

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG093021\
 Data File : BG050398.D
 Acq On : 2 Oct 2021 21:03
 Operator : CG/JU
 Sample : M3877-14
 Misc :
 ALS Vial : 84 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW2N0

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-BG093021.M
 Title : SVOA CALIBRATION

Signal : TIC: BG050398.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.561	10	12	21	rVB7	2691	5106	0.30%	0.028%
2	3.632	21	24	30	rVV3	7810	13798	0.82%	0.077%
3	4.067	93	98	107	rBV2	45107	85917	5.11%	0.478%
4	4.173	111	116	124	rVV	47580	74253	4.42%	0.413%
5	4.243	125	128	134	rVV5	5717	9886	0.59%	0.055%
6	4.302	134	138	144	rVB3	12799	21106	1.26%	0.117%
7	4.408	152	156	159	rVB4	3450	5852	0.35%	0.033%
8	4.778	212	219	225	rVB8	2735	6272	0.37%	0.035%
9	4.972	247	252	257	rVV4	5666	9105	0.54%	0.051%
10	5.118	273	277	280	rBV4	2726	4680	0.28%	0.026%
11	5.330	307	313	321	rBV6	5235	11577	0.69%	0.064%
12	5.418	324	328	336	rVB4	8068	15897	0.95%	0.088%
13	7.404	658	666	676	rVV	151273	274464	16.33%	1.528%
14	7.586	690	697	709	rBV	108847	190442	11.33%	1.060%
15	7.798	724	733	739	rBV	137719	250007	14.87%	1.392%
16	8.274	803	814	824	rVB	232005	414326	24.65%	2.306%
17	8.491	845	851	857	rVB6	4558	8045	0.48%	0.045%
18	8.797	899	903	907	rBV5	2635	4855	0.29%	0.027%
19	8.961	921	931	939	rBV	142479	258701	15.39%	1.440%
20	9.443	1004	1013	1022	rVV	140991	264940	15.76%	1.475%
21	9.819	1072	1077	1082	rBV7	2492	5887	0.35%	0.033%
22	10.166	1129	1136	1144	rBV2	116405	212197	12.62%	1.181%
23	10.700	1220	1227	1235	rVV	169304	318614	18.95%	1.774%
24	10.841	1245	1251	1260	rBV2	36426	63914	3.80%	0.356%
25	11.094	1287	1294	1303	rVV	310502	582895	34.68%	3.245%
26	11.223	1309	1316	1327	rBV	72682	142992	8.51%	0.796%
27	12.663	1556	1561	1564	rBV3	4503	5607	0.33%	0.031%
28	13.256	1656	1662	1667	rBV4	2591	5374	0.32%	0.030%
29	13.691	1733	1736	1740	rBV3	3755	6531	0.39%	0.036%
30	14.202	1818	1823	1828	rBV5	3496	6383	0.38%	0.036%
31	14.278	1828	1836	1844	rVV	303663	455330	27.09%	2.535%
32	14.507	1872	1875	1881	rVV5	3710	6357	0.38%	0.035%
33	14.584	1881	1888	1900	rVV	348759	578883	34.44%	3.223%
34	14.754	1914	1917	1922	rVB5	3303	5020	0.30%	0.028%
35	14.889	1932	1940	1947	rVB	470752	766370	45.59%	4.266%
36	14.948	1947	1950	1956	rVB4	5084	8559	0.51%	0.048%

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Integrator: RTE

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG093021.M
 Title : SVOA CALIBRATION

37	15.054	1960	1968	1977	rBV	128812	219175	13.04%	1.220%
38	15.213	1989	1995	2000	rBV5	12977	24817	1.48%	0.138%
39	15.265	2000	2004	2005	rBV2	4273	5289	0.31%	0.029%
40	15.700	2076	2078	2083	rVV4	3316	4752	0.28%	0.026%
41	15.876	2100	2108	2114	rBV	440501	696488	41.43%	3.877%
42	15.923	2114	2116	2120	rVV4	9068	13502	0.80%	0.075%
43	15.976	2120	2125	2134	rVB	78363	125473	7.46%	0.698%
44	16.111	2145	2148	2154	rVB4	5703	8965	0.53%	0.050%
45	16.223	2162	2167	2172	rBV5	4061	8435	0.50%	0.047%
46	16.370	2188	2192	2196	rBV4	4038	6065	0.36%	0.034%
47	16.505	2210	2215	2217	rBV4	3600	5373	0.32%	0.030%
48	16.564	2219	2225	2226	rBV5	6411	8212	0.49%	0.046%
49	16.582	2226	2228	2235	rVB7	6515	12903	0.77%	0.072%
50	16.699	2242	2248	2250	rBV6	5046	5915	0.35%	0.033%
51	16.787	2260	2263	2265	rBV3	4983	5811	0.35%	0.032%
52	16.922	2280	2286	2287	rVV5	5800	9548	0.57%	0.053%
53	16.940	2287	2289	2292	rVV4	6646	5995	0.36%	0.033%
54	16.993	2292	2298	2301	rVV6	8939	16016	0.95%	0.089%
55	17.081	2306	2313	2314	rBV5	4585	6684	0.40%	0.037%
56	17.099	2314	2316	2320	rVB5	6076	7640	0.45%	0.043%
57	17.163	2322	2327	2330	rBV7	3720	6598	0.39%	0.037%
58	17.357	2357	2360	2363	rBV5	7766	10583	0.63%	0.059%
59	17.445	2373	2375	2379	rVB2	4537	5577	0.33%	0.031%
60	17.492	2379	2383	2389	rBV2	54461	77320	4.60%	0.430%
61	17.557	2390	2394	2398	rVV3	24201	30921	1.84%	0.172%
62	17.633	2401	2407	2411	rVV	545818	875549	52.09%	4.874%
63	17.674	2411	2414	2418	rVV	98565	152104	9.05%	0.847%
64	17.727	2418	2423	2435	rVB2	434007	722687	42.99%	4.023%
65	18.092	2483	2485	2490	rVV6	5106	7578	0.45%	0.042%
66	18.144	2492	2494	2498	rVB5	6022	6696	0.40%	0.037%
67	18.215	2503	2506	2507	rBV3	7265	8414	0.50%	0.047%
68	18.274	2513	2516	2520	rVB3	20414	22163	1.32%	0.123%
69	18.374	2525	2533	2538	rBV6	35265	71237	4.24%	0.397%
70	18.432	2540	2543	2547	rVV4	12284	19629	1.17%	0.109%
71	18.491	2548	2553	2559	rVV2	165783	257404	15.31%	1.433%
72	18.550	2559	2563	2569	rVB2	34868	72045	4.29%	0.401%
73	18.597	2569	2571	2573	rBV3	4362	5351	0.32%	0.030%
74	18.632	2574	2577	2581	rVB2	12593	15642	0.93%	0.087%
75	18.697	2582	2588	2591	rBV3	34260	64216	3.82%	0.357%
76	18.738	2591	2595	2598	rVB3	18040	26996	1.61%	0.150%

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 Stop Thrs : 0 Peak Location: TOP

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77	18.785	2599	2603	2610	rBV2	39416	69222	4.12%	0.385%
78	18.914	2623	2625	2630	rBV6	7642	6873	0.41%	0.038%
79	18.990	2632	2638	2642	rBV2	23780	45249	2.69%	0.252%
80	19.079	2651	2653	2656	rBV3	9845	11590	0.69%	0.065%
81	19.408	2705	2709	2714	rBV4	27063	44840	2.67%	0.250%
82	19.637	2745	2748	2749	rBV3	24023	24444	1.45%	0.136%
83	19.672	2749	2754	2761	rVB	259003	413333	24.59%	2.301%
84	19.748	2764	2767	2774	rBV4	43792	68362	4.07%	0.381%
85	19.801	2774	2776	2777	rBV2	13585	9909	0.59%	0.055%
86	20.001	2805	2810	2814	rBV	534266	872661	51.91%	4.858%
87	20.412	2875	2880	2881	rBV5	41574	71707	4.27%	0.399%
88	20.618	2912	2915	2919	rVB2	31722	37757	2.25%	0.210%
89	21.535	3067	3071	3077	rBV2	203846	289156	17.20%	1.610%
90	21.934	3131	3139	3144	rBV2	483362	1035205	61.58%	5.763%
91	21.981	3145	3147	3154	rVB	108919	155215	9.23%	0.864%
92	22.786	3276	3284	3291	rBV4	167800	444343	26.43%	2.474%
93	23.879	3464	3470	3478	rVB2	229611	482766	28.72%	2.687%
94	24.372	3548	3554	3558	rBV	166075	346897	20.64%	1.931%
95	24.437	3558	3565	3578	rVB2	655270	1680948	100.00%	9.357%
96	25.124	3676	3682	3690	rVB3	168914	423167	25.17%	2.356%
97	25.365	3715	3723	3736	rVB2	267958	813192	48.38%	4.527%
98	26.623	3929	3937	3947	rVB2	198311	625652	37.22%	3.483%
99	26.752	3952	3959	3969	rVV5	126298	413641	24.61%	2.303%
100	31.106	4688	4700	4721	rVB9	156199	821645	48.88%	4.574%

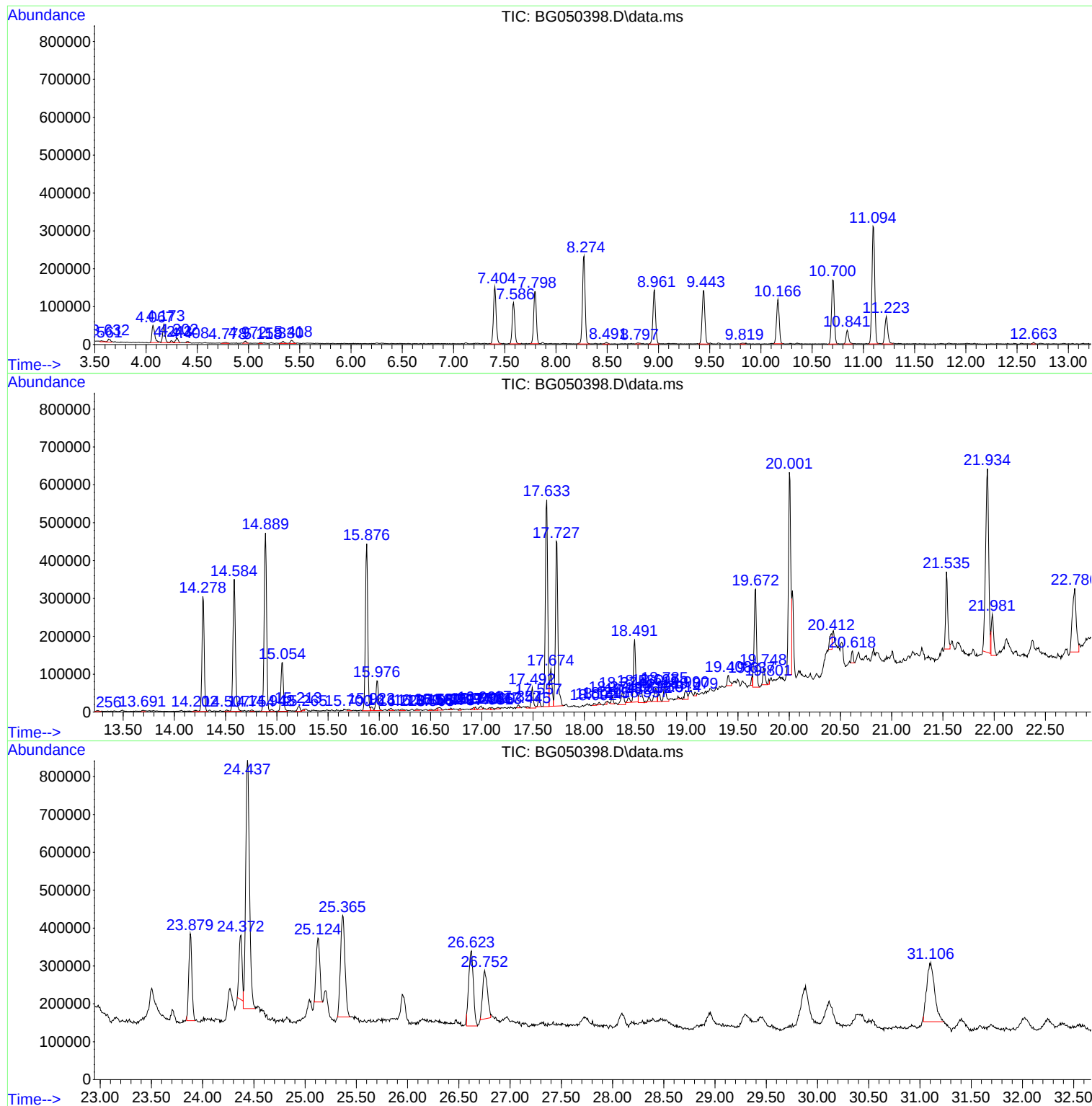
Sum of corrected areas: 17963684

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG093021.M
Quant Title : SVOA CALIBRATION

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



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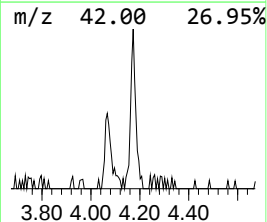
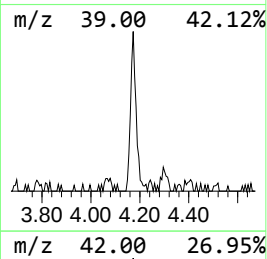
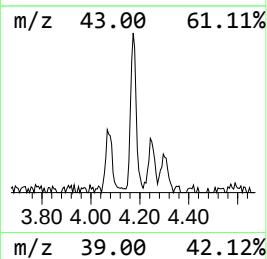
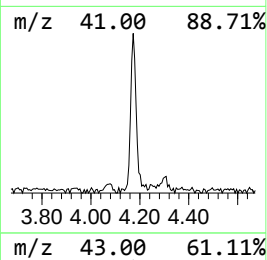
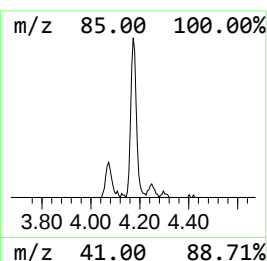
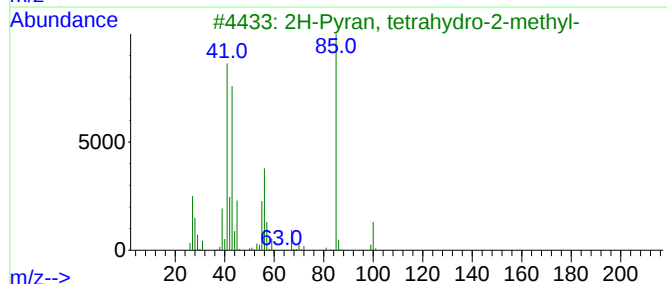
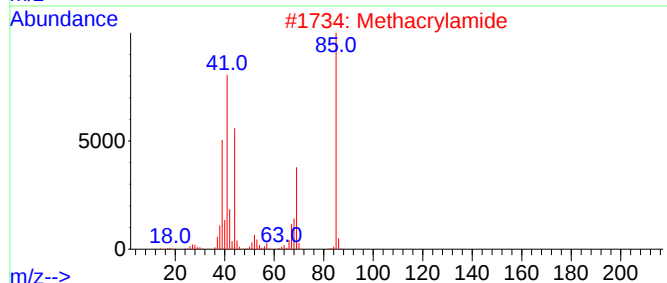
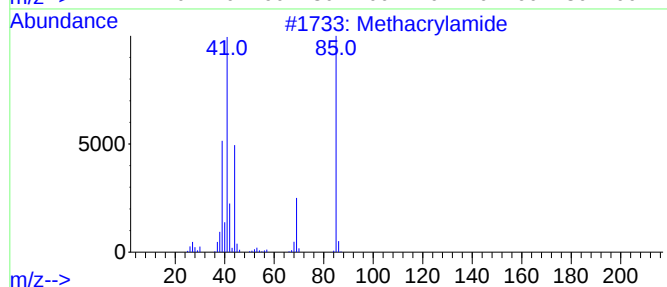
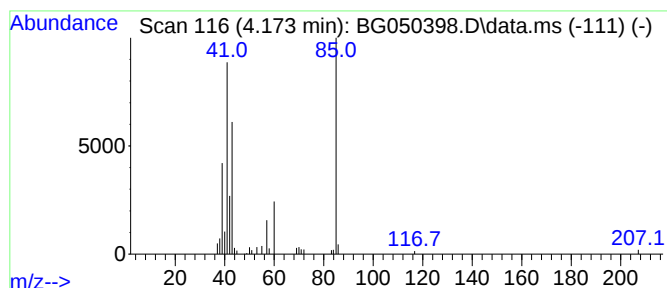
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG093021.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Methacrylamide Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.173	3.58 ng/ul	74253	1,4-Dichlorobenzene-d4	8.274

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methacrylamide	85	C4H7NO	000079-39-0	59
2			Methacrylamide	85	C4H7NO	000079-39-0	45
3			2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	40
4			2-(Bromomethyl)acrylic acid	164	C4H5BrO2	072707-66-5	40
5			Pentane, 2-bromo-4-methyl-	164	C6H13Br	030310-22-6	33



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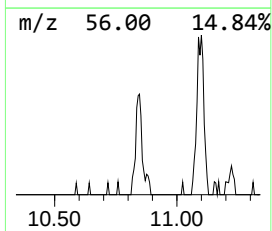
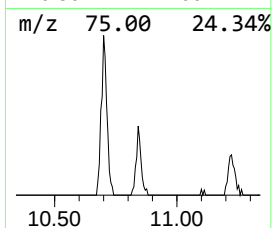
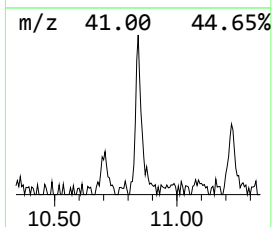
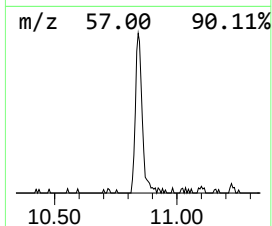
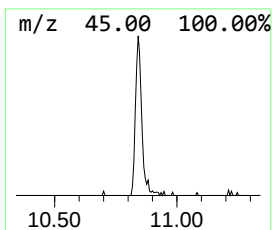
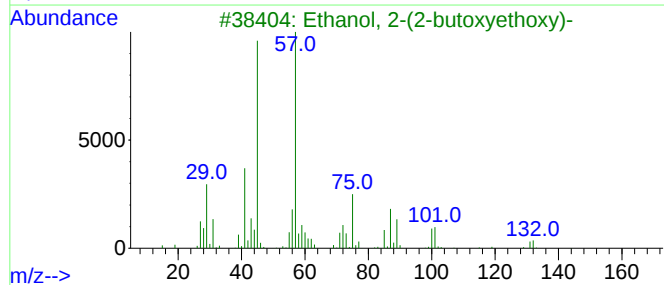
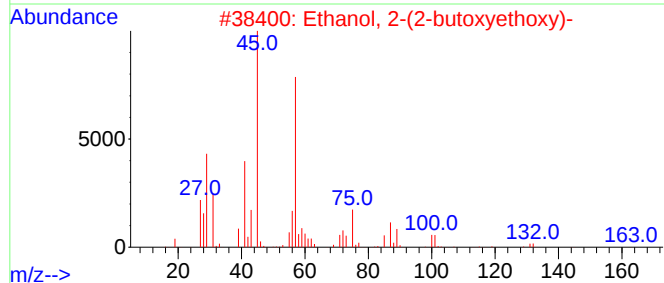
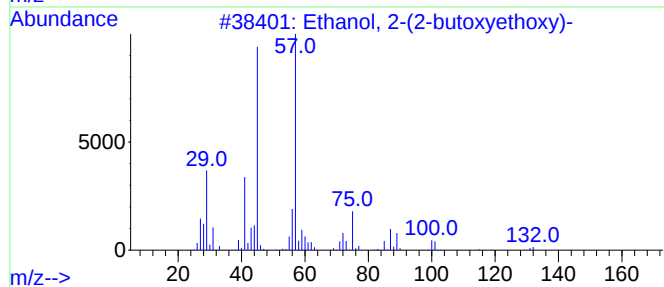
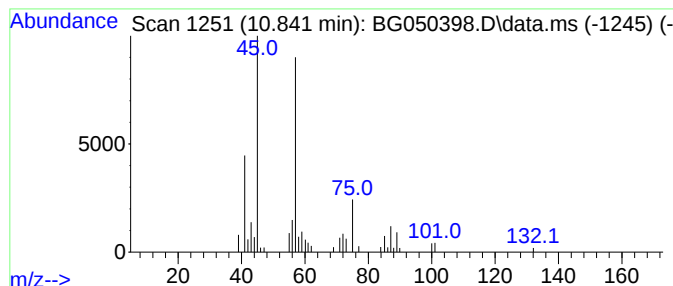
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG093021.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.841	2.19 ng/ul	63914	Naphthalene-d8	11.100

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	86
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72
5			Di-sec-Butyl ether	130	C8H18O	006863-58-7	50



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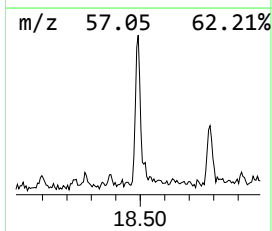
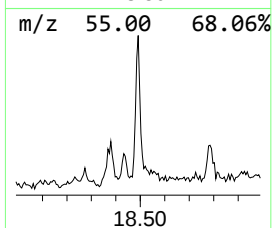
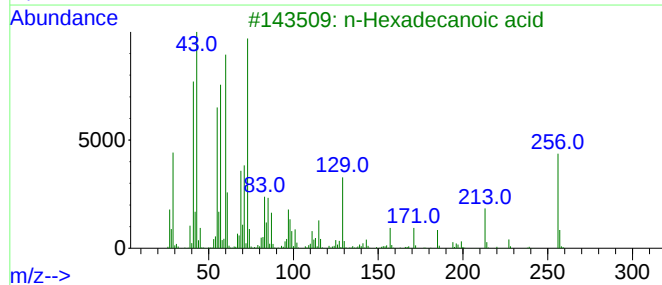
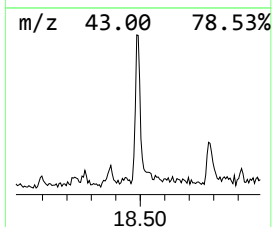
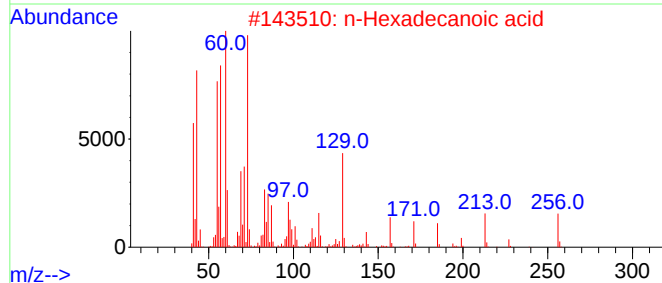
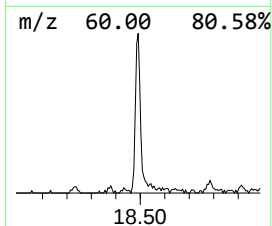
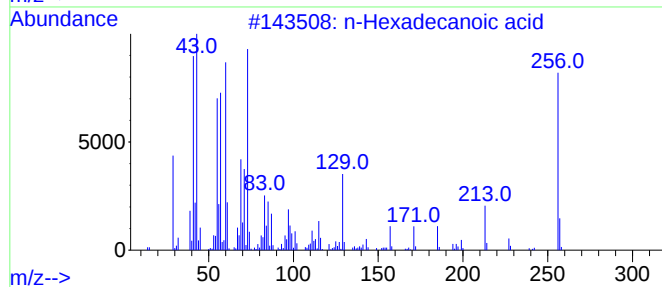
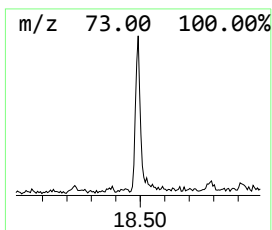
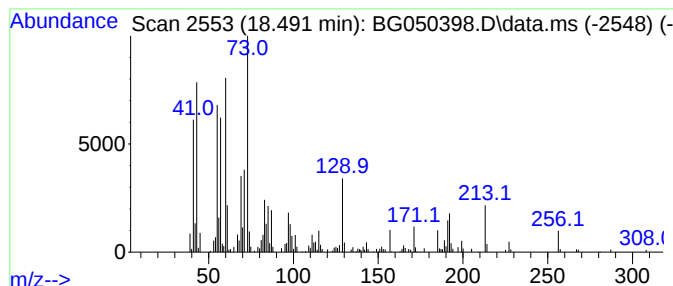
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG093021.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.491	5.88 ng/ul	257404	Phenanthrene-d10	17.633

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	91
5			Tridecanoic acid	214	C13H26O2	000638-53-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG093021\
Data File : BG050398.D
Acq On : 2 Oct 2021 21:03
Operator : CG/JU
Sample : M3877-14
Misc :
ALS Vial : 84 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW2N0

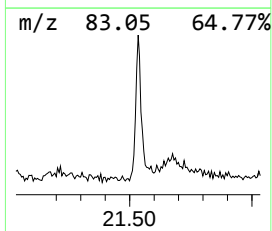
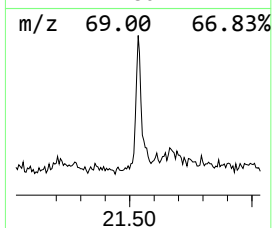
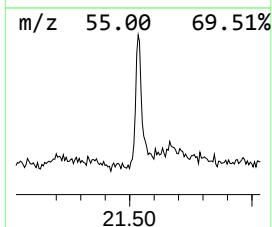
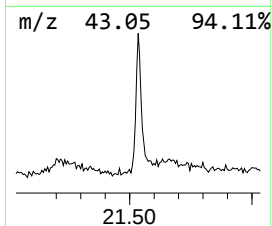
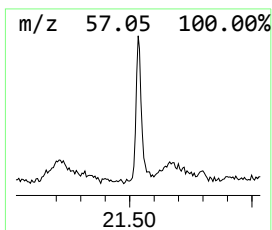
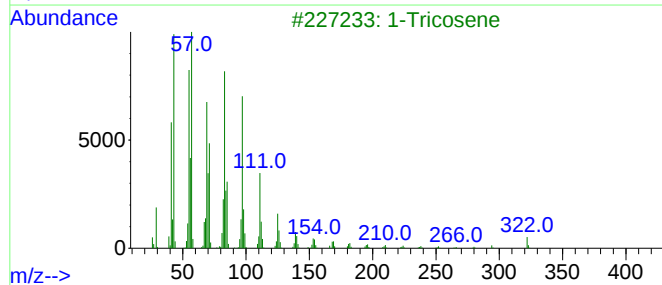
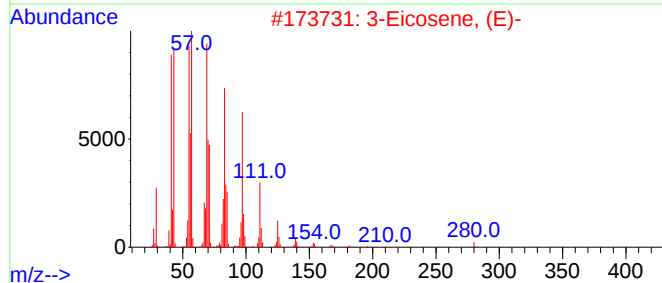
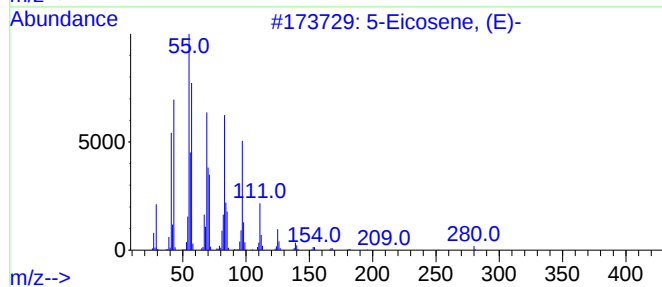
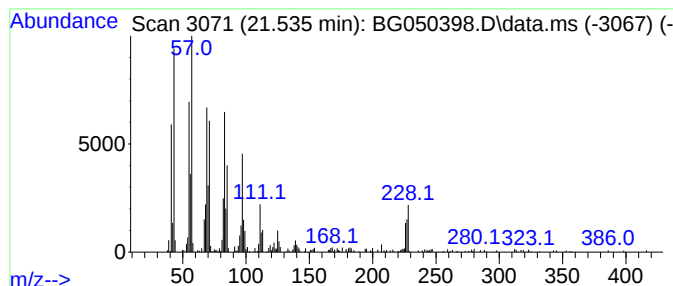
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.535	5.59 ng/ul	289156	Chrysene-d12	21.934

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-		280	C20H40	074685-30-6	96
2	3-Eicosene, (E)-		280	C20H40	074685-33-9	95
3	1-Tricosene		322	C23H46	018835-32-0	93
4	2-Chloropropionic acid, hexadec...		332	C19H37ClO2	086711-81-1	92
5	9-Eicosene, (E)-		280	C20H40	074685-29-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG093021\
Data File : BG050398.D
Acq On : 2 Oct 2021 21:03
Operator : CG/JU
Sample : M3877-14
Misc :
ALS Vial : 84 Sample Multiplier: 1

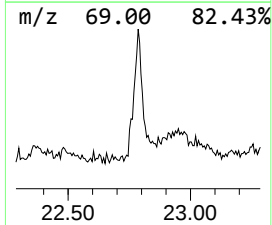
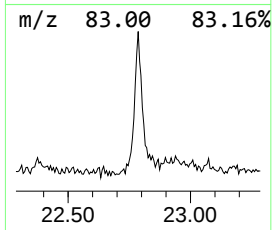
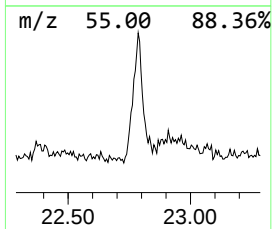
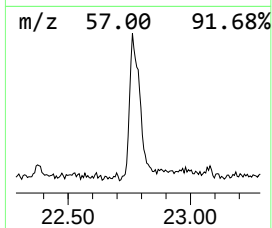
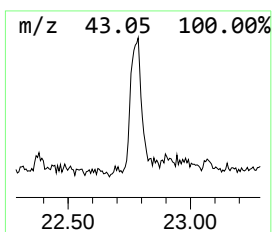
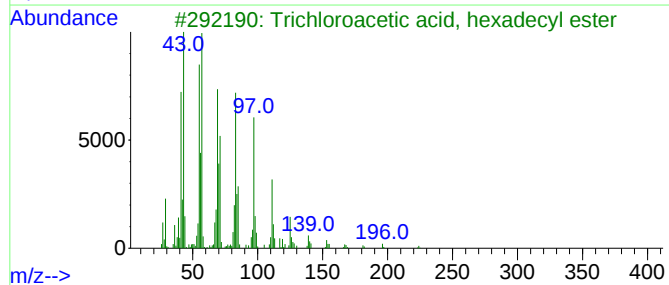
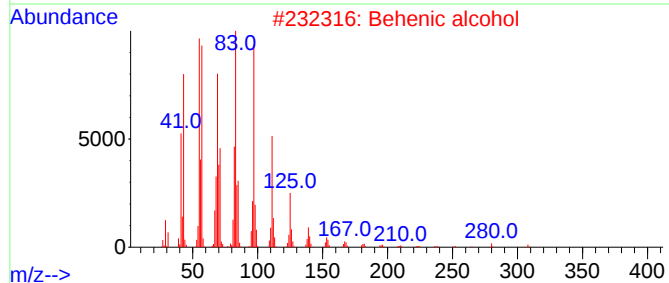
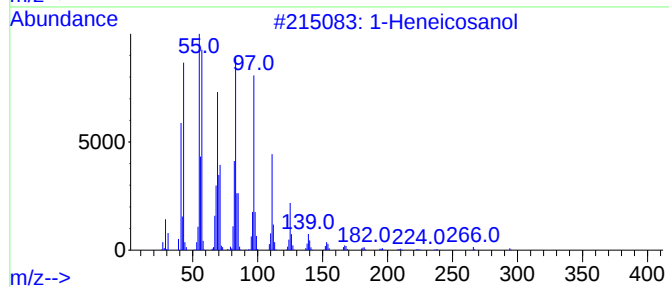
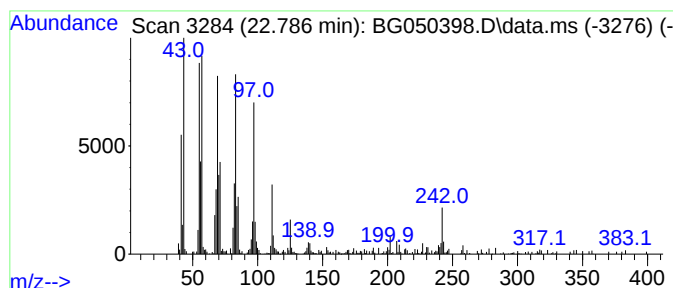
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG093021.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 1-Heneicosanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.786	8.58 ng/ul	444343	Chrysene-d12	21.934	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	93
2	Behenic alcohol	326	C22H46O	000661-19-8	93
3	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	93
4	1-Tricosene	322	C23H46	018835-32-0	93
5	1-Heneicosyl formate	340	C22H44O2	077899-03-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG093021\
Data File : BG050398.D
Acq On : 2 Oct 2021 21:03
Operator : CG/JU
Sample : M3877-14
Misc :
ALS Vial : 84 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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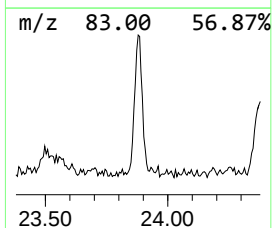
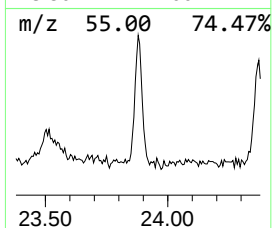
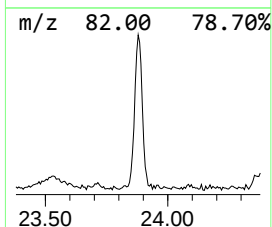
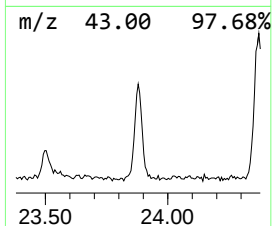
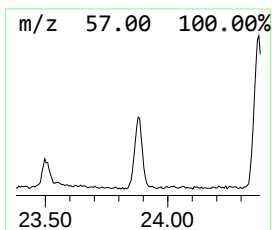
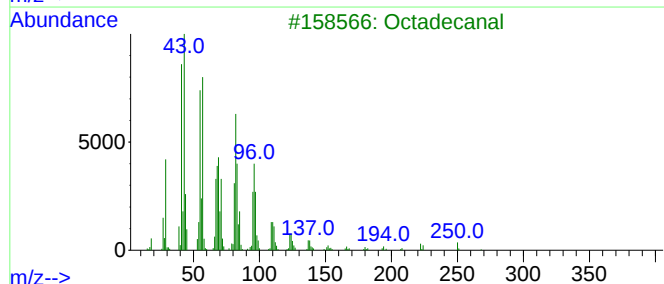
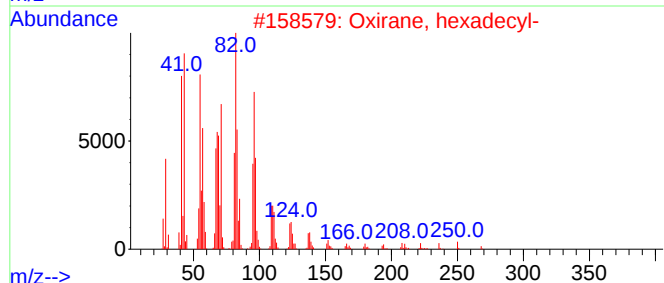
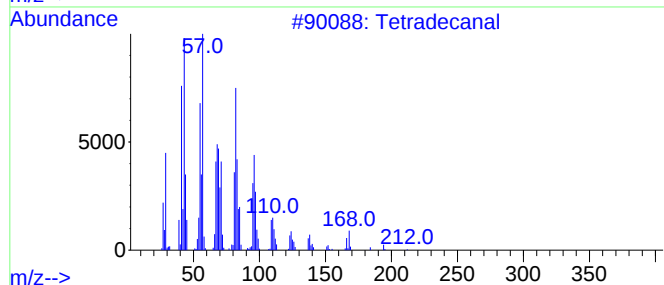
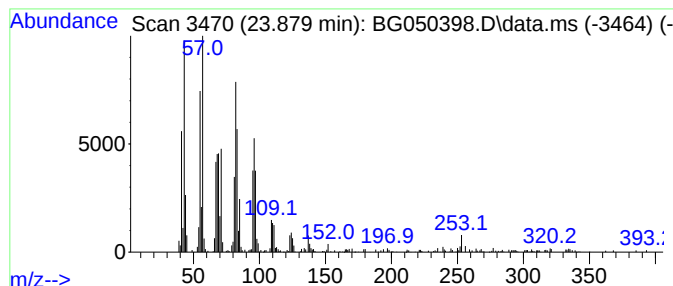
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Tetradecanal Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.879	11.87 ng/ul	482766	Perylene-d12	25.365

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecanal	212	C14H28O	000124-25-4	94
2			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	93
3			Octadecanal	268	C18H36O	000638-66-4	91
4			Hexacosanal	380	C26H52O	026627-85-0	91
5			Docosanal	324	C22H44O	057402-36-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG093021\
Data File : BG050398.D
Acq On : 2 Oct 2021 21:03
Operator : CG/JU
Sample : M3877-14
Misc :
ALS Vial : 84 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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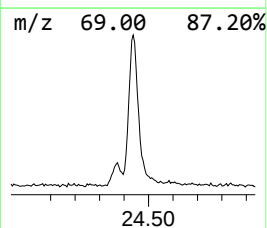
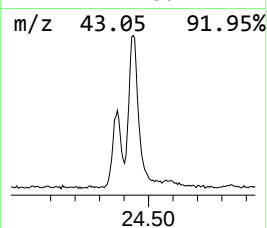
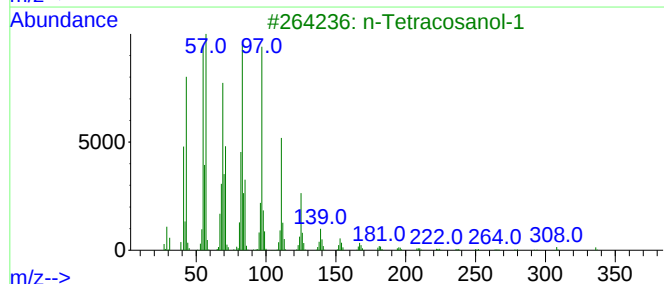
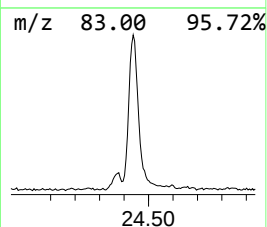
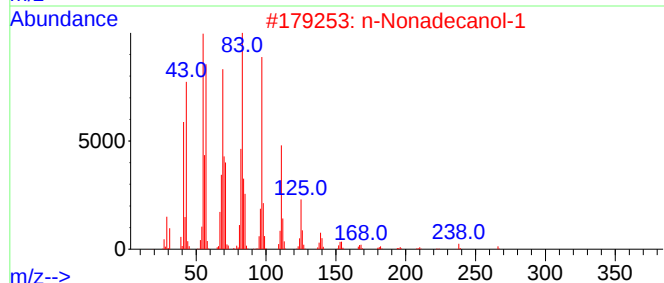
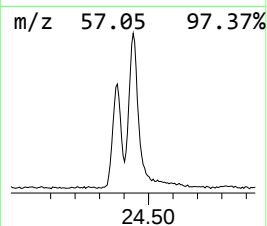
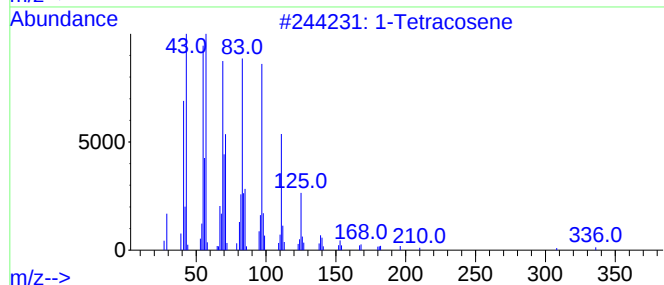
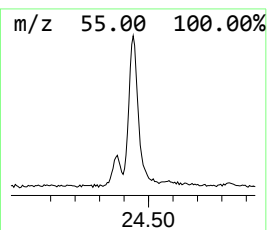
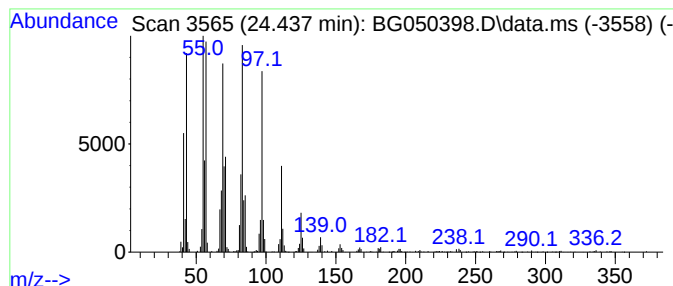
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.437	41.34 ng/ul	1680950	Perylene-d12	25.365

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene			336	C24H48	010192-32-2	99
2	n-Nonadecanol-1			284	C19H40O	001454-84-8	94
3	n-Tetracosanol-1			354	C24H50O	000506-51-4	94
4	Behenic alcohol			326	C22H46O	000661-19-8	94
5	Trifluoroacetoxy hexadecane			338	C18H33F3O2	006222-03-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG093021\
Data File : BG050398.D
Acq On : 2 Oct 2021 21:03
Operator : CG/JU
Sample : M3877-14
Misc :
ALS Vial : 84 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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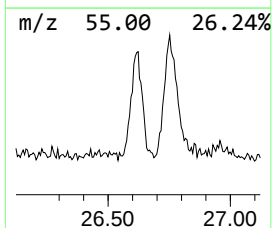
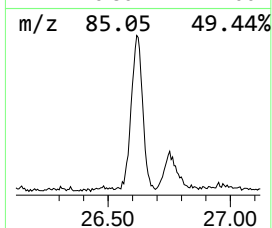
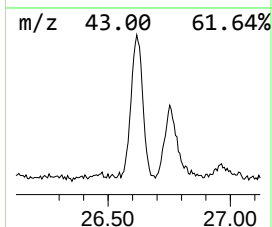
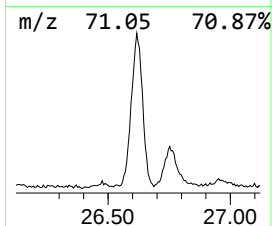
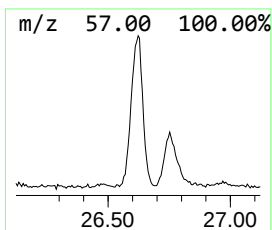
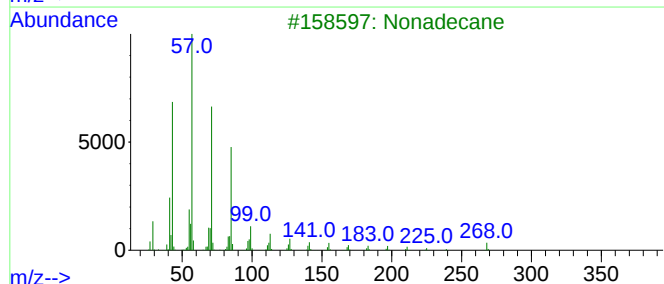
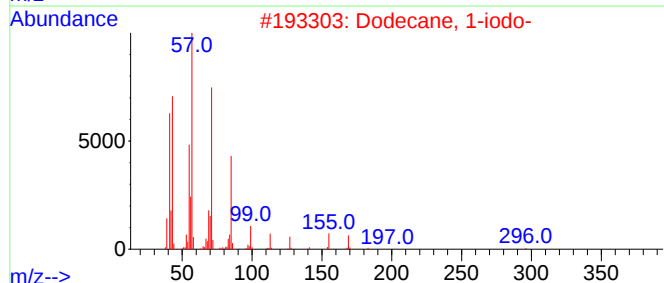
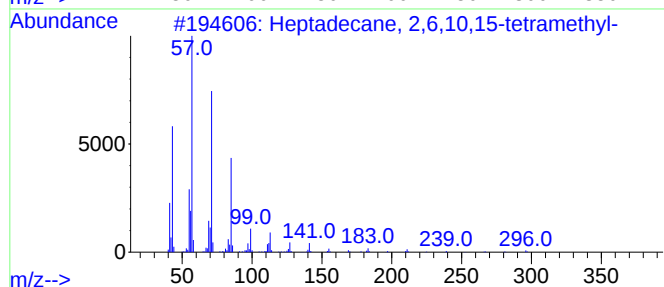
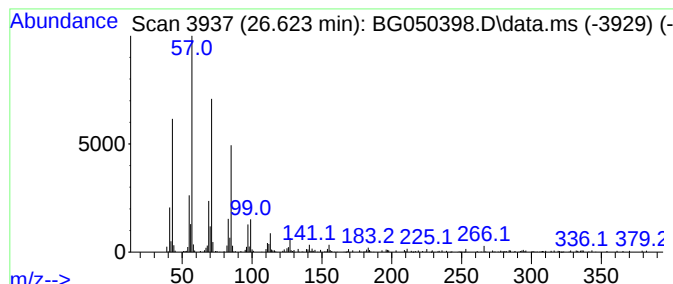
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.623	15.39 ng/ul	625652	Perylene-d12	25.365

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	96
2			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	91
3			Nonadecane	268	C19H40	000629-92-5	87
4			Tetracosane	338	C24H50	000646-31-1	87
5			Pentacosane	352	C25H52	000629-99-2	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG093021\
Data File : BG050398.D
Acq On : 2 Oct 2021 21:03
Operator : CG/JU
Sample : M3877-14
Misc :
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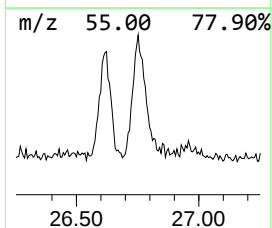
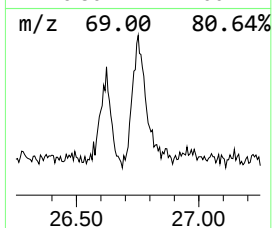
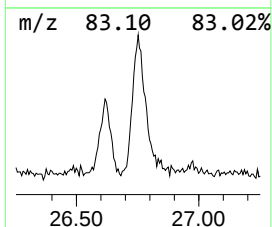
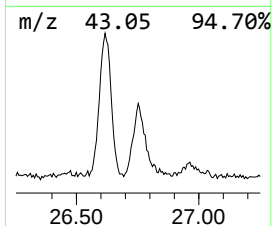
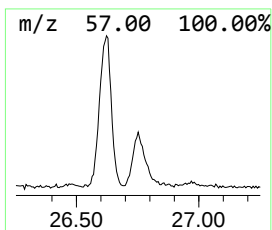
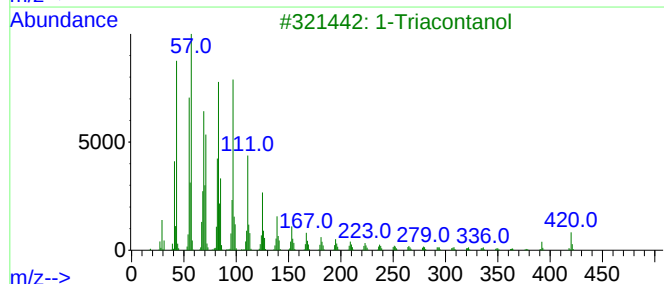
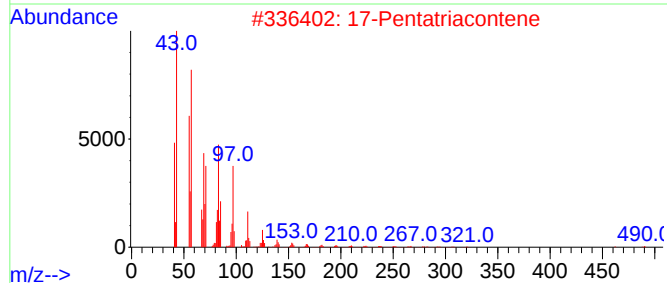
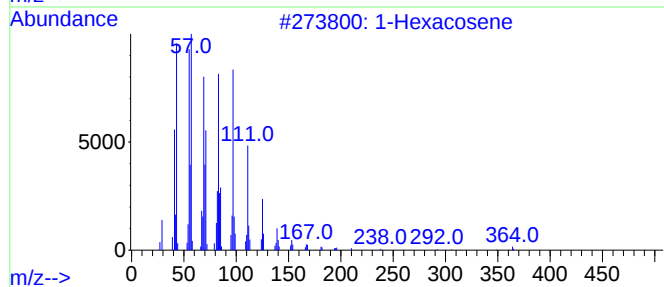
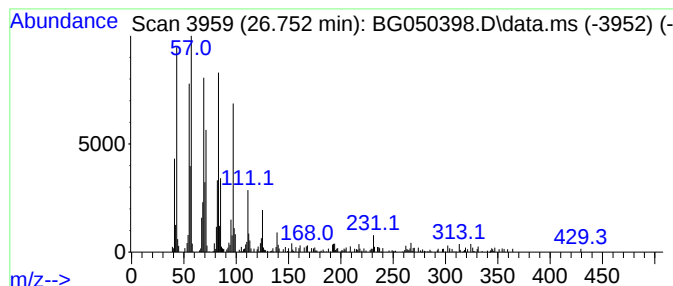
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG093021.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
26.752	10.17 ng/ul	413641	Perylene-d12		25.365	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Hexacosene		364	C26H52	018835-33-1	93
2	17-Pentatriacontene		491	C35H70	006971-40-0	91
3	1-Triacontanol		438	C30H62O	000593-50-0	64
4	1-Nonadecene		266	C19H38	018435-45-5	58
5	Heneicosyl heptafluorobutyrate		508	C25H43F7O2	1000351-83-8	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG093021\
Data File : BG050398.D
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Sample : M3877-14
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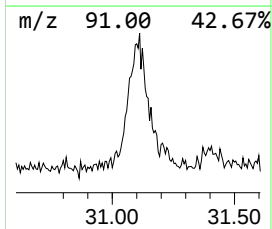
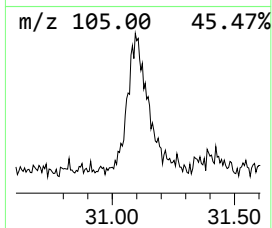
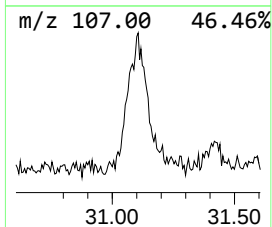
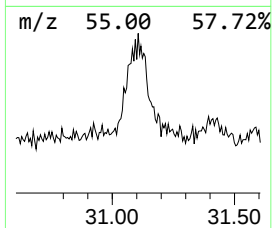
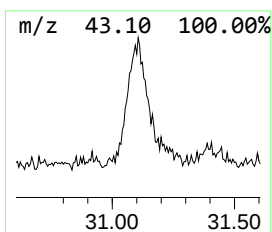
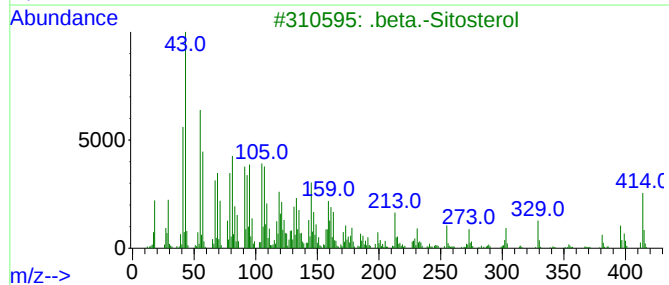
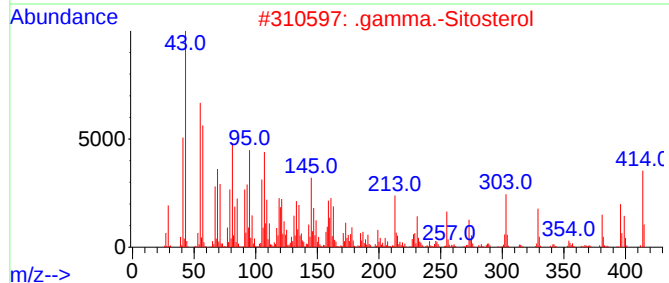
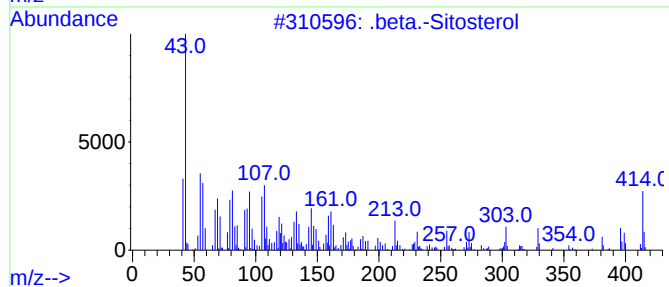
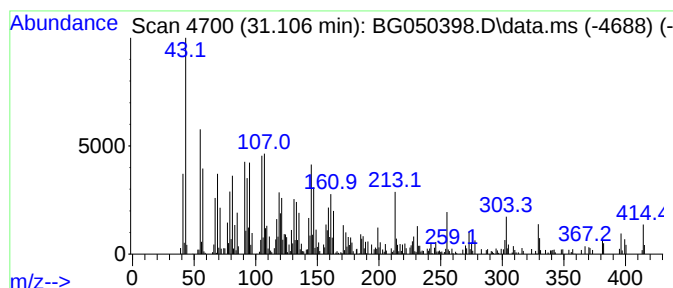
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ClientSampleId :
EW2N0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG093021.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 .beta.-Sitosterol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
31.106	20.21 ng/ul	821645	Perylene-d12		25.365	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	.beta.-Sitosterol		414	C29H50O	000083-46-5	94
2	.gamma.-Sitosterol		414	C29H50O	000083-47-6	90
3	.beta.-Sitosterol		414	C29H50O	000083-46-5	90
4	Campesterol		400	C28H48O	000474-62-4	58
5	Ergost-5-en-3-ol, (3.beta.)-		400	C28H48O	004651-51-8	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG093021\
 Data File : BG050398.D
 Acq On : 2 Oct 2021 21:03
 Operator : CG/JU
 Sample : M3877-14
 Misc :
 ALS Vial : 84 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW2N0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG093021.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
Methacrylamide	4.173	3.6	ng/ul	74253	1	8.274	414326	20.0
Ethanol, 2-(2-b...	10.841	2.2	ng/ul	63914	2	11.100	582895	20.0
n-Hexadecanoic ...	18.491	5.9	ng/ul	257404	4	17.633	875549	20.0
(DEL) Alkane: S...	21.535	5.6	ng/ul	289156	5	21.934	1035210	20.0
1-Heneicosanol	22.786	8.6	ng/ul	444343	5	21.934	1035210	20.0
Tetradecanal	23.879	11.9	ng/ul	482766	6	25.365	813192	20.0
(DEL) Alkane: S...	24.437	41.3	ng/ul	1680950	6	25.365	813192	20.0
(DEL) Alkane: S...	26.623	15.4	ng/ul	625652	6	25.365	813192	20.0
(DEL) Alkane: S...	26.752	10.2	ng/ul	413641	6	25.365	813192	20.0
.beta.-Sitosterol	31.106	20.2	ng/ul	821645	6	25.365	813192	20.0