LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
Data File : BG050065.D
Acq On : 31 Aug 2021 15:02
Operator : CG/JU
Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKG0

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-BG081921.M
Title : SVOA CALIBRATION

Signal : TIC: BG050065.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.435	401	408	417	rBV	524643	881303	4.67%	0.894%
2	5.570	424	431	443	rVB2	155846	357702	1.89%	0.363%
3	5.946	487	495	502	rBV	119469	202720	1.07%	0.206%
4	7.285	716	723	729	rBV	85887	151126	0.80%	0.153%
5	7.497	747	759	771	rBV2	80399	190325	1.01%	0.193%
6	7.603	771	777	784	rVV	188174	319126	1.69%	0.324%
7	7.679	784	790	797	rVB2	136766	236138	1.25%	0.240%
8	7.961	832	838	851	rVB	95333	172642	0.91%	0.175%
9	8.349	894	904	910	rVV	3753127	6745408	35.73%	6.842%
10	8.419	910	916	924	rVV3	111151	205300	1.09%	0.208%
11	8.501	924	930	938	rVV	292799	512882	2.72%	0.520%
12	9.142	1032	1039	1046	rBV2	142295	297075	1.57%	0.301%
13	9.324	1063	1070	1077	rBV	96760	179330	0.95%	0.182%
14	9.400	1077	1083	1086	rBV	98152	172975	0.92%	0.175%
15	9.453	1086	1092	1098	rVV	234038	515074	2.73%	0.522%
16	9.512	1098	1102	1109	rVB	89755	153168	0.81%	0.155%
17	9.864	1152	1162	1172	rBV	100849	202157	1.07%	0.205%
18	9.964	1172	1179	1183	rBV	269811	512745	2.72%	0.520%
19	10.023	1186	1189	1199	rVB	139039	243435	1.29%	0.247%
20	10.176	1203	1215	1227	rBV5	240193	730007	3.87%	0.740%
21	10.276	1227	1232	1244	rVB2	221574	455775	2.41%	0.462%
22	10.417	1247	1256	1264	rBV2	596383	1157224	6.13%	1.174%
23	10.675	1293	1300	1307	rBV2	193010	378117	2.00%	0.384%
24	10.875	1312	1334	1339	rVV	7082593	18880984	100.00%	19.150%
25	10.928	1339	1343	1345	rVV2	189676	322589	1.71%	0.327%
26	10.969	1345	1350	1358	rVB	863875	1506039	7.98%	1.528%
27	11.627	1455	1462	1468	rBV	169317	314603	1.67%	0.319%
28	11.815	1486	1494	1499	rBV2	74749	155702	0.82%	0.158%
29	11.873	1499	1504	1517	rVV3	106419	333299	1.77%	0.338%
30	12.179	1551	1556	1563	rVB4	104077	242572	1.28%	0.246%
31	12.343	1574	1584	1589	rBV3	154899	299534	1.59%	0.304%
32	12.455	1595	1603	1610	rBV	2903705	5166063	27.36%	5.240%
33	12.572	1615	1623	1629	rVB2	660681	1241523	6.58%	1.259%
34	12.678	1629	1641	1648	rVB	2168006	3950656	20.92%	4.007%
35	13.037	1696	1702	1708	rBV4	126916	243741	1.29%	0.247%
36	13.454	1765	1773	1783	rVB	634097	1150708	6.09%	1.167%

LSC Area Percent Report

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Acq On : 31 Aug 2021 15:02
Operator : CG/JU
Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKG0

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-BG081921.M
Title : SVOA CALIBRATION

37	13.659	1802	1808	1818	rVB4	271809	686445	3.64%	0.696%
38	13.800	1819	1832	1840	rBV4	254485	817838	4.33%	0.829%
39	13.941	1851	1856	1862	rBV2	372941	722926	3.83%	0.733%
40	14.000	1862	1866	1868	rBV	119124	185928	0.98%	0.189%
41	14.029	1868	1871	1878	rVB	298155	423243	2.24%	0.429%
42	14.182	1891	1897	1900	rBV	160050	277493	1.47%	0.281%
43	14.335	1914	1923	1930	rVB3	375837	1061245	5.62%	1.076%
44	14.628	1959	1973	1978	rBV4	212471	555894	2.94%	0.564%
45	14.699	1978	1985	1991	rBV	3081645	5037387	26.68%	5.109%
46	14.764	1991	1996	2001	rBV	455132	696665	3.69%	0.707%
47	14.822	2001	2006	2014	rBV2	292965	550345	2.91%	0.558%
48	14.899	2014	2019	2028	rVB3	279301	576019	3.05%	0.584%
49	15.034	2031	2042	2050	rVB2	1430960	2681992	14.20%	2.720%
50	15.339	2089	2094	2098	rBV	119644	177014	0.94%	0.180%
51	15.527	2116	2126	2132	rBV2	130019	291441	1.54%	0.296%
52	15.621	2134	2142	2146	rBV	421993	705749	3.74%	0.716%
53	15.680	2146	2152	2160	rVB	1399094	2145124	11.36%	2.176%
54	15.762	2160	2166	2168	rBV3	117527	181067	0.96%	0.184%
55	15.821	2168	2176	2183	rVB5	310690	866474	4.59%	0.879%
56	15.897	2183	2189	2190	rBV4	166474	256137	1.36%	0.260%
57	16.109	2221	2225	2232	rVV3	147705	274230	1.45%	0.278%
58	16.197	2232	2240	2247	rVB	208604	403840	2.14%	0.410%
59	16.361	2260	2268	2271	rBV2	203387	401224	2.13%	0.407%
60	16.408	2272	2276	2280	rVB2	286638	481744	2.55%	0.489%
61	16.567	2293	2303	2309	rVV3	268665	585300	3.10%	0.594%
62	16.643	2309	2316	2320	rVV	477114	1119570	5.93%	1.136%
63	16.685	2320	2323	2327	rVV2	363326	562570	2.98%	0.571%
64	16.743	2327	2333	2338	rVV3	156755	308487	1.63%	0.313%
65	16.972	2362	2372	2379	rVB4	210372	620055	3.28%	0.629%
66	17.084	2384	2391	2394	rBV	140891	242395	1.28%	0.246%
67	17.166	2400	2405	2415	rVB	442871	782301	4.14%	0.793%
68	17.278	2415	2424	2430	rBV3	255310	642062	3.40%	0.651%
69	17.337	2430	2434	2437	rVV2	157153	214923	1.14%	0.218%
70	17.384	2437	2442	2445	rVV2	279103	457133	2.42%	0.464%
71	17.431	2445	2450	2455	rVV2	1178633	2013913	10.67%	2.043%
72	17.478	2455	2458	2462	rVV	453671	660719	3.50%	0.670%
73	17.525	2462	2466	2470	rVV5	219465	417502	2.21%	0.423%
74	17.589	2473	2477	2481	rVV3	246303	392732	2.08%	0.398%
75	17.630	2481	2484	2491	rVB2	144736	262639	1.39%	0.266%
76	17.812	2501	2515	2520	rBV	3627827	7299000	38.66%	7.403%

LSC Area Percent Report

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Acq On : 31 Aug 2021 15:02
Operator : CG/JU
Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKG0

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
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Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Title : SVOA CALIBRATION

77	17.889	2520	2528	2535	rBV4	555667	1195239	6.33%	1.212%
78	17.959	2535	2540	2546	rVB	809791	1214742	6.43%	1.232%
79	18.030	2546	2552	2558	rBV3	160784	350873	1.86%	0.356%
80	18.141	2565	2571	2577	rVB3	131592	232362	1.23%	0.236%
81	18.229	2581	2586	2591	rBV	359419	521982	2.76%	0.529%
82	18.312	2595	2600	2608	rVB2	536464	942916	4.99%	0.956%
83	18.388	2608	2613	2619	rVB	508453	726115	3.85%	0.736%
84	18.453	2619	2624	2633	rVB4	144206	265406	1.41%	0.269%
85	18.553	2633	2641	2648	rBV3	250776	812312	4.30%	0.824%
86	18.647	2654	2657	2662	rVB2	109395	151777	0.80%	0.154%
87	18.705	2662	2667	2672	rVV2	180761	312828	1.66%	0.317%
88	18.758	2672	2676	2681	rVB	147557	204241	1.08%	0.207%
89	19.187	2744	2749	2757	rVB5	95388	193805	1.03%	0.197%
90	19.504	2796	2803	2806	rVV2	93660	166125	0.88%	0.168%
91	19.769	2843	2848	2853	rVV	545330	861115	4.56%	0.873%
92	19.816	2853	2856	2865	rVB	149532	282658	1.50%	0.287%
93	20.180	2913	2918	2921	rBV	172185	232092	1.23%	0.235%
94	20.233	2921	2927	2929	rVV	277131	475152	2.52%	0.482%
95	20.286	2929	2936	2941	rVB	755625	1609571	8.52%	1.633%
96	20.867	3030	3035	3038	rBV2	134262	253922	1.34%	0.258%
97	21.654	3163	3169	3174	rVB	232131	377789	2.00%	0.383%
98	22.300	3272	3279	3286	rVB	121070	238903	1.27%	0.242%
99	24.662	3672	3681	3693	rVB	262826	726224	3.85%	0.737%
100	24.879	3710	3718	3730	rVB3	139745	427590	2.26%	0.434%

Sum of corrected areas: 98594271

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
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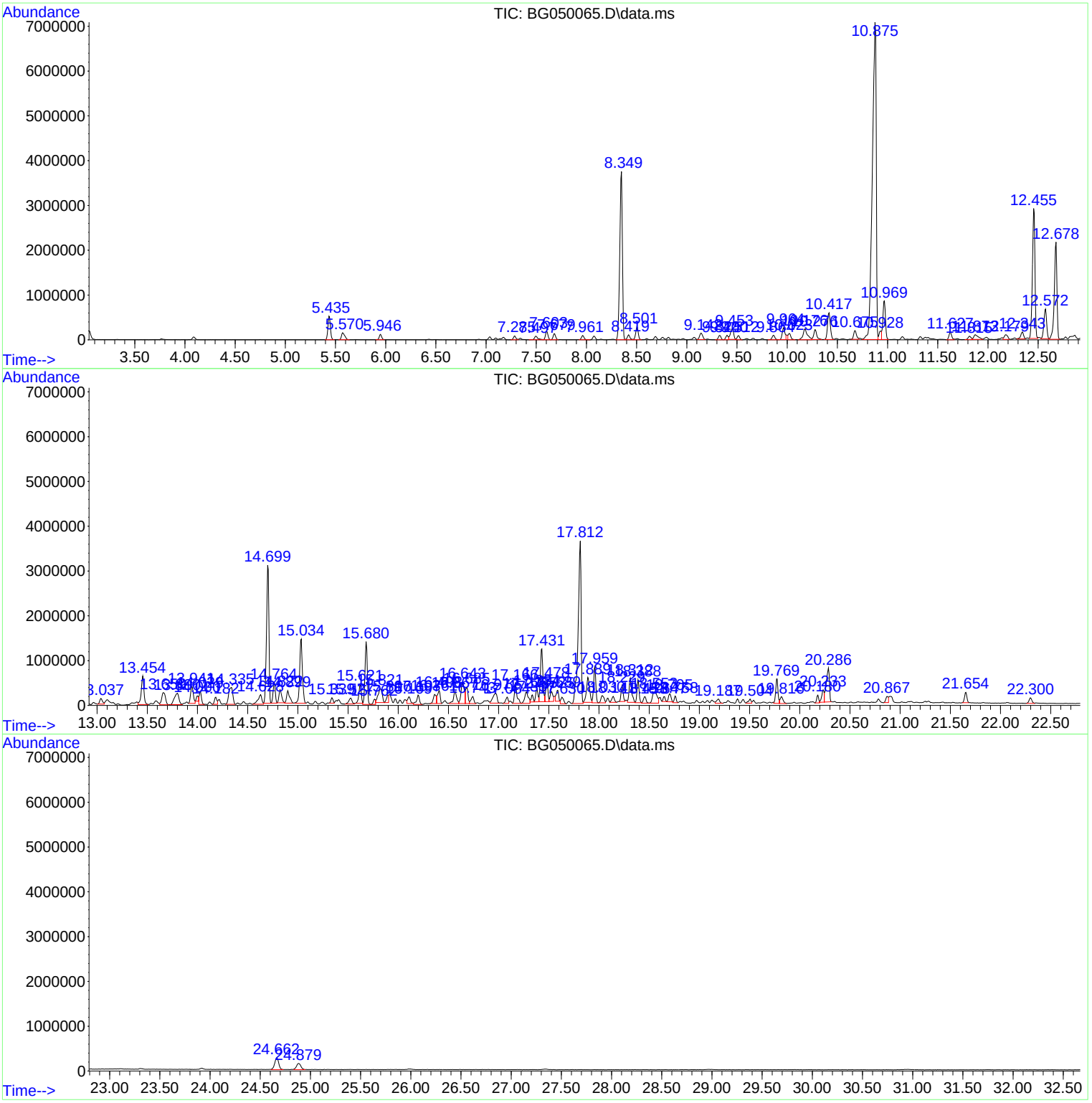
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKG0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
Data File : BG050065.D
Acq On : 31 Aug 2021 15:02
Operator : CG/JU
Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKG0

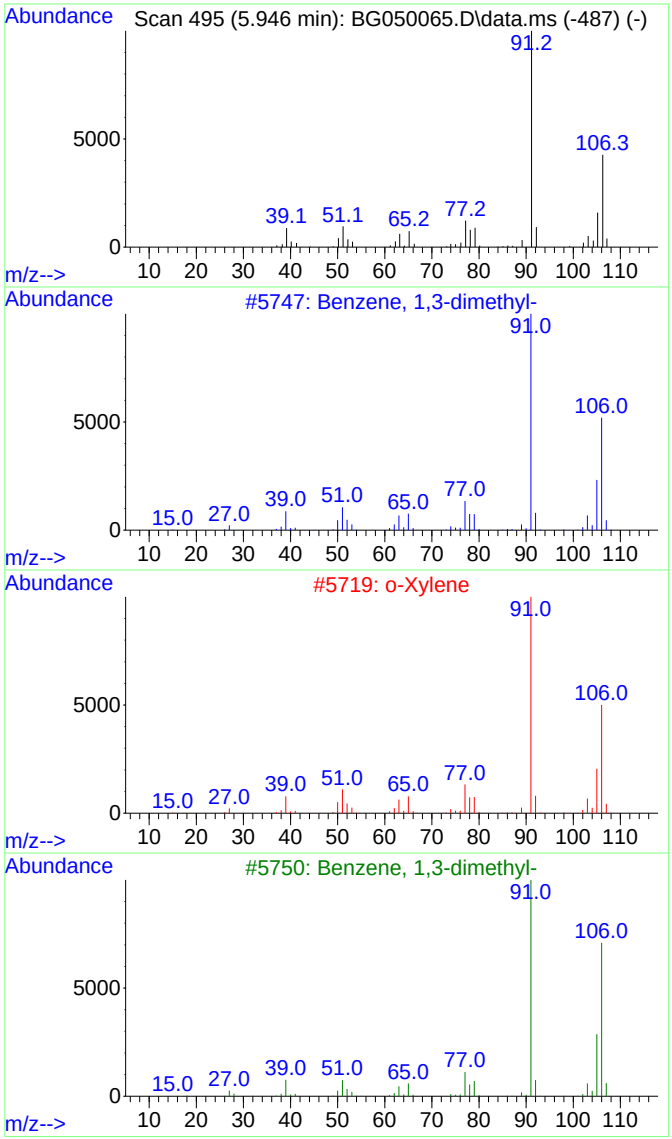
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

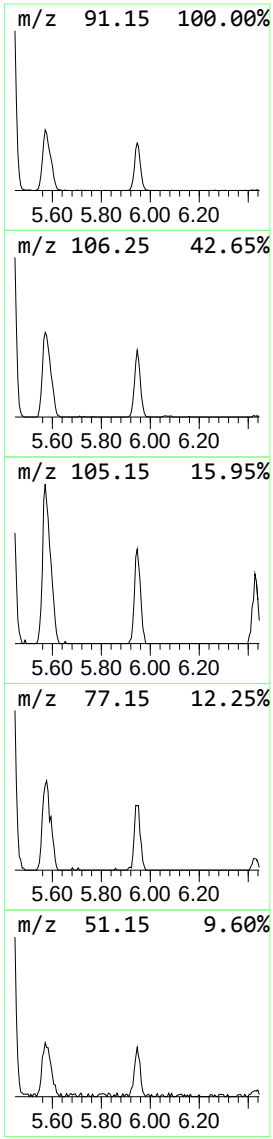
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1,3-dimethyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.946	23.48 ng/ul	202720	1,4-Dichlorobenzene-d4	7.961

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95
2	o-Xylene	106	C8H10	000095-47-6	95
3	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95
4	p-Xylene	106	C8H10	000106-42-3	95
5	o-Xylene	106	C8H10	000095-47-6	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
Data File : BG050065.D
Acq On : 31 Aug 2021 15:02
Operator : CG/JU
Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKG0

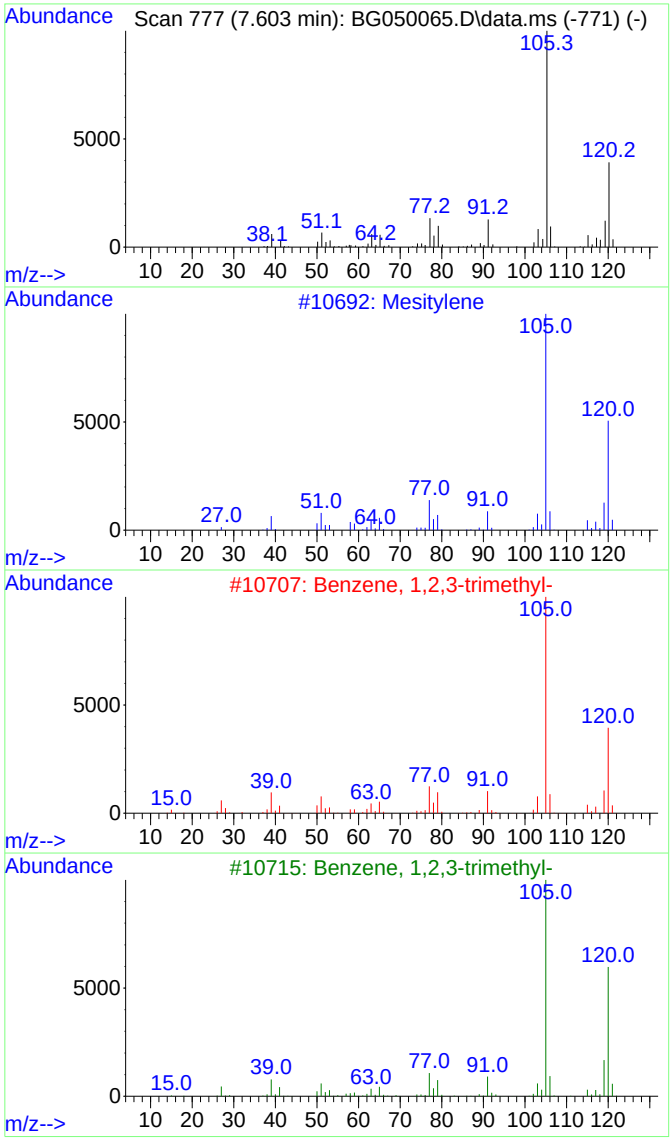
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

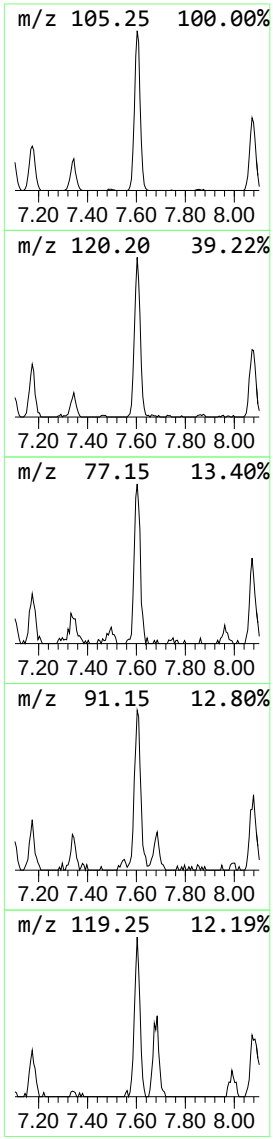
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Mesitylene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.603	36.97 ng/ul	319126	1,4-Dichlorobenzene-d4	7.961

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
Data File : BG050065.D
Acq On : 31 Aug 2021 15:02
Operator : CG/JU
Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKG0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

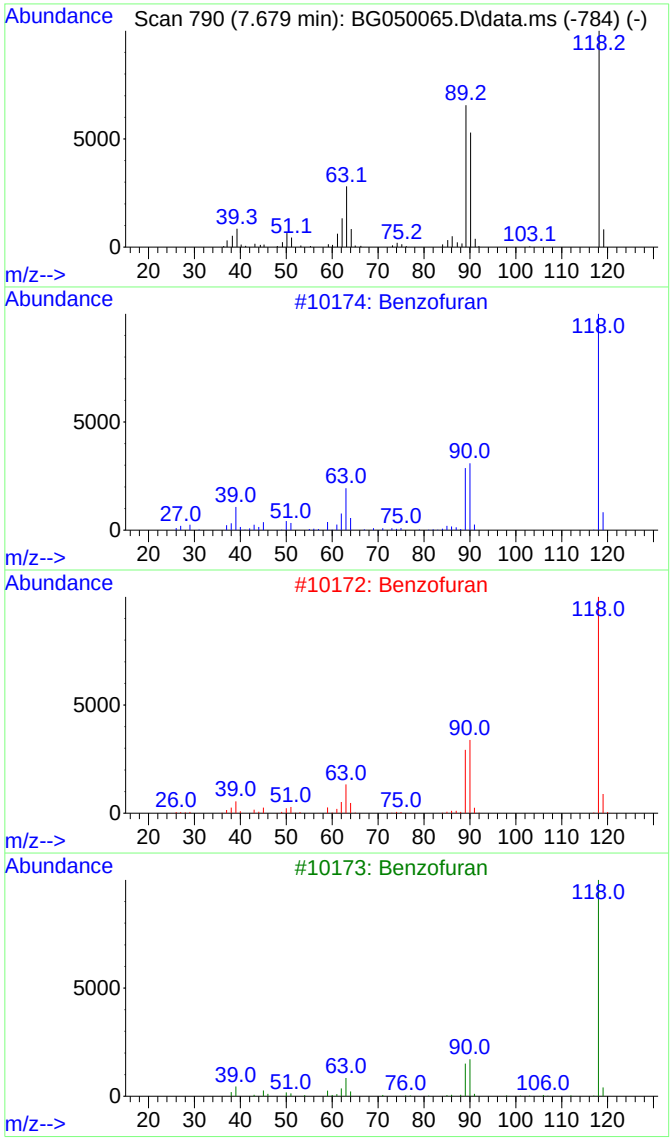
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

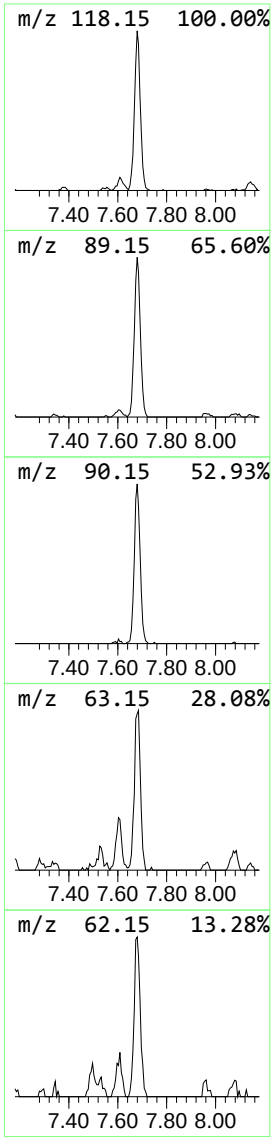
****

Peak Number 5 Benzofuran Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.679	27.36 ng/ul	236138	1,4-Dichlorobenzene-d4	7.961

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran	118	C8H6O	000271-89-6	91
2	Benzofuran	118	C8H6O	000271-89-6	91
3	Benzofuran	118	C8H6O	000271-89-6	90
4	3-Hydroxyphenylacetylene	118	C8H6O	010401-11-3	87
5	Benzocyclobuten-1(2H)-one	118	C8H6O	003469-06-5	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
Data File : BG050065.D
Acq On : 31 Aug 2021 15:02
Operator : CG/JU
Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKG0

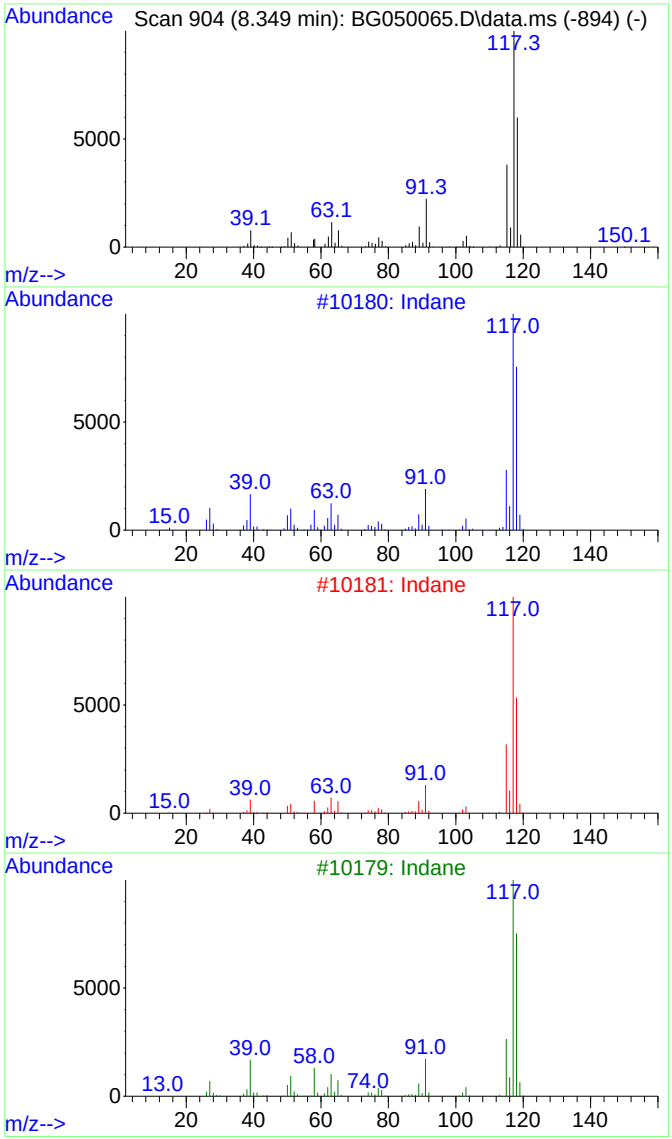
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

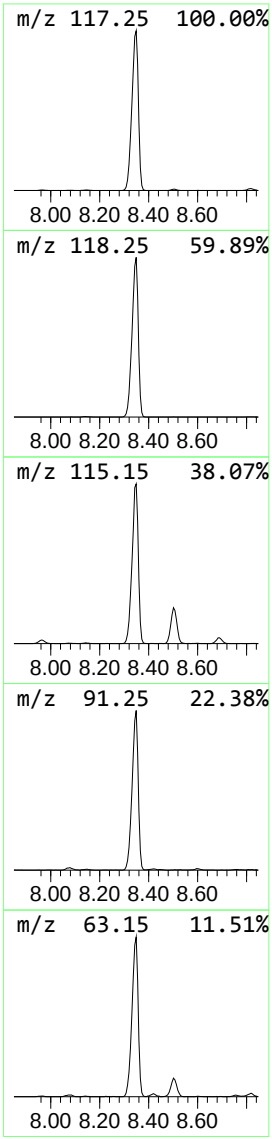
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.349	781.43 ng/ul	6745410	1,4-Dichlorobenzene-d4	7.961

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	94
2	Indane	118	C9H10	000496-11-7	94
3	Indane	118	C9H10	000496-11-7	81
4	Indane	118	C9H10	000496-11-7	70
5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	70





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
Data File : BG050065.D
Acq On : 31 Aug 2021 15:02
Operator : CG/JU
Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKG0

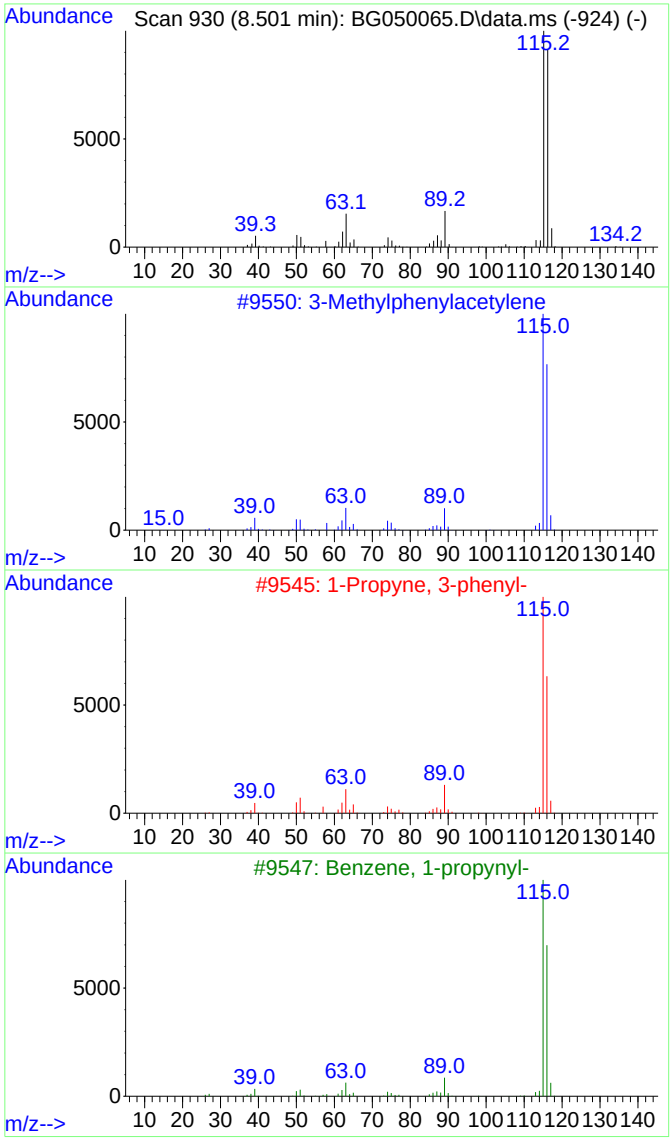
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

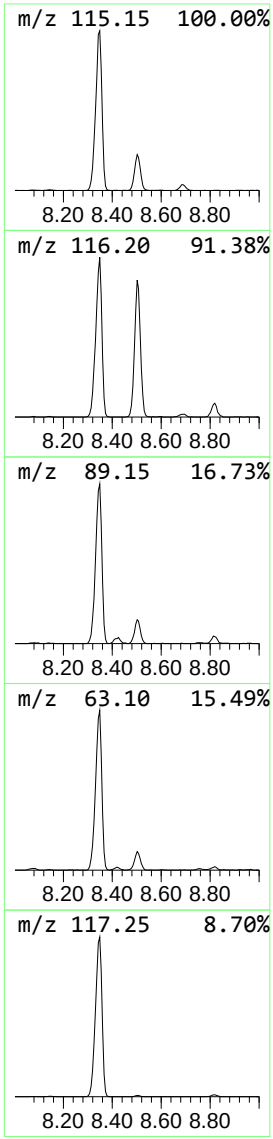
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 3-Methylphenylacetylene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.501	59.42 ng/ul	512882	1,4-Dichlorobenzene-d4	7.961

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Methylphenylacetylene	116	C9H8	000766-82-5	94
2	1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	94
3	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
4	Indene	116	C9H8	000095-13-6	91
5	Indene	116	C9H8	000095-13-6	91





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Acq On : 31 Aug 2021 15:02
Operator : CG/JU
Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKG0

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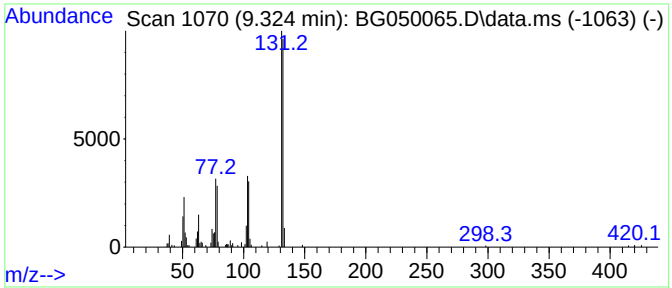
Quant Title : SVOA CALIBRATION

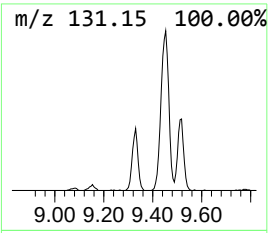
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

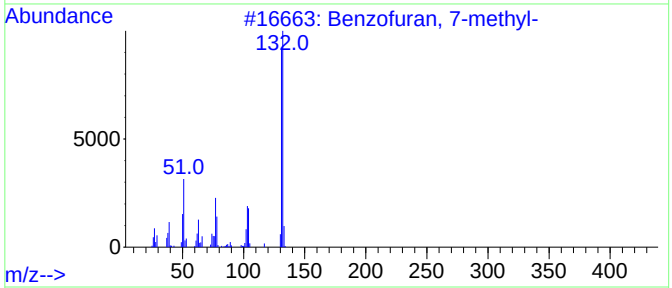
****
Peak Number 8 Benzofuran, 7-methyl- Concentration Rank 28

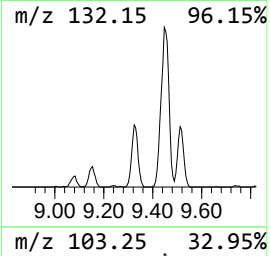
R.T.	EstConc	Area	Relative to ISTD	R.T.
9.324	20.77 ng/ul	179330	1,4-Dichlorobenzene-d4	7.961

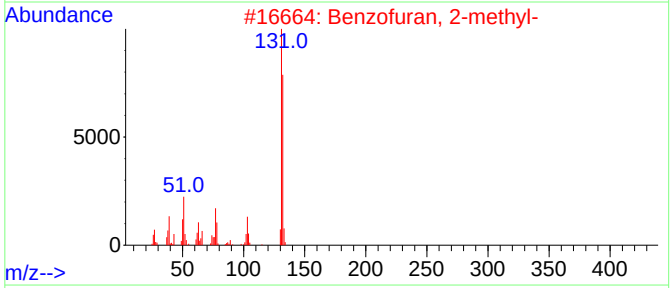
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	94
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	93
3	2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	90
4	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
5	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90

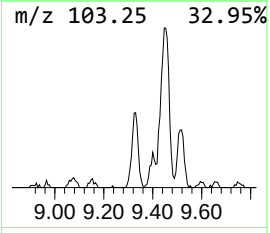


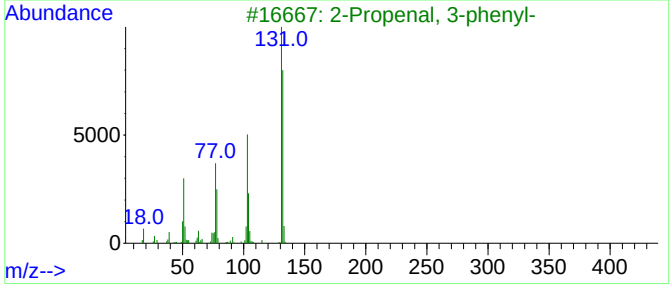


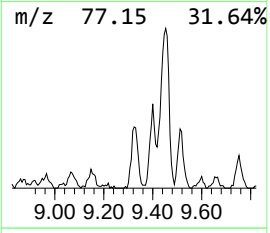


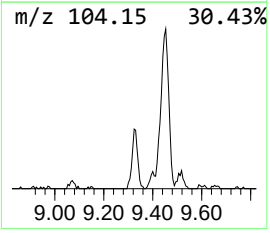












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Acq On : 31 Aug 2021 15:02
Operator : CG/JU
Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKG0

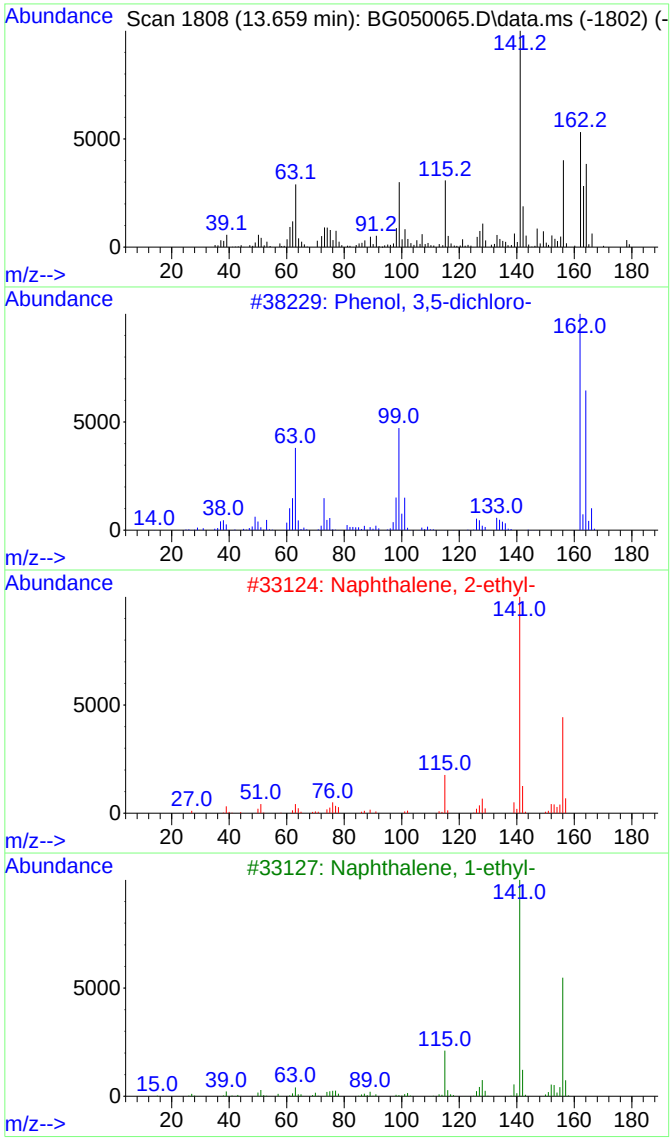
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

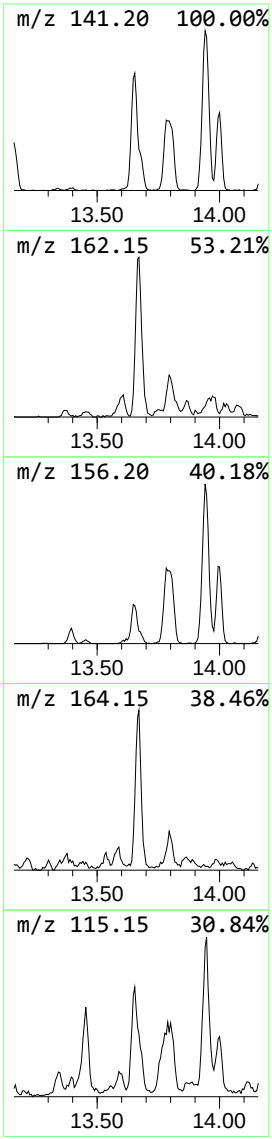
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Phenol, 3,5-dichloro- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.659	24.70 ng/ul	686445	Acenaphthene-d10	14.629

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	56
2	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	50
3	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	50
4	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	46
5	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	46





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Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

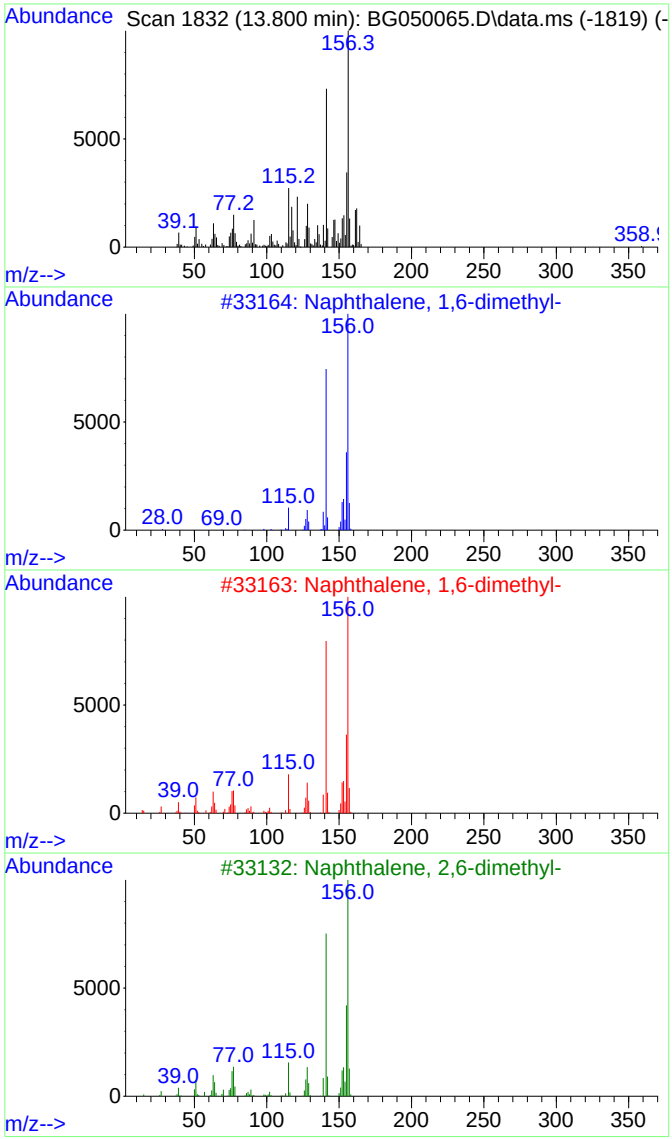
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

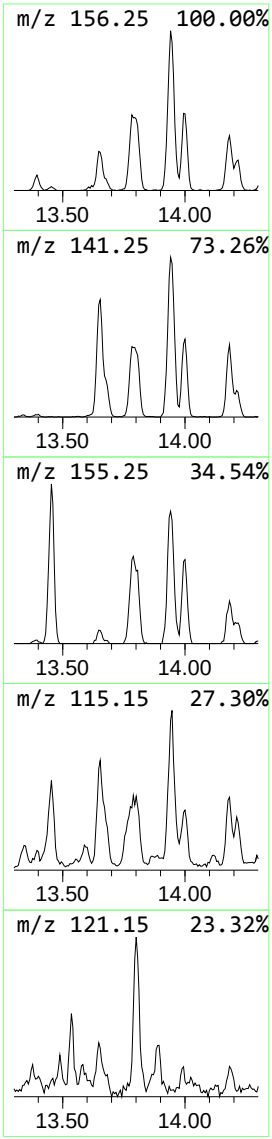
****

Peak Number 11 Naphthalene, 1,6-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.800	29.42 ng/ul	817838	Acenaphthene-d10	14.629

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
2	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
3	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95





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ALS Vial : 39 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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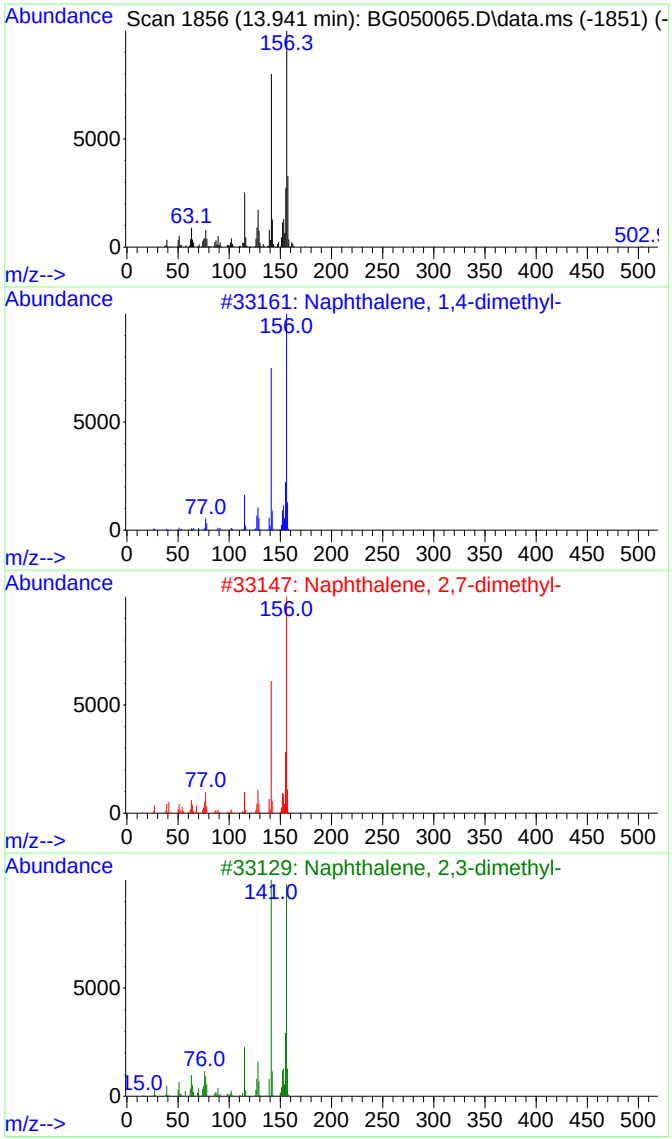
Quant Title : SVOA CALIBRATION

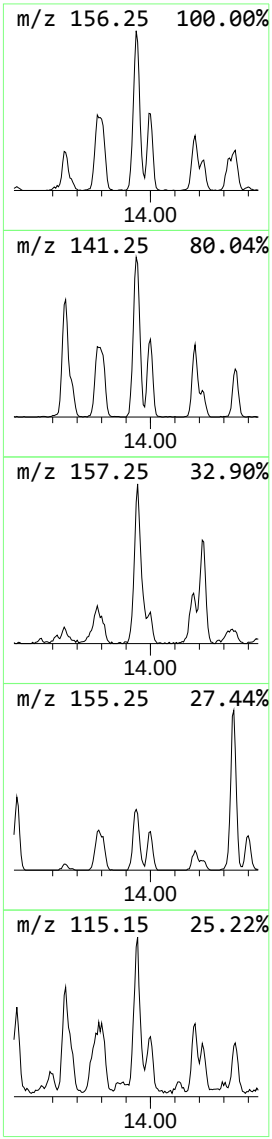
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Naphthalene, 1,4-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.941	26.01 ng/ul	722926	Acenaphthene-d10	14.629

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	94
2	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	94
3	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	93
4	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	93
5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	93





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Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKG0

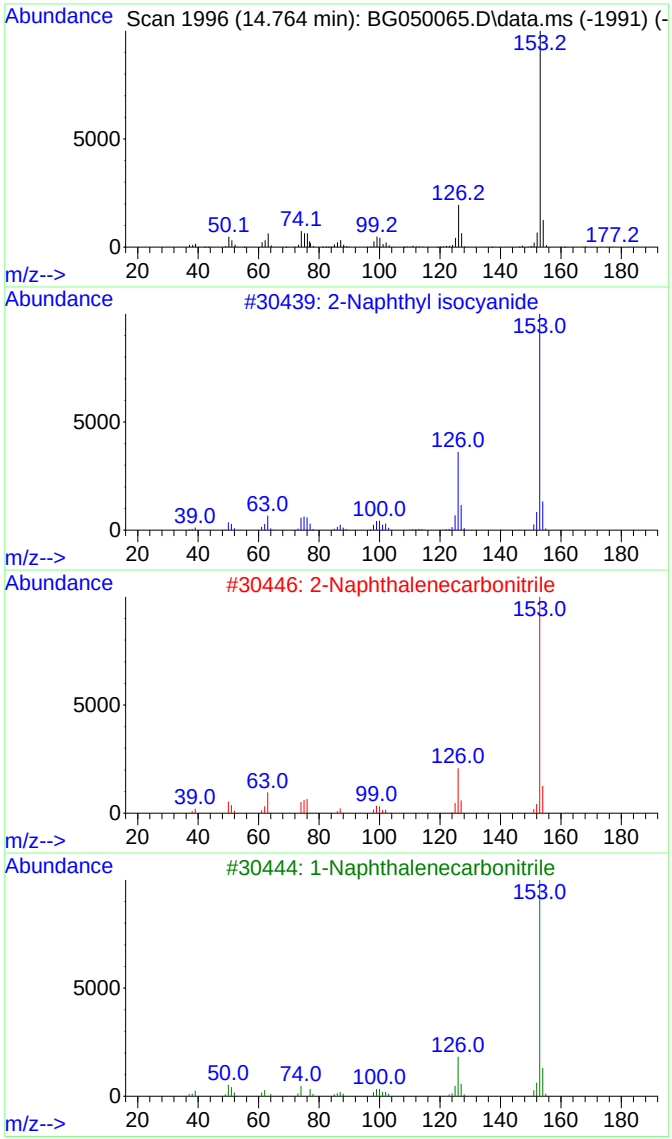
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

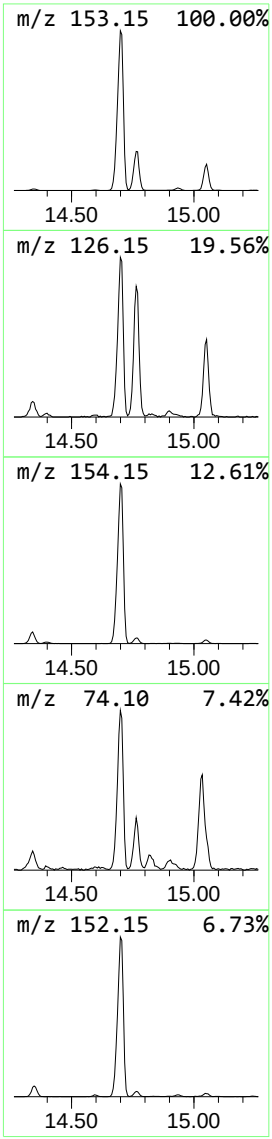
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 2-Naphthyl isocyanide Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.764	25.06 ng/ul	696665	Acenaphthene-d10	14.629

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthyl isocyanide	153	C11H7N	010124-78-4	95
2	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	94
3	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	93
4	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	91
5	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	91





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Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

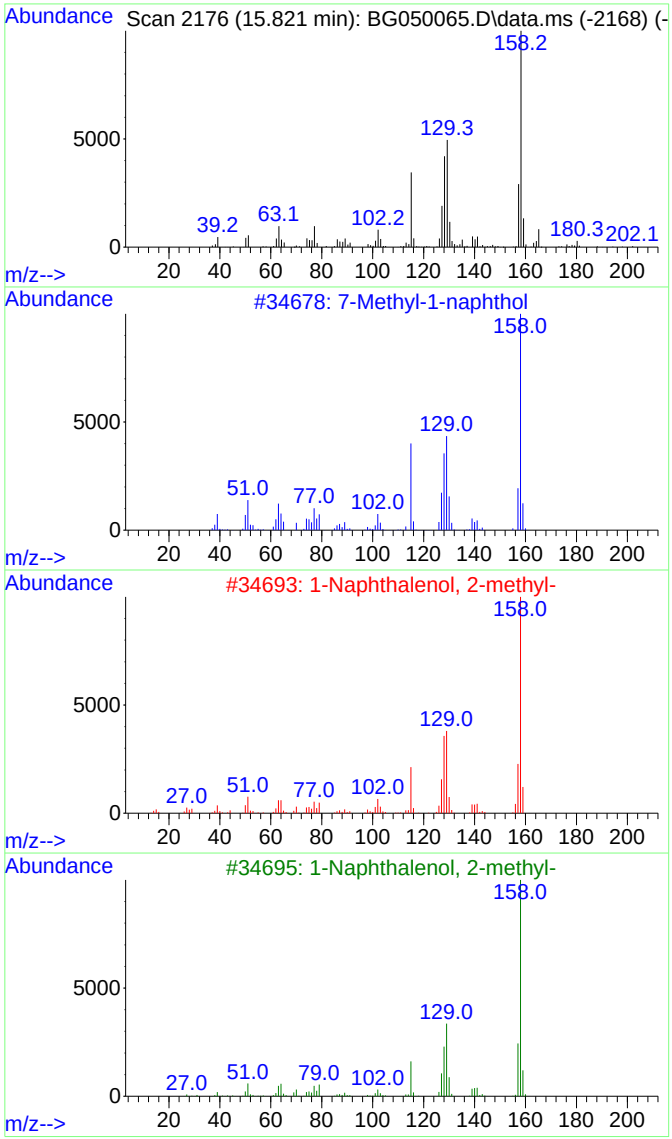
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

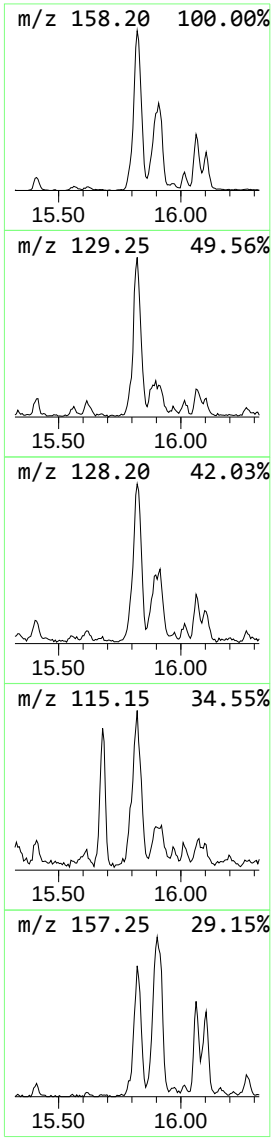
****

Peak Number 17 7-Methyl-1-naphthol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.821	31.17 ng/ul	866474	Acenaphthene-d10	14.629

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Methyl-1-naphthol	158	C11H10O	006939-33-9	94
2	1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	94
3	1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	94
4	1-Methyl-2-naphthol	158	C11H10O	001076-26-2	93
5	Phenol, 3-chloro-5-methoxy-	158	C7H7ClO2	065262-96-6	50





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ALS Vial : 39 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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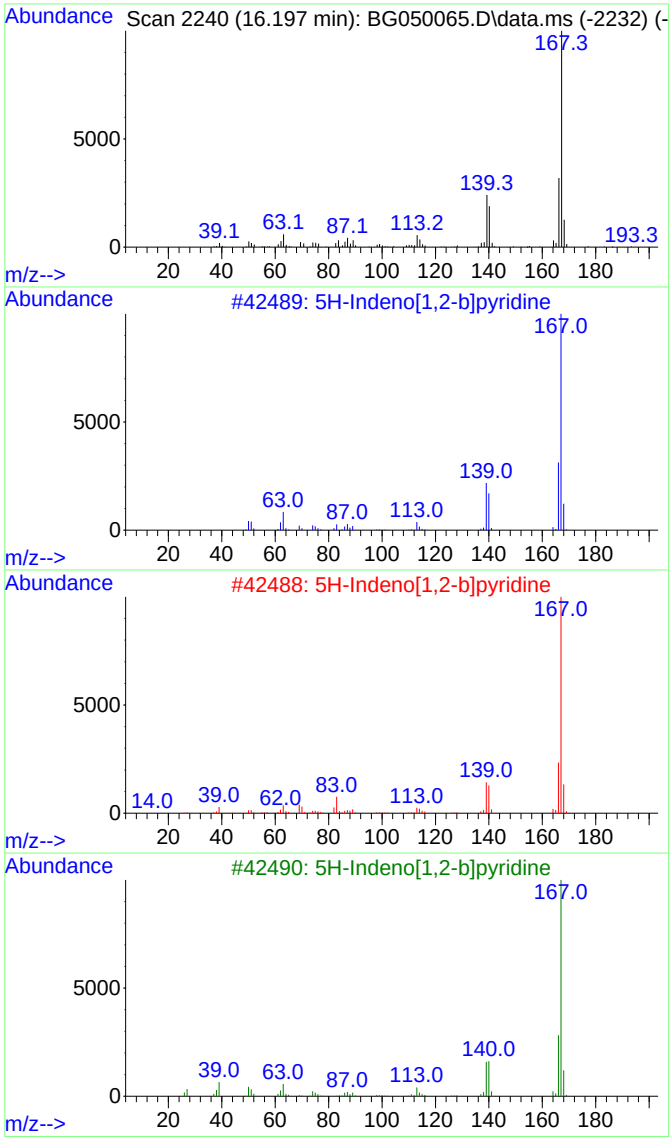
Quant Title : SVOA CALIBRATION

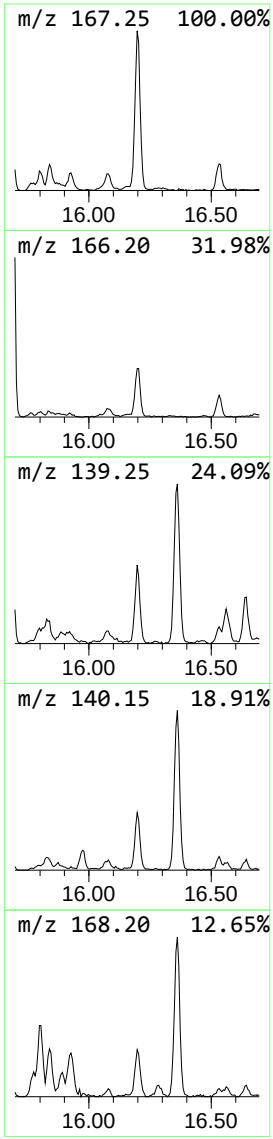
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 5H-Indeno[1,2-b]pyridine Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.197	17.67 ng/ul	403840	Phenanthrene-d10	17.384

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	91
2	5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	91
3	5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	90
4	Carbazole	167	C12H9N	000086-74-8	90
5	2-Naphthylacetonitrile	167	C12H9N	007498-57-9	87





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Sample : M3464-07
Misc :
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Instrument : BNA_G
ClientSampleId : DBKG0

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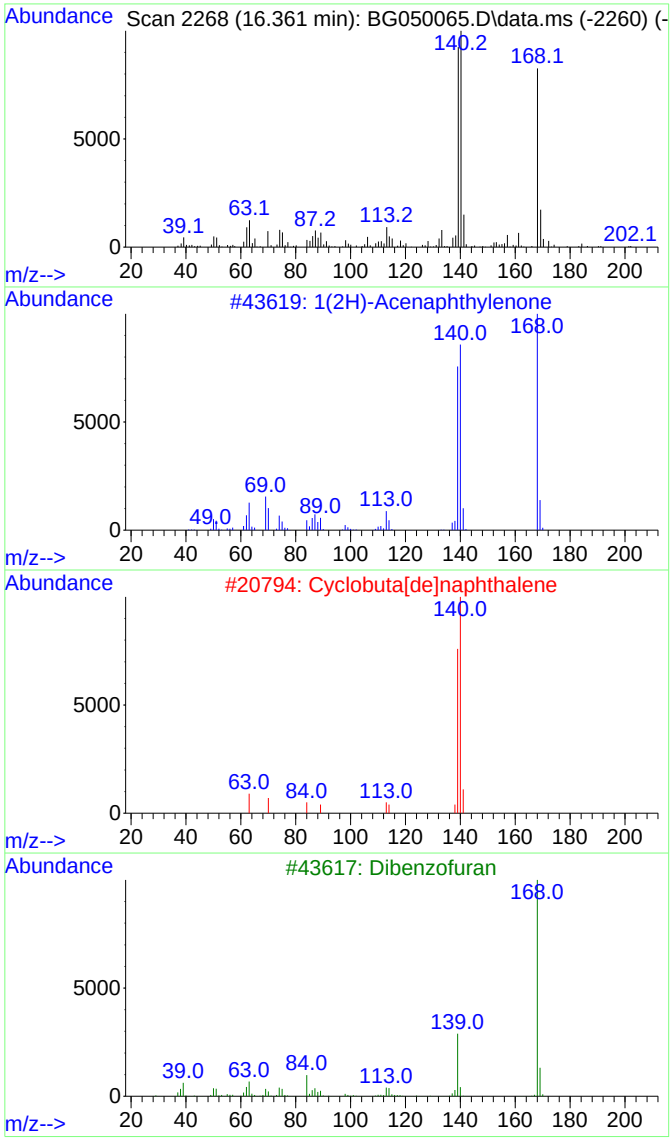
Quant Title : SVOA CALIBRATION

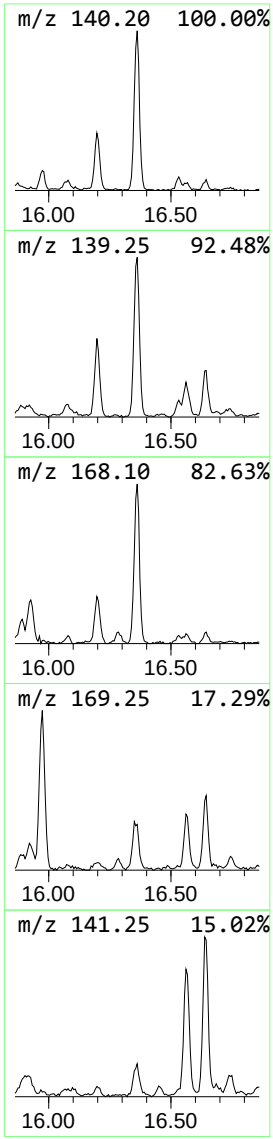
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 1(2H)-Acenaphthylenone Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.361	17.55 ng/ul	401224	Phenanthrene-d10	17.384

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(2H)-Acenaphthylenone	168	C12H8O	002235-15-6	93
2	Cyclobuta[de]naphthalene	140	C11H8	024973-91-9	64
3	Dibenzofuran	168	C12H8O	000132-64-9	49
4	Thiophene, 2,5-dipropyl-	168	C10H16S	054411-07-3	46
5	Benzaldehyde, 2-fluoro-3-hydroxy-	140	C7H5FO2	103438-86-4	43





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Instrument : BNA_G
ClientSampleId : DBKG0

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Quant Title : SVOA CALIBRATION

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

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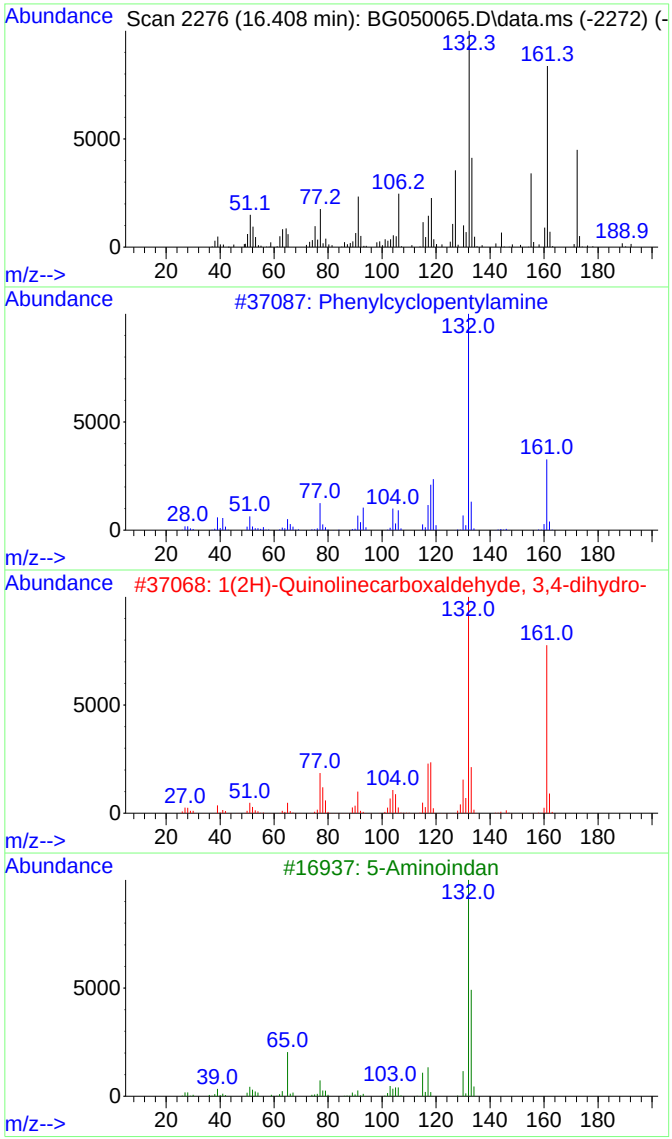
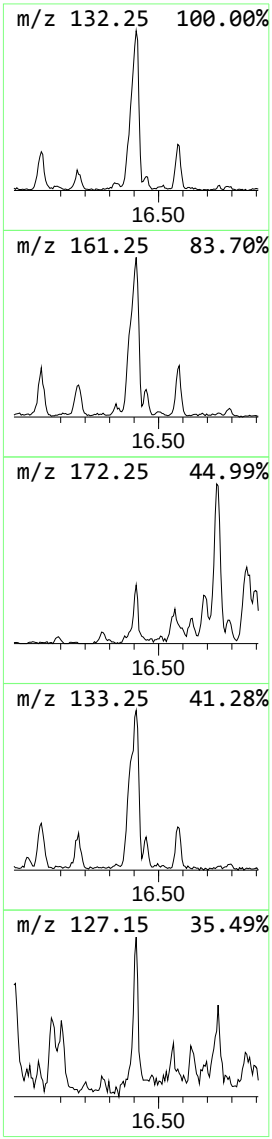
Peak Number 22 unknown-01 Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.408	21.08 ng/ul	481744	Phenanthrene-d10	17.384

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenylcyclopentylamine	161	C11H15N	040649-26-1	47
2	1(2H)-Quinolinecarboxaldehyde, 3...	161	C10H11NO	002739-16-4	43
3	5-Aminoindan	133	C9H11N	024425-40-9	38
4	3-Ethyl-5,6,7,8-tetrahydroquinoline	161	C11H15N	028971-02-0	35
5	3-Aminocoumarin	161	C9H7NO2	001635-31-0	30





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Quant Title : SVOA CALIBRATION

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

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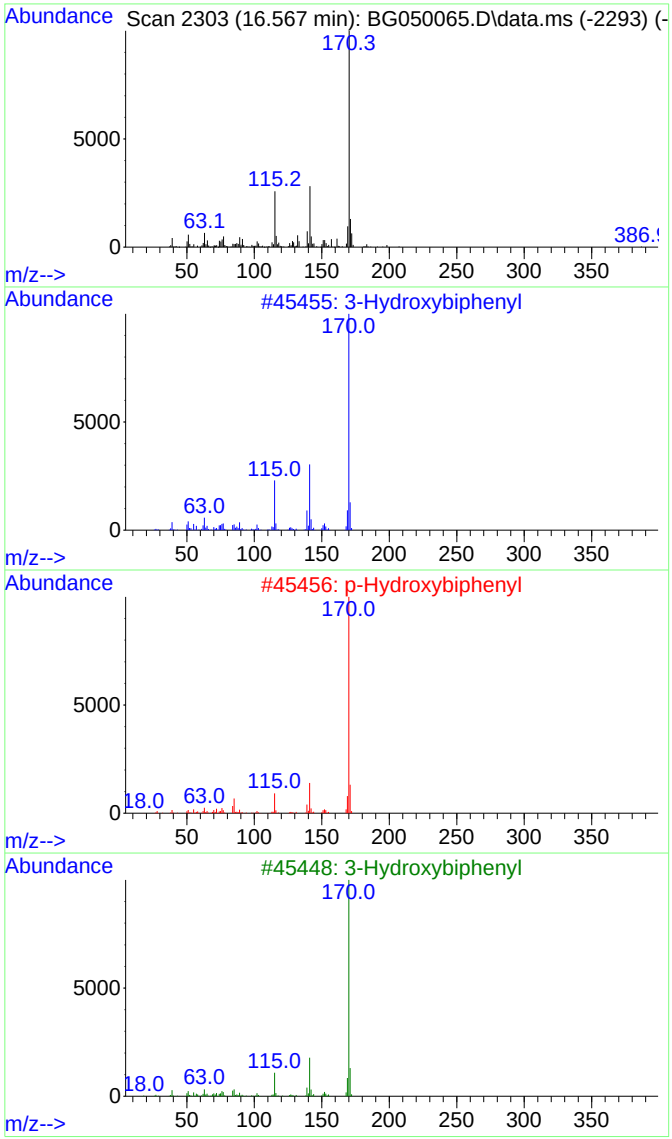
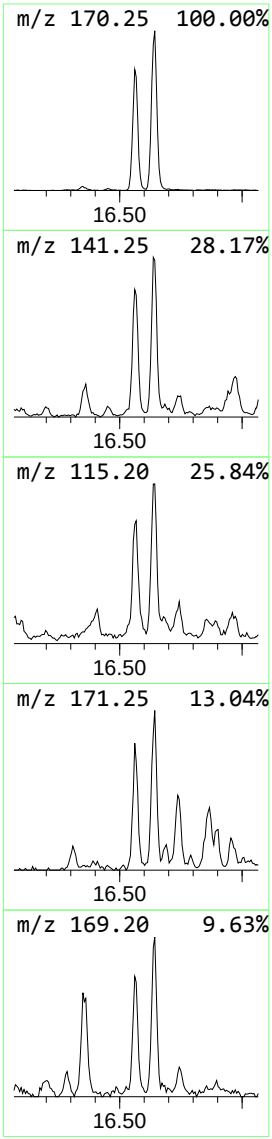
Peak Number 23 3-Hydroxybiphenyl Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.567	25.61 ng/ul	585300	Phenanthrene-d10	17.384

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hydroxybiphenyl	170	C12H10O	000580-51-8	96
2	p-Hydroxybiphenyl	170	C12H10O	000092-69-3	94
3	3-Hydroxybiphenyl	170	C12H10O	000580-51-8	94
4	3-Hydroxybiphenyl	170	C12H10O	000580-51-8	93
5	p-Hydroxybiphenyl	170	C12H10O	000092-69-3	91





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Acq On : 31 Aug 2021 15:02
Operator : CG/JU
Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKG0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

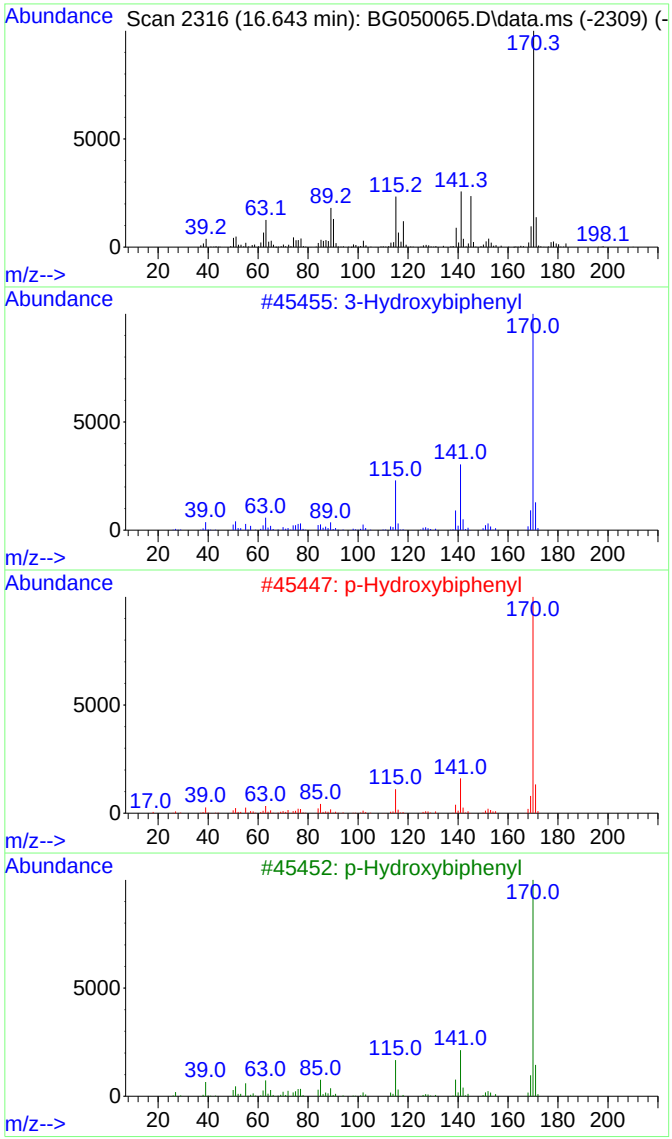
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

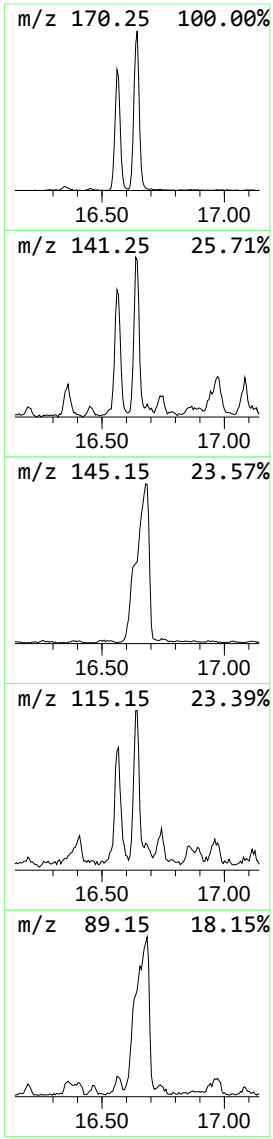
****

Peak Number 24 p-Hydroxybiphenyl Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.643	48.98 ng/ul	1119570	Phenanthrene-d10	17.384

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hydroxybiphenyl	170	C12H10O	000580-51-8	95
2	p-Hydroxybiphenyl	170	C12H10O	000092-69-3	93
3	p-Hydroxybiphenyl	170	C12H10O	000092-69-3	89
4	p-Hydroxybiphenyl	170	C12H10O	000092-69-3	76
5	3-Hydroxybiphenyl	170	C12H10O	000580-51-8	76





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Operator : CG/JU
Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKG0

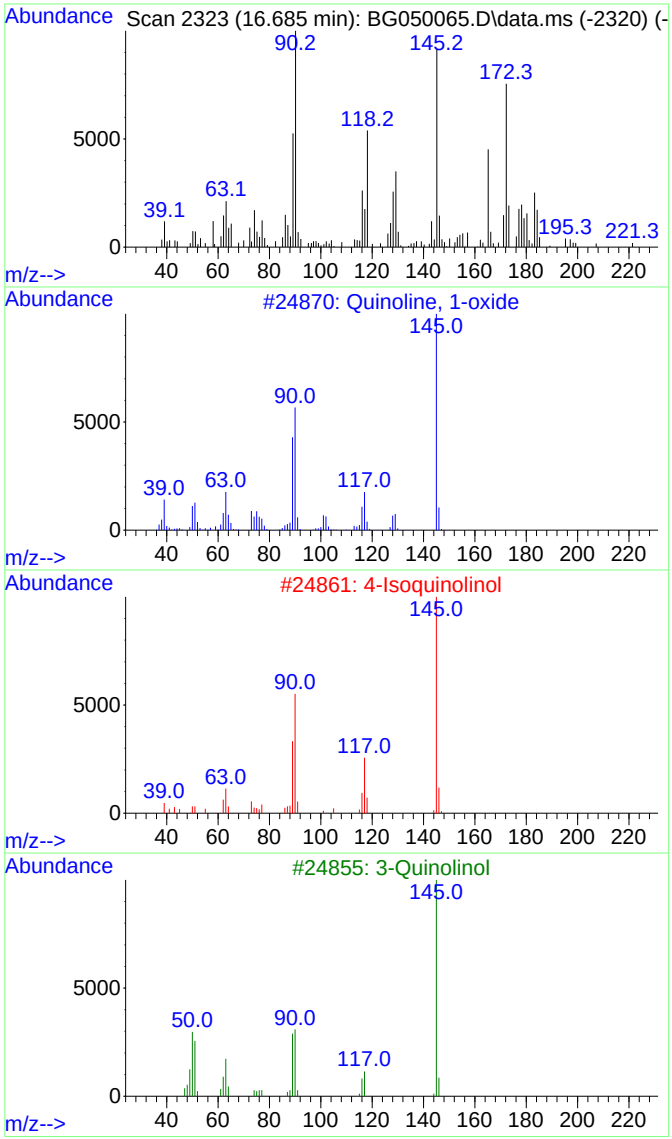
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

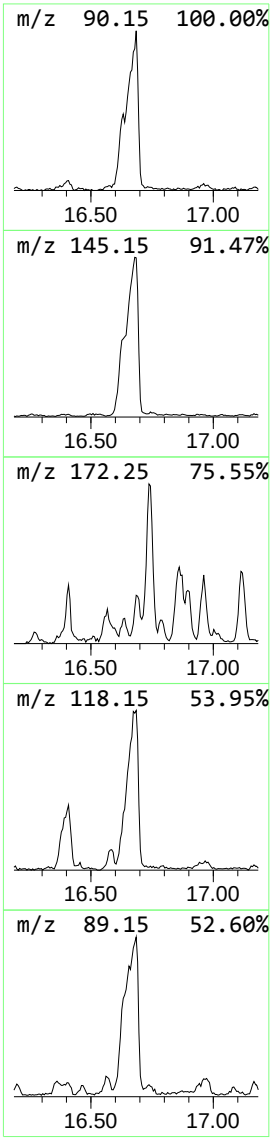
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 Quinoline, 1-oxide Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.685	24.61 ng/ul	562570	Phenanthrene-d10	17.384

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Quinoline, 1-oxide	145	C9H7NO	001613-37-2	64
2	4-Isoquinolinol	145	C9H7NO	003336-49-0	62
3	3-Quinolinol	145	C9H7NO	000580-18-7	47
4	Bicyclo[4.1.0]hepta-1,3,5-triene	90	C7H6	004646-69-9	46
5	Isoquinoline, 2-oxide	145	C9H7NO	001532-72-5	38





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Operator : CG/JU
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Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKG0

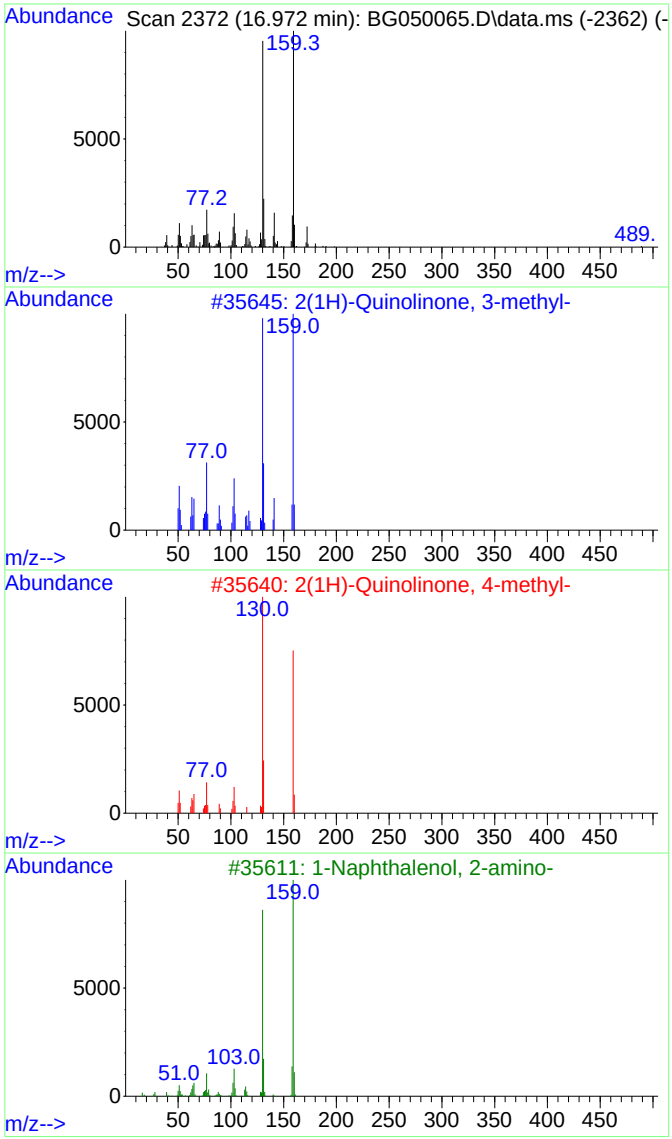
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

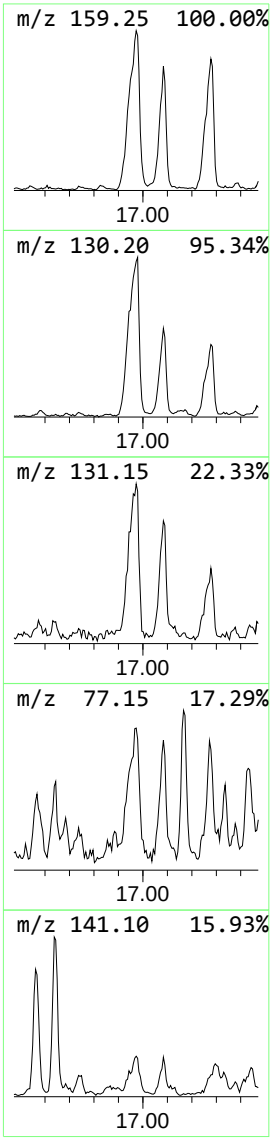
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 2(1H)-Quinolinone, 3-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.972	27.13 ng/ul	620055	Phenanthrene-d10	17.384

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(1H)-Quinolinone, 3-methyl-	159	C10H9NO	002721-59-7	97
2	2(1H)-Quinolinone, 4-methyl-	159	C10H9NO	000607-66-9	96
3	1-Naphthalenol, 2-amino-	159	C10H9NO	000606-41-7	94
4	1-Naphthalenol, 6-amino-	159	C10H9NO	023894-12-4	91
5	6-Methyl-4(1H)-quinolinone	159	C10H9NO	023432-40-8	91





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Operator : CG/JU
Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKG0

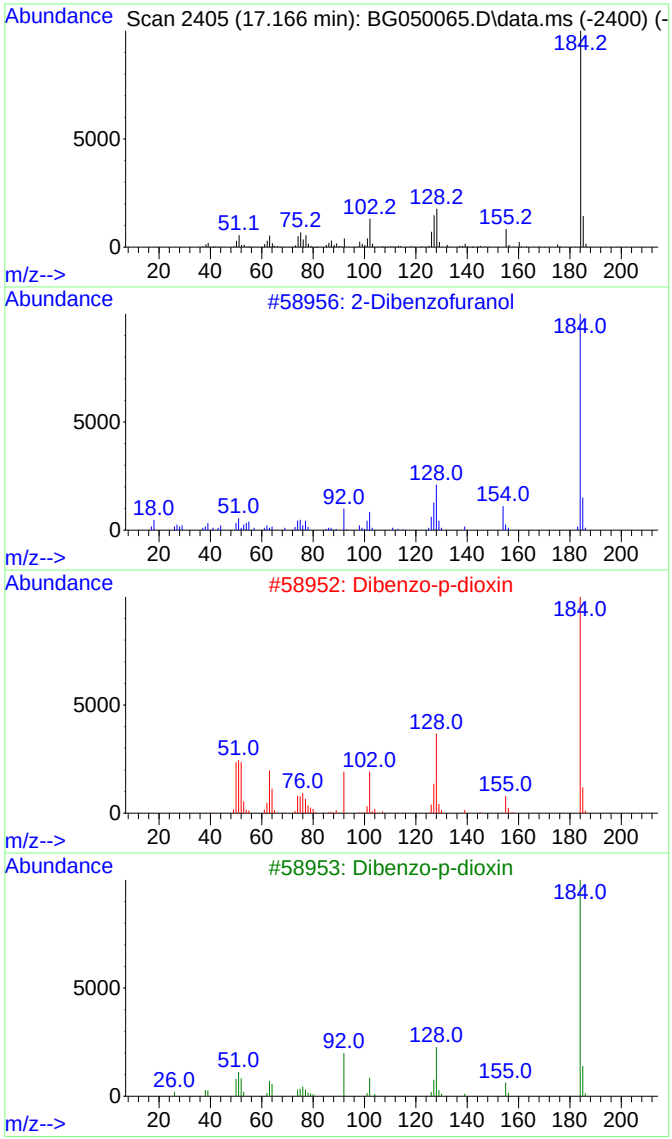
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

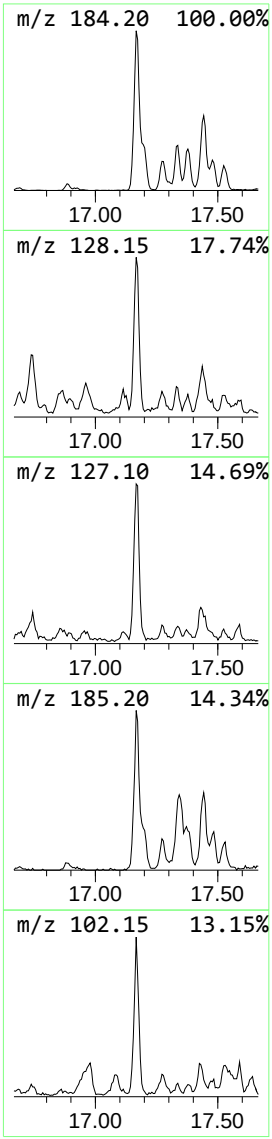
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 29 2-Dibenzofuranol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.166	34.23 ng/ul	782301	Phenanthrene-d10	17.384

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Dibenzofuranol	184	C12H8O2	000086-77-1	91
2	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	87
3	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	87
4	2-Dibenzofuranol	184	C12H8O2	000086-77-1	86
5	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	64





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Instrument :
BNA_G
ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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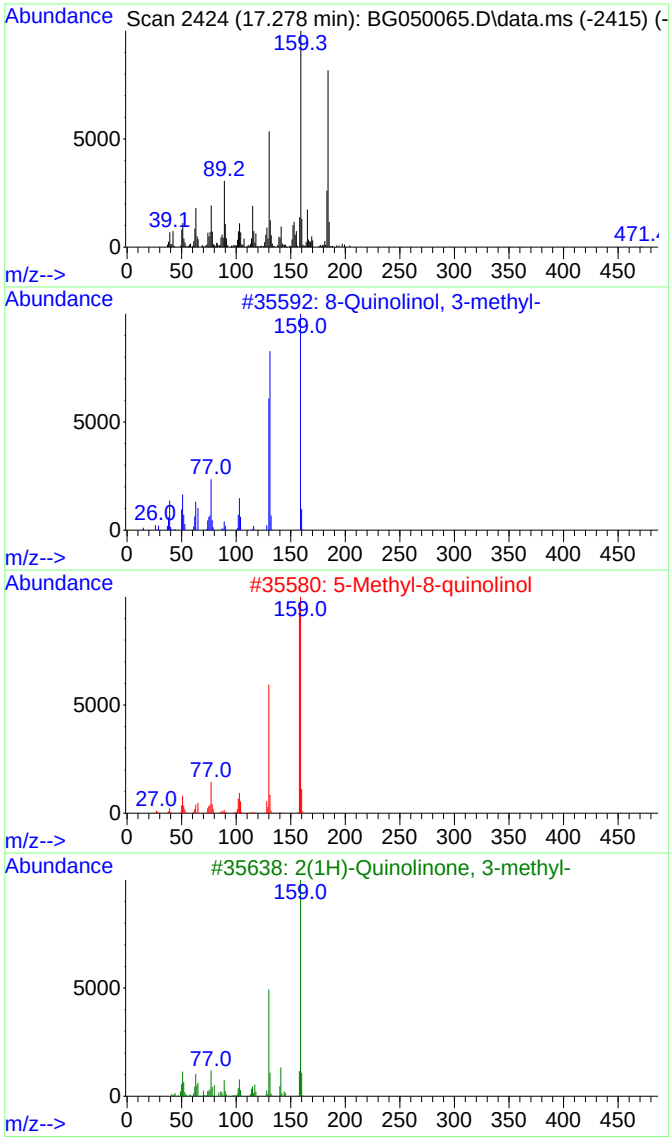
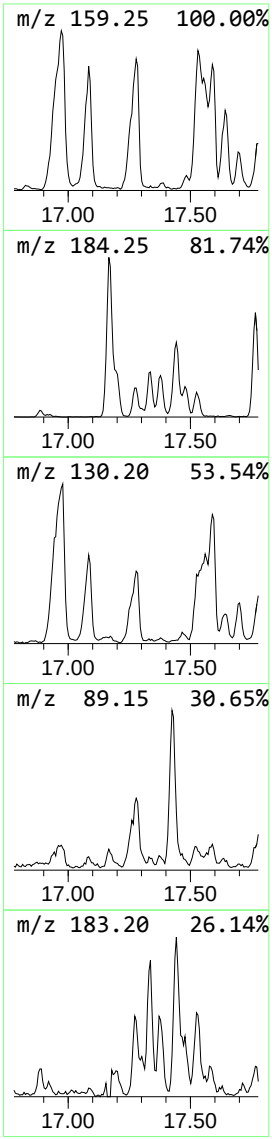
Peak Number 30 8-Quinolinol, 3-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.278	28.09 ng/ul	642062	Phenanthrene-d10	17.384

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	8-Quinolinol, 3-methyl-	159	C10H9NO	075457-13-5	62
2	5-Methyl-8-quinolinol	159	C10H9NO	005541-67-3	60
3	2(1H)-Quinolinone, 3-methyl-	159	C10H9NO	002721-59-7	50
4	2(1H)-Quinolinone, 1-methyl-	159	C10H9NO	000606-43-9	50
5	2(1H)-Quinolinone, 1-methyl-	159	C10H9NO	000606-43-9	50





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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

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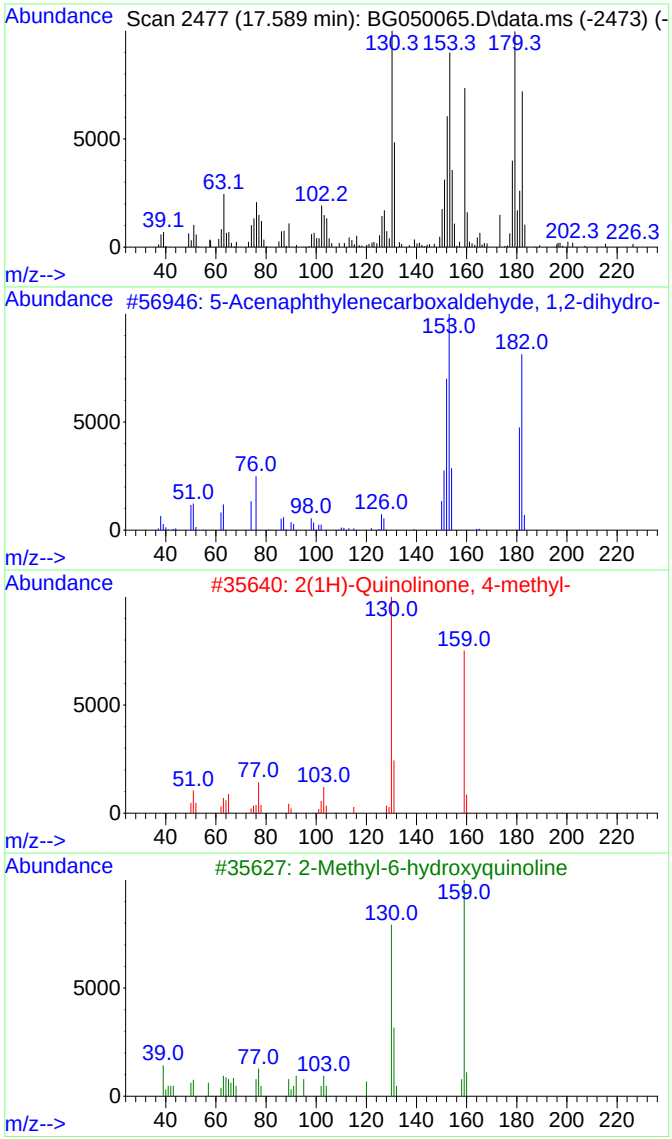
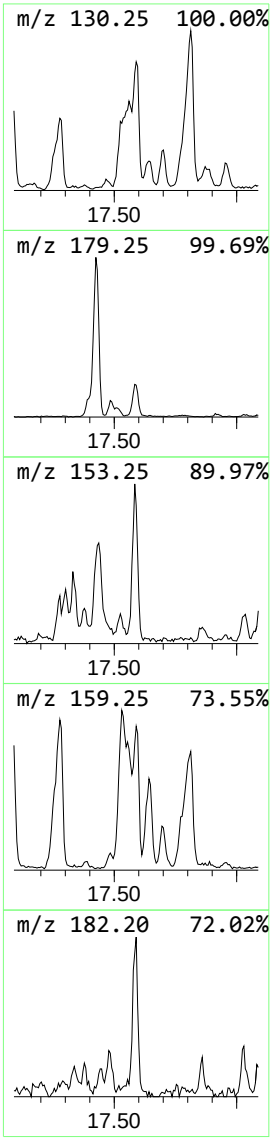
Peak Number 31 unknown-02 Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.589	17.18 ng/ul	392732	Phenanthrene-d10	17.384

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Acenaphthylenecarboxaldehyde, ...	182	C13H10O	1000362-64-3	47
2	2(1H)-Quinolinone, 4-methyl-	159	C10H9NO	000607-66-9	38
3	2-Methyl-6-hydroxyquinoline	159	C10H9NO	000613-21-8	30
4	Benzo[g]quinoline	179	C13H9N	000260-36-6	25
5	4-Amino-1-naphthol	159	C10H9NO	002834-90-4	18





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Operator : CG/JU
Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKG0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

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

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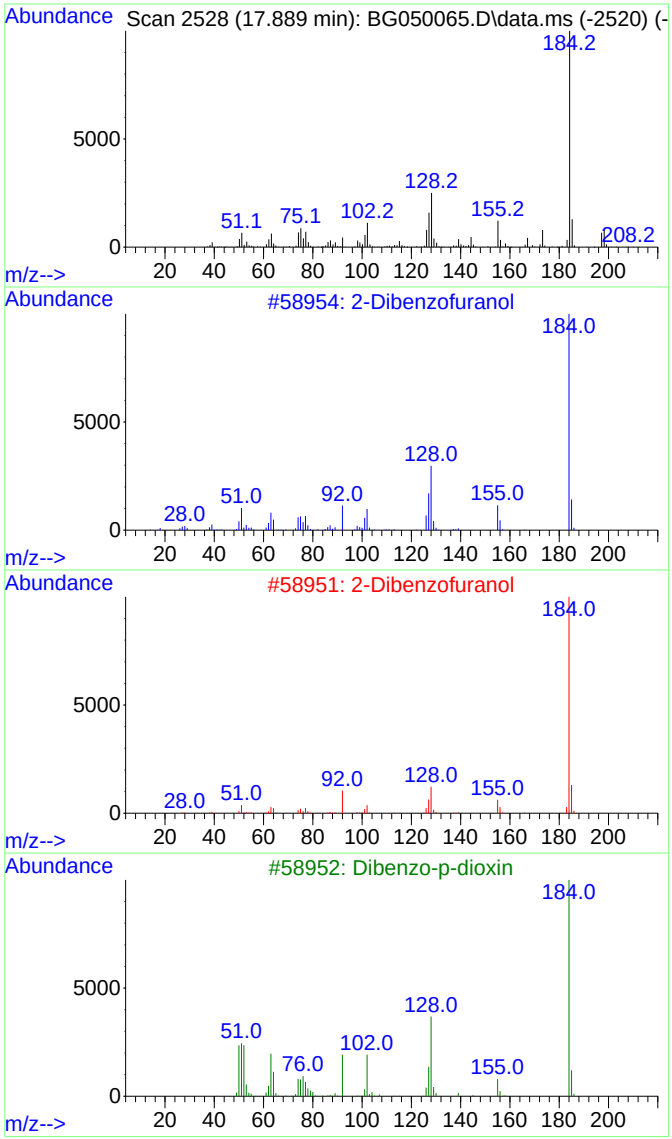
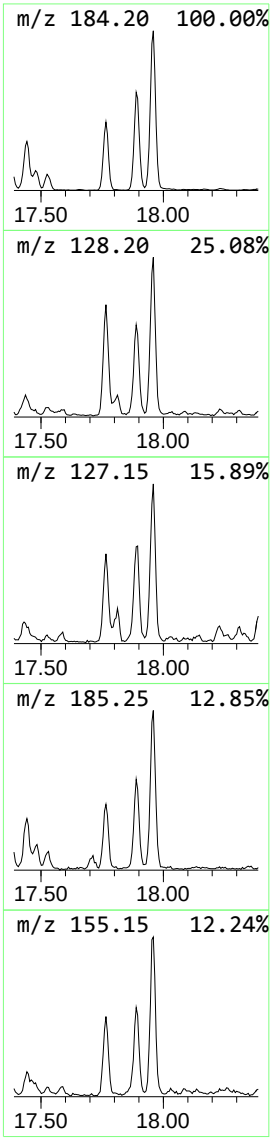
Peak Number 33 Dibenzo-p-dioxin Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.889	52.29 ng/ul	1195240	Phenanthrene-d10	17.384

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Dibenzofuranol	184	C12H8O2	000086-77-1	87
2	2-Dibenzofuranol	184	C12H8O2	000086-77-1	87
3	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	83
4	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	83
5	1H-Pyrazolo[3,4-b]quinolin-3-amine	184	C10H8N4	1000319-54-1	80





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Operator : CG/JU
Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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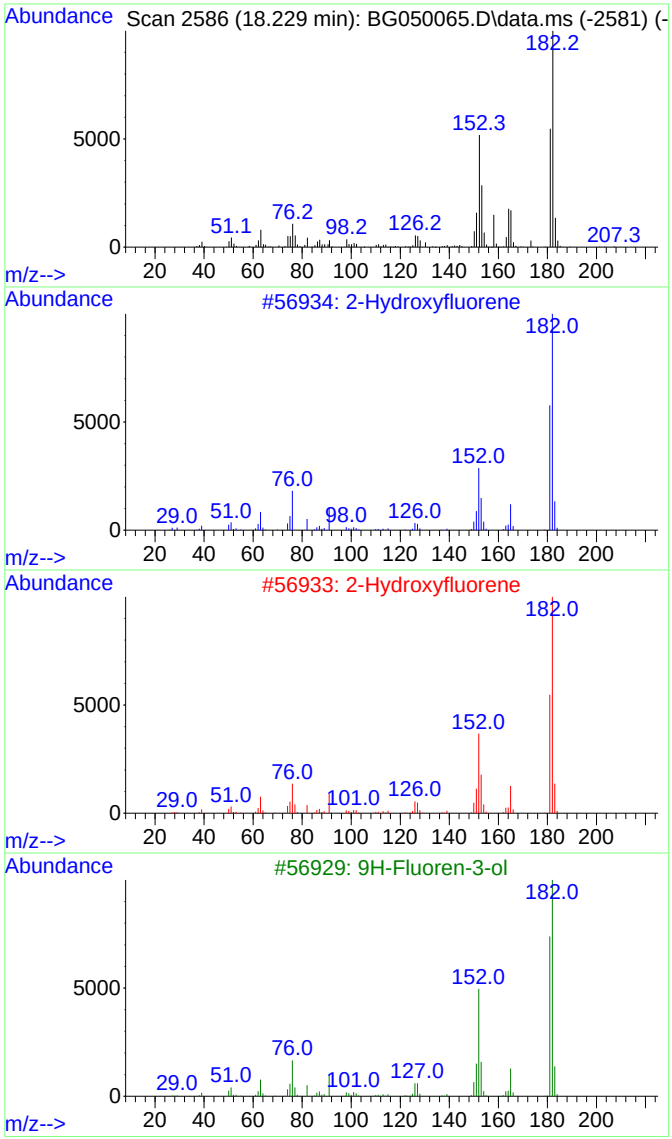
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

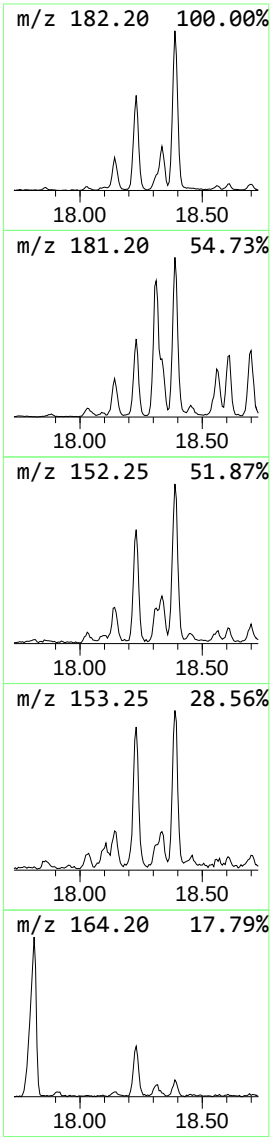
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 37 9H-Fluoren-3-ol Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.230	22.84 ng/ul	521982	Phenanthrene-d10	17.384

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hydroxyfluorene	182	C13H10O	002443-58-5	93
2	2-Hydroxyfluorene	182	C13H10O	002443-58-5	91
3	9H-Fluoren-3-ol	182	C13H10O	006344-67-8	87
4	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	70
5	3-(1-Naphthyl)acrylaldehyde	182	C13H10O	001504-73-0	64





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Instrument : BNA_G
ClientSampleId : DBKG0

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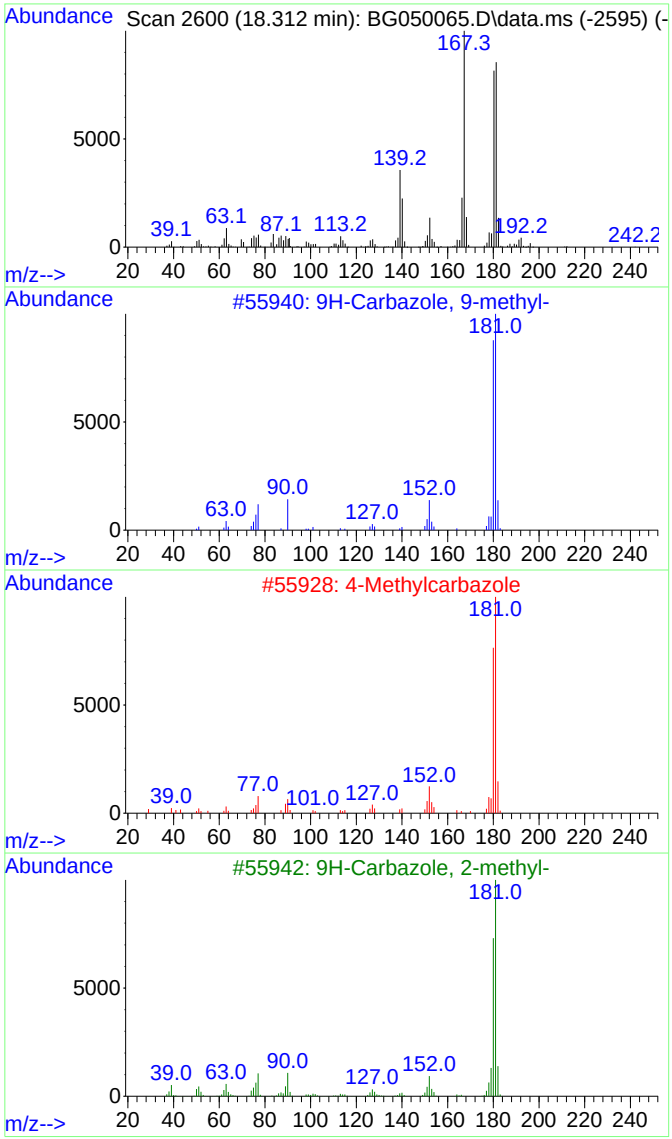
Quant Title : SVOA CALIBRATION

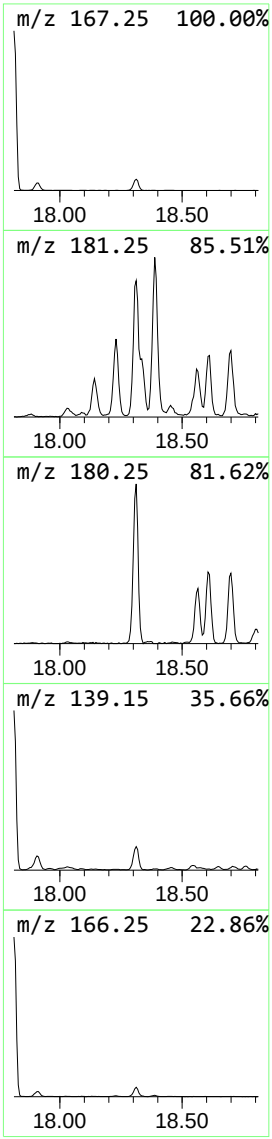
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 38 9H-Carbazole, 9-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.312	41.25 ng/ul	942916	Phenanthrene-d10	17.384

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Carbazole, 9-methyl-	181	C13H11N	001484-12-4	70
2	4-Methylcarbazole	181	C13H11N	003770-48-7	64
3	9H-Carbazole, 2-methyl-	181	C13H11N	003652-91-3	55
4	9H-Carbazole, 2-methyl-	181	C13H11N	003652-91-3	55
5	3-Methylcarbazole	181	C13H11N	004630-20-0	55





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ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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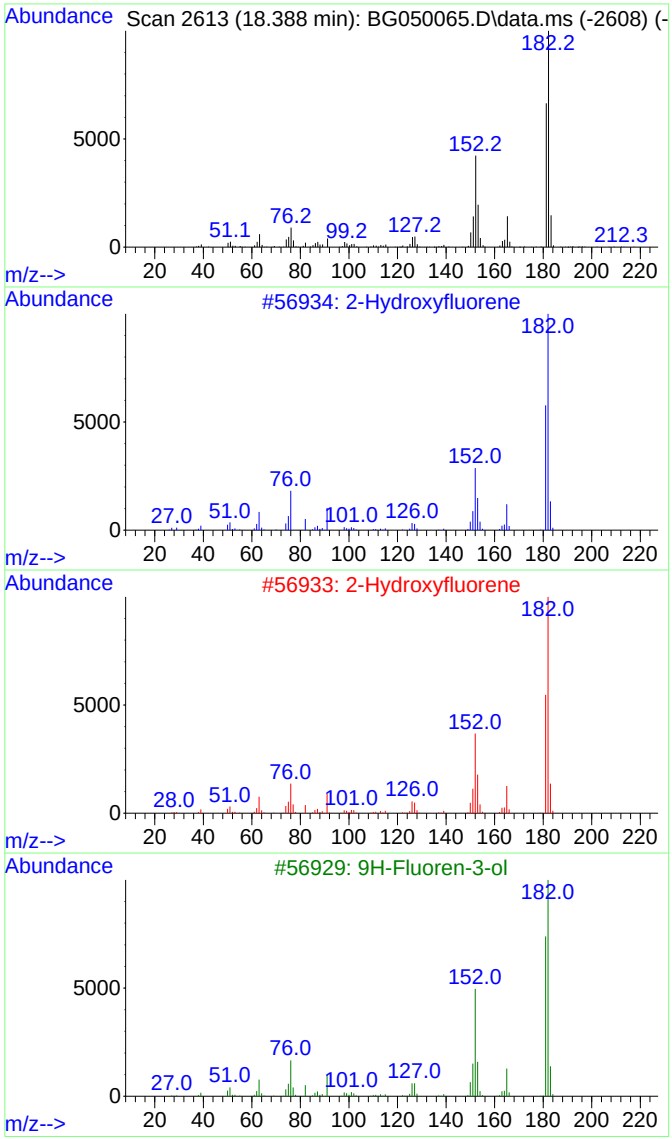
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

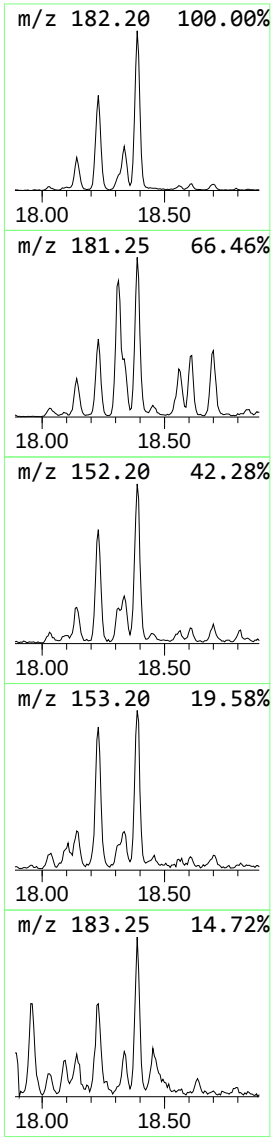
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 39 2-Hydroxyfluorene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.388	31.77 ng/ul	726115	Phenanthrene-d10	17.384

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hydroxyfluorene	182	C13H10O	002443-58-5	97
2	2-Hydroxyfluorene	182	C13H10O	002443-58-5	97
3	9H-Fluoren-3-ol	182	C13H10O	006344-67-8	93
4	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	90
5	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	72





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Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKG0

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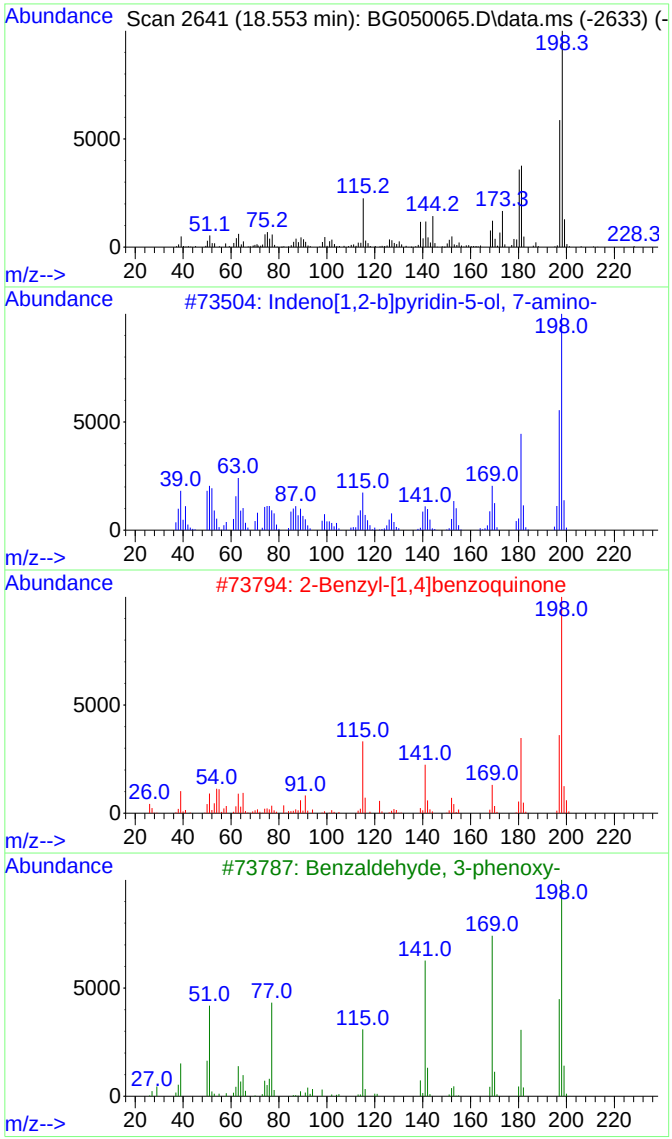
Quant Title : SVOA CALIBRATION

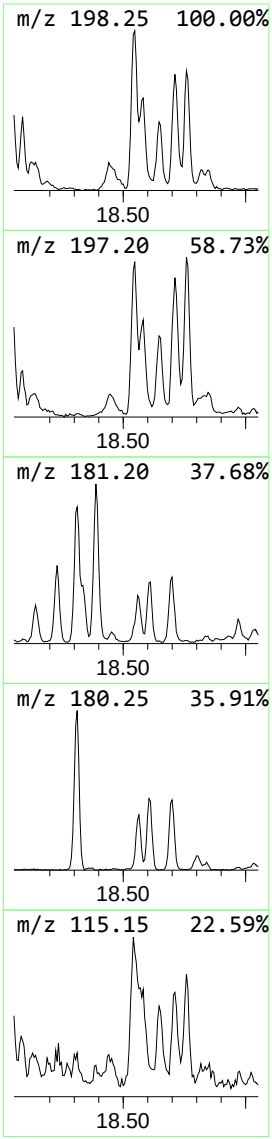
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 41 Indeno[1,2-b]pyridin-5-ol, ... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.553	35.54 ng/ul	812312	Phenanthrene-d10	17.384

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indeno[1,2-b]pyridin-5-ol, 7-amino-	198	C12H10N2O	339336-23-1	64
2	2-Benzyl-[1,4]benzoquinone	198	C13H10O2	038940-10-2	62
3	Benzaldehyde, 3-phenoxy-	198	C13H10O2	039515-51-0	62
4	Tacrine	198	C13H14N2	000321-64-2	58
5	1H-Imidazo[4,5-g]quinoxaline, 1,...	198	C11H10N4	1000319-12-2	49





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Acq On : 31 Aug 2021 15:02
Operator : CG/JU
Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKG0

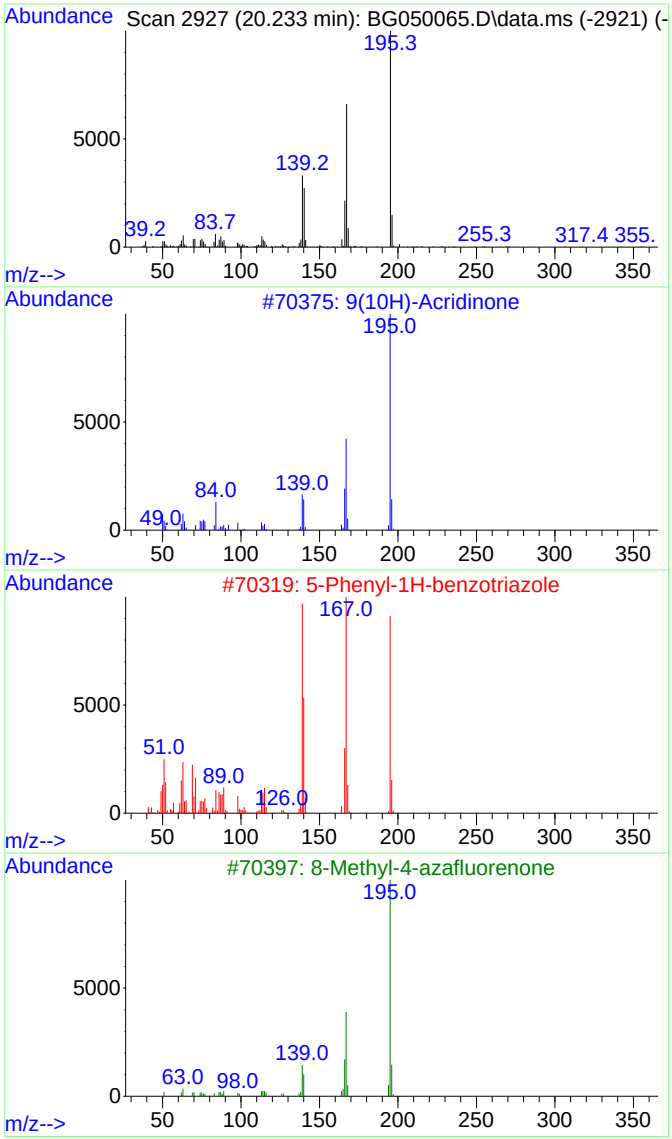
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

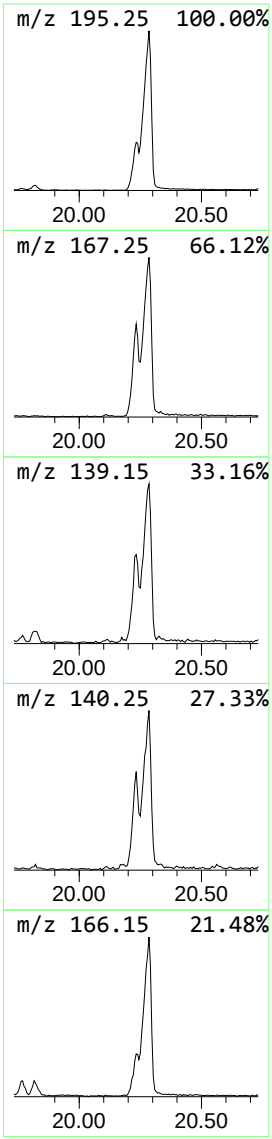
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 48 9(10H)-Acridinone Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.233	25.15 ng/ul	475152	Chrysene-d12	21.654

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9(10H)-Acridinone	195	C13H9NO	000578-95-0	93
2	5-Phenyl-1H-benzotriazole	195	C12H9N3	025877-73-0	91
3	8-Methyl-4-azafluorenone	195	C13H9NO	064292-03-1	90
4	3-Acridinol	195	C13H9NO	007132-70-9	90
5	7-methyl-4-azafluorenone	195	C13H9NO	064292-02-0	90





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Acq On : 31 Aug 2021 15:02
Operator : CG/JU
Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKG0

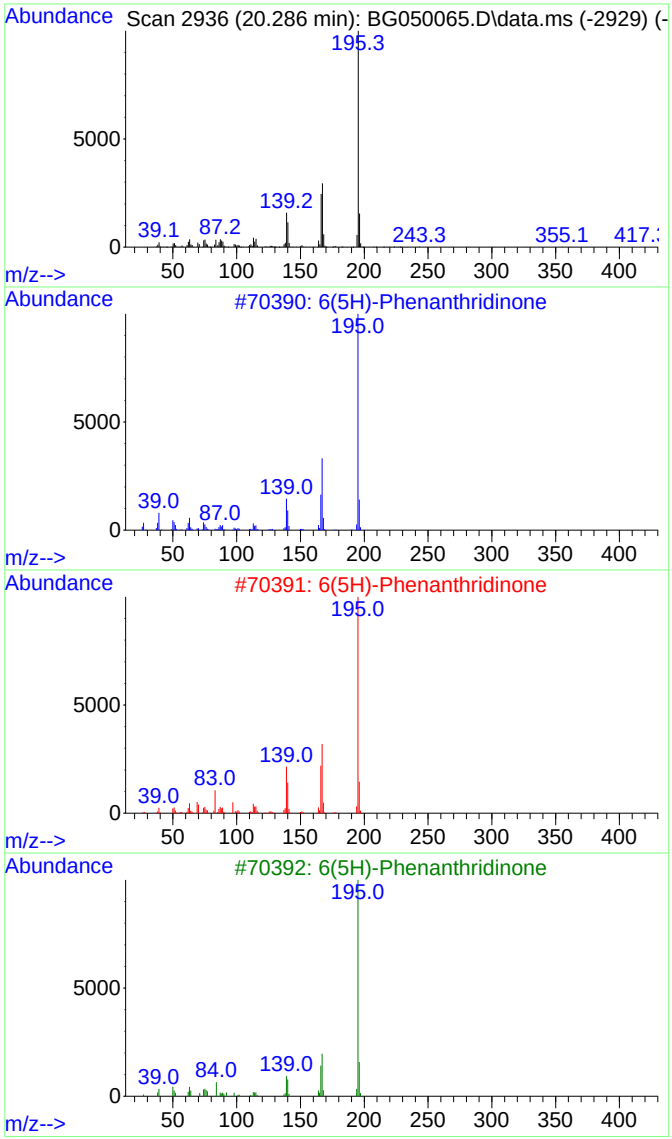
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

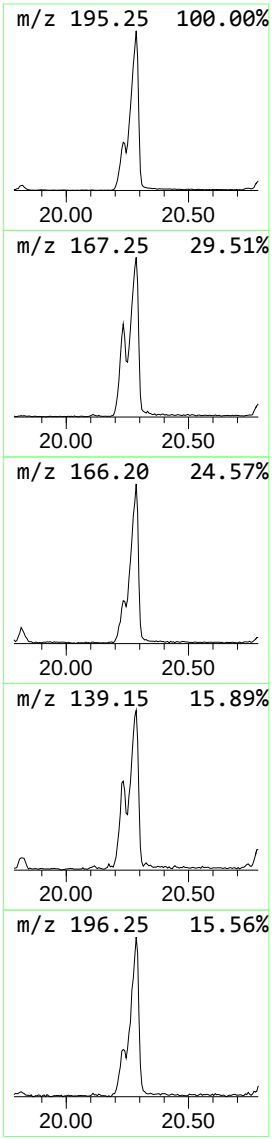
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 49 6(5H)-Phenanthridinone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.285	85.21 ng/ul	1609570	Chrysene-d12	21.654

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6(5H)-Phenanthridinone	195	C13H9NO	001015-89-0	96
2	6(5H)-Phenanthridinone	195	C13H9NO	001015-89-0	94
3	6(5H)-Phenanthridinone	195	C13H9NO	001015-89-0	94
4	3-Acridinol	195	C13H9NO	007132-70-9	93
5	9-Acridinol	195	C13H9NO	000643-62-9	93





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Data File : BG050065.D
Acq On : 31 Aug 2021 15:02
Operator : CG/JU
Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKG0

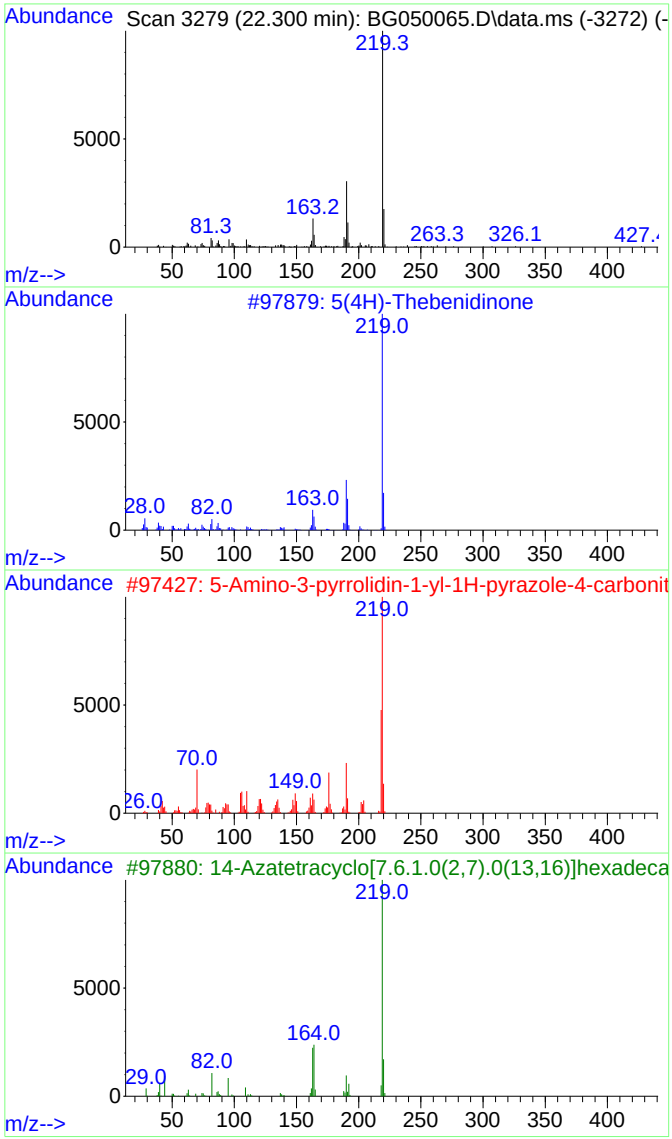
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

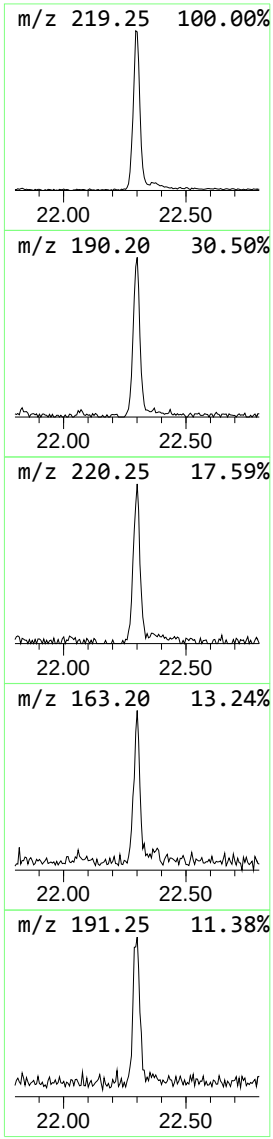
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 51 5(4H)-Thebenidinone Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.301	12.65 ng/ul	238903	Chrysene-d12	21.654

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5(4H)-Thebenidinone	219	C15H9NO	064884-40-8	87
2	5-Amino-3-pyrrolidin-1-yl-1H-pyr...	219	C11H17N5	1000499-90-1	64
3	14-Azatetracyclo[7.6.1.0(2,7).0(...	219	C15H9NO	042326-31-8	59
4	2-Phenyl-6-vinylindolizine	219	C16H13N	077652-49-4	53
5	2-Naphthalenamine, N-phenyl-	219	C16H13N	000135-88-6	52





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
Data File : BG050065.D
Acq On : 31 Aug 2021 15:02
Operator : CG/JU
Sample : M3464-07
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKG0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.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...	5.946	23.5	ng/ul	202720	1	7.961	172642	20.0
Mesitylene	7.603	37.0	ng/ul	319126	1	7.961	172642	20.0
Benzofuran	7.679	27.4	ng/ul	236138	1	7.961	172642	20.0
Indane	8.349	781.4	ng/ul	6745410	1	7.961	172642	20.0
3-Methylphenyla...	8.501	59.4	ng/ul	512882	1	7.961	172642	20.0
Benzofuran, 7-m...	9.324	20.8	ng/ul	179330	1	7.961	172642	20.0
Phenol, 3,5-dic...	13.659	24.7	ng/ul	686445	3	14.629	555894	20.0
Naphthalene, 1,...	13.800	29.4	ng/ul	817838	3	14.629	555894	20.0
Naphthalene, 1,...	13.941	26.0	ng/ul	722926	3	14.629	555894	20.0
2-Naphthyl isoc...	14.764	25.1	ng/ul	696665	3	14.629	555894	20.0
7-Methyl-1-naph...	15.821	31.2	ng/ul	866474	3	14.629	555894	20.0
5H-Indeno[1,2-b...	16.197	17.7	ng/ul	403840	4	17.384	457133	20.0
1(2H)-Acenaphth...	16.361	17.6	ng/ul	401224	4	17.384	457133	20.0
unknown-01	16.408	21.1	ng/ul	481744	4	17.384	457133	20.0
3-Hydroxybiphenyl	16.567	25.6	ng/ul	585300	4	17.384	457133	20.0
p-Hydroxybiphenyl	16.643	49.0	ng/ul	1119570	4	17.384	457133	20.0
Quinoline, 1-oxide	16.685	24.6	ng/ul	562570	4	17.384	457133	20.0
2(1H)-Quinolino...	16.972	27.1	ng/ul	620055	4	17.384	457133	20.0
2-Dibenzofuranol	17.166	34.2	ng/ul	782301	4	17.384	457133	20.0
8-Quinolinol, 3...	17.278	28.1	ng/ul	642062	4	17.384	457133	20.0
unknown-02	17.589	17.2	ng/ul	392732	4	17.384	457133	20.0
Dibenzo-p-dioxin	17.889	52.3	ng/ul	1195240	4	17.384	457133	20.0
9H-Fluoren-3-ol	18.230	22.8	ng/ul	521982	4	17.384	457133	20.0
9H-Carbazole, 9...	18.312	41.3	ng/ul	942916	4	17.384	457133	20.0
2-Hydroxyfluorene	18.388	31.8	ng/ul	726115	4	17.384	457133	20.0
Indeno[1,2-b]py...	18.553	35.5	ng/ul	812312	4	17.384	457133	20.0
9(10H)-Acridinone	20.233	25.1	ng/ul	475152	5	21.654	377789	20.0
6(5H)-Phenanthr...	20.285	85.2	ng/ul	1609570	5	21.654	377789	20.0
5(4H)-Thebenidi...	22.301	12.7	ng/ul	238903	5	21.654	377789	20.0