LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

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-BG090721.M
Title : SVOA CALIBRATION

Signal : TIC: BG050243.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.065	169	175	182	rBV2	21574	35367	0.78%	0.240%
2	5.416	401	405	411	rVB	22132	32843	0.72%	0.223%
3	5.551	422	428	438	rBV4	32265	86705	1.91%	0.588%
4	5.927	487	492	498	rVB2	29051	42793	0.94%	0.290%
5	7.155	694	701	707	rVV5	8323	15636	0.34%	0.106%
6	7.278	716	722	726	rBV4	8346	14831	0.33%	0.101%
7	7.490	751	758	764	rVB2	9815	16608	0.37%	0.113%
8	7.584	767	774	781	rBV3	22633	37418	0.83%	0.254%
9	7.666	781	788	798	rVB	46604	81339	1.79%	0.552%
10	7.942	829	835	844	rBV	104583	178204	3.93%	1.209%
11	8.060	850	855	859	rVB5	9280	14059	0.31%	0.095%
12	8.318	892	899	909	rVB	140212	257792	5.69%	1.749%
13	8.483	920	927	935	rBV	256634	449331	9.91%	3.048%
14	8.682	956	961	967	rVV3	15632	27252	0.60%	0.185%
15	8.806	975	982	988	rBV	28798	50793	1.12%	0.345%
16	9.129	1029	1037	1045	rBV4	30992	57982	1.28%	0.393%
17	9.311	1064	1068	1078	rVB4	17813	31340	0.69%	0.213%
18	9.434	1080	1089	1095	rVV4	43252	102824	2.27%	0.697%
19	9.499	1096	1100	1105	rVB3	10757	18205	0.40%	0.123%
20	9.851	1153	1160	1166	rBV4	17994	33076	0.73%	0.224%
21	10.010	1182	1187	1191	rBV2	13228	24355	0.54%	0.165%
22	10.157	1206	1212	1216	rBV3	20940	39139	0.86%	0.265%
23	10.398	1247	1253	1263	rVB6	18036	47498	1.05%	0.322%
24	10.674	1294	1300	1309	rBV6	16893	42702	0.94%	0.290%
25	10.768	1309	1316	1318	rVV	137603	250646	5.53%	1.700%
26	10.821	1318	1325	1336	rVV	2423274	4533004	100.00%	30.747%
27	10.932	1336	1344	1354	rVB	170135	323002	7.13%	2.191%
28	12.319	1574	1580	1587	rBV4	13530	27006	0.60%	0.183%
29	12.430	1593	1599	1605	rVB	102868	164610	3.63%	1.117%
30	12.560	1615	1621	1626	rVV7	7418	17708	0.39%	0.120%
31	12.648	1626	1636	1645	rVB	212854	366082	8.08%	2.483%
32	12.836	1663	1668	1674	rVB7	13395	31951	0.70%	0.217%
33	13.018	1695	1699	1703	rVV4	10269	13981	0.31%	0.095%
34	13.082	1703	1710	1718	rVB7	13878	34650	0.76%	0.235%
35	13.347	1749	1755	1762	rBV3	23293	43773	0.97%	0.297%
36	13.435	1762	1770	1780	rVB2	62136	120185	2.65%	0.815%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

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-BG090721.M
Title : SVOA CALIBRATION

37	13.576	1788	1794	1799	rBV6	10256	21503	0.47%	0.146%
38	13.652	1799	1807	1817	rVB3	32195	73931	1.63%	0.501%
39	13.787	1817	1830	1838	rBV8	31995	92589	2.04%	0.628%
40	13.922	1847	1853	1858	rVV3	28438	51701	1.14%	0.351%
41	13.981	1858	1863	1864	rVV4	17911	26469	0.58%	0.180%
42	14.005	1864	1867	1874	rVB	35111	55828	1.23%	0.379%
43	14.163	1889	1894	1898	rBV3	12198	19220	0.42%	0.130%
44	14.298	1910	1917	1930	rVB3	41064	102860	2.27%	0.698%
45	14.604	1958	1969	1975	rVV	242351	412556	9.10%	2.798%
46	14.668	1975	1980	1987	rVB	237855	374360	8.26%	2.539%
47	14.745	1987	1993	1999	rBV2	66814	107563	2.37%	0.730%
48	14.809	1999	2004	2011	rVB2	18227	33218	0.73%	0.225%
49	14.880	2011	2016	2024	rBV	44585	76654	1.69%	0.520%
50	15.003	2031	2037	2048	rVB2	127301	224644	4.96%	1.524%
51	15.162	2059	2064	2072	rVV2	107951	181242	4.00%	1.229%
52	15.244	2072	2078	2086	rVV7	23942	55909	1.23%	0.379%
53	15.602	2129	2139	2143	rBV	52588	80377	1.77%	0.545%
54	15.655	2143	2148	2157	rVB	120059	188265	4.15%	1.277%
55	15.773	2160	2168	2172	rBV5	46731	105194	2.32%	0.714%
56	15.884	2181	2187	2188	rBV4	19745	30510	0.67%	0.207%
57	15.955	2196	2199	2204	rVB4	10383	15398	0.34%	0.104%
58	16.049	2209	2215	2218	rBV3	10076	19313	0.43%	0.131%
59	16.096	2220	2223	2232	rVB6	15762	31315	0.69%	0.212%
60	16.337	2254	2264	2272	rVB7	30000	72771	1.61%	0.494%
61	16.548	2297	2300	2304	rVV2	13298	19157	0.42%	0.130%
62	16.619	2307	2312	2318	rVV	45032	77953	1.72%	0.529%
63	16.719	2324	2329	2336	rVB6	9801	19223	0.42%	0.130%
64	16.860	2348	2353	2360	rBV7	15524	31020	0.68%	0.210%
65	17.030	2377	2382	2390	rVB3	45824	80391	1.77%	0.545%
66	17.147	2399	2402	2406	rBV	19553	27748	0.61%	0.188%
67	17.259	2415	2421	2422	rBV	12845	16145	0.36%	0.110%
68	17.312	2427	2430	2433	rVV3	20162	30662	0.68%	0.208%
69	17.359	2433	2438	2443	rVV	332062	525708	11.60%	3.566%
70	17.400	2443	2445	2451	rVV2	92967	162746	3.59%	1.104%
71	17.459	2451	2455	2459	rVV	60566	98600	2.18%	0.669%
72	17.506	2459	2463	2468	rVV6	14719	26886	0.59%	0.182%
73	17.770	2497	2508	2519	rBV	395318	636062	14.03%	4.314%
74	17.864	2519	2524	2531	rBV2	41634	83572	1.84%	0.567%
75	17.929	2531	2535	2544	rVB	61322	99412	2.19%	0.674%
76	18.011	2544	2549	2555	rVB3	14194	24720	0.55%	0.168%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

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-BG090721.M
Title : SVOA CALIBRATION

77	18.070	2555	2559	2563	rBV4	9841	16147	0.36%	0.110%
78	18.123	2563	2568	2571	rBV6	9375	15414	0.34%	0.105%
79	18.205	2578	2582	2590	rVB2	31314	54944	1.21%	0.373%
80	18.287	2592	2596	2599	rBV3	35181	51203	1.13%	0.347%
81	18.369	2605	2610	2617	rVB2	36033	57872	1.28%	0.393%
82	18.440	2617	2622	2627	rVB5	10040	16749	0.37%	0.114%
83	18.540	2631	2639	2644	rBV5	19479	47151	1.04%	0.320%
84	18.587	2644	2647	2651	rVV2	17758	24709	0.55%	0.168%
85	18.687	2658	2664	2670	rVV2	16403	35827	0.79%	0.243%
86	18.739	2670	2673	2679	rVV3	10839	14367	0.32%	0.097%
87	19.750	2840	2845	2851	rVV	75856	105981	2.34%	0.719%
88	19.809	2851	2855	2860	rVB4	8436	13825	0.30%	0.094%
89	20.161	2911	2915	2921	rBV	11154	14061	0.31%	0.095%
90	20.226	2921	2926	2932	rBV	28594	42386	0.94%	0.287%
91	20.884	3033	3038	3045	rVV5	12234	23691	0.52%	0.161%
92	21.630	3159	3165	3172	rVB2	375419	602285	13.29%	4.085%
93	22.787	3361	3362	3366	rBV4	10602	17012	0.38%	0.115%
94	23.081	3408	3412	3420	rVB7	11827	25589	0.56%	0.174%
95	23.880	3541	3548	3555	rBV4	22849	50778	1.12%	0.344%
96	24.626	3665	3675	3685	rVB3	48021	128371	2.83%	0.871%
97	24.837	3700	3711	3724	rVB3	366621	1110356	24.49%	7.531%
98	25.936	3890	3898	3906	rVB5	38159	108678	2.40%	0.737%
99	27.269	4115	4125	4135	rVV5	17555	63958	1.41%	0.434%
100	28.890	4392	4401	4402	rBV6	11099	23525	0.52%	0.160%

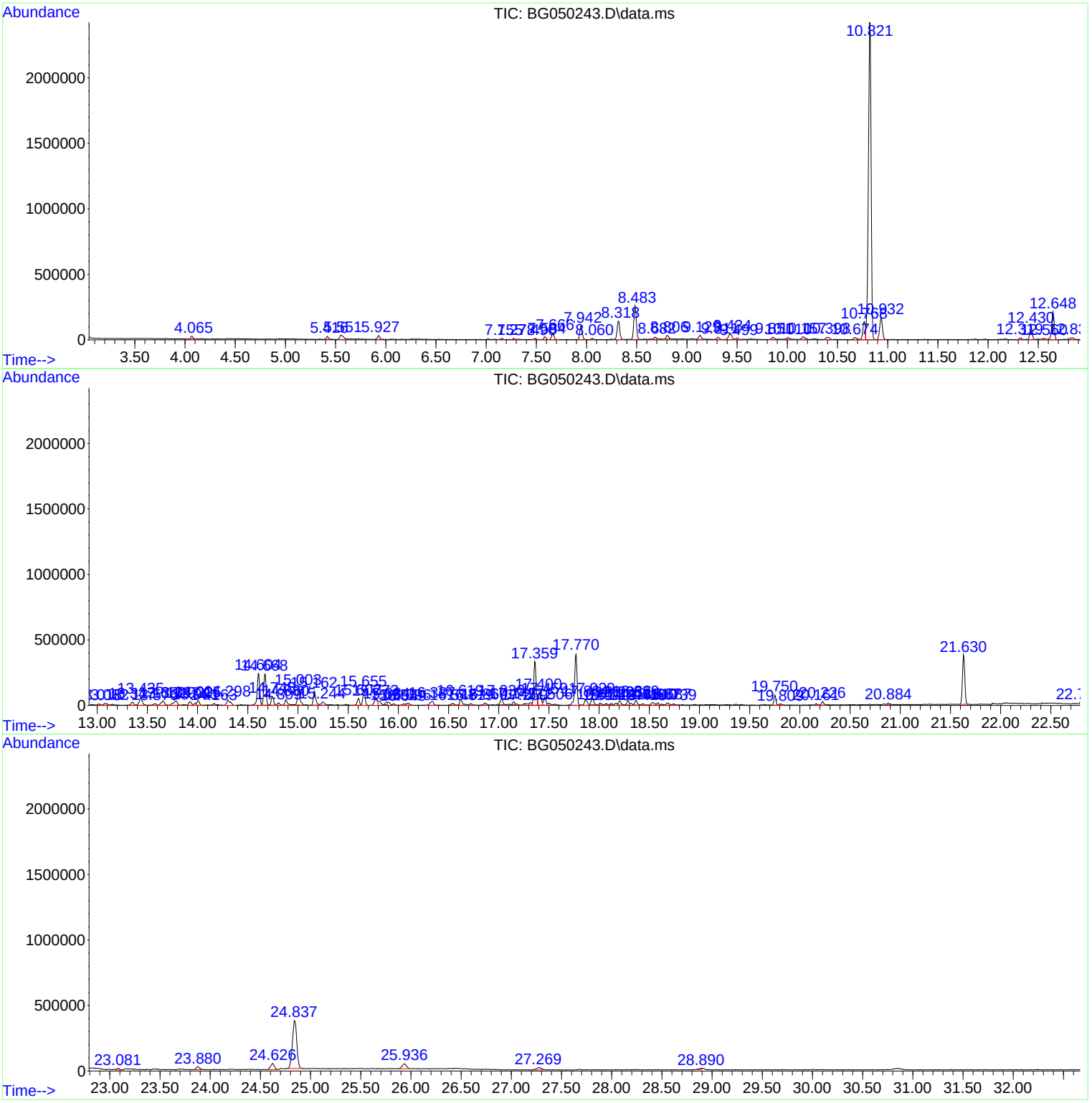
Sum of corrected areas: 14742964

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKK2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKK2DL

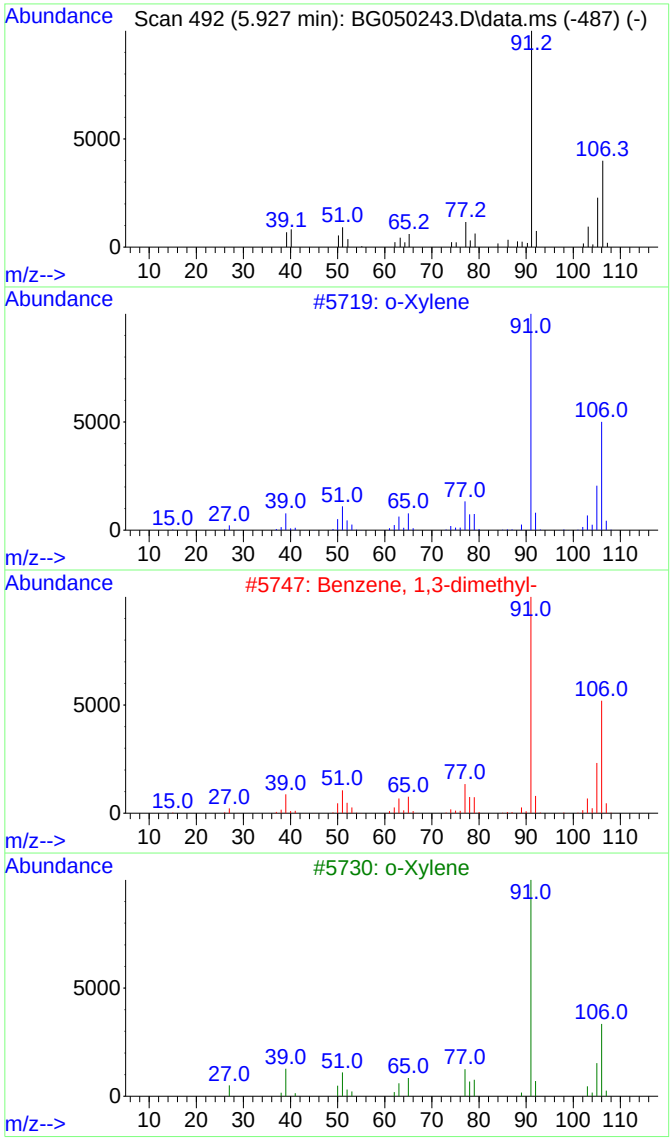
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

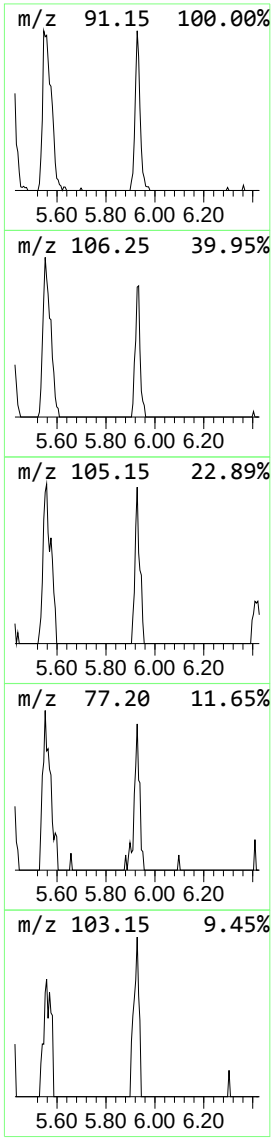
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.927	4.80 ng/ul	42793	1,4-Dichlorobenzene-d4	7.942

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Xylene	106	C8H10	000095-47-6	90
2	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	90
3	o-Xylene	106	C8H10	000095-47-6	87
4	p-Xylene	106	C8H10	000106-42-3	83
5	p-Xylene	106	C8H10	000106-42-3	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

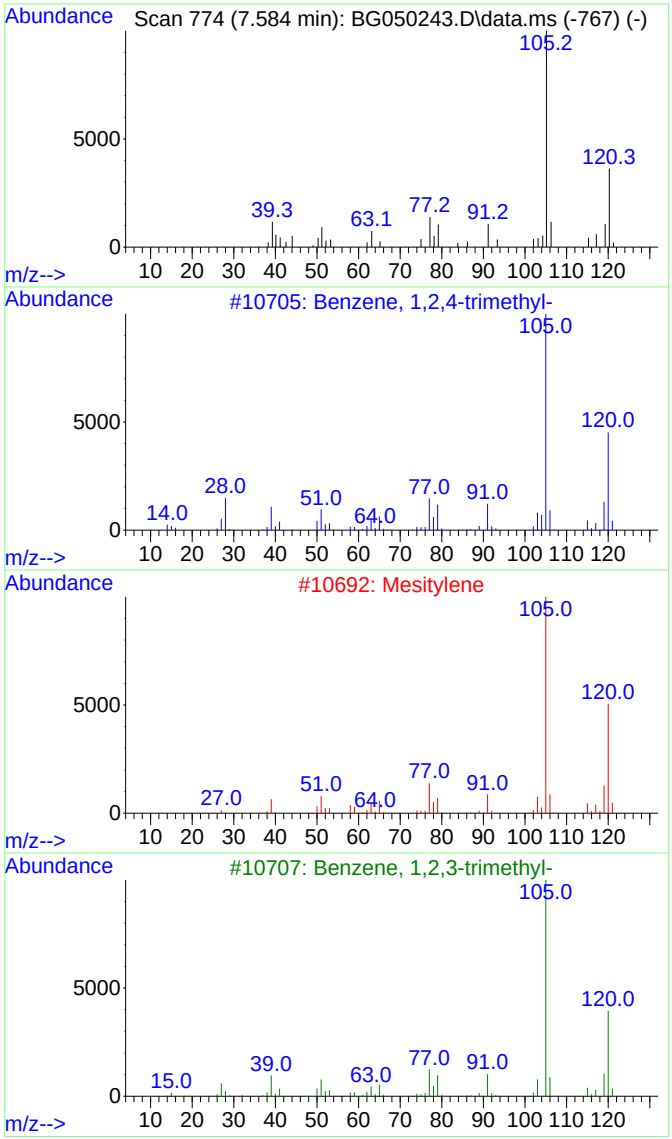
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

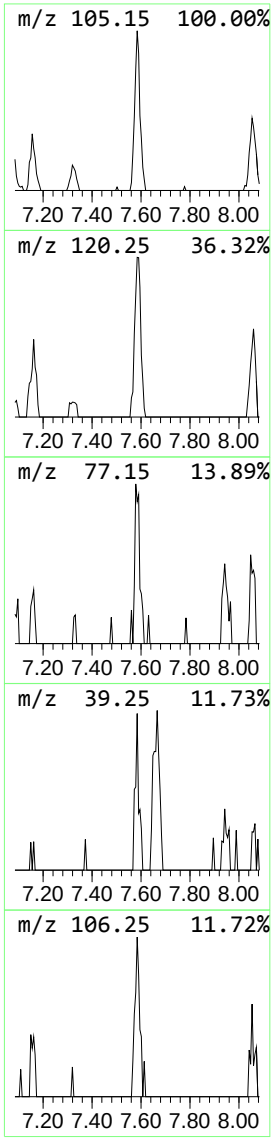
****

Peak Number 5 Benzene, 1,2,4-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.584	4.20 ng/ul	37418	1,4-Dichlorobenzene-d4	7.942

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	87
2	Mesitylene	120	C9H12	000108-67-8	86
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	83
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	72
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKK2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

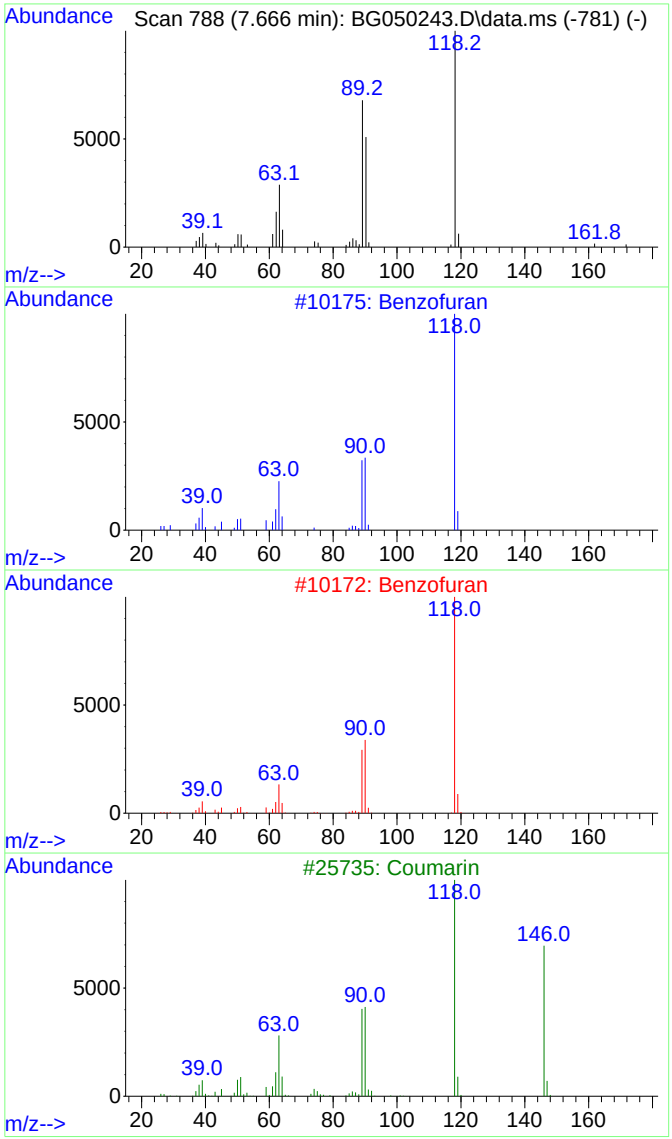
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

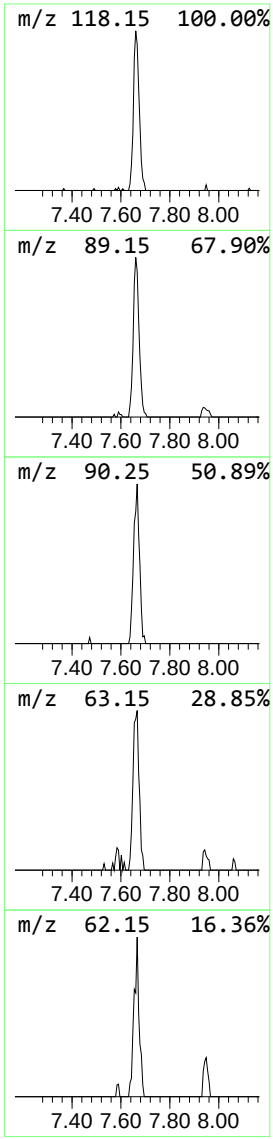
****

Peak Number 6 Benzofuran Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.666	9.13 ng/ul	81339	1,4-Dichlorobenzene-d4	7.942

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran	118	C8H6O	000271-89-6	94
2	Benzofuran	118	C8H6O	000271-89-6	91
3	Coumarin	146	C9H6O2	000091-64-5	83
4	4-Methylphthalic anhydride	162	C9H6O3	019438-61-0	83
5	Benzofuran	118	C8H6O	000271-89-6	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

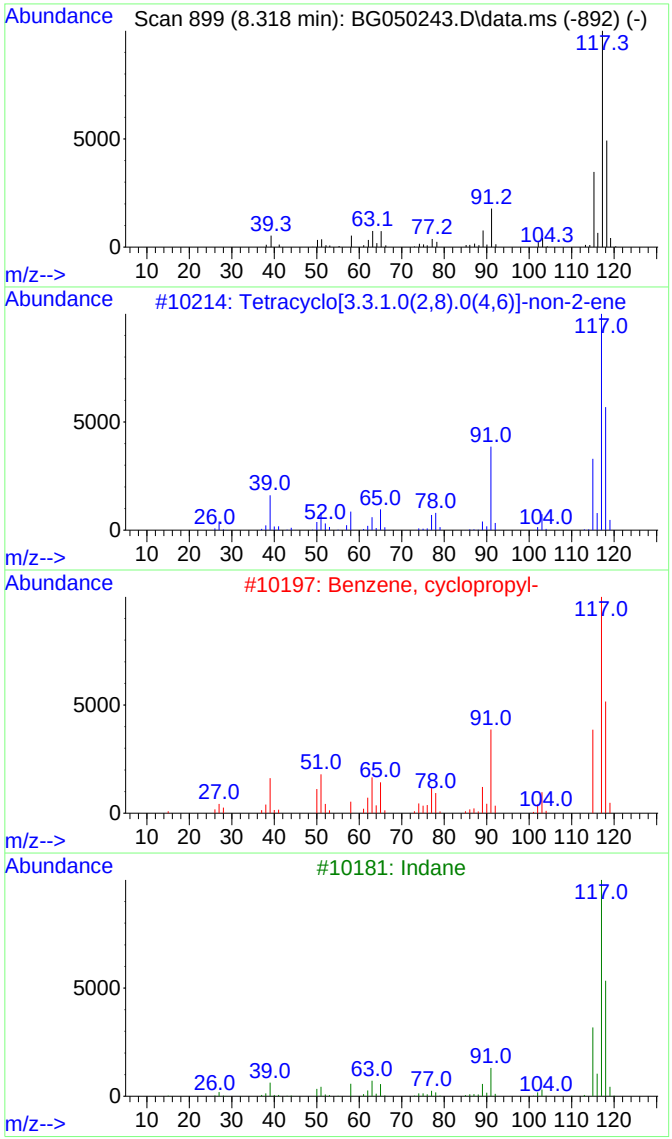
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

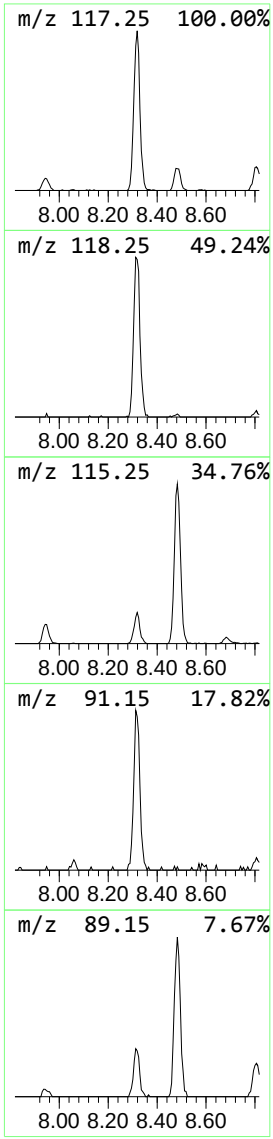
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Tetracyclo[3.3.1.0(2,8).0(4... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.318	28.93 ng/ul	257792	1,4-Dichlorobenzene-d4	7.942

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
3	Indane	118	C9H10	000496-11-7	60
4	Indane	118	C9H10	000496-11-7	60
5	Indane	118	C9H10	000496-11-7	60





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKK2DL

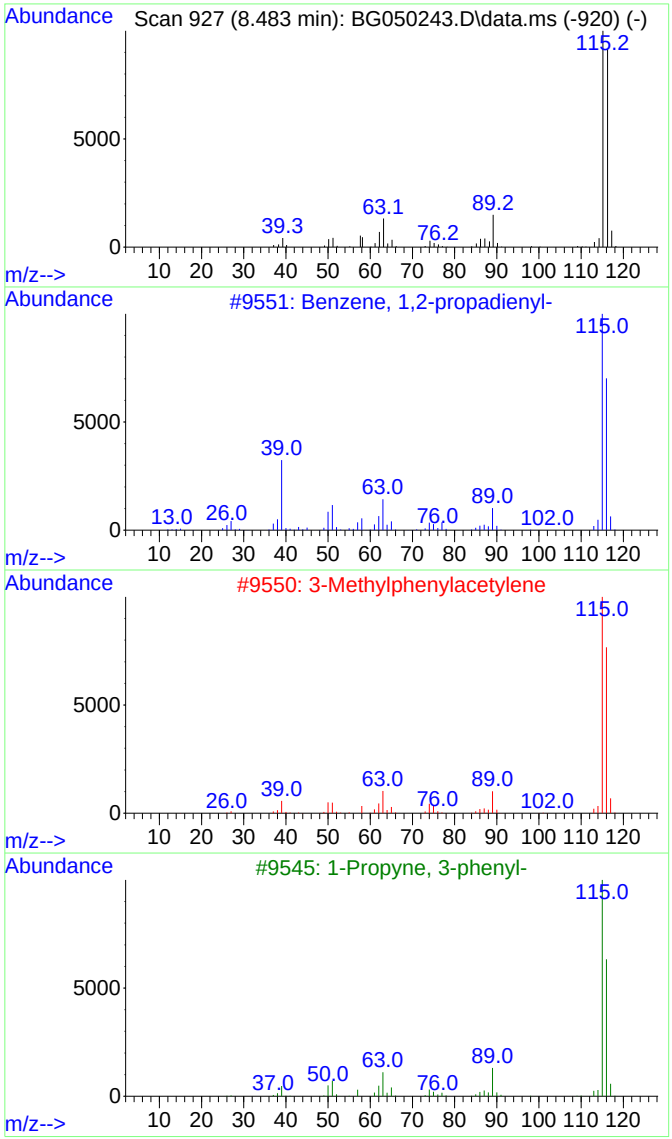
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

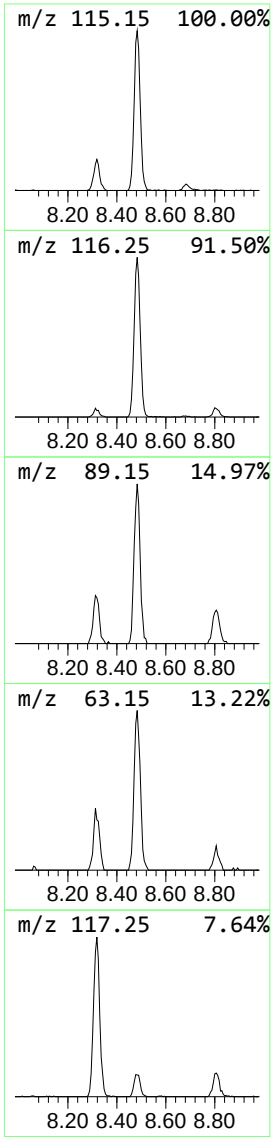
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1,2-propadienyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.483	50.43 ng/ul	449331	1,4-Dichlorobenzene-d4	7.942

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	94
2	3-Methylphenylacetylene	116	C9H8	000766-82-5	94
3	1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	94
4	Indene	116	C9H8	000095-13-6	93
5	Indene	116	C9H8	000095-13-6	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

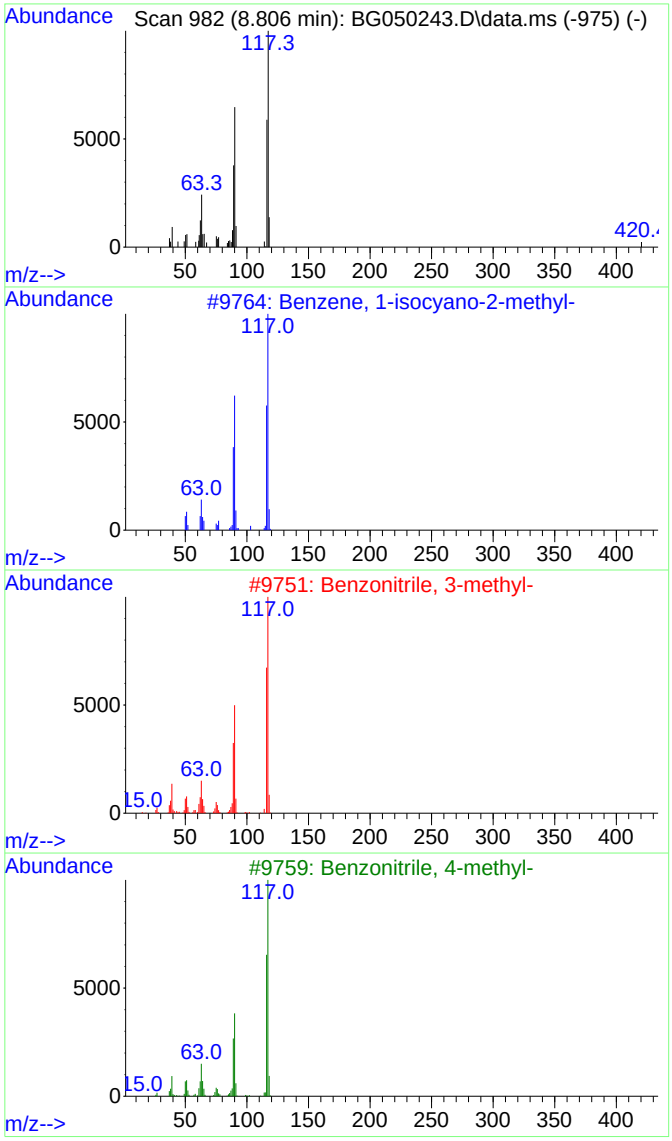
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

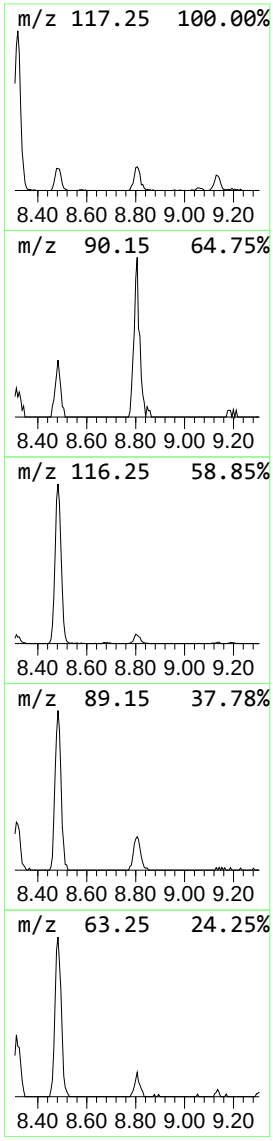
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1-isocyano-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.806	5.70 ng/ul	50793	1,4-Dichlorobenzene-d4	7.942

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-isocyano-2-methyl-	117	C8H7N	010468-64-1	95
2	Benzonitrile, 3-methyl-	117	C8H7N	000620-22-4	95
3	Benzonitrile, 4-methyl-	117	C8H7N	000104-85-8	94
4	Benzonitrile, 4-methyl-	117	C8H7N	000104-85-8	94
5	Benzonitrile, 4-methyl-	117	C8H7N	000104-85-8	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

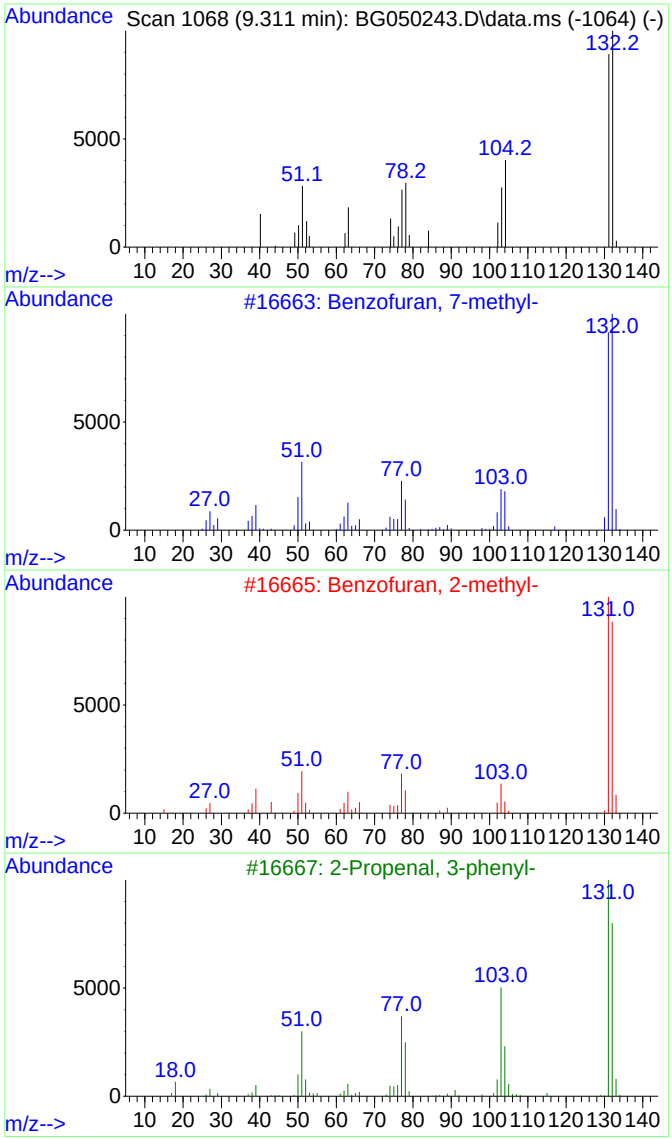
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

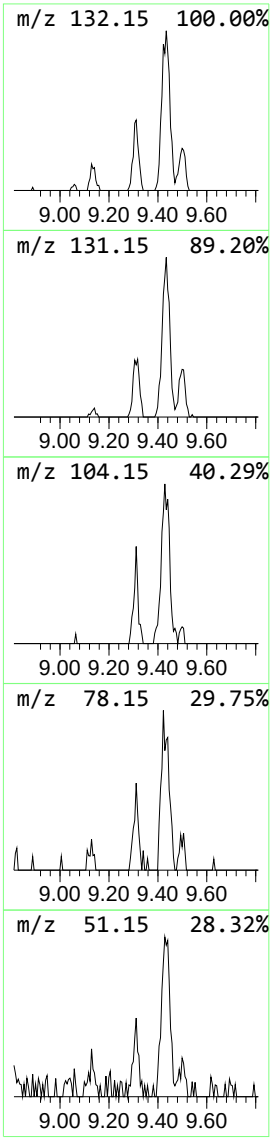
****

Peak Number 10 Benzofuran, 7-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.311	3.52 ng/ul	31340	1,4-Dichlorobenzene-d4	7.942

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	87
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	81
3	2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	68
4	Benzene, (2-propynyloxy)-	132	C9H8O	013610-02-1	64
5	Benzonitrile, 2-(methylamino)-	132	C8H8N2	017583-40-3	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

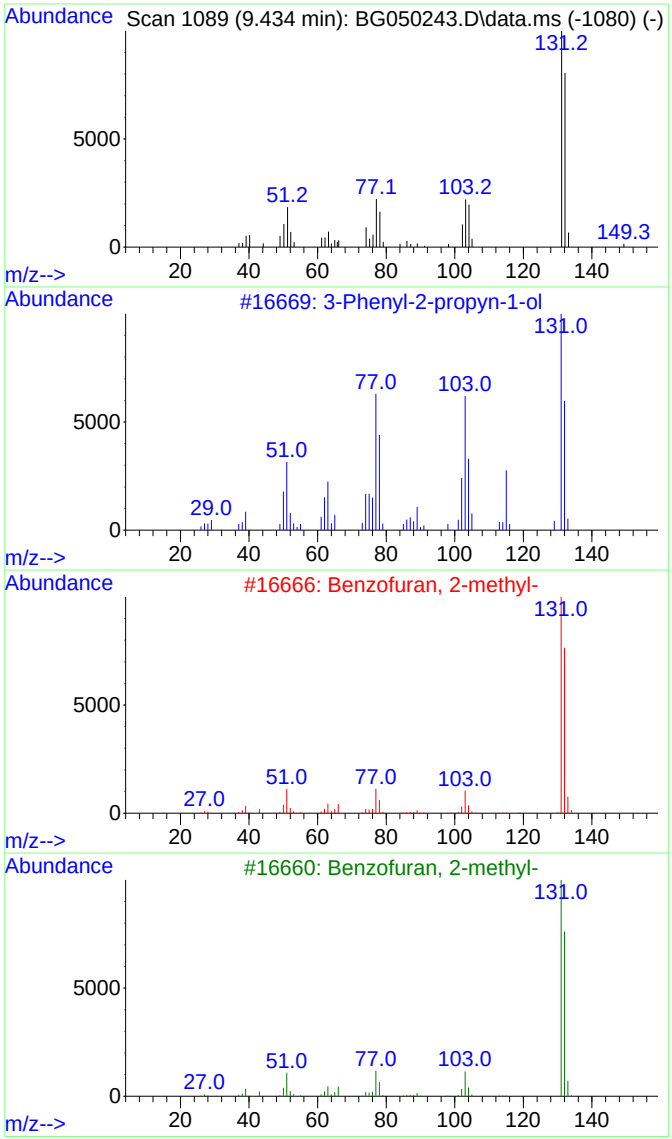
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

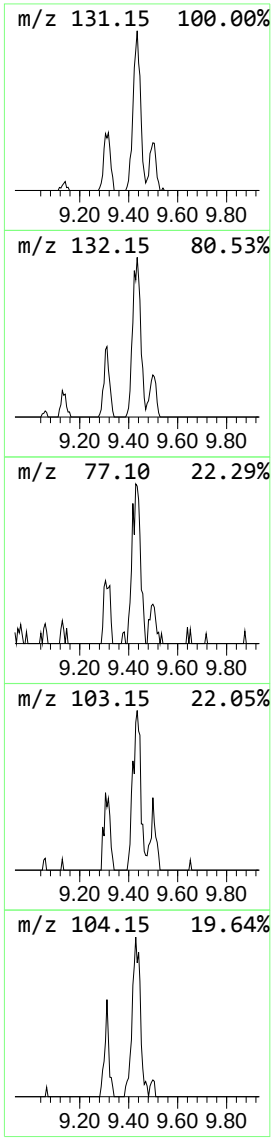
****

Peak Number 11 3-Phenyl-2-propyn-1-ol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.434	8.20 ng/ul	102824	Naphthalene-d8	10.768

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	91
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
3	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
4	Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	90
5	2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

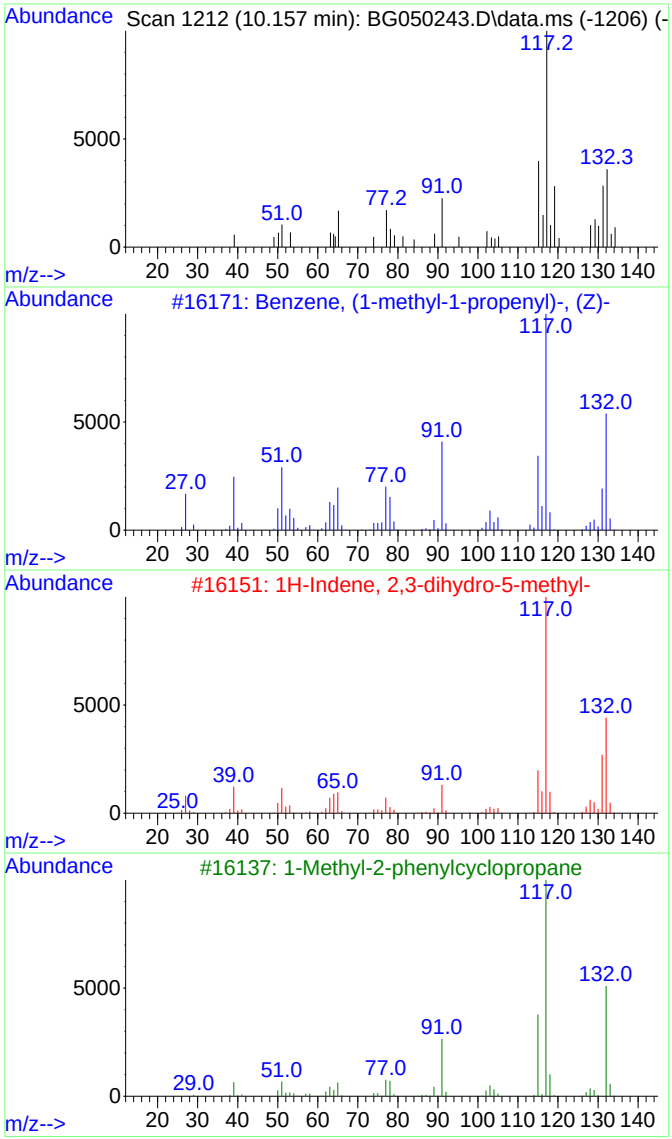
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

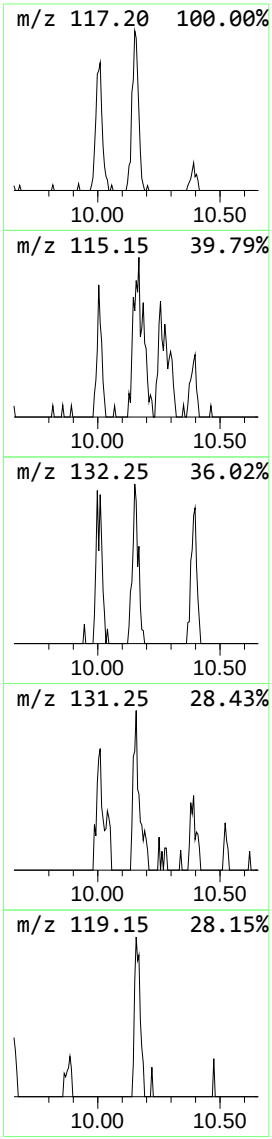
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Benzene, (1-methyl-1-propen... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.157	3.12 ng/ul	39139	Naphthalene-d8	10.768

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	70
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	64
3	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	64
4	(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	64
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

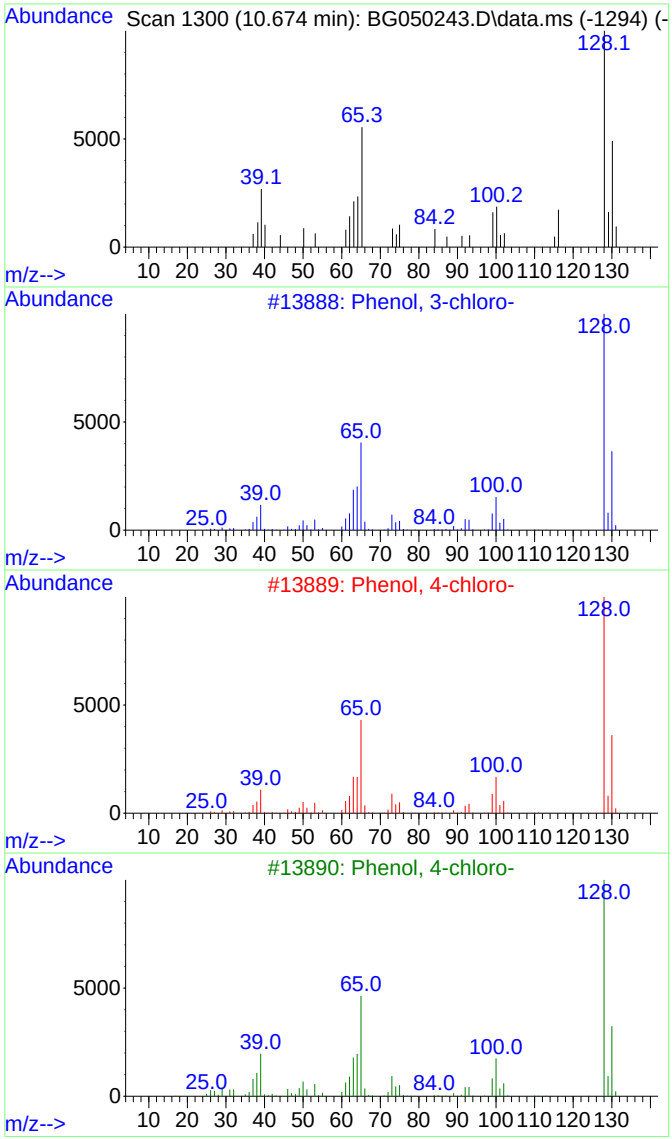
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

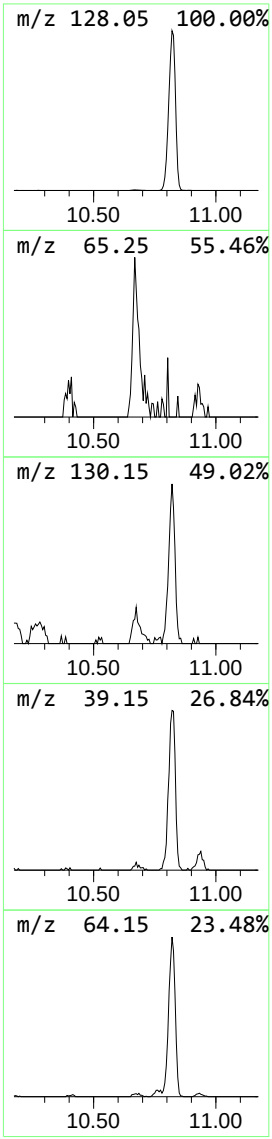
****

Peak Number 13 Phenol, 3-chloro- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.674	3.41 ng/ul	42702	Naphthalene-d8	10.768

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3-chloro-	128	C6H5ClO	000108-43-0	97
2	Phenol, 4-chloro-	128	C6H5ClO	000106-48-9	96
3	Phenol, 4-chloro-	128	C6H5ClO	000106-48-9	94
4	Phenol, 3-chloro-	128	C6H5ClO	000108-43-0	93
5	Phenol, 4-chloro-	128	C6H5ClO	000106-48-9	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

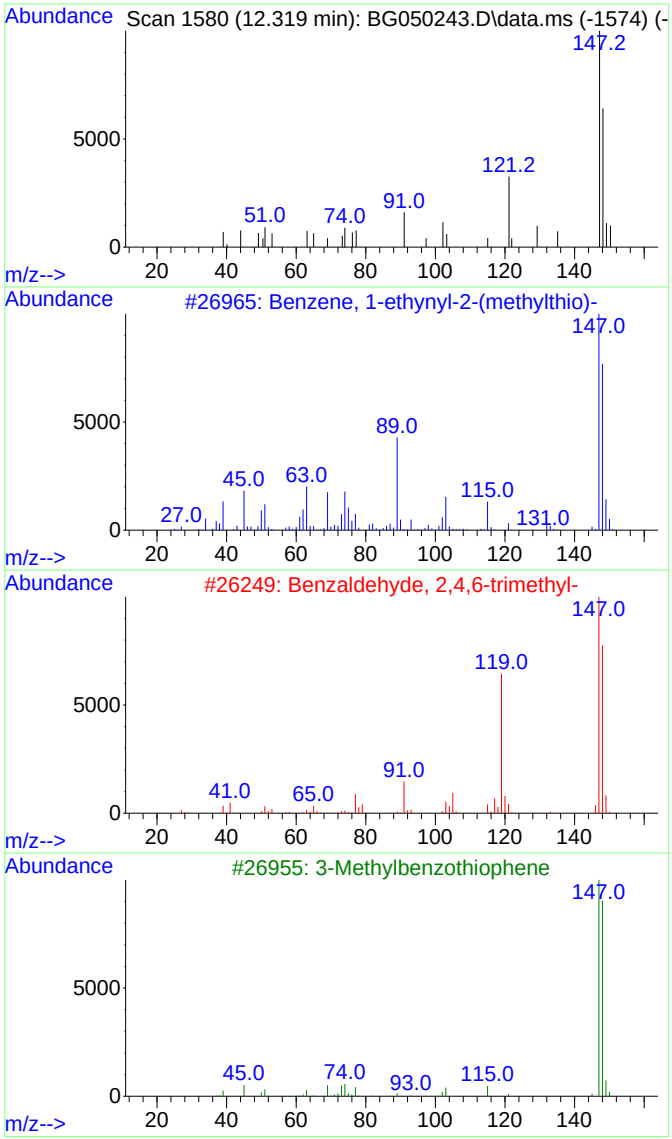
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

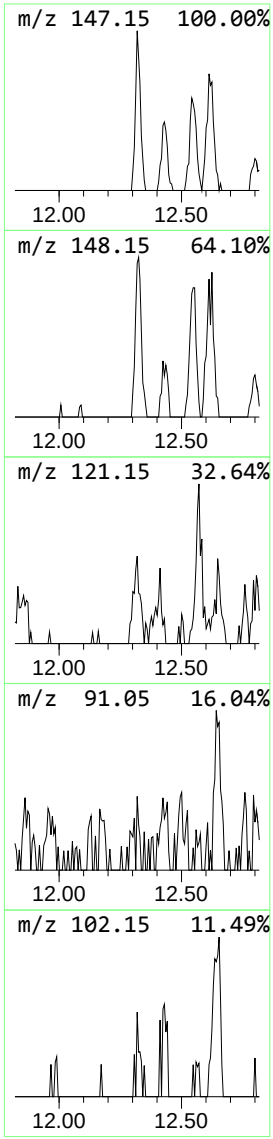
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Benzene, 1-ethynyl-2-(methy... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.319	2.15 ng/ul	27006	Naphthalene-d8	10.768

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethynyl-2-(methylthio)-	148	C9H8S	078905-08-5	64
2	Benzaldehyde, 2,4,6-trimethyl-	148	C10H12O	000487-68-3	64
3	3-Methylbenzothiophene	148	C9H8S	001455-18-1	64
4	3-Methylbenzothiophene	148	C9H8S	001455-18-1	64
5	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

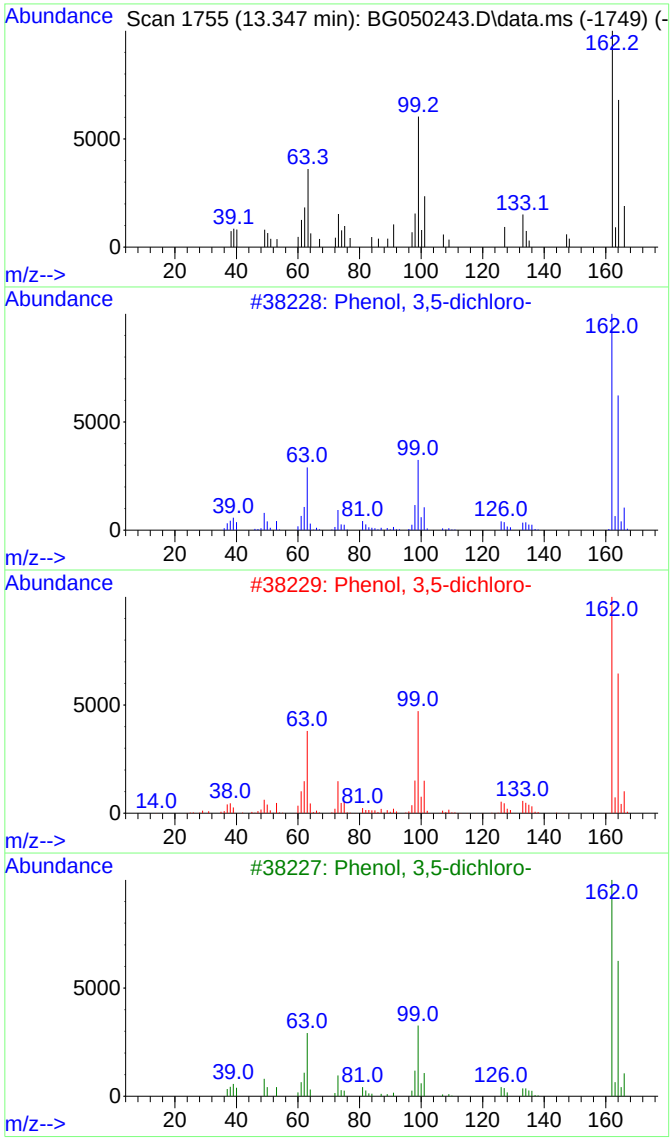
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

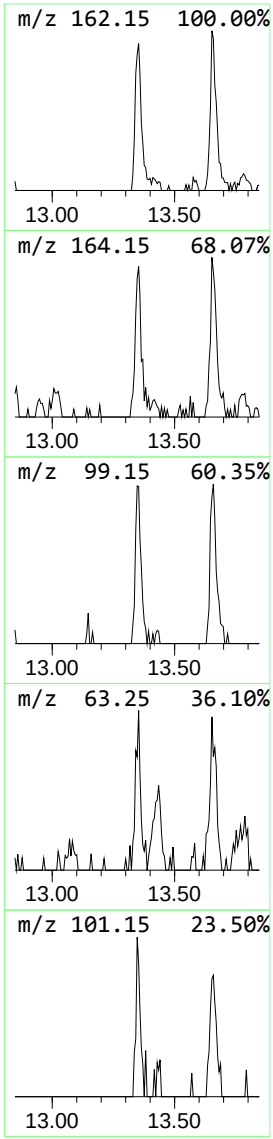
****

Peak Number 15 Phenol, 3,5-dichloro- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.347	2.12 ng/ul	43773	Acenaphthene-d10	14.610

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	91
2	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	91
3	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	91
4	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	91
5	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKK2DL

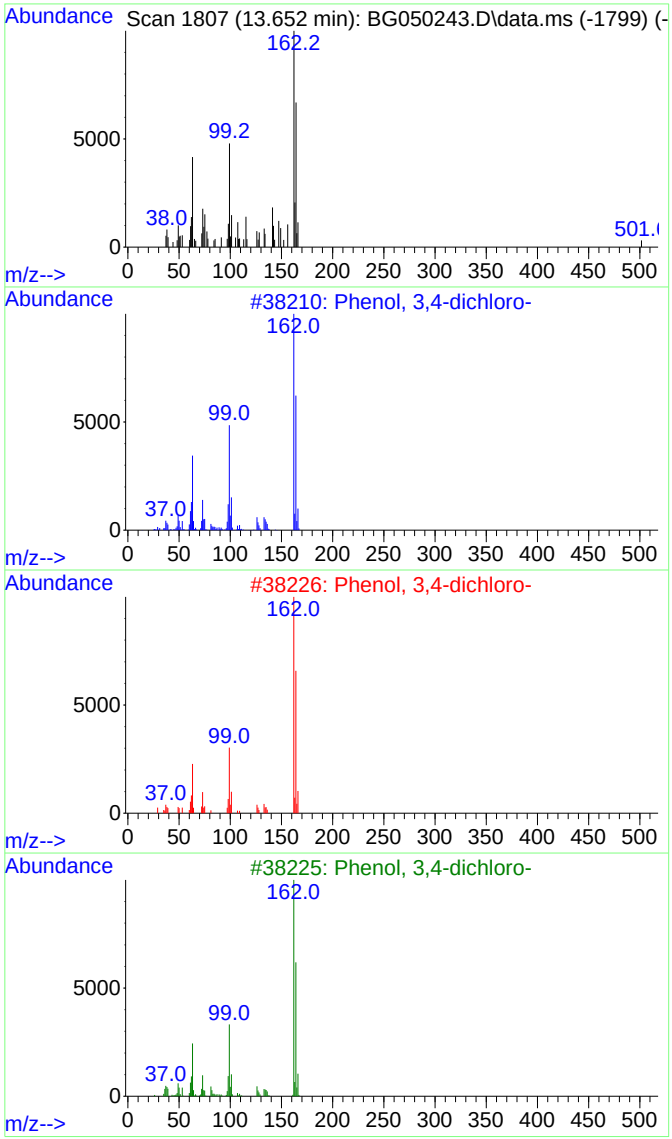
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

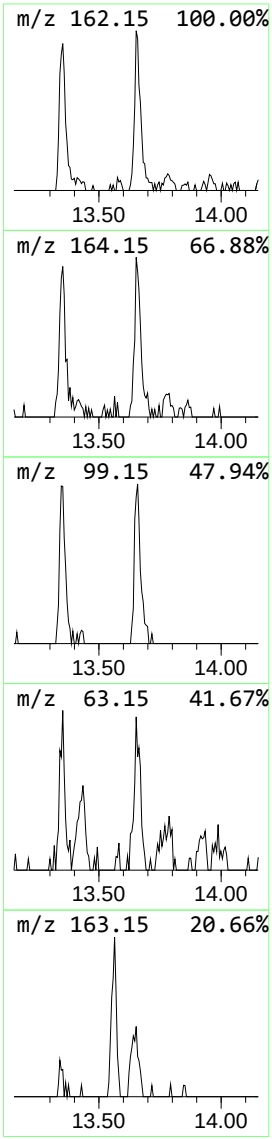
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Phenol, 3,4-dichloro- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.652	3.58 ng/ul	73931	Acenaphthene-d10	14.610

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97
2	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	96
3	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	96
4	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	95
5	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

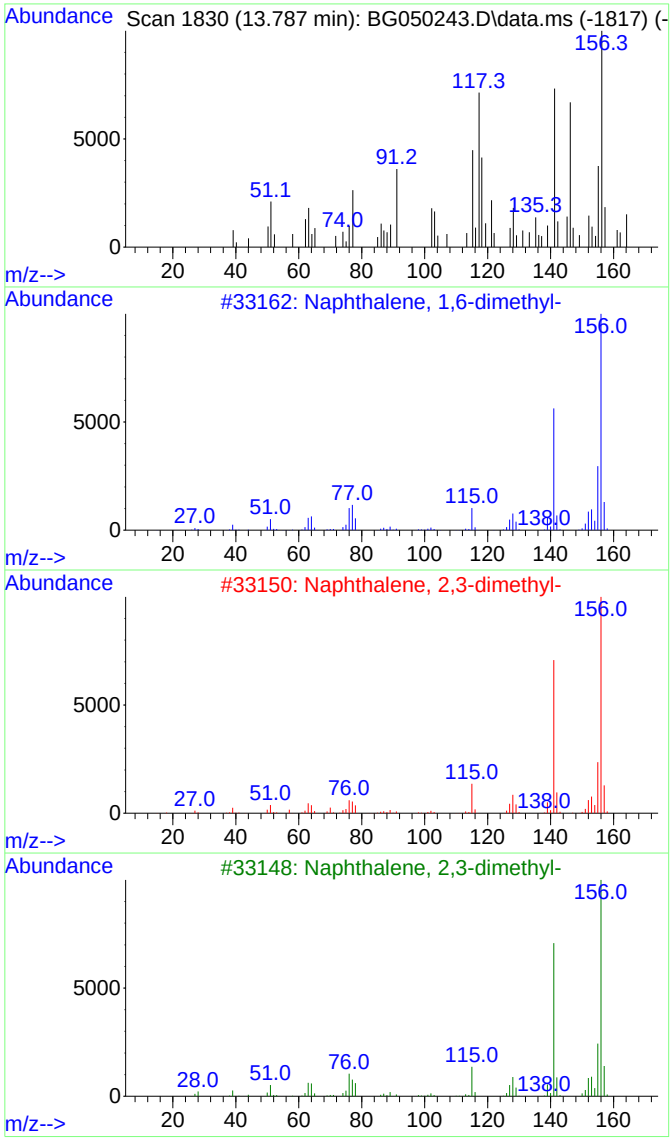
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

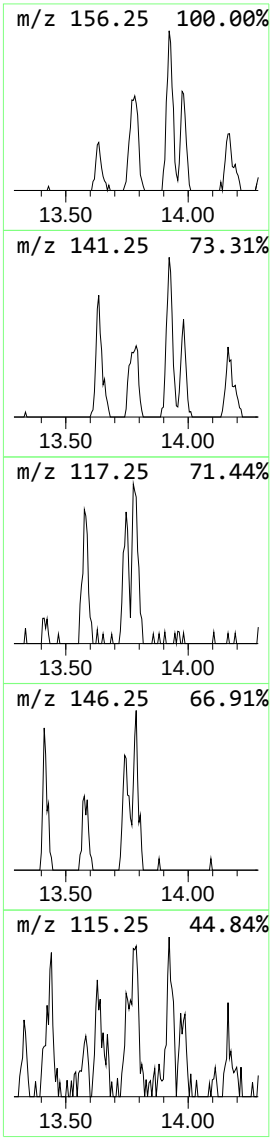
****

Peak Number 17 Naphthalene, 1,6-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.787	4.49 ng/ul	92589	Acenaphthene-d10	14.610

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	90
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	89
3	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	89
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	89
5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

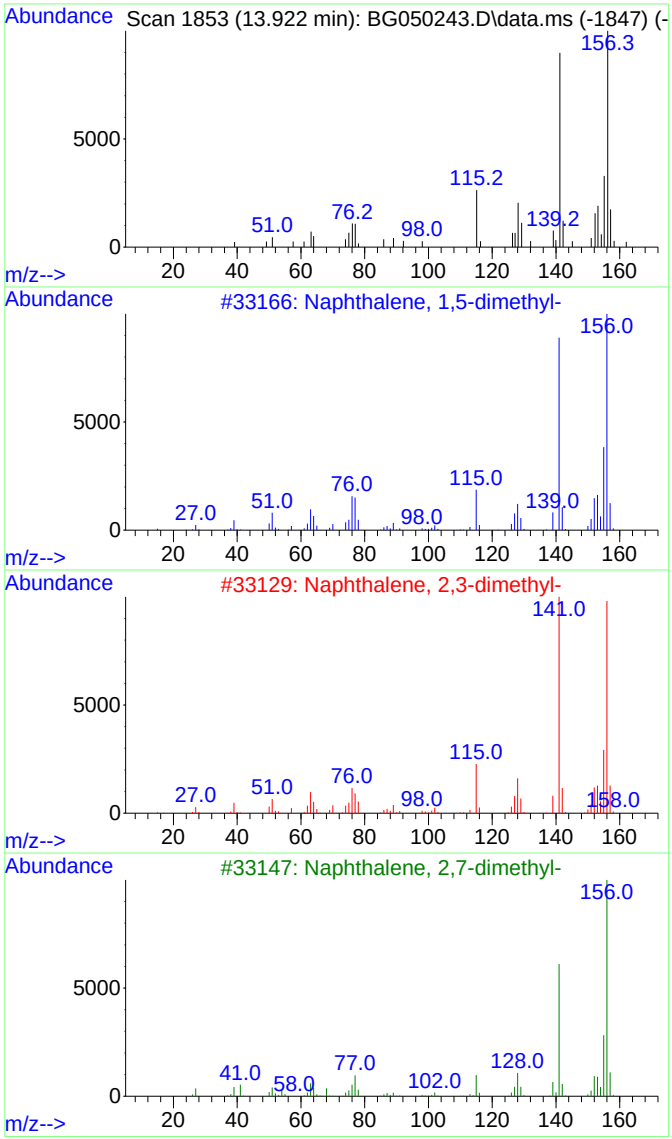
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

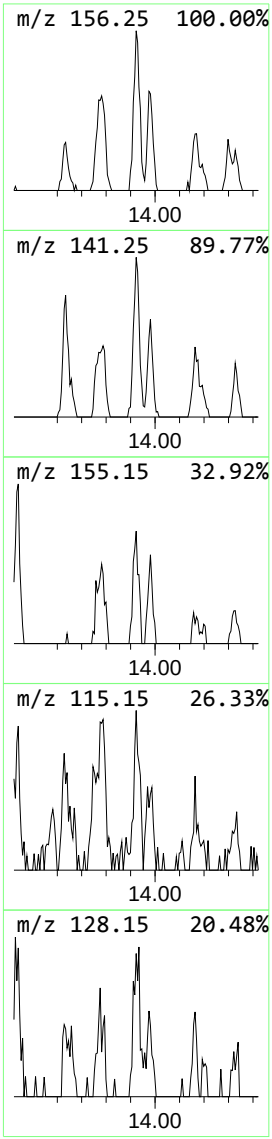
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Naphthalene, 1,5-dimethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.922	2.51 ng/ul	51701	Acenaphthene-d10	14.610

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	95
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
4	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	94
5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

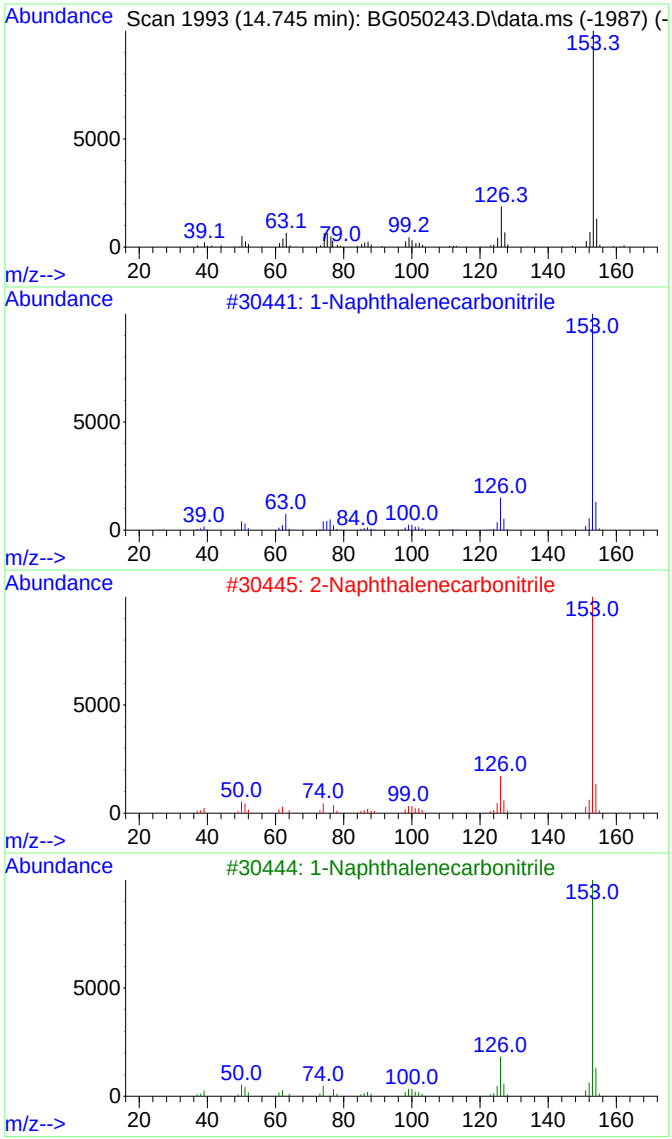
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

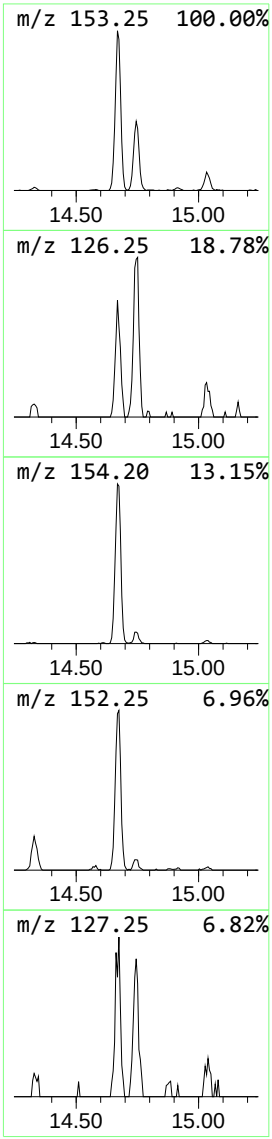
****

Peak Number 19 1-Naphthalenecarbonitrile Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.745	5.21 ng/ul	107563	Acenaphthene-d10	14.610

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	97
2	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	94
3	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	94
4	2-Naphthyl isocyanide	153	C11H7N	010124-78-4	94
5	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

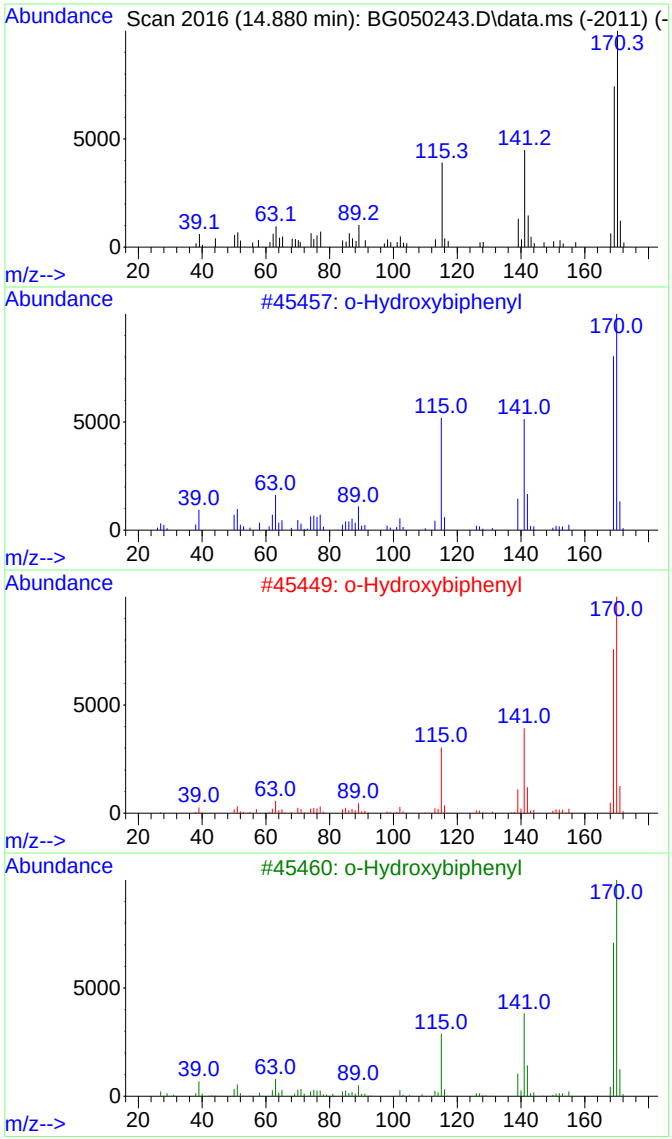
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

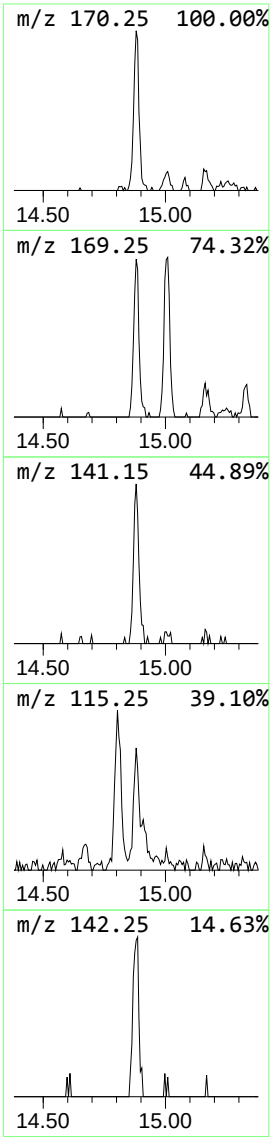
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 o-Hydroxybiphenyl Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.880	3.72 ng/ul	76654	Acenaphthene-d10	14.610

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	96
2	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	95
3	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	95
4	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	95
5	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

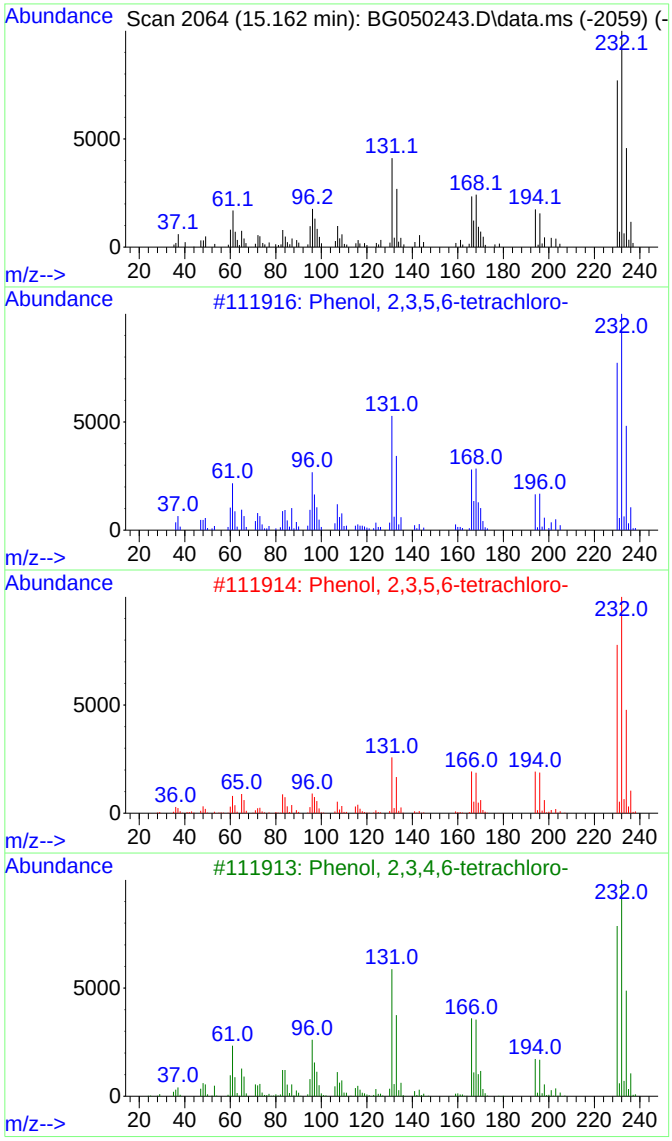
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

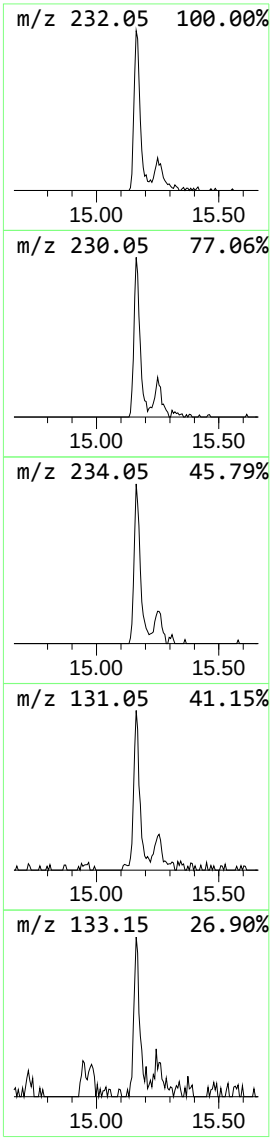
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Phenol, 2,3,5,6-tetrachloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.162	8.79 ng/ul	181242	Acenaphthene-d10	14.610

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





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

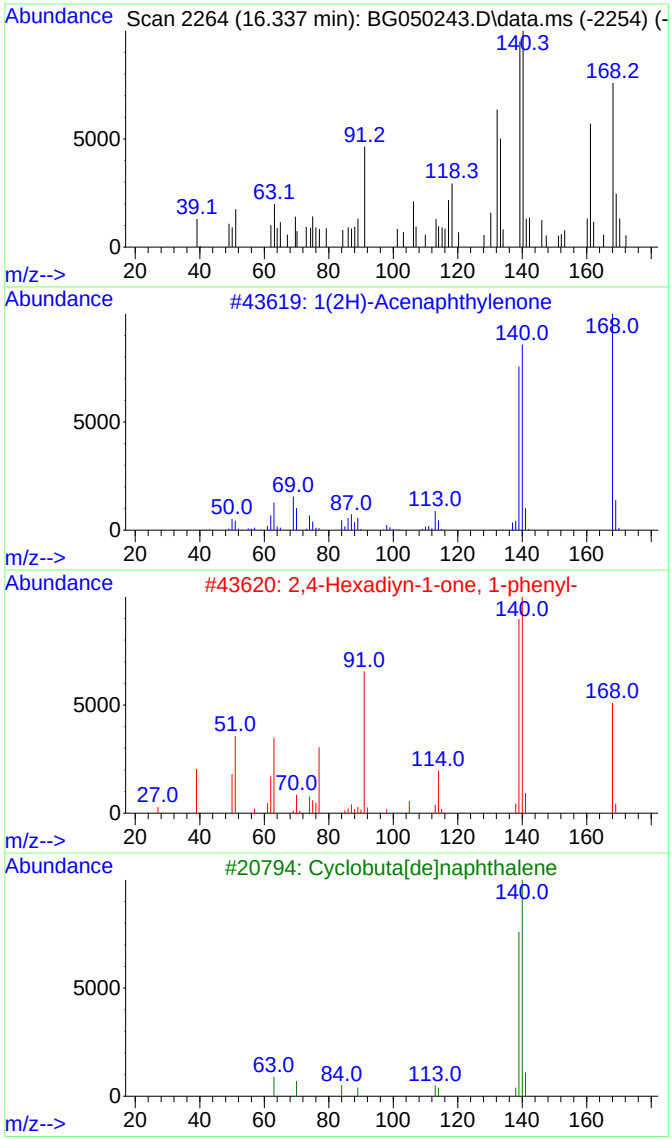
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

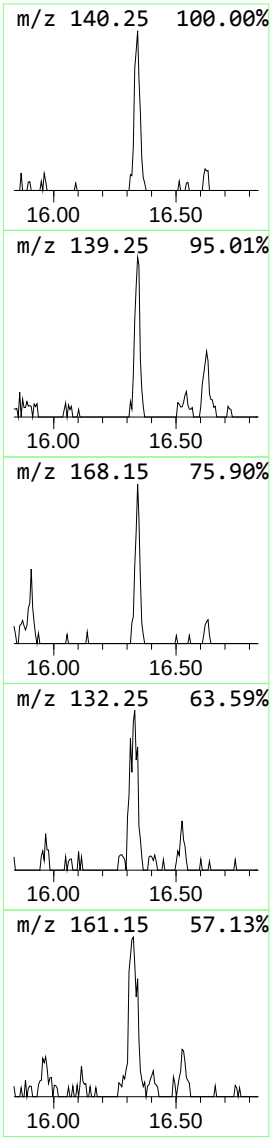
****

Peak Number 22 1(2H)-Acenaphthylenone Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.337	2.77 ng/ul	72771	Phenanthrene-d10	17.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(2H)-Acenaphthylenone	168	C12H8O	002235-15-6	55
2	2,4-Hexadiyn-1-one, 1-phenyl-	168	C12H8O	000495-74-9	53
3	Cyclobuta[de]naphthalene	140	C11H8	024973-91-9	38
4	Benzaldehyde, 2-fluoro-4-hydroxy-	140	C7H5FO2	000348-27-6	38
5	5,6,7,8-Tetrahydro-thiazolo[5,4-...	168	C7H8N2OS	1000317-75-4	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

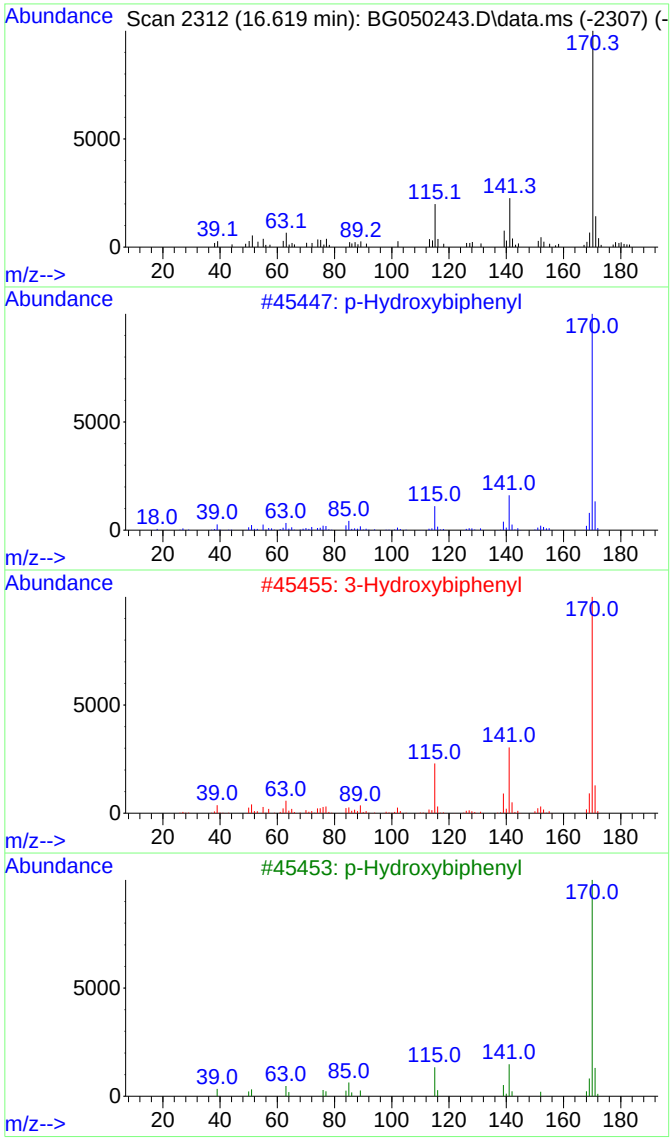
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

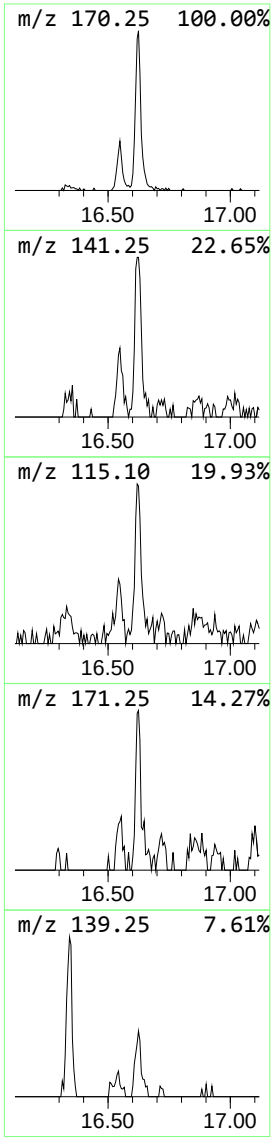
****

Peak Number 23 p-Hydroxybiphenyl Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.619	2.97 ng/ul	77953	Phenanthrene-d10	17.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Hydroxybiphenyl	170	C12H10O	000092-69-3	94
2	3-Hydroxybiphenyl	170	C12H10O	000580-51-8	93
3	p-Hydroxybiphenyl	170	C12H10O	000092-69-3	91
4	p-Hydroxybiphenyl	170	C12H10O	000092-69-3	90
5	p-Hydroxybiphenyl	170	C12H10O	000092-69-3	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

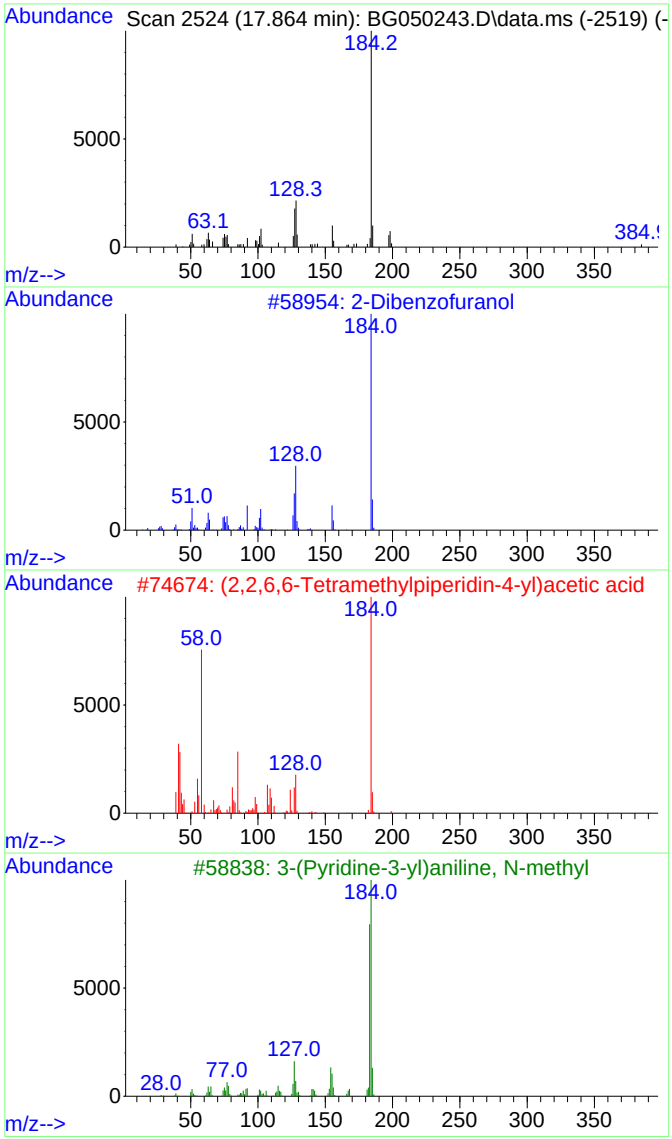
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

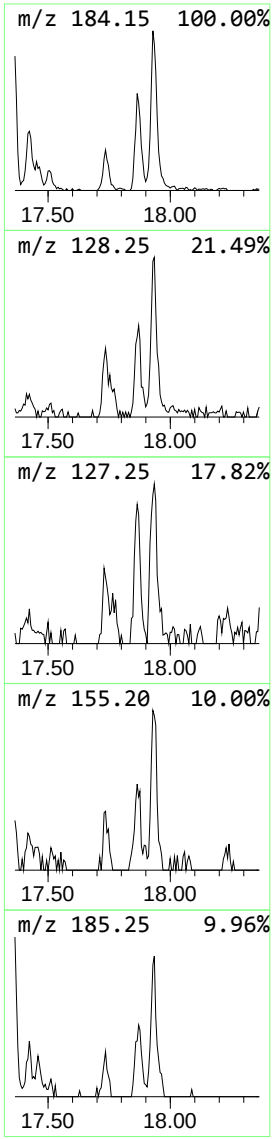
****

Peak Number 24 2-Dibenzofuranol Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.864	3.18 ng/ul	83572	Phenanthrene-d10	17.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Dibenzofuranol	184	C12H8O2	000086-77-1	74
2	(2,2,6,6-Tetramethylpiperidin-4-...	199	C11H21NO2	1000310-71-9	64
3	3-(Pyridine-3-yl)aniline, N-methyl	184	C12H12N2	1378740-24-9	64
4	2-Dibenzofuranol	184	C12H8O2	000086-77-1	64
5	1H-Pyrazolo[3,4-b]quinolin-3-amine	184	C10H8N4	1000319-54-1	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

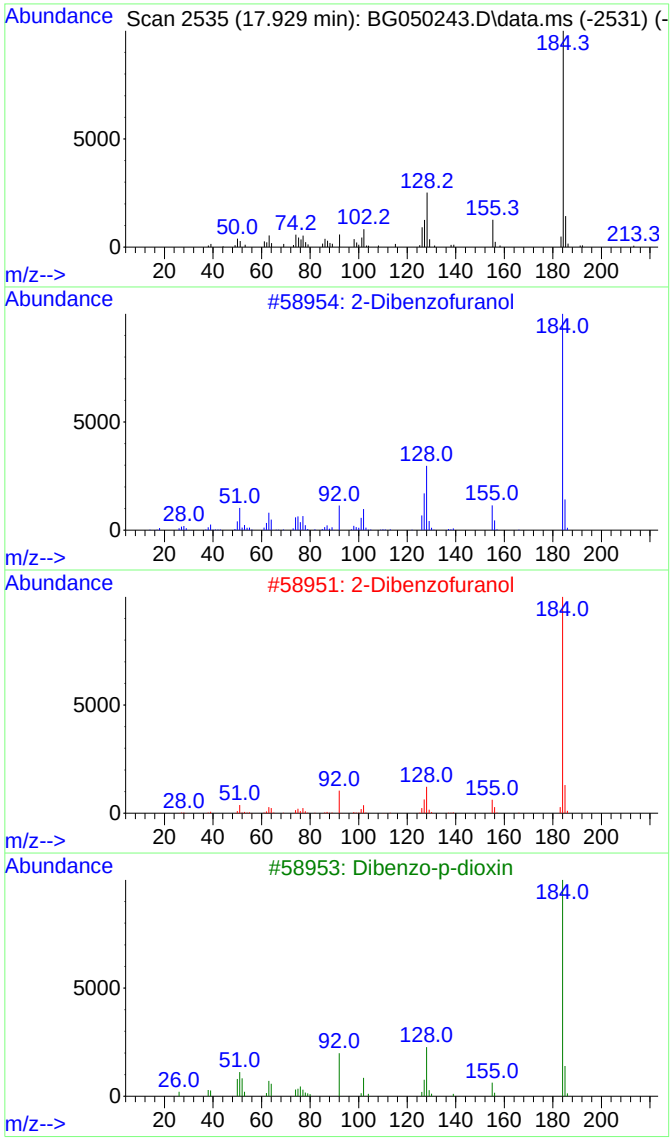
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

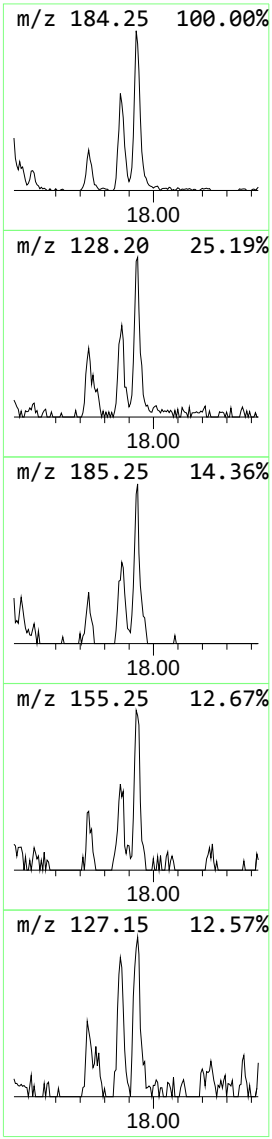
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 Dibenzo-p-dioxin Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.929	3.78 ng/ul	99412	Phenanthrene-d10	17.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Dibenzofuranol	184	C12H8O2	000086-77-1	91
2	2-Dibenzofuranol	184	C12H8O2	000086-77-1	91
3	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	90
4	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	87
5	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

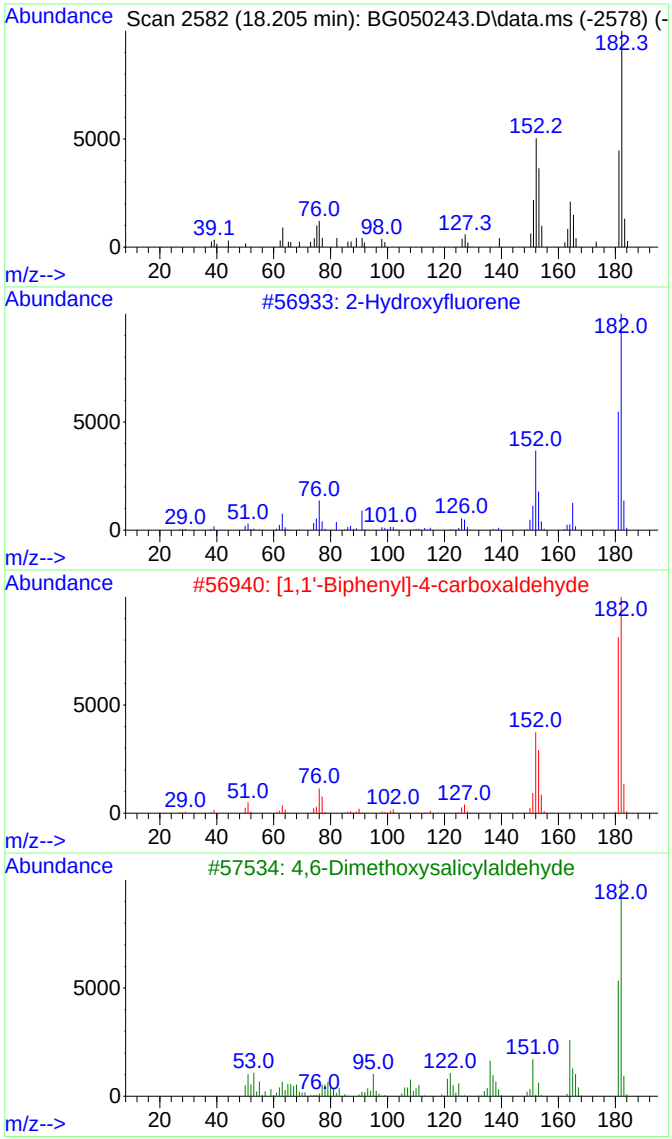
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

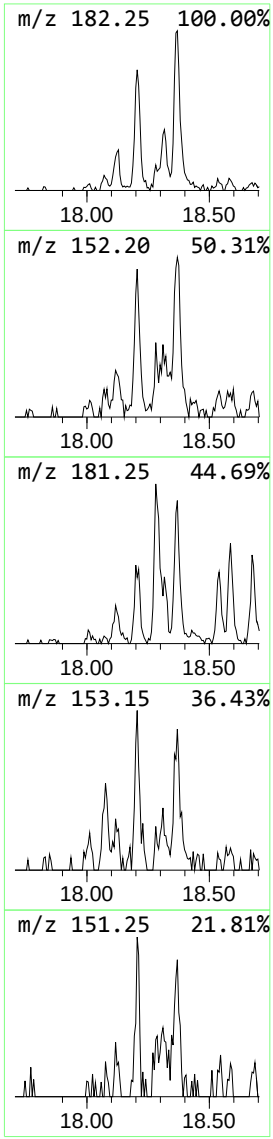
****

Peak Number 26 2-Hydroxyfluorene Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.205	2.09 ng/ul	54944	Phenanthrene-d10	17.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hydroxyfluorene	182	C13H10O	002443-58-5	87
2	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	52
3	4,6-Dimethoxysalicylaldehyde	182	C9H10O4	000708-76-9	50
4	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	50
5	9H-Fluoren-3-ol	182	C13H10O	006344-67-8	49





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

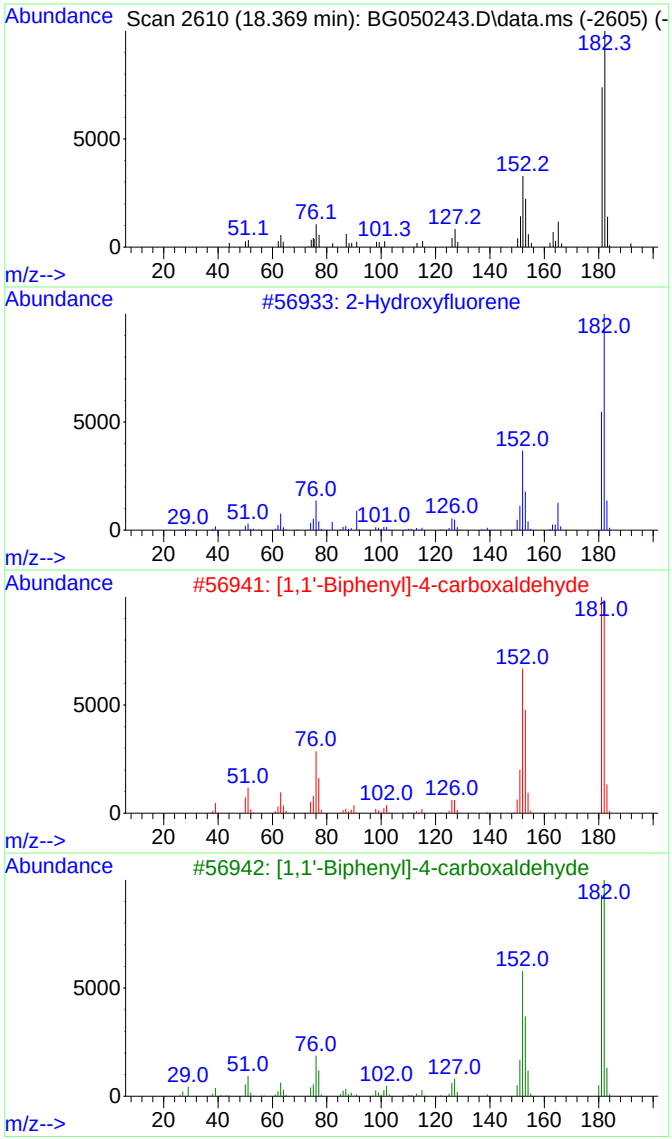
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

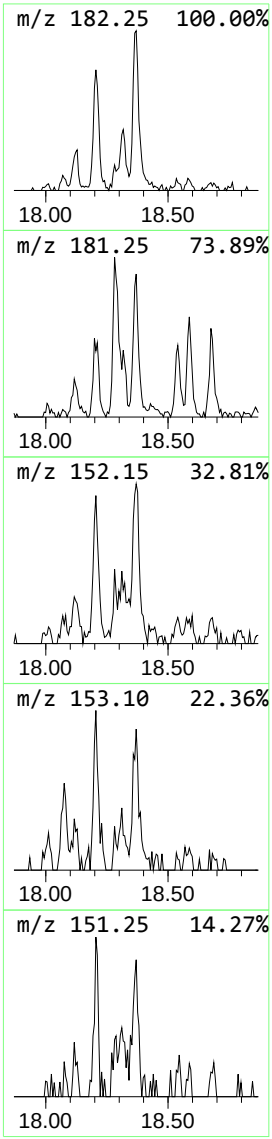
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 [1,1'-Biphenyl]-4-carboxald... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.369	2.20 ng/ul	57872	Phenanthrene-d10	17.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hydroxyfluorene	182	C13H10O	002443-58-5	93
2	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	90
3	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	90
4	2-Hydroxyfluorene	182	C13H10O	002443-58-5	90
5	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050243.D
Acq On : 10 Sep 2021 00:13
Operator : CG/JU
Sample : M3608-06DL 10X
Misc :
ALS Vial : 71 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKK2DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.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, 1,3-di...	5.927	4.8	ng/ul	42793	1	7.942	178204	20.0
Benzene, 1,2,4-...	7.584	4.2	ng/ul	37418	1	7.942	178204	20.0
Benzofuran	7.666	9.1	ng/ul	81339	1	7.942	178204	20.0
Tetracyclo[3.3....	8.318	28.9	ng/ul	257792	1	7.942	178204	20.0
Benzene, 1,2-pr...	8.483	50.4	ng/ul	449331	1	7.942	178204	20.0
Benzene, 1-isoc...	8.806	5.7	ng/ul	50793	1	7.942	178204	20.0
Benzofuran, 7-m...	9.311	3.5	ng/ul	31340	1	7.942	178204	20.0
3-Phenyl-2-prop...	9.434	8.2	ng/ul	102824	2	10.768	250646	20.0
Benzene, (1-met...	10.157	3.1	ng/ul	39139	2	10.768	250646	20.0
Phenol, 3-chloro-	10.674	3.4	ng/ul	42702	2	10.768	250646	20.0
Benzene, 1-ethy...	12.319	2.1	ng/ul	27006	2	10.768	250646	20.0
Phenol, 3,5-dic...	13.347	2.1	ng/ul	43773	3	14.610	412556	20.0
Phenol, 3,4-dic...	13.652	3.6	ng/ul	73931	3	14.610	412556	20.0
Naphthalene, 1,...	13.787	4.5	ng/ul	92589	3	14.610	412556	20.0
Naphthalene, 1,...	13.922	2.5	ng/ul	51701	3	14.610	412556	20.0
1-Naphthaleneca...	14.745	5.2	ng/ul	107563	3	14.610	412556	20.0
o-Hydroxybiphenyl	14.880	3.7	ng/ul	76654	3	14.610	412556	20.0
Phenol, 2,3,5,6...	15.162	8.8	ng/ul	181242	3	14.610	412556	20.0
1(2H)-Acenaphth...	16.337	2.8	ng/ul	72771	4	17.359	525708	20.0
p-Hydroxybiphenyl	16.619	3.0	ng/ul	77953	4	17.359	525708	20.0
2-Dibenzofuranol	17.864	3.2	ng/ul	83572	4	17.359	525708	20.0
Dibenzo-p-dioxin	17.929	3.8	ng/ul	99412	4	17.359	525708	20.0
2-Hydroxyfluorene	18.205	2.1	ng/ul	54944	4	17.359	525708	20.0
[1,1'-Biphenyl]...	18.369	2.2	ng/ul	57872	4	17.359	525708	20.0