

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
 Data File : BG056119.D
 Acq On : 25 Dec 2022 12:04
 Operator : CG/JU
 Sample : N6017-18
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBYG4

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-BG122422.M
 Title : SVOA CALIBRATION

Signal : TIC: BG056119.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.371	10	14	23	rBV	97914	144282	8.97%	0.901%
2	3.429	23	24	29	rVB3	3996	3490	0.22%	0.022%
3	3.647	58	61	68	rVB2	7873	10077	0.63%	0.063%
4	4.082	132	135	147	rBV	63765	104212	6.48%	0.651%
5	4.334	175	178	183	rBV2	1045	1967	0.12%	0.012%
6	4.434	191	195	200	rBV	3244	5434	0.34%	0.034%
7	4.663	230	234	237	rVB2	2712	3017	0.19%	0.019%
8	5.380	353	356	363	rVB2	2070	2849	0.18%	0.018%
9	7.472	702	712	726	rBV	175866	306807	19.07%	1.916%
10	7.648	735	742	750	rBV2	101829	173603	10.79%	1.084%
11	7.871	772	780	790	rVB	136311	236084	14.67%	1.475%
12	8.347	854	861	871	rVB	118739	194465	12.09%	1.215%
13	9.035	971	978	986	rBV	199130	347252	21.58%	2.169%
14	9.516	1052	1060	1068	rVB	129482	231925	14.42%	1.449%
15	9.904	1121	1126	1130	rBB3	6385	9119	0.57%	0.057%
16	10.245	1176	1184	1192	rVB	125453	216239	13.44%	1.351%
17	10.345	1196	1201	1209	rBB2	2797	5491	0.34%	0.034%
18	10.786	1268	1276	1285	rVB	244101	425023	26.42%	2.655%
19	11.179	1335	1343	1352	rVB	167032	301940	18.77%	1.886%
20	11.303	1355	1364	1373	rBV	163947	307702	19.12%	1.922%
21	12.730	1601	1607	1613	rBV2	4285	7004	0.44%	0.044%
22	12.807	1613	1620	1628	rVV2	97000	156380	9.72%	0.977%
23	13.759	1777	1782	1785	rVB2	3615	5223	0.32%	0.033%
24	14.234	1858	1863	1866	rBV3	7186	9404	0.58%	0.059%
25	14.276	1866	1870	1875	rVV4	2812	4699	0.29%	0.029%
26	14.340	1875	1881	1887	rVV	486493	710965	44.19%	4.441%
27	14.405	1887	1892	1894	rVV	1860	2374	0.15%	0.015%
28	14.452	1896	1900	1903	rVV	1124	1872	0.12%	0.012%
29	14.481	1903	1905	1909	rVB	1786	1909	0.12%	0.012%
30	14.652	1927	1934	1941	rVB	551542	873765	54.31%	5.458%
31	14.951	1976	1985	1993	rBV	348876	548655	34.10%	3.427%
32	15.116	2005	2013	2023	rBV	180490	295252	18.35%	1.844%
33	15.204	2023	2028	2031	rVB4	4345	7326	0.46%	0.046%
34	15.316	2043	2047	2049	rBV	1642	1966	0.12%	0.012%
35	15.415	2061	2064	2068	rBV	1482	1960	0.12%	0.012%
36	15.938	2146	2153	2160	rVB	780180	1193819	74.20%	7.457%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
 Data File : BG056119.D
 Acq On : 25 Dec 2022 12:04
 Operator : CG/JU
 Sample : N6017-18
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBYG4

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-BG122422.M
 Title : SVOA CALIBRATION

37	16.038	2163	2170	2176	rBV	212534	300067	18.65%	1.874%
38	16.179	2188	2194	2197	rVB	3585	5509	0.34%	0.034%
39	16.226	2197	2202	2206	rBV3	9868	14267	0.89%	0.089%
40	16.285	2206	2212	2217	rBV2	39898	56624	3.52%	0.354%
41	16.432	2230	2237	2239	rBV2	1461	2918	0.18%	0.018%
42	16.508	2245	2250	2252	rVB	1750	2083	0.13%	0.013%
43	16.714	2281	2285	2288	rBV2	2497	2897	0.18%	0.018%
44	16.955	2324	2326	2328	rBV	2128	2612	0.16%	0.016%
45	17.002	2331	2334	2337	rVV	2076	2260	0.14%	0.014%
46	17.072	2345	2346	2351	rVB3	2222	2563	0.16%	0.016%
47	17.331	2389	2390	2393	rBV	2374	2413	0.15%	0.015%
48	17.366	2393	2396	2399	rVV	2794	3245	0.20%	0.020%
49	17.419	2401	2405	2407	rVV	2399	2483	0.15%	0.016%
50	17.472	2410	2414	2418	rVV2	2271	3264	0.20%	0.020%
51	17.536	2422	2425	2429	rVV2	2088	2714	0.17%	0.017%
52	17.619	2436	2439	2443	rBV4	2408	2496	0.16%	0.016%
53	17.689	2443	2451	2457	rBV	479186	733782	45.61%	4.584%
54	17.736	2457	2459	2461	rVV3	4050	3914	0.24%	0.024%
55	17.789	2461	2468	2474	rVV	870963	1295267	80.51%	8.091%
56	18.289	2550	2553	2555	rBV2	1861	2197	0.14%	0.014%
57	18.330	2555	2560	2563	rBV3	5234	6956	0.43%	0.043%
58	18.400	2569	2572	2574	rBV2	2378	3546	0.22%	0.022%
59	18.424	2574	2576	2579	rVB	2670	2921	0.18%	0.018%
60	18.477	2582	2585	2590	rBV4	4833	7456	0.46%	0.047%
61	18.535	2590	2595	2604	rVV	133323	186486	11.59%	1.165%
62	18.606	2606	2607	2611	rVV4	4453	5499	0.34%	0.034%
63	18.735	2626	2629	2632	rBV4	3673	4753	0.30%	0.030%
64	18.829	2640	2645	2648	rVV5	3356	6174	0.38%	0.039%
65	18.900	2652	2657	2660	rVV6	3052	5911	0.37%	0.037%
66	18.952	2662	2666	2671	rVV4	12923	15074	0.94%	0.094%
67	19.052	2679	2683	2686	rBV5	3086	4487	0.28%	0.028%
68	19.123	2692	2695	2696	rBV2	2283	2157	0.13%	0.013%
69	19.176	2703	2704	2707	rVB3	4039	2308	0.14%	0.014%
70	19.246	2713	2716	2717	rBV3	1920	1950	0.12%	0.012%
71	19.264	2717	2719	2722	rVB4	3047	2098	0.13%	0.013%
72	19.452	2748	2751	2753	rBV4	4846	6013	0.37%	0.038%
73	19.516	2760	2762	2764	rBV2	3402	3097	0.19%	0.019%
74	19.581	2769	2773	2774	rBV4	3092	4219	0.26%	0.026%
75	19.622	2774	2780	2782	rBV2	17993	29490	1.83%	0.184%
76	19.675	2782	2789	2795	rBV6	74318	167856	10.43%	1.048%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
 Data File : BG056119.D
 Acq On : 25 Dec 2022 12:04
 Operator : CG/JU
 Sample : N6017-18
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBYG4

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-BG122422.M
 Title : SVOA CALIBRATION

77	19.793	2805	2809	2812	rBV3	65768	106574	6.62%	0.666%
78	20.051	2847	2853	2863	rVB	1097655	1577916	98.07%	9.856%
79	20.180	2872	2875	2881	rBV2	32689	44025	2.74%	0.275%
80	20.233	2882	2884	2888	rVB4	5063	6232	0.39%	0.039%
81	20.545	2931	2937	2941	rBV8	13771	25986	1.62%	0.162%
82	20.615	2948	2949	2950	rBV	7414	3964	0.25%	0.025%
83	20.903	2994	2998	3002	rBV3	41911	61754	3.84%	0.386%
84	20.938	3002	3004	3007	rVB4	5516	5648	0.35%	0.035%
85	21.538	3104	3106	3107	rBV2	5637	4218	0.26%	0.026%
86	21.579	3108	3113	3123	rVV4	73860	145777	9.06%	0.911%
87	21.884	3162	3165	3166	rBV3	8446	9670	0.60%	0.060%
88	21.984	3176	3182	3189	rVB2	465215	814479	50.62%	5.088%
89	22.454	3261	3262	3263	rBV	6604	2438	0.15%	0.015%
90	22.807	3320	3322	3328	rVB6	11051	14099	0.88%	0.088%
91	24.428	3592	3598	3606	rVB3	106578	239161	14.86%	1.494%
92	25.222	3721	3733	3744	rBV3	525552	1608911	100.00%	10.050%
93	25.457	3762	3773	3786	rVB2	283669	879637	54.67%	5.495%
94	26.003	3864	3866	3868	rBV3	6989	8372	0.52%	0.052%
95	26.696	3974	3984	3997	rBV3	168039	532290	33.08%	3.325%
96	29.029	4379	4381	4382	rBV2	9316	6782	0.42%	0.042%
97	29.975	4530	4542	4544	rBV7	45660	131288	8.16%	0.820%
98	30.674	4659	4661	4662	rBV2	5157	3045	0.19%	0.019%
99	31.044	4720	4724	4725	rBV4	6684	7233	0.45%	0.045%
100	31.450	4791	4793	4795	rBV3	7781	4088	0.25%	0.026%

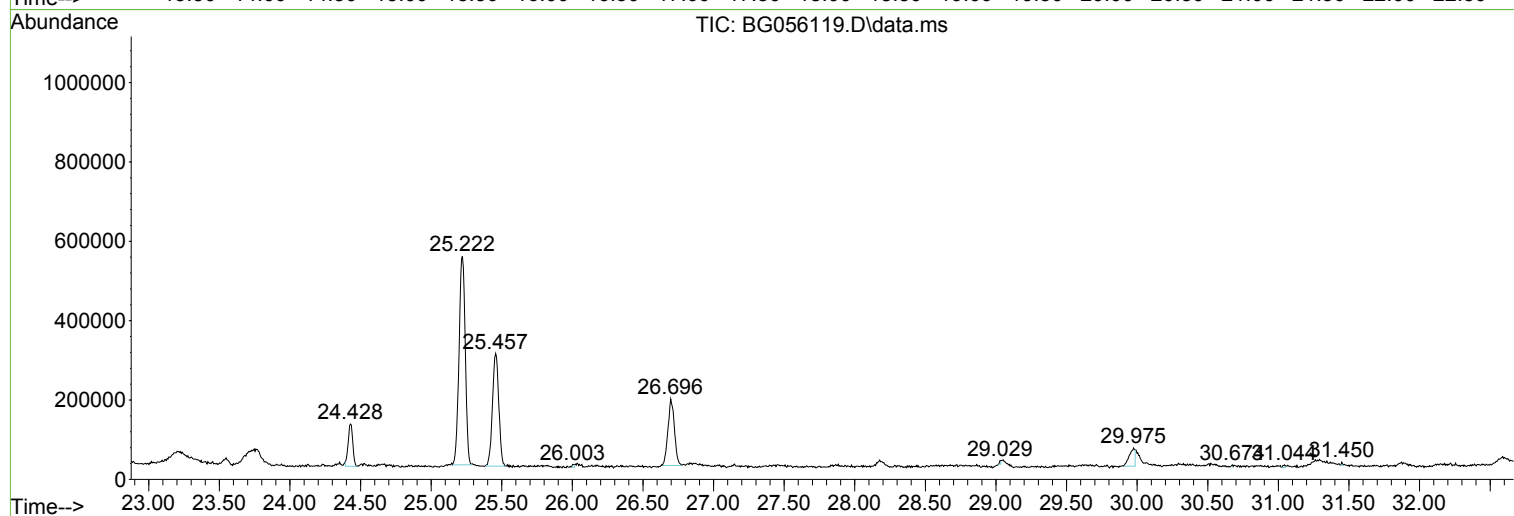
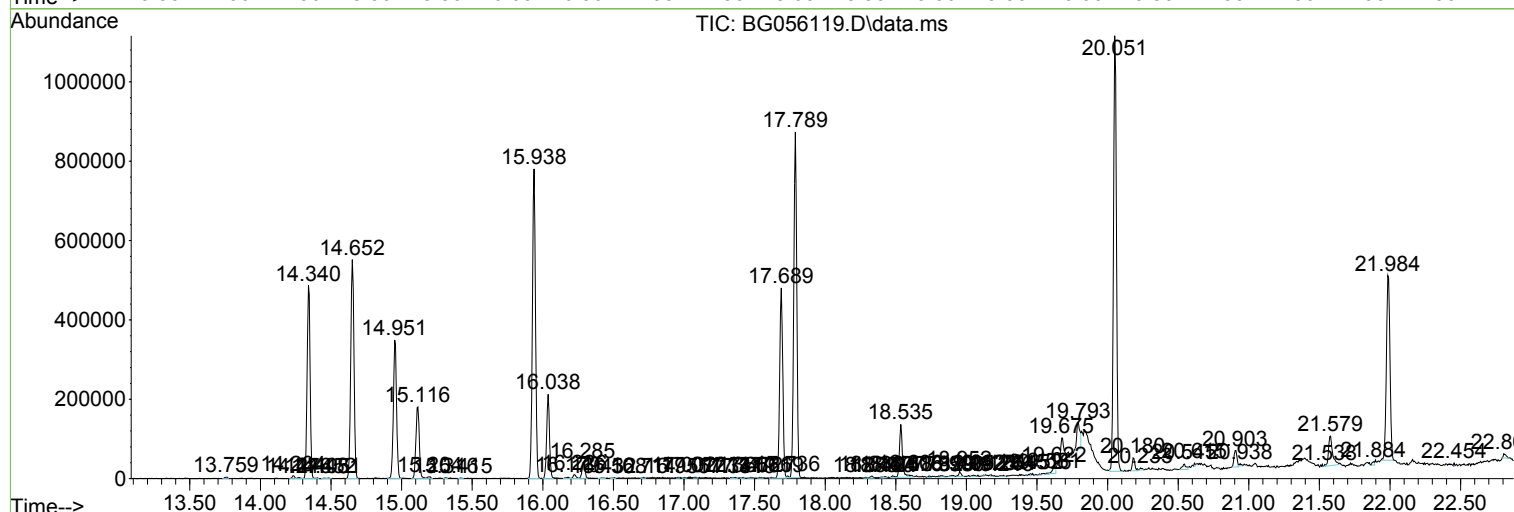
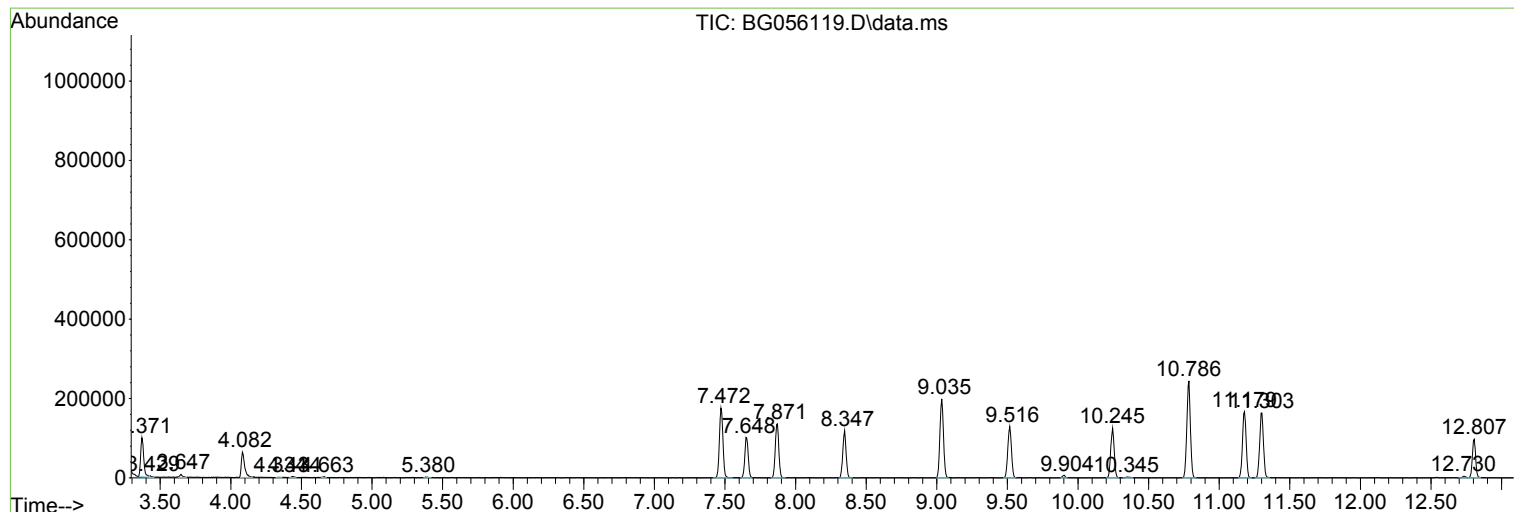
Sum of corrected areas: 16009164

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056119.D
Acq On : 25 Dec 2022 12:04
Operator : CG/JU
Sample : N6017-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056119.D
Acq On : 25 Dec 2022 12:04
Operator : CG/JU
Sample : N6017-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG4

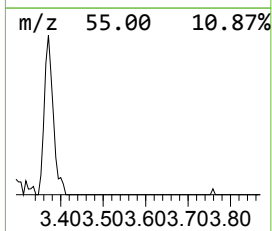
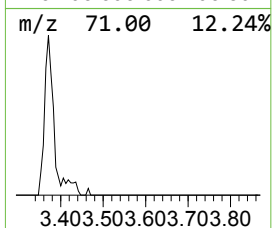
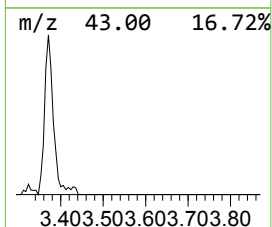
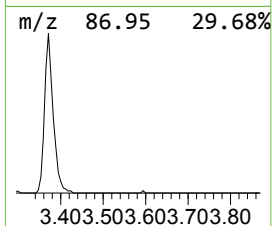
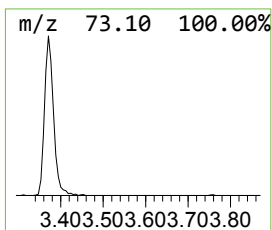
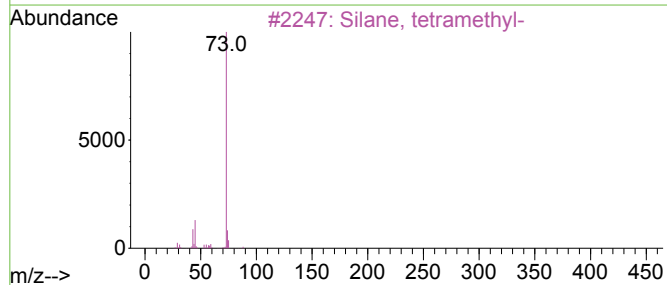
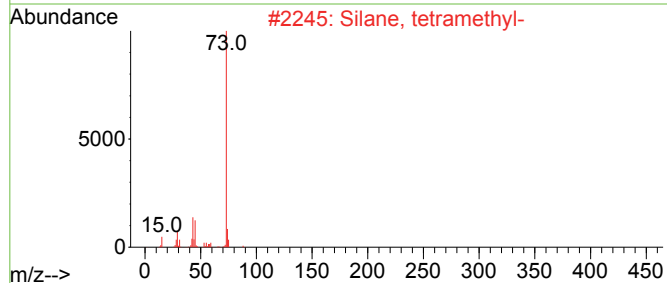
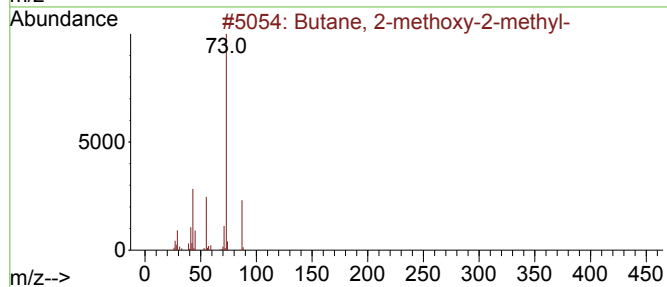
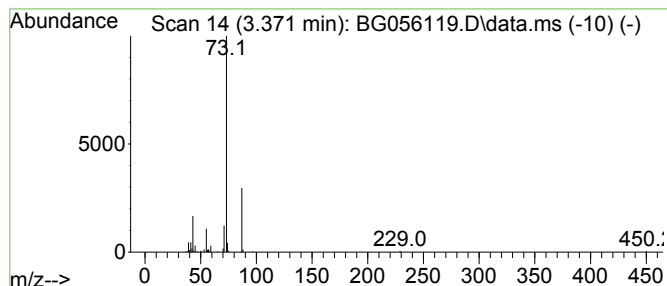
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.371	14.84 ng/ul	144282	1,4-Dichlorobenzene-d4	8.347

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
4			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	9
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056119.D
Acq On : 25 Dec 2022 12:04
Operator : CG/JU
Sample : N6017-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG4

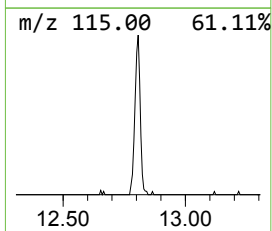
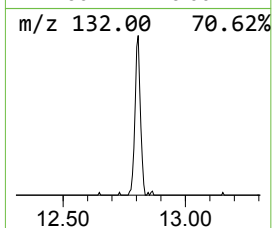
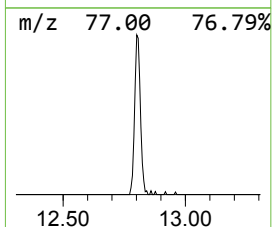
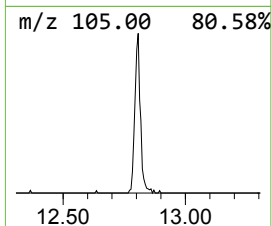
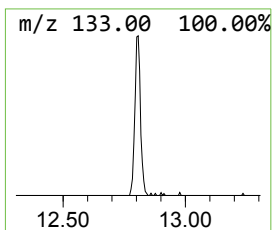
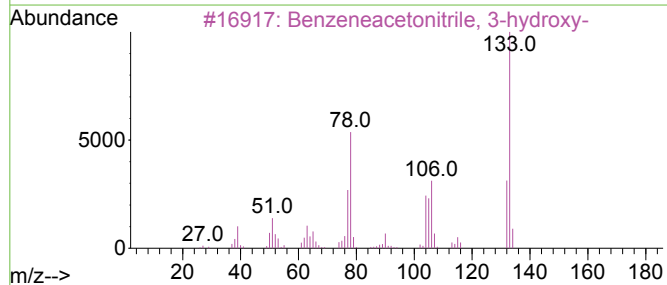
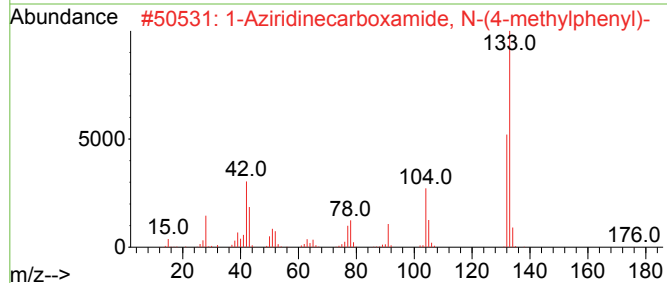
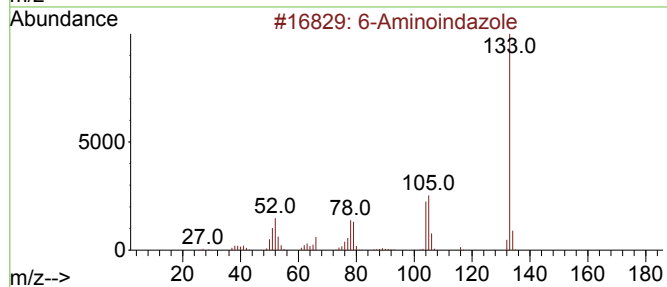
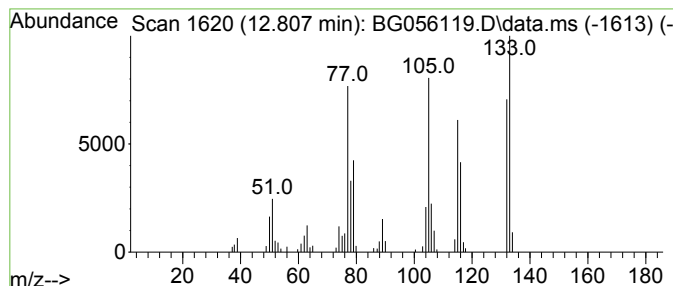
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.807	10.36 ng/ul	156380	Naphthalene-d8	11.179

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Aminoindazole	133	C7H7N3	006967-12-0	35
2			1-Aziridinecarboxamide, N-(4-met...	176	C10H12N2O	000829-65-2	27
3			Benzeneacetonitrile, 3-hydroxy-	133	C8H7NO	025263-44-9	27
4			1H-1,3-Benzodiazol-7-amine	133	C7H7N3	004331-29-7	25
5			3-Methylphenylacetylene	116	C9H8	000766-82-5	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056119.D
Acq On : 25 Dec 2022 12:04
Operator : CG/JU
Sample : N6017-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG4

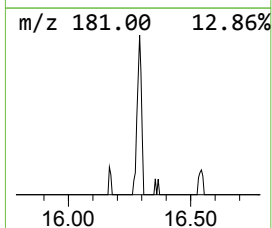
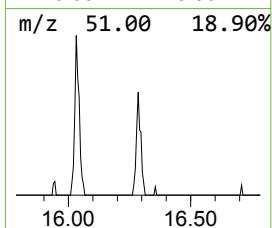
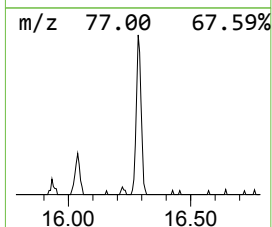
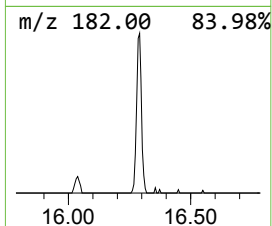
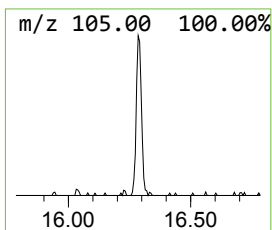
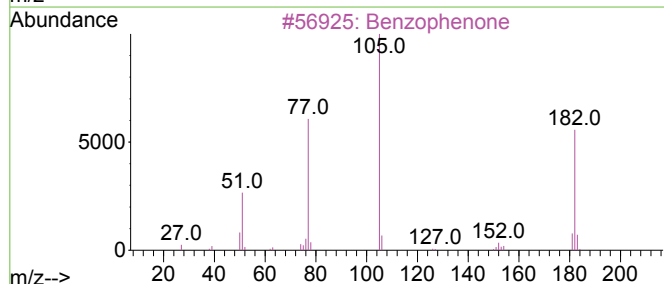
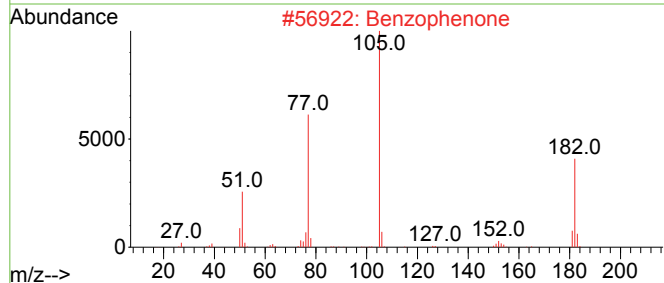
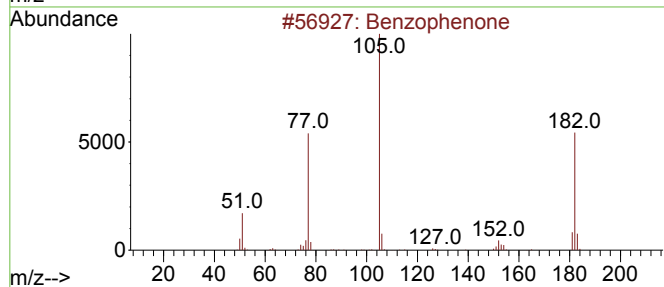
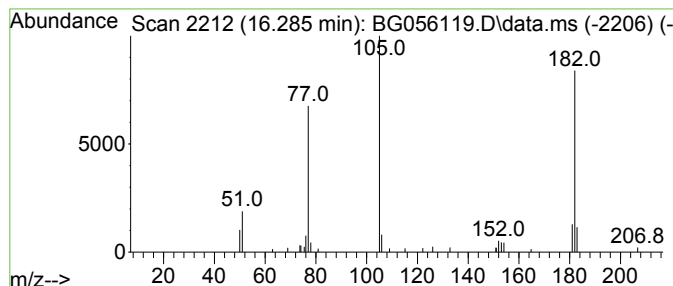
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzophenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.285	2.06 ng/ul	56624	Acenaphthene-d10	14.951

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	94
2			Benzophenone	182	C13H10O	000119-61-9	91
3			Benzophenone	182	C13H10O	000119-61-9	91
4			Benzophenone	182	C13H10O	000119-61-9	74
5			2,4,6-Trimethoxytoluene	182	C10H14O3	014107-97-2	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056119.D
Acq On : 25 Dec 2022 12:04
Operator : CG/JU
Sample : N6017-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG4

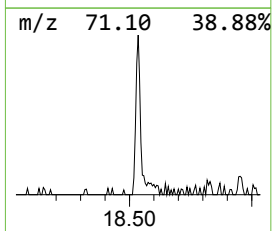
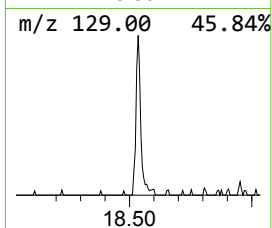
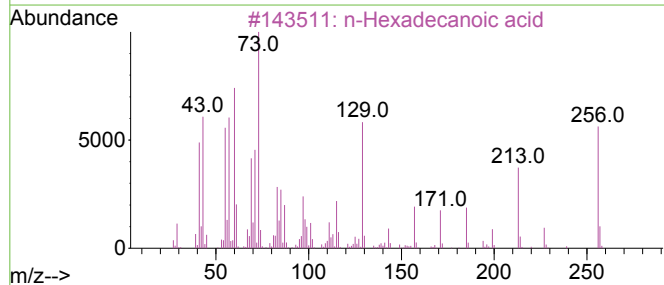
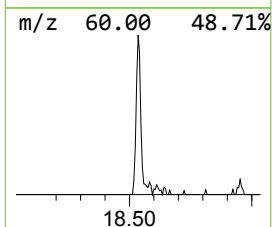
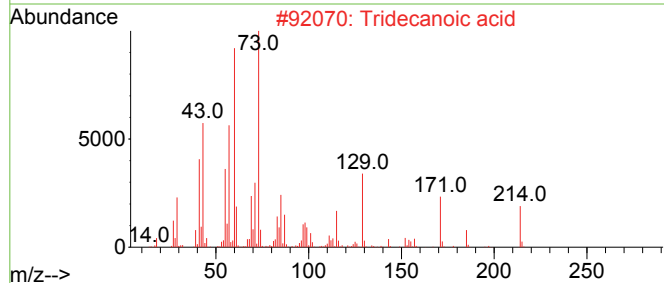
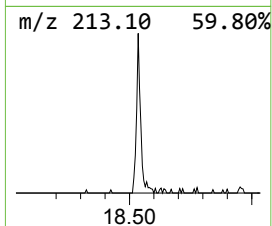
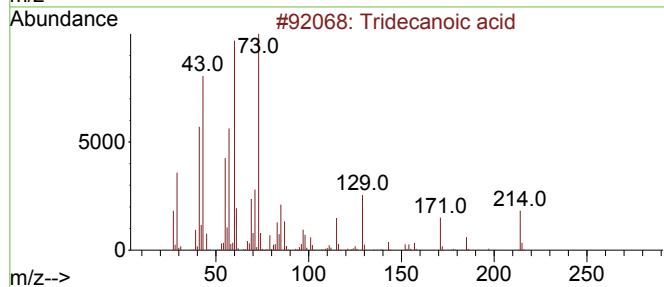
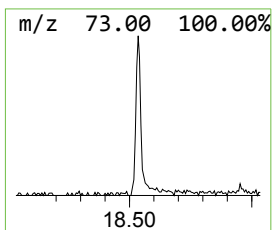
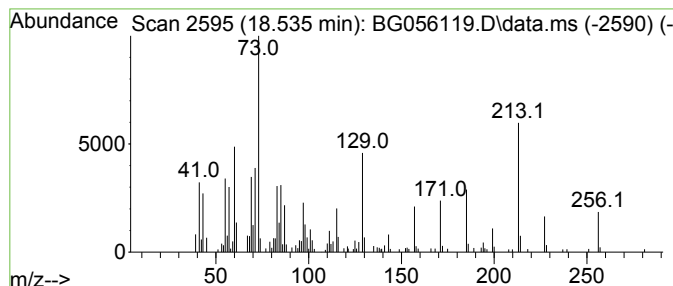
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Tridecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.535	5.08 ng/ul	186486	Phenanthrene-d10	17.689

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecanoic acid	214	C13H26O2	000638-53-9	83
2			Tridecanoic acid	214	C13H26O2	000638-53-9	53
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	53
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	49
5			Tridecanoic acid	214	C13H26O2	000638-53-9	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056119.D
Acq On : 25 Dec 2022 12:04
Operator : CG/JU
Sample : N6017-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG4

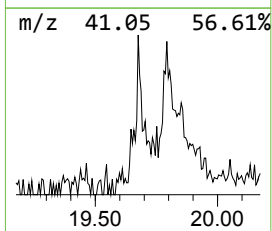
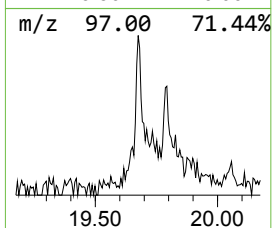
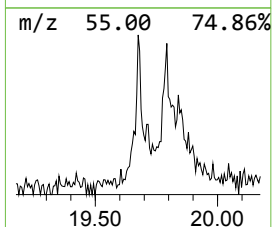
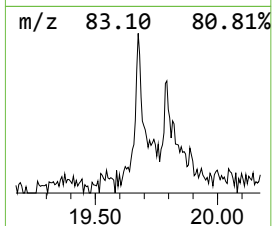
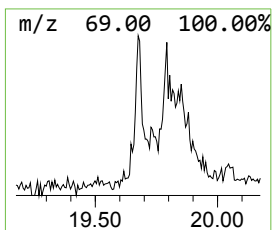
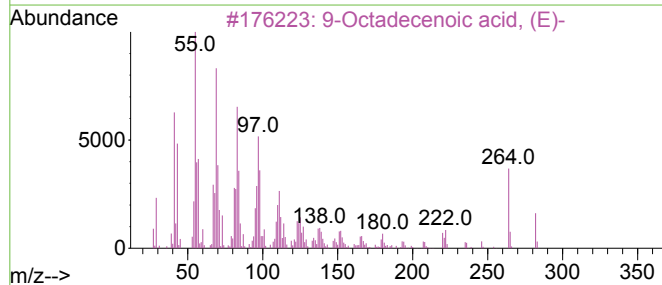
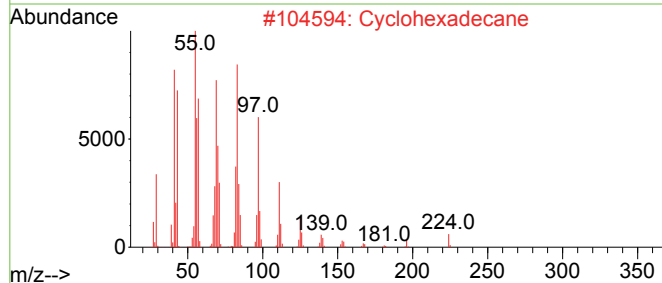
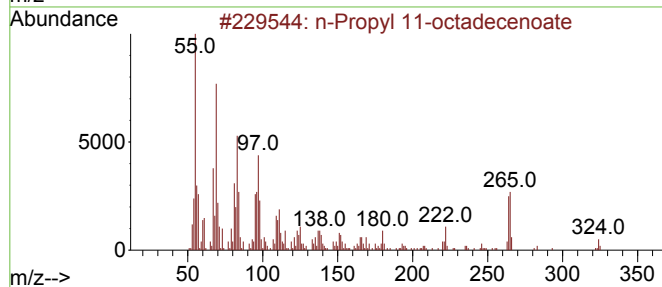
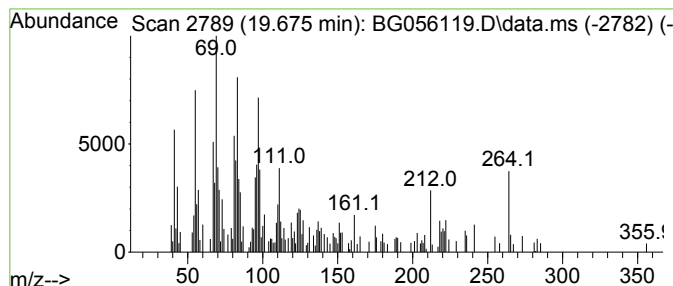
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Propyl 11-octadecenoate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.675	4.58 ng/ul	167856	Phenanthrene-d10	17.689

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Propyl 11-octadecenoate	324	C21H40O2	1000336-71-7	83
2			Cyclohexadecane	224	C16H32	000295-65-8	83
3			9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	83
4			9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	83
5			9-Octadecenoic acid (Z)-, 2,3-di...	356	C21H40O4	000111-03-5	66



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056119.D
Acq On : 25 Dec 2022 12:04
Operator : CG/JU
Sample : N6017-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG4

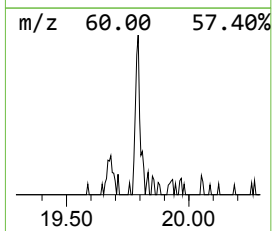
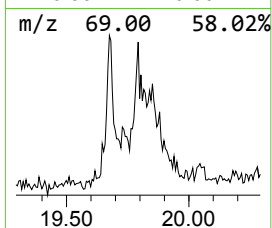
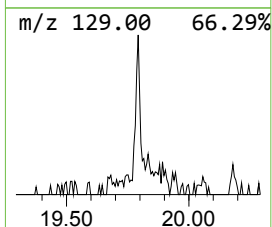
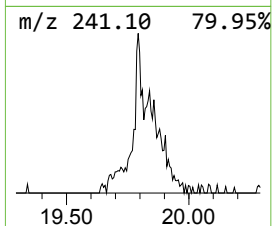
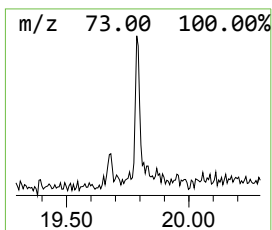
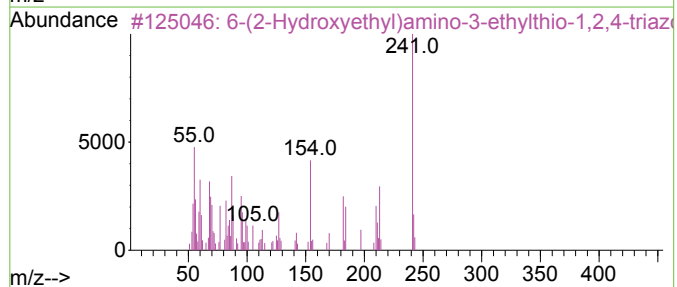
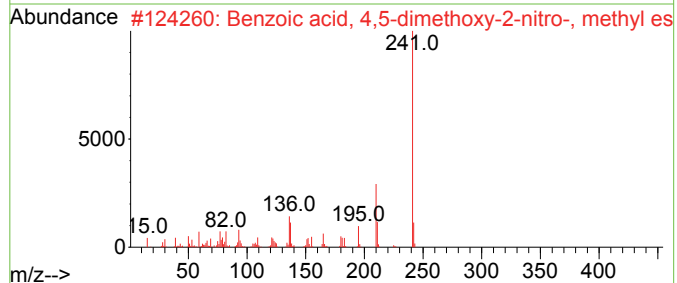
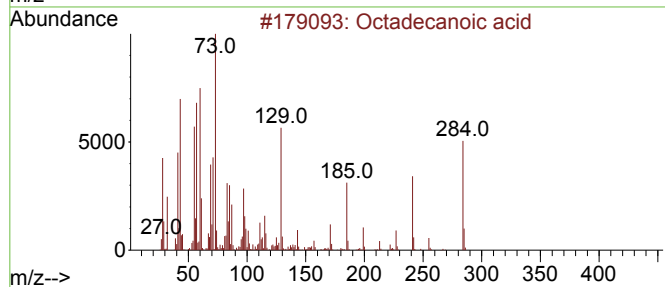
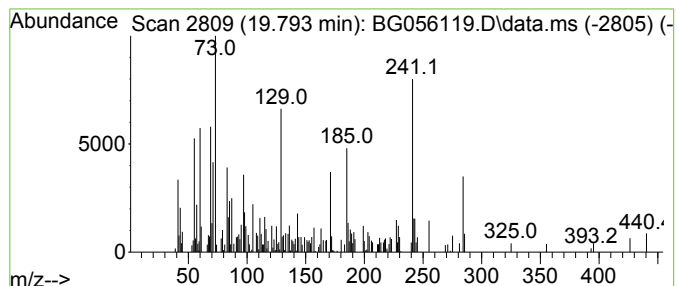
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.793	2.90 ng/ul	106574	Phenanthrene-d10	17.689

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	89
2			Benzoic acid, 4,5-dimethoxy-2-ni...	241	C10H11NO6	026791-93-5	35
3			6-(2-Hydroxyethyl)amino-3-ethylt...	241	C7H11N7O5	1000284-82-7	35
4			n-Decanoic acid	172	C10H20O2	000334-48-5	35
5			Benzaldehyde, (4-nitrophenyl)hyd...	241	C13H11N3O2	003078-09-9	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056119.D
Acq On : 25 Dec 2022 12:04
Operator : CG/JU
Sample : N6017-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG4

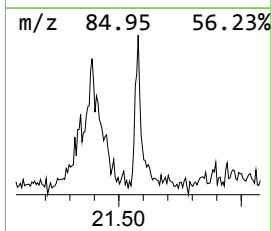
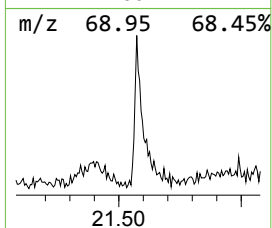
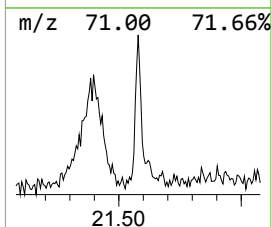
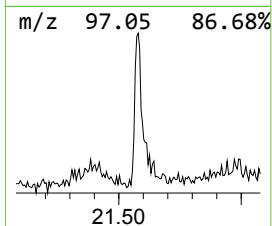
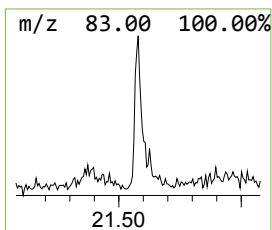
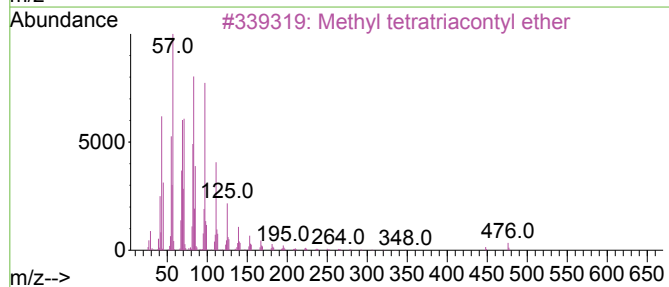
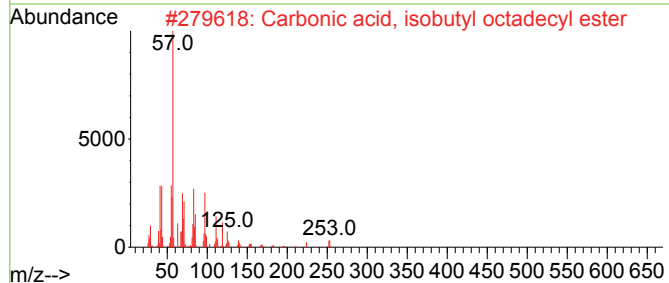
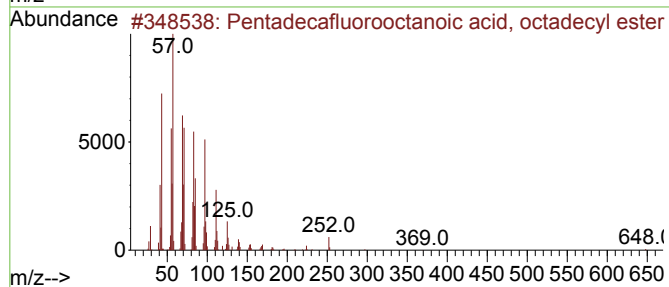
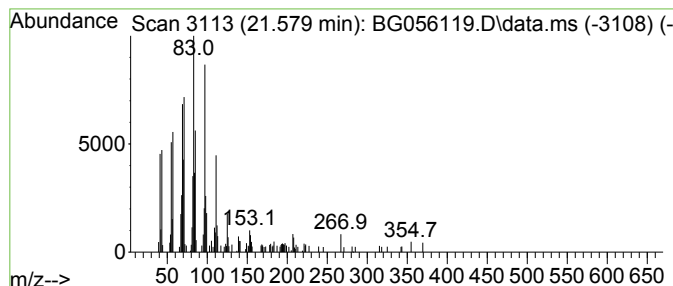
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Pentadecafluorooctanoic aci... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.579	3.58 ng/ul	145777	Chrysene-d12	21.984

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	91
2			Carbonic acid, isobutyl octadecy...	370	C23H46O3	1000314-61-5	91
3			Methyl tetratriacontyl ether	509	C35H72O	1000406-30-1	90
4			Oxirane, [(dodecyloxy)methyl]-	242	C15H30O2	002461-18-9	90
5			2- Chloropropionic acid, pentade...	318	C18H35ClO2	1000292-44-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056119.D
Acq On : 25 Dec 2022 12:04
Operator : CG/JU
Sample : N6017-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG4

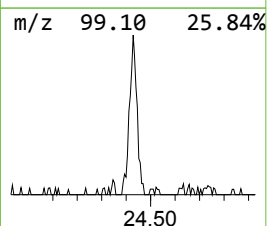
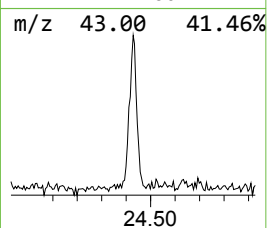
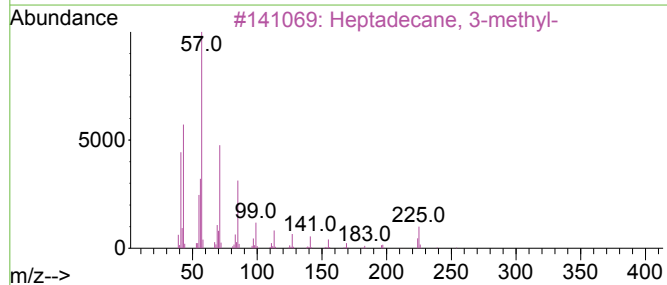
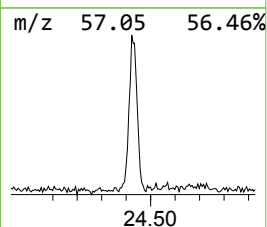
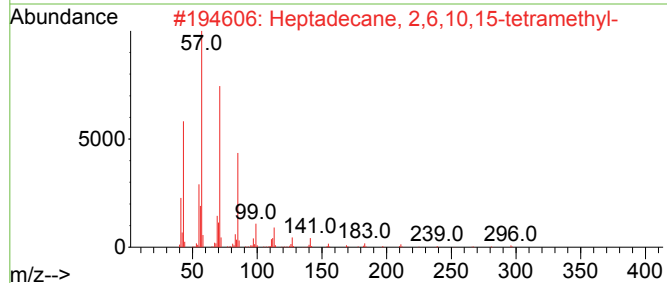
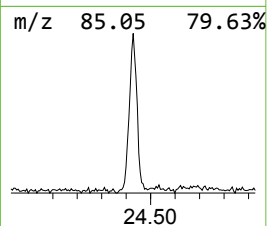
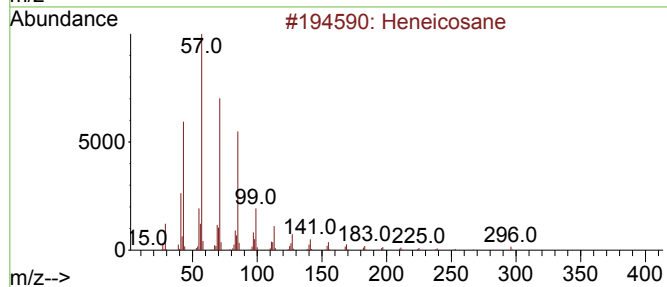
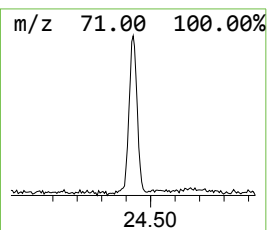
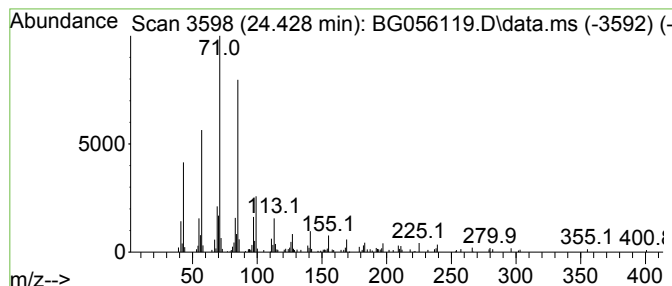
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.428	5.44 ng/ul	239161	Perylene-d12	25.457

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	94
2			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
3			Heptadecane, 3-methyl-	254	C18H38	006418-44-6	91
4			Octacosane	394	C28H58	000630-02-4	91
5			Docosane, 1-iodo-	436	C22H45I	1000406-31-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056119.D
Acq On : 25 Dec 2022 12:04
Operator : CG/JU
Sample : N6017-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG4

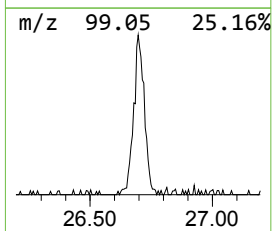
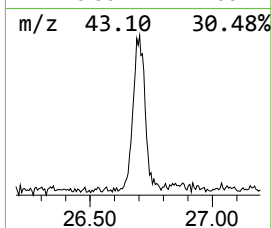
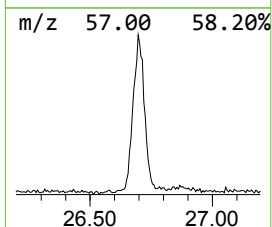
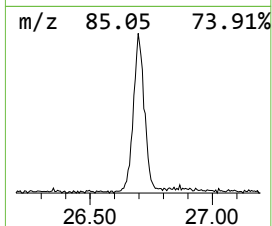
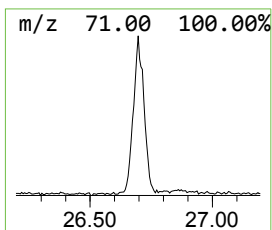
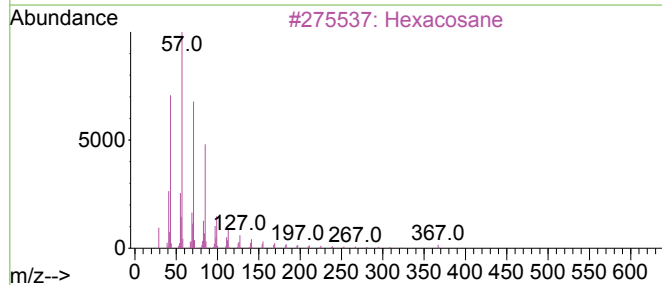
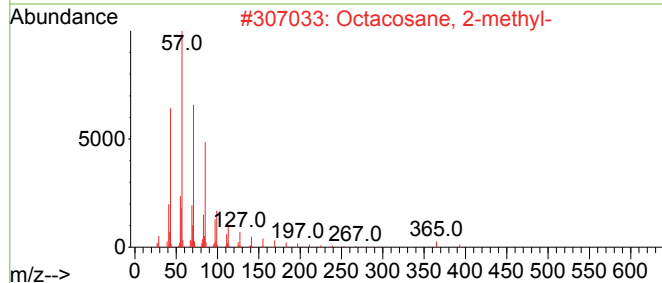
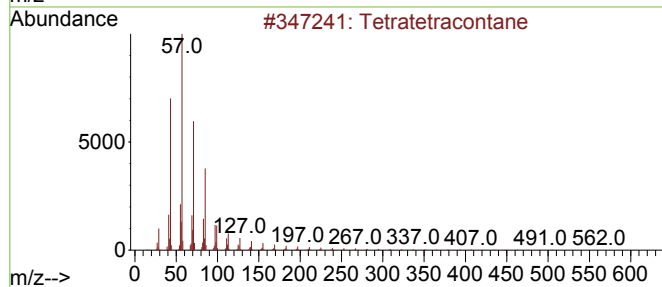
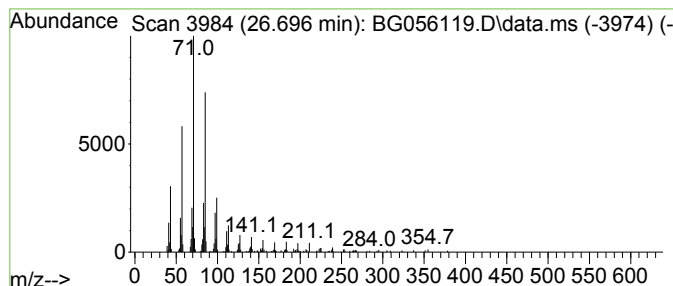
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.696	12.10 ng/ul	532290	Perylene-d12	25.457

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetratetracontane	619	C44H90	007098-22-8	91
2			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
3			Hexacosane	366	C26H54	000630-01-3	91
4			2-Methyltetracosane	352	C25H52	001560-78-7	90
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056119.D
Acq On : 25 Dec 2022 12:04
Operator : CG/JU
Sample : N6017-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG4

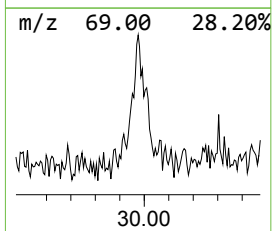
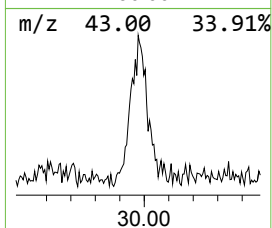
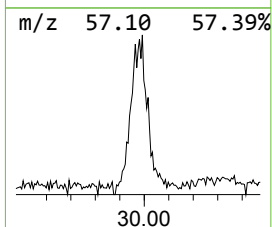
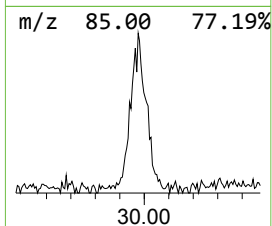
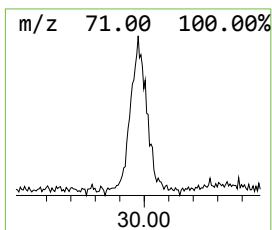
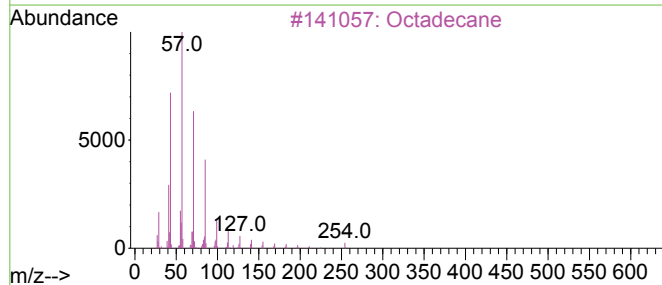
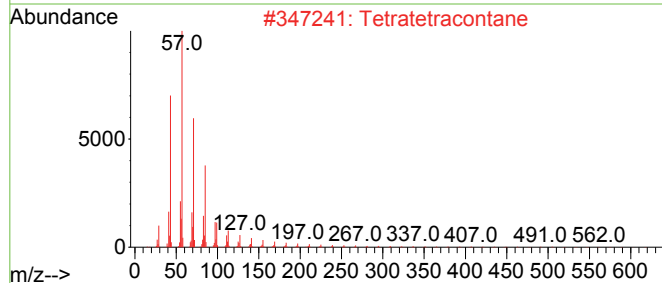
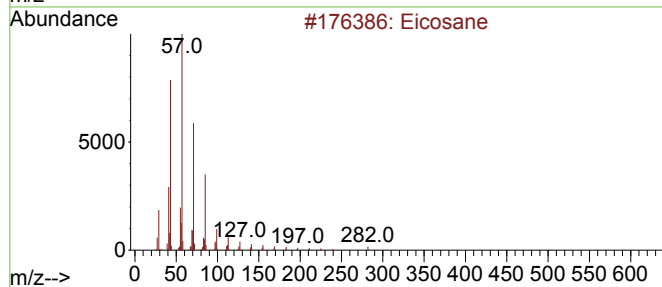
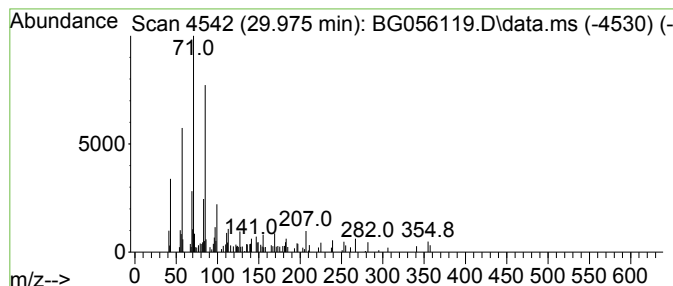
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.975	2.99 ng/ul	131288	Perylene-d12	25.457

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	90
2		Tetratetracontane	619	C44H90	007098-22-8	83
3		Octadecane	254	C18H38	000593-45-3	80
4		Eicosane	282	C20H42	000112-95-8	80
5		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056119.D
Acq On : 25 Dec 2022 12:04
Operator : CG/JU
Sample : N6017-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.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
Butane, 2-metho...	3.371	14.8	ng/ul	144282	1	8.347	194465	20.0
unknown-01	12.807	10.4	ng/ul	156380	2	11.179	301940	20.0
Benzophenone	16.285	2.1	ng/ul	56624	3	14.951	548655	20.0
Tridecanoic acid	18.535	5.1	ng/ul	186486	4	17.689	733782	20.0
n-Propyl 11-oct...	19.675	4.6	ng/ul	167856	4	17.689	733782	20.0
Octadecanoic acid	19.793	2.9	ng/ul	106574	4	17.689	733782	20.0
Pentadecafluoro...	21.579	3.6	ng/ul	145777	5	21.984	814479	20.0
(DEL) Alkane: S...	24.428	5.4	ng/ul	239161	6	25.457	879637	20.0
(DEL) Alkane: S...	26.696	12.1	ng/ul	532290	6	25.457	879637	20.0
(DEL) Alkane: S...	29.975	3.0	ng/ul	131288	6	25.457	879637	20.0