433%
70	20.473	2972	2976	2980	rBV2	20800	32126	3.17%	0.221%
71	20.515	2980	2983	2986	rVV3	24182	35338	3.49%	0.243%
72	20.567	2987	2992	2997	rVB	173973	212094	20.93%	1.461%
73	20.685	3008	3012	3015	rVB3	28397	32023	3.16%	0.221%
74	20.726	3017	3019	3022	rVB4	13431	12634	1.25%	0.087%
75	20.926	3048	3053	3060	rVB	49349	76751	7.57%	0.529%
76	21.102	3078	3083	3087	rBV3	30274	61782	6.10%	0.425%

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77	21.149	3087	3091	3092	rVV	34166	47441	4.68%	0.327%
78	21.178	3092	3096	3104	rVV4	77590	202202	19.95%	1.392%
79	21.249	3106	3108	3115	rVB	51944	80717	7.97%	0.556%
80	21.408	3130	3135	3141	rVB	186294	243888	24.07%	1.680%
81	21.507	3144	3152	3157	rVV2	396537	871757	86.02%	6.003%
82	21.554	3157	3160	3170	rVB	167176	267991	26.44%	1.846%
83	21.701	3182	3185	3190	rVB3	39148	56033	5.53%	0.386%
84	21.760	3193	3195	3201	rVB5	16241	26553	2.62%	0.183%
85	22.218	3269	3273	3278	rVB3	31209	50148	4.95%	0.345%
86	22.283	3280	3284	3289	rVV6	45678	72225	7.13%	0.497%
87	22.612	3338	3340	3341	rBV2	11627	8649	0.85%	0.060%
88	22.923	3387	3393	3403	rVB5	87943	214360	21.15%	1.476%
89	23.488	3487	3489	3496	rBV7	9685	20938	2.07%	0.144%
90	23.623	3505	3512	3519	rBV3	122988	376969	37.20%	2.596%
91	23.864	3550	3553	3560	rVB2	23113	52229	5.15%	0.360%
92	24.310	3623	3629	3633	rBV2	58804	130668	12.89%	0.900%
93	24.386	3635	3642	3648	rVV2	348135	881570	86.99%	6.071%
94	24.451	3650	3653	3666	rVB	112822	241754	23.86%	1.665%
95	24.598	3669	3678	3687	rBV	215993	570880	56.33%	3.931%
96	25.591	3845	3847	3848	rBV2	9262	6018	0.59%	0.041%
97	26.684	4031	4033	4036	rBV3	7258	8853	0.87%	0.061%
98	28.100	4261	4274	4288	rVB4	41556	200465	19.78%	1.380%
99	29.163	4449	4455	4456	rBV5	20366	33617	3.32%	0.232%
100	30.285	4644	4646	4648	rBV3	6663	8791	0.87%	0.061%

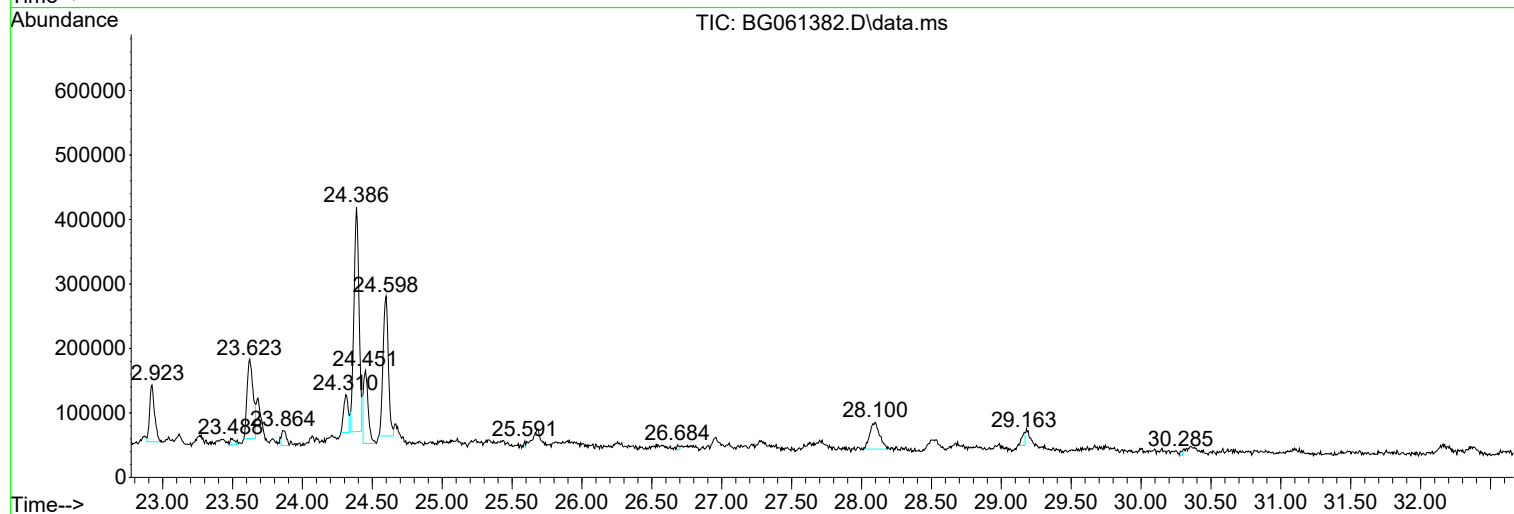
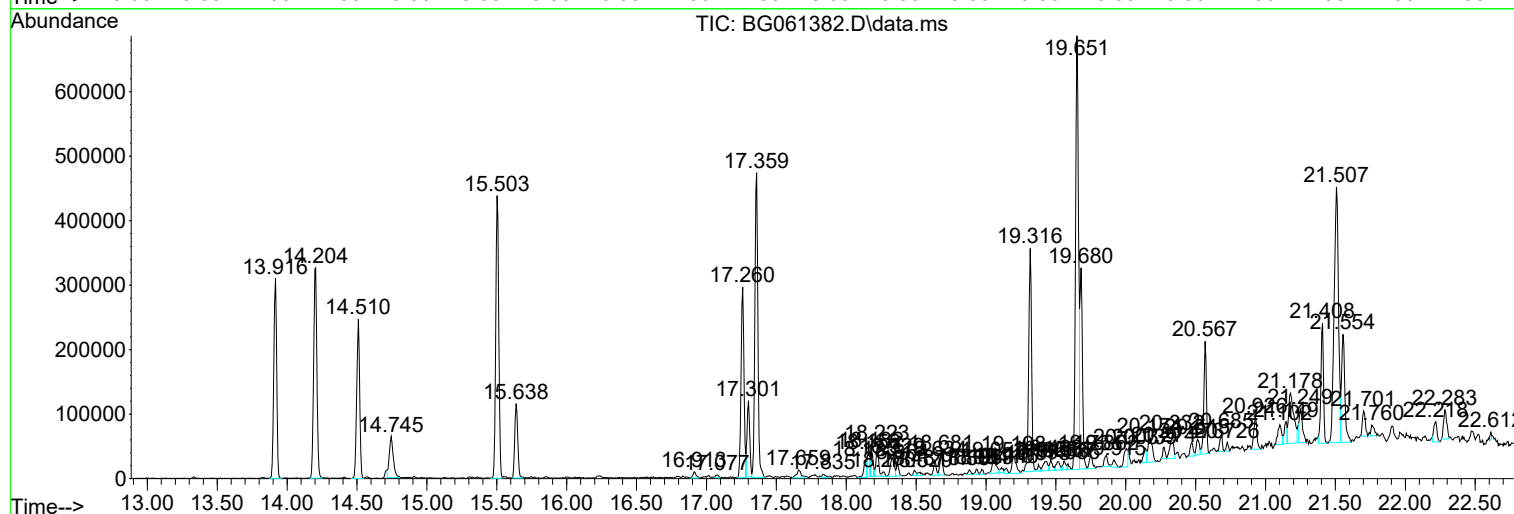
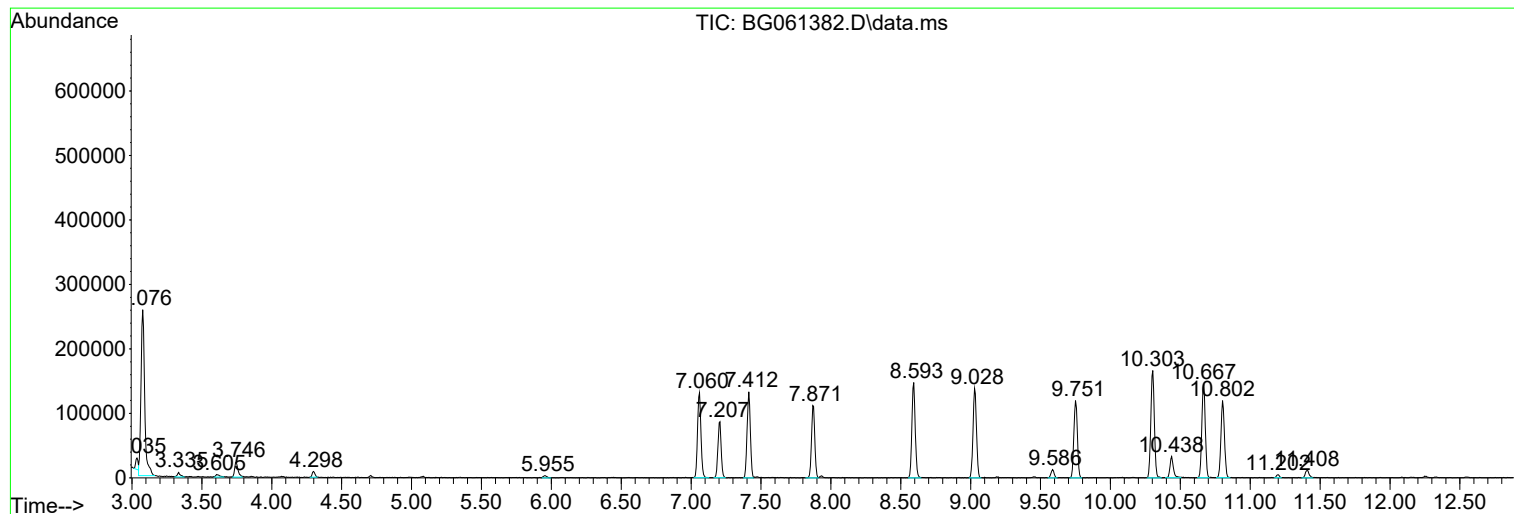
Sum of corrected areas: 14521274

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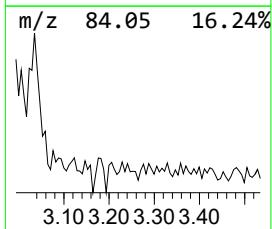
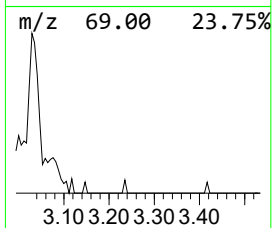
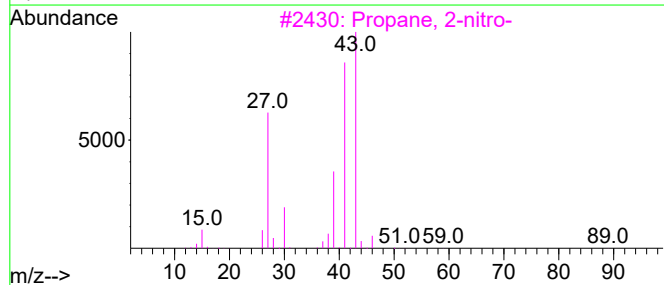
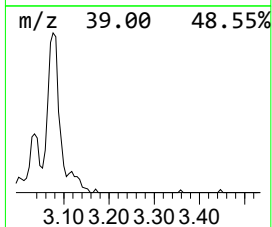
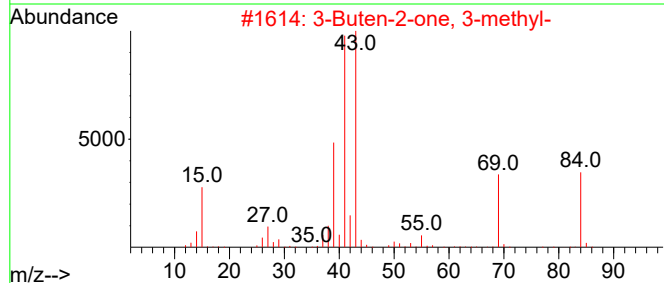
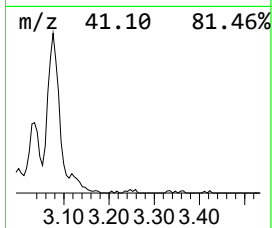
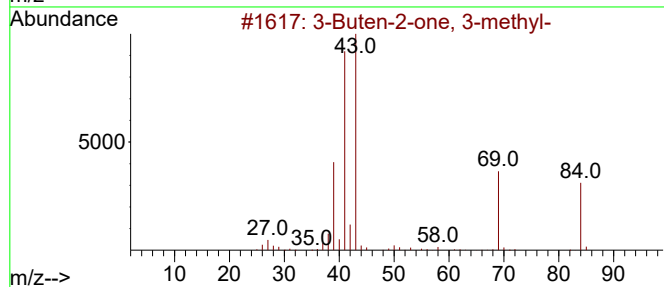
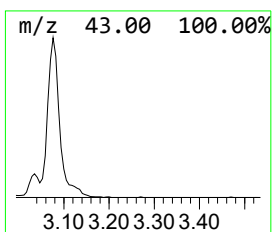
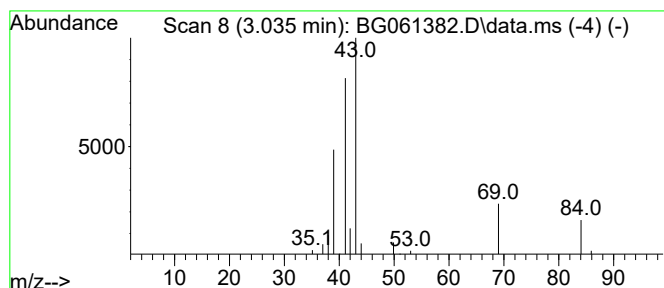
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 3-Buten-2-one, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.035	2.34 ng/ul	21686	1,4-Dichlorobenzene-d4	7.876

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Buten-2-one, 3-methyl-	84	C5H8O	000814-78-8	90
2		3-Buten-2-one, 3-methyl-	84	C5H8O	000814-78-8	83
3		Propane, 2-nitro-	89	C3H7NO2	000079-46-9	38
4		4-Penten-2-one	84	C5H8O	013891-87-7	33
5		Ethanone, 1-cyclopropyl-	84	C5H8O	000765-43-5	25



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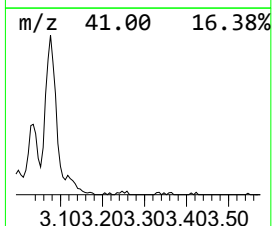
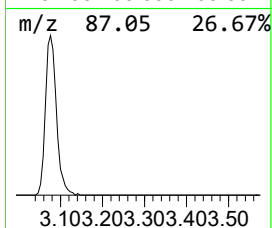
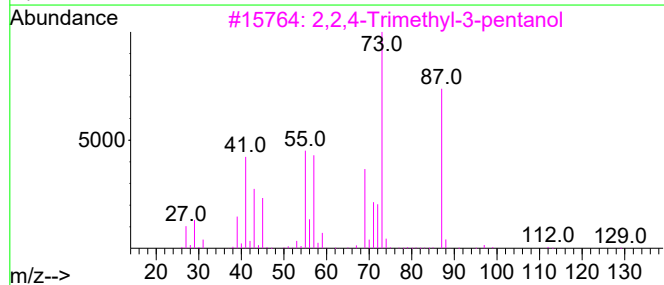
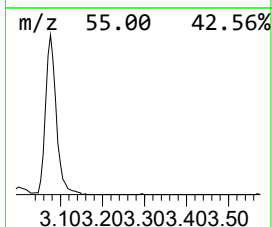
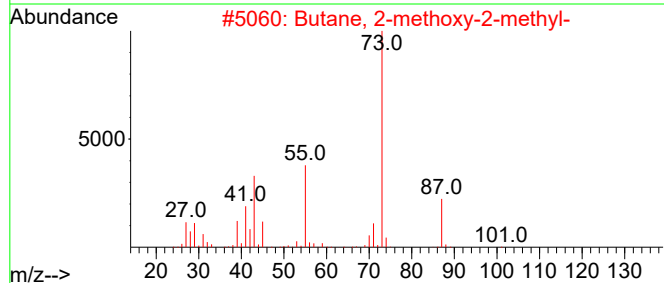
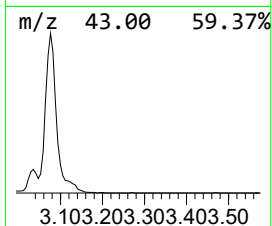
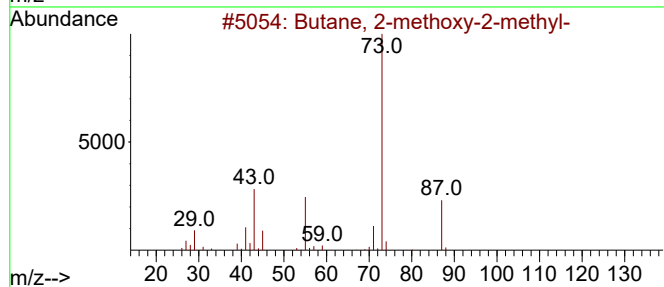
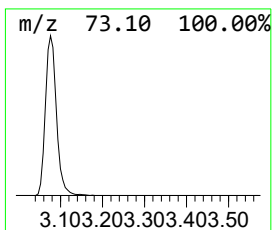
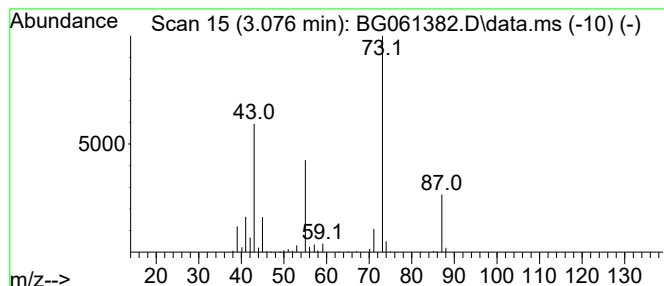
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.076	51.42 ng/ul	476569	1,4-Dichlorobenzene-d4	7.876

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	83
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4			3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	38
5			3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	28



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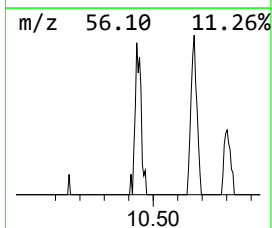
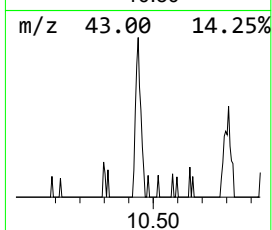
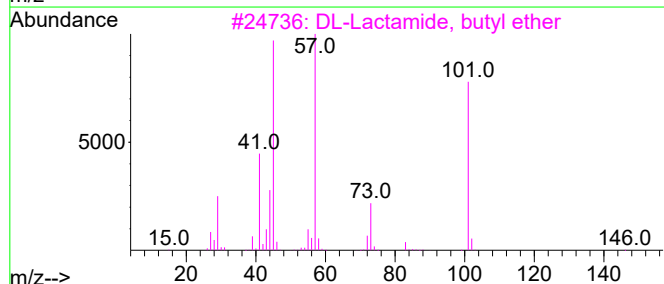
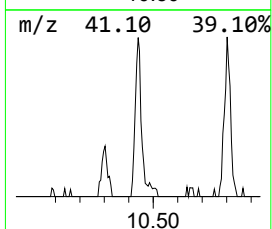
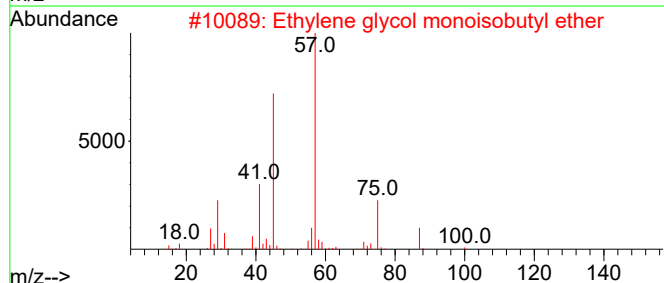
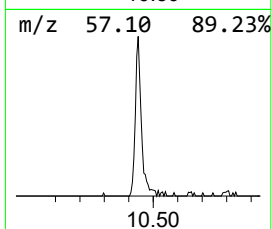
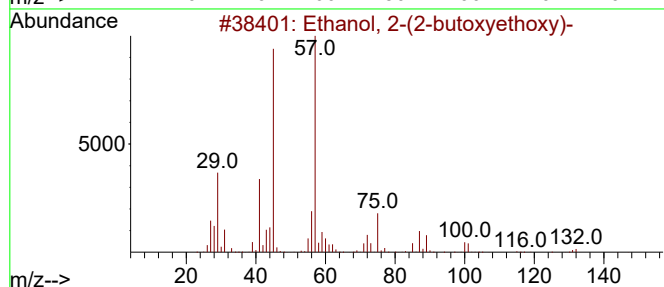
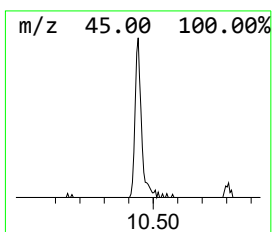
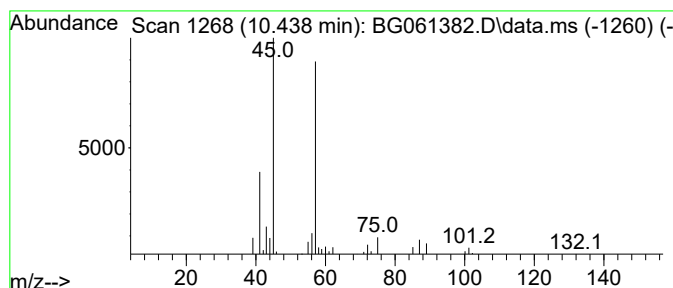
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.438	4.35 ng/ul	55208	Naphthalene-d8	10.667

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	64
2			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	59
3			DL-Lactamide, butyl ether	145	C7H15NO2	1000452-56-5	59
4			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	56
5			Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053024\
Data File : BG061382.D
Acq On : 31 May 2024 6:49
Operator : MA/JU
Sample : P2570-10
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH6B9

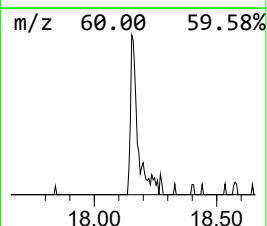
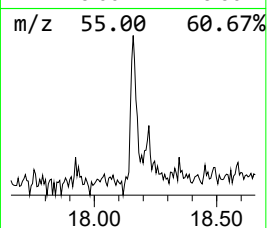
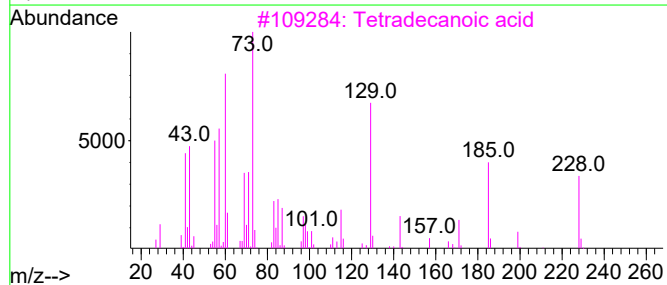
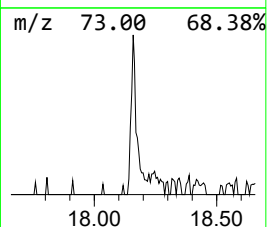
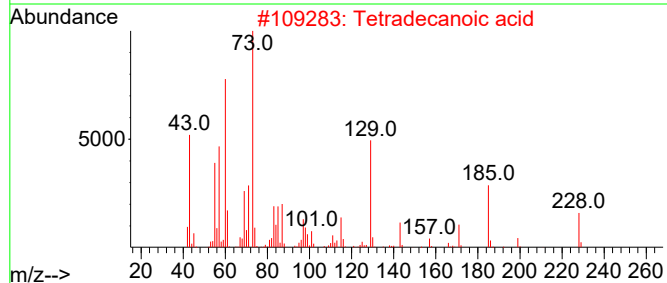
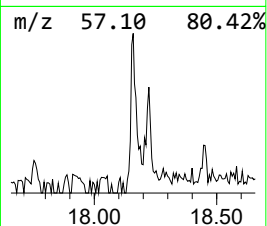
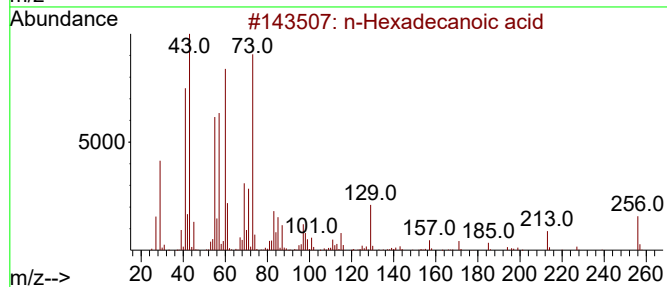
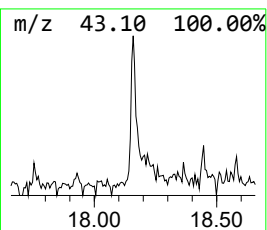
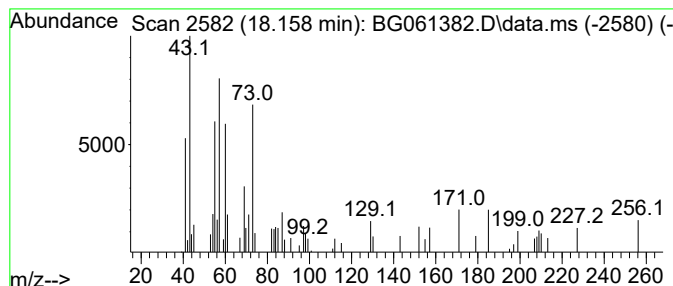
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.159	2.05 ng/ul	45737	Phenanthrene-d10	17.259

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	56
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	35
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	35
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	22
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	12



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053024\
Data File : BG061382.D
Acq On : 31 May 2024 6:49
Operator : MA/JU
Sample : P2570-10
Misc :
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Instrument :
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ClientSampleId :
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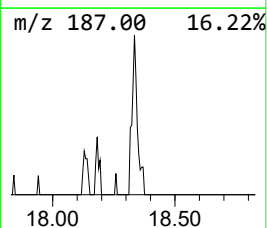
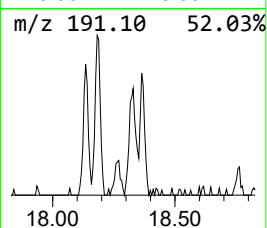
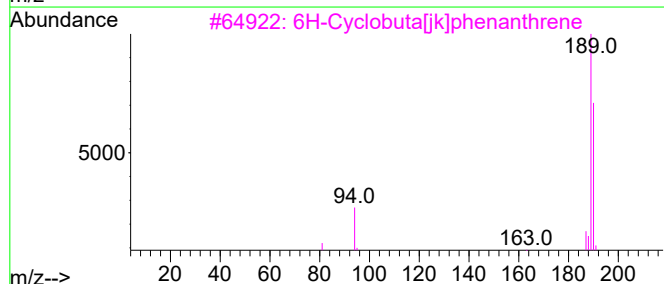
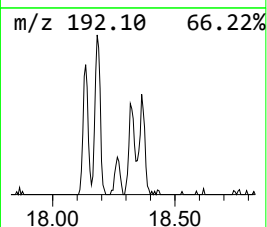
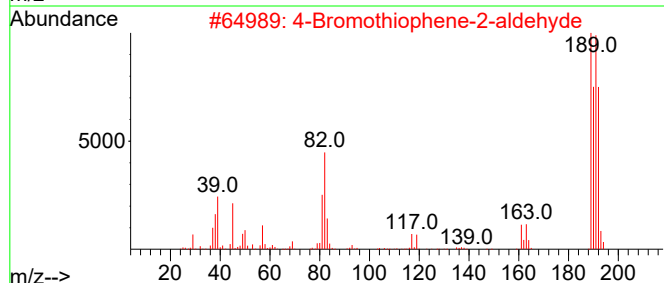
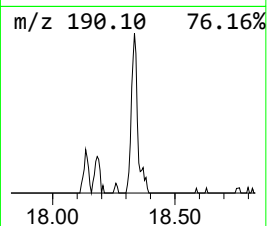
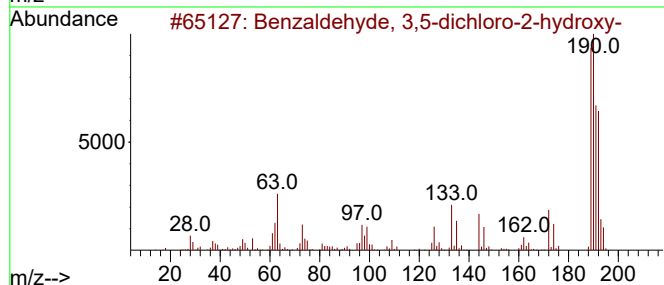
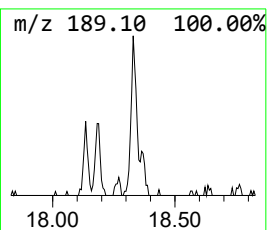
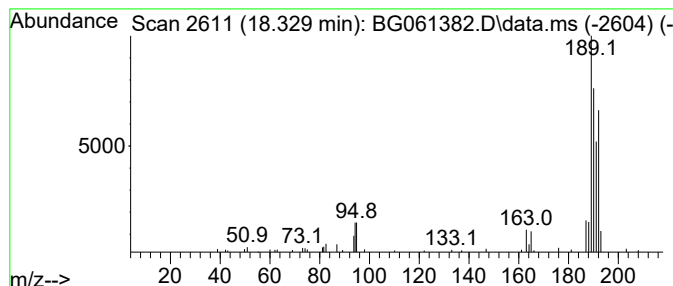
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzaldehyde, 3,5-dichloro-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.329	2.49 ng/ul	55571	Phenanthrene-d10	17.259

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	64
2			4-Bromothiophene-2-aldehyde	190	C5H3BrOS	018791-75-8	58
3			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	58
4			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	43
5			5-Amino-3-bromo-2-fluoropyridine	190	C5H4BrFN2	209328-99-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053024\
Data File : BG061382.D
Acq On : 31 May 2024 6:49
Operator : MA/JU
Sample : P2570-10
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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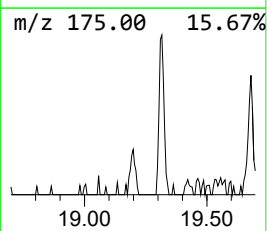
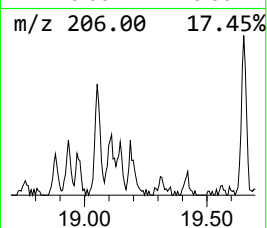
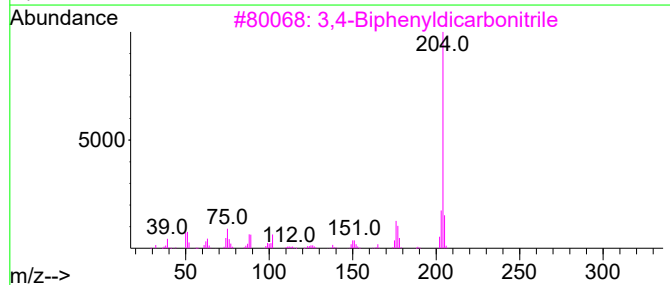
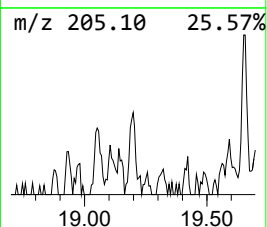
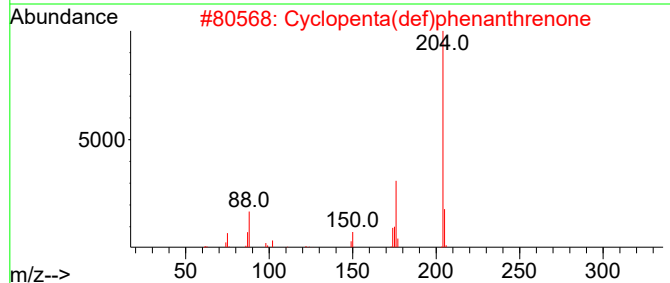
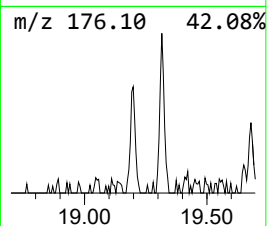
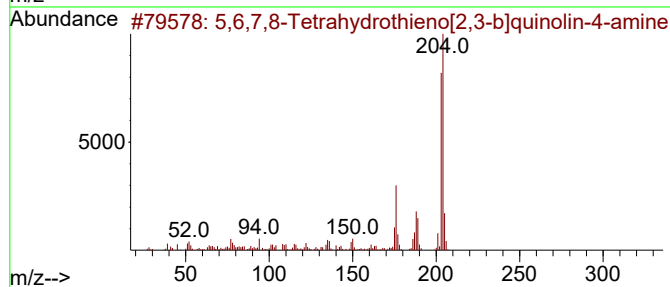
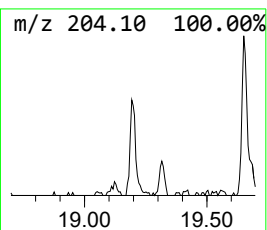
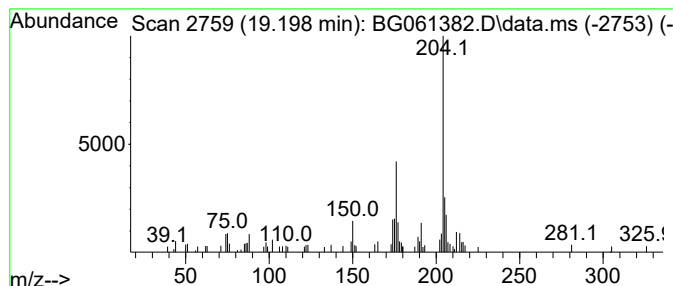
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 5,6,7,8-Tetrahydrothieno[2,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.198	2.53 ng/ul	56482	Phenanthrene-d10	17.259

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,6,7,8-Tetrahydrothieno[2,3-b]q...	204	C11H12N2S	122914-50-5	59	
2	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	58	
3	3,4-Biphenyldicarbonitrile	204	C14H8N2	004128-63-6	50	
4	[1,2,4]Triazolo[5,1-b]quinazolin...	204	C10H12N4O	1000337-56-5	50	
5	1-Phenyl-1,2,3,4-tetrahydroisoqu...	281	C18H23NSi	1000484-76-2	47	



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053024\
Data File : BG061382.D
Acq On : 31 May 2024 6:49
Operator : MA/JU
Sample : P2570-10
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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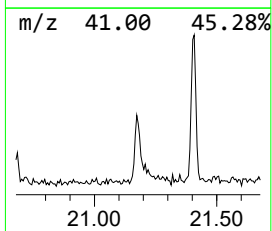
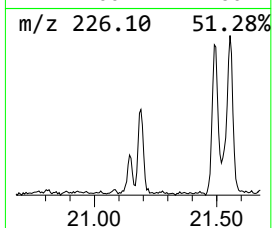
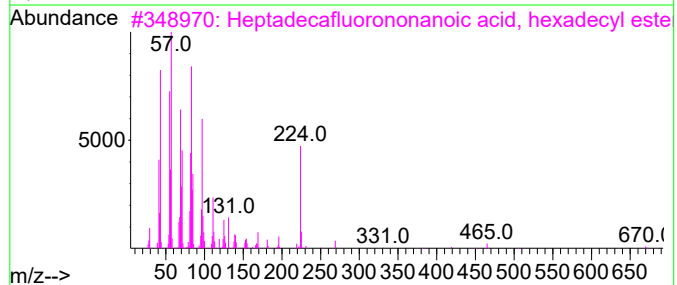
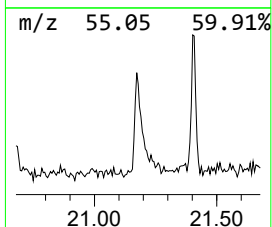
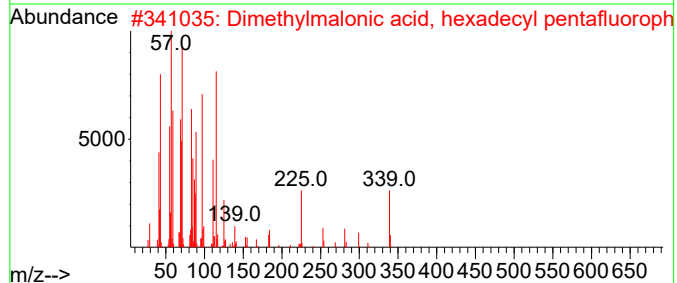
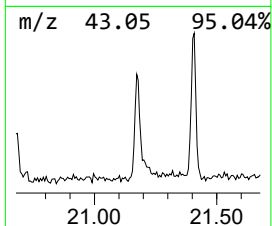
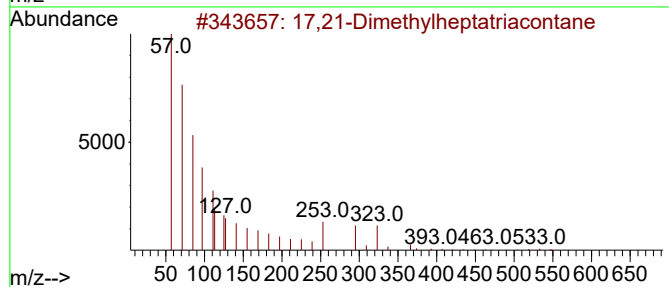
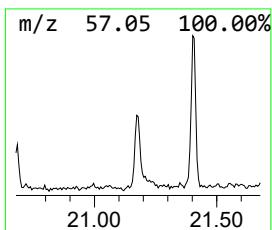
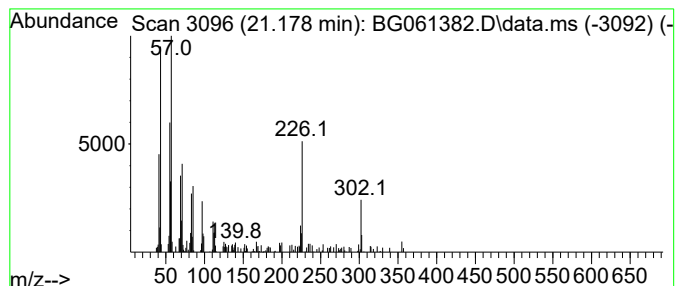
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.179	4.64 ng/ul	202202	Chrysene-d12	21.513

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	17,21-Dimethylheptatriacontane	549	C39H80	067979-79-7	27		
2	Dimethylmalonic acid, hexadecyl ...	522	C27H39F5O4	1000363-67-3	27		
3	Heptadecafluorononanoic acid, he...	688	C25H33F17O2	1000356-03-2	16		
4	Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	14		
5	Dihydrophytol, TMS derivative	370	C23H50OSi	078695-20-2	14		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053024\
Data File : BG061382.D
Acq On : 31 May 2024 6:49
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Sample : P2570-10
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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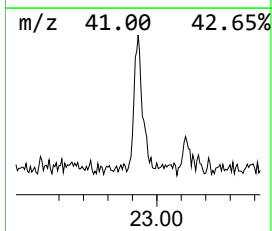
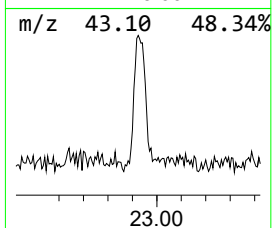
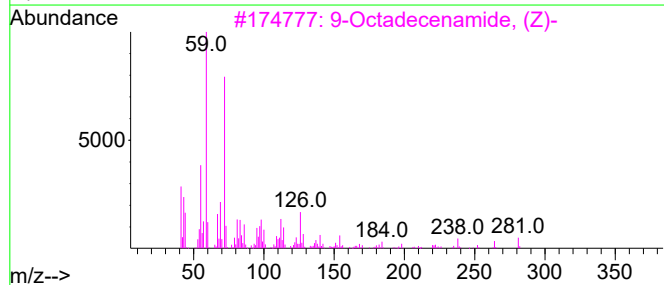
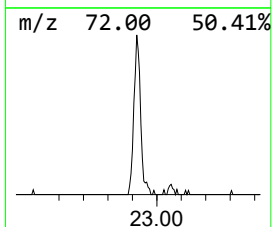
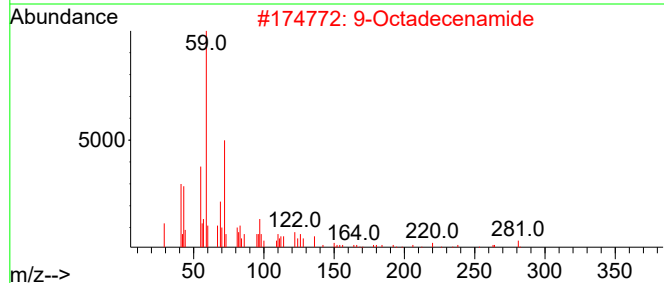
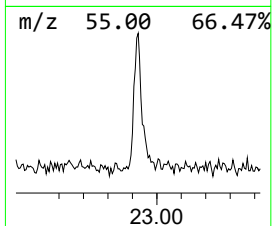
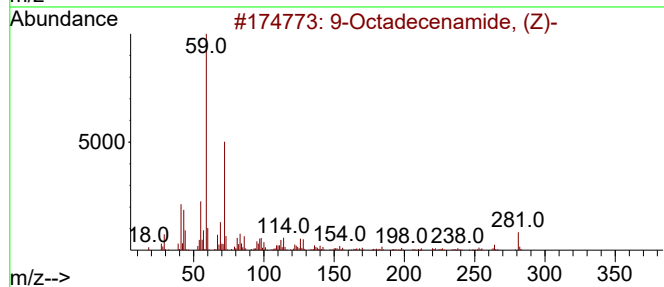
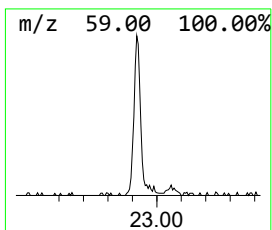
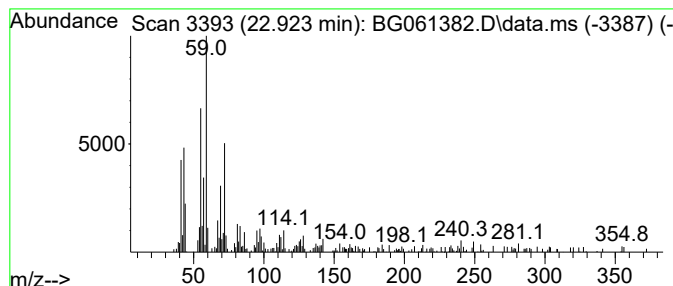
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 9-Octadecenamide, (Z)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.924	4.92 ng/ul	214360	Chrysene-d12	21.513

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	95
2	9-Octadecenamide			281	C18H35NO	003322-62-1	90
3	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	83
4	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	81
5	Tetradecanamide			227	C14H29NO	000638-58-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053024\
Data File : BG061382.D
Acq On : 31 May 2024 6:49
Operator : MA/JU
Sample : P2570-10
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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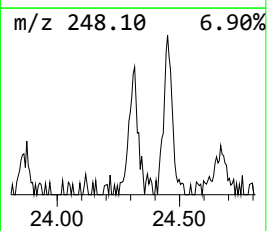
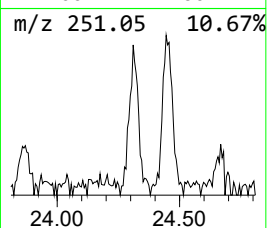
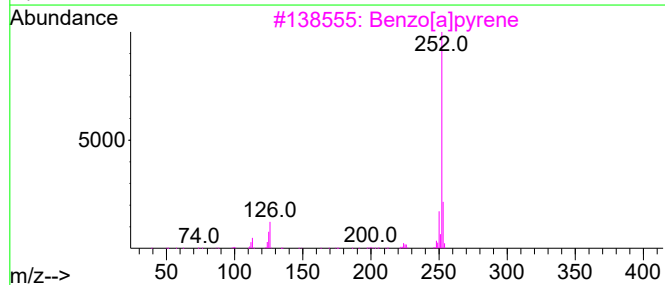
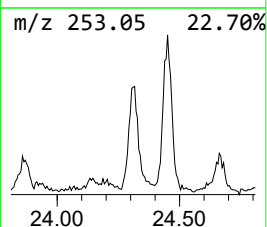
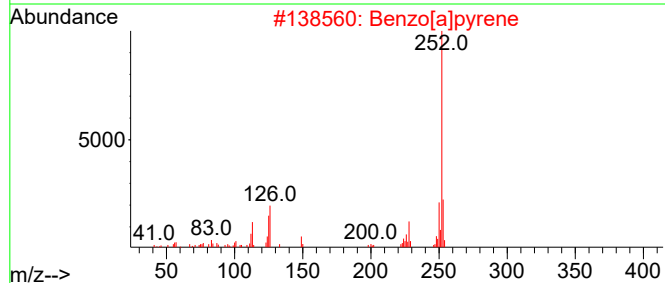
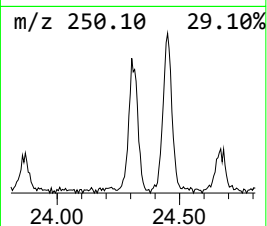
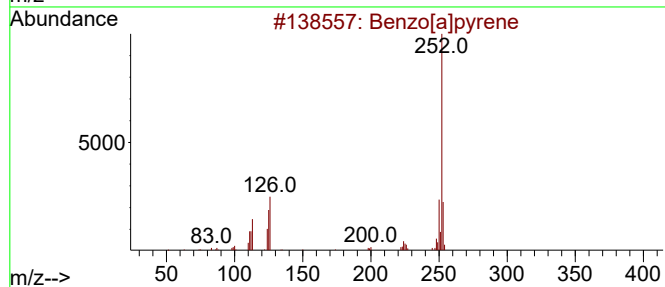
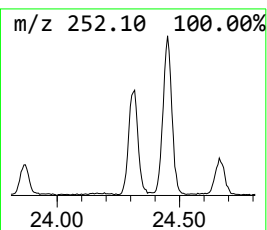
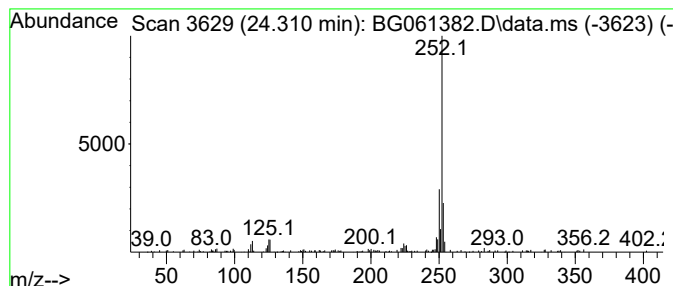
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Perylene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.310	4.58 ng/ul	130668	Perylene-d12	24.598

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[a]pyrene	252	C20H12	000050-32-8	96
2			Benzo[a]pyrene	252	C20H12	000050-32-8	93
3			Benzo[a]pyrene	252	C20H12	000050-32-8	90
4			Perylene	252	C20H12	000198-55-0	90
5			Benzo[k]fluoranthene	252	C20H12	000207-08-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053024\
Data File : BG061382.D
Acq On : 31 May 2024 6:49
Operator : MA/JU
Sample : P2570-10
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH6B9

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
3-Buten-2-one, ...	3.035	2.3	ng/ul	21686	1	7.876	185368	20.0
Butane, 2-metho...	3.076	51.4	ng/ul	476569	1	7.876	185368	20.0
Ethanol, 2-(2-b...	10.438	4.3	ng/ul	55208	2	10.667	253643	20.0
n-Hexadecanoic ...	18.159	2.0	ng/ul	45737	4	17.259	446512	20.0
Benzaldehyde, 3...	18.329	2.5	ng/ul	55571	4	17.259	446512	20.0
5,6,7,8-Tetrahy...	19.198	2.5	ng/ul	56482	4	17.259	446512	20.0
unknown-01	21.179	4.6	ng/ul	202202	5	21.513	871757	20.0
9-Octadecenamid...	22.924	4.9	ng/ul	214360	5	21.513	871757	20.0
Perylene	24.310	4.6	ng/ul	130668	6	24.598	570880	20.0