LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050931.D
Acq On : 10 Nov 2021 6:41
Operator : CG/JU
Sample : M4532-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB8E8

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-BG110321.M
Title : SVOA CALIBRATION

Signal : TIC: BG050931.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.590	12	17	19	rBV	13630	20026	0.72%	0.110%
2	3.625	19	23	36	rVB	78660	128133	4.61%	0.703%
3	3.860	60	63	71	rVB6	2765	5202	0.19%	0.029%
4	4.019	85	90	101	rBV	60775	110444	3.97%	0.606%
5	4.254	124	130	136	rVB2	6210	11911	0.43%	0.065%
6	4.359	143	148	154	rVB3	2996	5974	0.21%	0.033%
7	4.647	193	197	205	rVB	7641	11631	0.42%	0.064%
8	5.511	337	344	355	rVB	401681	687139	24.72%	3.767%
9	5.693	371	375	385	rBV5	4265	8109	0.29%	0.044%
10	5.864	396	404	408	rBV5	2295	4499	0.16%	0.025%
11	6.692	542	545	550	rVB2	7477	10722	0.39%	0.059%
12	7.368	653	660	670	rBV	54324	107699	3.87%	0.590%
13	7.544	681	690	699	rBV	170431	289305	10.41%	1.586%
14	7.755	719	726	735	rBV	164747	290400	10.45%	1.592%
15	8.120	779	788	793	rBV3	10954	17752	0.64%	0.097%
16	8.231	800	807	810	rBV	126272	230934	8.31%	1.266%
17	8.261	810	812	822	rVB	56327	90425	3.25%	0.496%
18	8.596	862	869	874	rBV5	3373	6621	0.24%	0.036%
19	8.866	909	915	918	rBV3	8549	11764	0.42%	0.064%
20	8.925	918	925	936	rVB	140625	254435	9.15%	1.395%
21	9.066	945	949	953	rVB4	4539	6524	0.23%	0.036%
22	9.336	991	995	999	rBV5	3800	5753	0.21%	0.032%
23	9.401	999	1006	1017	rVB	204616	374810	13.48%	2.055%
24	9.571	1029	1035	1041	rVB6	3267	7539	0.27%	0.041%
25	9.659	1044	1050	1051	rBV	5883	9016	0.32%	0.049%
26	9.859	1081	1084	1090	rVB4	3914	5419	0.19%	0.030%
27	9.929	1090	1096	1101	rBV5	2852	5451	0.20%	0.030%
28	10.035	1107	1114	1120	rVB2	6577	11902	0.43%	0.065%
29	10.123	1120	1129	1139	rBV	159854	305124	10.98%	1.673%
30	10.211	1139	1144	1150	rBV3	2070	4835	0.17%	0.027%
31	10.664	1214	1221	1231	rBV	213207	397260	14.29%	2.178%
32	10.805	1239	1245	1251	rBV3	10418	19668	0.71%	0.108%
33	10.934	1261	1267	1275	rBV3	15196	31082	1.12%	0.170%
34	11.052	1279	1287	1294	rBV	179942	328651	11.82%	1.802%
35	11.187	1301	1310	1323	rBV	207584	394786	14.20%	2.165%
36	11.839	1416	1421	1426	rVB3	12891	19459	0.70%	0.107%

LSC Area Percent Report

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Operator : CG/JU
Sample : M4532-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB8E8

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-BG110321.M
Title : SVOA CALIBRATION

37	11.968	1440	1443	1444	rVV3	4759	5355	0.19%	0.029%
38	11.980	1444	1445	1450	rVV2	5156	6543	0.24%	0.036%
39	12.432	1516	1522	1527	rBV7	4346	6947	0.25%	0.038%
40	12.620	1544	1554	1559	rBV5	7594	19382	0.70%	0.106%
41	13.566	1710	1715	1725	rVB3	9020	15323	0.55%	0.084%
42	13.901	1765	1772	1780	rBV	977911	1491466	53.66%	8.177%
43	14.242	1823	1830	1836	rBV	474863	699230	25.16%	3.834%
44	14.289	1836	1838	1843	rVB4	7025	8842	0.32%	0.048%
45	14.448	1857	1865	1875	rVV4	14857	33685	1.21%	0.185%
46	14.547	1875	1882	1892	rVV	495880	808325	29.08%	4.432%
47	14.630	1892	1896	1902	rVV5	5254	8976	0.32%	0.049%
48	14.853	1922	1934	1941	rBV	276313	462681	16.65%	2.537%
49	14.994	1953	1958	1961	rBV4	5047	8080	0.29%	0.044%
50	15.041	1961	1966	1975	rVV	41135	82417	2.97%	0.452%
51	15.123	1975	1980	1990	rVB2	24995	49048	1.76%	0.269%
52	15.317	2006	2013	2016	rBV3	4261	6022	0.22%	0.033%
53	15.394	2021	2026	2032	rBV3	6862	10563	0.38%	0.058%
54	15.458	2032	2037	2044	rVB4	6362	9526	0.34%	0.052%
55	15.570	2049	2056	2060	rBV4	9941	20651	0.74%	0.113%
56	15.617	2060	2064	2069	rVV4	8749	16546	0.60%	0.091%
57	15.670	2069	2073	2077	rVB4	3427	6050	0.22%	0.033%
58	15.752	2082	2087	2088	rVV3	3301	4058	0.15%	0.022%
59	15.787	2088	2093	2096	rVV	17670	27794	1.00%	0.152%
60	15.840	2096	2102	2115	rVV	658160	1050402	37.79%	5.759%
61	15.952	2115	2121	2129	rVV	213282	335405	12.07%	1.839%
62	16.010	2129	2131	2136	rVV4	5397	10366	0.37%	0.057%
63	16.081	2137	2143	2152	rVV	224169	387550	13.94%	2.125%
64	16.146	2152	2154	2162	rVV5	7948	15745	0.57%	0.086%
65	16.222	2162	2167	2172	rVV4	16587	25033	0.90%	0.137%
66	16.510	2211	2216	2218	rBV4	6282	8502	0.31%	0.047%
67	16.580	2224	2228	2236	rVB6	10150	15210	0.55%	0.083%
68	16.939	2283	2289	2297	rBV4	15111	28470	1.02%	0.156%
69	17.280	2341	2347	2352	rBV	14024	20947	0.75%	0.115%
70	17.415	2361	2370	2375	rVB7	6470	13726	0.49%	0.075%
71	17.597	2391	2401	2408	rBV2	358349	680372	24.48%	3.730%
72	17.697	2411	2418	2425	rVB	723988	1152102	41.45%	6.317%
73	17.979	2463	2466	2473	rVB6	3048	4585	0.16%	0.025%
74	18.225	2505	2508	2513	rBV4	6959	9977	0.36%	0.055%
75	18.461	2542	2548	2550	rBV3	3980	7048	0.25%	0.039%
76	18.519	2553	2558	2561	rVV3	30446	45205	1.63%	0.248%

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Operator : CG/JU
Sample : M4532-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB8E8

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-BG110321.M
Title : SVOA CALIBRATION

77	18.566	2561	2566	2574	rVV	78741	127549	4.59%	0.699%
78	18.719	2587	2592	2594	rBV	76300	115757	4.16%	0.635%
79	18.748	2594	2597	2608	rVB	90501	143255	5.15%	0.785%
80	18.966	2630	2634	2639	rVB6	2992	5238	0.19%	0.029%
81	19.389	2703	2706	2711	rVB6	5863	7862	0.28%	0.043%
82	19.542	2723	2732	2735	rBV2	11997	21699	0.78%	0.119%
83	19.606	2738	2743	2749	rVV	11038	18426	0.66%	0.101%
84	19.976	2798	2806	2819	rBV2	1758278	2779587	100.00%	15.240%
85	20.094	2821	2826	2841	rVB	360601	593520	21.35%	3.254%
86	20.258	2849	2854	2867	rVB6	11878	28567	1.03%	0.157%
87	20.546	2901	2903	2908	rBV6	4458	5040	0.18%	0.028%
88	21.181	3008	3011	3015	rVB5	4726	5048	0.18%	0.028%
89	21.369	3040	3043	3049	rVB9	6948	13379	0.48%	0.073%
90	21.504	3060	3066	3069	rBV7	14073	31905	1.15%	0.175%
91	21.739	3102	3106	3110	rBV7	5762	7993	0.29%	0.044%
92	21.892	3125	3132	3140	rVB	317229	555300	19.98%	3.045%
93	22.050	3155	3159	3164	rBV7	5973	9590	0.35%	0.053%
94	22.691	3263	3268	3272	rBV4	5790	9022	0.32%	0.049%
95	23.414	3385	3391	3399	rVB7	11636	24185	0.87%	0.133%
96	24.260	3530	3535	3545	rVB3	12769	24654	0.89%	0.135%
97	25.053	3656	3670	3683	rBV2	403768	1218856	43.85%	6.683%
98	25.288	3698	3710	3722	rVB2	191154	569283	20.48%	3.121%
99	26.457	3901	3909	3919	rVB6	13830	40886	1.47%	0.224%
100	27.902	4146	4155	4164	rVB7	10865	33737	1.21%	0.185%

Sum of corrected areas: 18238897

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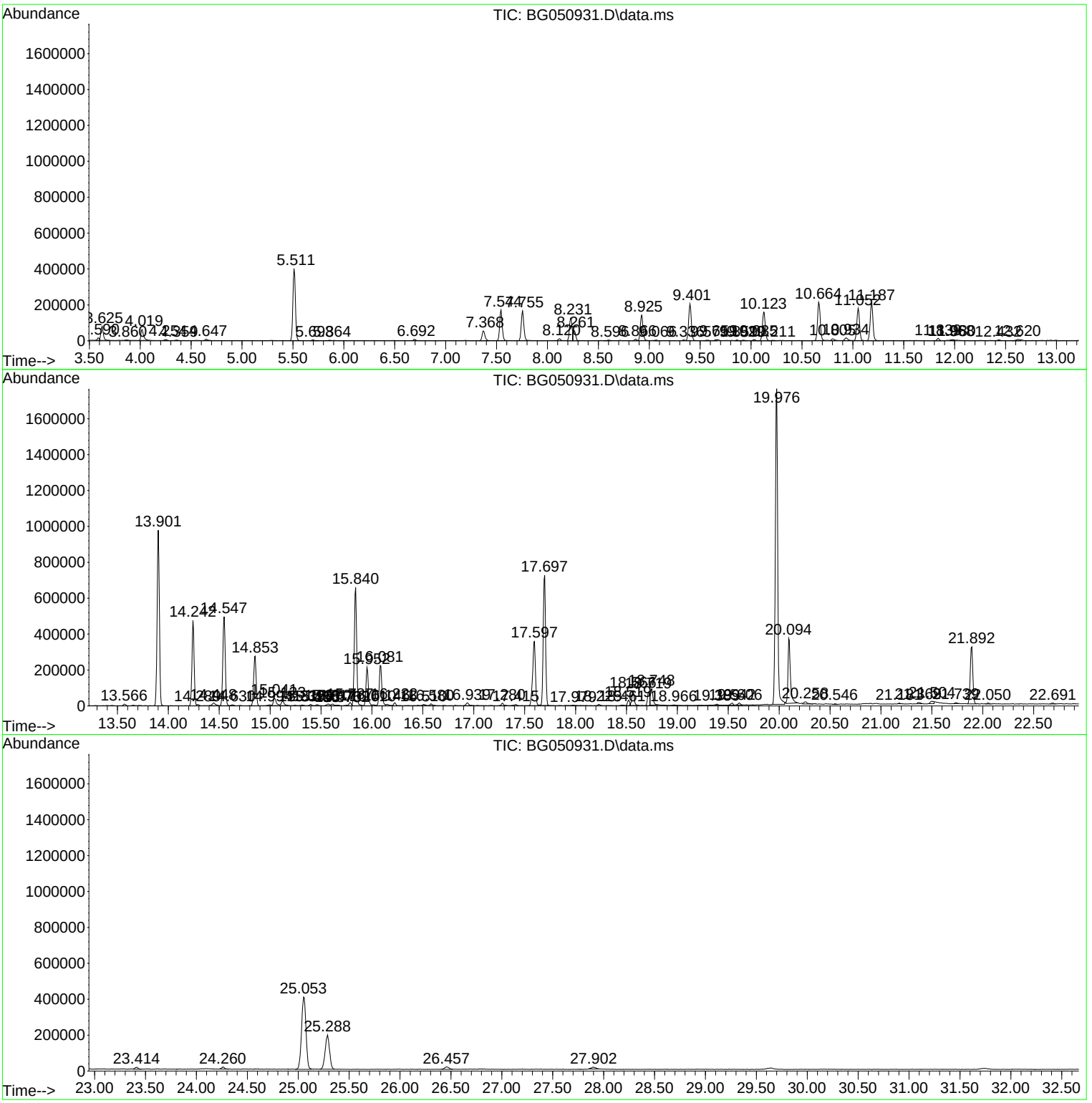
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB8E8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050931.D
Acq On : 10 Nov 2021 6:41
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Instrument :
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ClientSampleId :
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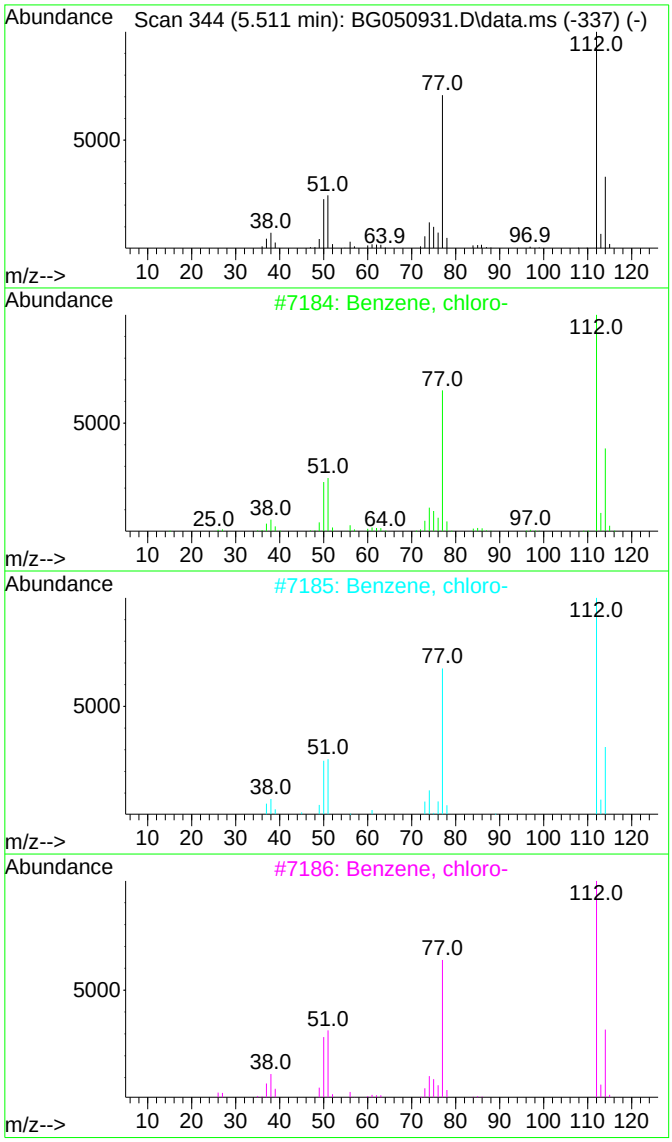
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

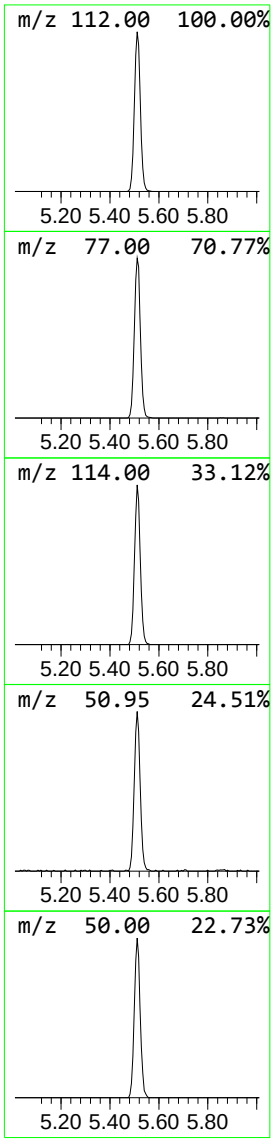
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Benzene, chloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.511	59.51 ng/ul	687139	1,4-Dichlorobenzene-d4	8.225

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, chloro-	112	C6H5Cl	000108-90-7	96
2	Benzene, chloro-	112	C6H5Cl	000108-90-7	96
3	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
4	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
5	Benzene, chloro-	112	C6H5Cl	000108-90-7	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050931.D
Acq On : 10 Nov 2021 6:41
Operator : CG/JU
Sample : M4532-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB8E8

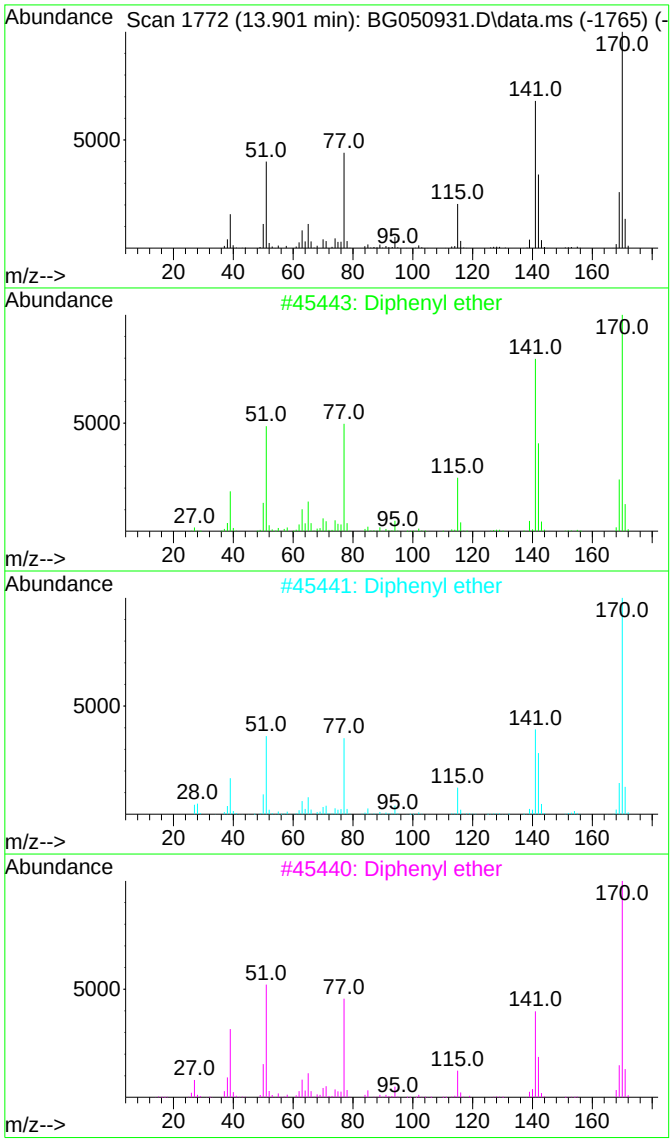
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

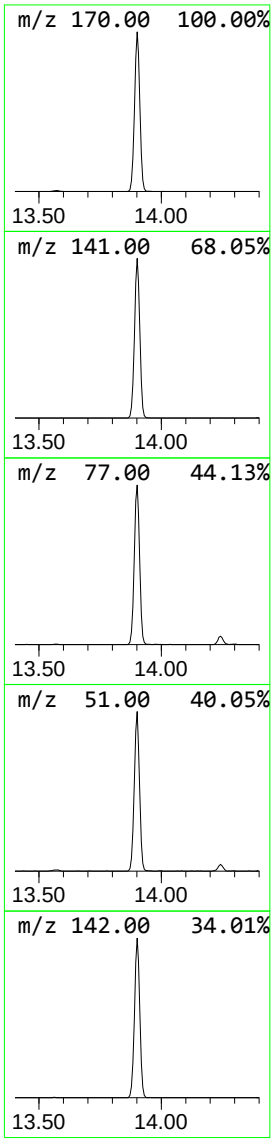
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Diphenyl ether Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.901	64.47 ng/ul	1491470	Acenaphthene-d10	14.853

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diphenyl ether	170	C12H10O	000101-84-8	98
2	Diphenyl ether	170	C12H10O	000101-84-8	90
3	Diphenyl ether	170	C12H10O	000101-84-8	87
4	Diphenyl ether	170	C12H10O	000101-84-8	87
5	Diphenyl ether	170	C12H10O	000101-84-8	81





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Sample : M4532-11
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ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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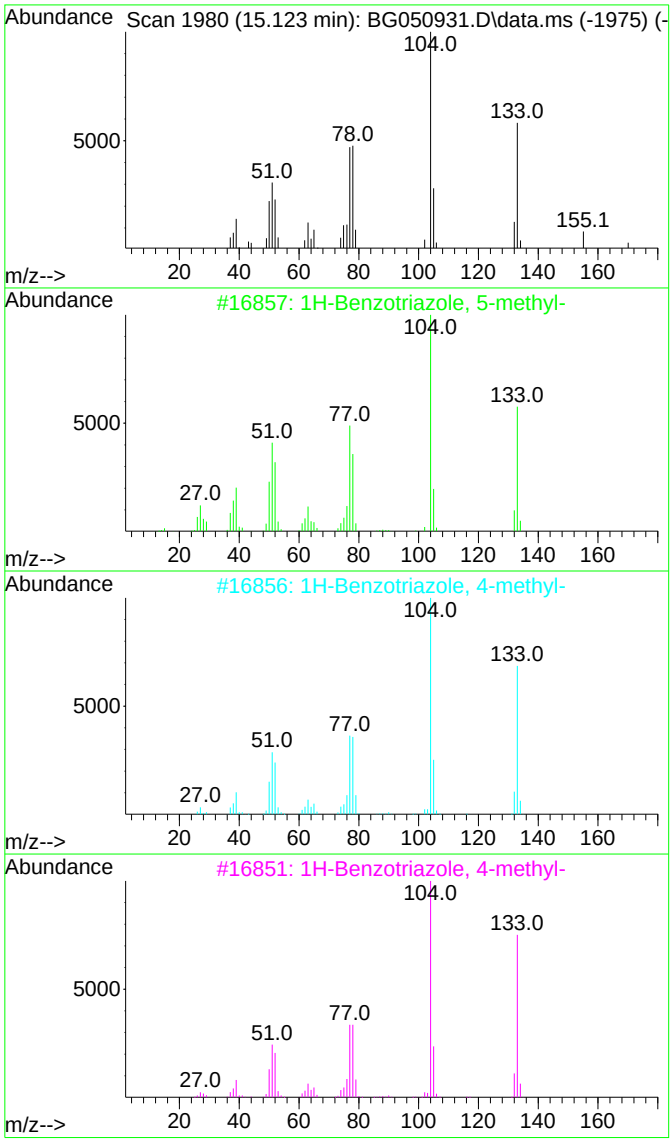
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

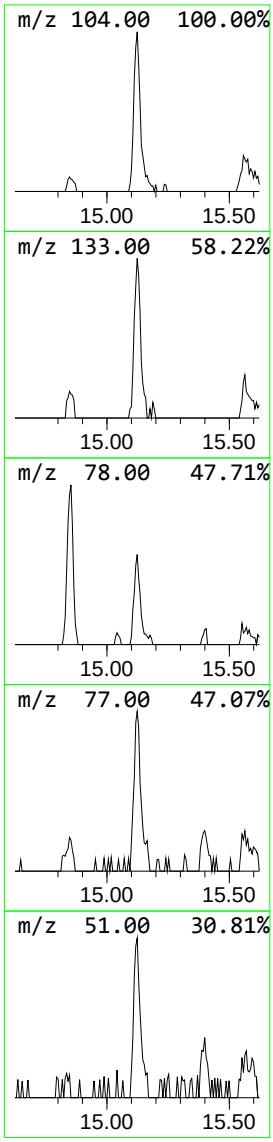
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1H-Benzotriazole, 5-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.123	2.12 ng/ul	49048	Acenaphthene-d10	14.853

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Benzotriazole, 5-methyl-	133	C7H7N3	000136-85-6	96
2	1H-Benzotriazole, 4-methyl-	133	C7H7N3	029878-31-7	95
3	1H-Benzotriazole, 4-methyl-	133	C7H7N3	029878-31-7	94
4	1H-Benzotriazole, 4-methyl-	133	C7H7N3	029878-31-7	93
5	1H-Benzotriazole, 5-methyl-	133	C7H7N3	000136-85-6	91





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BNA_G
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

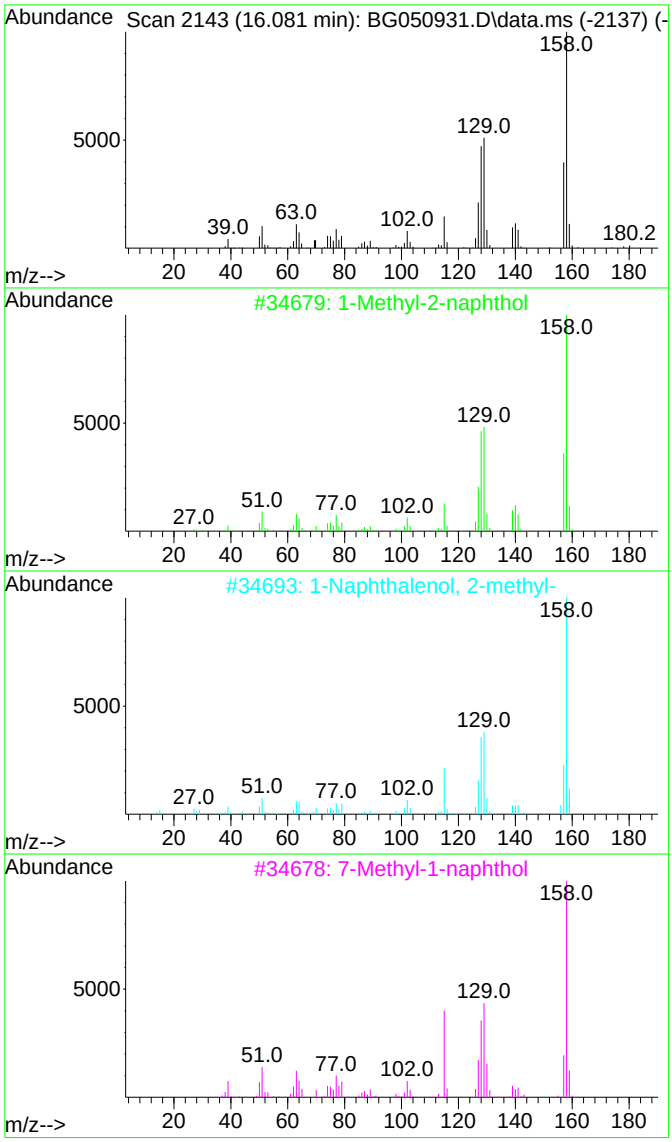
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

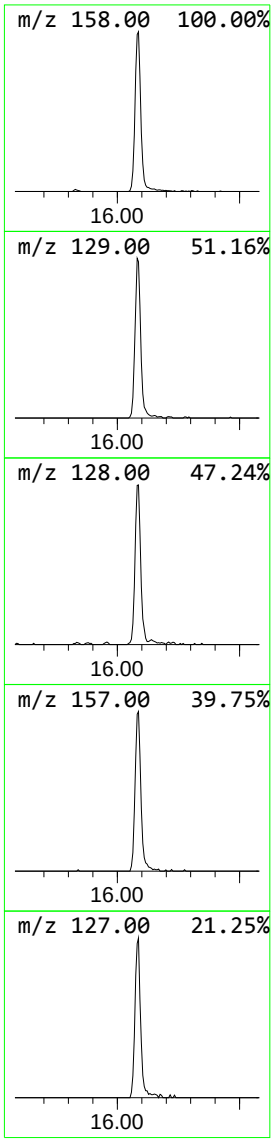
****

Peak Number 4 1-Methyl-2-naphthol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.081	16.75 ng/ul	387550	Acenaphthene-d10	14.853

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Methyl-2-naphthol	158	C11H10O	001076-26-2	98
2	1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	94
3	7-Methyl-1-naphthol	158	C11H10O	006939-33-9	87
4	1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	83
5	2-Naphthalenemethanol	158	C11H10O	001592-38-7	58





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050931.D
Acq On : 10 Nov 2021 6:41
Operator : CG/JU
Sample : M4532-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB8E8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

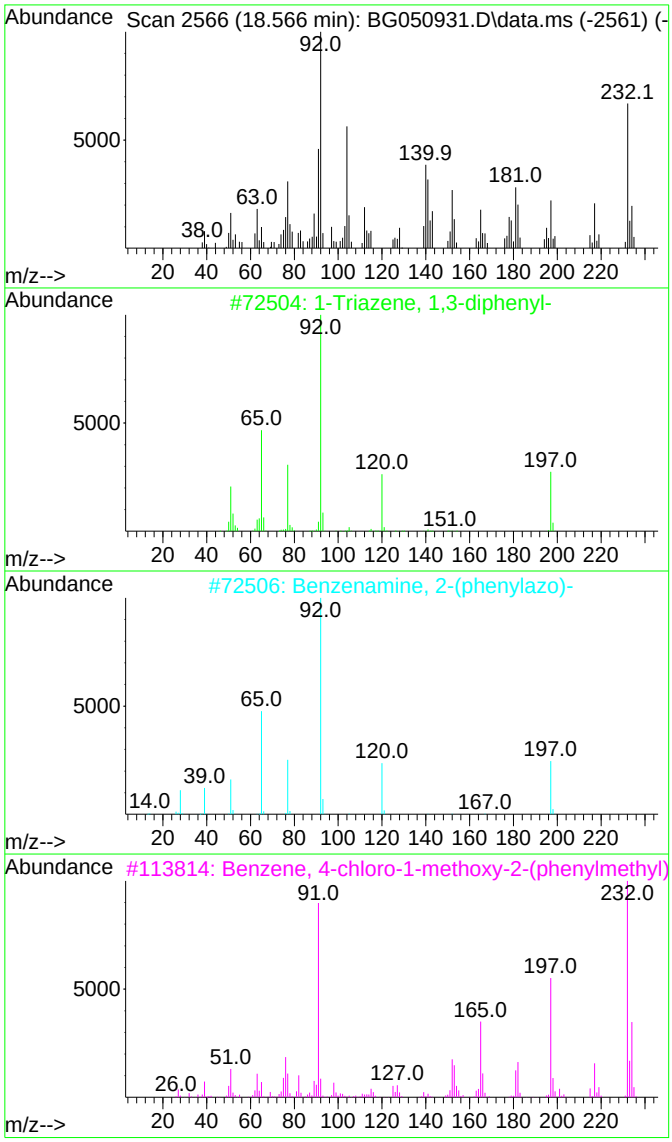
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

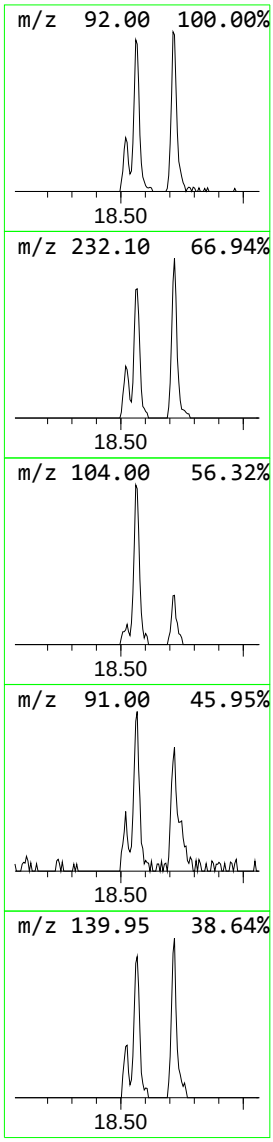
****

Peak Number 5 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.566	3.75 ng/ul	127549	Phenanthrene-d10	17.597

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Triazene, 1,3-diphenyl-	197	C12H11N3	000136-35-6	18
2	Benzenamine, 2-(phenylazo)-	197	C12H11N3	002835-58-7	18
3	Benzene, 4-chloro-1-methoxy-2-(p...	232	C14H13ClO	040395-13-9	14
4	Benzenamine, 4-(phenylazo)-	197	C12H11N3	000060-09-3	14
5	Benzenamine, 2-(phenylazo)-	197	C12H11N3	002835-58-7	14





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Data File : BG050931.D
Acq On : 10 Nov 2021 6:41
Operator : CG/JU
Sample : M4532-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB8E8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

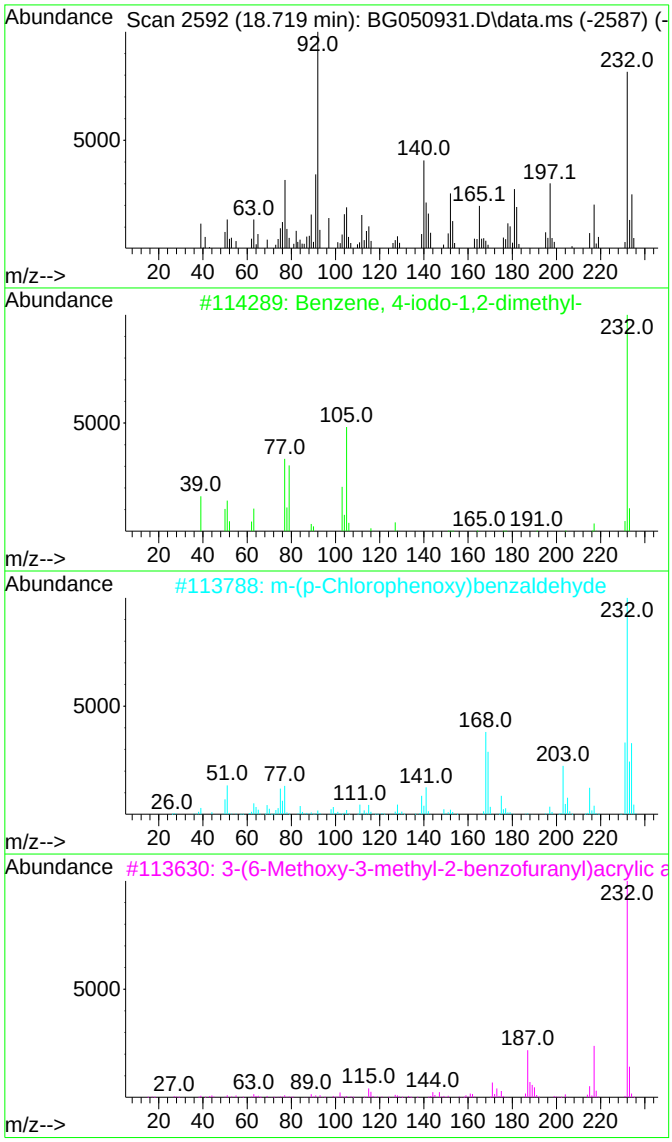
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

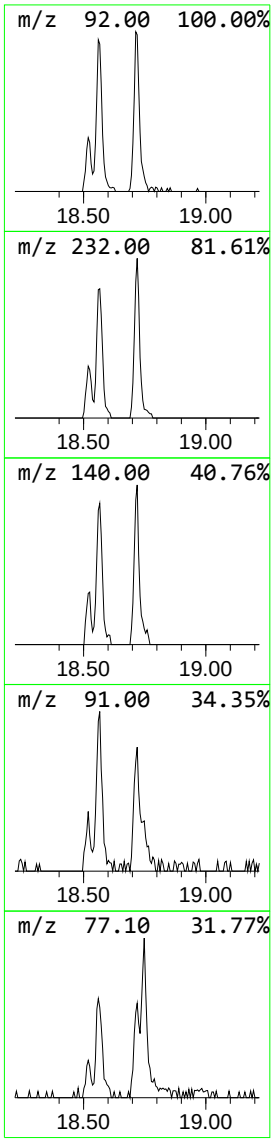
****

Peak Number 6 unknown-02 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.719	3.40 ng/ul	115757	Phenanthrene-d10	17.597

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-iodo-1,2-dimethyl-	232	C8H9I	031599-61-8	44
2	m-(p-Chlorophenoxy)benzaldehyde	232	C13H9ClO2	069770-20-3	25
3	3-(6-Methoxy-3-methyl-2-benzofur...	232	C13H12O4	010410-32-9	22
4	6-Chlorothiocarbonyl-3-nitro-1-m...	232	C7H5ClN2O3S	1000216-40-1	22
5	Benzene, 4-chloro-1-methoxy-2-(p...	232	C14H13ClO	040395-13-9	22





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050931.D
Acq On : 10 Nov 2021 6:41
Operator : CG/JU
Sample : M4532-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB8E8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

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

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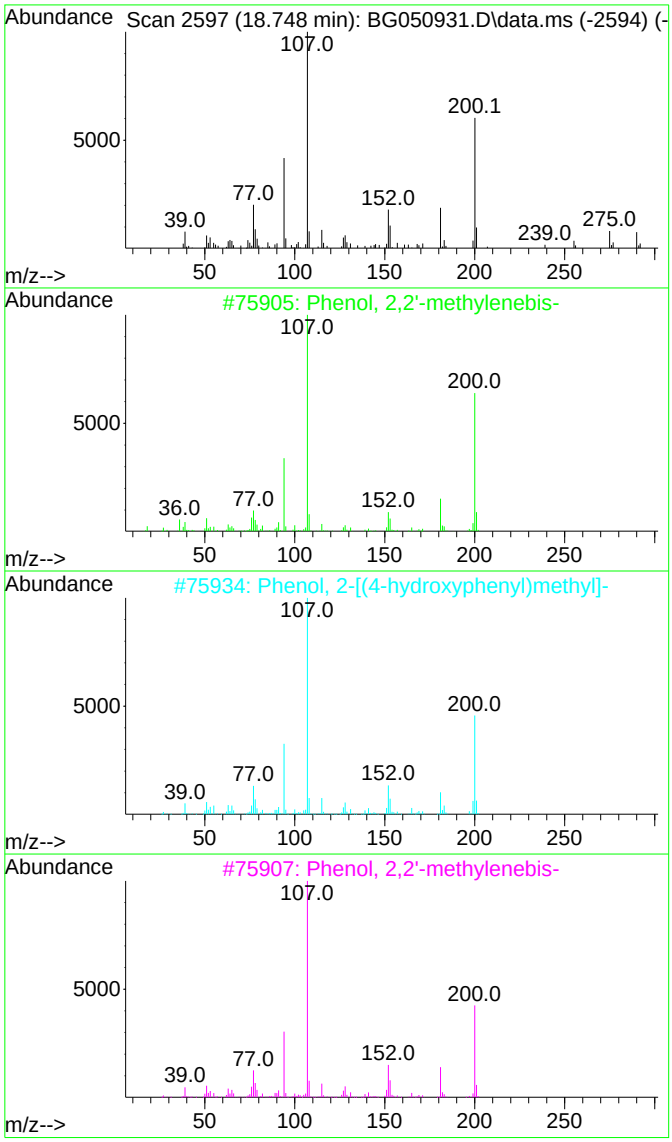
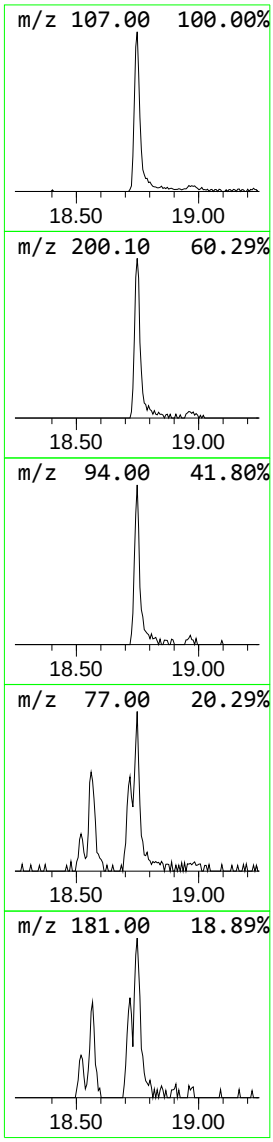
Peak Number 7 Phenol, 2,2'-methylenebis- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.748	4.21 ng/ul	143255	Phenanthrene-d10	17.597

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	93
2	Phenol, 2-[(4-hydroxyphenyl)meth...	200	C13H12O2	002467-03-0	90
3	Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	90
4	Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	68
5	Phenol, 2-[(4-hydroxyphenyl)meth...	200	C13H12O2	002467-03-0	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050931.D
Acq On : 10 Nov 2021 6:41
Operator : CG/JU
Sample : M4532-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB8E8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

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

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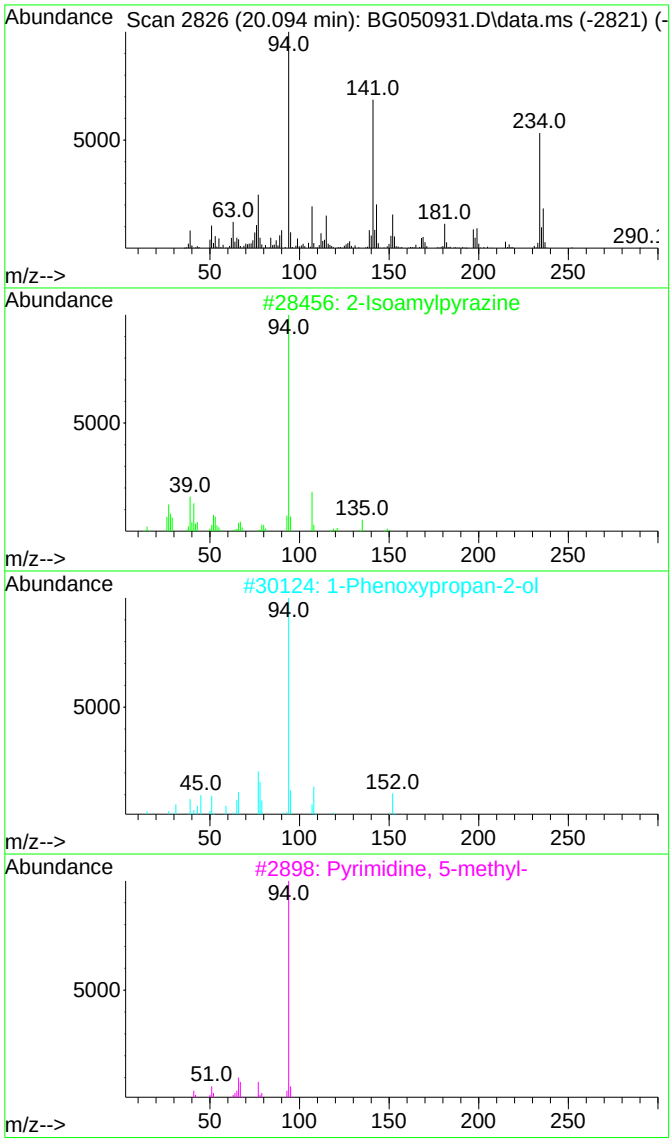
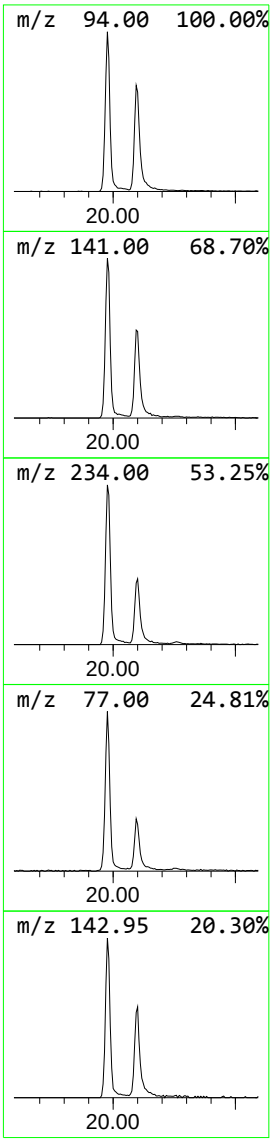
Peak Number 8 unknown-03 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.094	21.38 ng/ul	593520	Chrysene-d12	21.892

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Isoamylpyrazine	150	C9H14N2	040790-22-5	27
2	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	22
3	Pyrimidine, 5-methyl-	94	C5H6N2	002036-41-1	18
4	4H-1,2,4-Triazole-3-carbonitrile	94	C3H2N4	003641-10-9	14
5	Benzenemethanol, 3-iodo-	234	C7H7IO	057455-06-8	11





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050931.D
Acq On : 10 Nov 2021 6:41
Operator : CG/JU
Sample : M4532-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB8E8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, chloro-	5.511	59.5	ng/ul	687139	1	8.225	230934	20.0
Diphenyl ether	13.901	64.5	ng/ul	1491470	3	14.853	462681	20.0
1H-Benzotriazol...	15.123	2.1	ng/ul	49048	3	14.853	462681	20.0
1-Methyl-2-naph...	16.081	16.8	ng/ul	387550	3	14.853	462681	20.0
unknown-01	18.566	3.8	ng/ul	127549	4	17.597	680372	20.0
unknown-02	18.719	3.4	ng/ul	115757	4	17.597	680372	20.0
Phenol, 2,2'-me...	18.748	4.2	ng/ul	143255	4	17.597	680372	20.0
unknown-03	20.094	21.4	ng/ul	593520	5	21.892	555300	20.0