

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100523\
 Data File : BG059323.D
 Acq On : 6 Oct 2023 9:53
 Operator : MA/JU
 Sample : 04592-04 10X
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EZYY2

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-BG092723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG059323.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.559	25	29	35	rVB	68191	97768	5.41%	0.558%
2	3.888	83	85	88	rVB2	11341	10568	0.58%	0.060%
3	3.964	95	98	109	rBV2	75834	154999	8.57%	0.884%
4	4.070	115	116	120	rVB4	5390	4662	0.26%	0.027%
5	4.270	146	150	154	rBV2	9874	14229	0.79%	0.081%
6	4.557	193	199	207	rBV	18110	38209	2.11%	0.218%
7	4.634	210	212	217	rVB3	4050	5358	0.30%	0.031%
8	4.751	227	232	239	rVB	17335	22783	1.26%	0.130%
9	4.822	239	244	251	rBV4	12238	19433	1.07%	0.111%
10	5.210	304	310	315	rBV	32802	53615	2.97%	0.306%
11	5.550	366	368	372	rBV3	3543	5413	0.30%	0.031%
12	6.197	476	478	480	rVB2	3810	3256	0.18%	0.019%
13	6.232	480	484	485	rBV2	2895	4410	0.24%	0.025%
14	6.302	493	496	498	rBV2	3040	3237	0.18%	0.018%
15	6.714	563	566	568	rBV2	3014	3245	0.18%	0.019%
16	7.360	671	676	684	rVV2	56989	100402	5.55%	0.573%
17	7.554	703	709	717	rVV	221060	366249	20.25%	2.089%
18	7.630	721	722	727	rBV2	3682	3404	0.19%	0.019%
19	7.748	735	742	756	rBV	194235	351047	19.41%	2.002%
20	7.842	756	758	760	rVV3	3944	3831	0.21%	0.022%
21	8.147	807	810	813	rVB3	2220	3136	0.17%	0.018%
22	8.235	818	825	834	rVB	172577	310076	17.15%	1.769%
23	8.559	878	880	883	rVB3	3293	3178	0.18%	0.018%
24	8.923	935	942	952	rBV2	160007	288912	15.98%	1.648%
25	9.076	961	968	972	rVB4	8051	13139	0.73%	0.075%
26	9.328	1004	1011	1012	rBV4	3768	7023	0.39%	0.040%
27	9.428	1018	1028	1037	rBV	269077	486902	26.93%	2.777%
28	9.581	1052	1054	1058	rVB4	5523	7095	0.39%	0.040%
29	9.640	1061	1064	1065	rBV3	7492	6183	0.34%	0.035%
30	9.886	1103	1106	1108	rBV2	3822	3816	0.21%	0.022%
31	10.151	1142	1151	1158	rVV	214102	389199	21.52%	2.220%
32	10.227	1161	1164	1166	rBV	3717	4147	0.23%	0.024%
33	10.639	1228	1234	1235	rBV3	8752	14711	0.81%	0.084%
34	10.680	1235	1241	1247	rVV2	290400	536633	29.68%	3.061%
35	11.073	1301	1308	1315	rBV	255294	451708	24.98%	2.577%
36	11.138	1317	1319	1324	rVB3	3533	4100	0.23%	0.023%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092723.MA.M
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37	11.220	1324	1333	1343	rBV	286112	530807	29.36%	3.028%
38	11.414	1365	1366	1370	rVB	3638	3658	0.20%	0.021%
39	11.538	1383	1387	1393	rBV2	29515	46873	2.59%	0.267%
40	11.596	1393	1397	1404	rVV	37872	59329	3.28%	0.338%

41	12.137	1486	1489	1492	rVB	2789	3954	0.22%	0.023%
42	12.372	1524	1529	1533	rBV3	3558	5489	0.30%	0.031%
43	12.630	1570	1573	1579	rBV2	8445	14185	0.78%	0.081%
44	12.718	1583	1588	1594	rVV3	6232	11859	0.66%	0.068%
45	12.859	1608	1612	1614	rBV4	4529	6438	0.36%	0.037%

46	12.918	1619	1622	1624	rBV4	3467	5117	0.28%	0.029%
47	13.142	1659	1660	1665	rBV2	2497	3639	0.20%	0.021%
48	13.359	1693	1697	1700	rBV3	3661	3770	0.21%	0.022%
49	13.476	1712	1717	1719	rVB2	2380	3286	0.18%	0.019%
50	13.506	1719	1722	1724	rBV2	3034	3216	0.18%	0.018%

51	13.600	1735	1738	1742	rBV2	3368	5164	0.29%	0.029%
52	13.694	1750	1754	1760	rVB3	21988	34878	1.93%	0.199%
53	13.835	1774	1778	1779	rBV	3667	4350	0.24%	0.025%
54	14.011	1803	1808	1809	rBV	3343	4201	0.23%	0.024%
55	14.105	1823	1824	1828	rBV	3000	3355	0.19%	0.019%

56	14.258	1844	1850	1856	rBV	564714	841338	46.53%	4.799%
57	14.305	1856	1858	1866	rVB3	14932	19100	1.06%	0.109%
58	14.569	1895	1903	1911	rBV	712700	1122078	62.05%	6.401%
59	14.863	1944	1953	1960	rBV	398991	624594	34.54%	3.563%
60	15.057	1982	1986	1993	rBV4	44249	73954	4.09%	0.422%

61	15.221	2007	2014	2016	rBV6	5679	9861	0.55%	0.056%
62	15.345	2033	2035	2037	rBV3	3750	4166	0.23%	0.024%
63	15.427	2047	2049	2052	rVB	4264	3506	0.19%	0.020%
64	15.621	2079	2082	2086	rVV4	5320	6520	0.36%	0.037%
65	15.856	2114	2122	2130	rBV	903667	1380991	76.37%	7.877%

66	15.968	2135	2141	2148	rBV2	259404	368285	20.37%	2.101%
67	16.097	2158	2163	2165	rVB3	6940	9279	0.51%	0.053%
68	16.467	2224	2226	2227	rBV	4031	4110	0.23%	0.023%
69	16.538	2237	2238	2241	rBV3	4237	3374	0.19%	0.019%
70	16.579	2241	2245	2246	rVV	3630	4257	0.24%	0.024%

71	16.737	2269	2272	2274	rBV	2771	3186	0.18%	0.018%
72	16.837	2286	2289	2291	rBV3	3144	4049	0.22%	0.023%
73	16.931	2303	2305	2306	rBV	6065	5726	0.32%	0.033%
74	17.119	2334	2337	2338	rBV	5232	4588	0.25%	0.026%
75	17.196	2348	2350	2353	rVB3	3980	3545	0.20%	0.020%

76	17.231	2353	2356	2358	rBV2	4894	5575	0.31%	0.032%
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 Title : SVOA CALIBRATION

77	17.278	2361	2364	2365	rBV2	4643	4221	0.23%	0.024%
78	17.413	2385	2387	2391	rVB3	4015	5102	0.28%	0.029%
79	17.613	2415	2421	2428	rVB	499389	754451	41.72%	4.304%
80	17.713	2432	2438	2448	rVB	995917	1540899	85.22%	8.790%
81	17.783	2448	2450	2451	rBV2	4980	4318	0.24%	0.025%
82	18.130	2507	2509	2513	rVB4	4055	4120	0.23%	0.024%
83	18.429	2556	2560	2568	rBV3	124035	185392	10.25%	1.058%
84	18.659	2597	2599	2600	rBV2	7627	5278	0.29%	0.030%
85	18.711	2605	2608	2611	rBV5	6933	8975	0.50%	0.051%
86	18.776	2617	2619	2622	rBV4	3794	3599	0.20%	0.021%
87	18.982	2650	2654	2655	rBV4	6286	8747	0.48%	0.050%
88	19.493	2739	2741	2743	rBV3	6793	5582	0.31%	0.032%
89	19.552	2749	2751	2754	rVV4	13404	13667	0.76%	0.078%
90	19.599	2754	2759	2766	rVB2	51241	77928	4.31%	0.445%
91	19.693	2771	2775	2779	rBV3	55248	78610	4.35%	0.448%
92	19.834	2795	2799	2804	rBV	128285	145968	8.07%	0.833%
93	19.945	2815	2818	2820	rBV3	18895	21432	1.19%	0.122%
94	19.986	2820	2825	2832	rVB	1234495	1775480	98.19%	10.128%
95	20.662	2938	2940	2944	rVB4	28377	25681	1.42%	0.146%
96	20.874	2972	2976	2981	rBV	97704	126399	6.99%	0.721%
97	21.914	3147	3153	3162	rVB2	484046	820033	45.35%	4.678%
98	23.600	3435	3440	3445	rVB3	95231	186078	10.29%	1.061%
99	25.157	3696	3705	3716	rVB	616637	1808206	100.00%	10.314%
100	25.404	3740	3747	3759	rVB	292036	868938	48.06%	4.957%

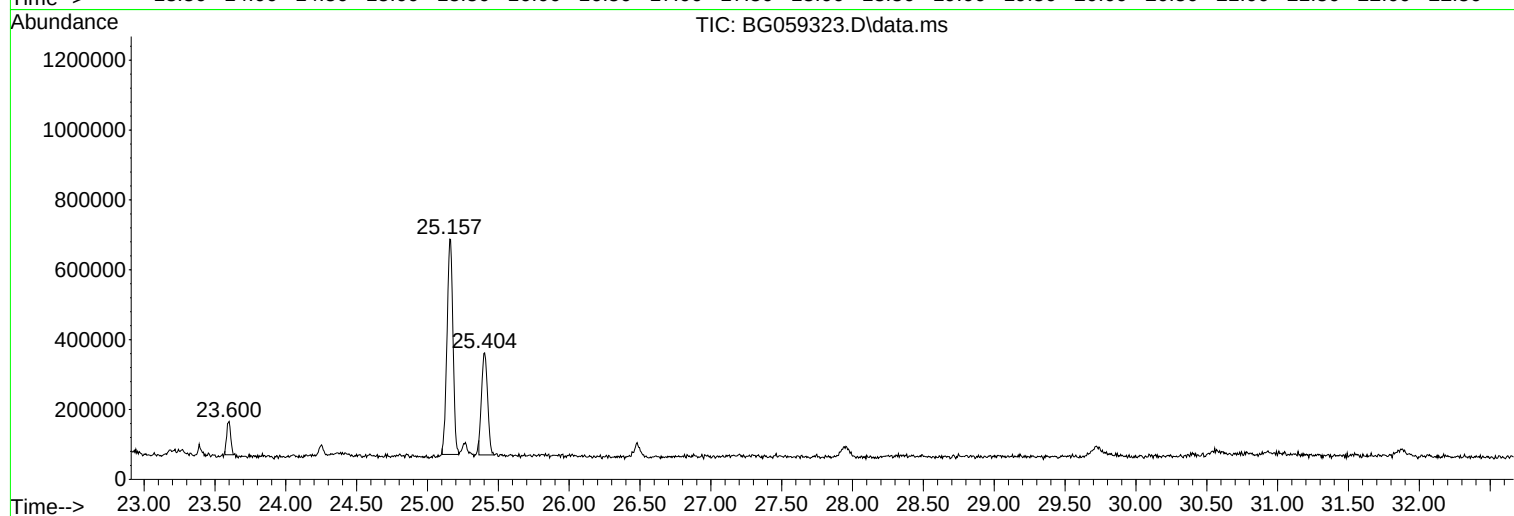
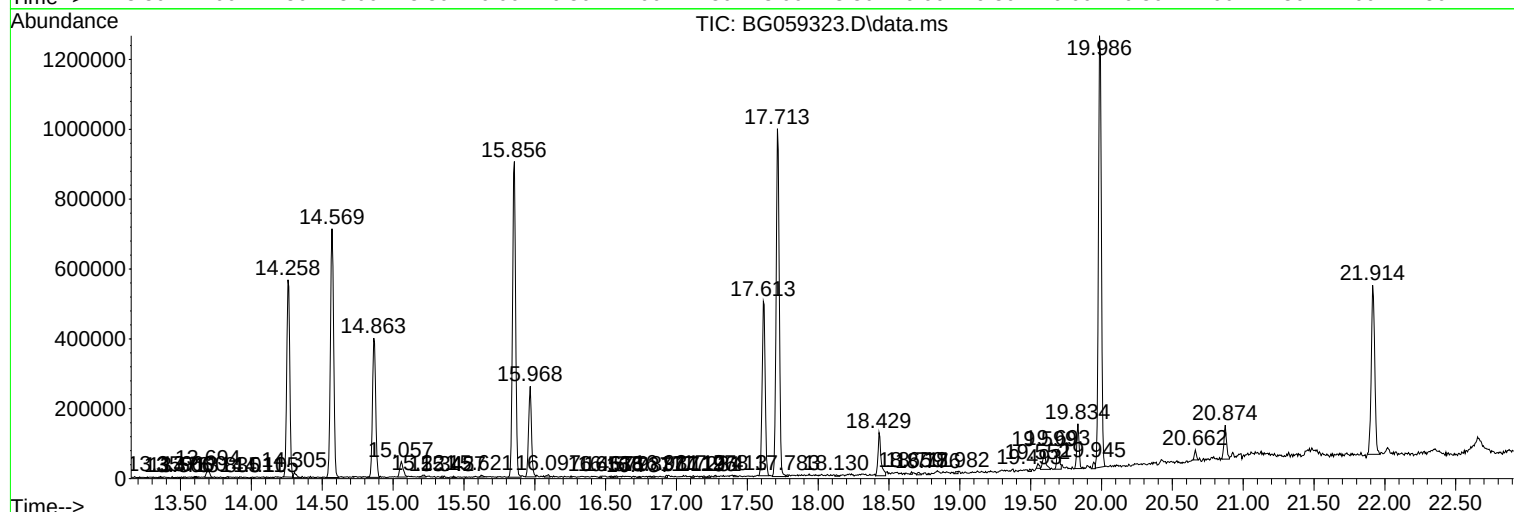
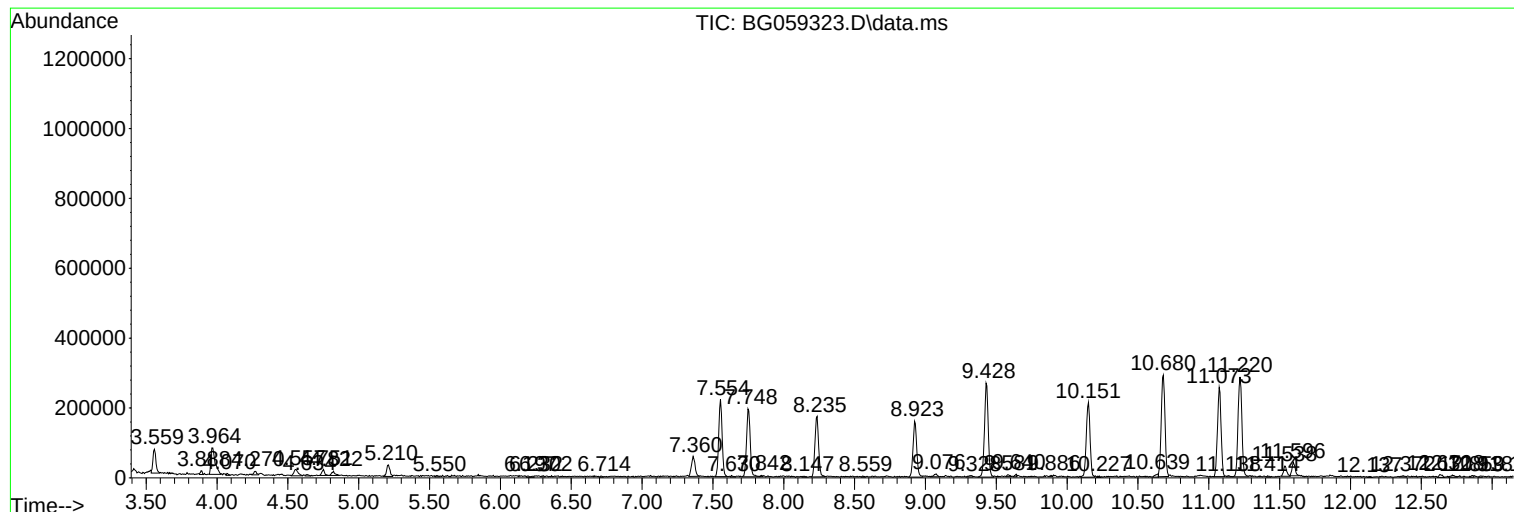
Sum of corrected areas: 17530830

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092723.MA.M
Quant Title : SVOA CALIBRATION

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



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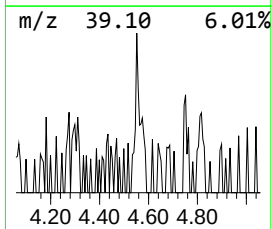
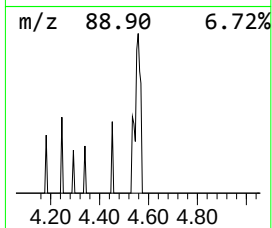
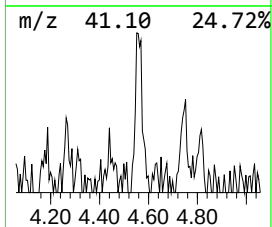
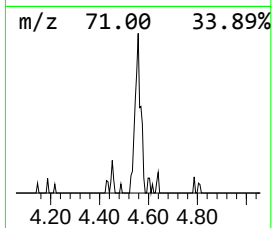
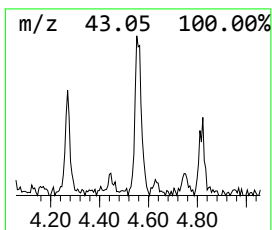
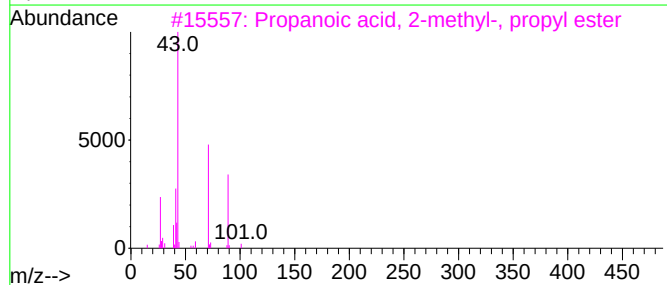
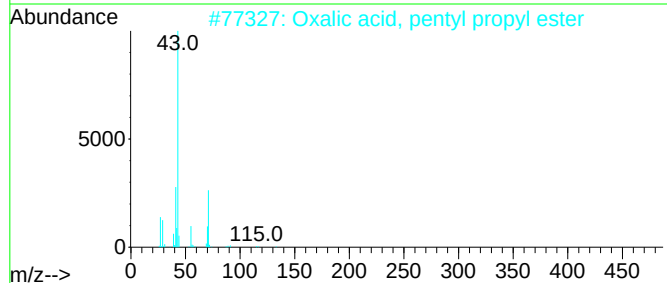
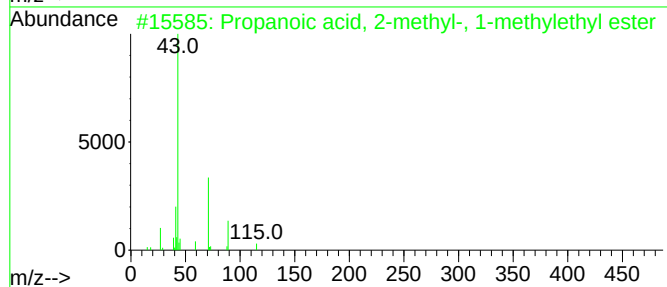
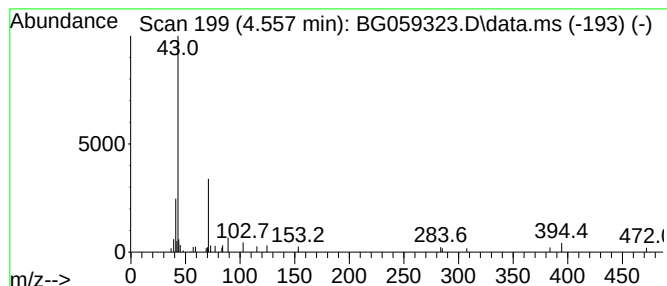
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.557	2.46 ng/ul	38209	1,4-Dichlorobenzene-d4	8.230

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	40
2			Oxalic acid, pentyl propyl ester	202	C10H18O4	1000309-25-7	9
3			Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	9
4			Oxalic acid, dipropyl ester	174	C8H14O4	1000309-25-4	9
5			dl-Mevalonic acid lactone	130	C6H10O3	000674-26-0	9



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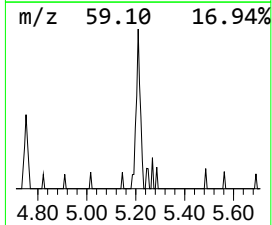
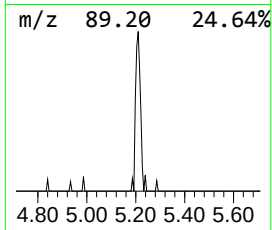
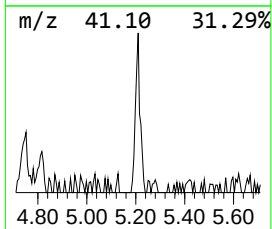
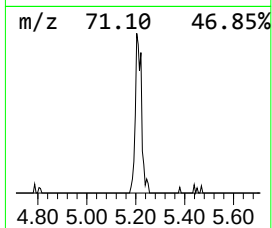
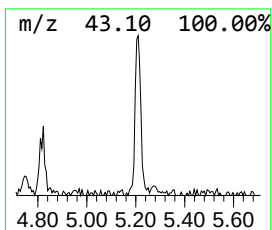
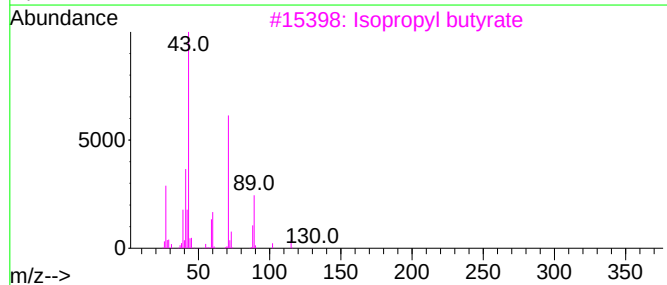
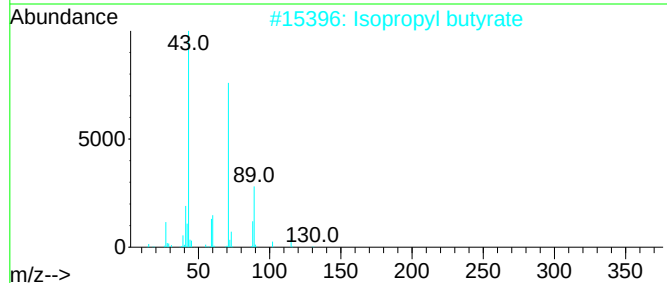
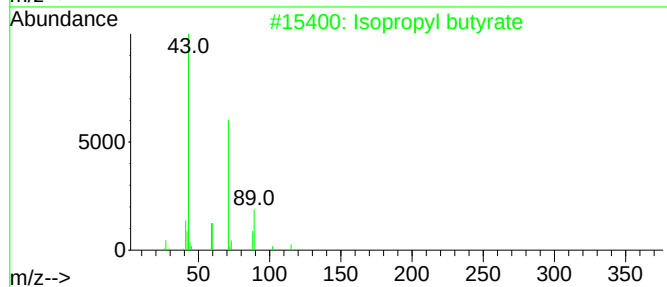
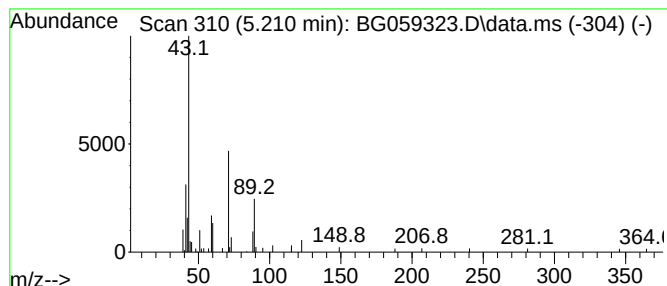
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Isopropyl butyrate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.210	3.46 ng/ul	53615	1,4-Dichlorobenzene-d4	8.230

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl butyrate	130	C7H14O2	000638-11-9	72
2			Isopropyl butyrate	130	C7H14O2	000638-11-9	72
3			Isopropyl butyrate	130	C7H14O2	000638-11-9	59
4			Isopropyl butyrate	130	C7H14O2	000638-11-9	59
5			Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	53



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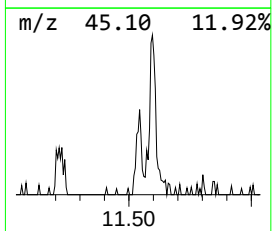
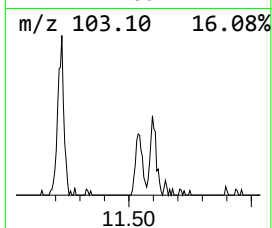
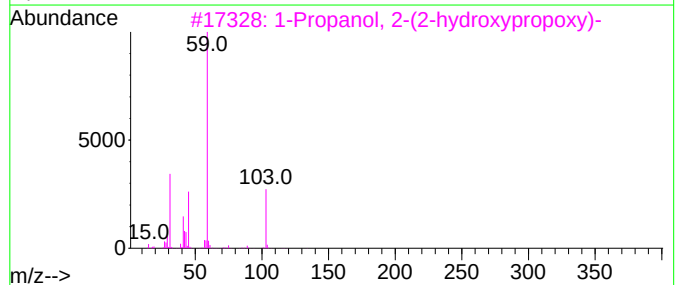
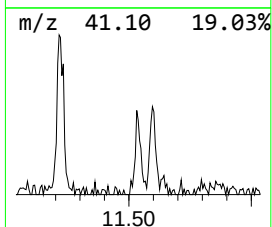
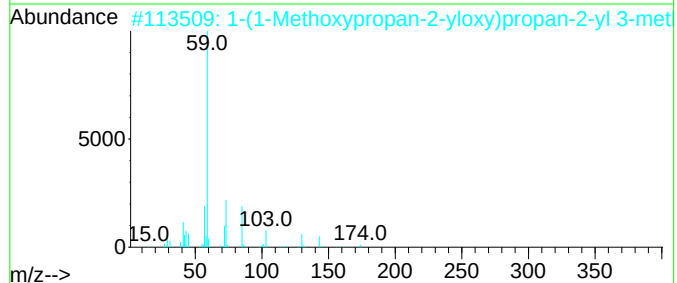
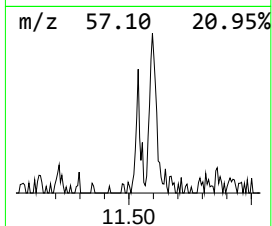
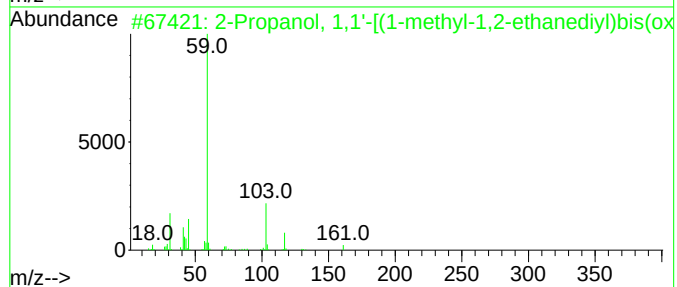
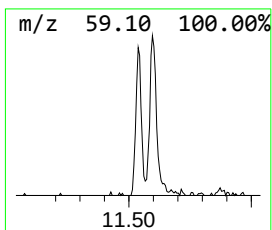
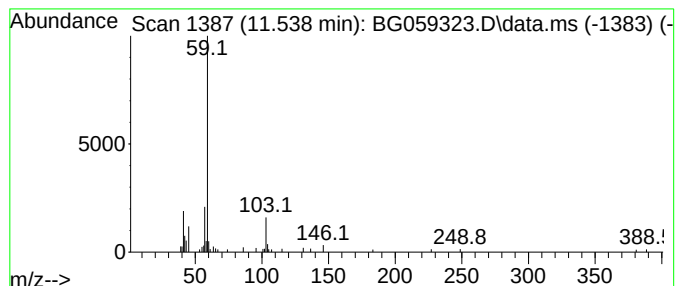
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 2-Propanol, 1,1'-[(1-methyl... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.538	2.08 ng/ul	46873	Naphthalene-d8	11.073

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1,1'-[(1-methyl-1,2-...	192	C9H20O4	001638-16-0	78		
2	1-(1-Methoxypropan-2-yloxy)propa...	232	C12H24O4	1000367-13-2	72		
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72		
4	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72		
5	4,8,12,16-tetraoxaicosan-1-ol	306	C16H34O5	1010397-63-6	64		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100523\
Data File : BG059323.D
Acq On : 6 Oct 2023 9:53
Operator : MA/JU
Sample : 04592-04 10X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZY2

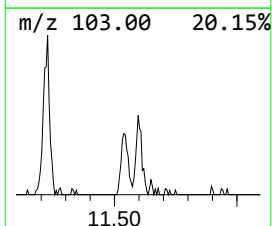
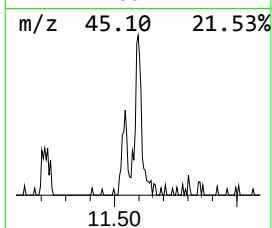
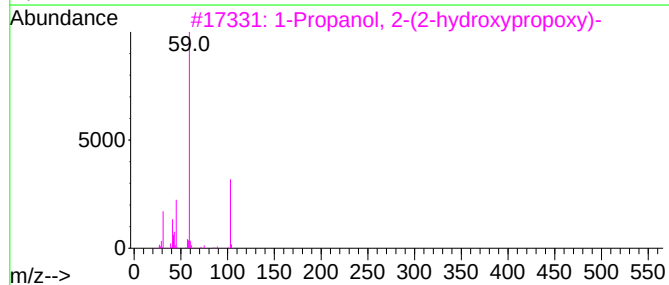
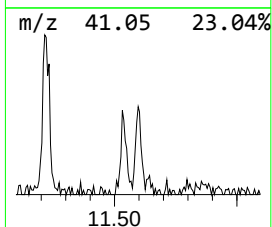
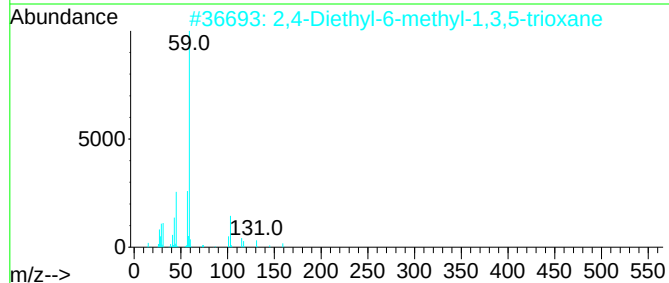
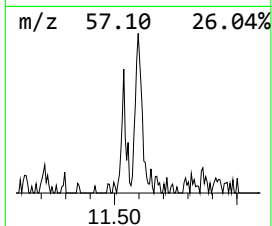
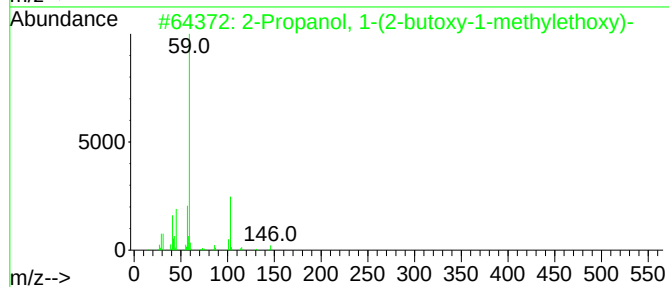
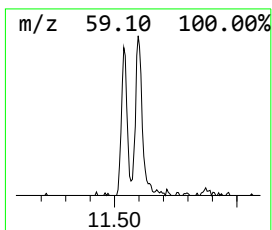
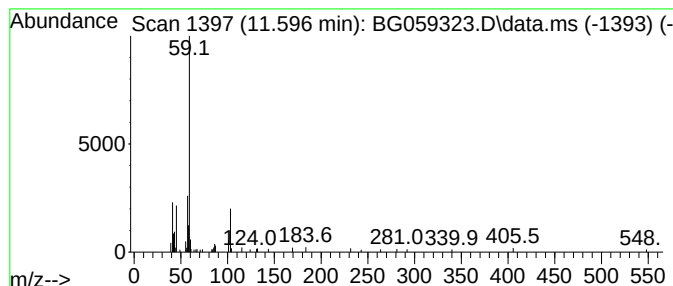
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 2-Propanol, 1-(2-butoxy-1-m... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.596	2.63 ng/ul	59329	Naphthalene-d8	11.073

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	86		
2	2,4-Diethyl-6-methyl-1,3,5-trioxane	160	C8H16O3	117888-04-7	78		
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72		
4	Dipropylene glycol (isomer 3)	134	C6H14O3	1000506-25-5	72		
5	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100523\
Data File : BG059323.D
Acq On : 6 Oct 2023 9:53
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Sample : 04592-04 10X
Misc :
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Instrument :
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ClientSampleId :
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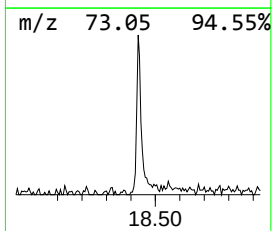
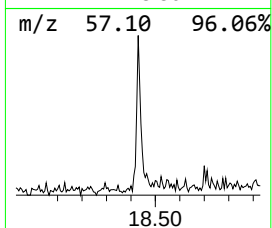
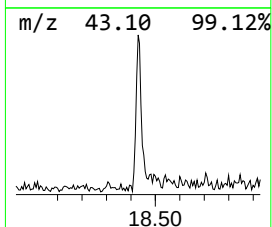
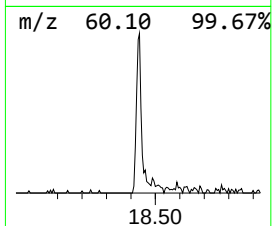
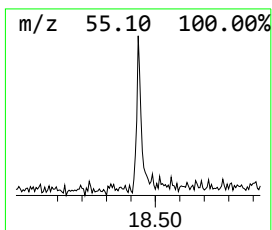
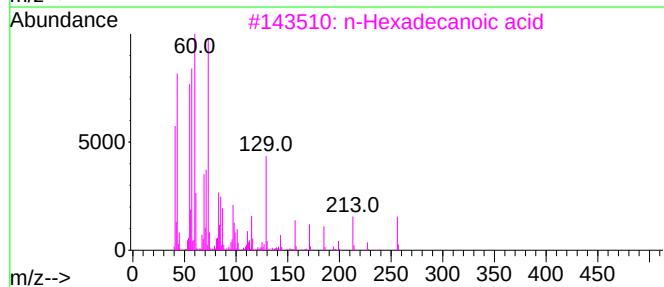
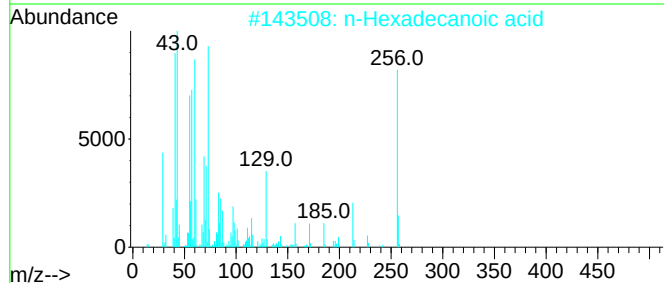
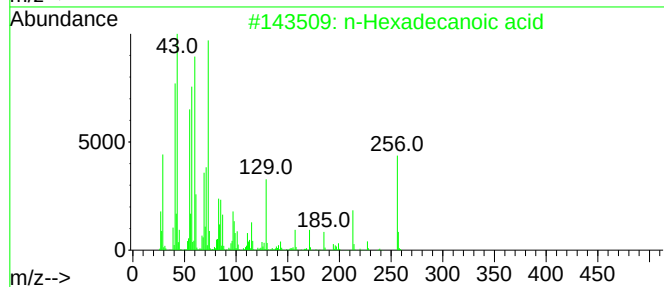
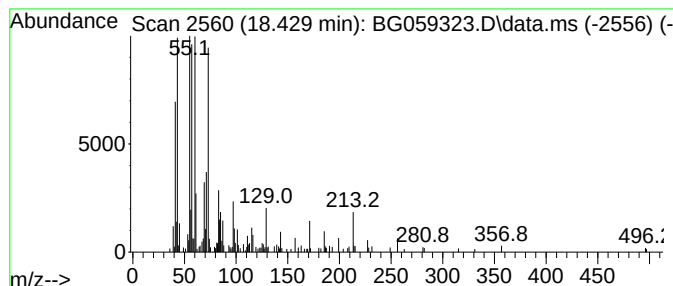
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.429	4.91 ng/ul	185392	Phenanthrene-d10	17.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	89
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	80
4			Isopropyl palmitate	298	C19H38O2	000142-91-6	72
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100523\
Data File : BG059323.D
Acq On : 6 Oct 2023 9:53
Operator : MA/JU
Sample : 04592-04 10X
Misc :
ALS Vial : 37 Sample Multiplier: 1

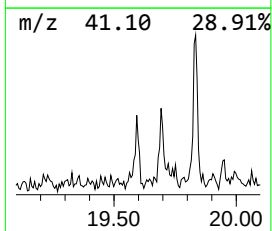
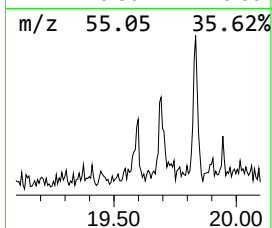
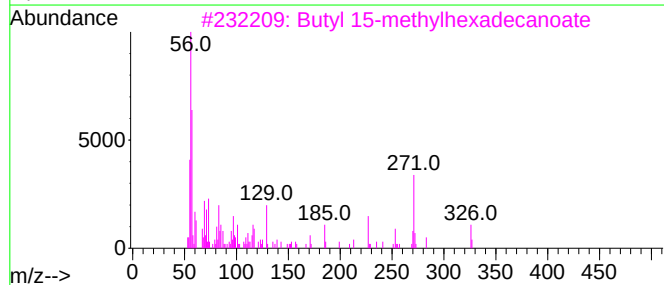
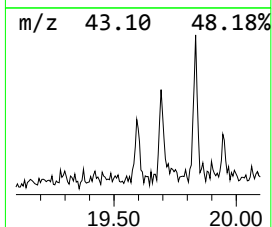
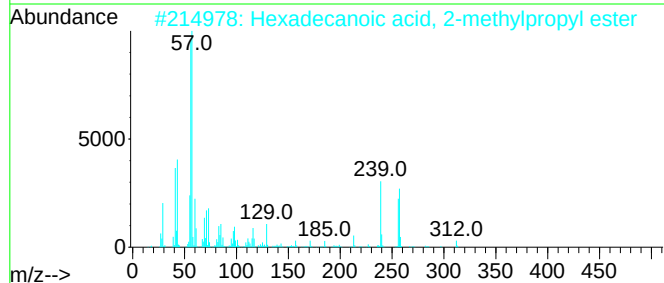
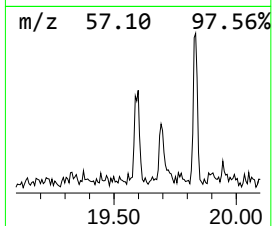
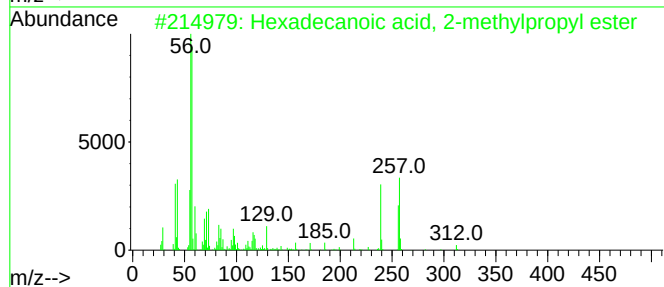
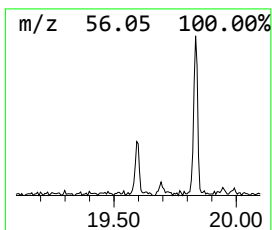
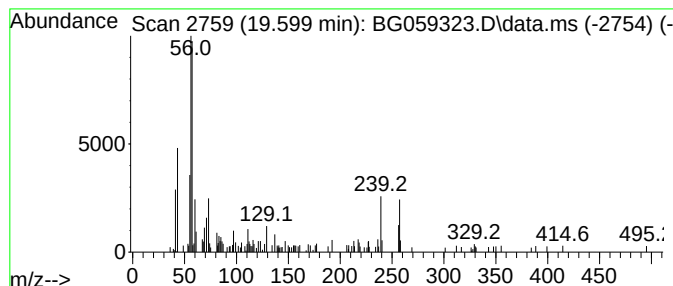
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Hexadecanoic acid, 2-methyl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.599	2.07 ng/ul	77928	Phenanthrene-d10		17.619	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, 2-methylpropy...		312	C20H40O2	000110-34-9	90
2	Hexadecanoic acid, 2-methylpropy...		312	C20H40O2	000110-34-9	90
3	Butyl 15-methylhexadecanoate		326	C21H42O2	1000336-52-4	43
4	Hexadecanoic acid, 1,1-dimethyle...		312	C20H40O2	031158-91-5	43
5	Hexadecanoic acid, butyl ester		312	C20H40O2	000111-06-8	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100523\
Data File : BG059323.D
Acq On : 6 Oct 2023 9:53
Operator : MA/JU
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ALS Vial : 37 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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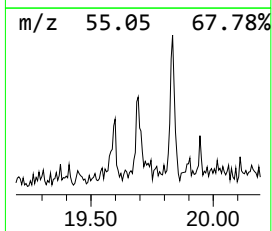
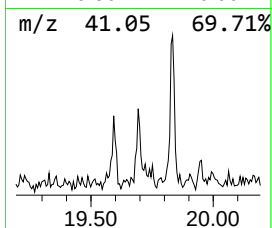
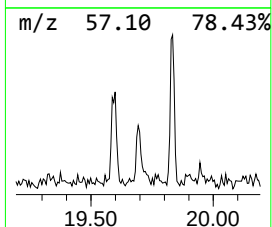
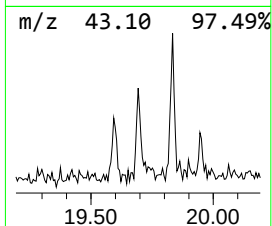
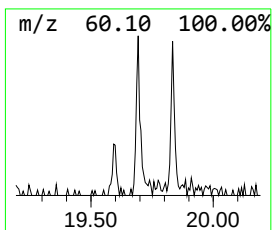
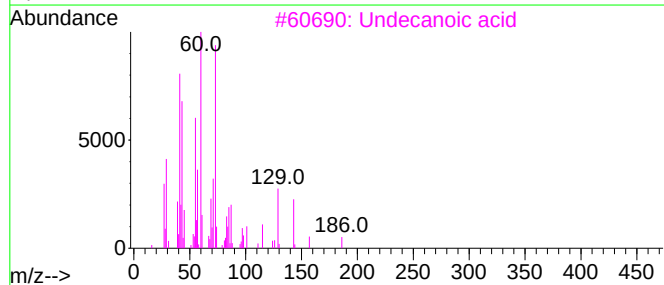
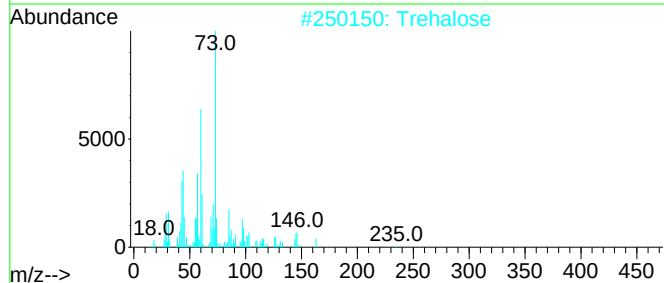
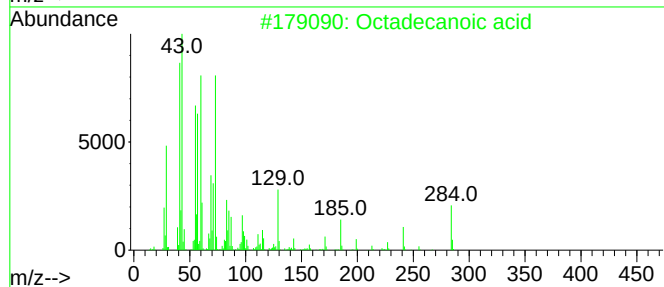
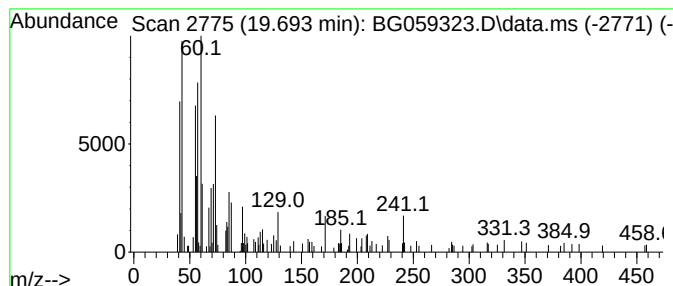
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.693	2.08 ng/ul	78610	Phenanthrene-d10	17.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	72
2			Trehalose	342	C12H22O11	000099-20-7	64
3			Undecanoic acid	186	C11H22O2	000112-37-8	60
4			Dodecanoic acid	200	C12H24O2	000143-07-7	47
5			Tridecanoic acid	214	C13H26O2	000638-53-9	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100523\
Data File : BG059323.D
Acq On : 6 Oct 2023 9:53
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Sample : 04592-04 10X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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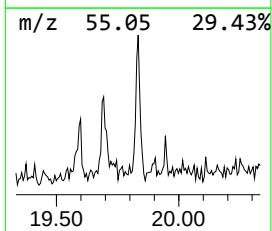
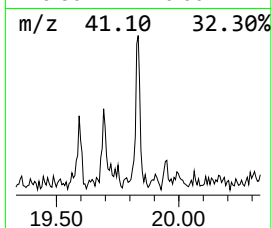
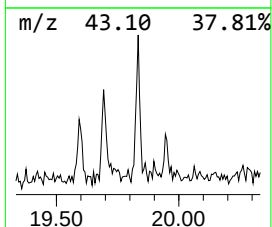
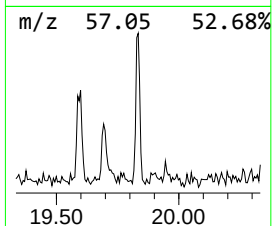
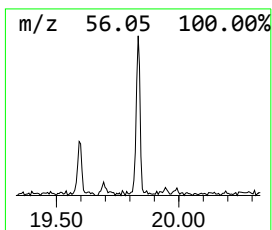
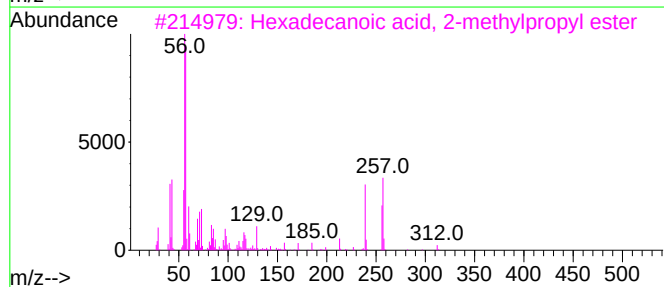
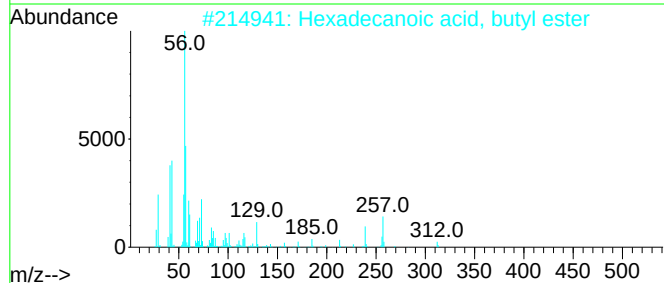
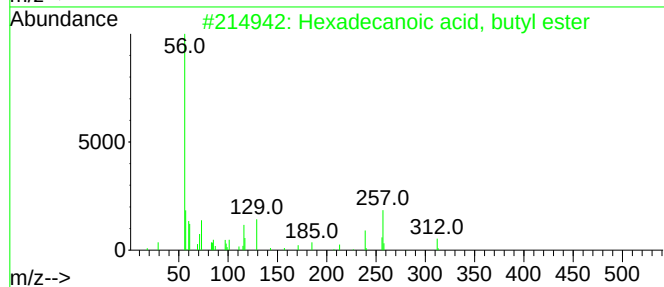
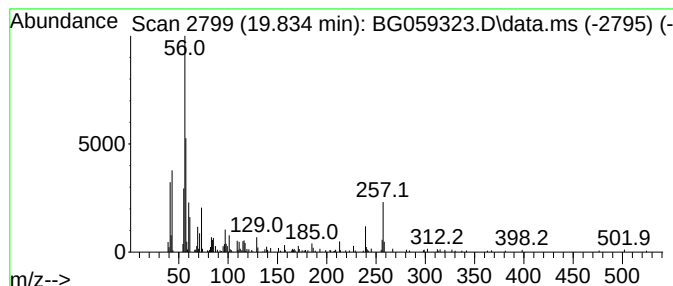
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Hexadecanoic acid, butyl ester Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.834	3.56 ng/ul	145968	Chrysene-d12	21.914

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	72
2			Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	46
3			Hexadecanoic acid, 2-methylpropy...	312	C20H40O2	000110-34-9	45
4			Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	38
5			Cyclobutane, 1-hexyl-2,3-dimethyl-	168	C12H24	055170-84-8	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100523\
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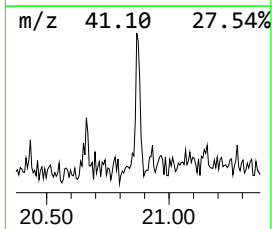
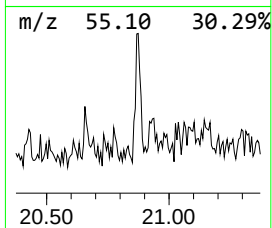
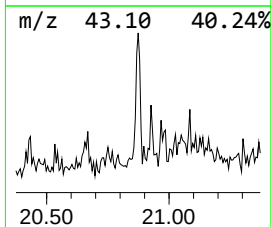
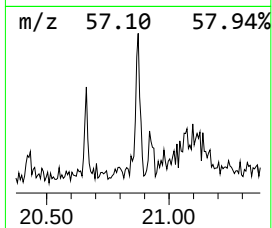
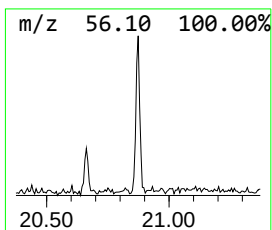
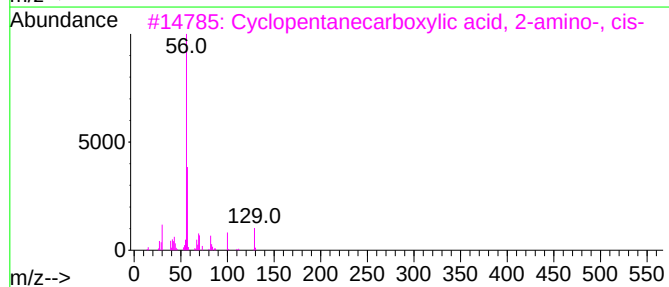
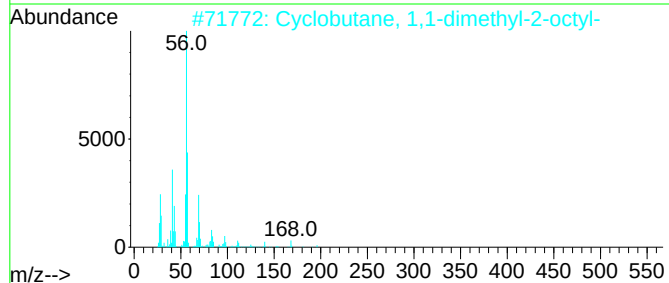
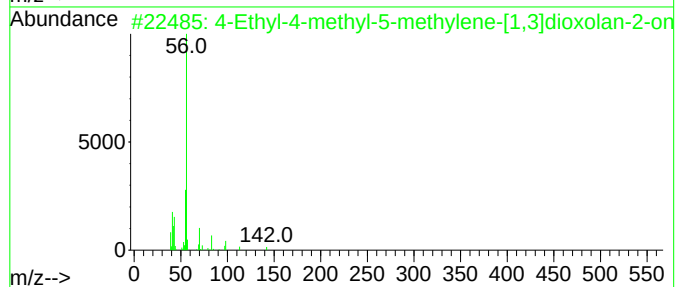
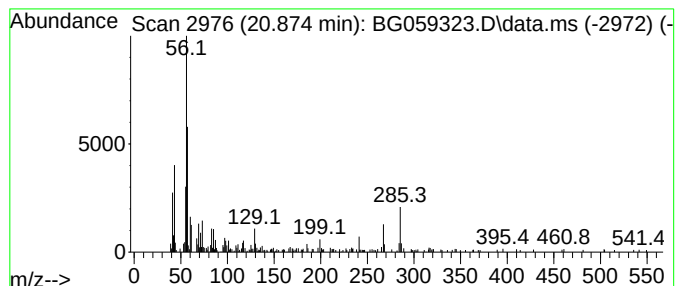
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 4-Ethyl-4-methyl-5-methylen... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.874	3.08 ng/ul	126399	Chrysene-d12		21.914	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	4-Ethyl-4-methyl-5-methylene-[1,...		142	C7H10O3	080956-52-1	52
2	Cyclobutane, 1,1-dimethyl-2-octyl-		196	C14H28	062338-30-1	50
3	Cyclopentanecarboxylic acid, 2-a...		129	C6H11NO2	037910-65-9	50
4	1-Hexene, 2,5-dimethyl-		112	C8H16	006975-92-4	50
5	1-Oxa-4-azaspiro[4.5]decan-4-oxy...		198	C10H16NO3	1000115-84-0	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100523\
Data File : BG059323.D
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Sample : 04592-04 10X
Misc :
ALS Vial : 37 Sample Multiplier: 1

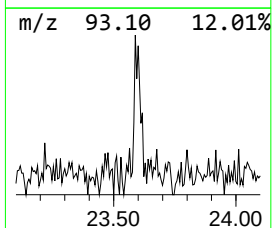
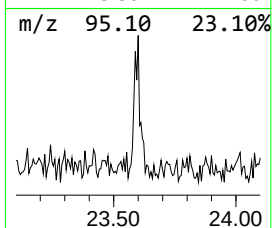
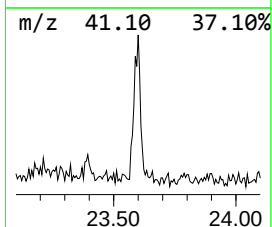
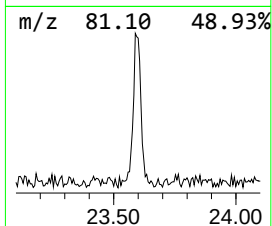
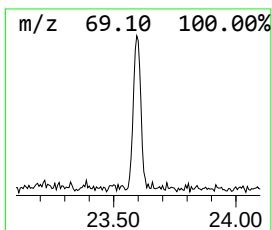
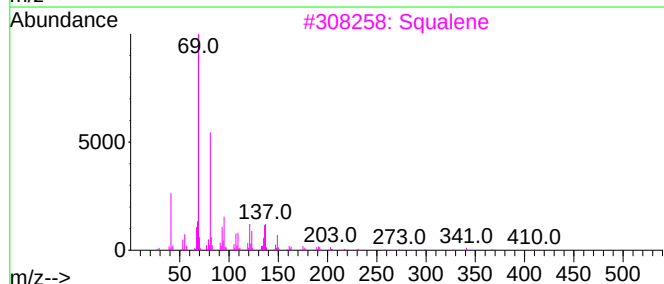
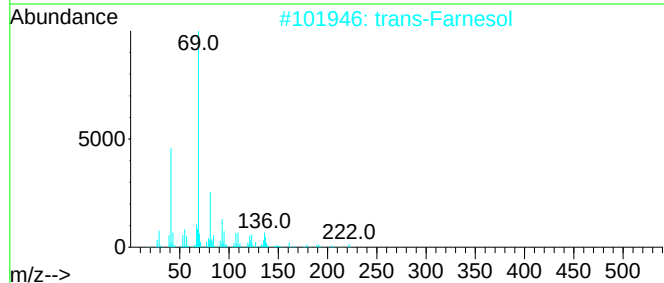
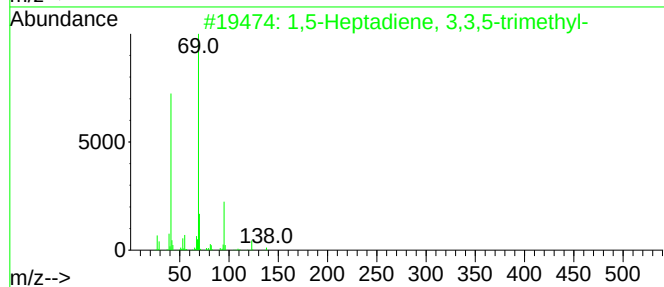
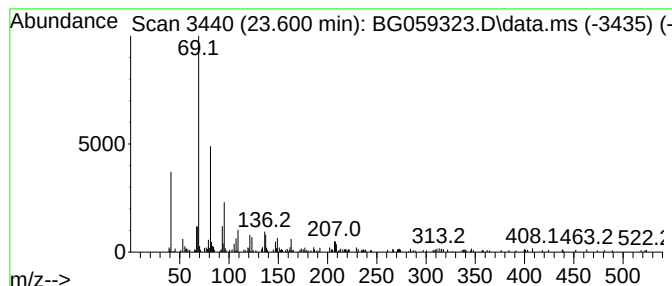
Instrument :
BNA_G
ClientSampleId :
EZY2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 1,5-Heptadiene, 3,3,5-trime... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.600	4.54 ng/ul	186078	Chrysene-d12	21.914	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,5-Heptadiene, 3,3,5-trimethyl-	138	C10H18	074630-29-8	72
2	trans-Farnesol	222	C15H26O	000106-28-5	70
3	Squalene	410	C30H50	000111-02-4	64
4	2,6-Dimethyl-2-trans-6-octadiene	138	C10H18	002609-23-6	64
5	1,2-Epoxy-3,7,11-trimethyl-(6E,1...	238	C15H26O2	1000432-46-2	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100523\
Data File : BG059323.D
Acq On : 6 Oct 2023 9:53
Operator : MA/JU
Sample : 04592-04 10X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZY2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092723.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	4.557	2.5	ng/ul	38209	1	8.230	310076	20.0
Isopropyl butyrate	5.210	3.5	ng/ul	53615	1	8.230	310076	20.0
2-Propanol, 1,1...	11.538	2.1	ng/ul	46873	2	11.073	451708	20.0
2-Propanol, 1-(...	11.596	2.6	ng/ul	59329	2	11.073	451708	20.0
n-Hexadecanoic ...	18.429	4.9	ng/ul	185392	4	17.619	754451	20.0
Hexadecanoic ac...	19.599	2.1	ng/ul	77928	4	17.619	754451	20.0
Octadecanoic acid	19.693	2.1	ng/ul	78610	4	17.619	754451	20.0
Hexadecanoic ac...	19.834	3.6	ng/ul	145968	5	21.914	820033	20.0
4-Ethyl-4-methy...	20.874	3.1	ng/ul	126399	5	21.914	820033	20.0
1,5-Heptadiene,...	23.600	4.5	ng/ul	186078	5	21.914	820033	20.0