LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
Data File : BG049333.D
Acq On : 14 Jul 2021 17:29
Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

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: BG049333.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.111	5	12	20	rBV	95417	234428	4.20%	0.786%
2	3.187	21	25	39	rVB2	79065	160234	2.87%	0.537%
3	3.892	142	145	152	rBV3	19573	40881	0.73%	0.137%
4	4.215	195	200	207	rBV6	7332	14766	0.26%	0.050%
5	7.223	705	712	722	rBV2	31220	65682	1.18%	0.220%
6	7.388	734	740	746	rBV	81124	142631	2.55%	0.478%
7	7.599	769	776	784	rVB	100579	187487	3.36%	0.629%
8	7.782	800	807	814	rBV5	8039	17849	0.32%	0.060%
9	8.011	841	846	850	rBV2	7632	14056	0.25%	0.047%
10	8.064	850	855	861	rVV	107009	191914	3.44%	0.644%
11	8.122	861	865	871	rVV3	20870	35772	0.64%	0.120%
12	8.181	871	875	883	rVB2	12830	21850	0.39%	0.073%
13	8.446	913	920	925	rBV	14246	25398	0.45%	0.085%
14	8.504	925	930	936	rVV4	19975	35549	0.64%	0.119%
15	8.769	968	975	982	rVV	86992	157629	2.82%	0.529%
16	8.916	995	1000	1009	rVB5	13191	23697	0.42%	0.079%
17	9.233	1048	1054	1064	rVV	125446	249920	4.47%	0.838%
18	9.356	1068	1075	1082	rBV3	20062	39012	0.70%	0.131%
19	9.956	1171	1177	1189	rVB	118413	217485	3.89%	0.729%
20	10.279	1226	1232	1241	rBV4	12116	22138	0.40%	0.074%
21	10.384	1241	1250	1254	rBV7	6717	14610	0.26%	0.049%
22	10.502	1263	1270	1278	rBV	168314	309625	5.54%	1.038%
23	10.649	1290	1295	1305	rVV2	34727	67363	1.21%	0.226%
24	10.884	1328	1335	1342	rBV	140675	259331	4.64%	0.870%
25	11.019	1351	1358	1371	rBV	106298	221229	3.96%	0.742%
26	11.501	1430	1440	1443	rBV8	11615	33233	0.59%	0.111%
27	11.994	1521	1524	1530	rBV4	11409	16151	0.29%	0.054%
28	12.271	1566	1571	1579	rVB9	11580	21901	0.39%	0.073%
29	12.429	1594	1598	1605	rBV2	14310	28655	0.51%	0.096%
30	12.758	1645	1654	1659	rVV8	18545	39111	0.70%	0.131%
31	12.840	1662	1668	1674	rBV8	8082	18659	0.33%	0.063%
32	12.911	1674	1680	1686	rVB6	12911	21919	0.39%	0.073%
33	13.146	1711	1720	1727	rBV4	53660	112722	2.02%	0.378%
34	13.222	1727	1733	1740	rVV	42835	76999	1.38%	0.258%
35	13.416	1760	1766	1779	rVB2	115983	226707	4.06%	0.760%
36	13.598	1791	1797	1805	rVB10	6627	18378	0.33%	0.062%

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Instrument : BNA_G
ClientSampleId : C0AT5

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

37	13.675	1805	1810	1813	rBV4	16163	30440	0.54%	0.102%
38	13.728	1813	1819	1831	rVB	86474	168656	3.02%	0.566%
39	13.839	1832	1838	1842	rBV5	24323	44435	0.80%	0.149%
40	14.110	1878	1884	1889	rBV	338801	482537	8.64%	1.618%
41	14.233	1900	1905	1908	rBV7	8237	14600	0.26%	0.049%
42	14.309	1914	1918	1924	rVB6	9135	15865	0.28%	0.053%
43	14.403	1927	1934	1945	rVB	339004	561436	10.05%	1.883%
44	14.503	1945	1951	1952	rBV6	13217	20063	0.36%	0.067%
45	14.709	1979	1986	1993	rVB	230209	373296	6.68%	1.252%
46	14.879	2010	2015	2017	rBV6	13574	23349	0.42%	0.078%
47	14.914	2017	2021	2031	rVB3	33216	60866	1.09%	0.204%
48	15.032	2033	2041	2044	rBV9	9233	19390	0.35%	0.065%
49	15.114	2049	2055	2061	rVB4	14904	27359	0.49%	0.092%
50	15.185	2061	2067	2069	rBV6	15201	21553	0.39%	0.072%
51	15.255	2074	2079	2082	rVV2	44882	70439	1.26%	0.236%
52	15.302	2082	2087	2089	rVV2	58980	103541	1.85%	0.347%
53	15.332	2089	2092	2101	rVB2	141000	224933	4.03%	0.754%
54	15.437	2106	2110	2116	rBV5	10897	16705	0.30%	0.056%
55	15.490	2116	2119	2122	rBV4	9960	15324	0.27%	0.051%
56	15.702	2148	2155	2161	rVV	478037	723686	12.96%	2.427%
57	15.761	2161	2165	2168	rVV2	24901	38170	0.68%	0.128%
58	15.837	2169	2178	2185	rVV	584692	1166099	20.88%	3.910%
59	15.913	2188	2191	2194	rVV5	16476	29284	0.52%	0.098%
60	15.949	2194	2197	2206	rVB6	25632	57014	1.02%	0.191%
61	16.048	2210	2214	2221	rVB10	13039	26675	0.48%	0.089%
62	16.119	2221	2226	2231	rBV9	14163	29855	0.53%	0.100%
63	16.442	2275	2281	2284	rBV8	9003	16754	0.30%	0.056%
64	16.607	2305	2309	2313	rVB2	21051	30697	0.55%	0.103%
65	16.736	2324	2331	2337	rBV4	17697	44243	0.79%	0.148%
66	16.947	2364	2367	2375	rVB5	8488	17089	0.31%	0.057%
67	17.118	2388	2396	2411	rVV	3371152	5585678	100.00%	18.729%
68	17.235	2413	2416	2423	rVB6	37340	68476	1.23%	0.230%
69	17.464	2449	2455	2465	rVB	289457	448673	8.03%	1.504%
70	17.564	2465	2472	2478	rBV2	458227	731775	13.10%	2.454%
71	17.641	2481	2485	2487	rBV2	29203	43724	0.78%	0.147%
72	17.817	2508	2515	2528	rBV	888867	1412566	25.29%	4.736%
73	18.117	2561	2566	2572	rVB2	44940	75541	1.35%	0.253%
74	18.263	2587	2591	2594	rBV4	12512	20603	0.37%	0.069%
75	18.305	2595	2598	2601	rVV4	14528	20512	0.37%	0.069%
76	18.346	2601	2605	2611	rVB8	21542	37762	0.68%	0.127%

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Misc :
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Instrument : BNA_G
ClientSampleId : C0AT5

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
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Title : SVOA CALIBRATION

77	18.651	2650	2657	2662	rBV2	39666	53975	0.97%	0.181%
78	19.262	2759	2761	2768	rVB7	16905	18800	0.34%	0.063%
79	19.556	2804	2811	2815	rBV	4555094	5568807	99.70%	18.673%
80	19.715	2832	2838	2852	rVV	1380065	2007319	35.94%	6.731%
81	19.850	2854	2861	2866	rBV	552670	816662	14.62%	2.738%
82	20.073	2894	2899	2904	rBV2	34536	55459	0.99%	0.186%
83	20.120	2904	2907	2913	rVB3	22338	29633	0.53%	0.099%
84	20.402	2951	2955	2960	rVB8	15868	23000	0.41%	0.077%
85	20.455	2960	2964	2966	rBV4	14055	19319	0.35%	0.065%
86	20.537	2974	2978	2985	rVB	319330	389904	6.98%	1.307%
87	20.796	3019	3022	3024	rBV4	29226	35572	0.64%	0.119%
88	20.819	3024	3026	3030	rVB3	26955	27519	0.49%	0.092%
89	21.219	3088	3094	3104	rVV	849295	1237560	22.16%	4.150%
90	21.377	3116	3121	3124	rBV5	25794	43038	0.77%	0.144%
91	21.618	3159	3162	3165	rBV2	16001	18141	0.32%	0.061%
92	21.742	3177	3183	3192	rVB	302394	512914	9.18%	1.720%
93	22.917	3376	3383	3396	rBV	305932	675913	12.10%	2.266%
94	23.416	3461	3468	3473	rBV9	20372	50452	0.90%	0.169%
95	23.734	3515	3522	3528	rVB3	24926	57972	1.04%	0.194%
96	24.080	3578	3581	3587	rVB5	11474	21962	0.39%	0.074%
97	24.809	3693	3705	3716	rBV2	312012	893176	15.99%	2.995%
98	25.032	3734	3743	3756	rVB2	181740	546893	9.79%	1.834%
99	25.385	3794	3803	3816	rBV2	100971	321016	5.75%	1.076%
100	29.303	4460	4470	4472	rBV6	35086	93574	1.68%	0.314%

Sum of corrected areas: 29823214

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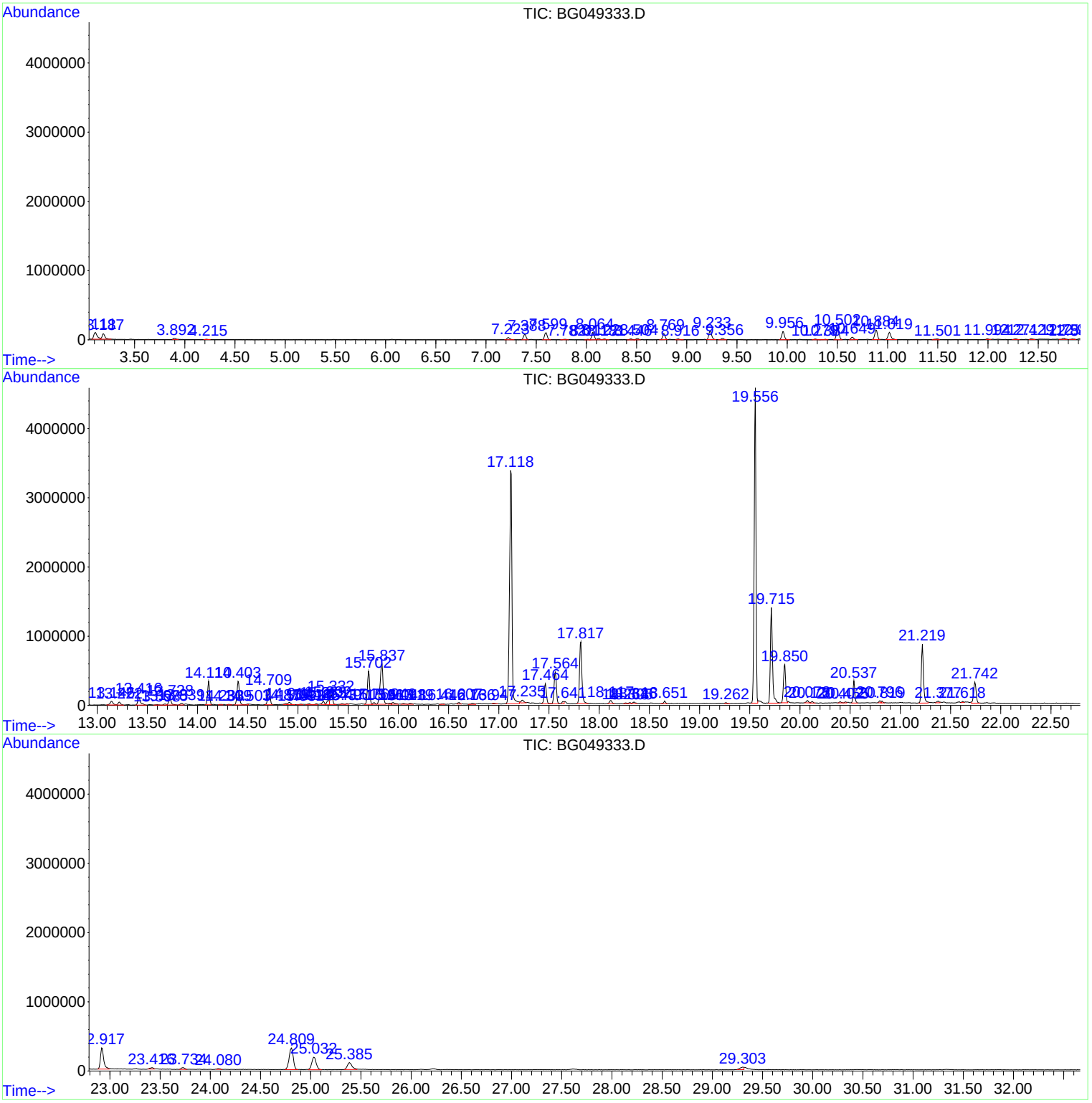
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AT5

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 : BG049333.D
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Misc :
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Instrument : BNA_G
ClientSampleId : C0AT5

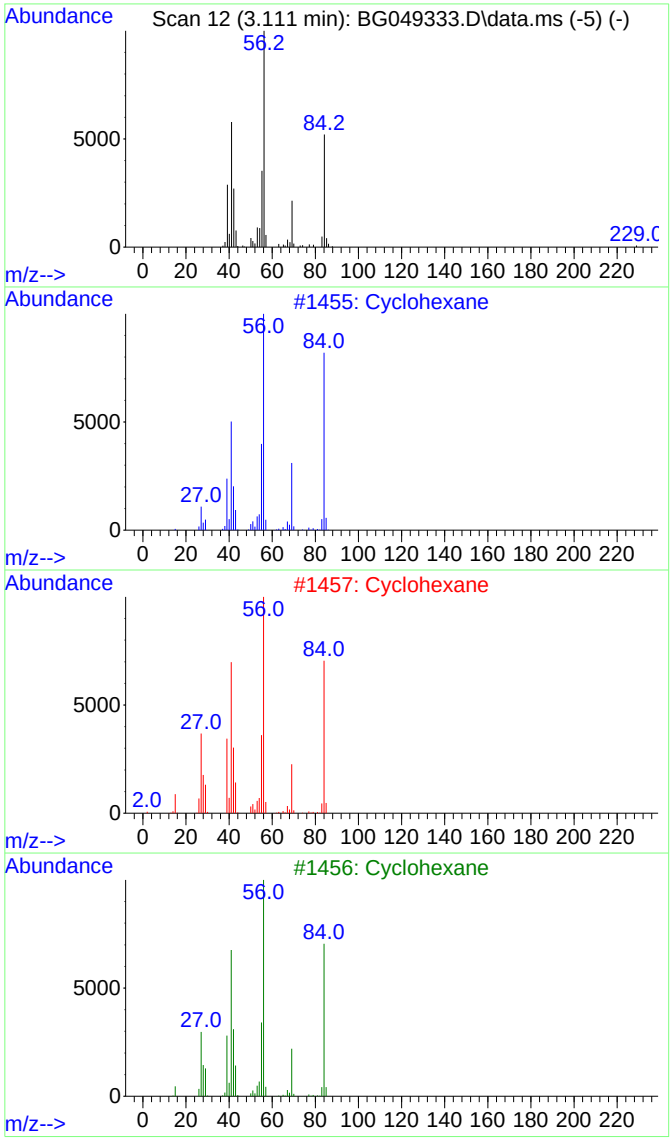
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

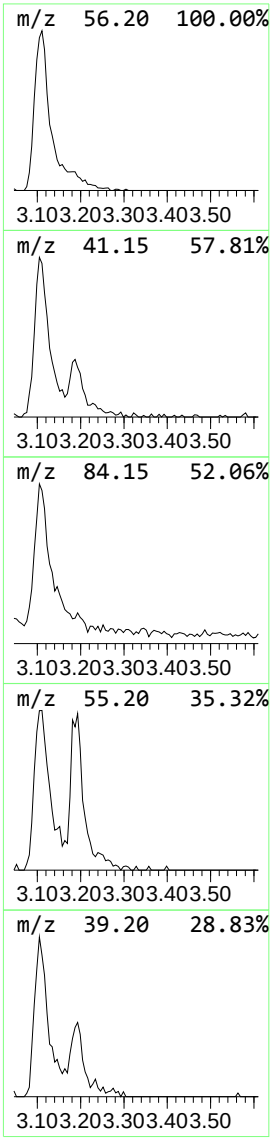
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Cyclic3.11 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.110	24.43 ng/ul	234428	1,4-Dichlorobenzene-d4	8.064

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane	84	C6H12	000110-82-7	91
2	Cyclohexane	84	C6H12	000110-82-7	90
3	Cyclohexane	84	C6H12	000110-82-7	90
4	2-Amino-oxazole	84	C3H4N2O	004570-45-0	87
5	Cyclohexane	84	C6H12	000110-82-7	83





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Acq On : 14 Jul 2021 17:29
Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AT5

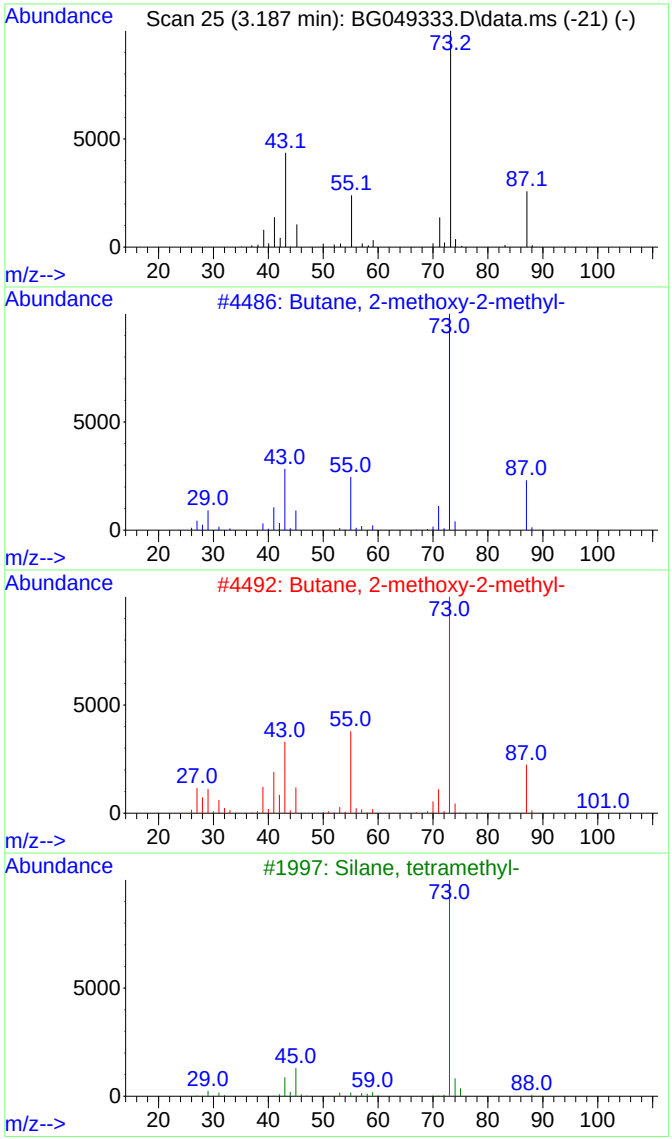
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

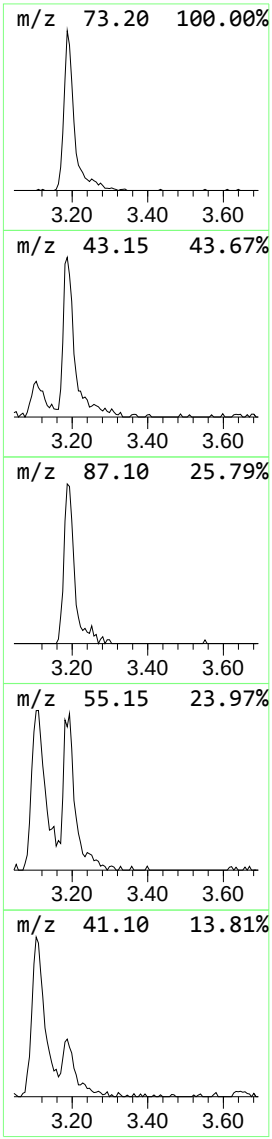
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.190	16.70 ng/ul	160234	1,4-Dichlorobenzene-d4	8.064

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	50
3	Silane, tetramethyl-	88	C4H12Si	000075-76-3	25
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
5	Silane, tetramethyl-	88	C4H12Si	000075-76-3	10





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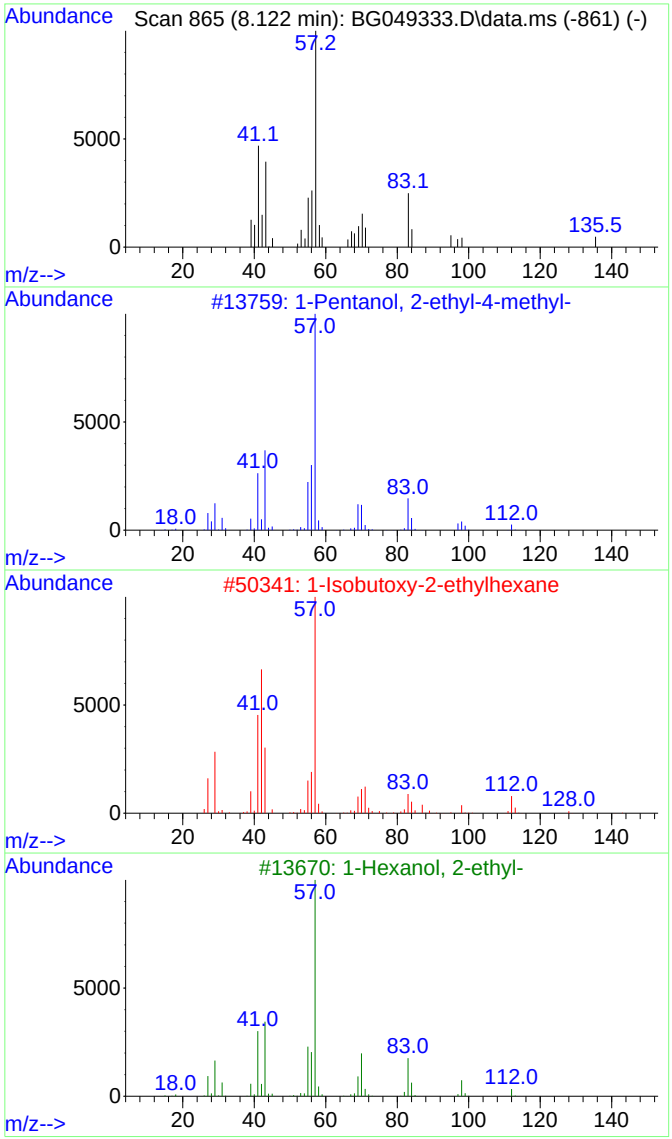
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

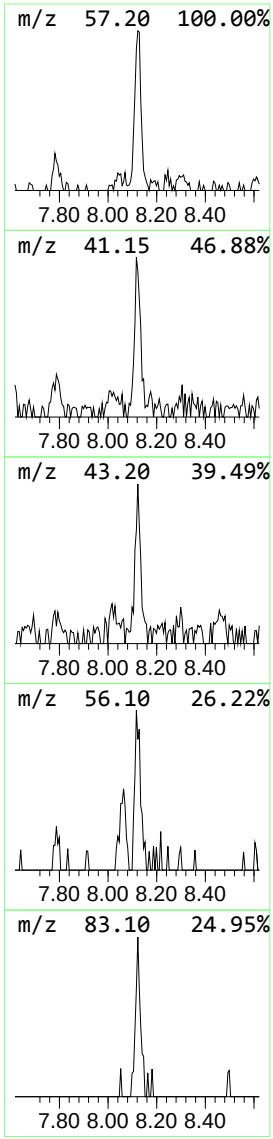
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1-Pentanol, 2-ethyl-4-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.120	3.73 ng/ul	35772	1,4-Dichlorobenzene-d4	8.064

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	64
2	1-Isobutoxy-2-ethylhexane	186	C12H26O	1000139-90-3	59
3	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	42
4	Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	40
5	Formic acid, 2-propylpentyl ester	158	C9H18O2	1000368-74-8	40





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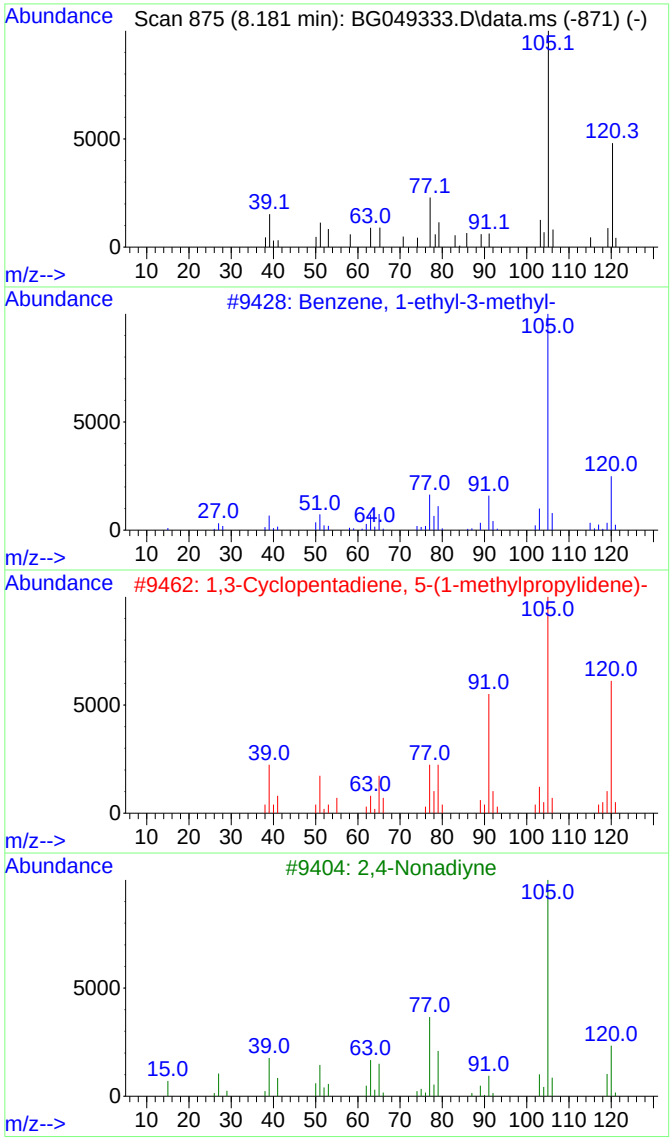
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

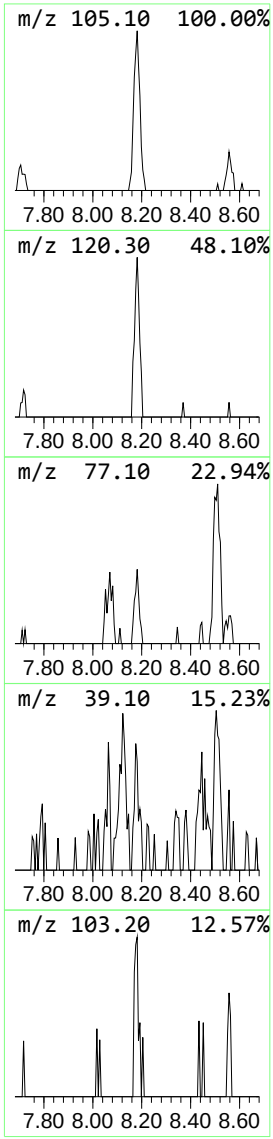
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1-ethyl-3-methyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.180	2.28 ng/ul	21850	1,4-Dichlorobenzene-d4	8.064

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	87
2	1,3-Cyclopentadiene, 5-(1-methyl...	120	C9H12	003141-02-4	86
3	2,4-Nonadiyne	120	C9H12	063621-15-8	83
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	80
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	80





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Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

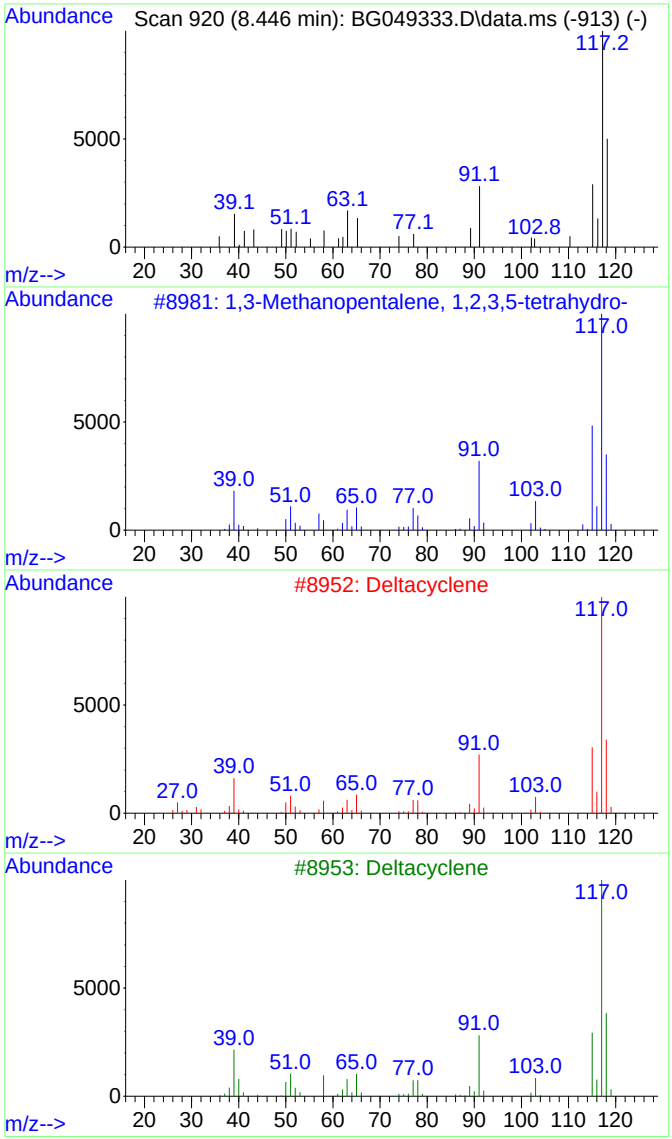
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

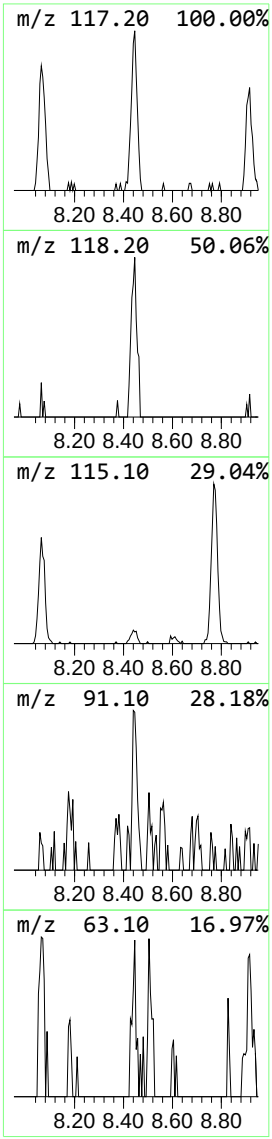
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 1,3-Methanopentalene, 1,2,3... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.450	2.65 ng/ul	25398	1,4-Dichlorobenzene-d4	8.064

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Methanopentalene, 1,2,3,5-te...	118	C9H10	128600-88-4	64
2	Deltacyclene	118	C9H10	007785-10-6	64
3	Deltacyclene	118	C9H10	007785-10-6	64
4	Indolizine	117	C8H7N	000274-40-8	43
5	5H-1-Pyrindine	117	C8H7N	000270-91-7	43





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Data File : BG049333.D
Acq On : 14 Jul 2021 17:29
Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AT5

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

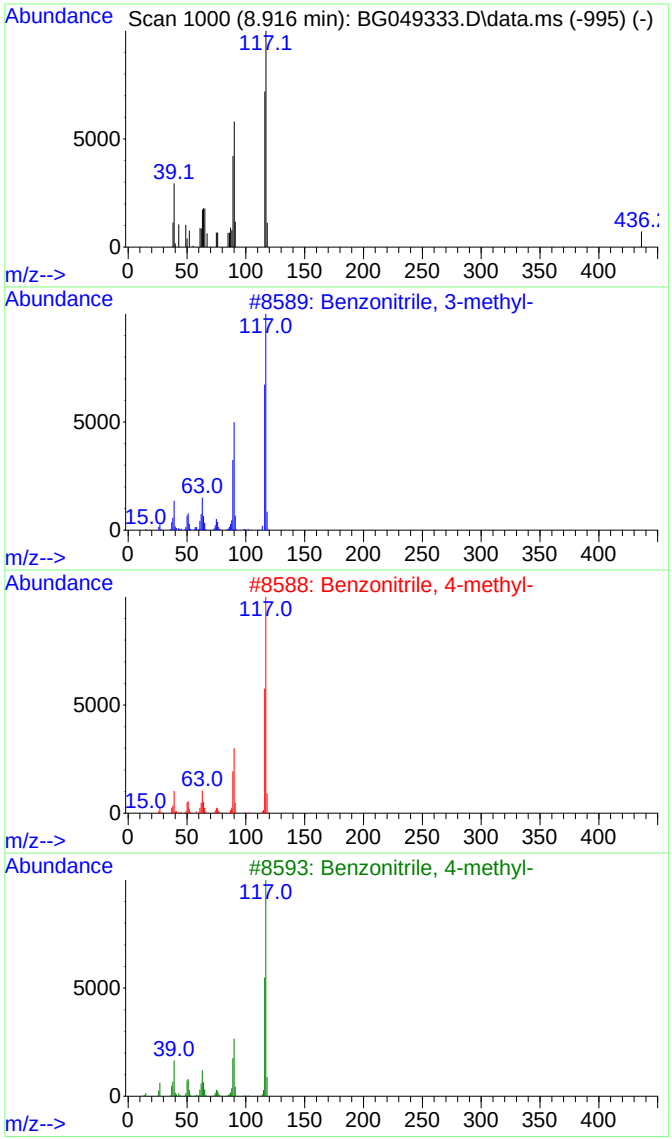
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

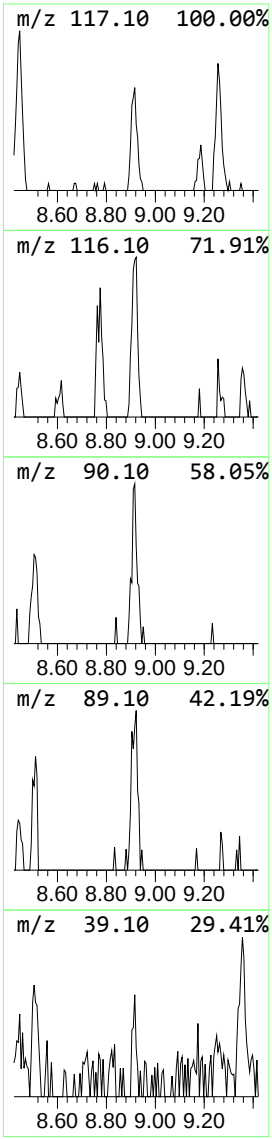
****

Peak Number 6 Benzonitrile, 3-methyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.920	2.47 ng/ul	23697	1,4-Dichlorobenzene-d4	8.064

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzonitrile, 3-methyl-	117	C8H7N	000620-22-4	93
2	Benzonitrile, 4-methyl-	117	C8H7N	000104-85-8	93
3	Benzonitrile, 4-methyl-	117	C8H7N	000104-85-8	93
4	Benzonitrile, 4-methyl-	117	C8H7N	000104-85-8	90
5	Benzonitrile, 3-methyl-	117	C8H7N	000620-22-4	90





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Acq On : 14 Jul 2021 17:29
Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

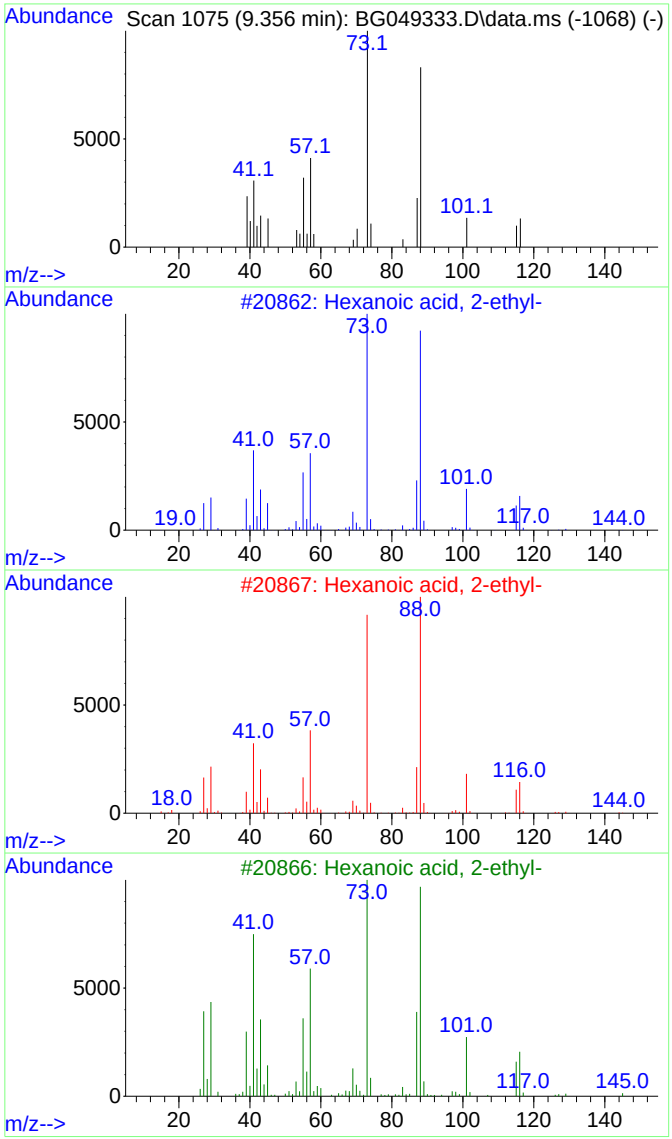
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

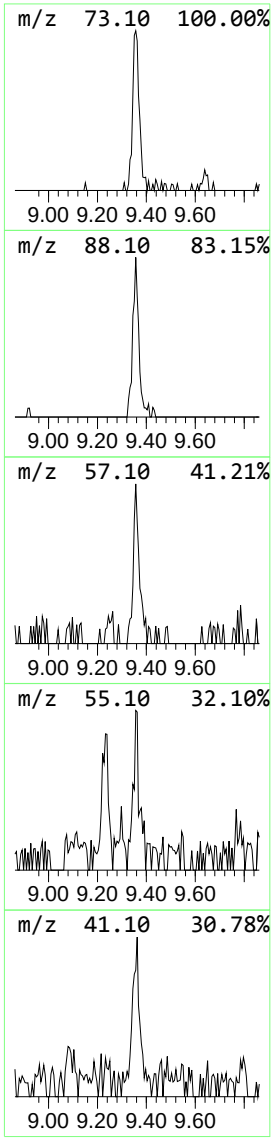
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Hexanoic acid, 2-ethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.360	4.07 ng/ul	39012	1,4-Dichlorobenzene-d4	8.064

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	72
2	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	50
3	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	40
4	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	12
5	Hydrazine, 1-butyl-1-ethyl-	116	C6H16N2	052728-56-0	9





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Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

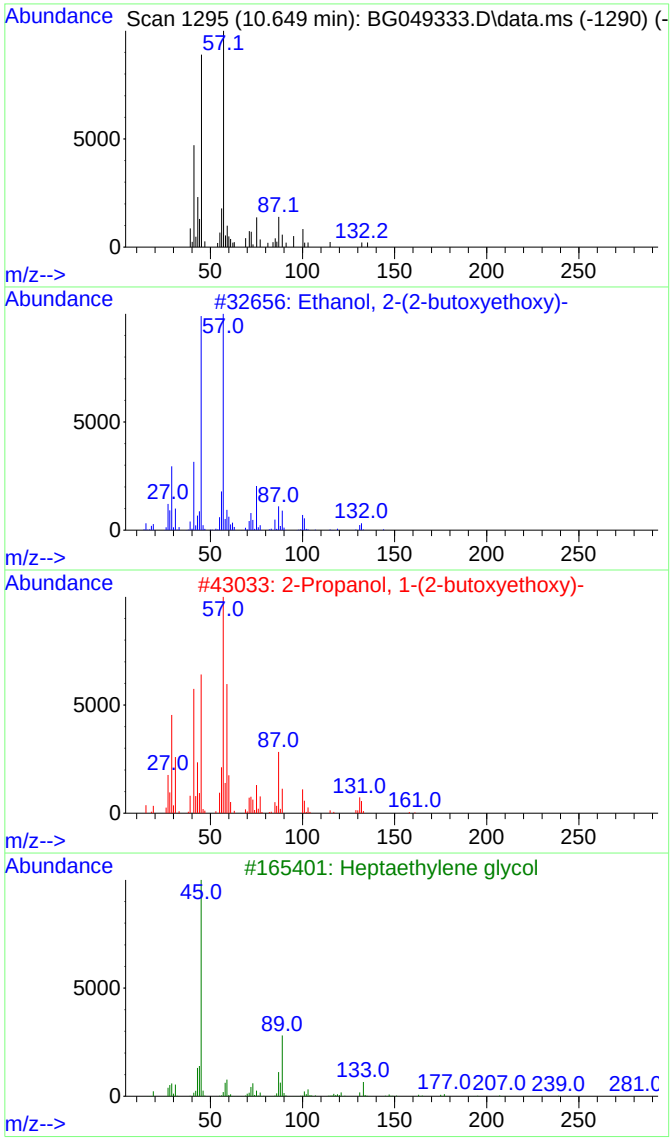
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

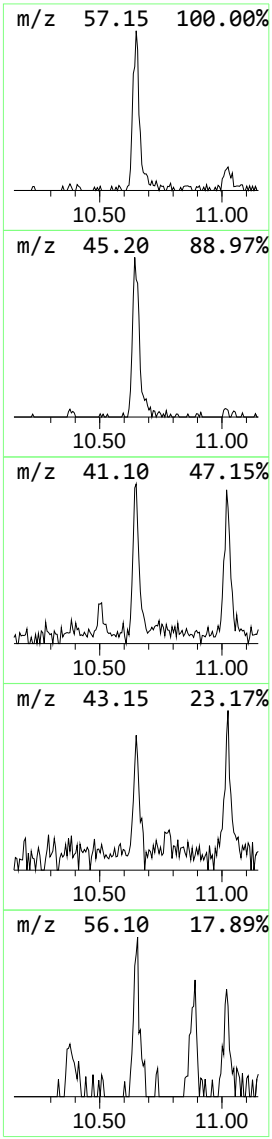
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.650	5.20 ng/ul	67363	Naphthalene-d8	10.884

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	56
2	2-Propanol, 1-(2-butoxyethoxy)-	176	C9H20O3	000124-16-3	56
3	Heptaethylene glycol	326	C14H30O8	005617-32-3	50
4	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	50
5	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	47





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Acq On : 14 Jul 2021 17:29
Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

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

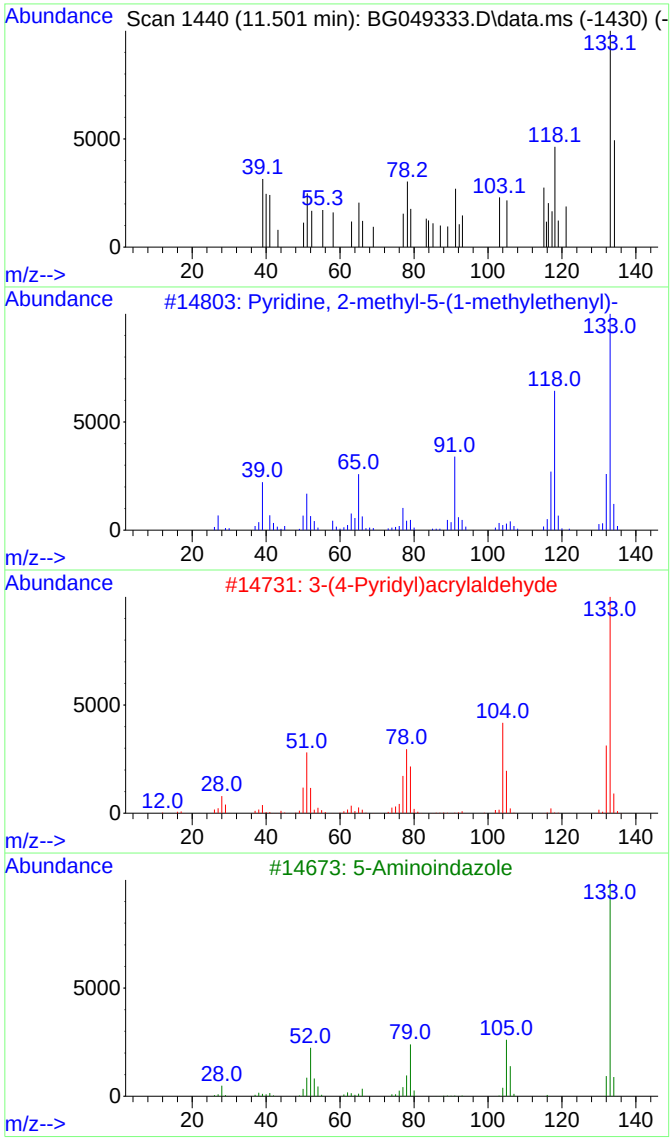
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

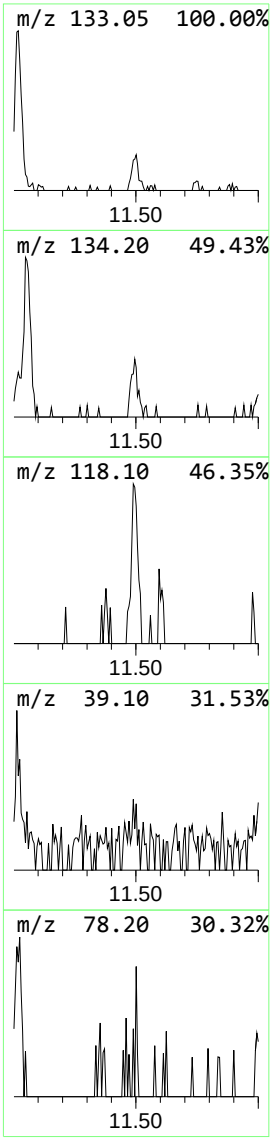
****

Peak Number 9 unknown-01 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.500	2.56 ng/ul	33233	Naphthalene-d8	10.884

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyridine, 2-methyl-5-(1-methylet...	133	C9H11N	056057-93-3	46
2	3-(4-Pyridyl)acrylaldehyde	133	C8H7NO	026505-36-2	30
3	5-Aminoindazole	133	C7H7N3	019335-11-6	25
4	[1,2,4]Triazolo[1,5-a]pyridine, ...	133	C7H7N3	000768-19-4	25
5	Benzene, 1-isocyanato-3-methyl-	133	C8H7NO	000621-29-4	25





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Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AT5

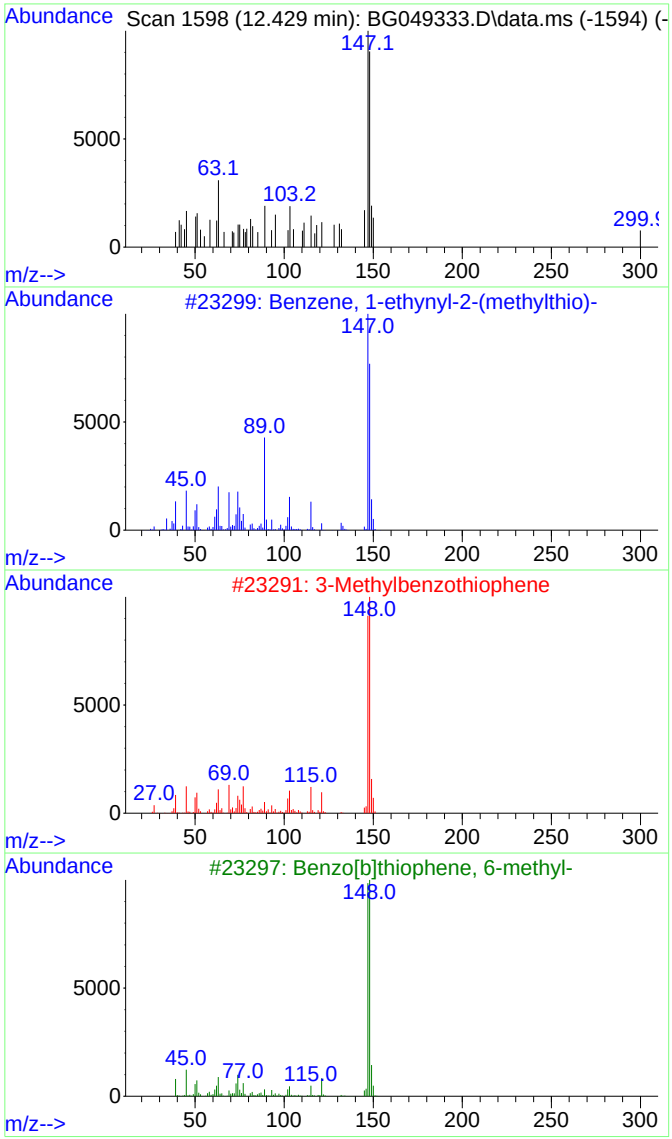
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

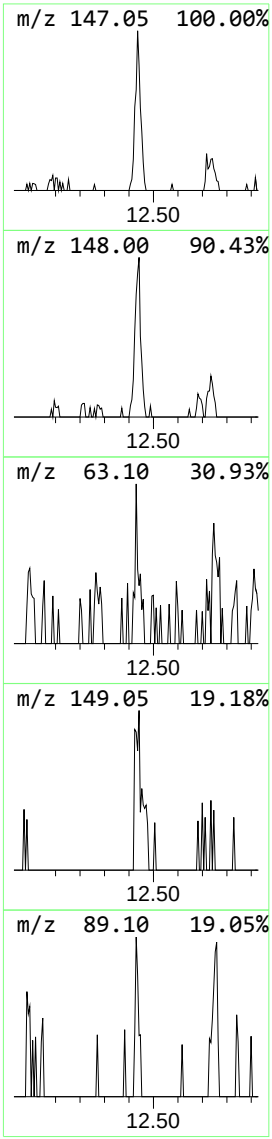
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1-ethynyl-2-(methy... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.430	2.21 ng/ul	28655	Naphthalene-d8	10.884

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethynyl-2-(methylthio)-	148	C9H8S	078905-08-5	64
2	3-Methylbenzothiophene	148	C9H8S	001455-18-1	62
3	Benzo[b]thiophene, 6-methyl-	148	C9H8S	016587-47-6	50
4	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	47
5	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	47





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Acq On : 14 Jul 2021 17:29
Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

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

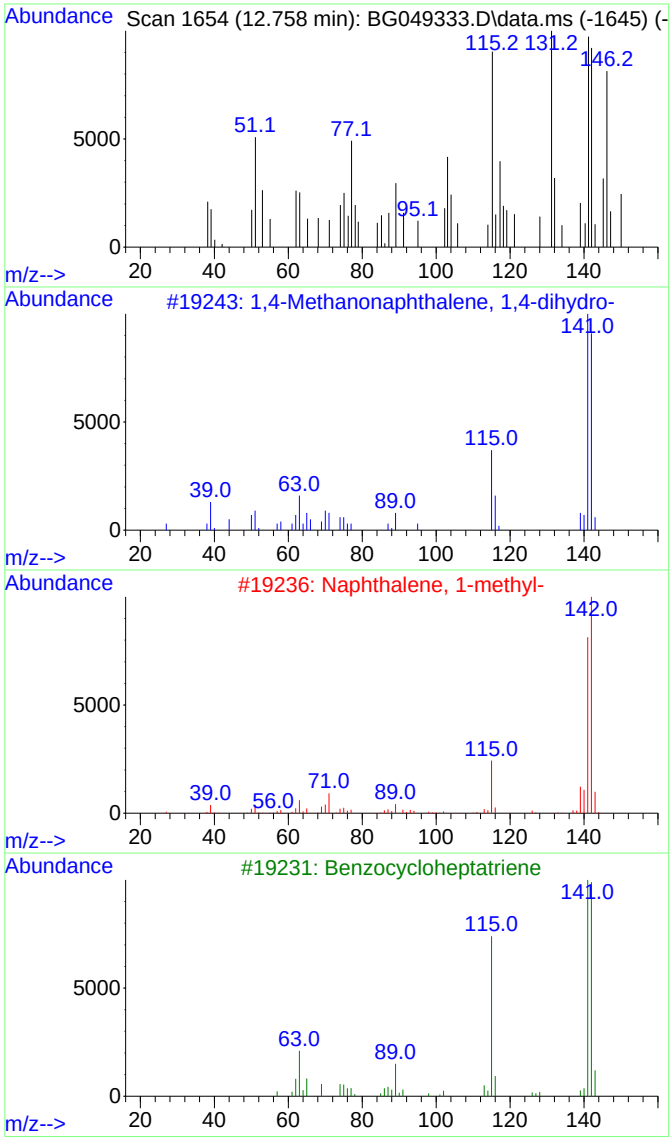
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

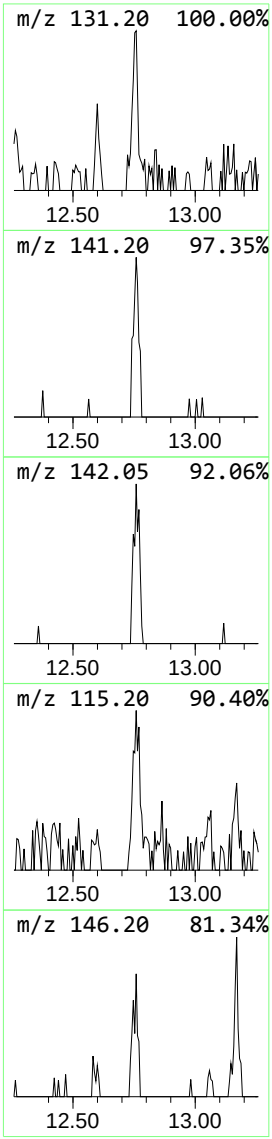
****

Peak Number 11 unknown-02 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.760	3.02 ng/ul	39111	Naphthalene-d8	10.884

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	38
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	30
3	Benzocycloheptatriene	142	C11H10	000264-09-5	30
4	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	30
5	1-Methylindan-2-one	146	C10H10O	035587-60-1	30





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Acq On : 14 Jul 2021 17:29
Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

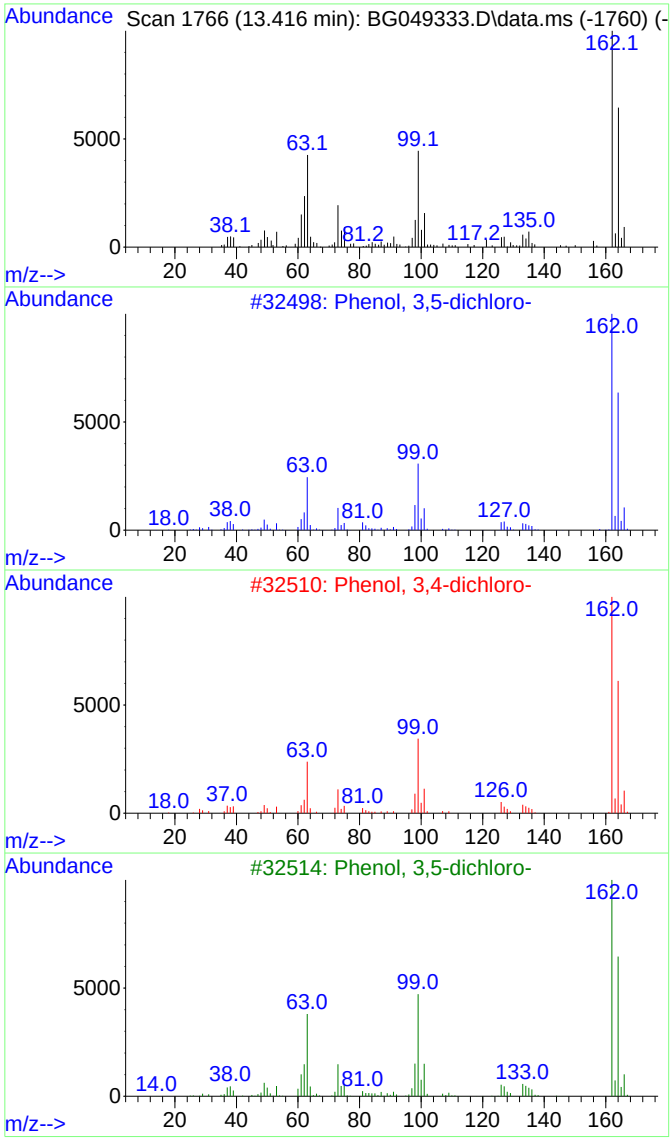
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

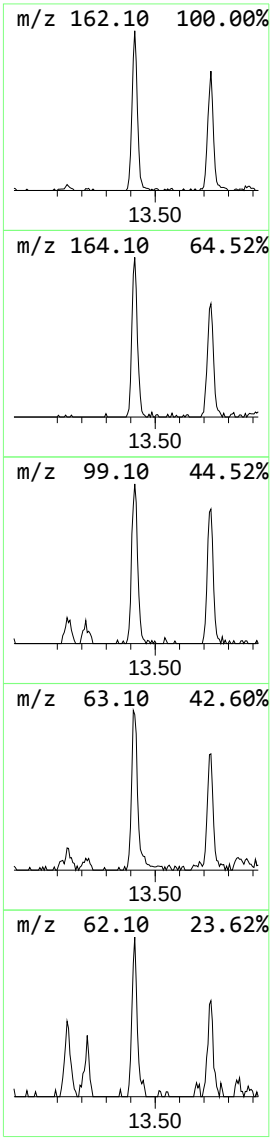
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Phenol, 3,5-dichloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.420	12.15 ng/ul	226707	Acenaphthene-d10	14.709

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	95
2	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	95
3	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	95
4	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	93
5	Phenol, 2,5-dichloro-	162	C6H4Cl2O	000583-78-8	90





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Acq On : 14 Jul 2021 17:29
Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

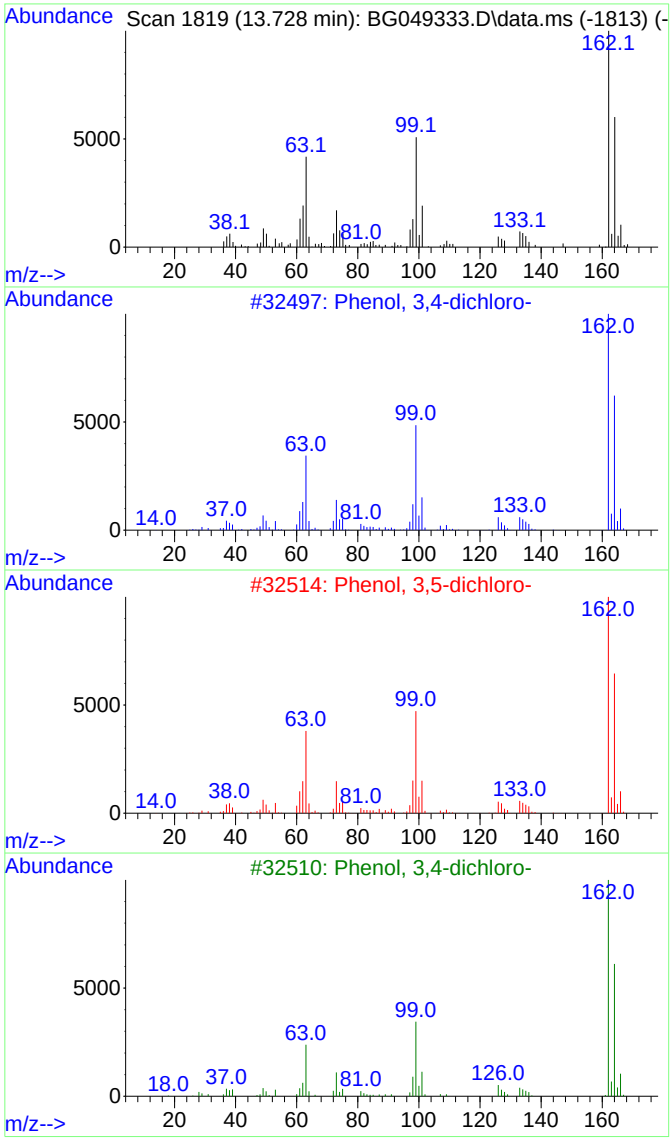
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

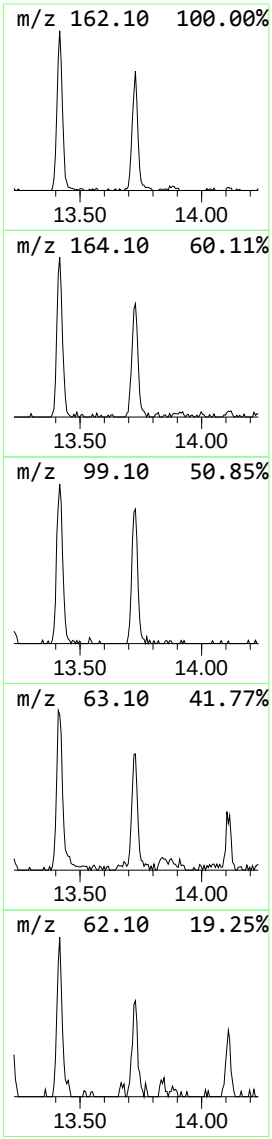
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Phenol, 3,4-dichloro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.730	9.04 ng/ul	168656	Acenaphthene-d10	14.709

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97
2	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	97
3	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	95
4	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	95
5	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	87





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Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

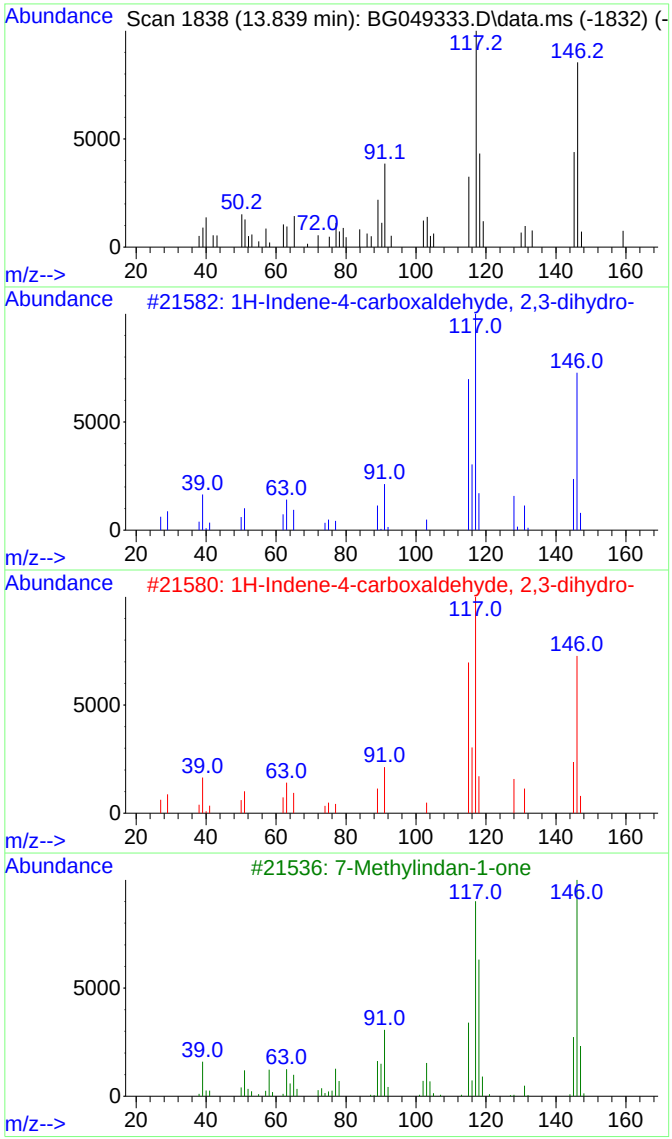
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

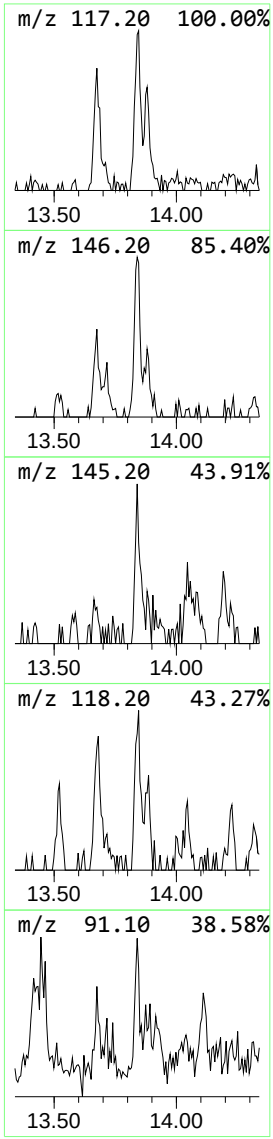
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 1H-Indene-4-carboxaldehyde,... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.840	2.38 ng/ul	44435	Acenaphthene-d10	14.709

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene-4-carboxaldehyde, 2,3-...	146	C10H10O	051932-70-8	58
2	1H-Indene-4-carboxaldehyde, 2,3-...	146	C10H10O	051932-70-8	58
3	7-Methylindan-1-one	146	C10H10O	039627-61-7	49
4	1,3-Methanopentalene, 1,2,3,5-te...	118	C9H10	128600-88-4	38
5	4H-1-Benzopyran-4-one	146	C9H6O2	000491-38-3	35





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Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AT5

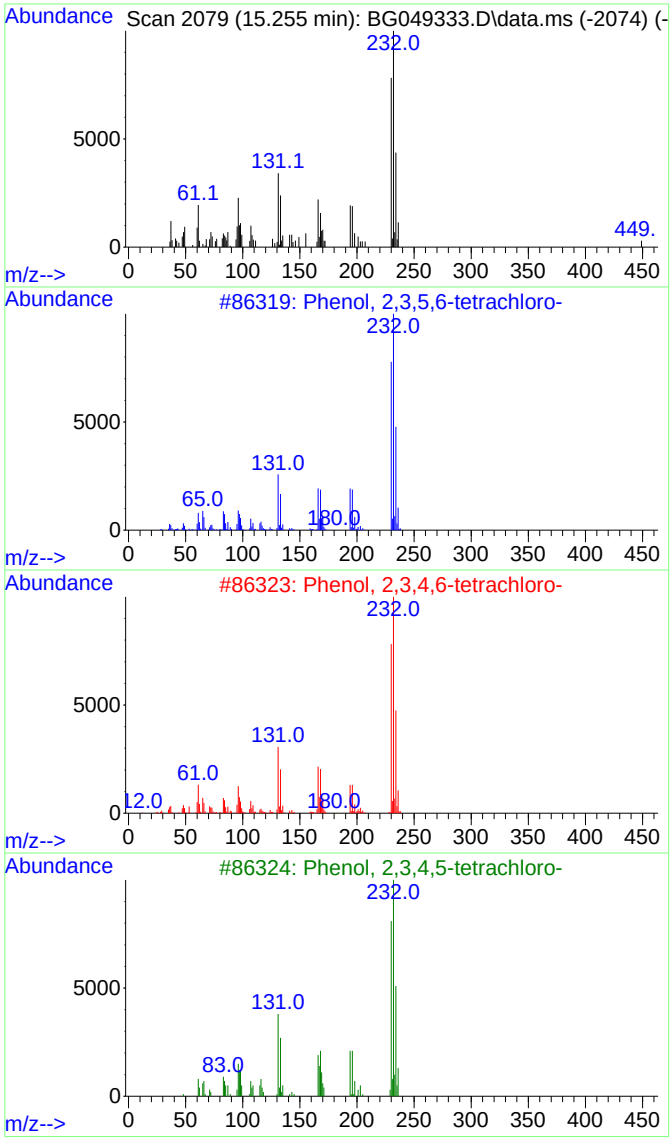
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

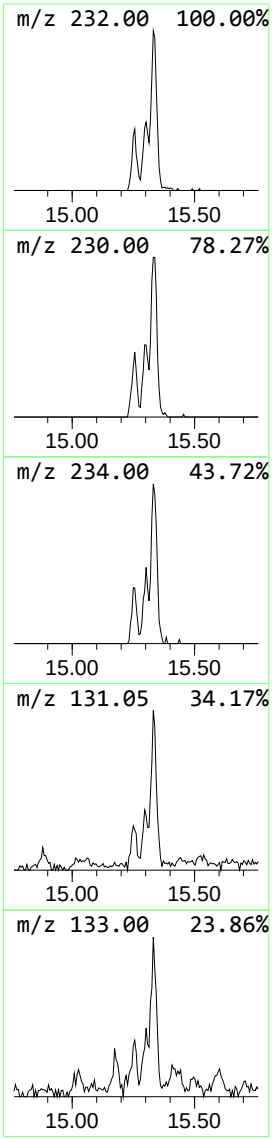
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Phenol, 2,3,5,6-tetrachloro- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.260	3.77 ng/ul	70439	Acenaphthene-d10	14.709

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	99
2	Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	97
3	Phenol, 2,3,4,5-tetrachloro-	230	C6H2Cl4O	004901-51-3	96
4	Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	95
5	Phenol, 2,3,4,5-tetrachloro-	230	C6H2Cl4O	004901-51-3	95





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ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

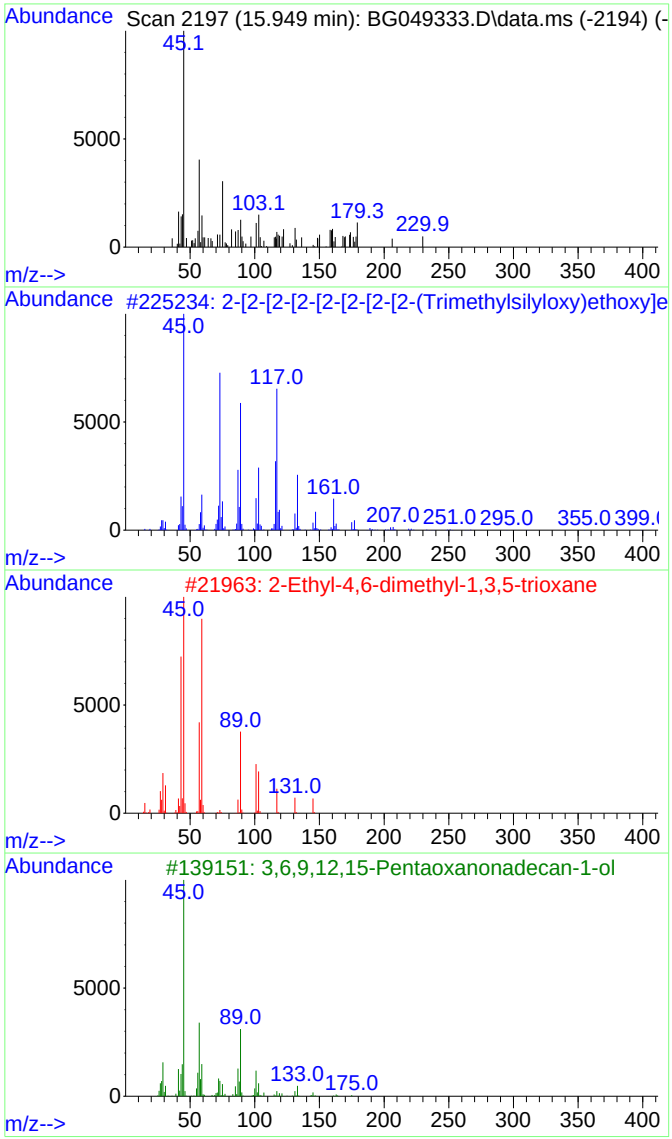
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

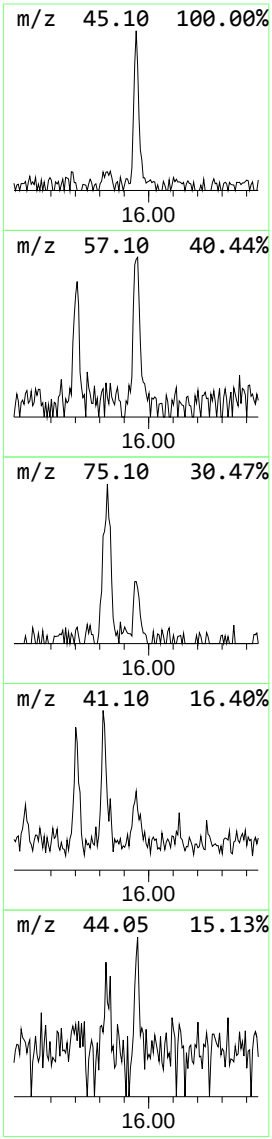
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 2-[2-[2-[2-[2-[2-[2-[2-(Tri... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.950	3.05 ng/ul	57014	Acenaphthene-d10	14.709

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-[2-[2-[2-[2-[2-[2-[2-(Trimethy...	442	C19H42O9Si	1000352-06-4	50
2	2-Ethyl-4,6-dimethyl-1,3,5-trioxane	146	C7H14O3	024261-86-7	47
3	3,6,9,12,15-Pentaoxanonadecan-1-ol	294	C14H30O6	001786-94-3	47
4	tert-Butyldimethylsilyl 2-(2-eth...	262	C12H26O4Si	1000366-92-5	45
5	Methoxyacetic acid, 2-butyl ester	146	C7H14O3	070159-90-9	43





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Acq On : 14 Jul 2021 17:29
Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

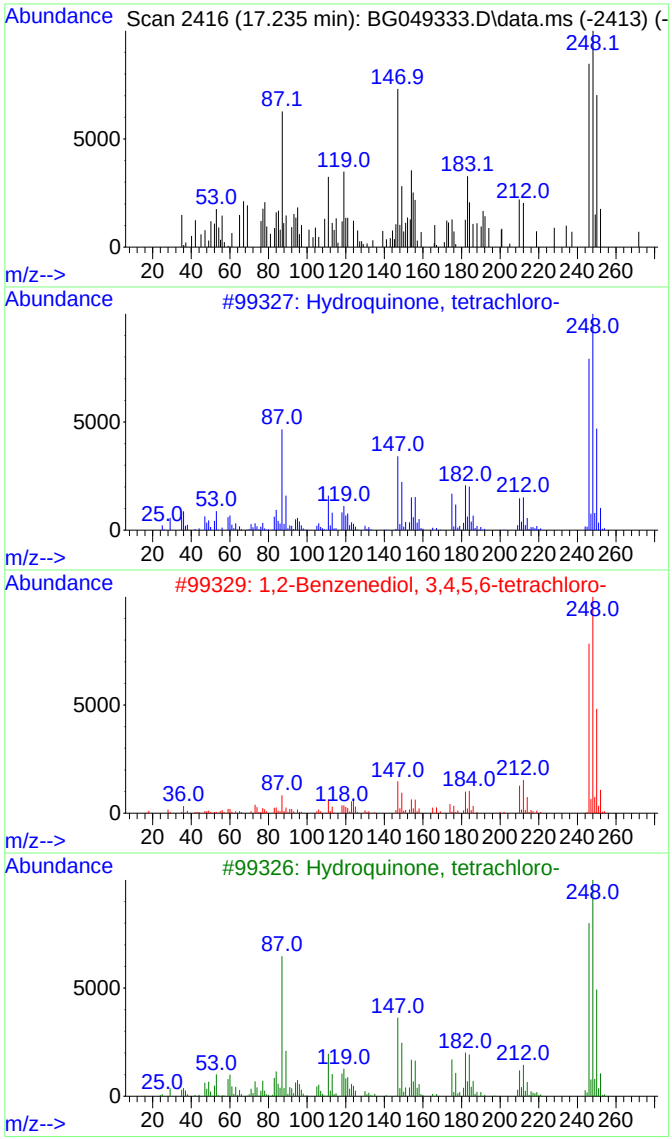
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

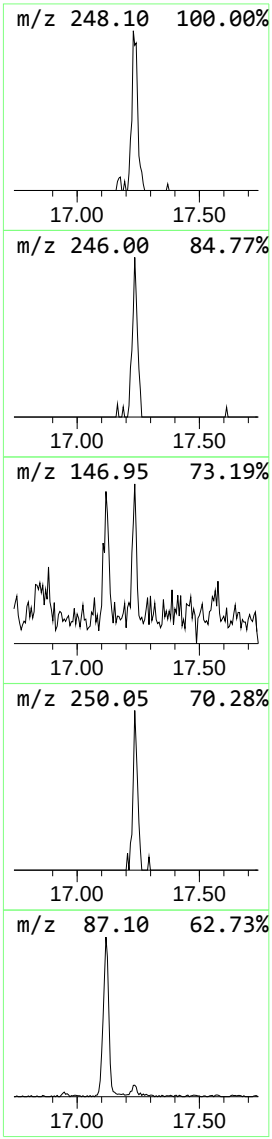
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Hydroquinone, tetrachloro- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.240	3.05 ng/ul	68476	Phenanthrene-d10	17.464

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hydroquinone, tetrachloro-	246	C6H2Cl4O2	000087-87-6	93
2	1,2-Benzenediol, 3,4,5,6-tetrach...	246	C6H2Cl4O2	001198-55-6	91
3	Hydroquinone, tetrachloro-	246	C6H2Cl4O2	000087-87-6	90
4	Hydroquinone, tetrachloro-	246	C6H2Cl4O2	000087-87-6	52
5	5-Bromo-3H-1,2-benzodithiol-3-one	246	C7H3BrOS2	115475-20-2	30





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Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

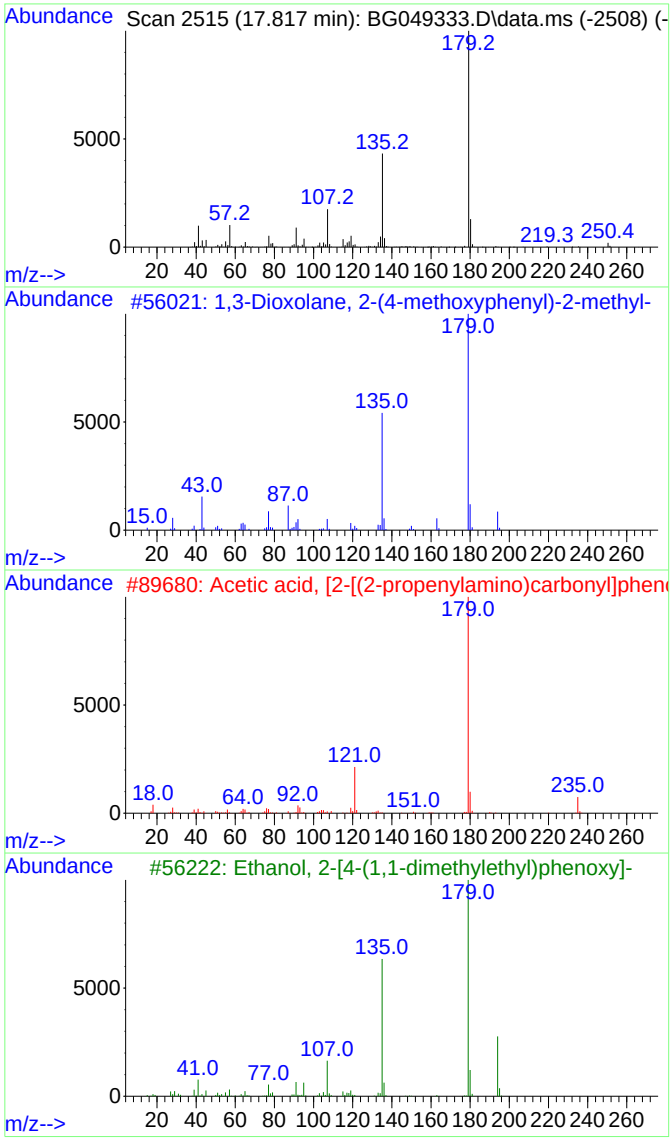
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

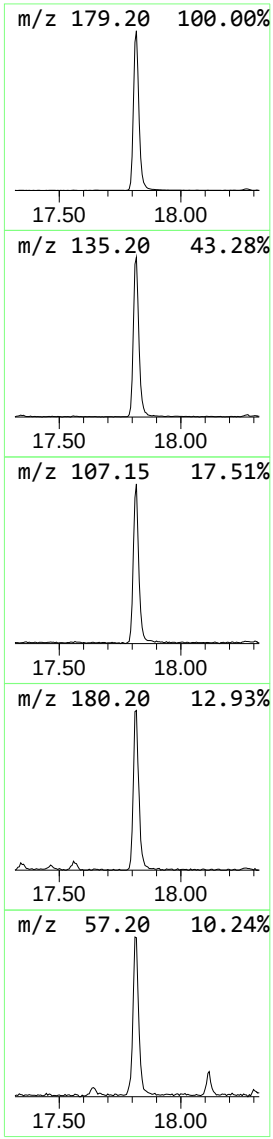
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 1,3-Dioxolane, 2-(4-methoxy... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.820	62.97 ng/ul	1412570	Phenanthrene-d10	17.464

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dioxolane, 2-(4-methoxypheny...	194	C11H14O3	036881-00-2	56
2	Acetic acid, [2-[(2-propenylamin...	235	C12H13NO4	000119-45-9	43
3	Ethanol, 2-[4-(1,1-dimethylethyl...	194	C12H18O2	000713-46-2	40
4	2-(4-Formyl-phenoxy)-acetamide	179	C9H9NO3	135857-20-4	38
5	Ethanol, 2-[4-(1,1-dimethylpropy...	208	C13H20O2	006382-07-6	33





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Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

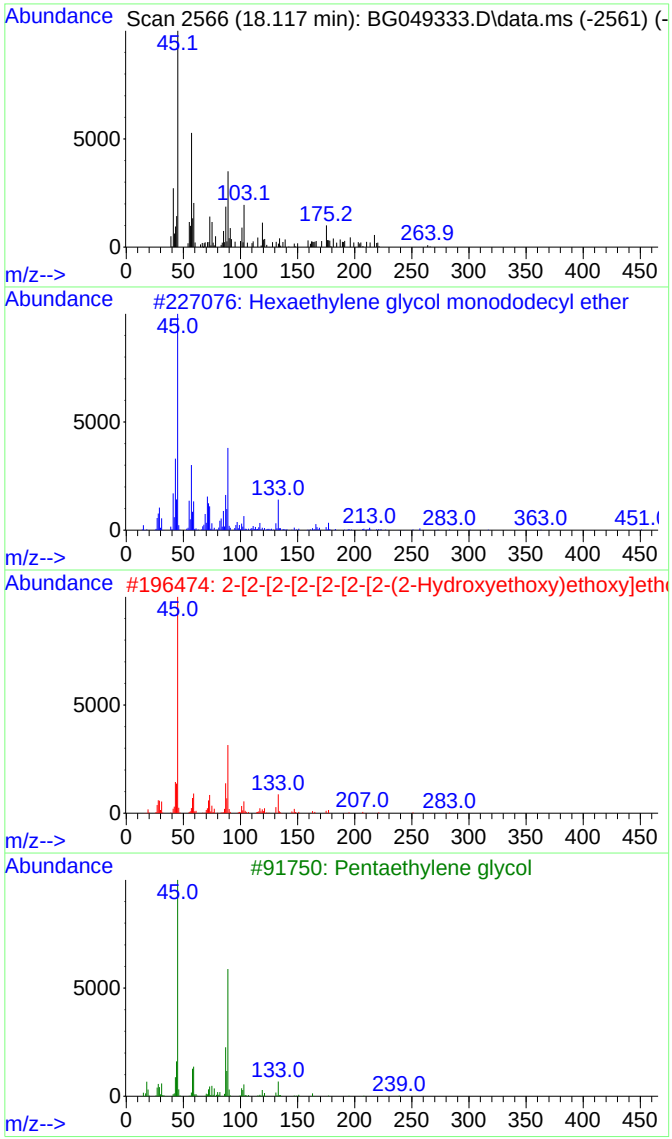
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

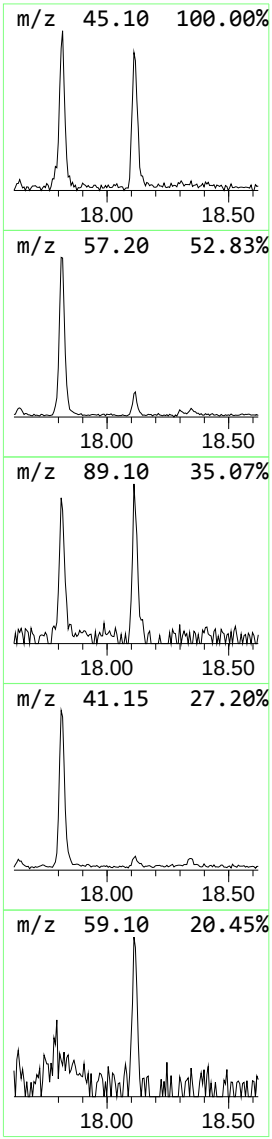
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Hexaethylene glycol monodod... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.120	3.37 ng/ul	75541	Phenanthrene-d10	17.464

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexaethylene glycol monododecyl ...	450	C24H50O7	003055-96-7	53
2	2-[2-[2-[2-[2-[2-(2-Hydroxyet...	370	C16H34O9	005117-19-1	53
3	Pentaethylene glycol	238	C10H22O6	004792-15-8	50
4	Tetraethylene glycol	194	C8H18O5	000112-60-7	50
5	Ethanol, 2-[2-(2-ethoxyethoxy)et...	178	C8H18O4	000112-50-5	50





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Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

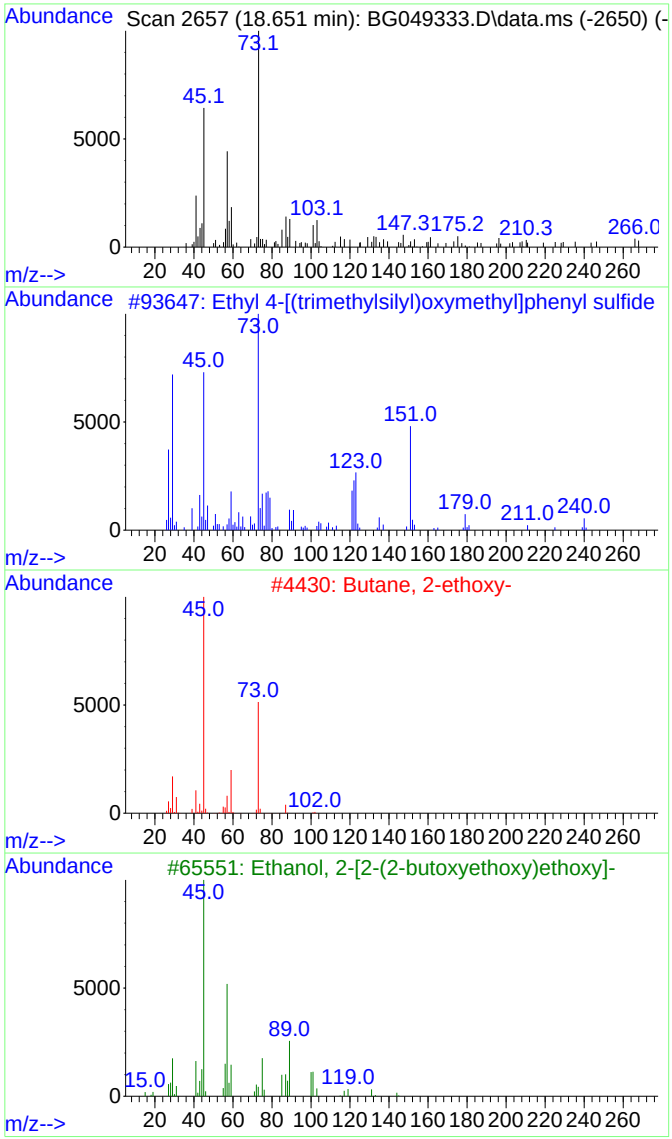
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

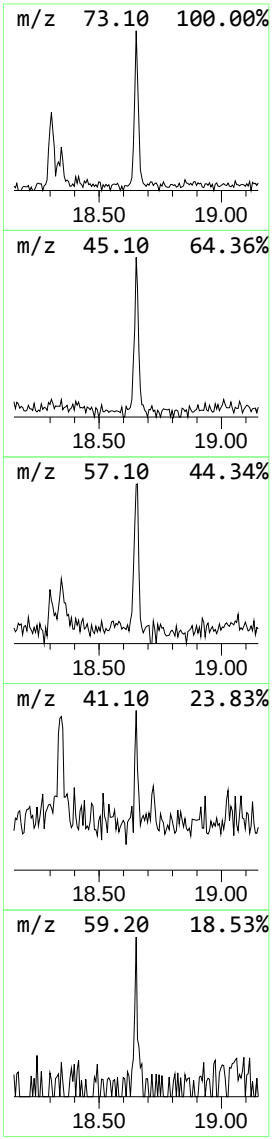
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Ethyl 4-[(trimethylsilyl)ox... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.650	2.41 ng/ul	53975	Phenanthrene-d10	17.464

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethyl 4-[(trimethylsilyl)oxymeth...	240	C12H20OSSi	1000100-02-3	50
2	Butane, 2-ethoxy-	102	C6H14O	002679-87-0	50
3	Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	47
4	3,6,9,12-Tetraoxahexadecan-1-ol	250	C12H26O5	001559-34-8	47
5	11,13-Dihydroxy-tetradec-5-enoic...	272	C15H28O4	1000193-81-4	42





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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

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

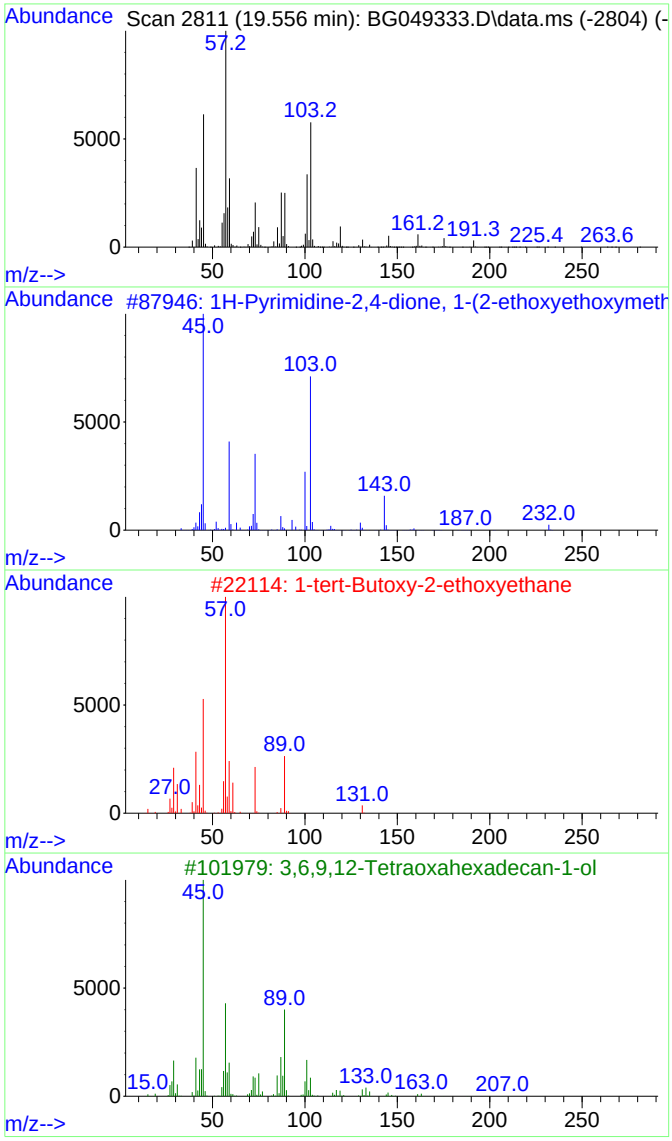
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

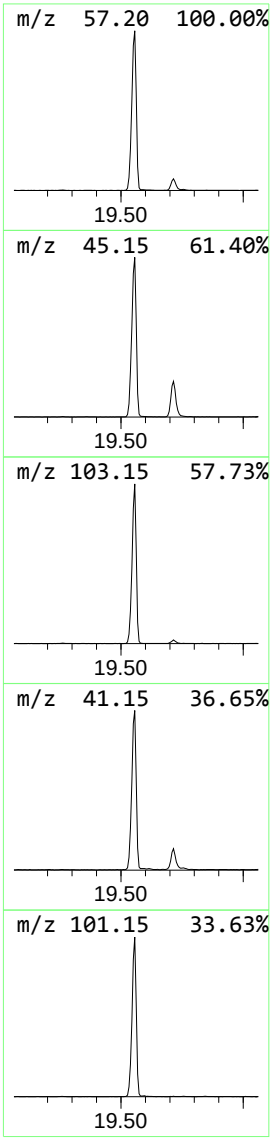
****

Peak Number 21 unknown-03 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.560	248.24 ng/ul	5568810	Phenanthrene-d10	17.464

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Pyrimidine-2,4-dione, 1-(2-et...	232	C9H13FN2O4	1000310-13-7	43
2	1-tert-Butoxy-2-ethoxyethane	146	C8H18O2	051422-54-9	43
3	3,6,9,12-Tetraoxahexadecan-1-ol	250	C12H26O5	001559-34-8	38
4	Di-sec-butyl ether	130	C8H18O	006863-58-7	35
5	3,6,9,12,15-Pentaoxanonadecan-1-ol	294	C14H30O6	001786-94-3	27





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Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

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

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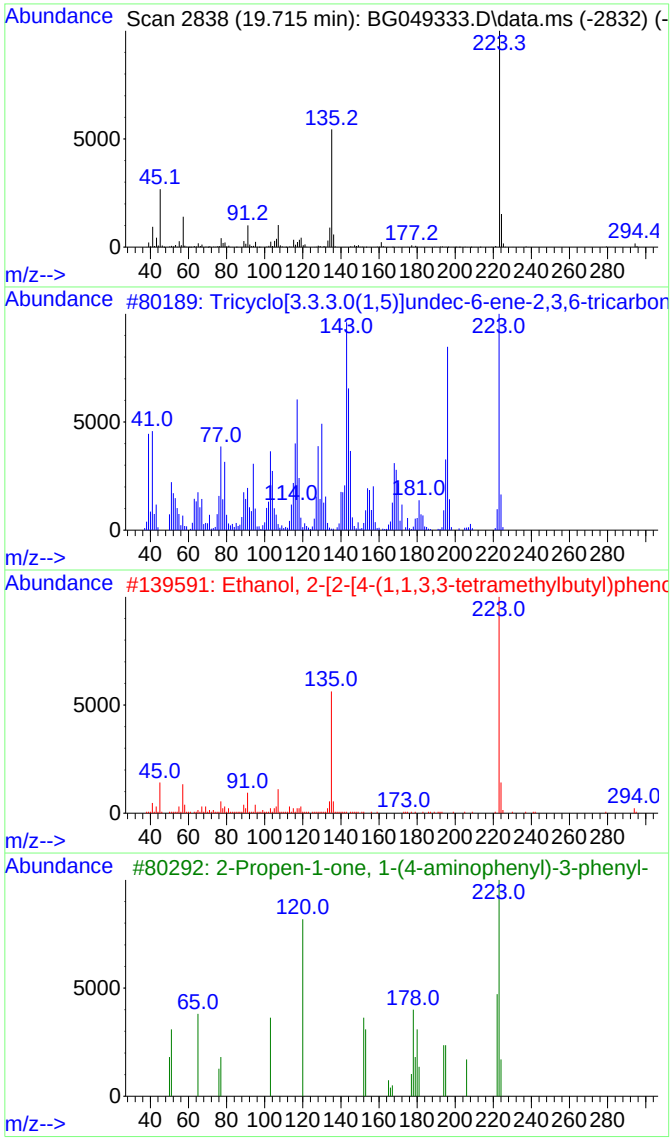
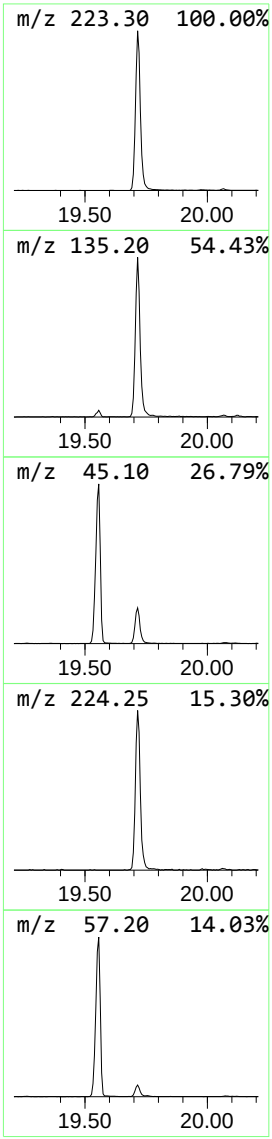
Peak Number 22 Tricyclo[3.3.3.0(1,5)]undec... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.710	78.27 ng/ul	2007320	Chrysene-d12	21.742

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricyclo[3.3.3.0(1,5)]undec-6-en...	223	C14H13N3	118894-71-6	86
2	Ethanol, 2-[2-[4-(1,1,3,3-tetram...	294	C18H30O3	002315-61-9	55
3	2-Propen-1-one, 1-(4-aminophenyl...	223	C15H13NO	002403-30-7	43
4	9-Acridinecarboxylic acid	223	C14H9NO2	005336-90-3	38
5	Benzenamine, 3-(5-methyl-1H-1,3-...	223	C14H13N3	1000319-43-9	38





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Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

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

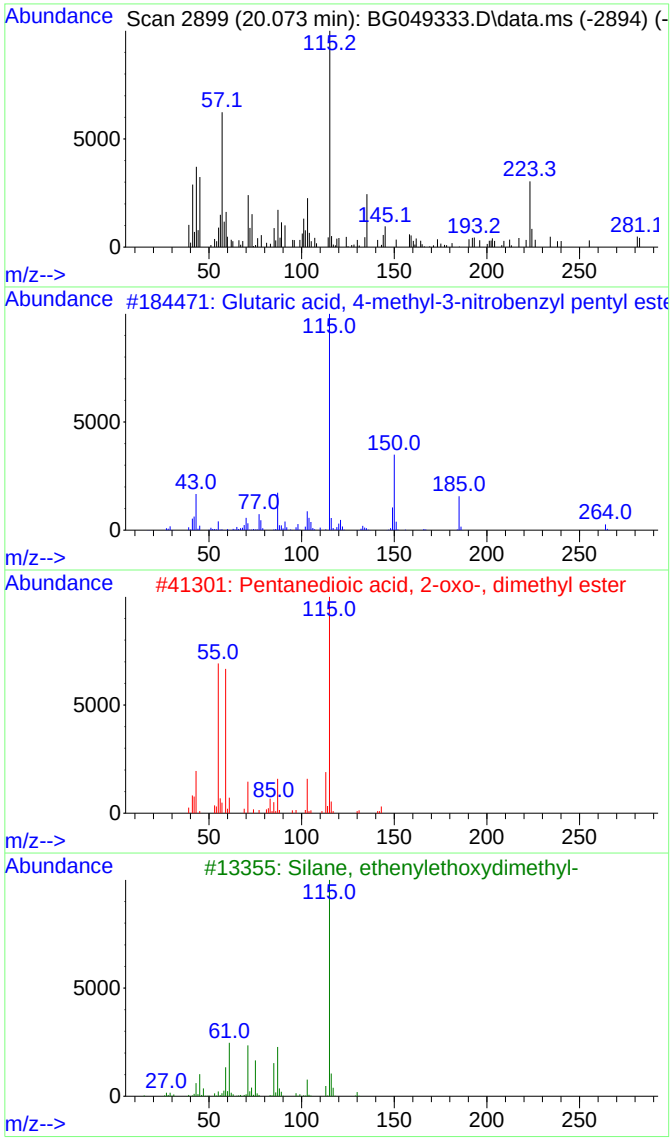
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

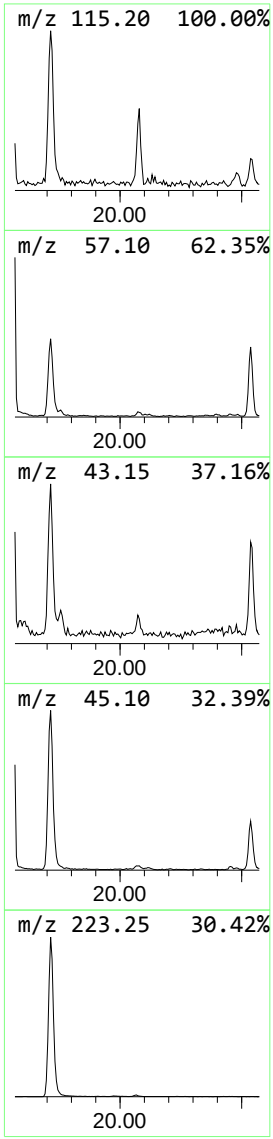
****

Peak Number 23 unknown-04 Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.070	2.16 ng/ul	55459	Chrysene-d12	21.742

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Glutaric acid, 4-methyl-3-nitrob...	351	C18H25NO6	1000376-78-1	38
2	Pentanedioic acid, 2-oxo-, dimet...	174	C7H10O5	013192-04-6	38
3	Silane, ethenylethoxydimethyl-	130	C6H14OSi	005356-83-2	38
4	Silane, ethenylethoxydimethyl-	130	C6H14OSi	005356-83-2	38
5	Morpholine, 2,6-dimethyl-	115	C6H13NO	000141-91-3	38





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Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

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

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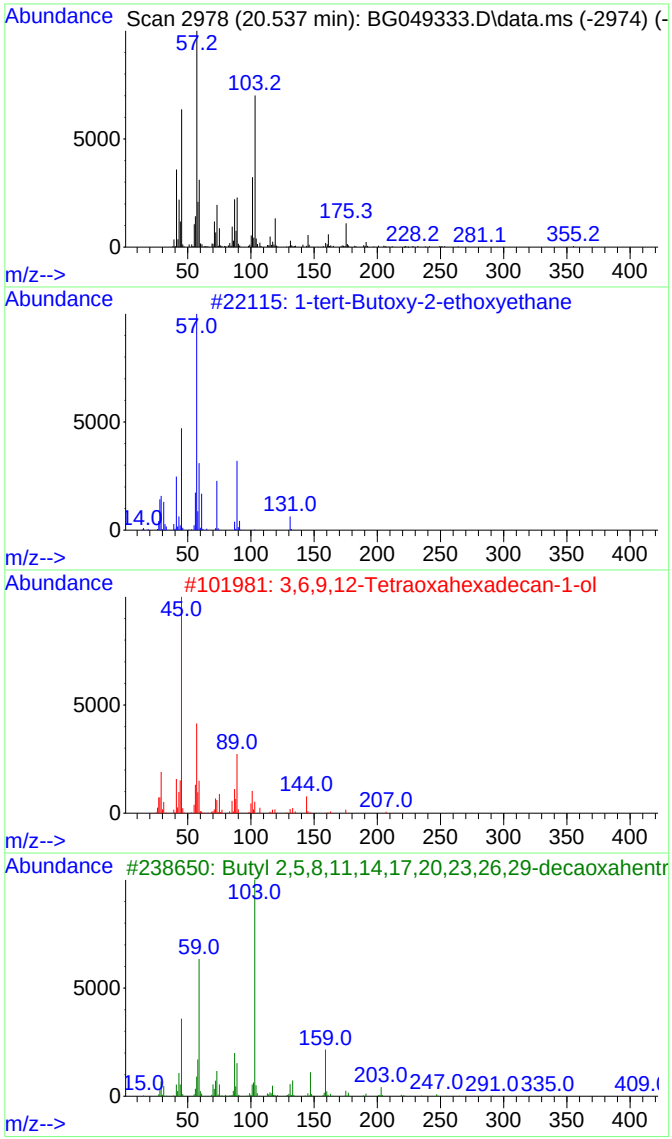
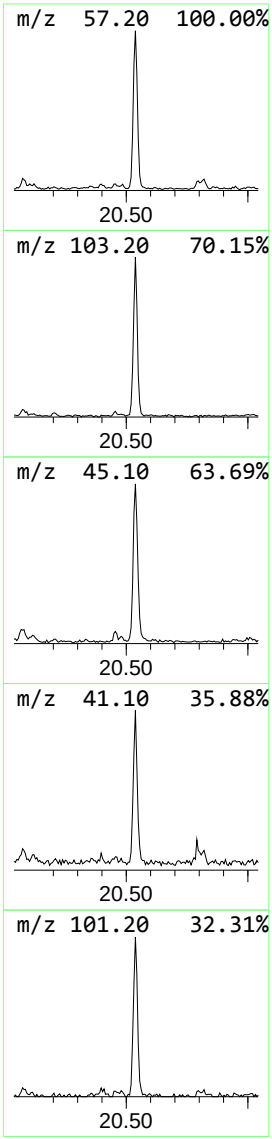
Peak Number 24 unknown-05 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.540	15.20 ng/ul	389904	Chrysene-d12	21.742

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-tert-Butoxy-2-ethoxyethane	146	C8H18O2	051422-54-9	38
2	3,6,9,12-Tetraoxahexadecan-1-ol	250	C12H26O5	001559-34-8	38
3	Butyl 2,5,8,11,14,17,20,23,26,29...	542	C25H50O12	1000366-81-6	35
4	3,6,9,12,15-Pentaoxanonadecan-1-ol	294	C14H30O6	001786-94-3	35
5	3,6,9,12-Tetraoxahexadecan-1-ol	250	C12H26O5	001559-34-8	35





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Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AT5

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

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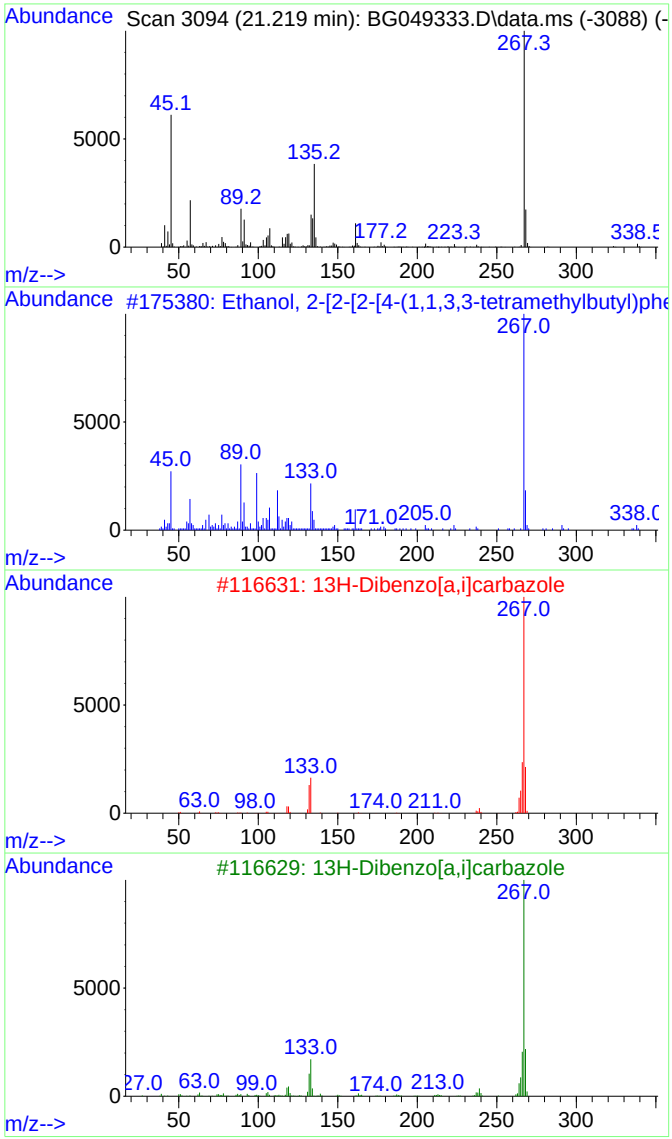
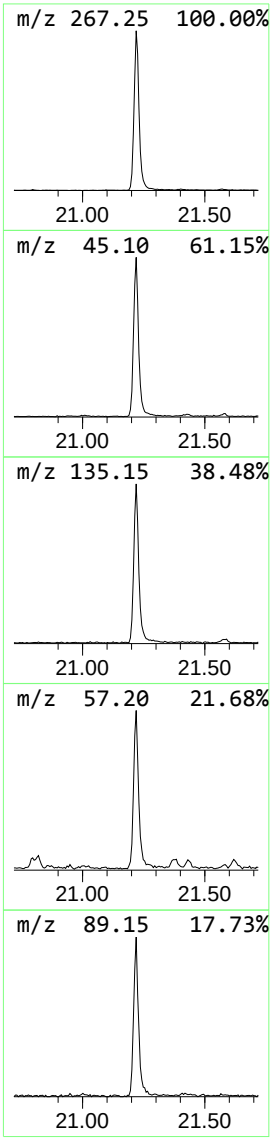
Peak Number 25 unknown-06 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.220	48.26 ng/ul	1237560	Chrysene-d12	21.742

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	46
2	13H-Dibenzo[a,i]carbazole	267	C20H13N	000239-64-5	43
3	13H-Dibenzo[a,i]carbazole	267	C20H13N	000239-64-5	35
4	Pyrido[2,3-d]pyrimidin-5(8H)-one...	267	C15H13N3O2	1000260-26-0	35
5	Benzaldehyde, 2,5-bis[(trimethyl...	282	C13H22O3Si2	056114-69-3	32





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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AT5

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Quant Title : SVOA CALIBRATION

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

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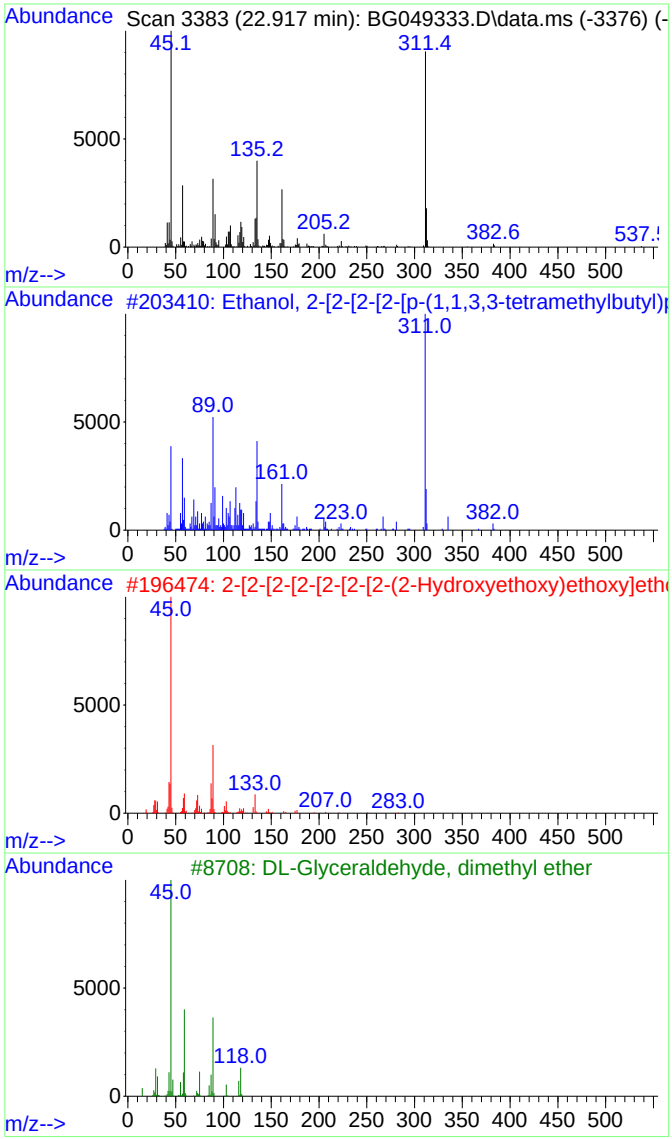
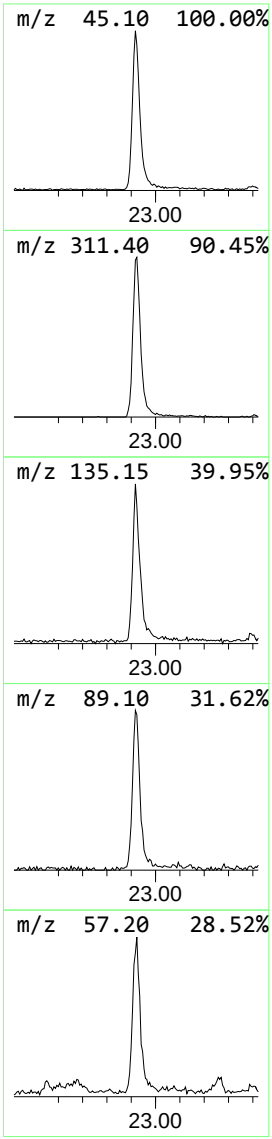
Peak Number 26 unknown-07 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.920	26.36 ng/ul	675913	Chrysene-d12	21.742

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-[2-[2-[2-[p-(1,1,3,3-...	382	C22H38O5	002315-63-1	35
2	2-[2-[2-[2-[2-[2-(2-Hydroxyet...	370	C16H34O9	005117-19-1	27
3	DL-Glyceraldehyde, dimethyl ether	118	C5H10O3	1000332-93-3	22
4	Hexaethylene glycol	282	C12H26O7	002615-15-8	22
5	Hexaethylene glycol	282	C12H26O7	002615-15-8	16





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ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

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Quant Title : SVOA CALIBRATION

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

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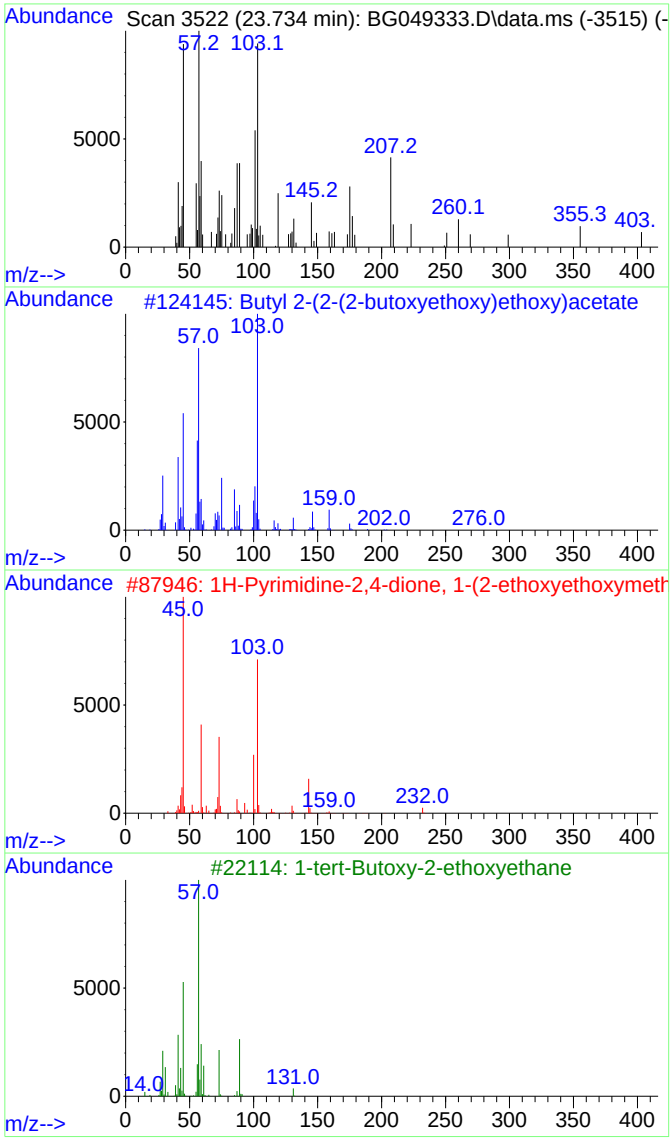
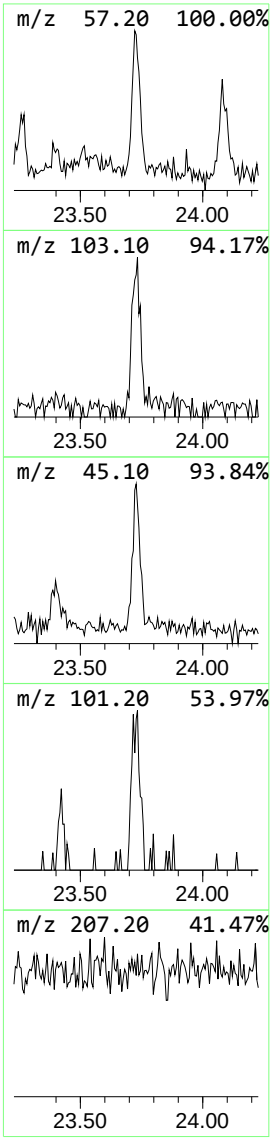
Peak Number 27 unknown-08 Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.730	2.12 ng/ul	57972	Perylene-d12	25.032

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butyl 2-(2-(2-butoxyethoxy)ethox...	276	C14H28O5	1000366-77-4	32
2	1H-Pyrimidine-2,4-dione, 1-(2-et...	232	C9H13FN2O4	1000310-13-7	32
3	1-tert-Butoxy-2-ethoxyethane	146	C8H18O2	051422-54-9	22
4	Di-sec-butyl ether	130	C8H18O	006863-58-7	16
5	Adonitol, pentamethyl ether	222	C10H22O5	1000332-81-0	16





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
Data File : BG049333.D
Acq On : 14 Jul 2021 17:29
Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

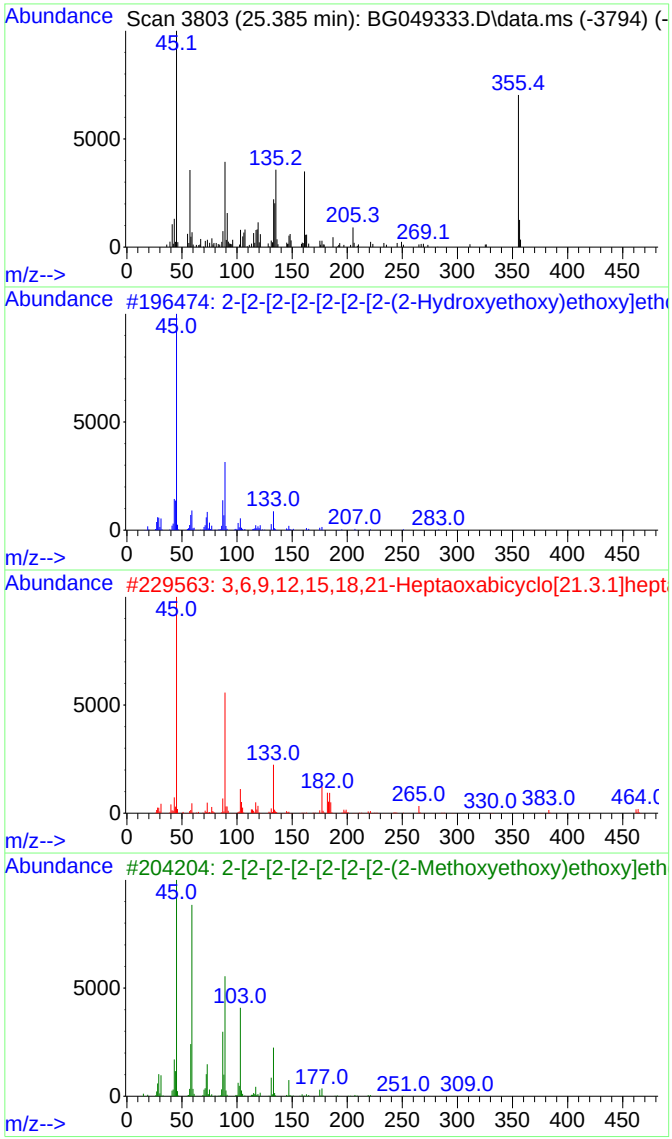
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

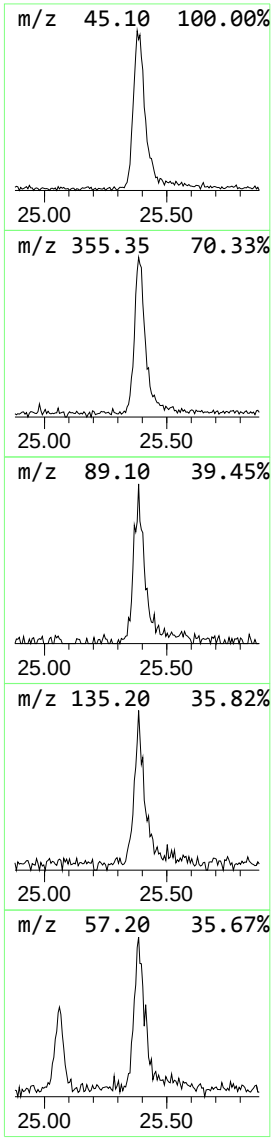
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 28 unknown-09 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.380	11.74 ng/ul	321016	Perylene-d12	25.032

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-[2-[2-[2-[2-[2-[2-(2-Hydroxyet...	370	C16H34O9	005117-19-1	38
2	3,6,9,12,15,18,21-Heptaoxabicycl...	462	C20H31BrO7	111982-23-1	32
3	2-[2-[2-[2-[2-[2-[2-(2-Methoxyet...	384	C17H36O9	1000352-04-2	32
4	Hexaethylene glycol	282	C12H26O7	002615-15-8	32
5	Cyclohexene, 4-isopropenyl-1-met...	196	C12H20O2	1000195-93-2	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
Data File : BG049333.D
Acq On : 14 Jul 2021 17:29
Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

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

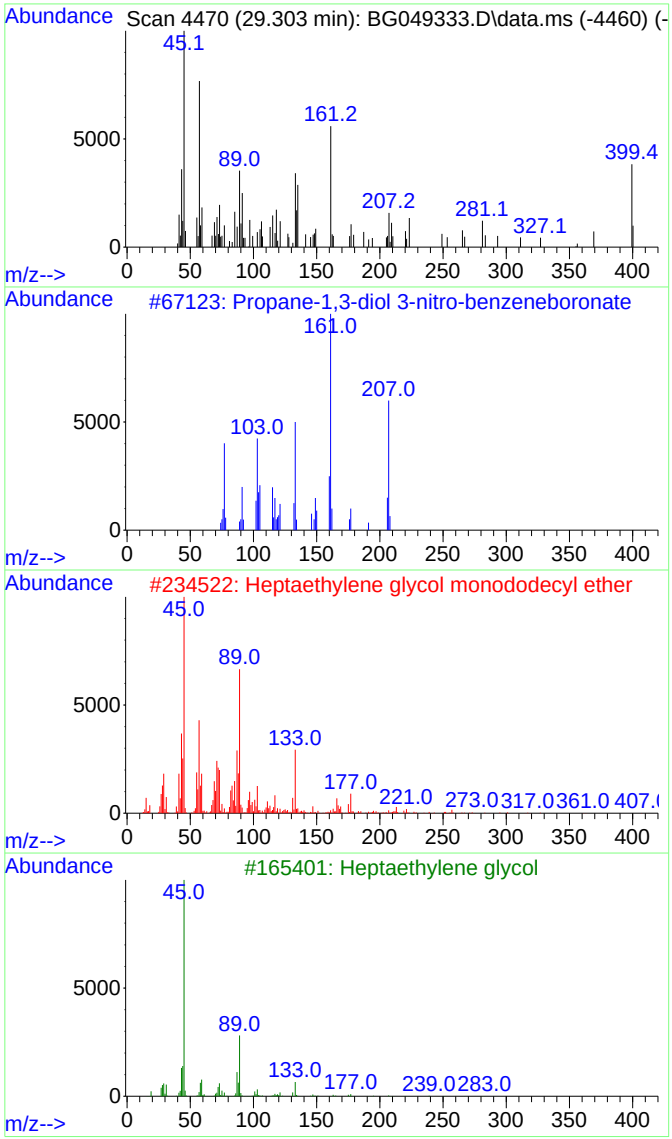
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

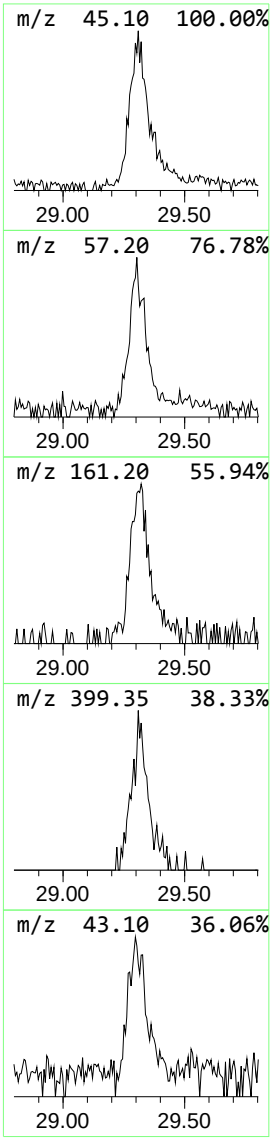
****

Peak Number 29 unknown-10 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.300	3.42 ng/ul	93574	Perylene-d12	25.032

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane-1,3-diol 3-nitro-benzene...	207	C9H10BNO4	085107-44-4	38
2	Heptaethylene glycol monododecyl...	494	C26H54O8	003055-97-8	35
3	Heptaethylene glycol	326	C14H30O8	005617-32-3	27
4	[2-[2-[2-[2-[2-[2-[2-(tert-Bu...	484	C22H48O9Si	1000352-10-5	25
5	2-[2-[2-[2-[2-[2-[2-[2-[2-(2-...	502	C22H46O12	1000352-13-2	25





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
Data File : BG049333.D
Acq On : 14 Jul 2021 17:29
Operator : CG/JU
Sample : M2996-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT5

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: C...	3.110	24.4	ng/ul	234428	1	8.064	191914	20.0
Butane, 2-metho...	3.190	16.7	ng/ul	160234	1	8.064	191914	20.0
1-Pentanol, 2-e...	8.120	3.7	ng/ul	35772	1	8.064	191914	20.0
Benzene, 1-ethy...	8.180	2.3	ng/ul	21850	1	8.064	191914	20.0
1,3-Methanopent...	8.450	2.6	ng/ul	25398	1	8.064	191914	20.0
Benzonitrile, 3...	8.920	2.5	ng/ul	23697	1	8.064	191914	20.0
Hexanoic acid, ...	9.360	4.1	ng/ul	39012	1	8.064	191914	20.0
Ethanol, 2-(2-b...	10.650	5.2	ng/ul	67363	2	10.884	259331	20.0
unknown-01	11.500	2.6	ng/ul	33233	2	10.884	259331	20.0
Benzene, 1-ethy...	12.430	2.2	ng/ul	28655	2	10.884	259331	20.0
unknown-02	12.760	3.0	ng/ul	39111	2	10.884	259331	20.0
Phenol, 3,5-dic...	13.420	12.2	ng/ul	226707	3	14.709	373296	20.0
Phenol, 3,4-dic...	13.730	9.0	ng/ul	168656	3	14.709	373296	20.0
1H-Indene-4-car...	13.840	2.4	ng/ul	44435	3	14.709	373296	20.0
Phenol, 2,3,5,6...	15.260	3.8	ng/ul	70439	3	14.709	373296	20.0
2-[2-[2-[2-[2-[...	15.950	3.0	ng/ul	57014	3	14.709	373296	20.0
Hydroquinone, t...	17.240	3.0	ng/ul	68476	4	17.464	448673	20.0
1,3-Dioxolane, ...	17.820	63.0	ng/ul	1412570	4	17.464	448673	20.0
Hexaethylene gl...	18.120	3.4	ng/ul	75541	4	17.464	448673	20.0
Ethyl 4-[(trime...	18.650	2.4	ng/ul	53975	4	17.464	448673	20.0
unknown-03	19.560	248.2	ng/ul	5568810	4	17.464	448673	20.0
Tricyclo[3.3.3....	19.710	78.3	ng/ul	2007320	5	21.742	512914	20.0
unknown-04	20.070	2.2	ng/ul	55459	5	21.742	512914	20.0
unknown-05	20.540	15.2	ng/ul	389904	5	21.742	512914	20.0
unknown-06	21.220	48.3	ng/ul	1237560	5	21.742	512914	20.0
unknown-07	22.920	26.4	ng/ul	675913	5	21.742	512914	20.0
unknown-08	23.730	2.1	ng/ul	57972	6	25.032	546893	20.0
unknown-09	25.380	11.7	ng/ul	321016	6	25.032	546893	20.0
unknown-10	29.300	3.4	ng/ul	93574	6	25.032	546893	20.0