

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
 Data File : BG049334.D
 Acq On : 14 Jul 2021 18:10
 Operator : CG/JU
 Sample : M2996-01DL 5X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AT5DL

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-BG070921.M
 Title : SVOA CALIBRATION

Signal : TIC: BG049334.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.109	6	12	19	rBV2	35331	77476	6.03%	0.954%
2	3.191	22	26	31	rVB3	16303	26587	2.07%	0.328%
3	3.908	146	148	151	rBV4	3758	4757	0.37%	0.059%
4	4.213	196	200	206	rBV4	3834	5479	0.43%	0.067%
5	4.525	250	253	256	rVB4	2142	2515	0.20%	0.031%
6	5.594	431	435	438	rBV3	1589	2352	0.18%	0.029%
7	5.770	463	465	470	rBV4	1669	2125	0.17%	0.026%
8	7.228	705	713	718	rBV3	6642	11514	0.90%	0.142%
9	7.386	735	740	746	rVB	14625	26355	2.05%	0.325%
10	7.598	769	776	781	rBV3	19498	36557	2.84%	0.450%
11	7.780	804	807	814	rBV4	2346	4520	0.35%	0.056%
12	8.021	845	848	849	rBV2	2008	2296	0.18%	0.028%
13	8.068	849	856	862	rVV	113953	197296	15.35%	2.431%
14	8.121	862	865	871	rVV4	4779	8093	0.63%	0.100%
15	8.179	874	875	882	rVB3	3403	3263	0.25%	0.040%
16	8.432	915	918	919	rBV	2513	2285	0.18%	0.028%
17	8.503	928	930	935	rVB4	4012	6321	0.49%	0.078%
18	8.773	971	976	983	rVB4	17058	29387	2.29%	0.362%
19	8.908	996	999	1001	rBV3	2289	2926	0.23%	0.036%
20	9.172	1040	1044	1047	rBV2	1882	3411	0.27%	0.042%
21	9.237	1049	1055	1064	rVB	26780	53425	4.16%	0.658%
22	9.349	1069	1074	1075	rBV2	3632	4666	0.36%	0.057%
23	9.960	1172	1178	1188	rVB3	20581	39744	3.09%	0.490%
24	10.271	1229	1231	1236	rBV5	2153	3306	0.26%	0.041%
25	10.506	1264	1271	1279	rBV2	32506	63172	4.91%	0.778%
26	10.653	1291	1296	1302	rBV5	7070	12959	1.01%	0.160%
27	10.882	1329	1335	1342	rBV	151518	276181	21.48%	3.402%
28	11.017	1352	1358	1364	rBV3	20426	39304	3.06%	0.484%
29	11.288	1401	1404	1409	rVB2	1318	2093	0.16%	0.026%
30	11.987	1521	1523	1529	rBV4	2021	3115	0.24%	0.038%
31	12.516	1608	1613	1615	rBV3	1676	2455	0.19%	0.030%
32	12.756	1651	1654	1660	rVB2	3313	5538	0.43%	0.068%
33	13.144	1715	1720	1727	rBV6	9917	17579	1.37%	0.217%
34	13.226	1729	1734	1741	rVV4	7833	13981	1.09%	0.172%
35	13.303	1743	1747	1753	rVB3	3054	4340	0.34%	0.053%
36	13.414	1761	1766	1775	rBV2	22955	39073	3.04%	0.481%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
 Data File : BG049334.D
 Acq On : 14 Jul 2021 18:10
 Operator : CG/JU
 Sample : M2996-01DL 5X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AT5DL

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-BG070921.M
 Title : SVOA CALIBRATION

37	13.667	1804	1809	1815	rBV4	3161	7396	0.58%	0.091%
38	13.726	1815	1819	1825	rVV4	16270	25907	2.02%	0.319%
39	13.843	1832	1839	1843	rBV6	4214	8629	0.67%	0.106%
40	13.879	1843	1845	1849	rVB4	3731	4022	0.31%	0.050%

41	14.108	1878	1884	1890	rBV	76912	111492	8.67%	1.373%
42	14.407	1929	1935	1942	rBV	73847	120579	9.38%	1.485%
43	14.713	1980	1987	1993	rBV	253141	406856	31.65%	5.012%
44	14.883	2012	2016	2018	rBV2	2580	2953	0.23%	0.036%
45	14.913	2018	2021	2026	rVB4	5524	8513	0.66%	0.105%

46	15.024	2036	2040	2043	rVB2	2988	4531	0.35%	0.056%
47	15.107	2051	2054	2057	rBV	3040	4146	0.32%	0.051%
48	15.248	2076	2078	2083	rVV4	4117	5630	0.44%	0.069%
49	15.306	2083	2088	2089	rVV2	10555	14837	1.15%	0.183%
50	15.336	2089	2093	2099	rVB3	30962	46199	3.59%	0.569%

51	15.430	2107	2109	2112	rBV2	2569	2705	0.21%	0.033%
52	15.459	2112	2114	2117	rVB3	2643	2496	0.19%	0.031%
53	15.606	2134	2139	2141	rBV	1696	3180	0.25%	0.039%
54	15.700	2150	2155	2161	rBV2	106070	163749	12.74%	2.017%
55	15.759	2161	2165	2170	rVB3	5367	8982	0.70%	0.111%

56	15.835	2170	2178	2188	rBV	142409	248566	19.33%	3.062%
57	15.912	2189	2191	2194	rBV4	2725	2853	0.22%	0.035%
58	15.941	2194	2196	2197	rBV2	4572	2870	0.22%	0.035%
59	16.041	2210	2213	2216	rBV4	3075	3332	0.26%	0.041%
60	16.117	2219	2226	2228	rBV4	2809	4863	0.38%	0.060%

61	16.211	2239	2242	2247	rBV3	2243	2908	0.23%	0.036%
62	16.288	2253	2255	2261	rBV4	2198	4107	0.32%	0.051%
63	16.382	2267	2271	2273	rBV3	3208	4748	0.37%	0.058%
64	16.523	2291	2295	2298	rBV3	2605	2685	0.21%	0.033%
65	16.593	2305	2307	2313	rVB3	4859	6374	0.50%	0.079%

66	16.740	2328	2332	2335	rBV3	3634	5799	0.45%	0.071%
67	16.805	2340	2343	2346	rBV2	3470	3959	0.31%	0.049%
68	16.951	2365	2368	2371	rVB4	2913	2987	0.23%	0.037%
69	17.110	2388	2395	2406	rBV	660937	1054666	82.03%	12.993%
70	17.404	2442	2445	2448	rBV3	1919	2081	0.16%	0.026%

71	17.463	2448	2455	2461	rVB	325245	498991	38.81%	6.147%
72	17.557	2465	2471	2479	rVB	112861	171904	13.37%	2.118%
73	17.633	2479	2484	2486	rBV5	3954	5657	0.44%	0.070%
74	17.727	2499	2500	2503	rBV2	2553	2279	0.18%	0.028%
75	17.809	2509	2514	2527	rBV	192407	314371	24.45%	3.873%

76	18.115	2560	2566	2572	rVB5	8714	18261	1.42%	0.225%
----	--------	------	------	------	------	------	-------	-------	--------

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
 Data File : BG049334.D
 Acq On : 14 Jul 2021 18:10
 Operator : CG/JU
 Sample : M2996-01DL 5X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AT5DL

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-BG070921.M
 Title : SVOA CALIBRATION

77	18.262	2588	2591	2594	rBV3	3988	6099	0.47%	0.075%
78	18.309	2595	2599	2601	rVB4	4836	6929	0.54%	0.085%
79	18.338	2601	2604	2607	rBV4	3615	5672	0.44%	0.070%
80	18.414	2615	2617	2622	rBV5	1962	3130	0.24%	0.039%
81	18.514	2631	2634	2635	rBV2	2969	3492	0.27%	0.043%
82	18.550	2639	2640	2643	rBV3	2626	2248	0.17%	0.028%
83	18.620	2650	2652	2653	rBV2	3249	2458	0.19%	0.030%
84	18.655	2655	2658	2660	rVB2	6385	6456	0.50%	0.080%
85	18.708	2665	2667	2669	rBV2	2930	2703	0.21%	0.033%
86	18.849	2687	2691	2692	rBV3	2037	2734	0.21%	0.034%
87	19.396	2782	2784	2787	rVB4	2796	2682	0.21%	0.033%
88	19.543	2804	2809	2815	rBV	1036662	1285636	100.00%	15.838%
89	19.707	2832	2837	2843	rBV	290551	407239	31.68%	5.017%
90	19.842	2855	2860	2870	rBV	130842	196373	15.27%	2.419%
91	20.071	2897	2899	2902	rVB3	7659	7657	0.60%	0.094%
92	20.536	2974	2978	2983	rVB2	64954	80468	6.26%	0.991%
93	20.794	3020	3022	3024	rBV3	8282	7206	0.56%	0.089%
94	21.217	3089	3094	3104	rBV	151086	244056	18.98%	3.007%
95	21.740	3177	3183	3192	rVB	328040	558364	43.43%	6.879%
96	22.915	3378	3383	3395	rVB5	55535	122943	9.56%	1.515%
97	24.801	3697	3704	3717	rVB	69374	194388	15.12%	2.395%
98	25.030	3734	3743	3754	rVB3	193720	571279	44.44%	7.038%
99	25.383	3798	3803	3815	rVB3	16750	43670	3.40%	0.538%

Sum of corrected areas: 8117391

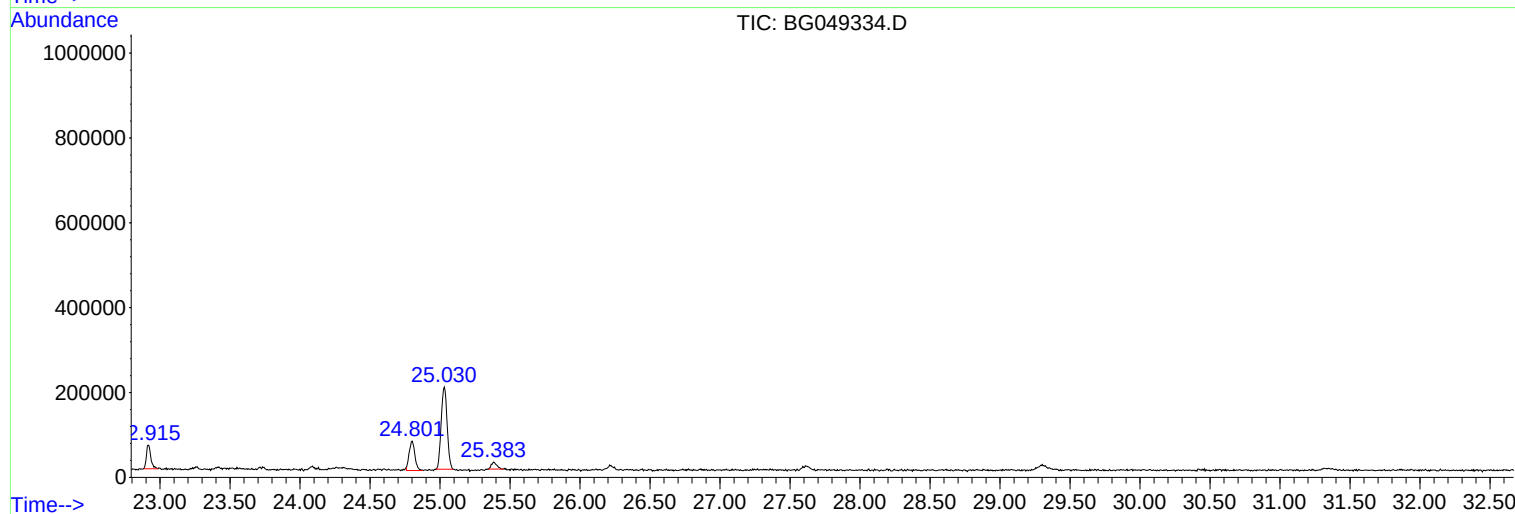
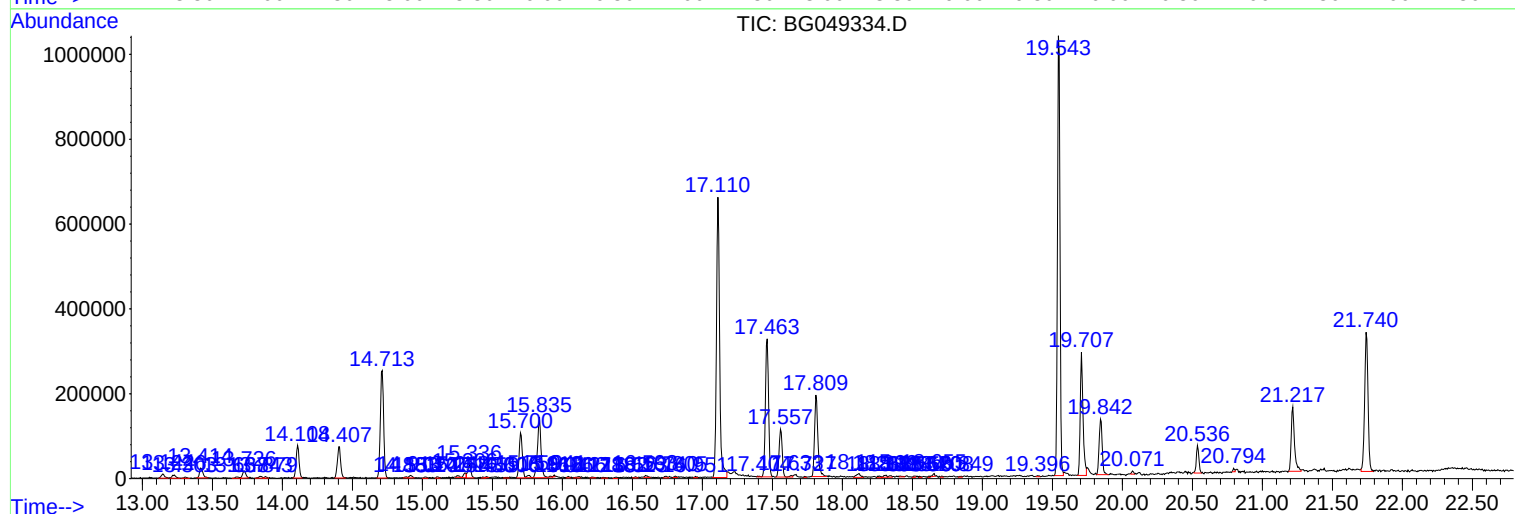
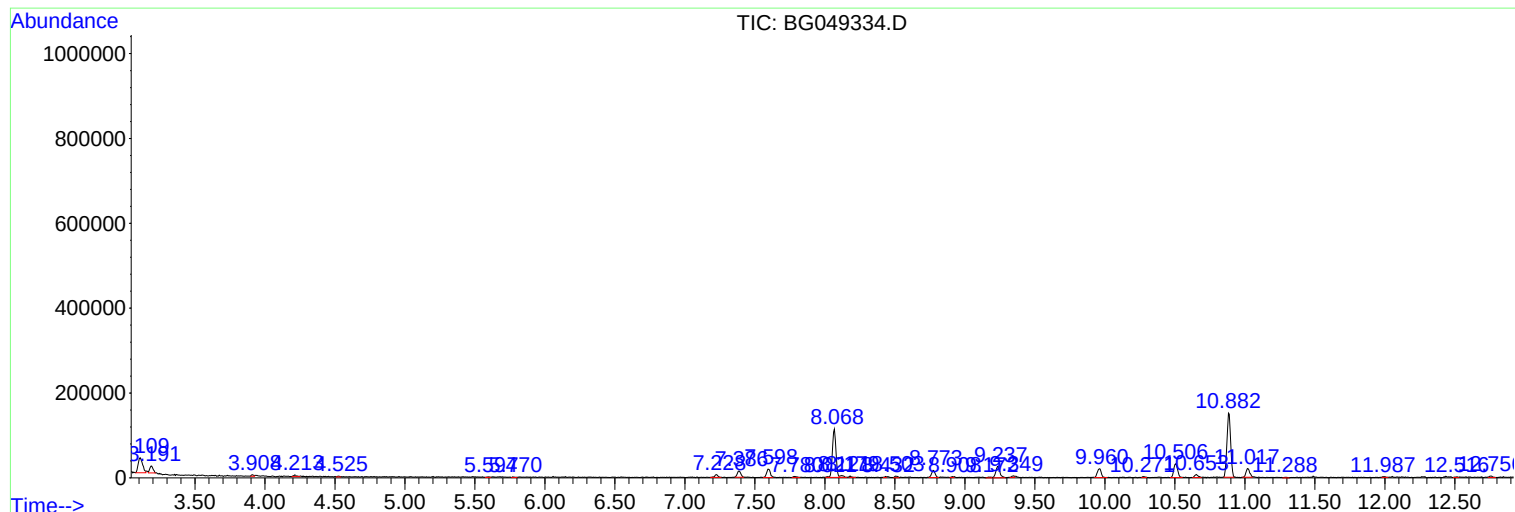
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
Data File : BG049334.D
Acq On : 14 Jul 2021 18:10
Operator : CG/JU
Sample : M2996-01DL 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AT5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L

TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
Data File : BG049334.D
Acq On : 14 Jul 2021 18:10
Operator : CG/JU
Sample : M2996-01DL 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AT5DL

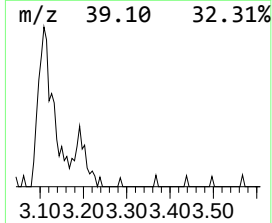
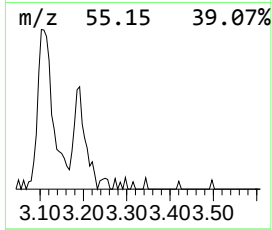
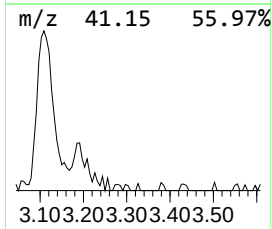
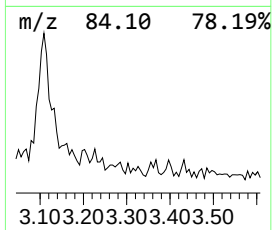
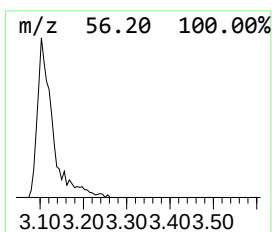
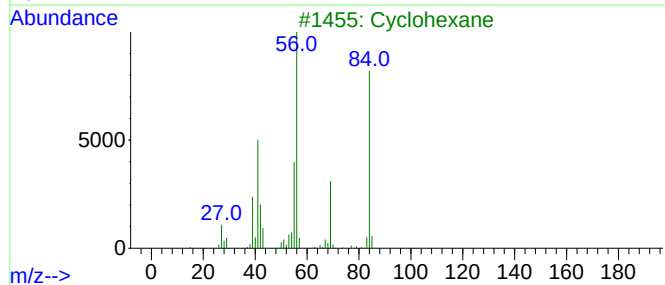
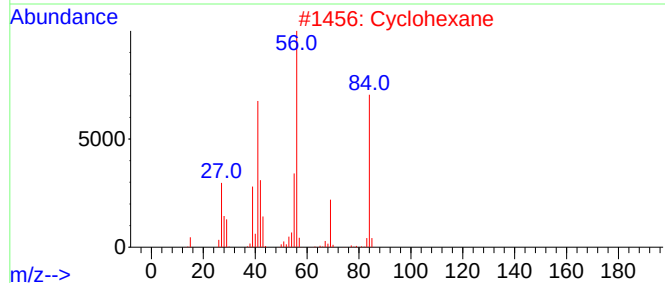
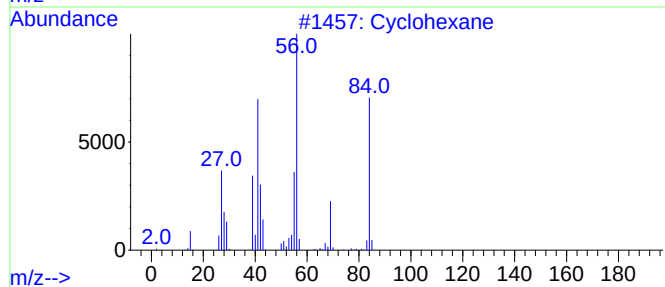
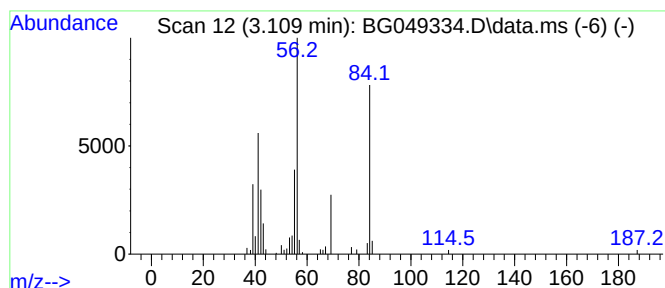
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Cyclic3.11 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.110	7.85 ng/ul	77476	1,4-Dichlorobenzene-d4	8.068

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane	84	C6H12	000110-82-7	91
2		Cyclohexane	84	C6H12	000110-82-7	91
3		Cyclohexane	84	C6H12	000110-82-7	90
4		Cyclohexane	84	C6H12	000110-82-7	90
5		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
Data File : BG049334.D
Acq On : 14 Jul 2021 18:10
Operator : CG/JU
Sample : M2996-01DL 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AT5DL

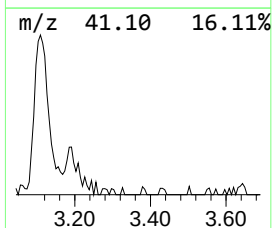
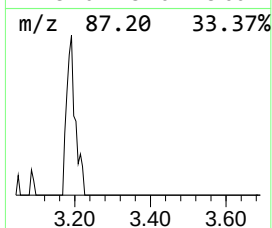
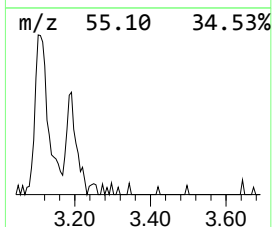
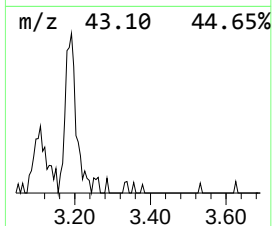
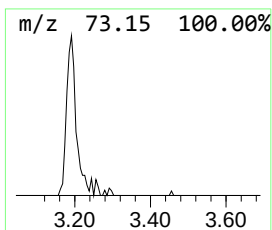
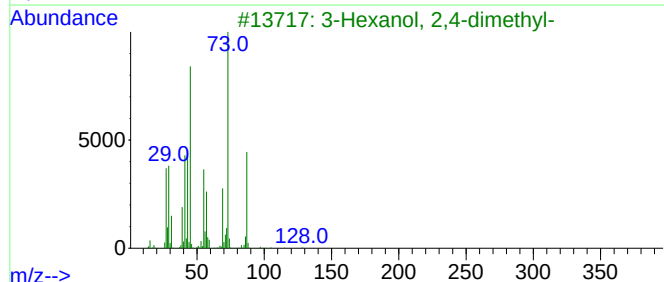
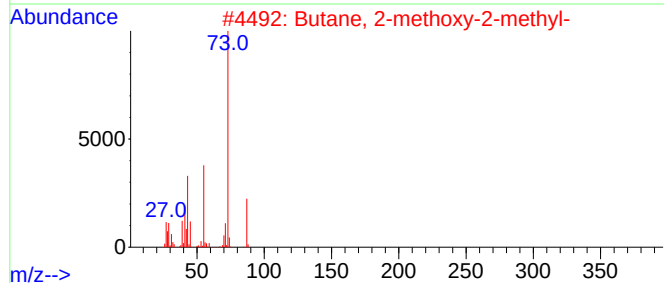
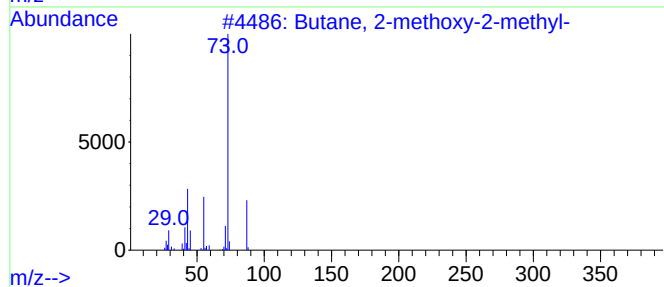
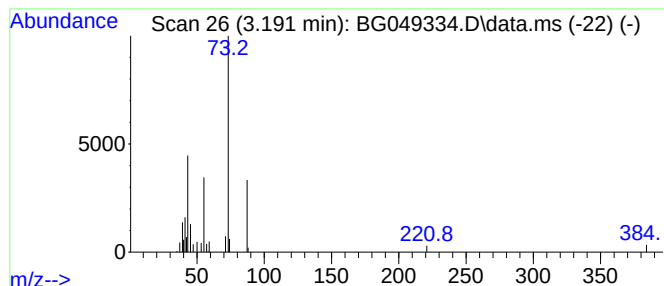
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.190	2.70 ng/ul	26587	1,4-Dichlorobenzene-d4	8.068

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	59
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	36
3			3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	28
4			2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	17
5			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	12



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
Data File : BG049334.D
Acq On : 14 Jul 2021 18:10
Operator : CG/JU
Sample : M2996-01DL 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AT5DL

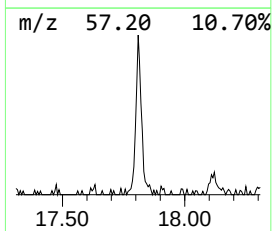
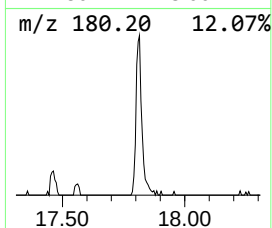
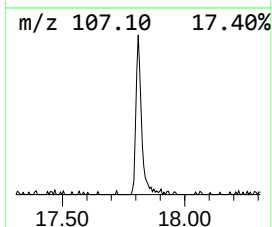
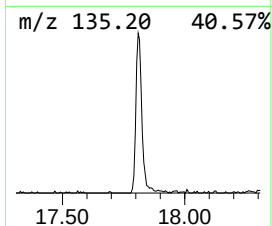
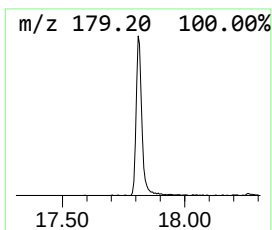
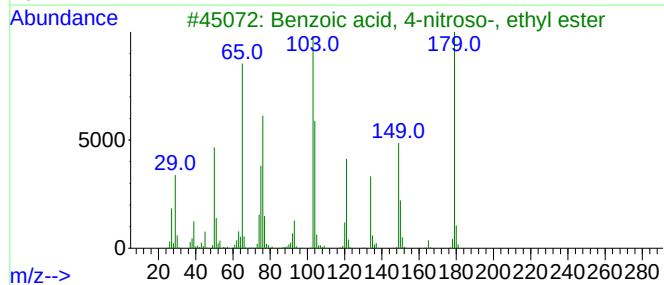
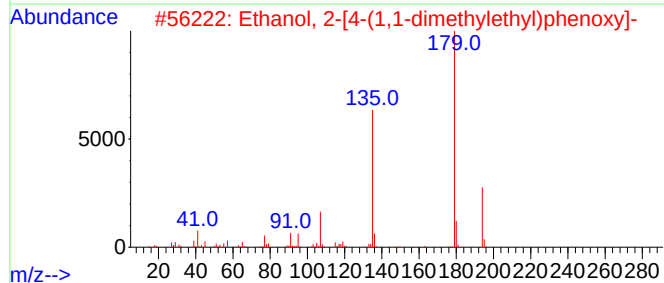
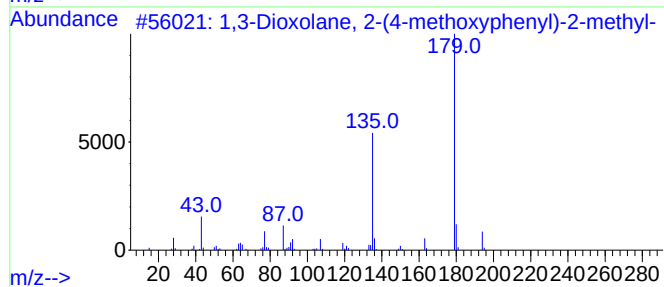
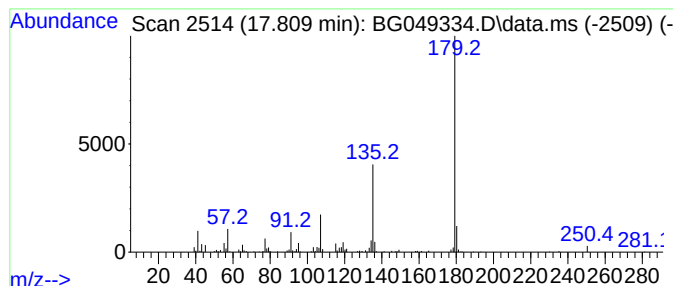
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1,3-Dioxolane, 2-(4-methoxy... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.810	12.60 ng/ul	314371	Phenanthrene-d10	17.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Dioxolane, 2-(4-methoxypheny...	194	C11H14O3	036881-00-2	56
2			Ethanol, 2-[4-(1,1-dimethylethyl...]	194	C12H18O2	000713-46-2	50
3			Benzoic acid, 4-nitroso-, ethyl ...	179	C9H9NO3	007476-79-1	43
4			2,4-Dimethoxyphenyl isocyanate	179	C9H9NO3	084370-87-6	37
5			Acetic acid, [2-[(2-propenylamin...	235	C12H13NO4	000119-45-9	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
Data File : BG049334.D
Acq On : 14 Jul 2021 18:10
Operator : CG/JU
Sample : M2996-01DL 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AT5DL

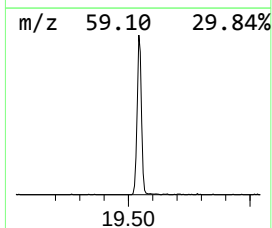
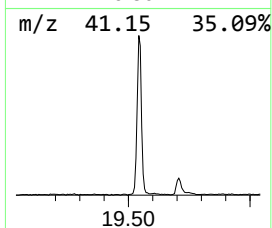
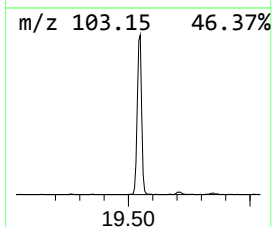
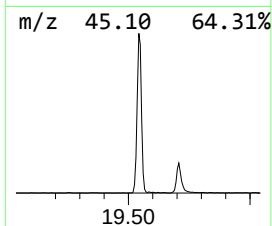
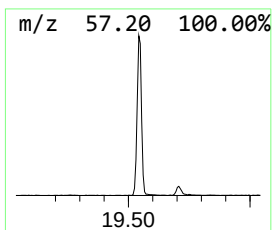
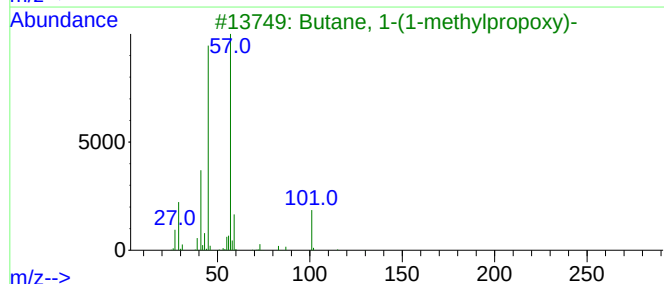
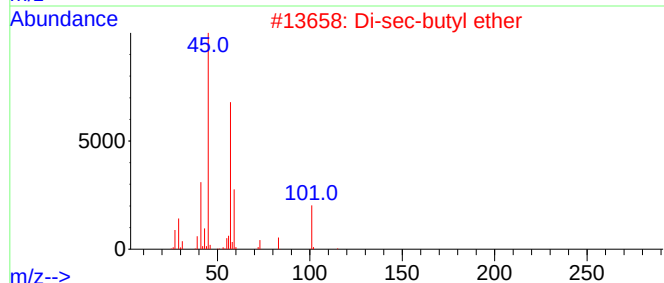
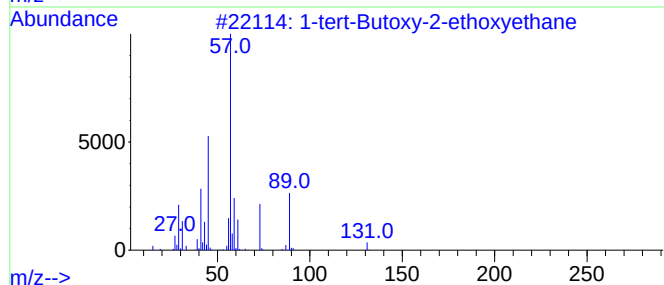
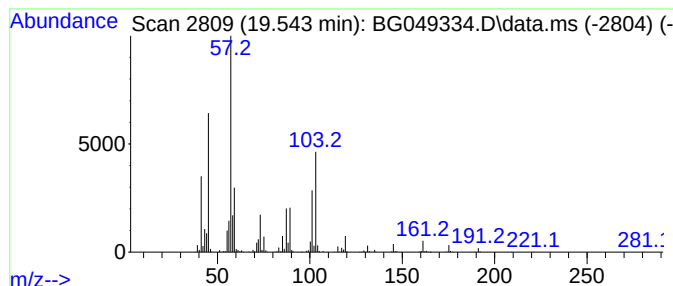
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.540	51.53 ng/ul	1285640	Phenanthrene-d10	17.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-tert-Butoxy-2-ethoxyethane	146	C8H18O2	051422-54-9	38
2			Di-sec-butyl ether	130	C8H18O	006863-58-7	38
3			Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	35
4			3,6,9,12-Tetraoxahexadecan-1-ol	250	C12H26O5	001559-34-8	35
5			2-[2-[2-[2-[2-[2-(2-Methoxyethox...	340	C15H32O8	1000352-04-1	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
Data File : BG049334.D
Acq On : 14 Jul 2021 18:10
Operator : CG/JU
Sample : M2996-01DL 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AT5DL

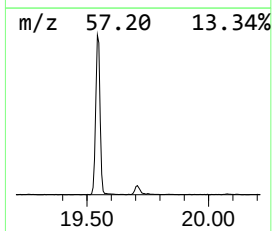
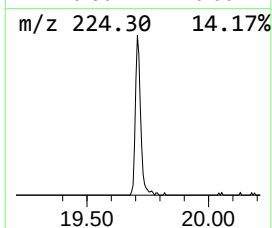
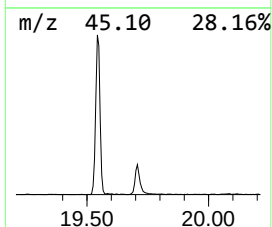
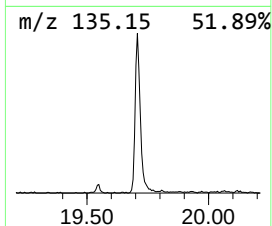
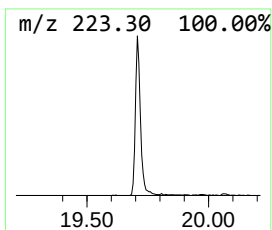
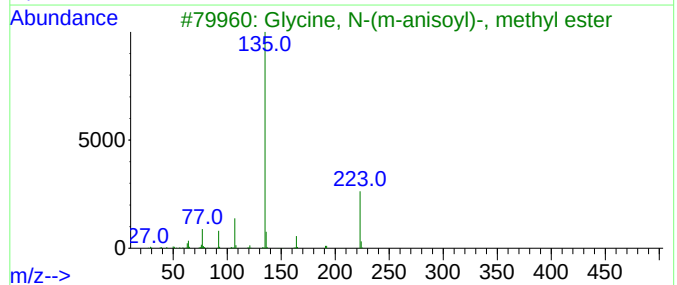
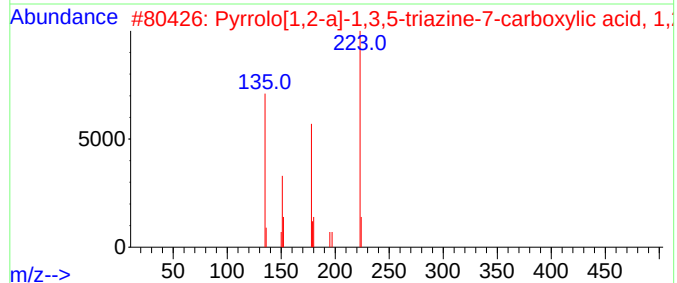
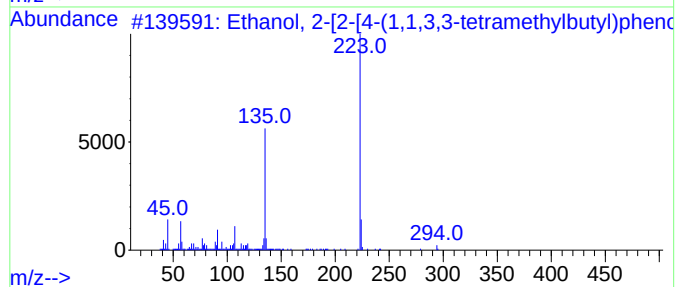
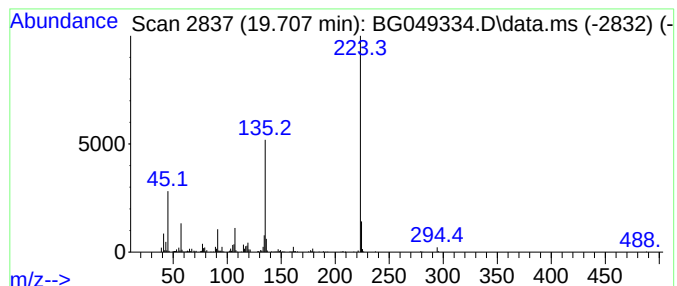
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Ethanol, 2-[2-[4-(1,1,3,3-t... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.710	14.59 ng/ul	407239	Chrysene-d12	21.740

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-[2-[4-(1,1,3,3-tetram...	294	C18H30O3	002315-61-9	93
2			Pyrrolo[1,2-a]-1,3,5-triazine-7-...	223	C9H9N3O4	054449-89-7	90
3			Glycine, N-(m-anisoyl)-, methyl ...	223	C11H13NO4	1000299-70-7	59
4			Tricyclo[3.3.3.0(1,5)]undec-6-en...	223	C14H13N3	118894-71-6	43
5			4'-Methoxybutyrophenone	178	C11H14O2	004160-51-4	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
Data File : BG049334.D
Acq On : 14 Jul 2021 18:10
Operator : CG/JU
Sample : M2996-01DL 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AT5DL

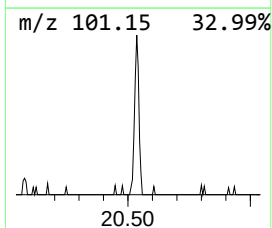
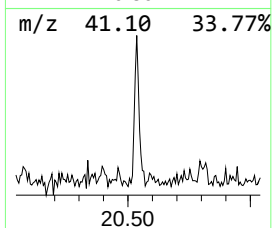
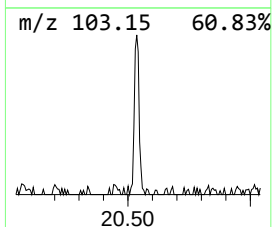
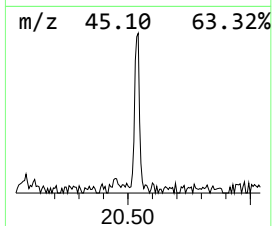
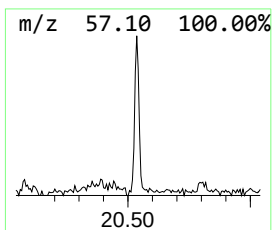
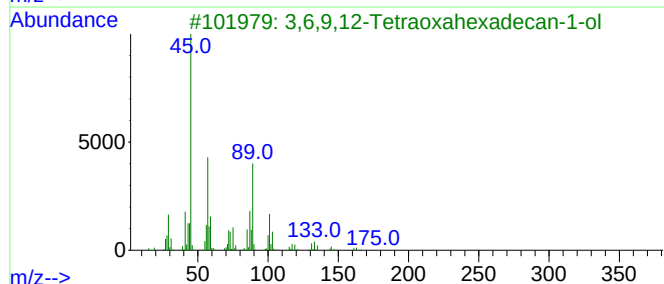
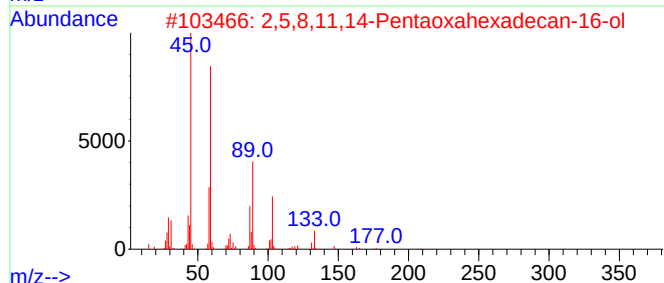
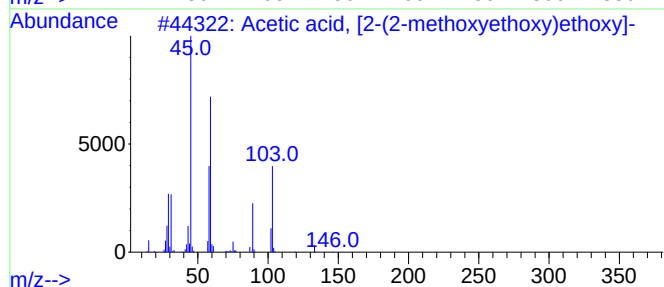
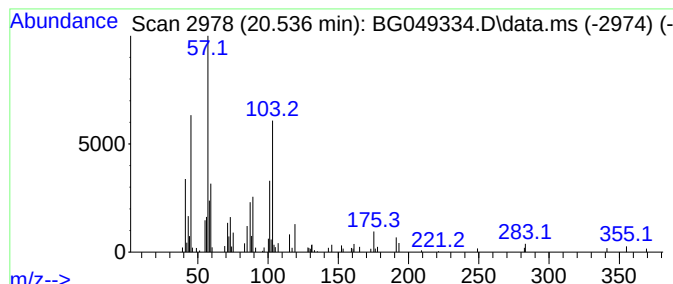
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown-02 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.540	2.88 ng/ul	80468	Chrysene-d12	21.740

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetic acid, [2-(2-methoxyethoxy)...	178	C7H14O5	016024-58-1	47
2			2,5,8,11,14-Pentaoxahexadecan-16-ol	252	C11H24O6	023778-52-1	43
3			3,6,9,12-Tetraoxahexadecan-1-ol	250	C12H26O5	001559-34-8	43
4			2,5,8,11,14-Pentaoxahexadecan-16-ol	252	C11H24O6	023778-52-1	38
5			1-(2-Hydroxyethoxy)-2-(vinylthio...	148	C6H12O2S	086241-10-3	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
Data File : BG049334.D
Acq On : 14 Jul 2021 18:10
Operator : CG/JU
Sample : M2996-01DL 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AT5DL

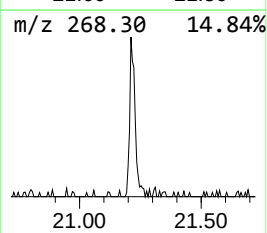
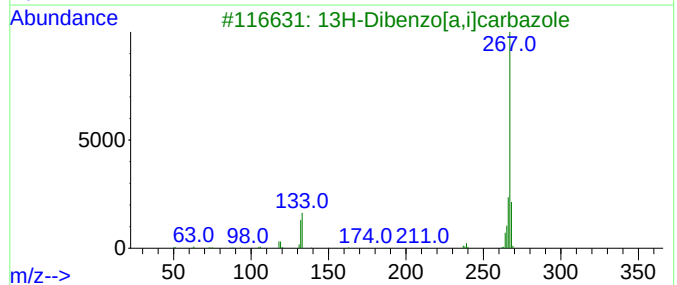
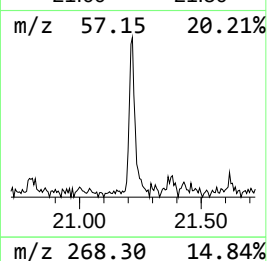
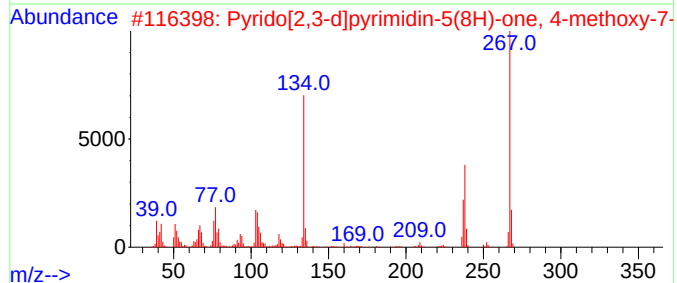
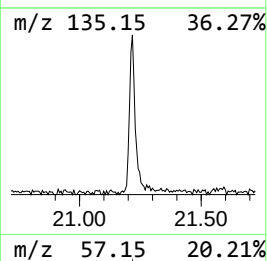
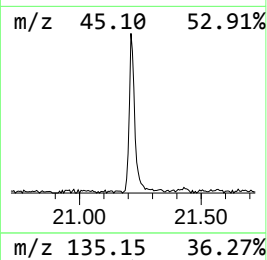
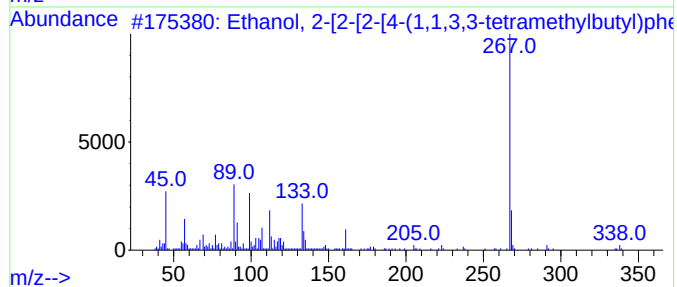
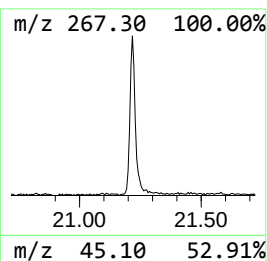
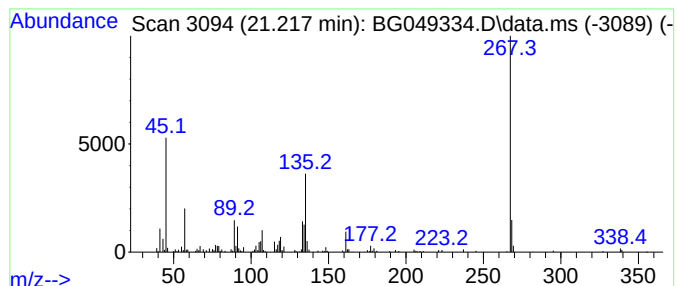
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Ethanol, 2-[2-[2-[4-(1,1,3,... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.220	8.74 ng/ul	244056	Chrysene-d12	21.740

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-[2-[2-[4-(1,1,3,3-tet...	338	C20H34O4	002315-62-0	50
2			Pyrido[2,3-d]pyrimidin-5(8H)-one...	267	C15H13N3O2	1000260-26-0	38
3			13H-Dibenzo[a,i]carbazole	267	C20H13N	000239-64-5	38
4			5H-Naphtho[2,3-c]carbazole	267	C20H13N	000198-96-9	35
5			Phthalic acid, di(2-propylphenyl...	402	C26H26O4	1000357-03-1	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
 Data File : BG049334.D
 Acq On : 14 Jul 2021 18:10
 Operator : CG/JU
 Sample : M2996-01DL 5X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AT5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: C...	3.110	7.8	ng/ul	77476	1	8.068	197296	20.0
Butane, 2-metho...	3.190	2.7	ng/ul	26587	1	8.068	197296	20.0
1,3-Dioxolane, ...	17.810	12.6	ng/ul	314371	4	17.463	498991	20.0
unknown-01	19.540	51.5	ng/ul	1285640	4	17.463	498991	20.0
Ethanol, 2-[2-[...	19.710	14.6	ng/ul	407239	5	21.740	558364	20.0
unknown-02	20.540	2.9	ng/ul	80468	5	21.740	558364	20.0
Ethanol, 2-[2-[...	21.220	8.7	ng/ul	244056	5	21.740	558364	20.0
unknown-03	22.920	4.4	ng/ul	122943	5	21.740	558364	20.0