

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071124\
 Data File : BG062031.D
 Acq On : 12 Jul 2024 6:40
 Operator : MA/JU
 Sample : P3124-01
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4CB3

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-BG070224.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG062031.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.440	23	26	28	rVB4	9572	10326	0.19%	0.030%
2	3.711	66	72	79	rBV	100310	166796	3.14%	0.480%
3	3.858	93	97	107	rBV	86122	154720	2.91%	0.445%
4	3.969	111	116	123	rVB2	47537	81130	1.53%	0.233%
5	4.040	126	128	132	rBV4	5445	7935	0.15%	0.023%
6	4.198	152	155	161	rVB6	4901	7922	0.15%	0.023%
7	4.627	223	228	231	rBV4	4045	8020	0.15%	0.023%
8	4.751	246	249	253	rBV3	11699	15868	0.30%	0.046%
9	5.003	288	292	297	rVB6	6265	9802	0.18%	0.028%
10	5.115	305	311	316	rVV2	20018	34665	0.65%	0.100%
11	5.209	319	327	333	rVB2	13479	26115	0.49%	0.075%
12	5.362	350	353	357	rVB3	5360	6560	0.12%	0.019%
13	6.090	475	477	480	rVB4	5282	5517	0.10%	0.016%
14	6.584	560	561	567	rVB3	4983	6207	0.12%	0.018%
15	7.124	651	653	657	rBV5	4224	6255	0.12%	0.018%
16	7.195	659	665	674	rVB	362300	645718	12.15%	1.858%
17	7.365	687	694	701	rBV	221726	359070	6.75%	1.033%
18	7.571	722	729	734	rBV	349643	584070	10.99%	1.681%
19	7.618	734	737	745	rVB2	46700	73917	1.39%	0.213%
20	7.888	782	783	789	rVB3	5204	5750	0.11%	0.017%
21	8.041	800	809	815	rBV	242793	417711	7.86%	1.202%
22	8.746	922	929	938	rBV	421436	712994	13.41%	2.052%
23	8.864	943	949	955	rBV4	17399	27896	0.52%	0.080%
24	9.204	1000	1007	1019	rBV	333932	588577	11.07%	1.694%
25	9.539	1062	1064	1069	rBV3	5322	6638	0.12%	0.019%
26	9.751	1095	1100	1105	rBV2	19717	34212	0.64%	0.098%
27	9.927	1123	1130	1140	rBV	317806	553899	10.42%	1.594%
28	10.144	1165	1167	1176	rVB7	7349	15355	0.29%	0.044%
29	10.468	1215	1222	1230	rBV	467119	843562	15.87%	2.428%
30	10.603	1243	1245	1251	rBV5	4425	9108	0.17%	0.026%
31	10.849	1280	1287	1294	rBV2	338127	620750	11.68%	1.786%
32	10.985	1304	1310	1322	rBV2	135613	262624	4.94%	0.756%
33	11.366	1371	1375	1383	rVB7	17073	27901	0.52%	0.080%
34	11.584	1406	1412	1426	rBV	133071	270822	5.09%	0.779%
35	12.424	1553	1555	1560	rVB2	7326	9732	0.18%	0.028%
36	12.500	1562	1568	1574	rBV8	6183	13315	0.25%	0.038%

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37	12.718	1603	1605	1610	rVB3	7008	9481	0.18%	0.027%
38	12.824	1616	1623	1629	rBV8	5261	10285	0.19%	0.030%
39	13.288	1697	1702	1709	rVB6	7041	16081	0.30%	0.046%
40	13.540	1743	1745	1749	rBV4	4945	6952	0.13%	0.020%
41	13.852	1794	1798	1802	rVB5	8418	13596	0.26%	0.039%
42	13.987	1817	1821	1822	rBV2	8021	9305	0.18%	0.027%
43	14.069	1827	1835	1842	rBV	773315	1167460	21.96%	3.360%
44	14.304	1873	1875	1879	rBV5	9289	11575	0.22%	0.033%
45	14.363	1879	1885	1899	rVB2	772921	1278141	24.04%	3.678%
46	14.545	1914	1916	1920	rVB3	11897	14105	0.27%	0.041%
47	14.668	1930	1937	1944	rBV	536419	834660	15.70%	2.402%
48	14.733	1944	1948	1954	rVB2	26434	41879	0.79%	0.121%
49	14.886	1963	1974	1985	rBV	2341752	5315709	100.00%	15.297%
50	15.015	1994	1996	2001	rBV3	24391	33663	0.63%	0.097%
51	15.062	2001	2004	2011	rVB4	32288	56816	1.07%	0.164%
52	15.268	2037	2039	2044	rVB5	8021	9186	0.17%	0.026%
53	15.374	2056	2057	2060	rBV3	5708	6370	0.12%	0.018%
54	15.438	2064	2068	2069	rBV4	9255	11014	0.21%	0.032%
55	15.661	2099	2106	2112	rBV	1077086	1600920	30.12%	4.607%
56	15.714	2112	2115	2119	rVB	33345	45757	0.86%	0.132%
57	15.773	2119	2125	2134	rBV	432180	656331	12.35%	1.889%
58	15.879	2142	2143	2145	rBV2	8309	6491	0.12%	0.019%
59	16.155	2187	2190	2194	rBV5	13041	19880	0.37%	0.057%
60	16.196	2194	2197	2200	rVV3	8763	11215	0.21%	0.032%
61	16.361	2220	2225	2228	rBV6	18734	33990	0.64%	0.098%
62	16.684	2277	2280	2283	rBV5	7484	11625	0.22%	0.033%
63	16.978	2328	2330	2334	rBV4	9653	7924	0.15%	0.023%
64	17.066	2340	2345	2350	rBV2	27664	49221	0.93%	0.142%
65	17.154	2356	2360	2364	rBV7	29782	43336	0.82%	0.125%
66	17.236	2370	2374	2379	rVB3	26424	39143	0.74%	0.113%
67	17.412	2398	2404	2408	rBV	675726	1046939	19.70%	3.013%
68	17.459	2408	2412	2416	rVV	541823	805002	15.14%	2.317%
69	17.512	2416	2421	2432	rVB	1241400	1925359	36.22%	5.541%
70	17.818	2469	2473	2476	rVB2	46337	64786	1.22%	0.186%
71	18.000	2499	2504	2511	rBV10	31685	70251	1.32%	0.202%
72	18.064	2511	2515	2519	rBV5	40489	50199	0.94%	0.144%
73	18.288	2549	2553	2558	rBV2	291891	429722	8.08%	1.237%
74	18.335	2558	2561	2570	rVB2	111147	182521	3.43%	0.525%
75	18.482	2581	2586	2590	rBV3	127097	234798	4.42%	0.676%
76	18.576	2600	2602	2607	rVB4	38983	45989	0.87%	0.132%

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77	18.781	2634	2637	2640	rVV2	52099	63471	1.19%	0.183%
78	18.822	2641	2644	2648	rVB	110184	149214	2.81%	0.429%
79	19.463	2748	2753	2759	rVB	1317896	1813963	34.12%	5.220%
80	19.557	2767	2769	2773	rVB5	56530	56946	1.07%	0.164%
81	19.798	2804	2810	2813	rBV	1356194	2222883	41.82%	6.397%
82	19.827	2813	2815	2820	rVB	890826	961783	18.09%	2.768%
83	20.139	2864	2868	2875	rBV5	75318	195056	3.67%	0.561%
84	20.309	2895	2897	2901	rVB	242048	299304	5.63%	0.861%
85	21.308	3063	3067	3079	rVB4	411206	807361	15.19%	2.323%
86	21.549	3104	3108	3113	rVB	432462	600384	11.29%	1.728%
87	21.672	3121	3129	3133	rBV4	798022	1971870	37.10%	5.675%
88	21.719	3134	3137	3148	rVB	543950	886220	16.67%	2.550%
89	22.483	3259	3267	3277	rBV7	247190	830909	15.63%	2.391%
90	24.663	3632	3638	3644	rBV2	445802	1030563	19.39%	2.966%

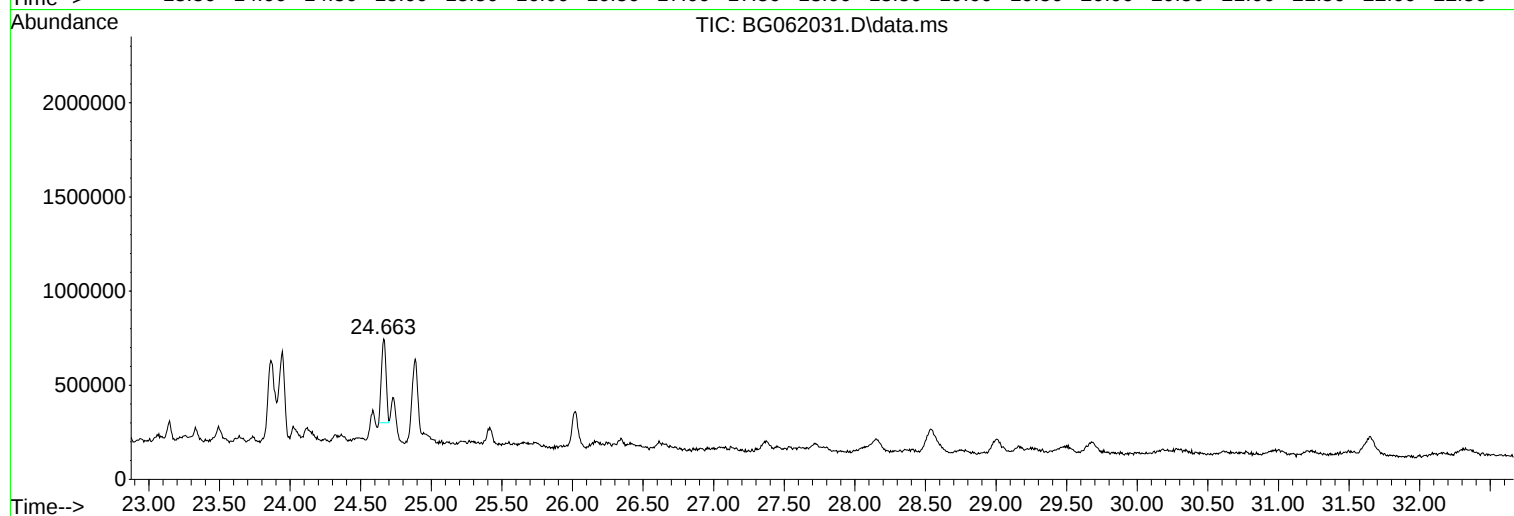
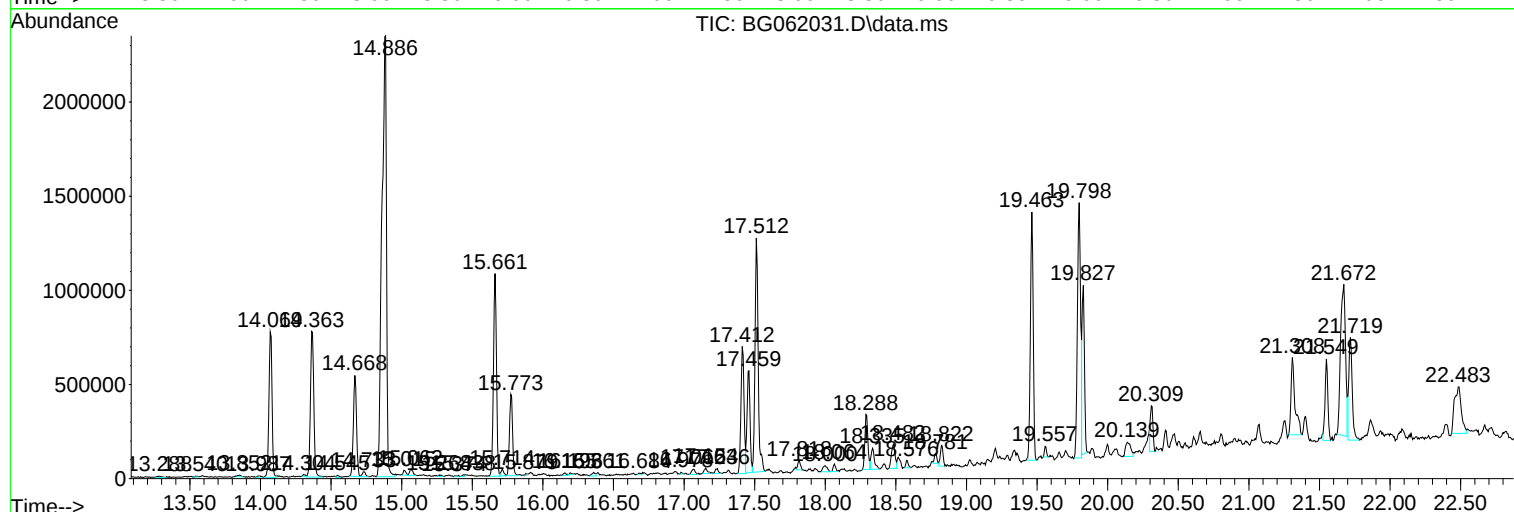
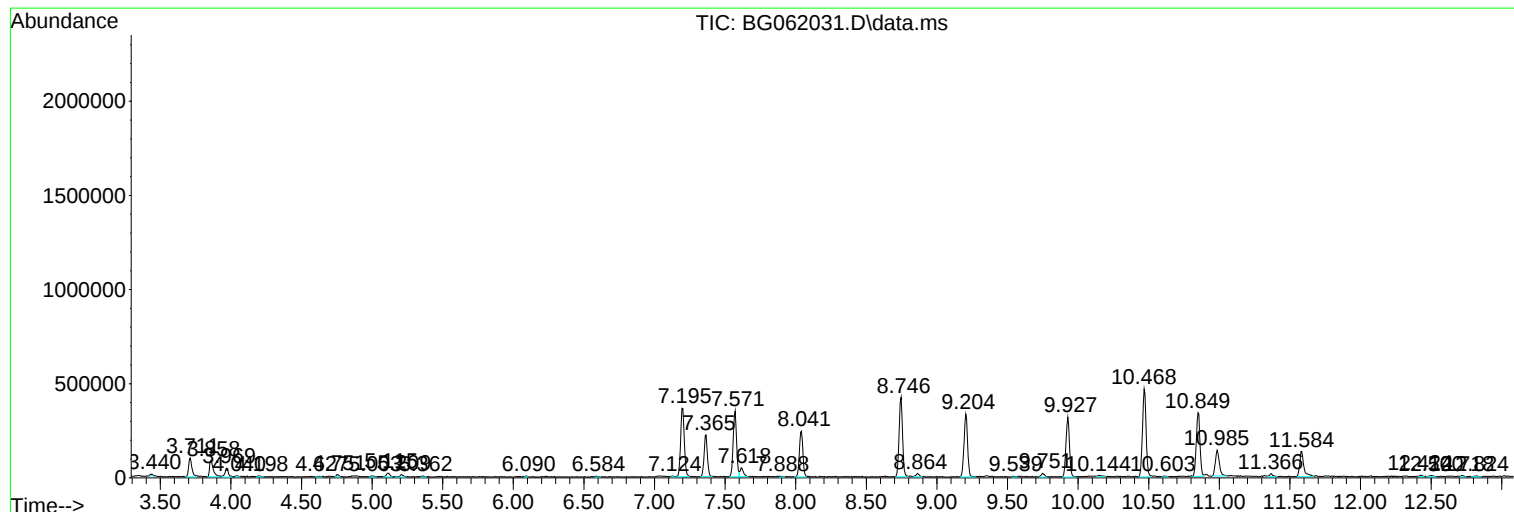
Sum of corrected areas: 34749063

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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



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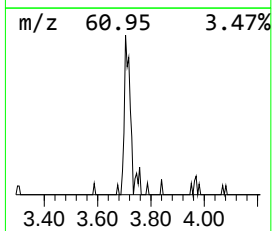
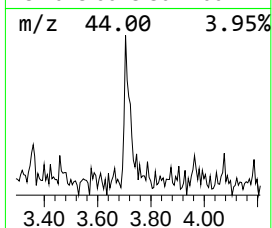
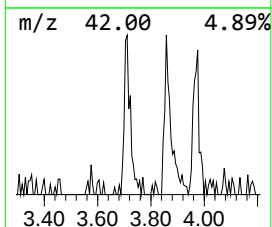
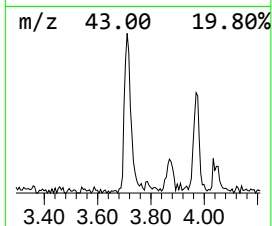
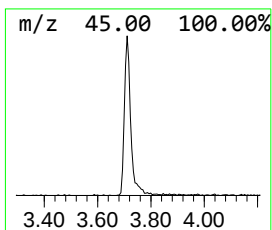
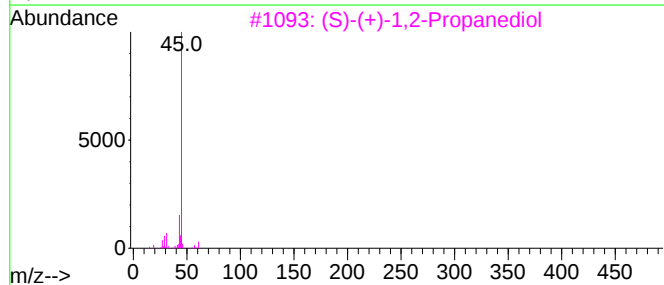
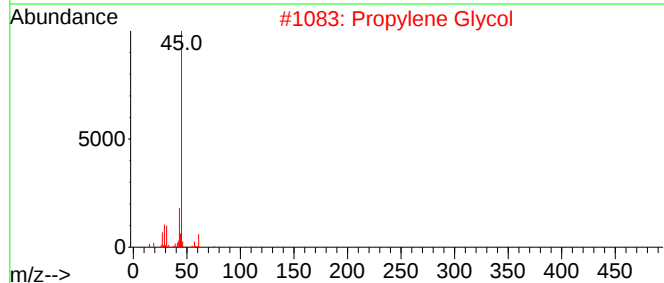
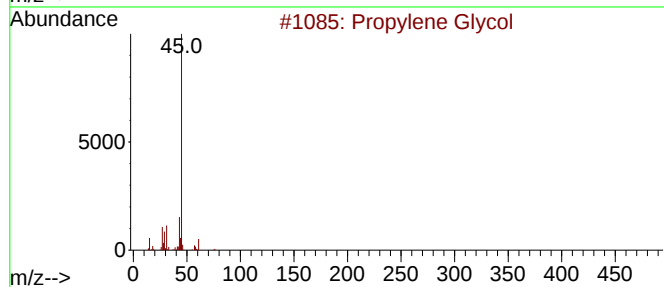
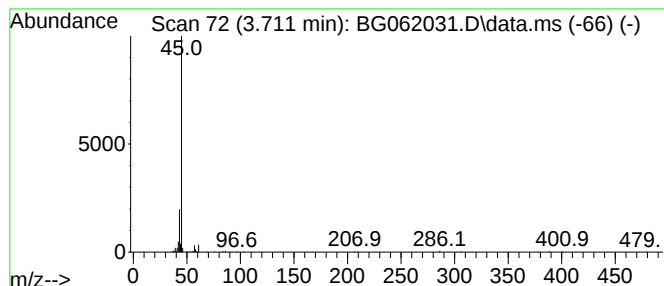
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.711	7.99 ng/ul	166796	1,4-Dichlorobenzene-d4	8.041

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	40
2			Propylene Glycol	76	C3H8O2	000057-55-6	9
3			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	9
4			2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	9
5			Isopropyl Alcohol	60	C3H8O	000067-63-0	9



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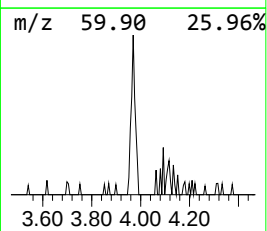
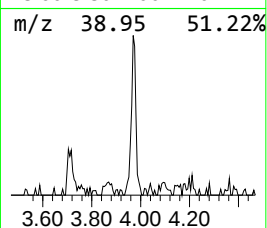
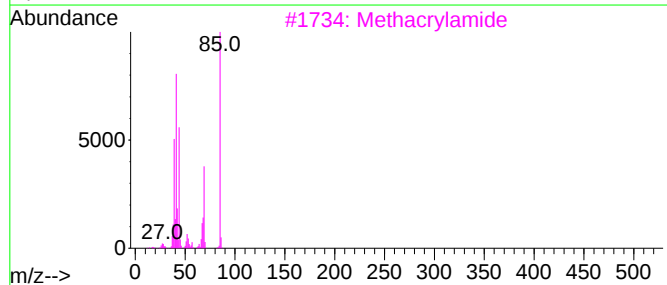
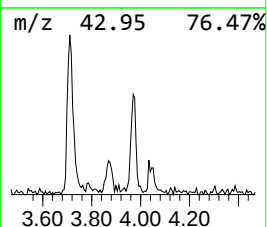
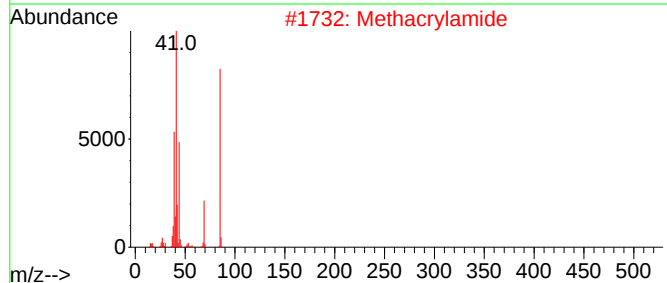
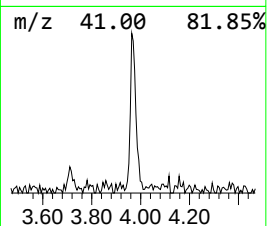
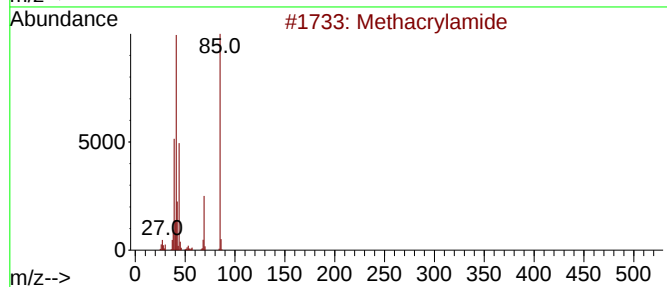
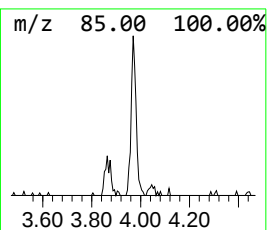
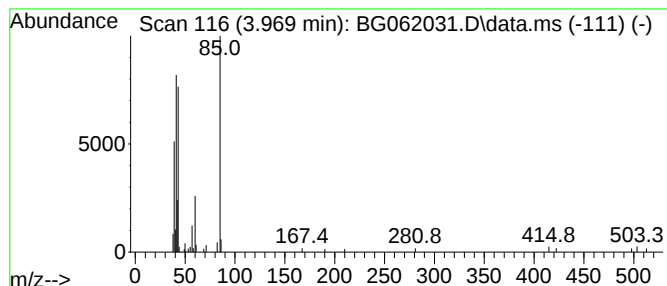
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Methacrylamide Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.969	3.88 ng/ul	81130	1,4-Dichlorobenzene-d4	8.041

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methacrylamide	85	C4H7NO	000079-39-0	72
2			Methacrylamide	85	C4H7NO	000079-39-0	72
3			Methacrylamide	85	C4H7NO	000079-39-0	59
4			2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	43
5			Cyclopropanecarboxylic acid	86	C4H6O2	001759-53-1	35



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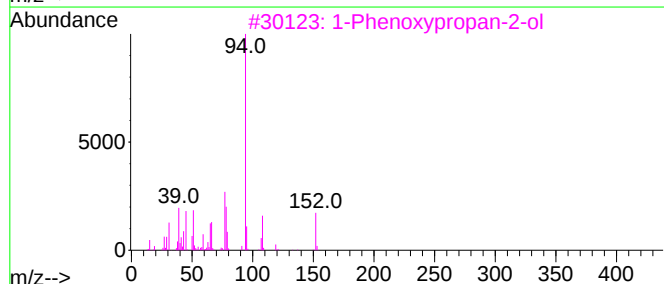
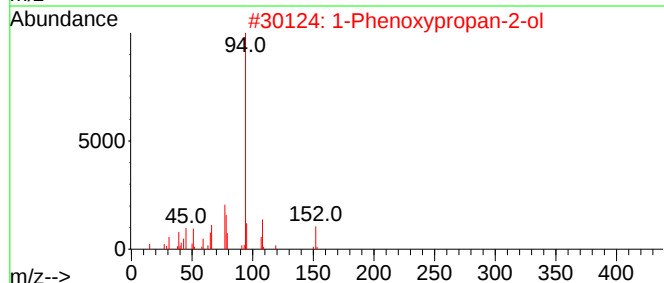
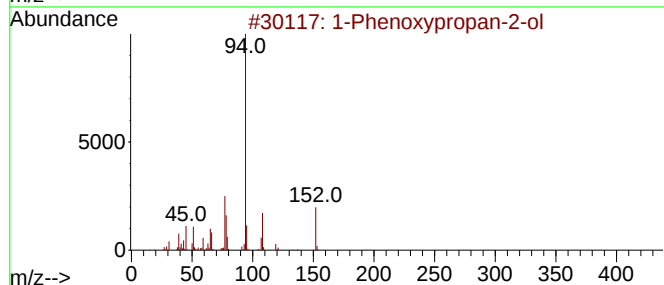
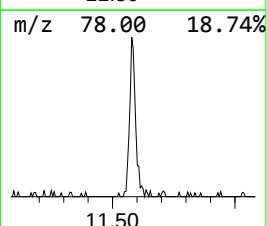
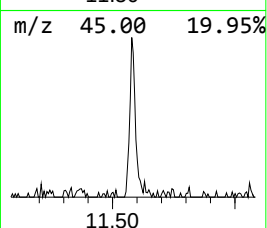
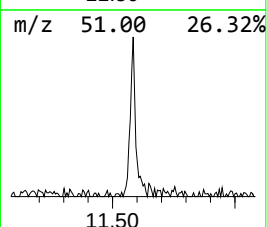
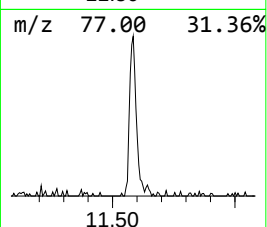
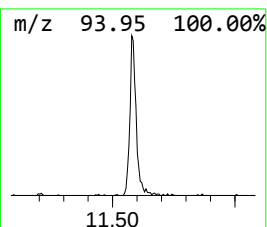
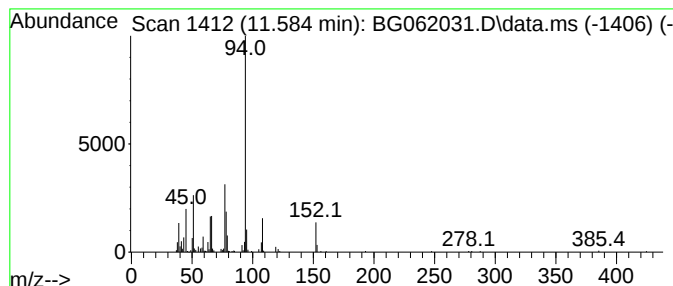
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.584	8.73 ng/ul	270822	Naphthalene-d8	10.849

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	90
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	90
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	90
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	81
5			2-Phenoxybutyric acid	180	C10H12O3	013794-14-4	72



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Sample : P3124-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4CB3

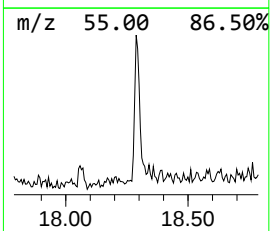
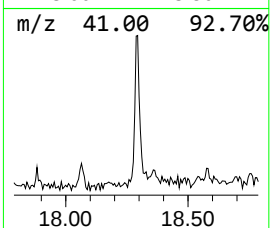
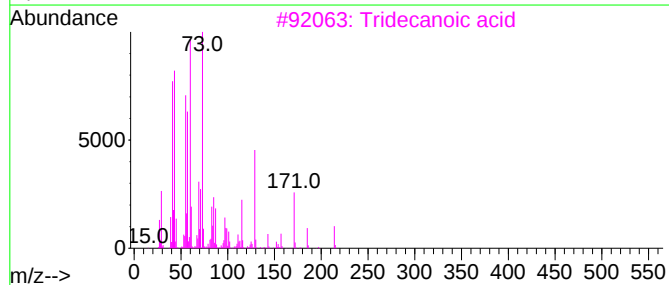
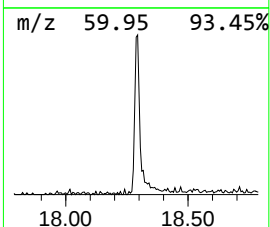
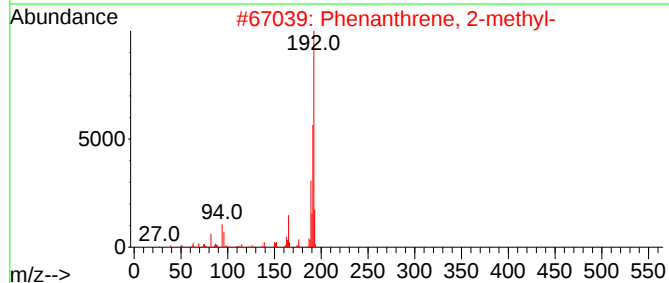
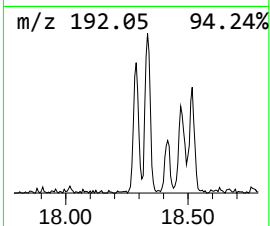
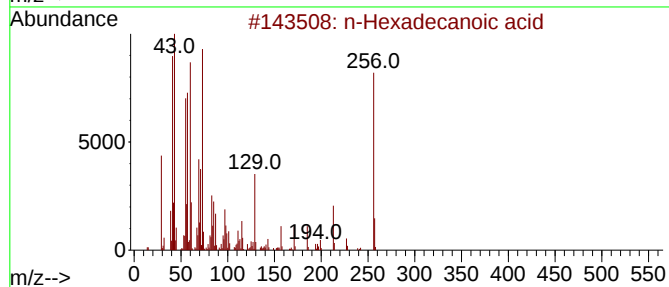
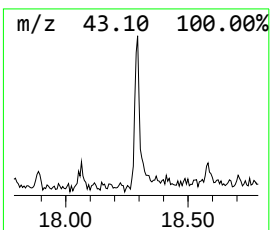
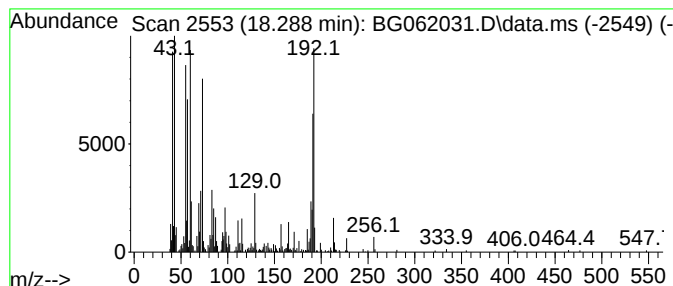
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.288	8.21 ng/ul	429722	Phenanthrene-d10	17.412

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	70
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	55
3			Tridecanoic acid	214	C13H26O2	000638-53-9	45
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	43
5			n-Decanoic acid	172	C10H20O2	000334-48-5	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071124\
Data File : BG062031.D
Acq On : 12 Jul 2024 6:40
Operator : MA/JU
Sample : P3124-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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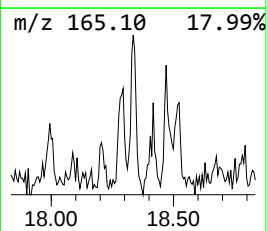
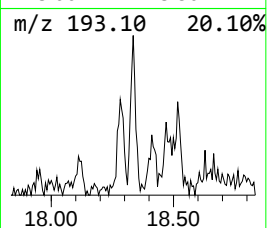
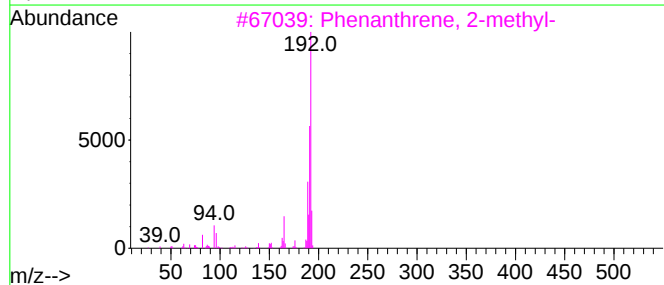
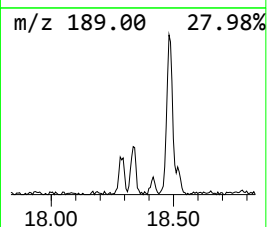
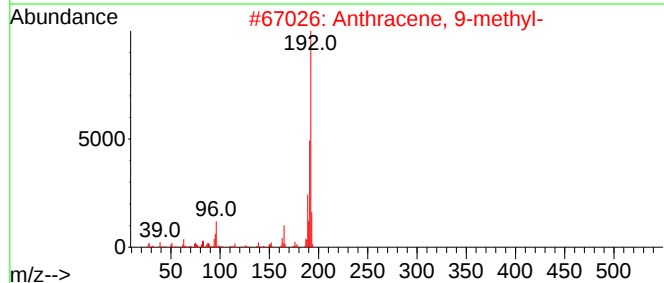
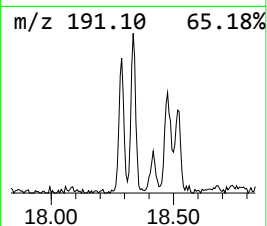
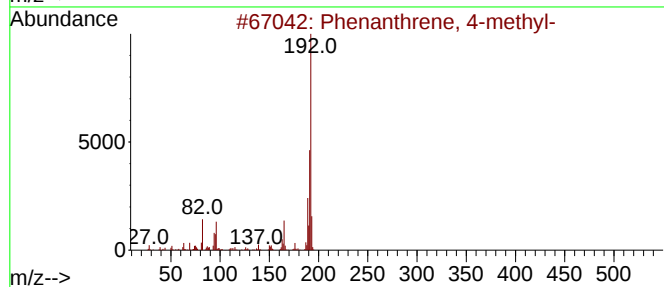
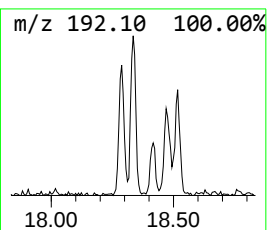
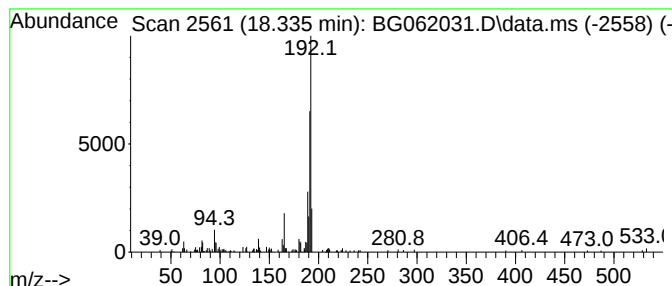
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Phenanthrene, 4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.335	3.49 ng/ul	182521	Phenanthrene-d10	17.412

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	87
2			Anthracene, 9-methyl-	192	C15H12	000779-02-2	87
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	83
4			Anthracene, 9-methyl-	192	C15H12	000779-02-2	81
5			Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071124\
Data File : BG062031.D
Acq On : 12 Jul 2024 6:40
Operator : MA/JU
Sample : P3124-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

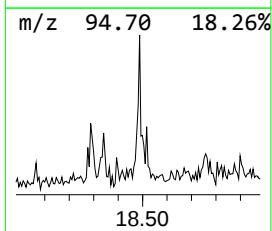
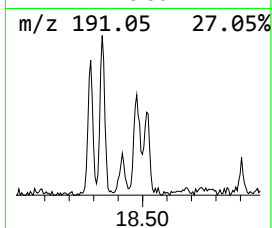
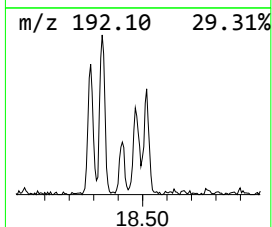
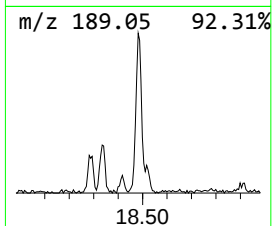
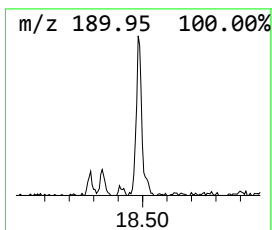
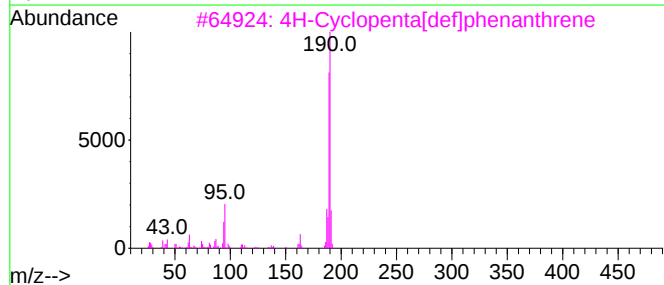
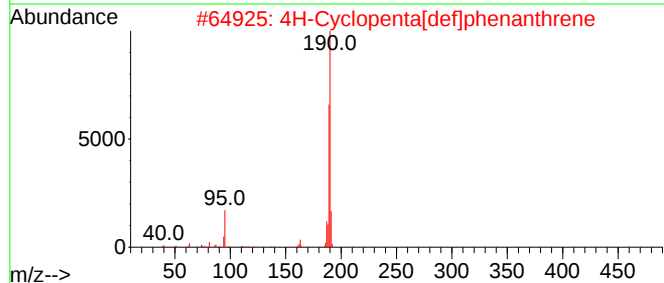
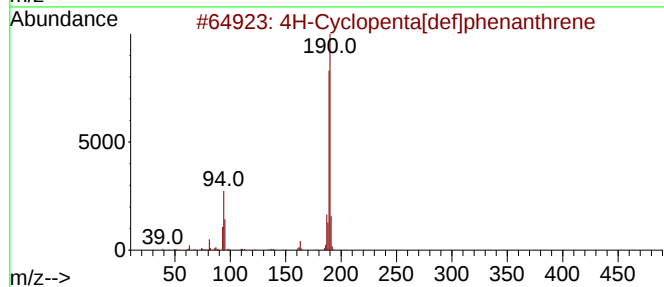
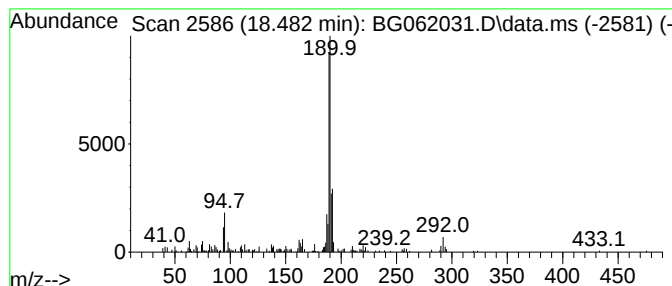
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.482	4.49 ng/ul	234798	Phenanthrene-d10	17.412	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	81
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	74
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	72
4	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	50
5	1H-Pyrazole-4-carboxaldehyde, 1-...	190	C10H7FN2O	1000338-05-6	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071124\
Data File : BG062031.D
Acq On : 12 Jul 2024 6:40
Operator : MA/JU
Sample : P3124-01
Misc :
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Instrument :
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ClientSampleId :
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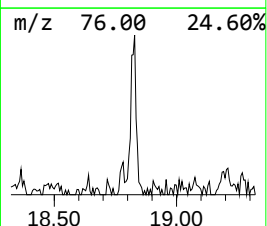
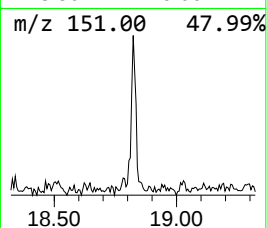
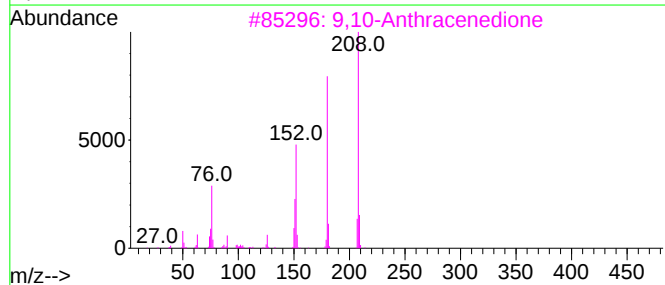
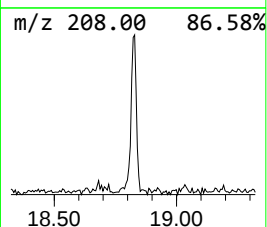
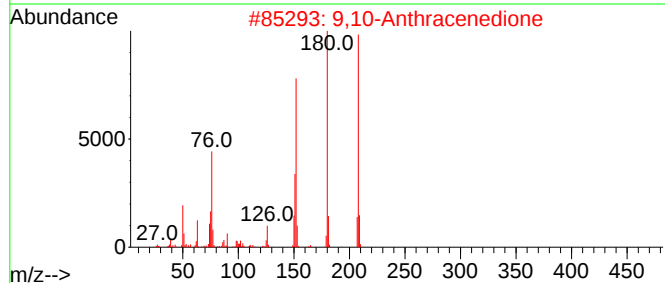
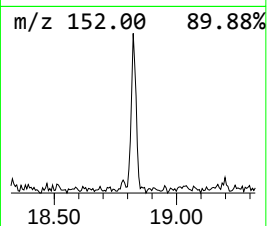
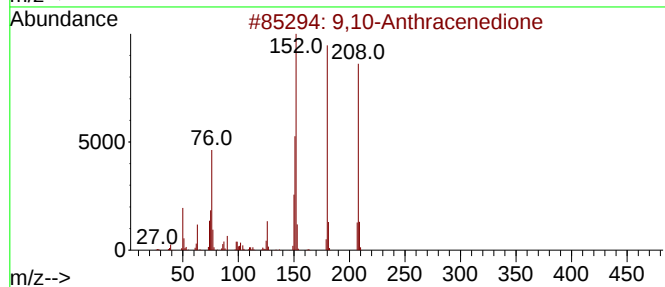
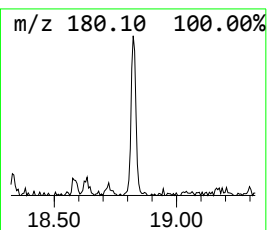
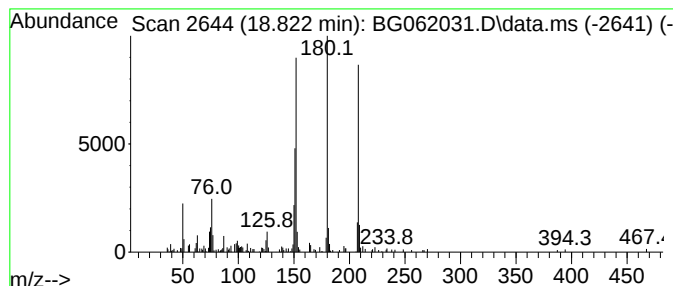
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9,10-Anthracenedione Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.822	2.85 ng/ul	149214	Phenanthrene-d10	17.412

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione			208	C14H8O2	000084-65-1	96
2	9,10-Anthracenedione			208	C14H8O2	000084-65-1	93
3	9,10-Anthracenedione			208	C14H8O2	000084-65-1	93
4	9,10-Phenanthrenedione			208	C14H8O2	000084-11-7	91
5	9,10-Anthracenedione			208	C14H8O2	000084-65-1	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071124\
Data File : BG062031.D
Acq On : 12 Jul 2024 6:40
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Sample : P3124-01
Misc :
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Instrument :
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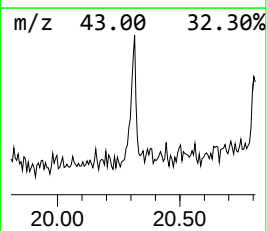
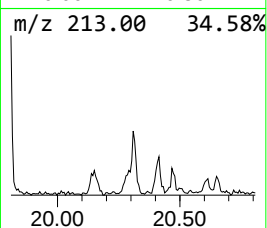
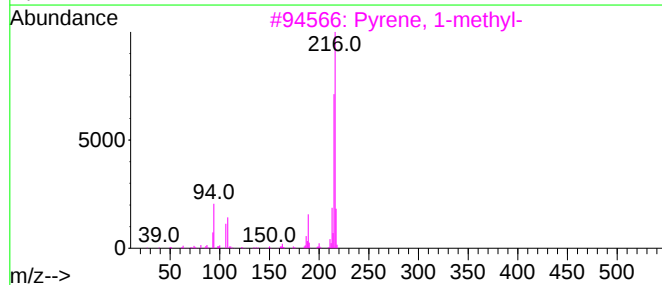
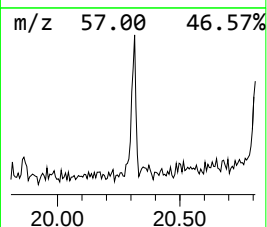
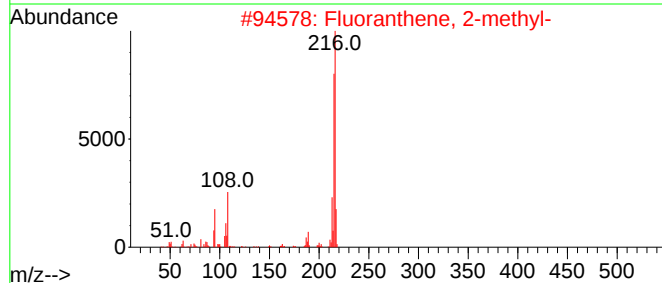
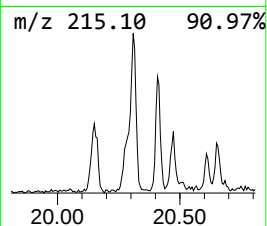
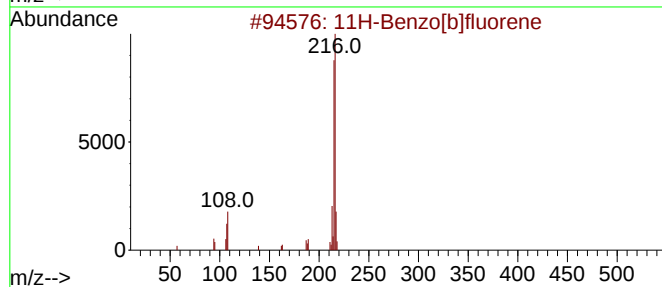
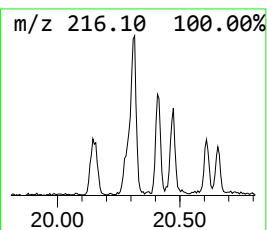
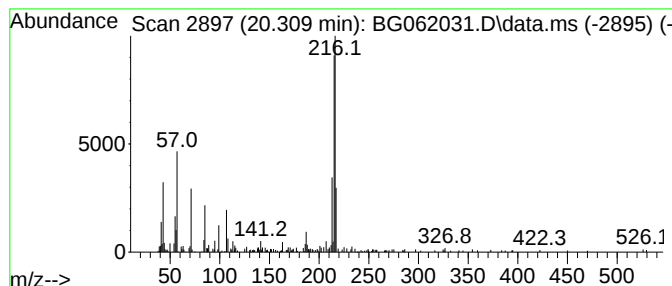
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 11H-Benzo[b]fluorene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.309	3.04 ng/ul	299304	Chrysene-d12	21.672

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	58
2			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	53
3			Pyrene, 1-methyl-	216	C17H12	002381-21-7	50
4			Pyrene, 2-methyl-	216	C17H12	003442-78-2	50
5			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071124\
Data File : BG062031.D
Acq On : 12 Jul 2024 6:40
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Sample : P3124-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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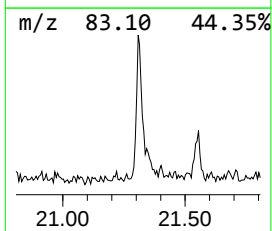
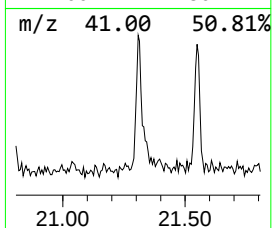
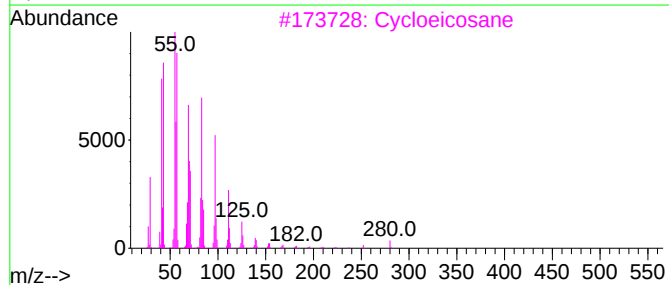
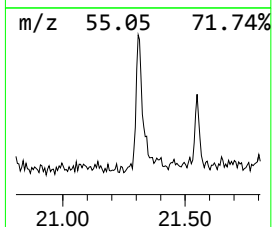
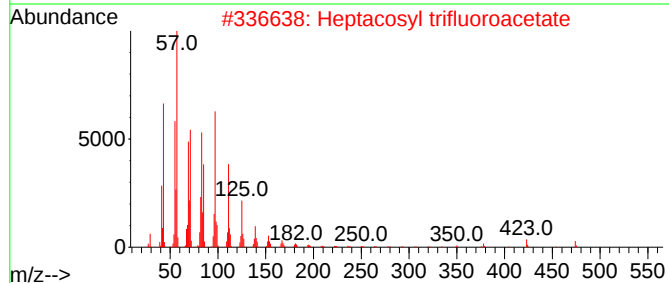
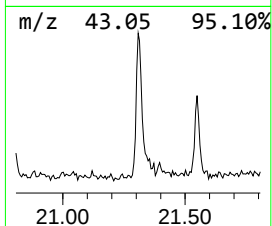
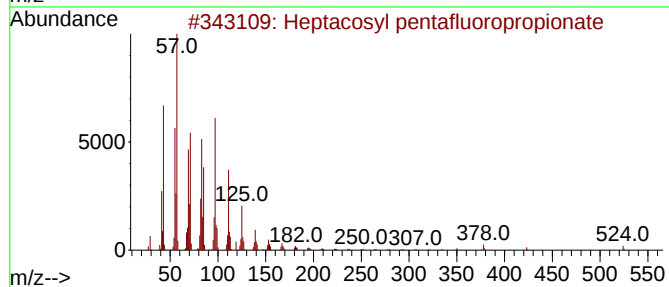
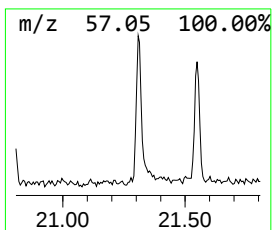
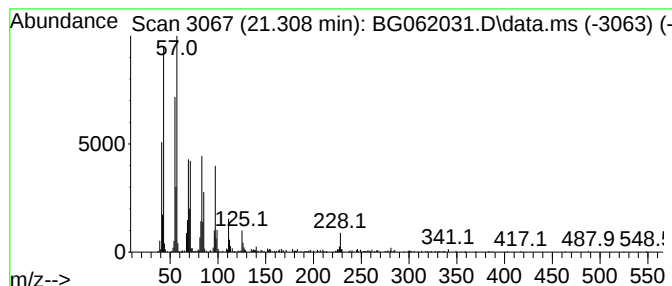
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Heptacosyl pentafluoropropi... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.308	8.19 ng/ul	807361	Chrysene-d12	21.672

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosyl pentafluoropropionate	542	C30H55F5O2	1000351-81-6	91
2			Heptacosyl trifluoroacetate	492	C29H55F3O2	1000351-75-8	91
3			Cycloeicosane	280	C20H40	000296-56-0	91
4			Triacontyl heptafluorobutyrate	634	C34H61F7O2	1000351-83-2	91
5			Octacosyl heptafluorobutyrate	606	C32H57F7O2	1010351-83-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071124\
Data File : BG062031.D
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Sample : P3124-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

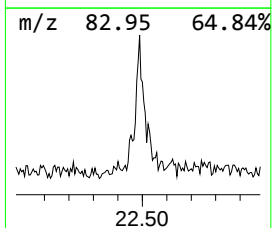
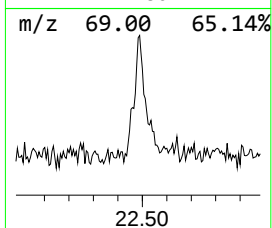
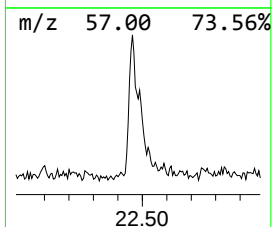
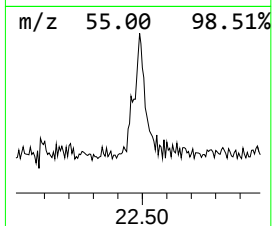
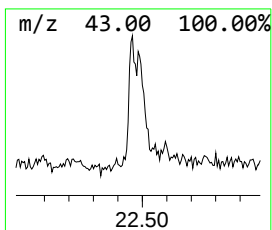
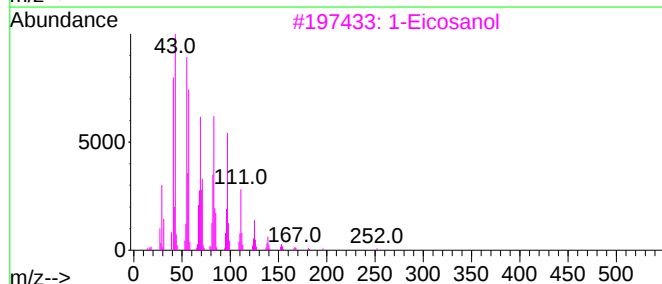
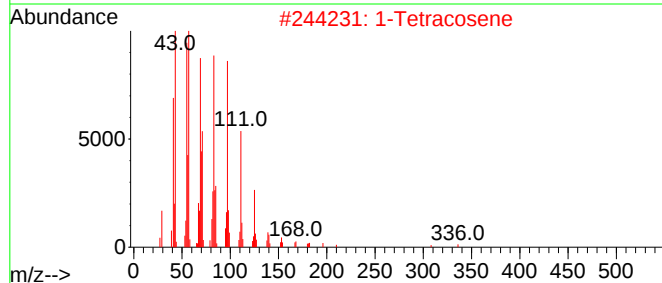
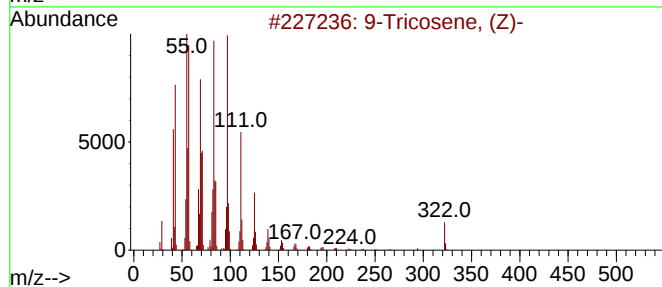
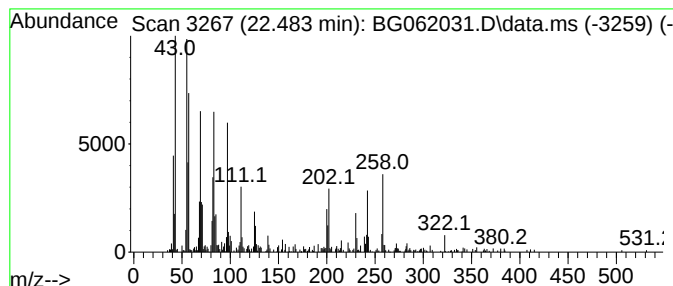
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.483	8.43 ng/ul	830909	Chrysene-d12	21.672	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Tricosene, (Z)-	322	C23H46	027519-02-4	64
2	1-Tetracosene	336	C24H48	010192-32-2	64
3	1-Eicosanol	298	C20H42O	000629-96-9	53
4	9-Tricosene, (Z)-	322	C23H46	027519-02-4	53
5	Cyclopentadecane	210	C15H30	000295-48-7	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071124\
Data File : BG062031.D
Acq On : 12 Jul 2024 6:40
Operator : MA/JU
Sample : P3124-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4CB3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.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
unknown-01	3.711	8.0	ng/ul	166796	1	8.041	417711	20.0
Methacrylamide	3.969	3.9	ng/ul	81130	1	8.041	417711	20.0
1-Phenoxypropan...	11.584	8.7	ng/ul	270822	2	10.849	620750	20.0
n-Hexadecanoic ...	18.288	8.2	ng/ul	429722	4	17.412	1046940	20.0
Phenanthrene, 4...	18.335	3.5	ng/ul	182521	4	17.412	1046940	20.0
4H-Cyclopenta[d...	18.482	4.5	ng/ul	234798	4	17.412	1046940	20.0
9,10-Anthracene...	18.822	2.9	ng/ul	149214	4	17.412	1046940	20.0
11H-Benzo[b]flu...	20.309	3.0	ng/ul	299304	5	21.672	1971870	20.0
Heptacosyl pent...	21.308	8.2	ng/ul	807361	5	21.672	1971870	20.0
(DEL) Alkane: S...	22.483	8.4	ng/ul	830909	5	21.672	1971870	20.0