LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062835.D
Acq On : 15 Sep 2024 7:08
Operator : JU/RC
Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ3

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-BG090924.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BG062835.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.769	112	116	125	rBV	47787	84319	4.04%	0.381%
2	4.545	241	248	256	rBV7	10473	21976	1.05%	0.099%
3	5.420	391	397	403	rVB	12208	17507	0.84%	0.079%
4	6.396	557	563	573	rBV	187575	292567	14.02%	1.321%
5	6.883	640	646	657	rVB	194007	310038	14.86%	1.400%
6	7.065	670	677	688	rVB	69782	124467	5.97%	0.562%
7	7.230	698	705	712	rVV	127036	201847	9.67%	0.912%
8	7.330	712	722	730	rVB3	19658	40763	1.95%	0.184%
9	7.436	732	740	750	rBV	167197	288033	13.81%	1.301%
10	7.900	811	819	829	rVV	131523	226762	10.87%	1.024%
11	8.017	833	839	848	rVB	84423	140139	6.72%	0.633%
12	8.135	850	859	868	rBV6	16482	38696	1.85%	0.175%
13	8.276	876	883	889	rBV	134085	223004	10.69%	1.007%
14	8.393	898	903	908	rBV2	21473	34751	1.67%	0.157%
15	8.446	908	912	918	rVB	14399	21466	1.03%	0.097%
16	8.540	922	928	933	rBV3	17071	29924	1.43%	0.135%
17	8.605	933	939	948	rVV	207012	360942	17.30%	1.630%
18	8.699	950	955	961	rVB4	9013	16371	0.78%	0.074%
19	8.851	976	981	988	rVV4	14863	34930	1.67%	0.158%
20	8.928	988	994	1001	rVV4	26007	53071	2.54%	0.240%
21	9.004	1001	1007	1011	rVV2	138686	230560	11.05%	1.041%
22	9.051	1011	1015	1027	rVB2	182704	372790	17.87%	1.684%
23	9.339	1058	1064	1070	rVB2	16298	30259	1.45%	0.137%
24	9.527	1082	1096	1102	rBV	123309	203481	9.75%	0.919%
25	9.774	1131	1138	1147	rBV	157006	286853	13.75%	1.296%
26	9.944	1162	1167	1175	rVB2	23244	43796	2.10%	0.198%
27	10.097	1184	1193	1200	rVB2	144379	269048	12.90%	1.215%
28	10.168	1200	1205	1211	rBV5	7469	14181	0.68%	0.064%
29	10.314	1222	1230	1238	rVV	281825	506785	24.29%	2.289%
30	10.391	1238	1243	1252	rVB3	13068	25421	1.22%	0.115%
31	10.691	1286	1294	1298	rBV	209238	373469	17.90%	1.687%
32	10.743	1298	1303	1310	rVV	371639	640288	30.69%	2.892%
33	10.826	1310	1317	1327	rVV	166778	310824	14.90%	1.404%
34	10.908	1327	1331	1337	rVB3	11642	18226	0.87%	0.082%
35	10.967	1337	1341	1346	rVB2	8875	13675	0.66%	0.062%
36	11.090	1357	1362	1372	rVB6	10639	27589	1.32%	0.125%

LSC Area Percent Report

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Misc :
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Instrument :
BNA_G
ClientSampleId :
C0AQ3

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-BG090924.MA.M
Title : SVOA CALIBRATION

37	11.360	1404	1408	1412	rVB5	9186	15727	0.75%	0.071%
38	11.513	1425	1434	1436	rBV7	7162	19996	0.96%	0.090%
39	11.795	1478	1482	1490	rVB7	7131	13795	0.66%	0.062%
40	12.018	1514	1520	1523	rBV2	7861	12266	0.59%	0.055%
41	12.071	1524	1529	1537	rVB4	15998	30750	1.47%	0.139%
42	12.300	1566	1568	1573	rVV6	8183	13533	0.65%	0.061%
43	12.406	1581	1586	1591	rVV4	19531	33001	1.58%	0.149%
44	12.565	1605	1613	1624	rVV3	50111	123177	5.90%	0.556%
45	12.671	1624	1631	1638	rVV2	80415	182667	8.76%	0.825%
46	12.870	1660	1665	1670	rVV8	7526	15713	0.75%	0.071%
47	12.941	1672	1677	1678	rVV5	8662	14520	0.70%	0.066%
48	12.982	1679	1684	1691	rVV2	72474	122705	5.88%	0.554%
49	13.058	1691	1697	1700	rVV6	9745	19662	0.94%	0.089%
50	13.105	1700	1705	1711	rVB2	26075	42742	2.05%	0.193%
51	13.246	1723	1729	1732	rBV3	10599	19840	0.95%	0.090%
52	13.487	1765	1770	1774	rBV	45219	70653	3.39%	0.319%
53	13.534	1774	1778	1785	rVV	119083	186651	8.95%	0.843%
54	13.652	1792	1798	1802	rBV2	73743	121954	5.85%	0.551%
55	13.693	1802	1805	1809	rVV2	41218	60834	2.92%	0.275%
56	13.758	1809	1816	1824	rVV2	142312	275695	13.21%	1.245%
57	13.828	1824	1828	1834	rVV3	29242	49239	2.36%	0.222%
58	13.934	1840	1846	1852	rVB	569755	808628	38.76%	3.652%
59	14.016	1856	1860	1863	rBV5	12137	19342	0.93%	0.087%
60	14.057	1863	1867	1872	rVB7	11274	19545	0.94%	0.088%
61	14.163	1879	1885	1890	rVV2	142100	266685	12.78%	1.204%
62	14.222	1890	1895	1902	rVB	616091	927697	44.47%	4.190%
63	14.292	1902	1907	1911	rBV2	22390	36676	1.76%	0.166%
64	14.345	1911	1916	1921	rVB	82560	131724	6.31%	0.595%
65	14.445	1929	1933	1939	rVB6	12579	21037	1.01%	0.095%
66	14.527	1939	1947	1955	rBV	420459	643492	30.84%	2.906%
67	14.609	1955	1961	1965	rBV7	7271	14029	0.67%	0.063%
68	14.692	1966	1975	1978	rVV4	85426	192972	9.25%	0.872%
69	14.727	1978	1981	1992	rVB	100612	168638	8.08%	0.762%
70	14.844	1997	2001	2003	rBV5	14049	18536	0.89%	0.084%
71	14.915	2009	2013	2016	rBV4	15379	23348	1.12%	0.105%
72	14.944	2016	2018	2023	rVB6	14712	16360	0.78%	0.074%
73	15.009	2023	2029	2032	rBV3	30020	56206	2.69%	0.254%
74	15.044	2033	2035	2040	rVB3	23160	27112	1.30%	0.122%
75	15.173	2049	2057	2062	rBV7	22831	52878	2.53%	0.239%
76	15.226	2062	2066	2075	rVB	36995	68966	3.31%	0.311%

LSC Area Percent Report

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Operator : JU/RC
Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ3

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

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Peak separation: 5

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Title : SVOA CALIBRATION

77	15.350	2083	2087	2092	rVV7	7972	12654	0.61%	0.057%
78	15.414	2096	2098	2103	rVB	16753	22125	1.06%	0.100%
79	15.520	2109	2116	2126	rBV	860788	1311504	62.86%	5.923%
80	15.638	2126	2136	2140	rBV	315134	473274	22.68%	2.138%
81	15.738	2148	2153	2157	rBV	66810	84985	4.07%	0.384%
82	15.849	2167	2172	2173	rBV2	11962	14552	0.70%	0.066%
83	15.884	2173	2178	2182	rVV	98472	151108	7.24%	0.682%
84	15.920	2182	2184	2188	rVV4	21636	27440	1.32%	0.124%
85	16.072	2206	2210	2217	rBV2	36507	54942	2.63%	0.248%
86	16.260	2238	2242	2250	rVV10	8775	19562	0.94%	0.088%
87	17.271	2408	2414	2422	rVB	574848	870817	41.74%	3.933%
88	17.371	2425	2431	2442	rVB2	1053908	1570223	75.26%	7.092%
89	18.164	2562	2566	2573	rBV2	56978	83381	4.00%	0.377%
90	18.652	2646	2649	2654	rVB7	11678	21135	1.01%	0.095%
91	19.134	2728	2731	2736	rVB	25275	33456	1.60%	0.151%
92	19.228	2744	2747	2752	rVB3	20774	27159	1.30%	0.123%
93	19.298	2756	2759	2766	rVB2	17762	33889	1.62%	0.153%
94	19.445	2781	2784	2793	rVB9	27987	38607	1.85%	0.174%
95	19.662	2815	2821	2826	rBV	1412107	1997957	95.76%	9.024%
96	19.709	2826	2829	2833	rVB	39224	54615	2.62%	0.247%
97	21.519	3131	3137	3144	rBV	673043	1033410	49.53%	4.667%
98	23.129	3407	3411	3418	rVB3	21156	40961	1.96%	0.185%
99	24.374	3612	3623	3635	rBV	801272	2086345	100.00%	9.423%
100	24.586	3649	3659	3675	rVB	417124	1189172	57.00%	5.371%

Sum of corrected areas: 22141082

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
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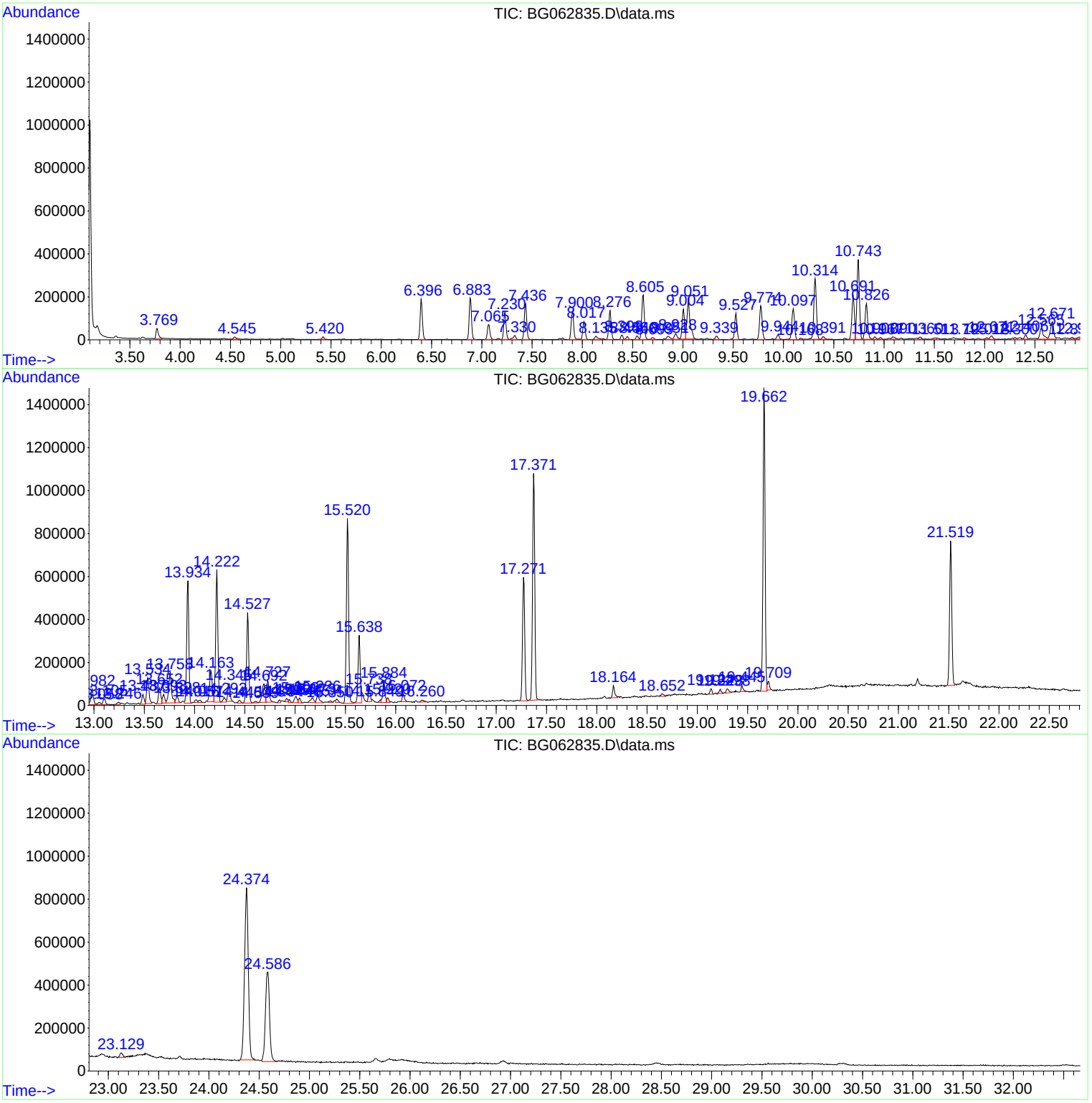
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062835.D
Acq On : 15 Sep 2024 7:08
Operator : JU/RC
Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ3

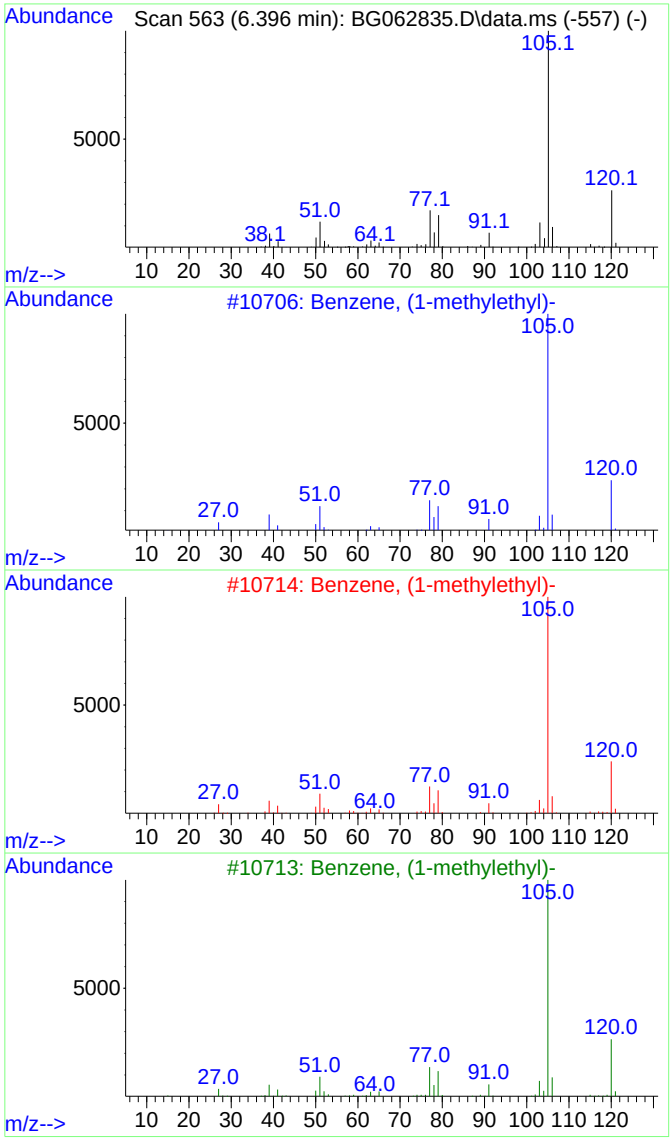
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

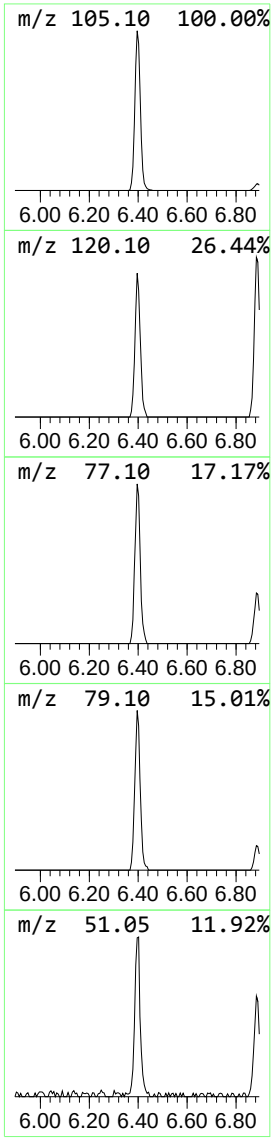
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Benzene, (1-methylethyl)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.396	25.80 ng/ul	292567	1,4-Dichlorobenzene-d4	7.900

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	95
2	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
3	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
4	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062835.D
Acq On : 15 Sep 2024 7:08
Operator : JU/RC
Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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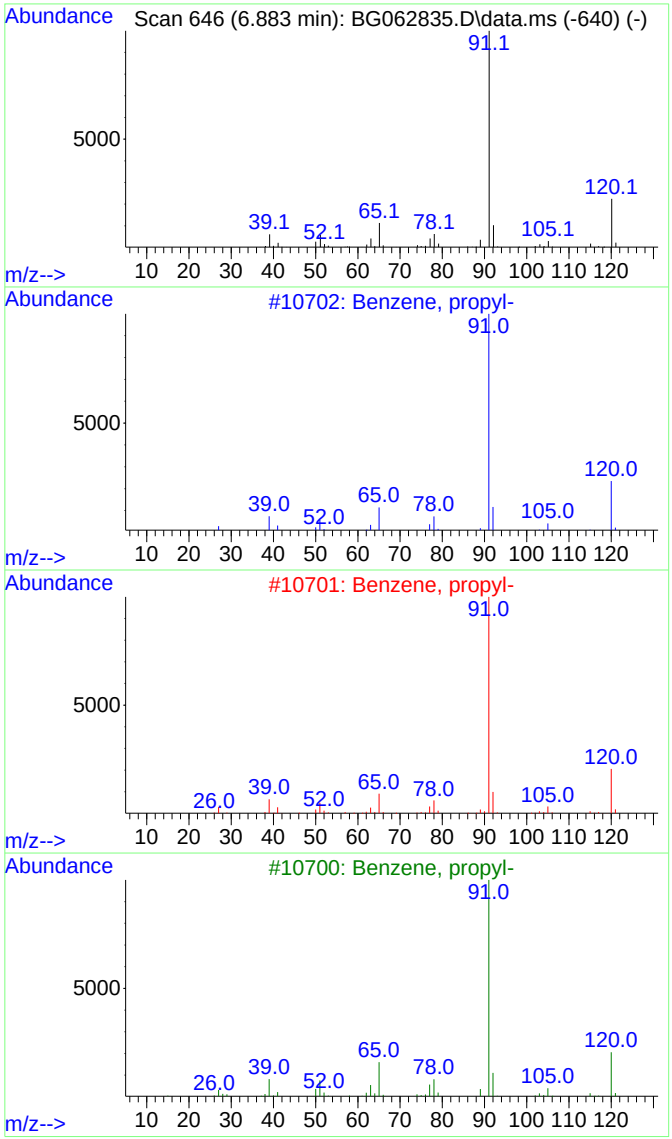
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

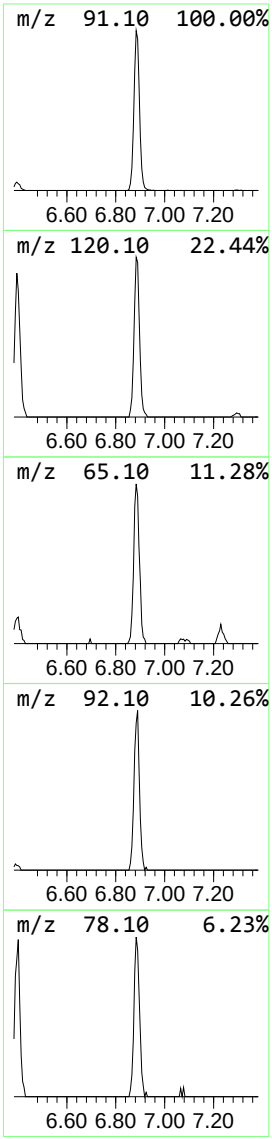
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.883	27.34 ng/ul	310038	1,4-Dichlorobenzene-d4	7.900

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	91
2	Benzene, propyl-	120	C9H12	000103-65-1	91
3	Benzene, propyl-	120	C9H12	000103-65-1	87
4	Benzene, propyl-	120	C9H12	000103-65-1	86
5	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
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ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ3

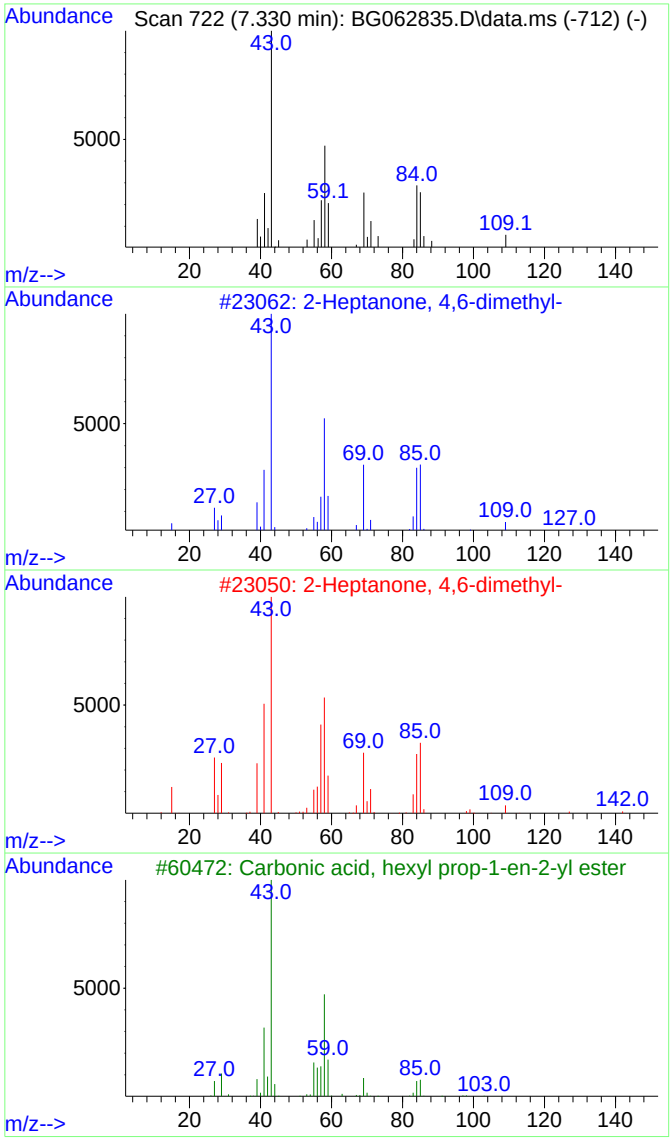
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

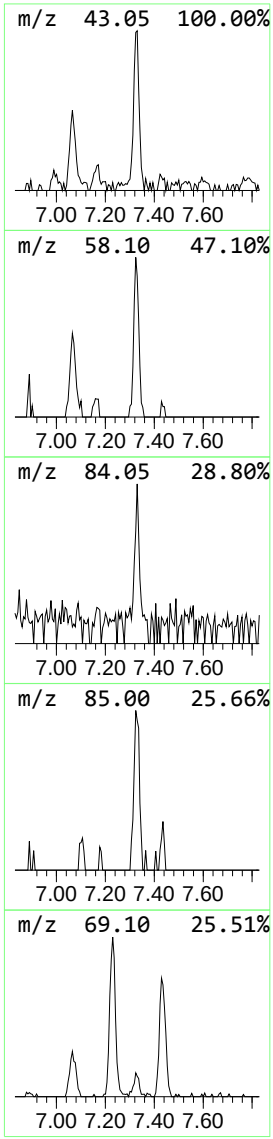
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 2-Heptanone, 4,6-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.330	3.60 ng/ul	40763	1,4-Dichlorobenzene-d4	7.900

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	64
2	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	56
3	Carbonic acid, hexyl prop-1-en-2...	186	C10H18O3	1000382-54-2	53
4	1,3-Dioxane-2-propanol, 2-methyl-	160	C8H16O3	036651-31-7	42
5	2-Hexanone, 4-methyl-	114	C7H14O	000105-42-0	40





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ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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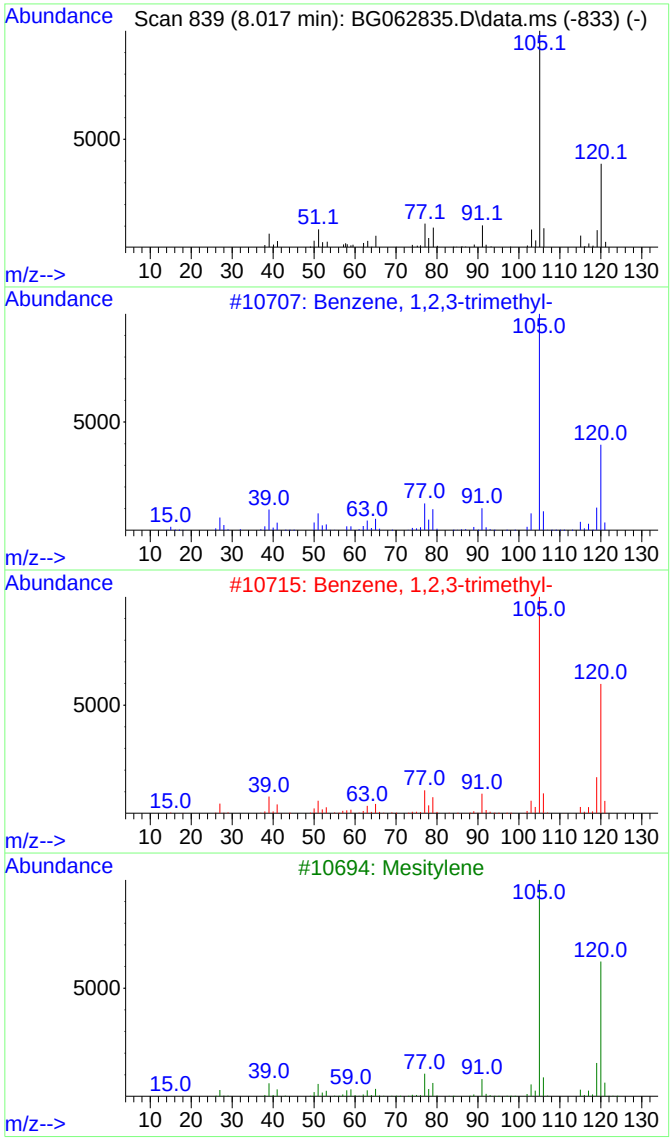
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

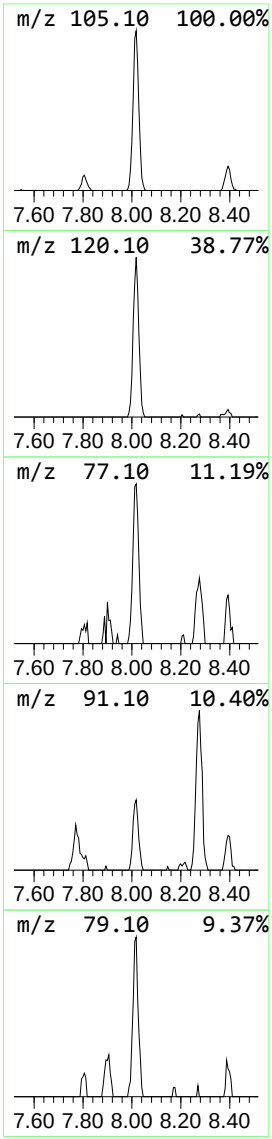
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1,2,3-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.017	12.36 ng/ul	140139	1,4-Dichlorobenzene-d4	7.900

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Mesitylene	120	C9H12	000108-67-8	94
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
5	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062835.D
Acq On : 15 Sep 2024 7:08
Operator : JU/RC
Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ3

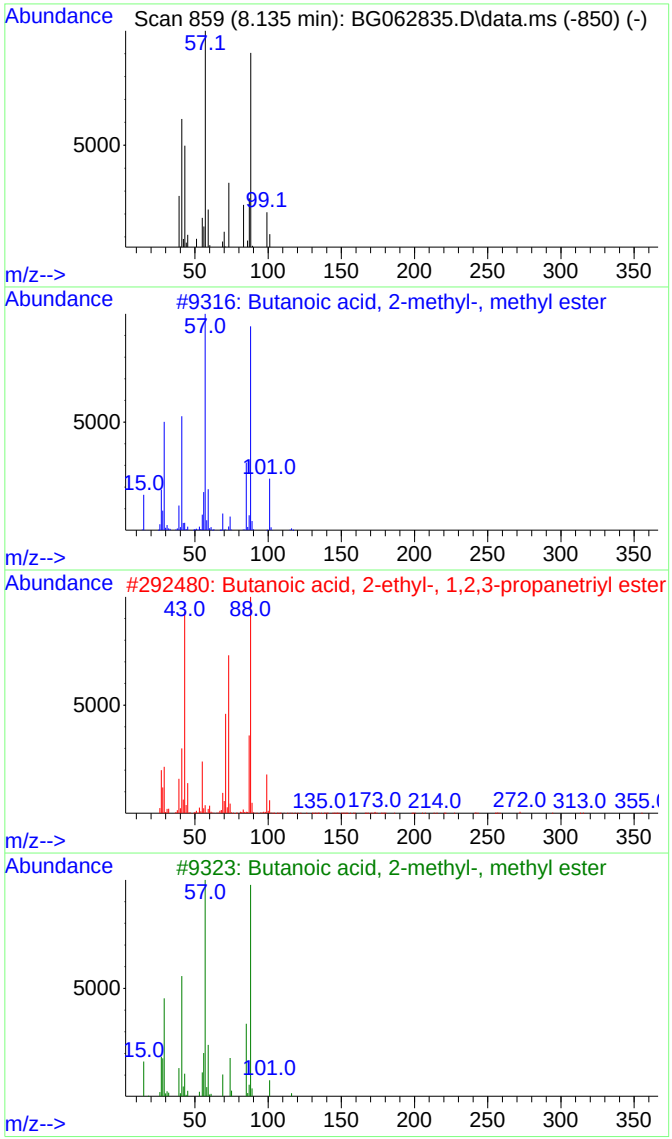
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

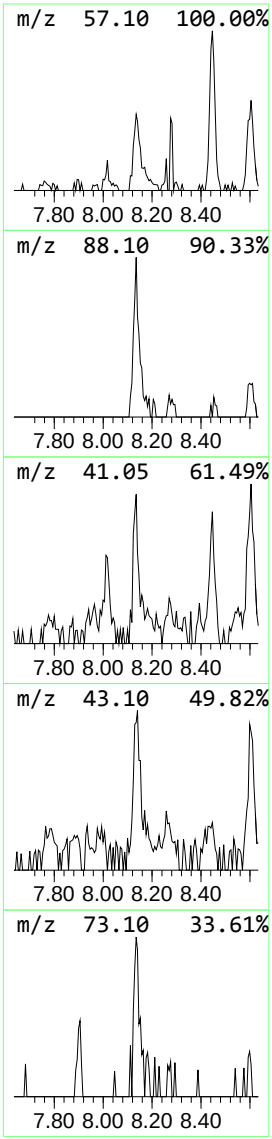
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Butanoic acid, 2-methyl-, m... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.135	3.41 ng/ul	38696	1,4-Dichlorobenzene-d4	7.900

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, 2-methyl-, methyl...	116	C6H12O2	000868-57-5	50
2	Butanoic acid, 2-ethyl-, 1,2,3-p...	386	C21H38O6	056554-54-2	42
3	Butanoic acid, 2-methyl-, methyl...	116	C6H12O2	000868-57-5	40
4	3-Furanol, tetrahydro-	88	C4H8O2	000453-20-3	38
5	Butanoic acid, 2-methyl-, methyl...	116	C6H12O2	000868-57-5	38





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Acq On : 15 Sep 2024 7:08
Operator : JU/RC
Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ3

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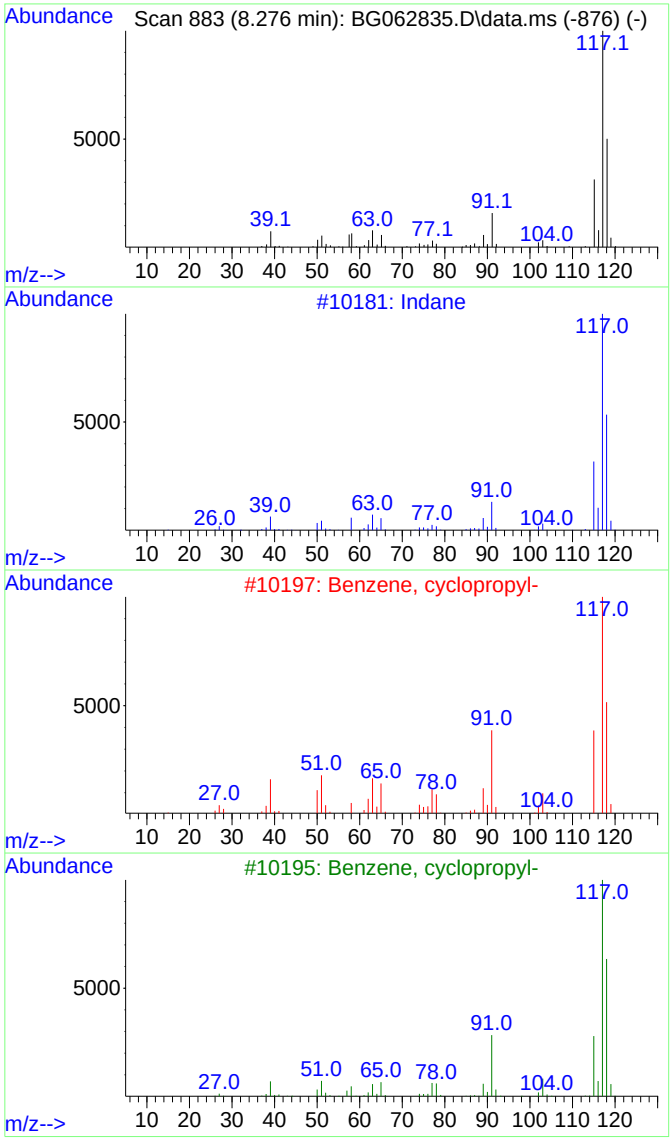
Quant Title : SVOA CALIBRATION

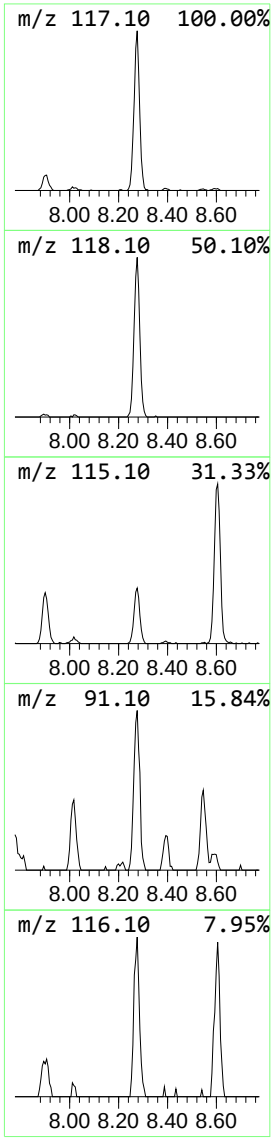
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.276	19.67 ng/ul	223004	1,4-Dichlorobenzene-d4	7.900

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	87
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	83
4	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	72
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	72





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Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ3

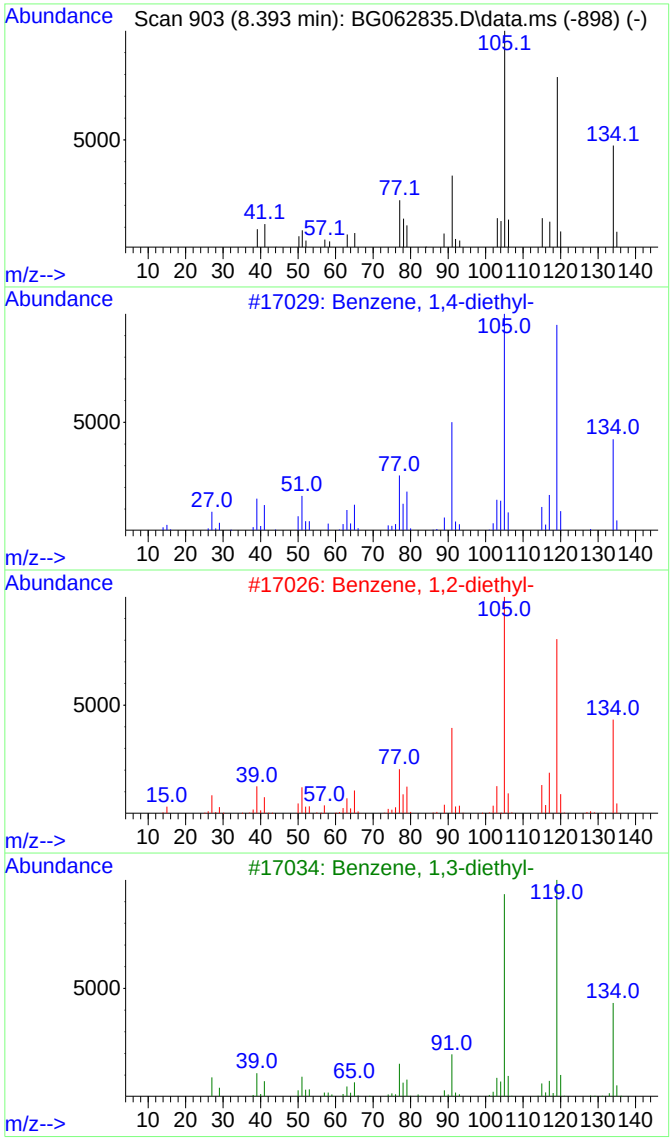
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

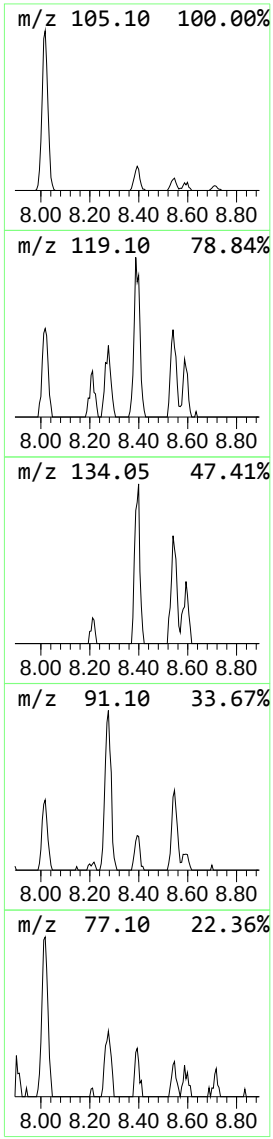
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1,4-diethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.393	3.06 ng/ul	34751	1,4-Dichlorobenzene-d4	7.900

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	95
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	91
4	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	91
5	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	91





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Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ3

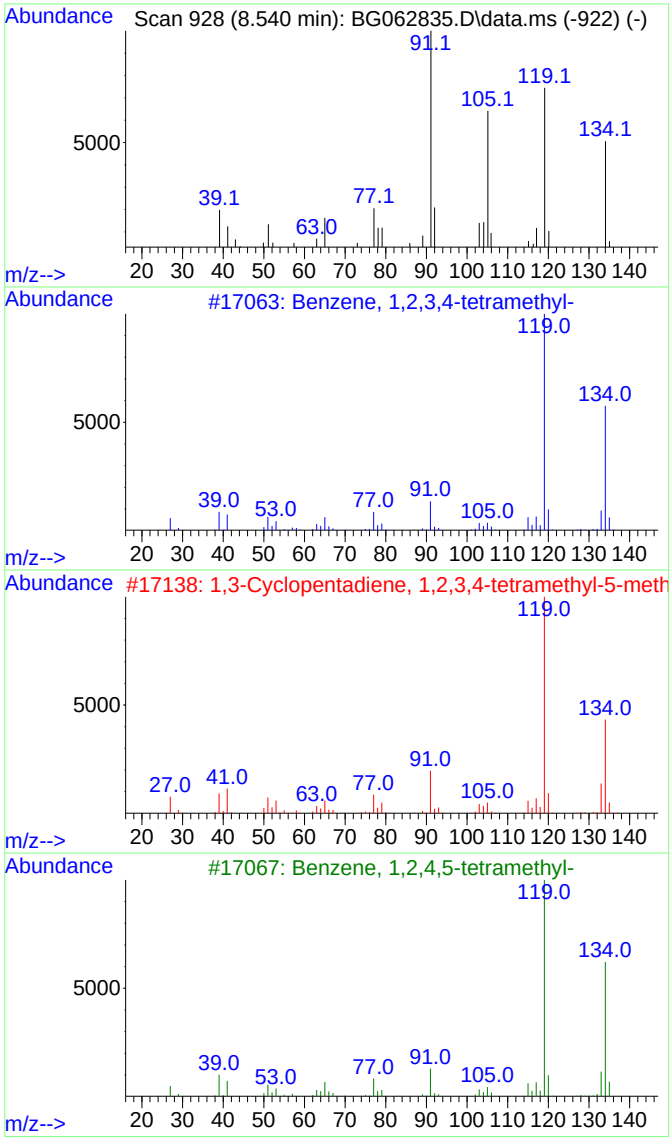
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

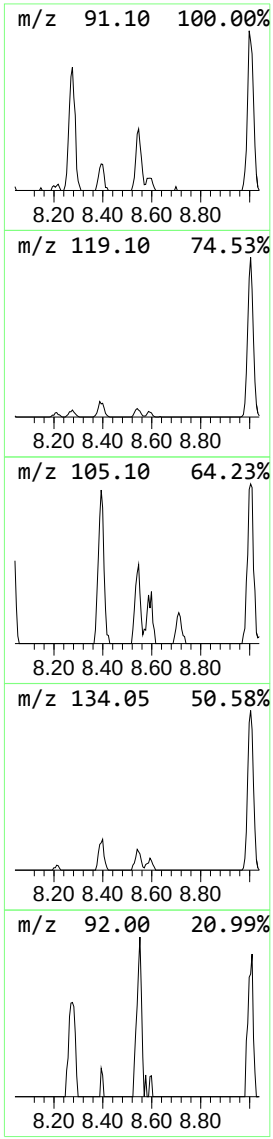
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.540	2.64 ng/ul	29924	1,4-Dichlorobenzene-d4	7.900

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	87
2	1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	81
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	81
4	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	81
5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	81





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Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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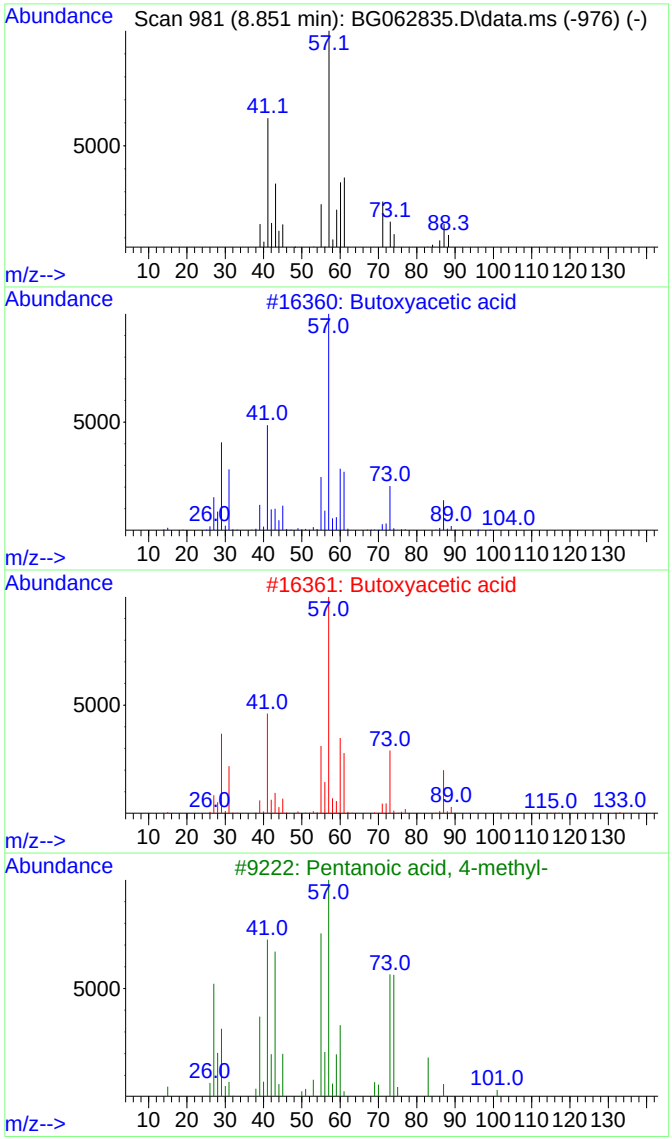
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

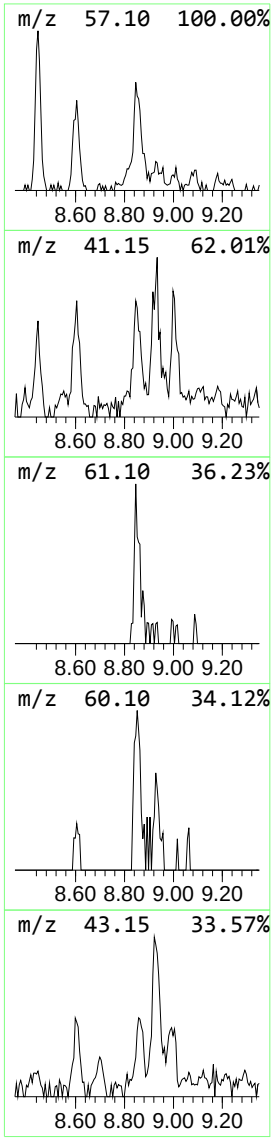
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 unknown-01 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.851	3.08 ng/ul	34930	1,4-Dichlorobenzene-d4	7.900

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butoxyacetic acid	132	C6H12O3	002516-93-0	37
2	Butoxyacetic acid	132	C6H12O3	002516-93-0	25
3	Pentanoic acid, 4-methyl-	116	C6H12O2	000646-07-1	17
4	3-Hydroxybutyric acid, t-butyl e...	160	C8H16O3	090435-23-7	10
5	3-Furanol, tetrahydro-	88	C4H8O2	000453-20-3	9





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Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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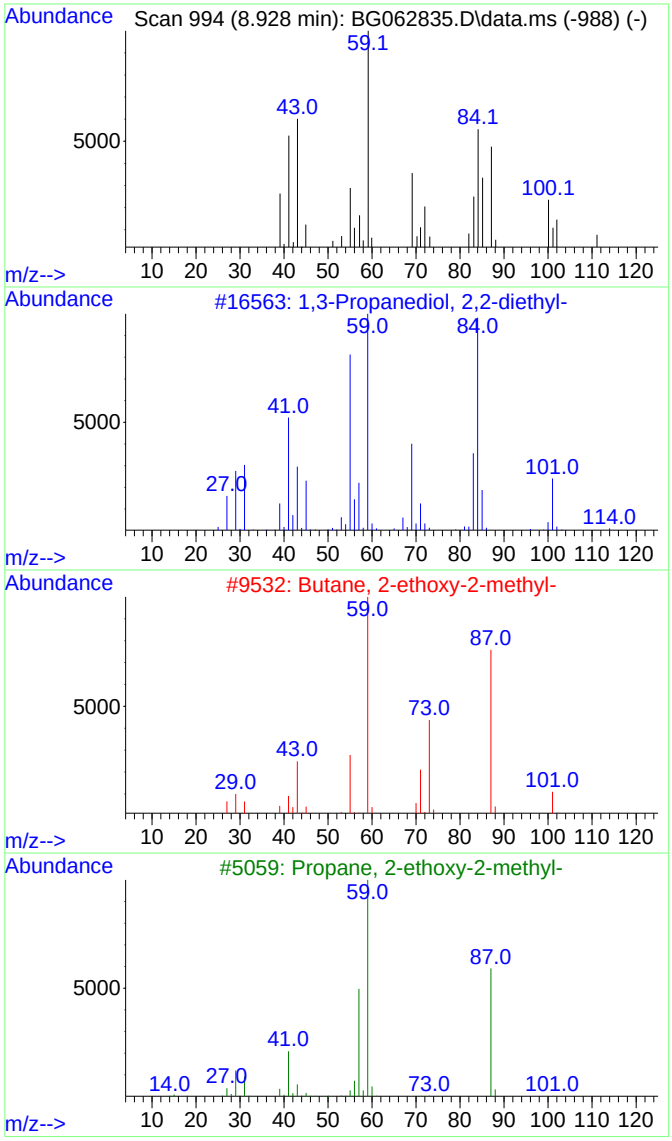
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

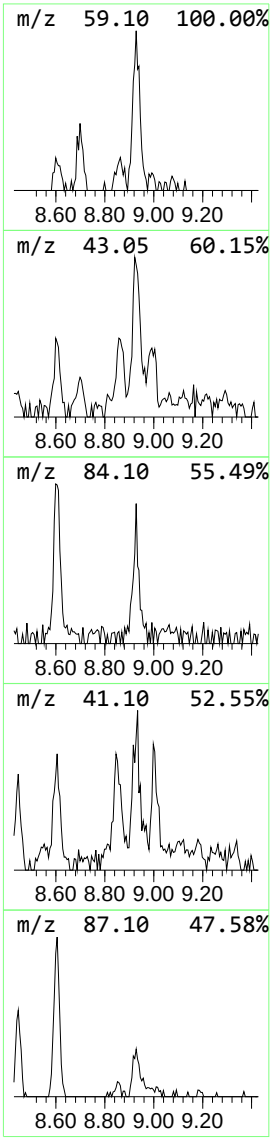
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 unknown-02 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.928	4.68 ng/ul	53071	1,4-Dichlorobenzene-d4	7.900

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Propanediol, 2,2-diethyl-	132	C7H16O2	000115-76-4	43
2	Butane, 2-ethoxy-2-methyl-	116	C7H16O	000919-94-8	38
3	Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	35
4	1,3-Propanediol, 2,2-diethyl-	132	C7H16O2	000115-76-4	25
5	3-Pentanol, 2,2-dimethyl-	116	C7H16O	003970-62-5	23





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Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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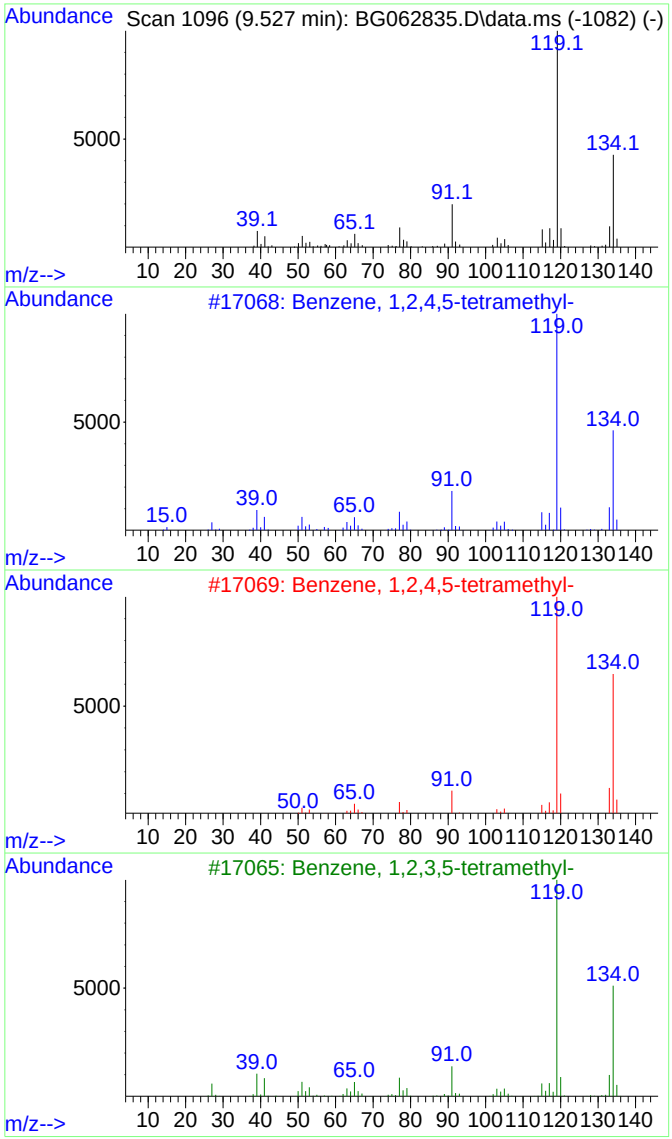
Quant Title : SVOA CALIBRATION

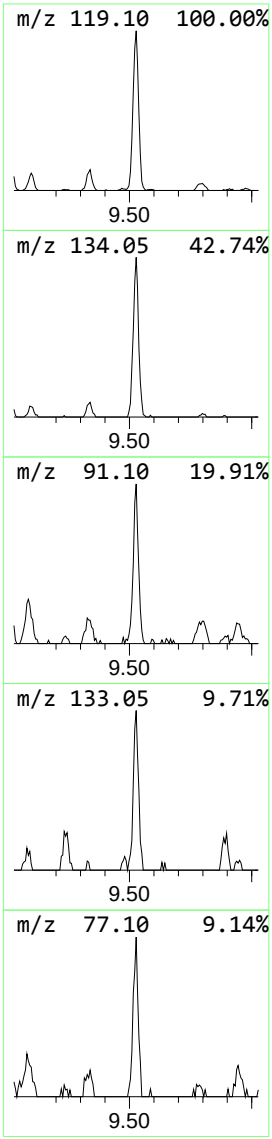
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.527	10.90 ng/ul	203481	Naphthalene-d8	10.691

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	93
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	93
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	93





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Operator : JU/RC
Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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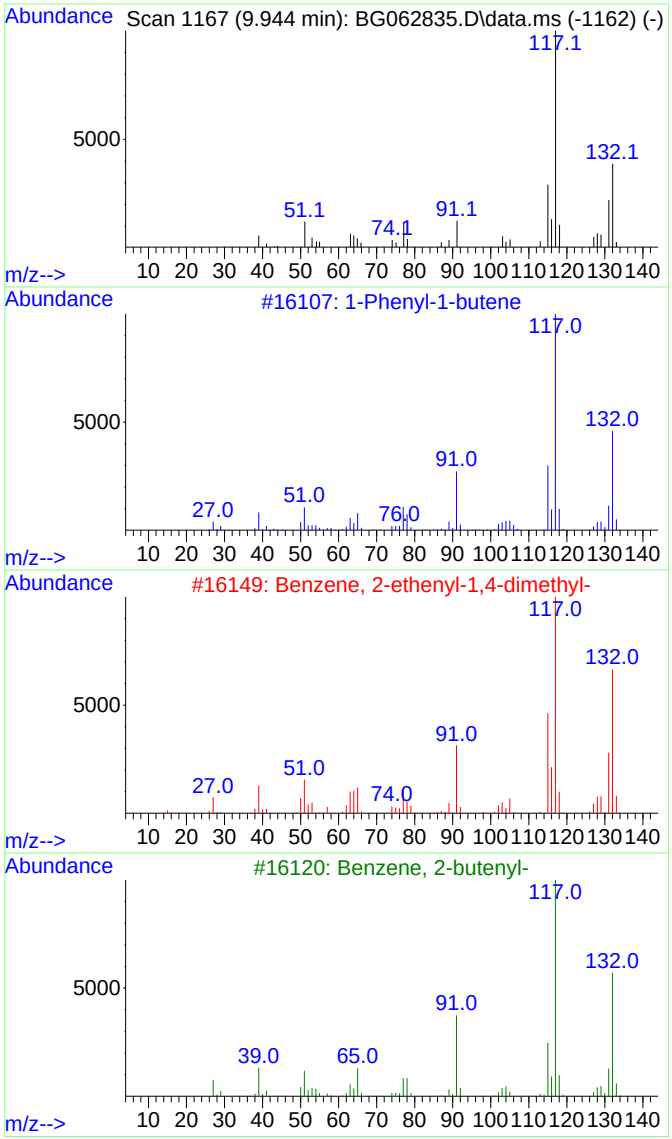
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

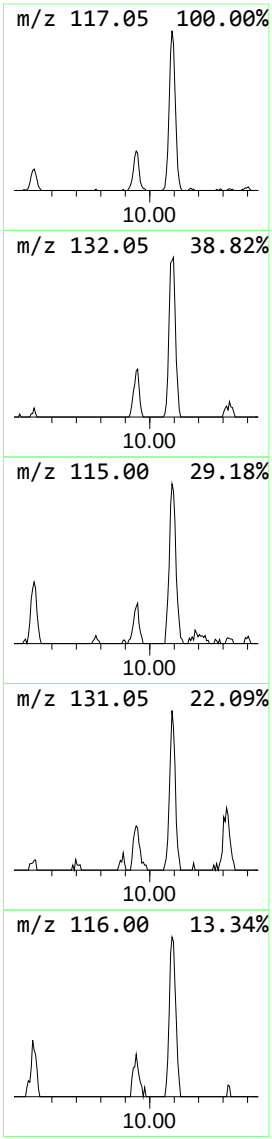
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 1-Phenyl-1-butene Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.944	2.35 ng/ul	43796	Naphthalene-d8	10.691

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	93
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	89
3	Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
5	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	86





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Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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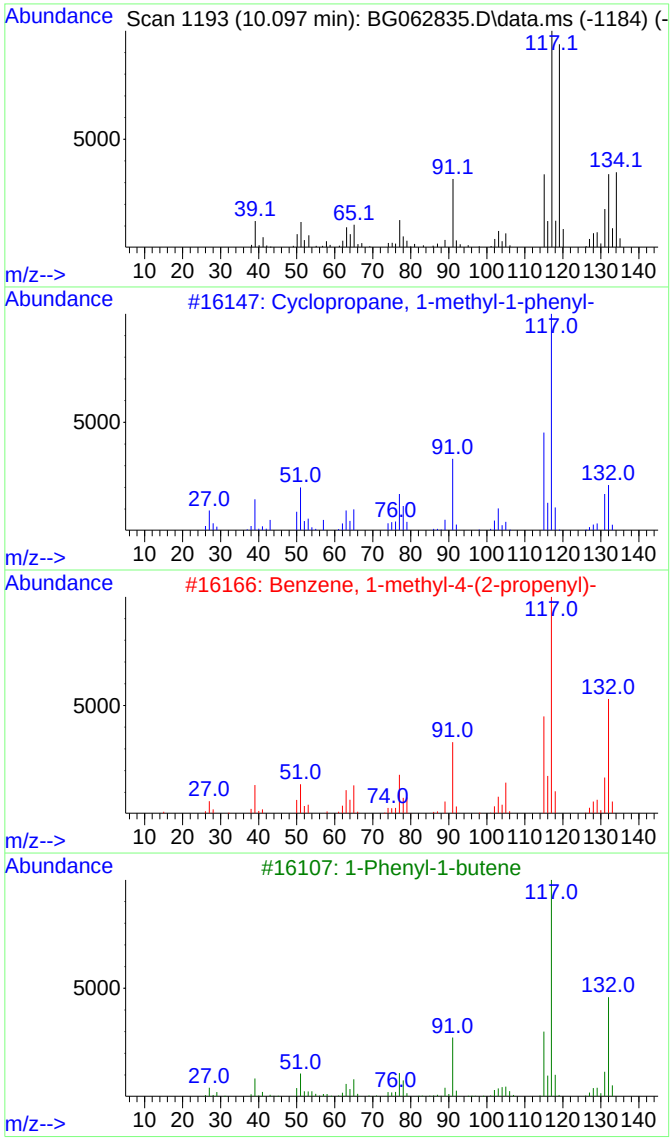
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

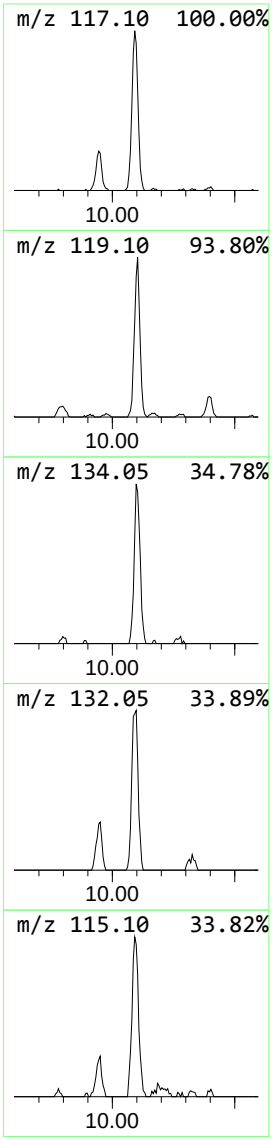
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Cyclopropane, 1-methyl-1-ph... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.097	14.41 ng/ul	269048	Naphthalene-d8	10.691

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, 1-methyl-1-phenyl-	132	C10H12	002214-14-4	70
2	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	70
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	60
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	55
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	55





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Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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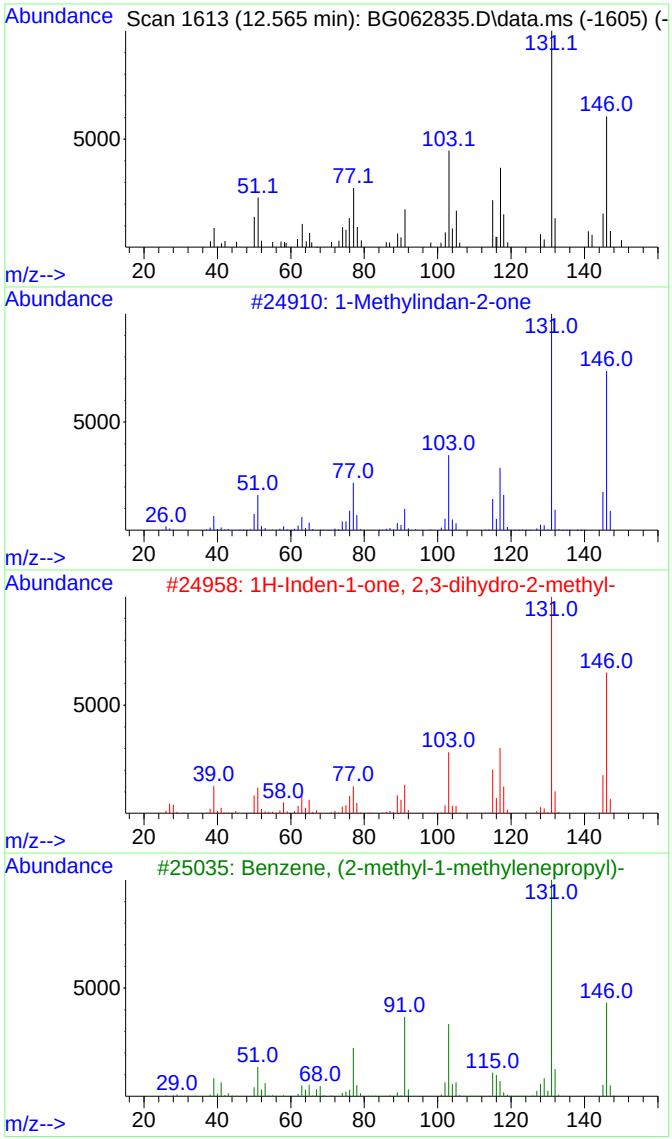
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

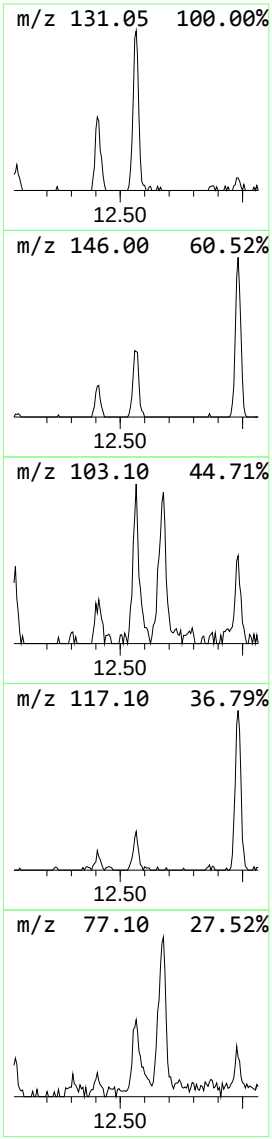
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 1-Methylindan-2-one Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.565	6.60 ng/ul	123177	Naphthalene-d8	10.691

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Methylindan-2-one	146	C10H10O	035587-60-1	97
2	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	91
3	Benzene, (2-methyl-1-methylenepr...	146	C11H14	017498-71-4	87
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	64
5	Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	64





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ALS Vial : 23 Sample Multiplier: 1

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ClientSampleId :
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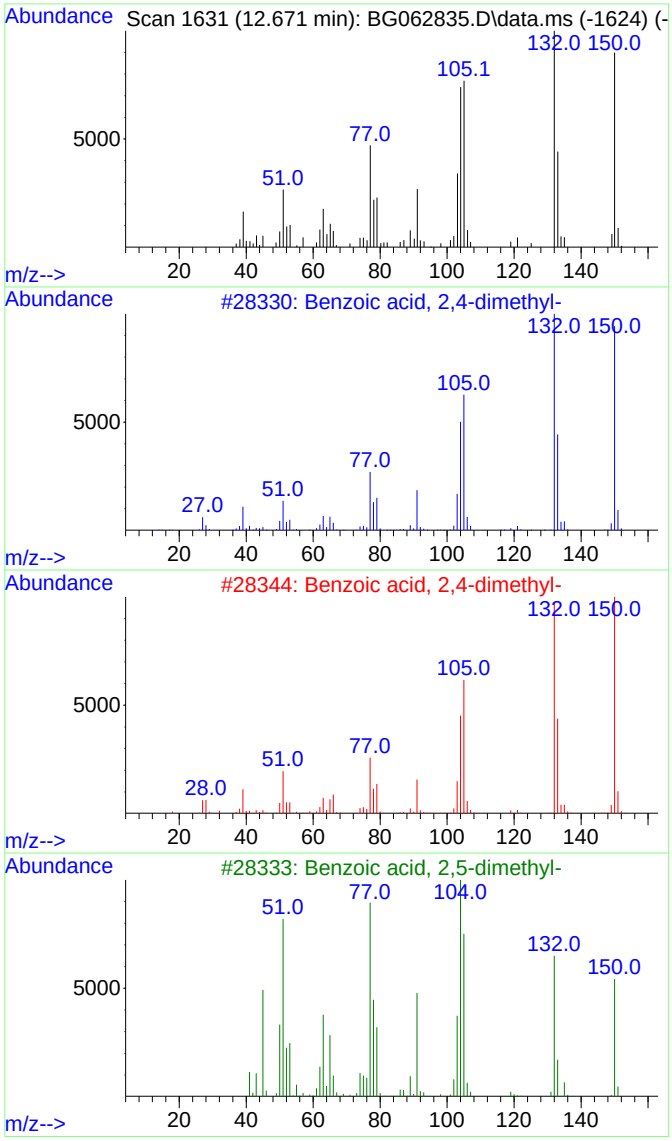
Quant Title : SVOA CALIBRATION

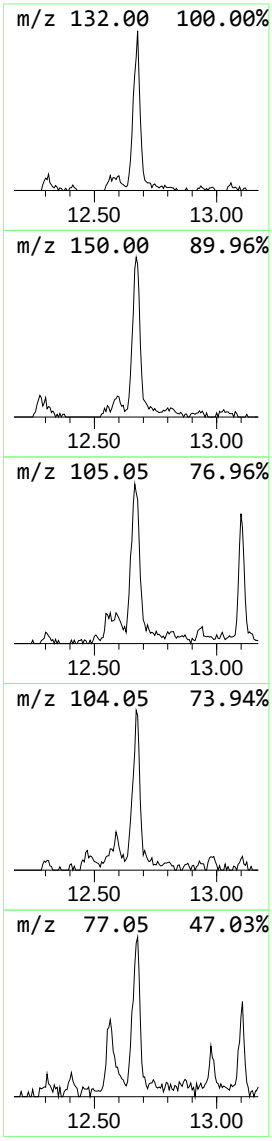
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Benzoic acid, 2,4-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.671	5.68 ng/ul	182667	Acenaphthene-d10	14.527

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 2,4-dimethyl-	150	C9H10O2	000611-01-8	97
2	Benzoic acid, 2,4-dimethyl-	150	C9H10O2	000611-01-8	96
3	Benzoic acid, 2,5-dimethyl-	150	C9H10O2	000610-72-0	94
4	Benzoic acid, 2,5-dimethyl-	150	C9H10O2	000610-72-0	94
5	Benzoic acid, 2,5-dimethyl-	150	C9H10O2	000610-72-0	93





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ClientSampleId :
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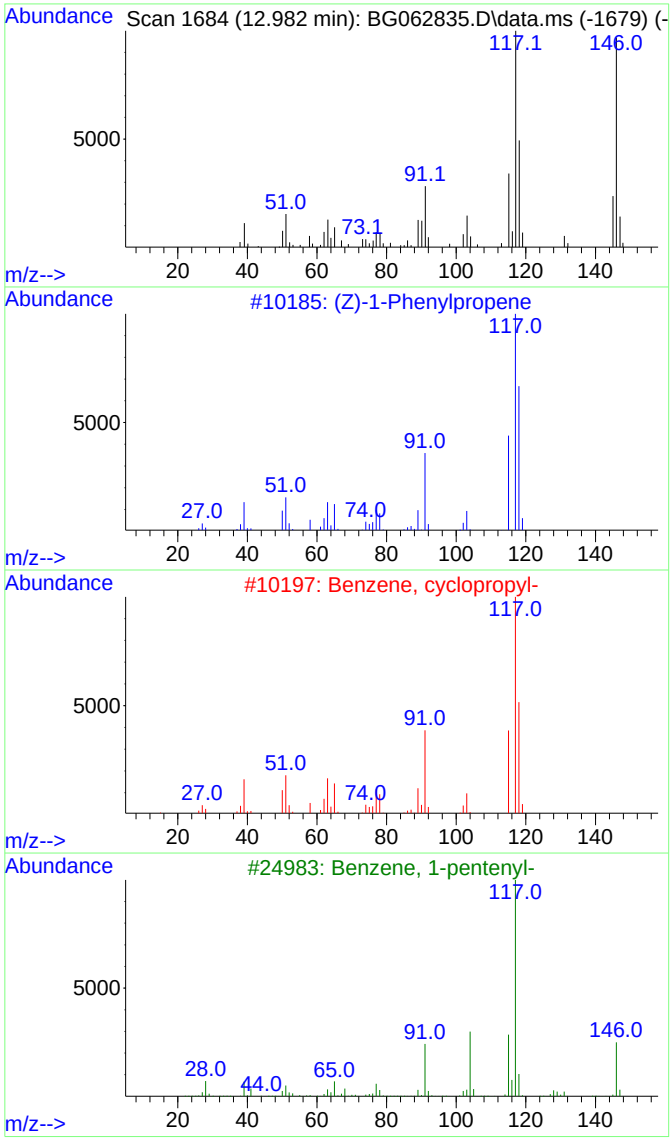
Quant Title : SVOA CALIBRATION

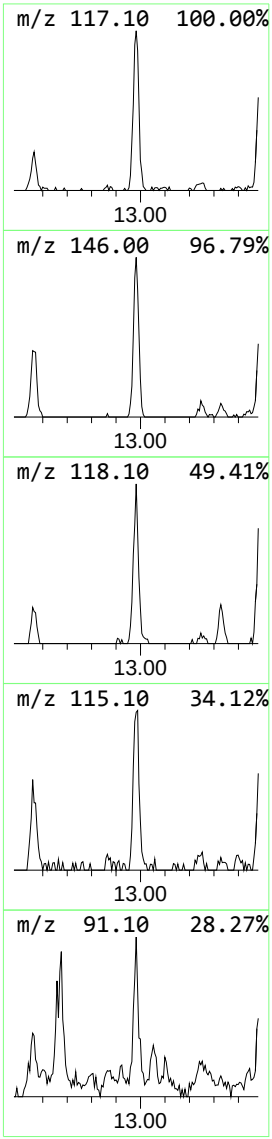
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 (Z)-1-Phenylpropene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.982	3.81 ng/ul	122705	Acenaphthene-d10	14.527

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(Z)-1-Phenylpropene	118	C9H10	000766-90-5	86
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
3	Benzene, 1-pentenyl-	146	C11H14	000826-18-6	62
4	7-Methylindan-1-one	146	C10H10O	039627-61-7	58
5	1H-Indol-4-ylmethylamine	146	C9H10N2	003468-18-6	53





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Data File : BG062835.D
Acq On : 15 Sep 2024 7:08
Operator : JU/RC
Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ3

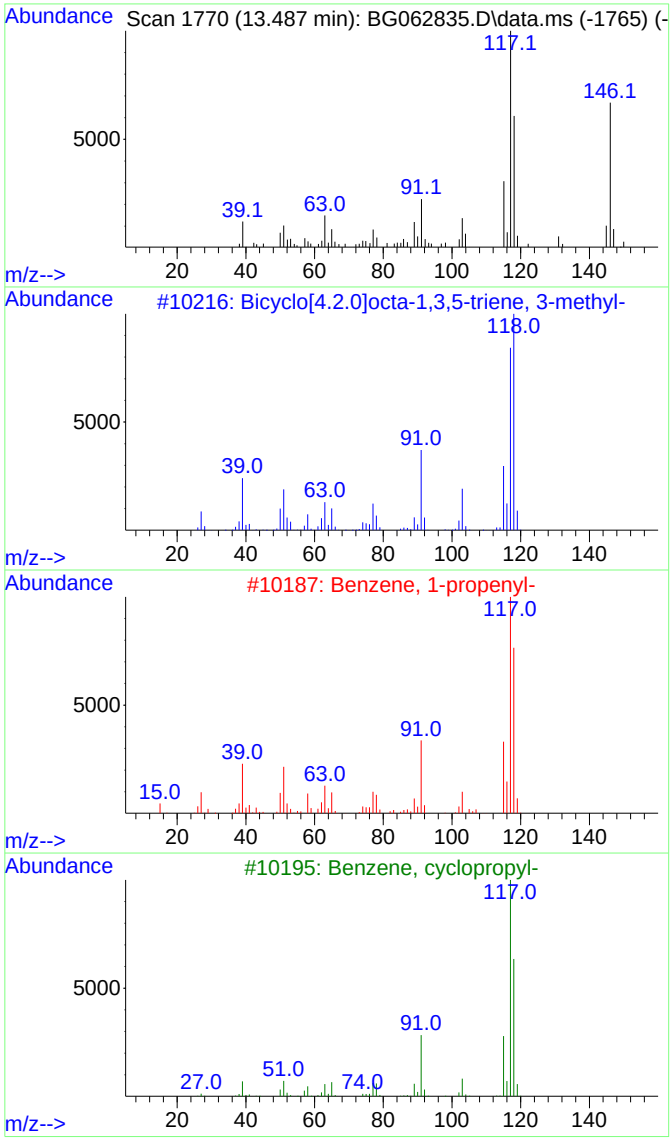
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

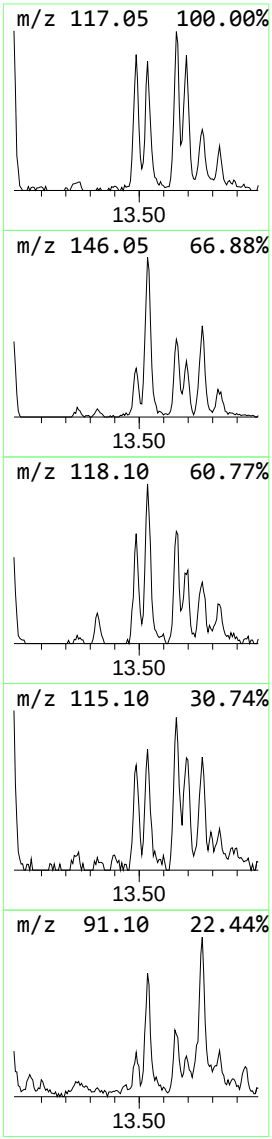
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.487	2.20 ng/ul	70653	Acenaphthene-d10	14.527

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	022250-74-4	80
2	Benzene, 1-propenyl-	118	C9H10	000637-50-3	64
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
4	Benzene, 1-pentenyl-	146	C11H14	000826-18-6	62
5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	60





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Acq On : 15 Sep 2024 7:08
Operator : JU/RC
Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ3

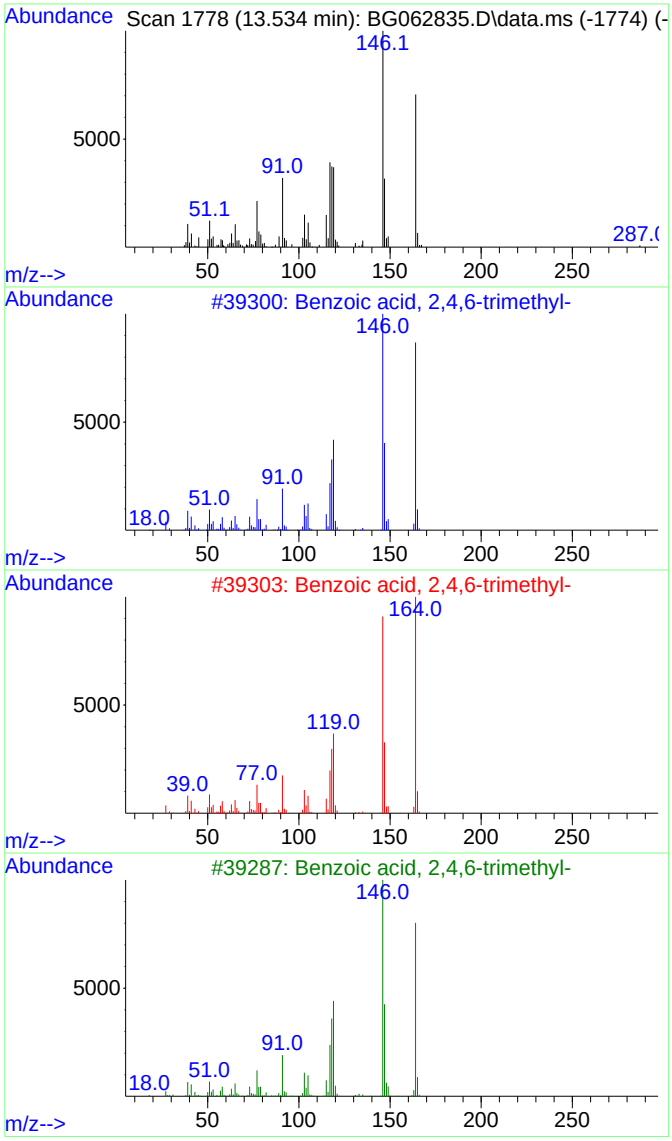
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

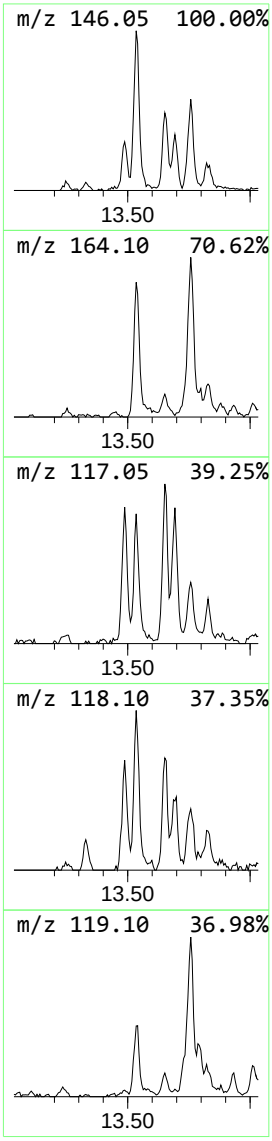
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Benzoic acid, 2,4,6-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.534	5.80 ng/ul	186651	Acenaphthene-d10	14.527

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 2,4,6-trimethyl-	164	C10H12O2	000480-63-7	87
2	Benzoic acid, 2,4,6-trimethyl-	164	C10H12O2	000480-63-7	83
3	Benzoic acid, 2,4,6-trimethyl-	164	C10H12O2	000480-63-7	64
4	4-Oxazolecarboxylic acid, 4,5-di...	233	C13H15NO3	037592-54-4	38
5	3-Ethyl-3-phenyl-1-methylazetidi...	203	C12H13NO2	1000305-99-8	38





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Operator : JU/RC
Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ3

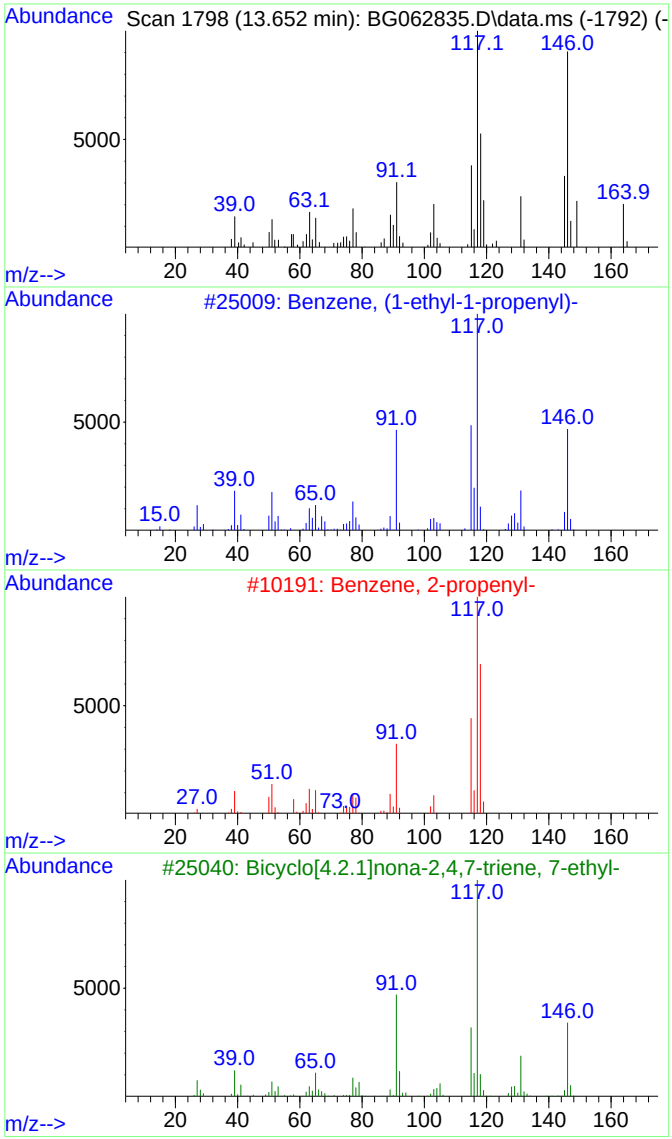
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

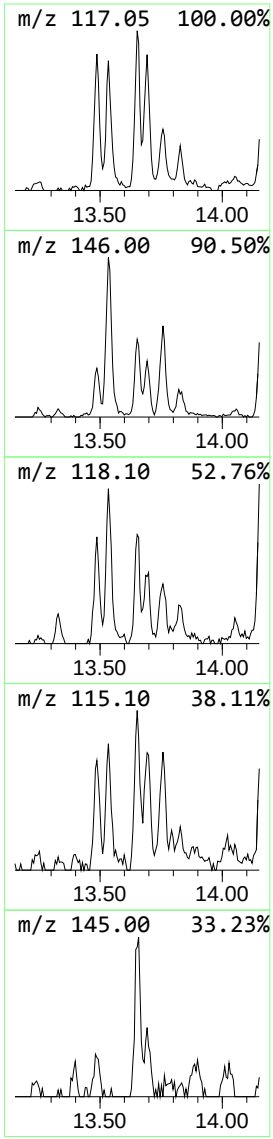
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Benzene, (1-ethyl-1-propenyl)- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.652	3.79 ng/ul	121954	Acenaphthene-d10	14.527

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-ethyl-1-propenyl)-	146	C11H14	004701-36-4	58
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	52
3	Bicyclo[4.2.1]nona-2,4,7-triene,...	146	C11H14	1000164-42-6	49
4	Benzene, 1-pentenyl-	146	C11H14	000826-18-6	49
5	1H-Imidazole, 4,5-dihydro-2-phenyl-	146	C9H10N2	000936-49-2	45





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Operator : JU/RC
Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ3

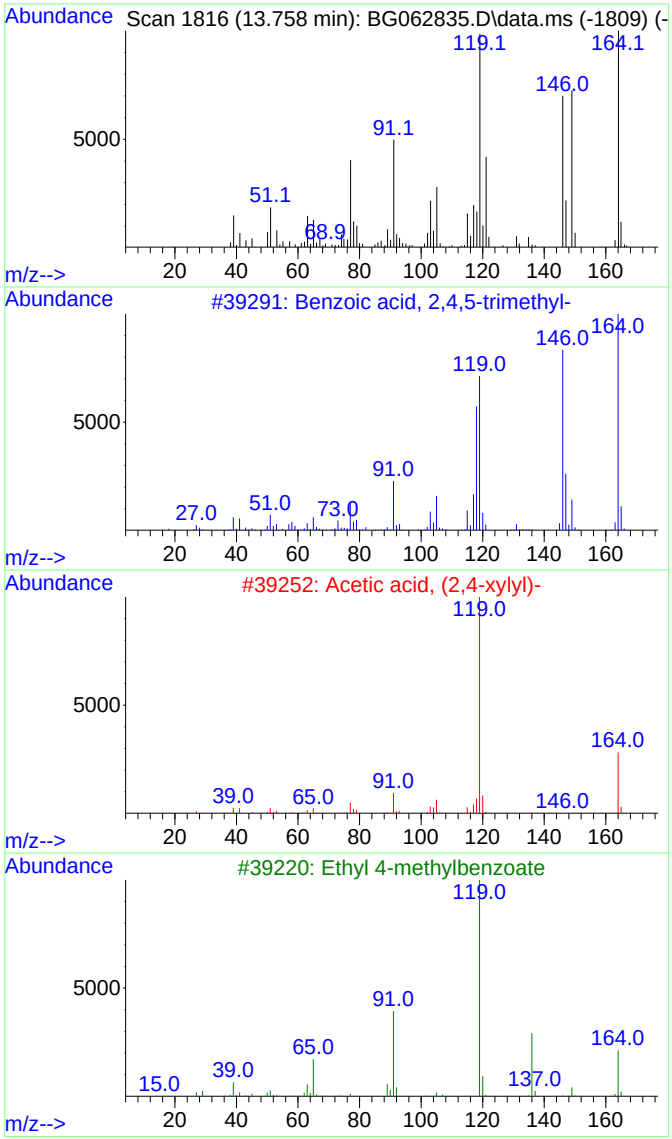
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

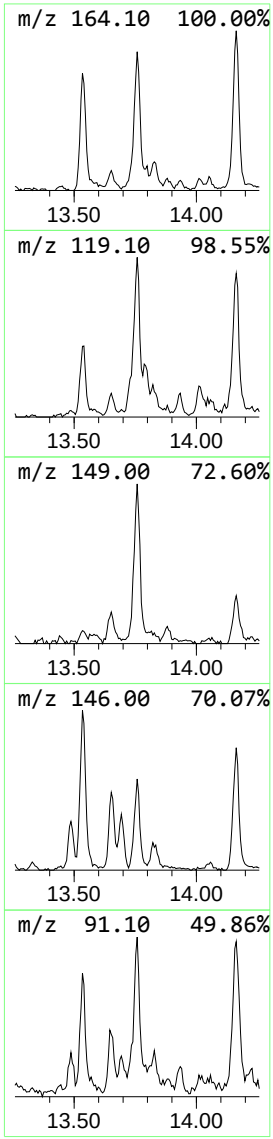
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Benzoic acid, 2,4,5-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.758	8.57 ng/ul	275695	Acenaphthene-d10	14.527

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 2,4,5-trimethyl-	164	C10H12O2	000528-90-5	91
2	Acetic acid, (2,4-xylyl)-	164	C10H12O2	006331-04-0	60
3	Ethyl 4-methylbenzoate	164	C10H12O2	000094-08-6	50
4	Ethyl m-methylbenzoate	164	C10H12O2	000120-33-2	45
5	Ethyl m-methylbenzoate	164	C10H12O2	000120-33-2	45





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Acq On : 15 Sep 2024 7:08
Operator : JU/RC
Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

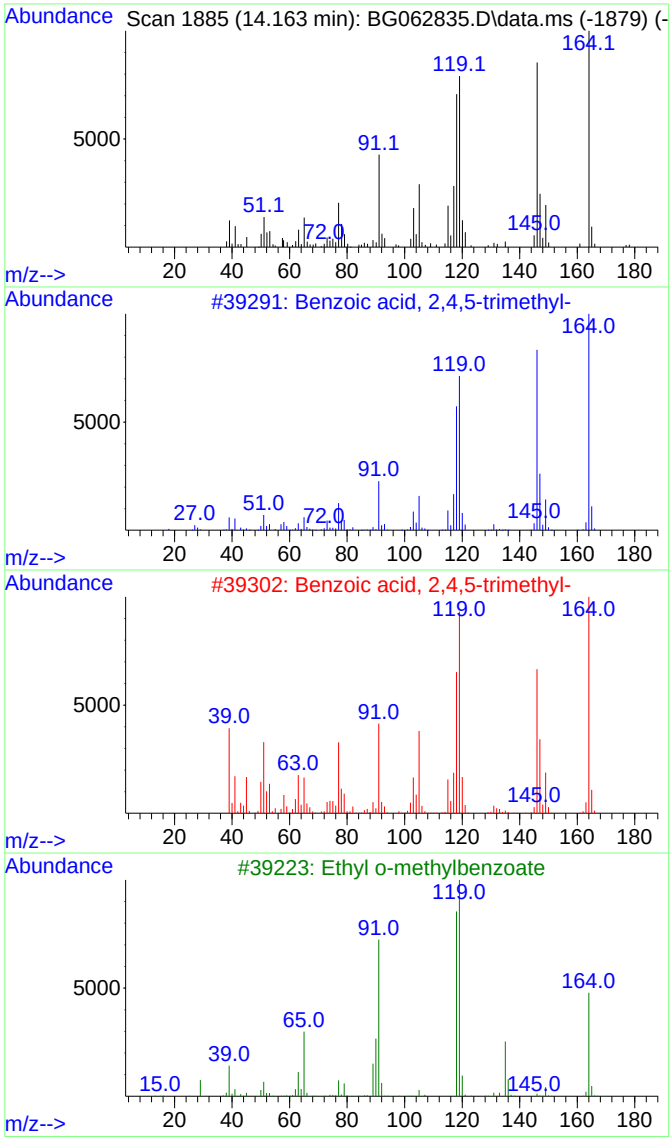
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

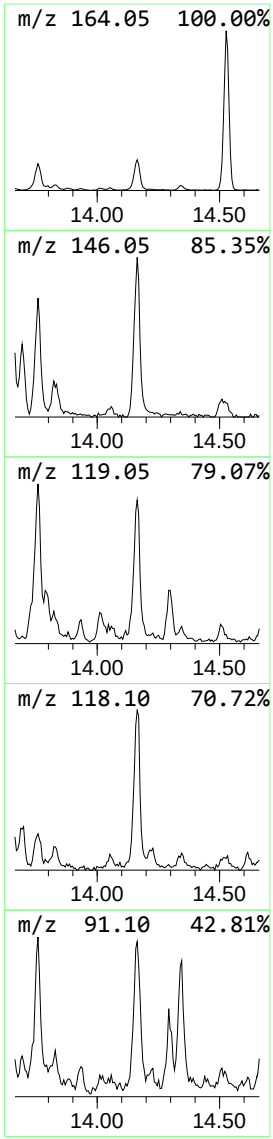
****

Peak Number 21 Ethyl o-methylbenzoate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.163	8.29 ng/ul	266685	Acenaphthene-d10	14.527

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 2,4,5-trimethyl-	164	C10H12O2	000528-90-5	90
2	Benzoic acid, 2,4,5-trimethyl-	164	C10H12O2	000528-90-5	74
3	Ethyl o-methylbenzoate	164	C10H12O2	000087-24-1	53
4	Ethyl o-methylbenzoate	164	C10H12O2	000087-24-1	43
5	3,5-Dimethylphenylacetic acid	164	C10H12O2	042288-46-0	43





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Acq On : 15 Sep 2024 7:08
Operator : JU/RC
Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ3

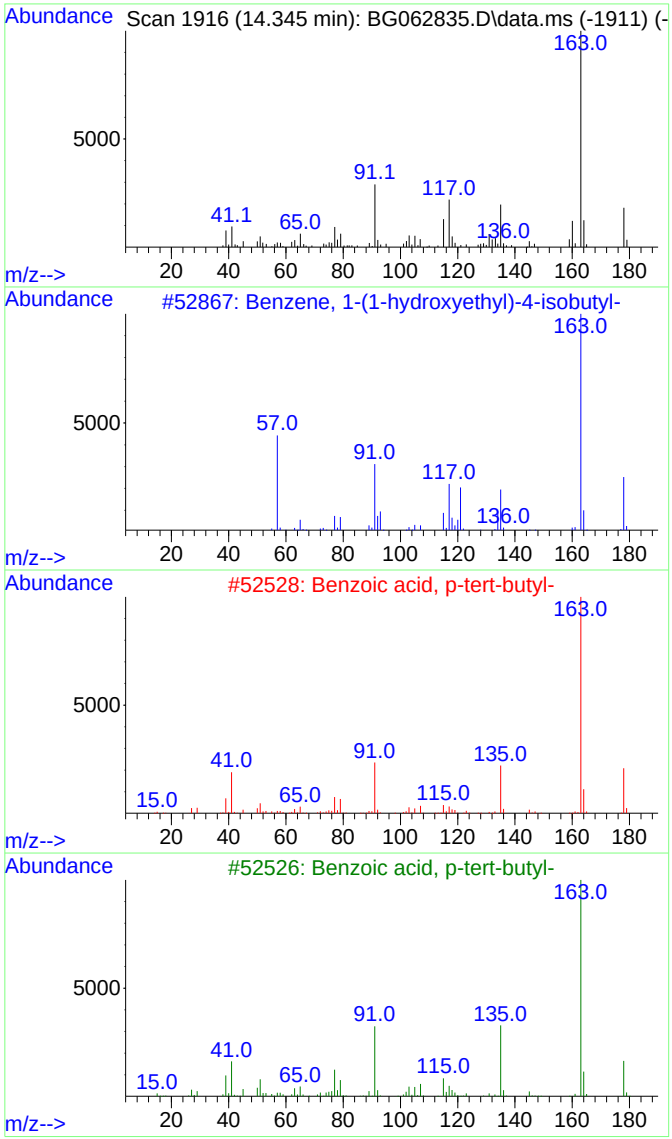
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

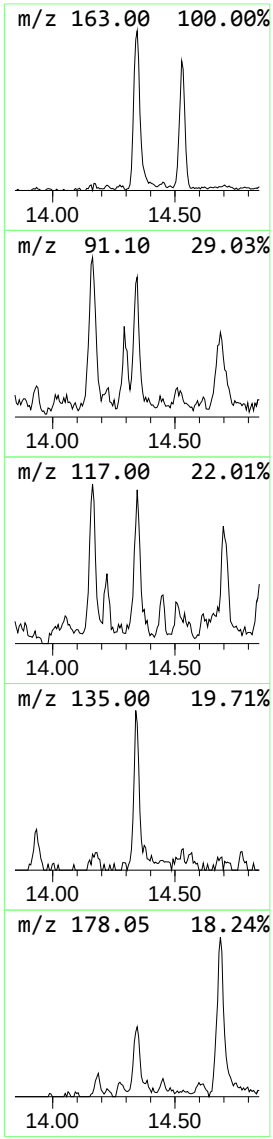
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 Benzene, 1-(1-hydroxyethyl)... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.345	4.09 ng/ul	131724	Acenaphthene-d10	14.527

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-(1-hydroxyethyl)-4-is...	178	C12H18O	040150-92-3	83
2	Benzoic acid, p-tert-butyl-	178	C11H14O2	000098-73-7	76
3	Benzoic acid, p-tert-butyl-	178	C11H14O2	000098-73-7	76
4	Benzene, 1-(1,1-dimethylethyl)-2...	178	C12H18O	000088-40-4	68
5	Benzene, 1-(1,1-dimethylethyl)-2...	178	C12H18O	000088-40-4	68





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Acq On : 15 Sep 2024 7:08
Operator : JU/RC
Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ3

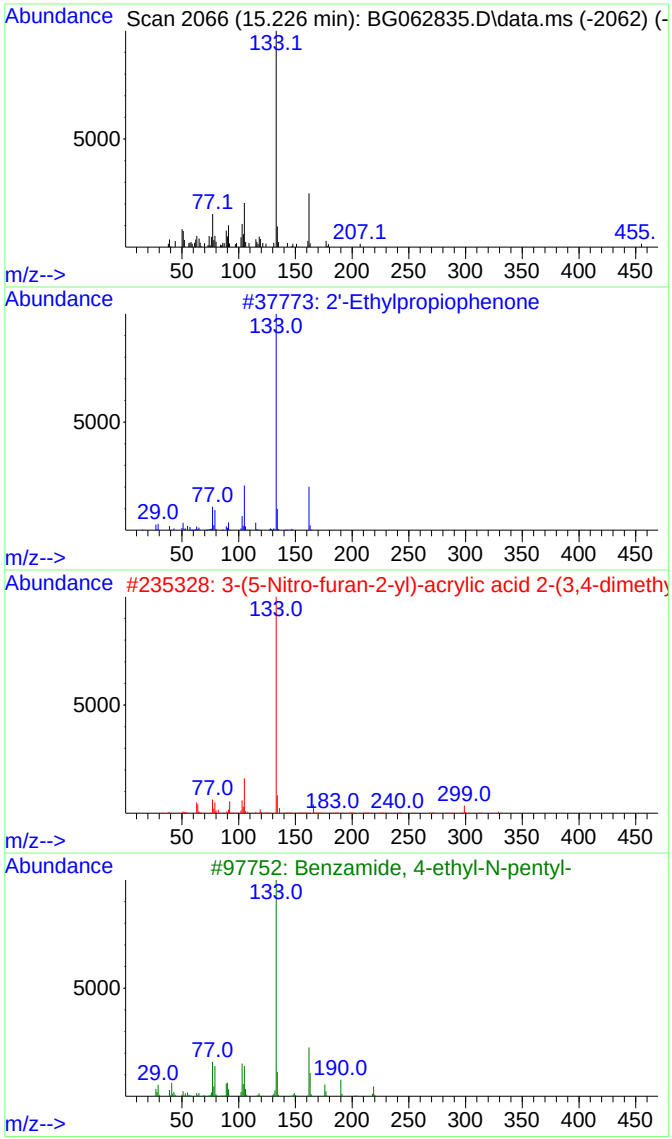
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

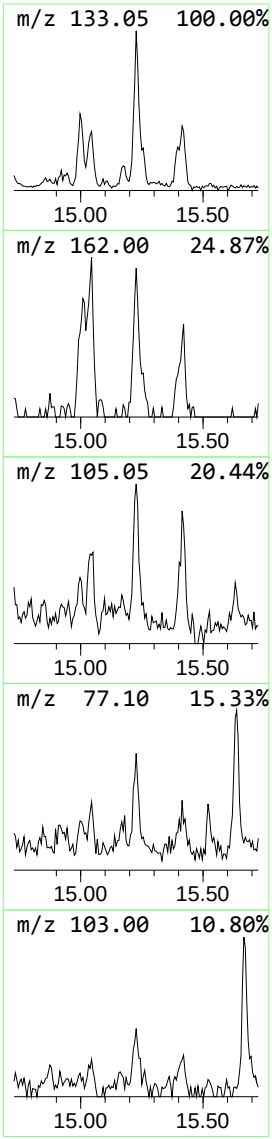
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 2'-Ethylpropiophenone Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.226	2.14 ng/ul	68966	Acenaphthene-d10	14.527

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2'-Ethylpropiophenone	162	C11H14O	016819-79-7	80
2	3-(5-Nitro-furan-2-yl)-acrylic a...	329	C17H15NO6	1000275-46-2	64
3	Benzamide, 4-ethyl-N-pentyl-	219	C14H21NO	1000419-66-4	59
4	3,4-Dimethylmethcathinone, N-tri...	287	C14H16F3NO2	1000448-10-9	59
5	(Benzothiazol-2-ylsulfanyl)aceti...	371	C19H17NO3S2	1000294-36-1	59





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Acq On : 15 Sep 2024 7:08
Operator : JU/RC
Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ3

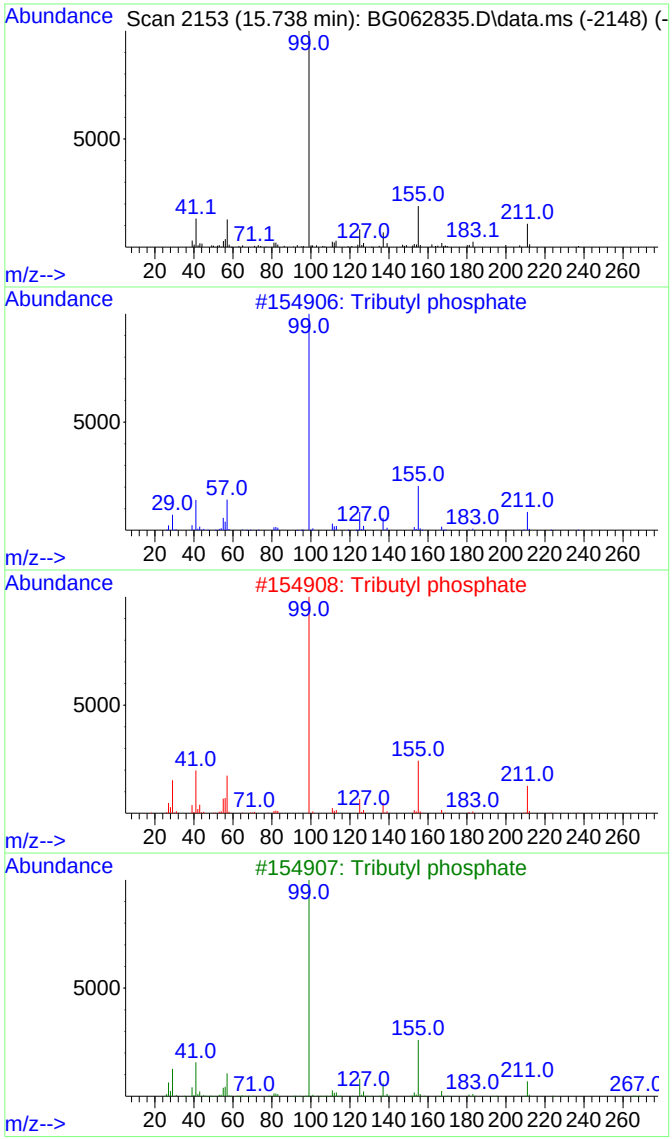
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

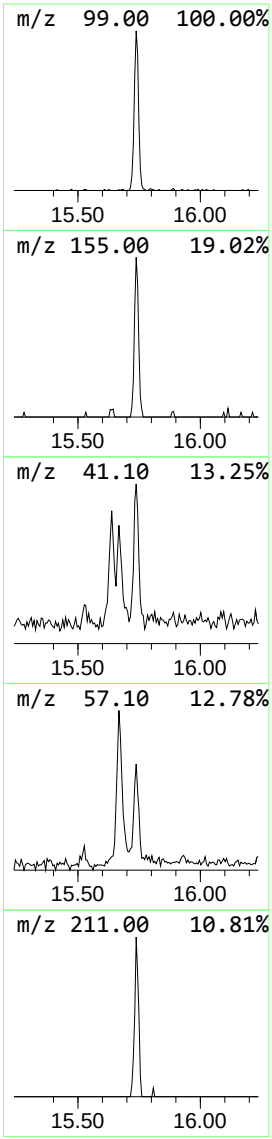
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Tributyl phosphate Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.738	2.64 ng/ul	84985	Acenaphthene-d10	14.527

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tributyl phosphate	266	C12H27O4P	000126-73-8	91
2	Tributyl phosphate	266	C12H27O4P	000126-73-8	90
3	Tributyl phosphate	266	C12H27O4P	000126-73-8	86
4	Tributyl phosphate	266	C12H27O4P	000126-73-8	83
5	Tributyl phosphate	266	C12H27O4P	000126-73-8	47





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Acq On : 15 Sep 2024 7:08
Operator : JU/RC
Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

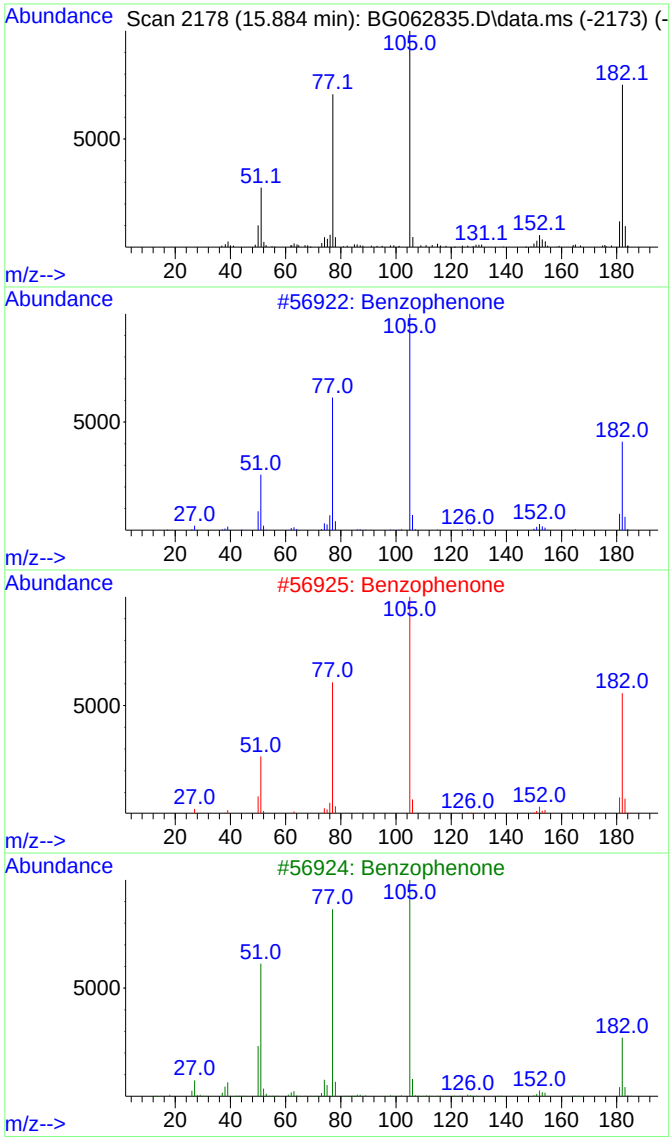
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

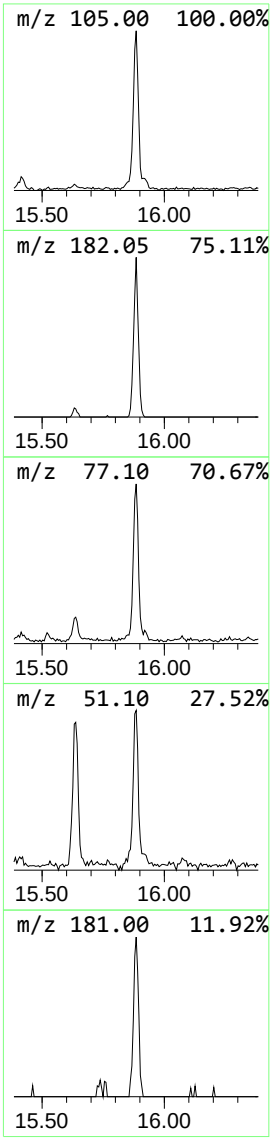
****

Peak Number 25 Benzophenone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.884	4.70 ng/ul	151108	Acenaphthene-d10	14.527

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	94
2	Benzophenone	182	C13H10O	000119-61-9	94
3	Benzophenone	182	C13H10O	000119-61-9	94
4	Benzophenone	182	C13H10O	000119-61-9	91
5	Benzophenone	182	C13H10O	000119-61-9	87





Tentatively Identified Compound (LSC) summary

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Data File : BG062835.D
Acq On : 15 Sep 2024 7:08
Operator : JU/RC
Sample : P3997-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AQ3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.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, (1-met...	6.396	25.8	ng/ul	292567	1	7.900	226762	20.0
Benzene, propyl-	6.883	27.3	ng/ul	310038	1	7.900	226762	20.0
2-Heptanone, 4,...	7.330	3.6	ng/ul	40763	1	7.900	226762	20.0
Benzene, 1,2,3-...	8.017	12.4	ng/ul	140139	1	7.900	226762	20.0
Butanoic acid, ...	8.135	3.4	ng/ul	38696	1	7.900	226762	20.0
Indane	8.276	19.7	ng/ul	223004	1	7.900	226762	20.0
Benzene, 1,4-di...	8.393	3.1	ng/ul	34751	1	7.900	226762	20.0
Benzene, 1,2,3,...	8.540	2.6	ng/ul	29924	1	7.900	226762	20.0
unknown-01	8.851	3.1	ng/ul	34930	1	7.900	226762	20.0
unknown-02	8.928	4.7	ng/ul	53071	1	7.900	226762	20.0
Benzene, 1,2,4,...	9.527	10.9	ng/ul	203481	2	10.691	373469	20.0
1-Phenyl-1-butene	9.944	2.4	ng/ul	43796	2	10.691	373469	20.0
Cyclopropane, 1...	10.097	14.4	ng/ul	269048	2	10.691	373469	20.0
1-Methylindan-2...	12.565	6.6	ng/ul	123177	2	10.691	373469	20.0
Benzoic acid, 2...	12.671	5.7	ng/ul	182667	3	14.527	643492	20.0
(Z)-1-Phenylpro...	12.982	3.8	ng/ul	122705	3	14.527	643492	20.0
Bicyclo[4.2.0]o...	13.487	2.2	ng/ul	70653	3	14.527	643492	20.0
Benzoic acid, 2...	13.534	5.8	ng/ul	186651	3	14.527	643492	20.0
Benzene, (1-eth...	13.652	3.8	ng/ul	121954	3	14.527	643492	20.0
Benzoic acid, 2...	13.758	8.6	ng/ul	275695	3	14.527	643492	20.0
Ethyl o-methylb...	14.163	8.3	ng/ul	266685	3	14.527	643492	20.0
Benzene, 1-(1-h...	14.345	4.1	ng/ul	131724	3	14.527	643492	20.0
2'-Ethylpropiop...	15.226	2.1	ng/ul	68966	3	14.527	643492	20.0
Tributyl phosphate	15.738	2.6	ng/ul	84985	3	14.527	643492	20.0
Benzophenone	15.884	4.7	ng/ul	151108	3	14.527	643492	20.0