

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
 Data File : BG058222.D
 Acq On : 14 Jul 2023 8:03
 Operator : CG/JU
 Sample : 03418-10
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4663

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-BG070123.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG058222.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.346	41	44	46	rBV3	12437	15269	0.44%	0.025%
2	3.381	46	50	58	rVB	61405	90720	2.63%	0.151%
3	3.804	119	122	127	rVB	13857	19699	0.57%	0.033%
4	4.092	167	171	177	rVB2	7813	12702	0.37%	0.021%
5	4.415	219	226	245	rBV	1048873	1683933	48.74%	2.800%
6	5.009	319	327	343	rBV	439876	680364	19.69%	1.131%
7	6.084	504	510	518	rVB2	8645	13596	0.39%	0.023%
8	7.042	665	673	676	rBV	341632	659847	19.10%	1.097%
9	7.095	680	682	698	rVV	376121	554017	16.03%	0.921%
10	7.230	698	705	718	rVB	207115	354288	10.25%	0.589%
11	7.435	733	740	743	rBV	250687	466914	13.51%	0.776%
12	7.465	743	745	759	rVV	223591	339634	9.83%	0.565%
13	7.905	813	820	831	rBV	193304	348197	10.08%	0.579%
14	8.611	933	940	950	rBV	316799	554162	16.04%	0.921%
15	8.705	950	956	965	rVB	258688	433331	12.54%	0.721%
16	9.063	1008	1017	1021	rBV	254292	469249	13.58%	0.780%
17	9.104	1021	1024	1038	rVV	189260	330534	9.57%	0.550%
18	9.216	1040	1043	1047	rVB4	2995	4806	0.14%	0.008%
19	9.786	1132	1140	1143	rBV	210688	409901	11.86%	0.682%
20	9.815	1143	1145	1156	rVB	191401	283535	8.21%	0.471%
21	10.326	1224	1232	1234	rBV	342675	605916	17.54%	1.008%
22	10.702	1285	1296	1301	rBV	303356	586891	16.99%	0.976%
23	10.755	1301	1305	1313	rVV	284269	489248	14.16%	0.814%
24	10.837	1313	1319	1337	rVB	269716	518849	15.02%	0.863%
25	11.037	1345	1353	1361	rVB	280347	485979	14.07%	0.808%
26	11.989	1507	1515	1533	rBV	558979	974647	28.21%	1.621%
27	12.224	1551	1555	1562	rVB3	11613	17133	0.50%	0.028%
28	12.289	1562	1566	1572	rVB3	4690	7724	0.22%	0.013%
29	12.582	1608	1616	1625	rBV	488063	799699	23.14%	1.330%
30	12.729	1634	1641	1649	rBV	322051	533368	15.44%	0.887%
31	12.970	1675	1682	1687	rBV	518528	859893	24.89%	1.430%
32	13.017	1687	1690	1699	rVB	85962	125475	3.63%	0.209%
33	13.141	1706	1711	1716	rVB5	3079	5205	0.15%	0.009%
34	13.370	1743	1750	1754	rBV	639342	1060857	30.70%	1.764%
35	13.417	1754	1758	1767	rVB	337826	541534	15.67%	0.900%
36	13.945	1837	1848	1852	rBV	667898	989302	28.63%	1.645%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
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37	13.992	1852	1856	1863	rVV	389324	581824	16.84%	0.967%
38	14.063	1864	1868	1871	rVV4	3520	4742	0.14%	0.008%
39	14.116	1871	1877	1884	rVV	387509	571564	16.54%	0.950%
40	14.233	1890	1897	1907	rVB	712418	1133209	32.80%	1.884%
41	14.427	1922	1930	1936	rVB3	2990	4855	0.14%	0.008%
42	14.539	1941	1949	1955	rBV	520668	812331	23.51%	1.351%
43	14.604	1955	1960	1969	rVB	599873	930082	26.92%	1.547%
44	14.750	1977	1985	2003	rBV3	385261	892976	25.84%	1.485%
45	14.938	2010	2017	2031	rVB	448053	715987	20.72%	1.191%
46	15.162	2050	2055	2069	rBV	398180	635127	18.38%	1.056%
47	15.320	2075	2082	2086	rBV5	4846	9540	0.28%	0.016%
48	15.379	2086	2092	2096	rVB2	21944	30912	0.89%	0.051%
49	15.532	2111	2118	2122	rBV	982047	1507558	43.63%	2.507%
50	15.579	2122	2126	2132	rVV	424256	608356	17.61%	1.012%
51	15.667	2132	2141	2151	rVV3	584695	1130547	32.72%	1.880%
52	15.896	2174	2180	2194	rVB	247138	374370	10.83%	0.622%
53	16.260	2240	2242	2248	rVV3	5950	8347	0.24%	0.014%
54	16.454	2270	2275	2282	rBV	76189	108303	3.13%	0.180%
55	16.589	2291	2298	2305	rBV	514932	763773	22.10%	1.270%
56	16.819	2333	2337	2341	rVB4	6955	9429	0.27%	0.016%
57	16.936	2350	2357	2369	rBV	608933	982416	28.43%	1.634%
58	17.030	2369	2373	2376	rVB	9254	13206	0.38%	0.022%
59	17.112	2385	2387	2392	rVB2	4551	5301	0.15%	0.009%
60	17.289	2410	2417	2420	rBV	636158	1033669	29.92%	1.719%
61	17.330	2420	2424	2428	rVV	673021	1016045	29.41%	1.689%
62	17.383	2428	2433	2437	rVV2	1074840	1842564	53.33%	3.064%
63	17.418	2437	2439	2453	rVB	771709	1020288	29.53%	1.697%
64	17.559	2460	2463	2470	rVB7	2567	4796	0.14%	0.008%
65	17.688	2478	2485	2491	rBV	626461	982217	28.43%	1.633%
66	17.759	2491	2497	2510	rVB	648317	951628	27.54%	1.582%
67	17.894	2517	2520	2523	rBV3	5418	6744	0.20%	0.011%
68	17.929	2523	2526	2531	rVB5	11523	15560	0.45%	0.026%
69	18.170	2562	2567	2573	rVV	195417	279882	8.10%	0.465%
70	18.240	2573	2579	2588	rVB	698093	936473	27.10%	1.557%
71	18.464	2613	2617	2621	rBV5	5958	10950	0.32%	0.018%
72	18.699	2653	2657	2662	rVB	15175	22176	0.64%	0.037%
73	19.039	2708	2715	2717	rBV5	5221	10903	0.32%	0.018%
74	19.092	2720	2724	2728	rVB5	5471	8862	0.26%	0.015%
75	19.239	2742	2749	2754	rBV	19288	31071	0.90%	0.052%
76	19.339	2754	2766	2778	rBV	1661859	2578518	74.63%	4.288%

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77	19.445	2779	2784	2793	rVV3	34428	60604	1.75%	0.101%
78	19.586	2803	2808	2814	rVB2	18950	30435	0.88%	0.051%
79	19.674	2816	2823	2826	rVV	1345606	2247913	65.06%	3.738%
80	19.703	2826	2828	2840	rVV	1274941	1659740	48.04%	2.760%
81	19.809	2844	2846	2851	rVV6	7422	11475	0.33%	0.019%
82	20.197	2907	2912	2917	rVB9	8990	16470	0.48%	0.027%
83	20.585	2972	2978	2983	rBV	712962	864882	25.03%	1.438%
84	20.690	2992	2996	2999	rBV6	11459	16408	0.47%	0.027%
85	21.184	3075	3080	3096	rBV	169973	386406	11.18%	0.643%
86	21.419	3114	3120	3127	rVB	993773	1300277	37.63%	2.162%
87	21.525	3130	3138	3144	rVV2	1515535	3455226	100.00%	5.745%
88	21.584	3144	3148	3163	rVB	1095981	1774412	51.35%	2.950%
89	21.719	3168	3171	3176	rVB6	12206	15109	0.44%	0.025%
90	22.306	3266	3271	3276	rBV4	23694	54134	1.57%	0.090%
91	22.671	3330	3333	3338	rVB4	18686	26019	0.75%	0.043%
92	22.970	3379	3384	3389	rVB5	25954	52950	1.53%	0.088%
93	23.681	3495	3505	3513	rBV	767262	1863125	53.92%	3.098%
94	23.757	3514	3518	3528	rVB3	39609	85869	2.49%	0.143%
95	24.468	3627	3639	3645	rBV2	766619	2249853	65.11%	3.741%
96	24.539	3645	3651	3663	rVV	625210	1650751	47.78%	2.745%
97	24.680	3665	3675	3690	rVB	453304	1307639	37.85%	2.174%
98	25.761	3852	3859	3860	rBV3	40844	63964	1.85%	0.106%
99	27.071	4077	4082	4084	rBV3	22252	39021	1.13%	0.065%
100	28.287	4267	4289	4290	rBV	484300	1994003	57.71%	3.316%

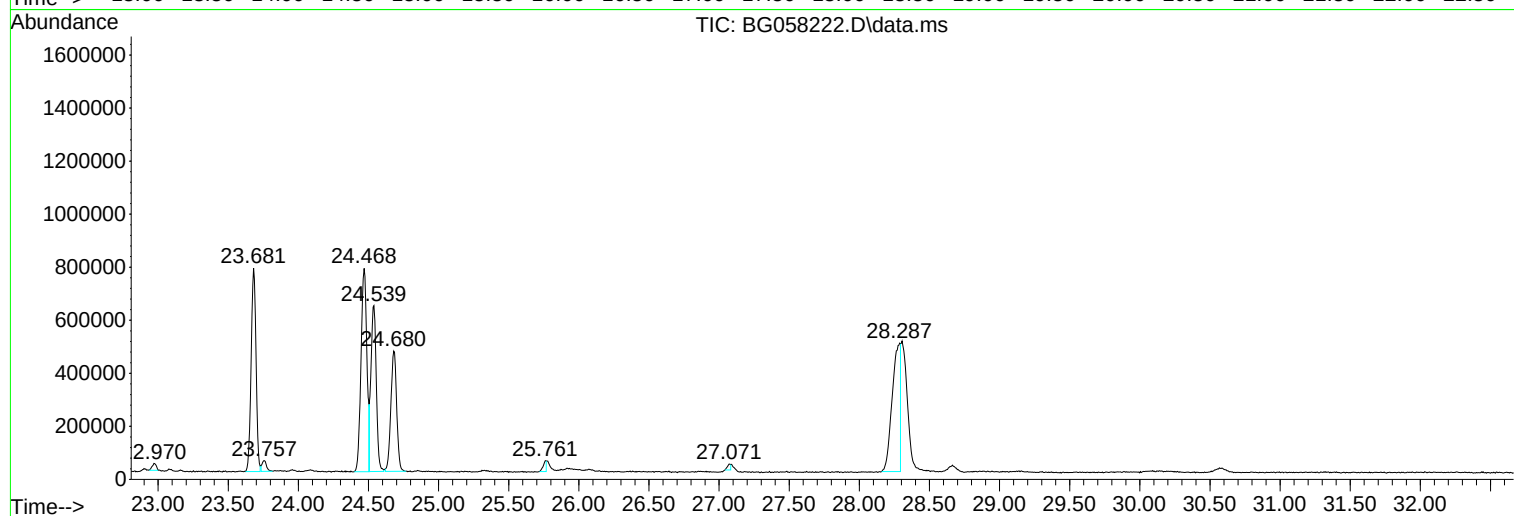
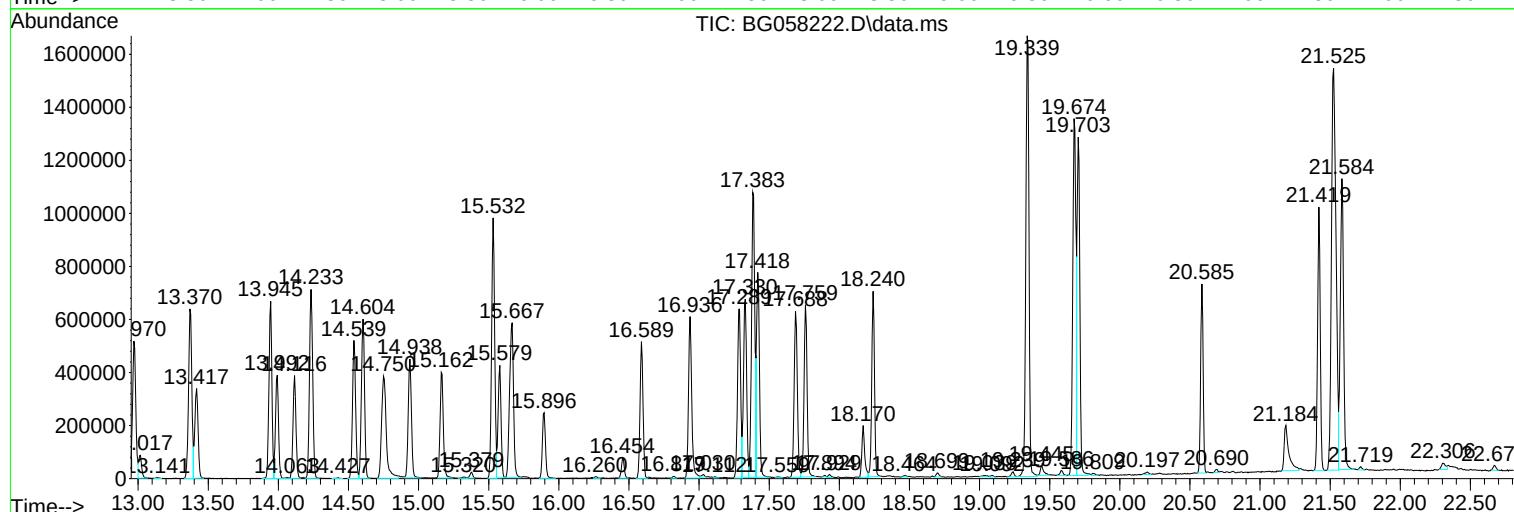
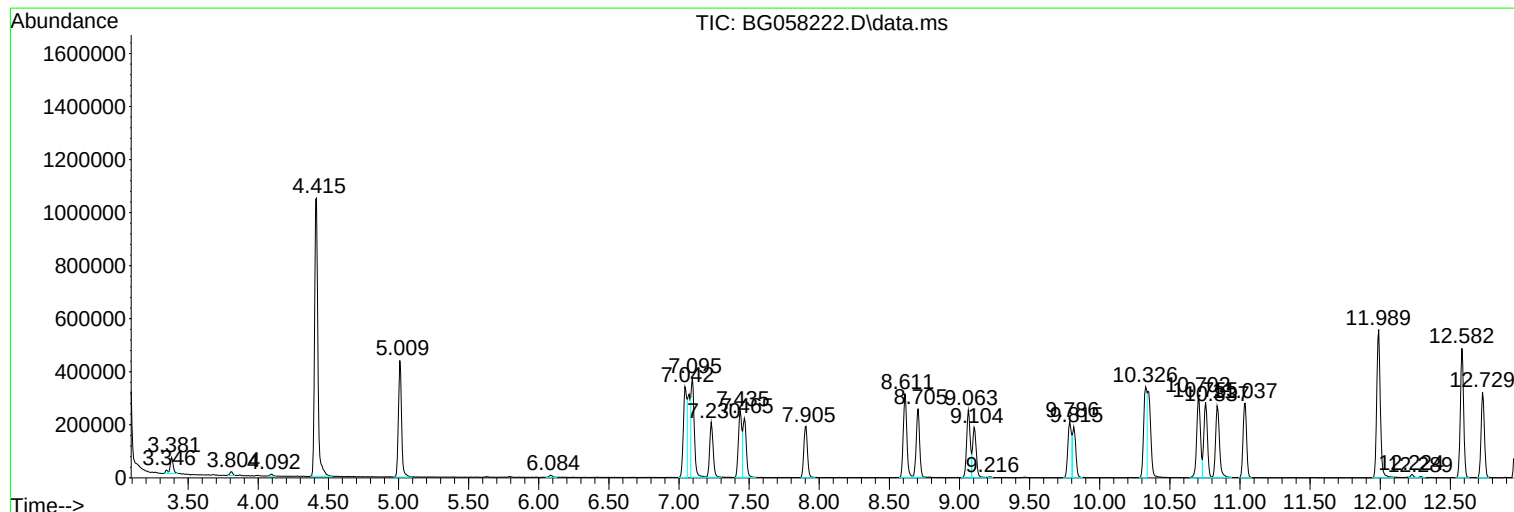
Sum of corrected areas: 60140204

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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



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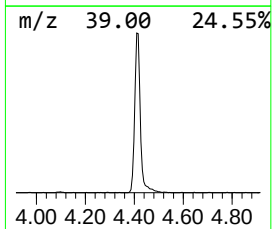
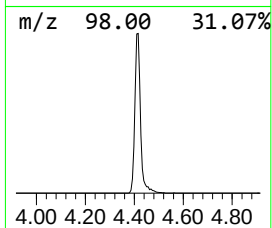
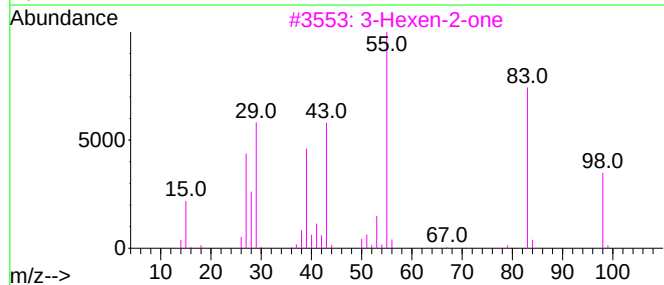
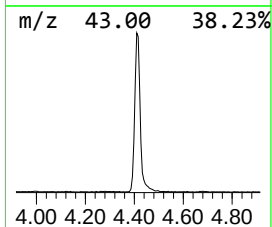
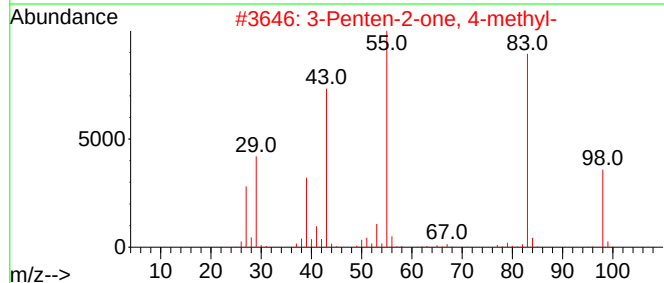
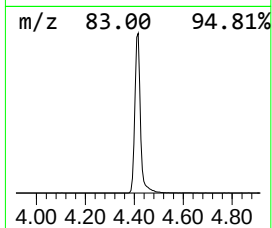
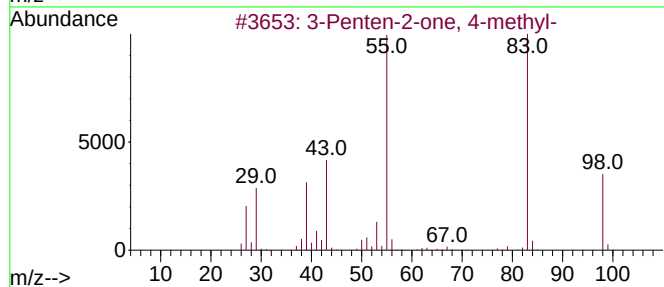
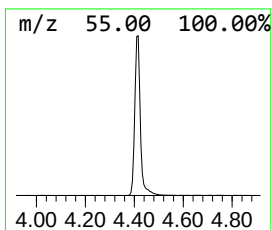
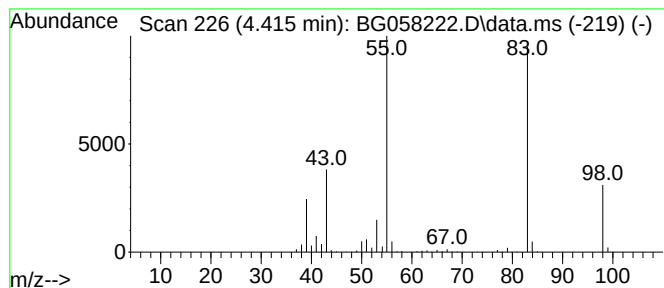
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.415	96.72 ng/ul	1683930	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Penten-2-one, 4-methyl-			98	C6H10O	000141-79-7	94
2	3-Penten-2-one, 4-methyl-			98	C6H10O	000141-79-7	91
3	3-Hexen-2-one			98	C6H10O	000763-93-9	91
4	3-Penten-2-one, 4-methyl-			98	C6H10O	000141-79-7	91
5	3-Hexen-2-one			98	C6H10O	000763-93-9	91



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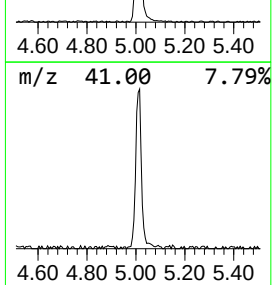
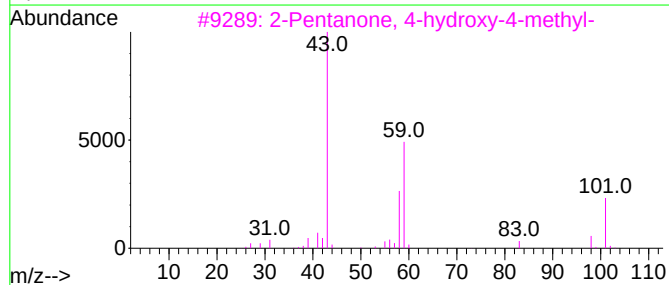
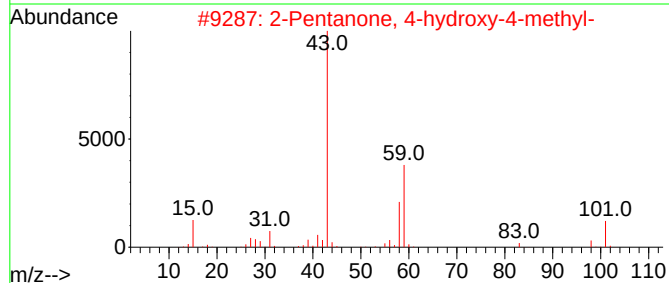
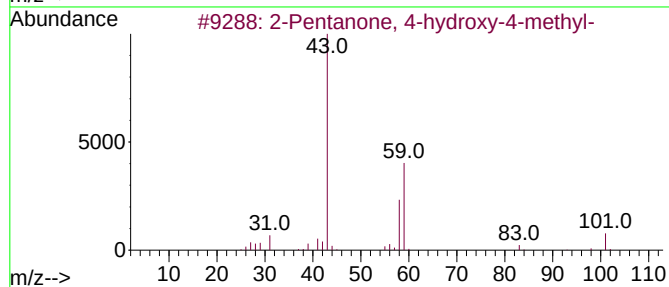
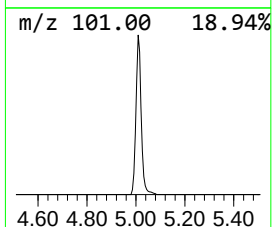
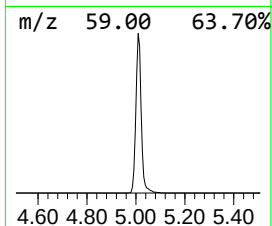
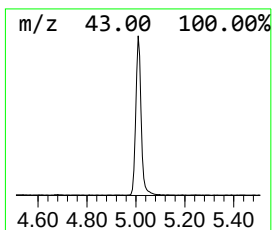
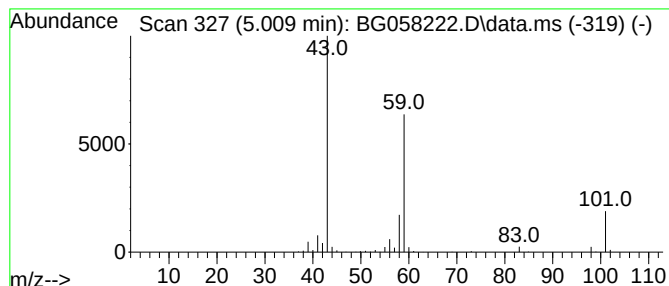
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.009	39.08 ng/ul	680364	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
3			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
4			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	37
5			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33



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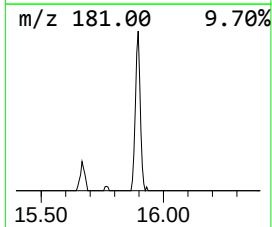
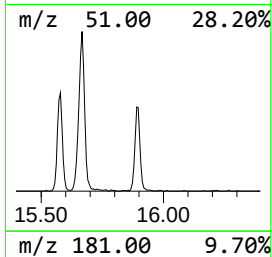
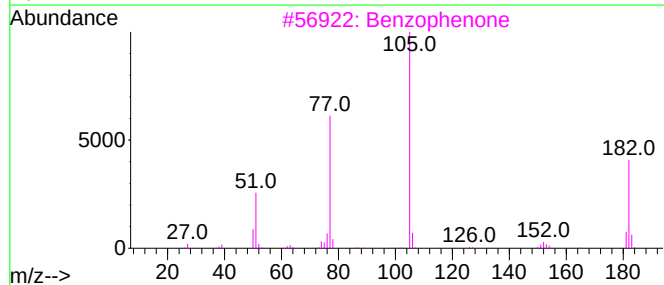
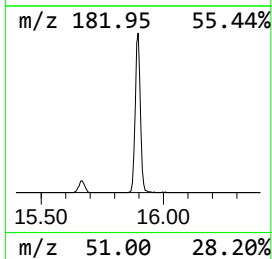
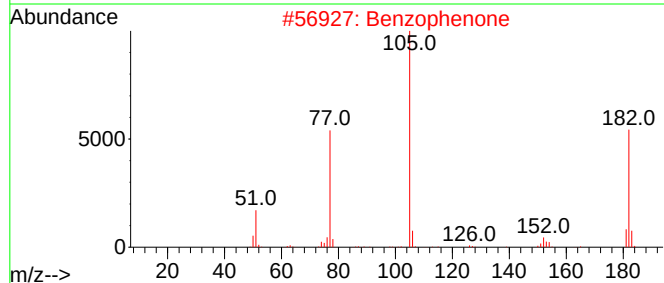
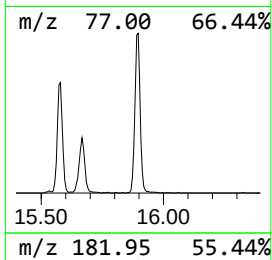
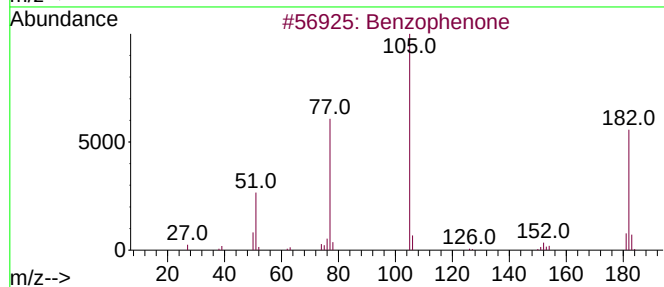
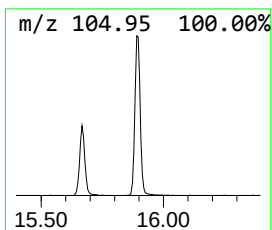
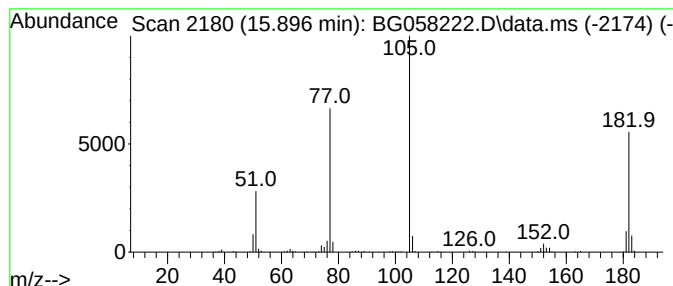
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.896	9.22 ng/ul	374370	Acenaphthene-d10	14.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	93
5			Azobenzene	182	C12H10N2	000103-33-3	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
Data File : BG058222.D
Acq On : 14 Jul 2023 8:03
Operator : CG/JU
Sample : 03418-10
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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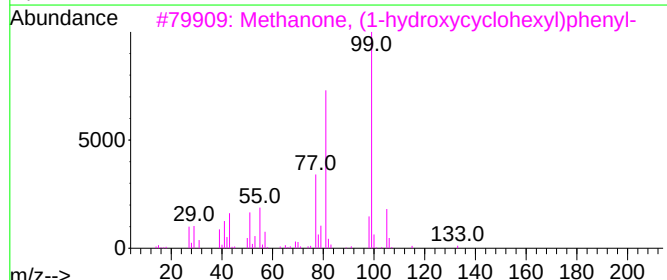
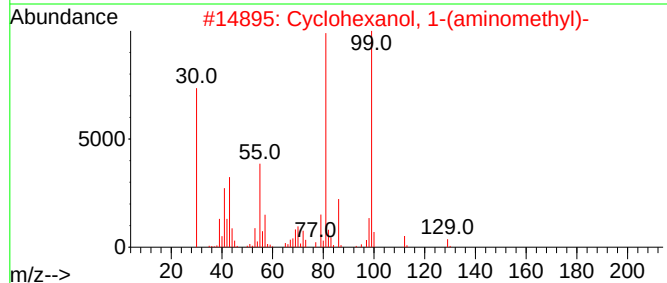
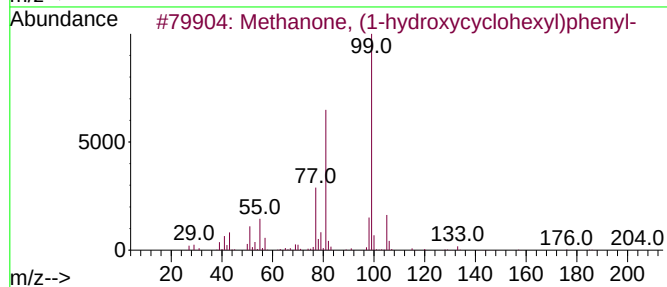
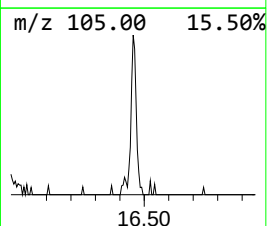
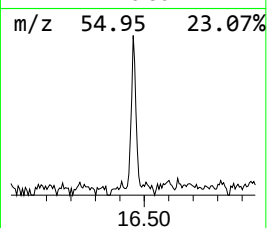
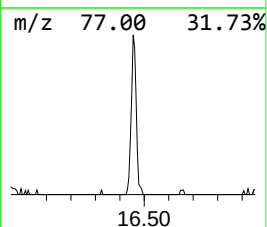
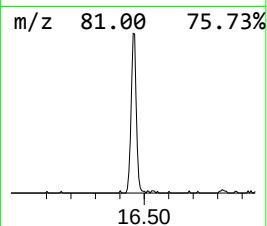
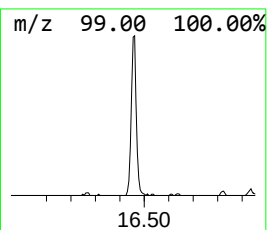
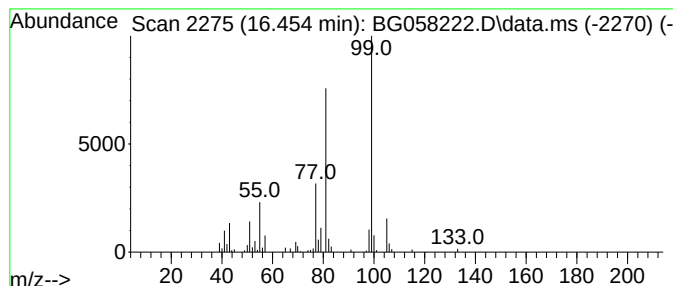
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Methanone, (1-hydroxycyclohexyl) Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.454	2.10 ng/ul	108303	Phenanthrene-d10	17.289

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methanone, (1-hydroxycyclohexyl)...	204	C13H16O2	000947-19-3	90
2			Cyclohexanol, 1-(aminomethyl)-	129	C7H15NO	004000-72-0	59
3			Methanone, (1-hydroxycyclohexyl)...	204	C13H16O2	000947-19-3	58
4			2-Methylcyclosarin	194	C8H16FO2P	085473-32-1	53
5			Cyclohexanol, 1-(aminomethyl)-	129	C7H15NO	004000-72-0	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
Data File : BG058222.D
Acq On : 14 Jul 2023 8:03
Operator : CG/JU
Sample : 03418-10
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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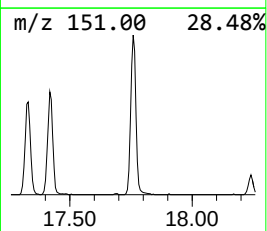
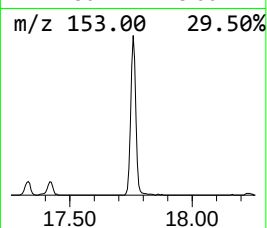
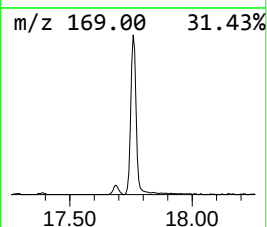
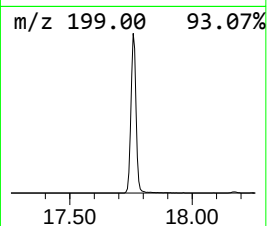
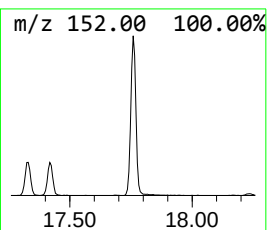
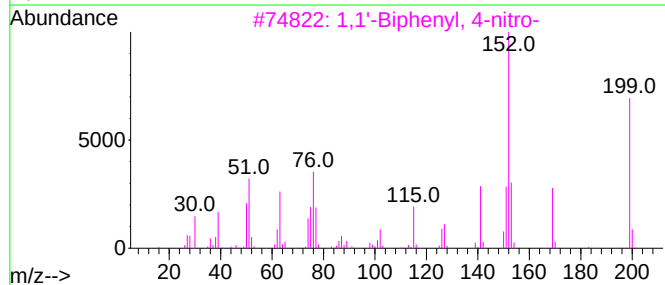
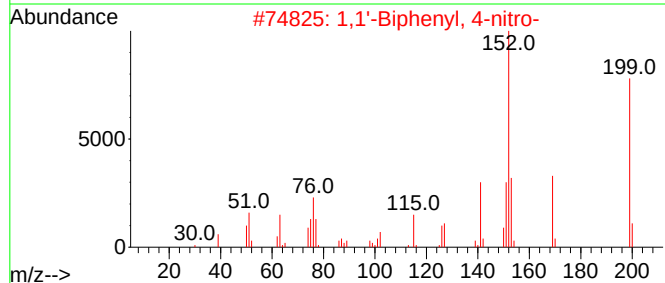
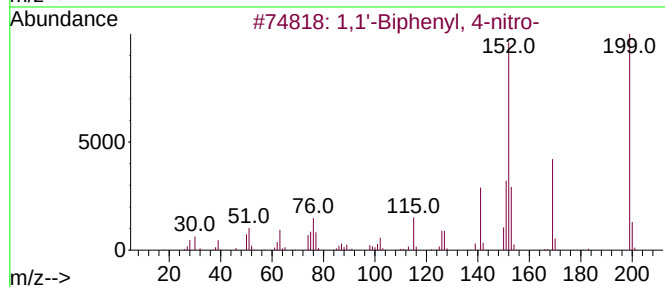
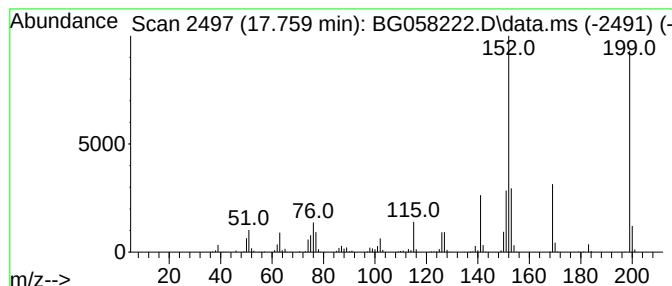
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,1'-Biphenyl, 4-nitro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.759	18.41 ng/ul	951628	Phenanthrene-d10	17.289

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 4-nitro-			199	C12H9NO2	000092-93-3	99
2	1,1'-Biphenyl, 4-nitro-			199	C12H9NO2	000092-93-3	98
3	1,1'-Biphenyl, 4-nitro-			199	C12H9NO2	000092-93-3	97
4	1,1'-Biphenyl, 3-nitro-			199	C12H9NO2	002113-58-8	90
5	1,1'-Biphenyl, 3-nitro-			199	C12H9NO2	002113-58-8	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
Data File : BG058222.D
Acq On : 14 Jul 2023 8:03
Operator : CG/JU
Sample : 03418-10
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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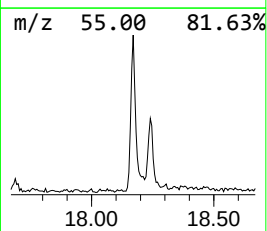
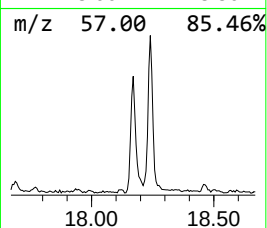
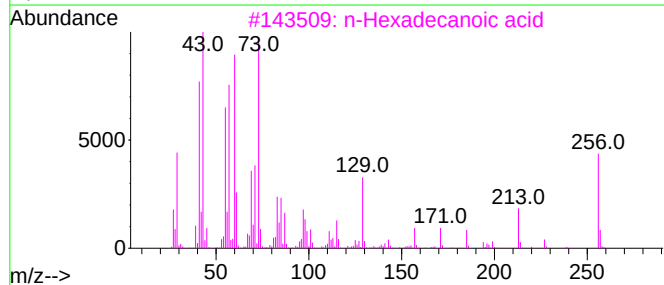
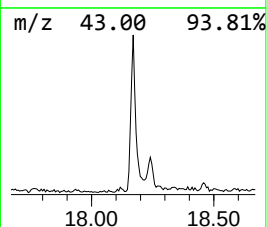
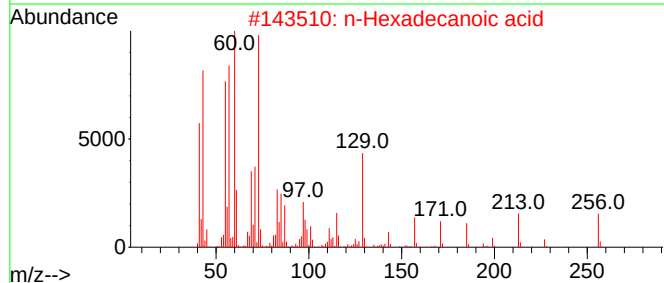
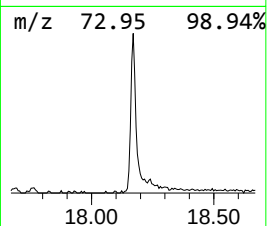
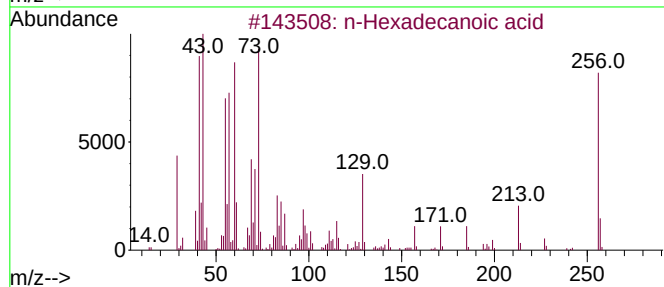
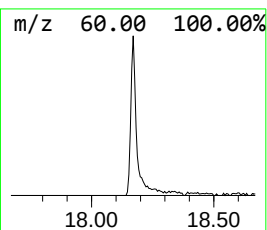
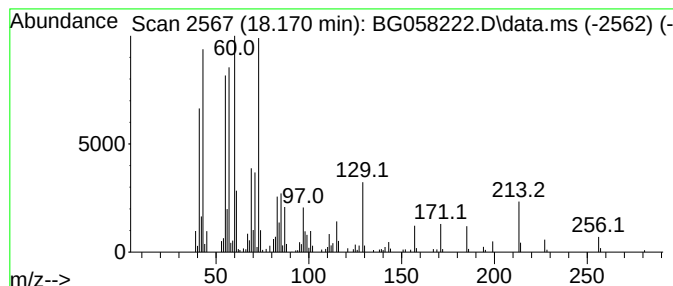
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.170	5.42 ng/ul	279882	Phenanthrene-d10	17.289

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
4			Tridecanoic acid	214	C13H26O2	000638-53-9	90
5			Tridecanoic acid	214	C13H26O2	000638-53-9	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
Data File : BG058222.D
Acq On : 14 Jul 2023 8:03
Operator : CG/JU
Sample : 03418-10
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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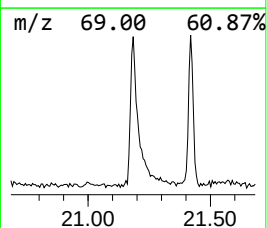
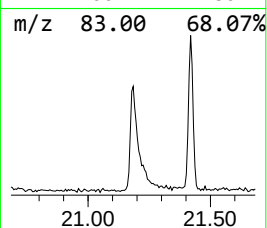
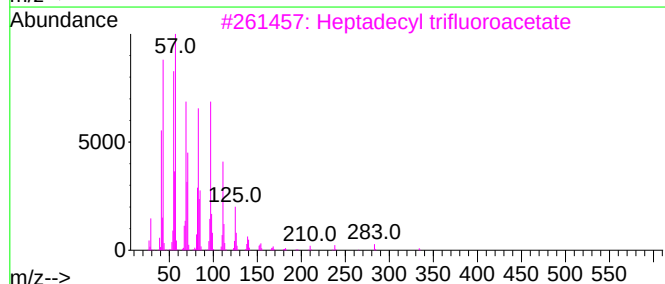
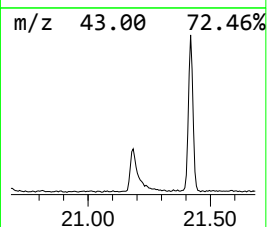
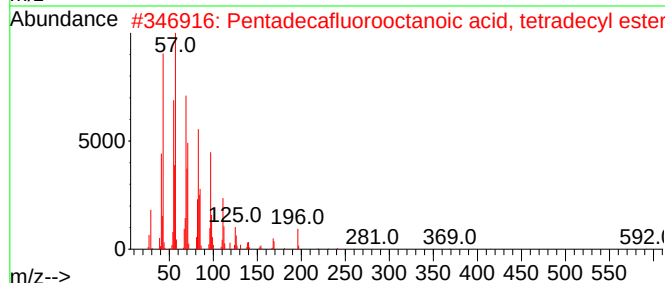
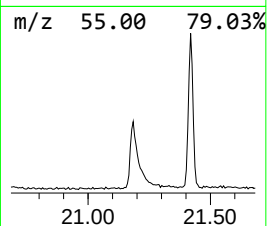
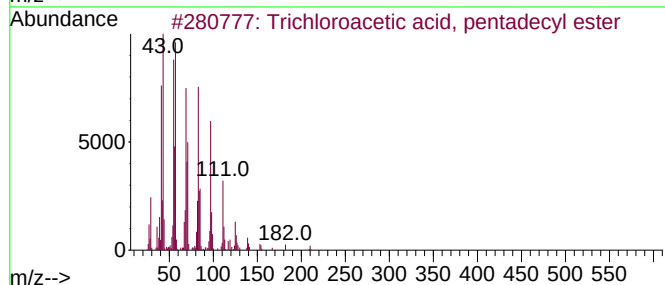
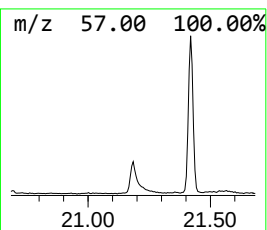
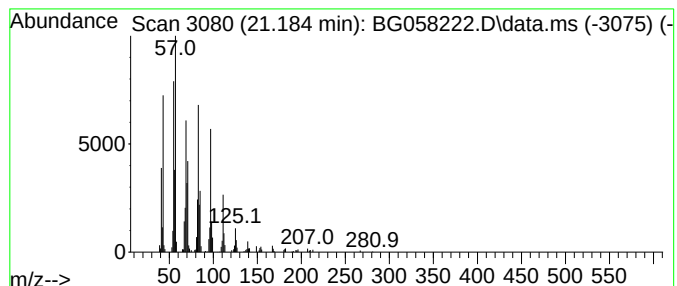
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Trichloroacetic acid, penta... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.184	2.24 ng/ul	386406	Chrysene-d12	21.542

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	94
2			Pentadecafluorooctanoic acid, te...	610	C22H29F15O2	1000406-04-4	93
3			Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	91
4			Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	91
5			1-Hexacosene	364	C26H52	018835-33-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
Data File : BG058222.D
Acq On : 14 Jul 2023 8:03
Operator : CG/JU
Sample : 03418-10
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Penten-2-one,...	4.415	96.7	ng/ul	1683930	1	7.905	348197	20.0
2-Pentanone, 4-...	5.009	39.1	ng/ul	680364	1	7.905	348197	20.0
Benzophenone	15.896	9.2	ng/ul	374370	3	14.539	812331	20.0
Methanone, (1-h...	16.454	2.1	ng/ul	108303	4	17.289	1033670	20.0
1,1'-Biphenyl, ...	17.759	18.4	ng/ul	951628	4	17.289	1033670	20.0
n-Hexadecanoic ...	18.170	5.4	ng/ul	279882	4	17.289	1033670	20.0
Trichloroacetic...	21.184	2.2	ng/ul	386406	5	21.542	3455230	20.0