LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
Data File : BG050067.D
Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YAS53

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: BG050067.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.438	399	408	416	rBV	440062	757444	0.78%	0.381%
2	5.590	424	434	442	rBV	129440	265142	0.27%	0.133%
3	5.949	486	495	504	rBV	386319	651342	0.67%	0.328%
4	6.430	569	577	587	rVB	105664	176744	0.18%	0.089%
5	7.094	684	690	695	rVV	211054	357738	0.37%	0.180%
6	7.177	699	704	711	rVB	148383	266347	0.28%	0.134%
7	7.288	715	723	727	rBV	92688	165168	0.17%	0.083%
8	7.341	727	732	738	rVB2	131397	213516	0.22%	0.107%
9	7.500	752	759	765	rBV	100512	188282	0.19%	0.095%
10	7.605	770	777	785	rVV	464625	807854	0.84%	0.406%
11	7.682	785	790	798	rVB	69293	119687	0.12%	0.060%
12	7.964	831	838	849	rVV2	93364	200363	0.21%	0.101%
13	8.075	849	857	865	rVV	457285	871641	0.90%	0.438%
14	8.146	865	869	879	rVB	61018	118041	0.12%	0.059%
15	8.363	892	906	912	rBV	7674044	16639019	17.21%	8.368%
16	8.504	925	930	938	rVB	238928	400800	0.41%	0.202%
17	8.604	941	947	952	rBV2	81235	137799	0.14%	0.069%
18	8.692	952	962	971	rVB	95989	193254	0.20%	0.097%
19	9.074	1015	1027	1033	rBV	240521	423143	0.44%	0.213%
20	9.150	1033	1040	1047	rVV3	481072	902398	0.93%	0.454%
21	9.332	1062	1071	1078	rVB	578667	1049111	1.09%	0.528%
22	9.462	1078	1093	1098	rBV	1332944	3094890	3.20%	1.557%
23	9.520	1098	1103	1111	rVV	474114	828287	0.86%	0.417%
24	9.603	1111	1117	1122	rVV	133477	230957	0.24%	0.116%
25	9.661	1122	1127	1136	rVB	150659	257716	0.27%	0.130%
26	9.873	1157	1163	1174	rVB3	101448	257287	0.27%	0.129%
27	10.026	1182	1189	1198	rVV	585882	1057543	1.09%	0.532%
28	10.184	1207	1216	1227	rBV3	951318	2840155	2.94%	1.428%
29	10.290	1227	1234	1250	rVB	389366	1259991	1.30%	0.634%
30	10.431	1250	1258	1275	rVB7	112196	450089	0.47%	0.226%
31	10.971	1312	1350	1353	rBV	17712677	96679523	100.00%	48.624%
32	11.030	1353	1360	1367	rVV	4688332	7620714	7.88%	3.833%
33	11.095	1367	1371	1377	rVB2	82165	142543	0.15%	0.072%
34	11.212	1385	1391	1397	rVB	182873	366055	0.38%	0.184%
35	11.447	1424	1431	1439	rVB2	78072	140465	0.15%	0.071%
36	11.706	1469	1475	1485	rBV2	65859	135122	0.14%	0.068%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
Data File : BG050067.D
Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YAS53

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	11.900	1499	1508	1512	rBV3	93841	185845	0.19%	0.093%
38	12.170	1549	1554	1564	rVB3	46186	97464	0.10%	0.049%
39	12.346	1578	1584	1594	rVV	221802	408376	0.42%	0.205%
40	12.481	1594	1607	1612	rVV	6708305	14532538	15.03%	7.309%
41	12.575	1616	1623	1629	rVV	251169	473531	0.49%	0.238%
42	12.687	1629	1642	1648	rVV	4036936	7502799	7.76%	3.773%
43	13.163	1717	1723	1728	rBV	62777	97151	0.10%	0.049%
44	13.456	1766	1773	1782	rVV	1077672	1651281	1.71%	0.830%
45	13.650	1800	1806	1816	rVB	271012	548157	0.57%	0.276%
46	13.803	1818	1832	1839	rBV3	310732	860170	0.89%	0.433%
47	13.944	1849	1856	1861	rVV	438921	852812	0.88%	0.429%
48	14.003	1861	1866	1868	rVV	353549	620753	0.64%	0.312%
49	14.026	1868	1870	1882	rVB	349386	416710	0.43%	0.210%
50	14.185	1890	1897	1900	rBV	181159	307877	0.32%	0.155%
51	14.214	1900	1902	1908	rVB	87728	99952	0.10%	0.050%
52	14.320	1913	1920	1930	rVB2	368730	826366	0.85%	0.416%
53	14.596	1957	1967	1969	rVV2	133848	236987	0.25%	0.119%
54	14.625	1969	1972	1977	rVV	196300	326093	0.34%	0.164%
55	14.702	1977	1985	1990	rVV	3728088	6135278	6.35%	3.086%
56	14.766	1990	1996	2001	rVV	522692	842257	0.87%	0.424%
57	14.819	2001	2005	2009	rVV2	85786	149854	0.16%	0.075%
58	14.937	2020	2025	2029	rVV	84404	149643	0.15%	0.075%
59	15.031	2033	2041	2049	rVV	1840272	2971359	3.07%	1.494%
60	15.618	2136	2141	2146	rVV	446744	732145	0.76%	0.368%
61	15.683	2146	2152	2158	rVB	1696838	2579267	2.67%	1.297%
62	15.759	2159	2165	2169	rBV3	149421	268898	0.28%	0.135%
63	15.800	2169	2172	2175	rVV2	117845	197347	0.20%	0.099%
64	15.835	2175	2178	2183	rVV3	144213	236536	0.24%	0.119%
65	15.924	2190	2193	2197	rVV	77107	119226	0.12%	0.060%
66	15.971	2197	2201	2205	rVB	86998	127684	0.13%	0.064%
67	16.111	2220	2225	2233	rVV3	100671	237826	0.25%	0.120%
68	16.188	2233	2238	2245	rVB2	51373	99685	0.10%	0.050%
69	16.464	2279	2285	2291	rBV	105506	163736	0.17%	0.082%
70	16.681	2318	2322	2326	rVV3	57702	99370	0.10%	0.050%
71	16.734	2326	2331	2337	rVB3	58547	111526	0.12%	0.056%
72	16.834	2343	2348	2363	rVB5	30451	110647	0.11%	0.056%
73	17.192	2405	2409	2416	rVV	128270	211723	0.22%	0.106%
74	17.380	2436	2441	2444	rVV	253552	444212	0.46%	0.223%
75	17.427	2444	2449	2454	rVV	1345204	2132318	2.21%	1.072%
76	17.480	2454	2458	2462	rVV2	564750	891588	0.92%	0.448%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
Data File : BG050067.D
Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YAS53

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

77	17.515	2462	2464	2469	rVV2	147268	182701	0.19%	0.092%
78	17.574	2469	2474	2483	rVB	226660	364247	0.38%	0.183%
79	17.797	2500	2512	2518	rVV	2466571	4258099	4.40%	2.142%
80	17.897	2518	2529	2534	rVV3	148389	314566	0.33%	0.158%
81	17.944	2534	2537	2544	rVB	78526	110739	0.11%	0.056%
82	18.032	2544	2552	2557	rBV2	69754	139322	0.14%	0.070%
83	18.220	2579	2584	2593	rVB2	79030	140108	0.14%	0.070%
84	18.303	2593	2598	2607	rBV	362989	557481	0.58%	0.280%
85	18.450	2618	2623	2629	rVB4	61048	116460	0.12%	0.059%
86	18.561	2634	2642	2646	rBV2	144000	313633	0.32%	0.158%
87	18.602	2646	2649	2653	rVV	131122	182754	0.19%	0.092%
88	18.696	2659	2665	2671	rVV2	127312	218118	0.23%	0.110%
89	19.771	2842	2848	2853	rBV	593074	883735	0.91%	0.444%
90	20.171	2911	2916	2923	rBV	94166	149007	0.15%	0.075%
91	21.651	3162	3168	3178	rVB	232642	381368	0.39%	0.192%
92	24.659	3671	3680	3692	rVB2	286785	791905	0.82%	0.398%
93	24.882	3706	3718	3728	rBV2	141459	404507	0.42%	0.203%

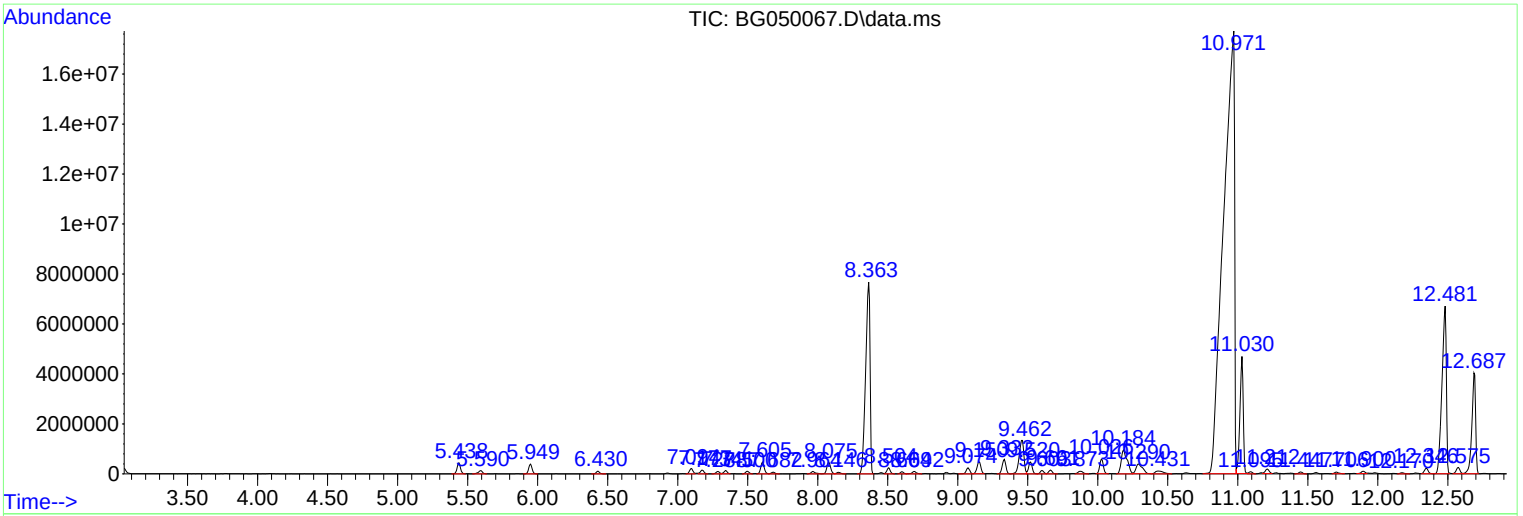
Sum of corrected areas: 198831907

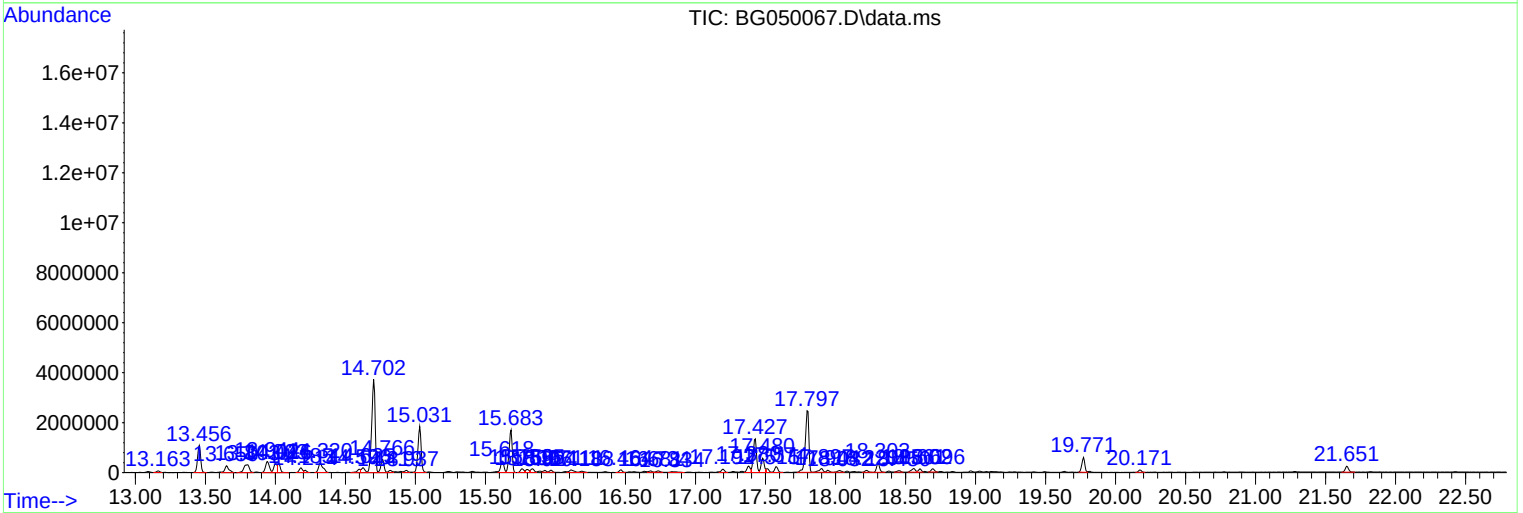
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
Data File : BG050067.D
Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

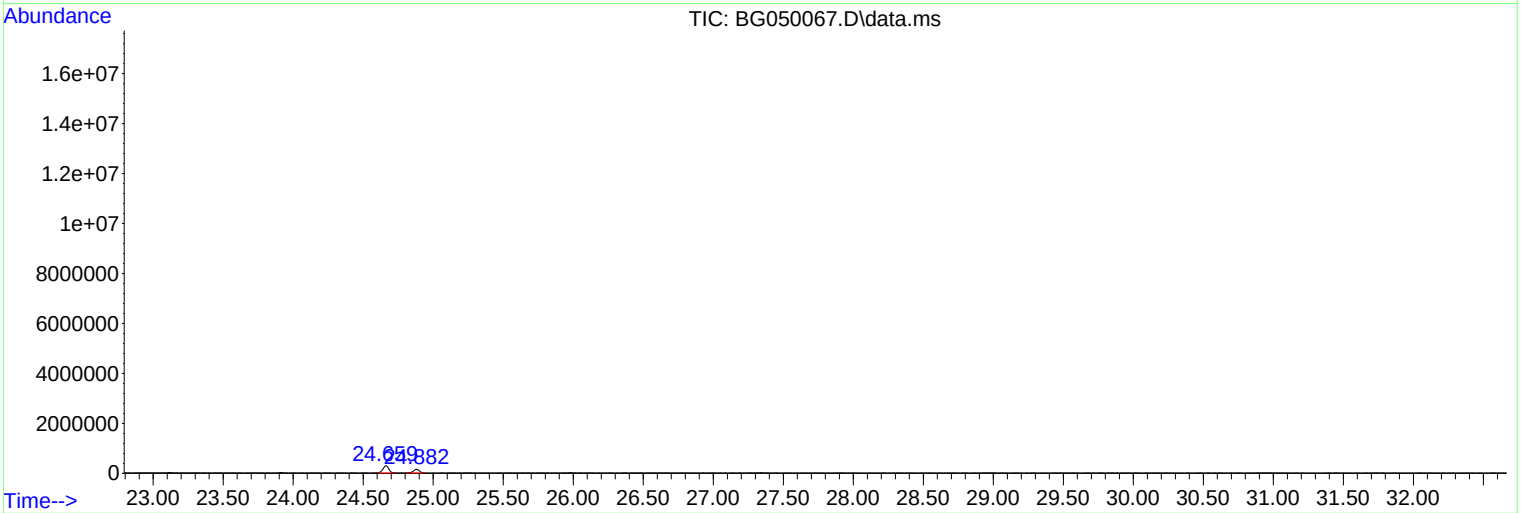
Instrument :
BNA_G
ClientSampleId :
YAS53

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 : BG050067.D
Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YAS53

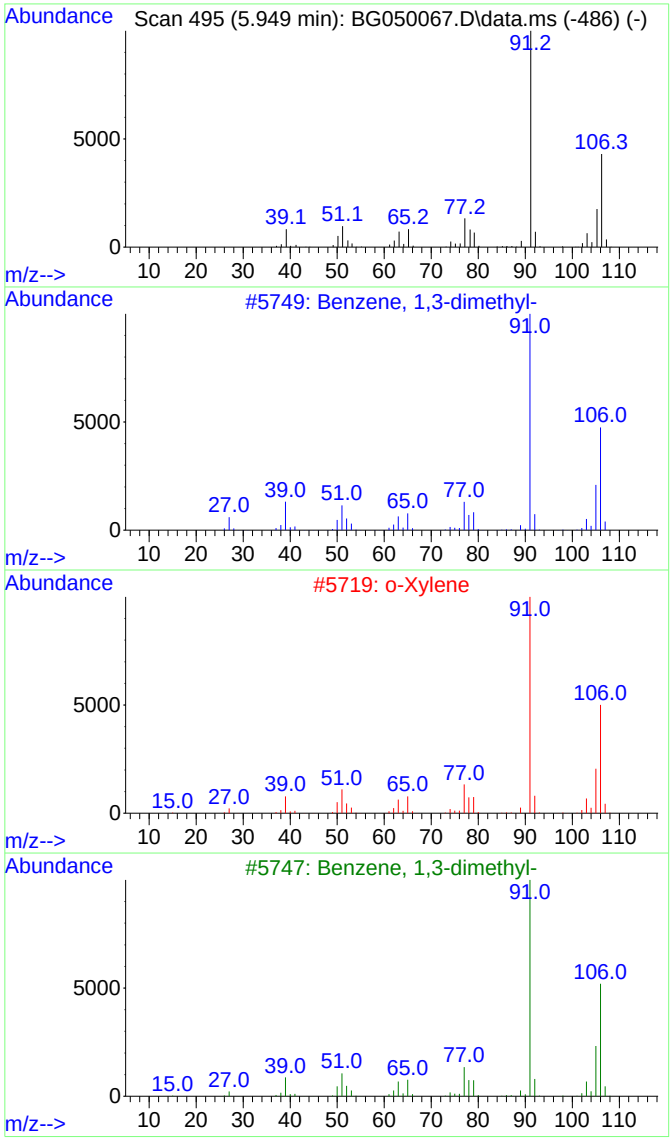
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

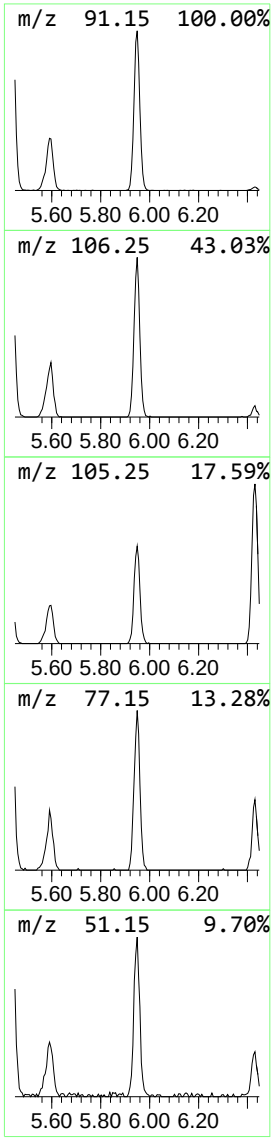
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.949	65.02 ng/ul	651342	1,4-Dichlorobenzene-d4	7.964

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2	o-Xylene	106	C8H10	000095-47-6	97
3	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
4	o-Xylene	106	C8H10	000095-47-6	95
5	o-Xylene	106	C8H10	000095-47-6	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
Data File : BG050067.D
Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YAS53

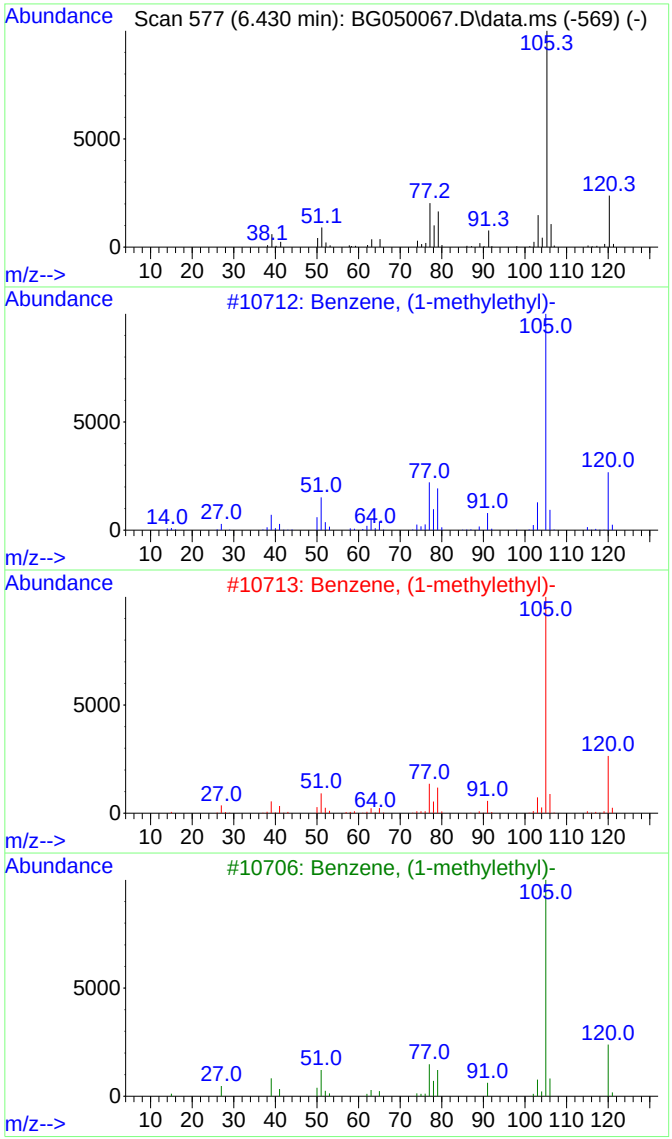
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

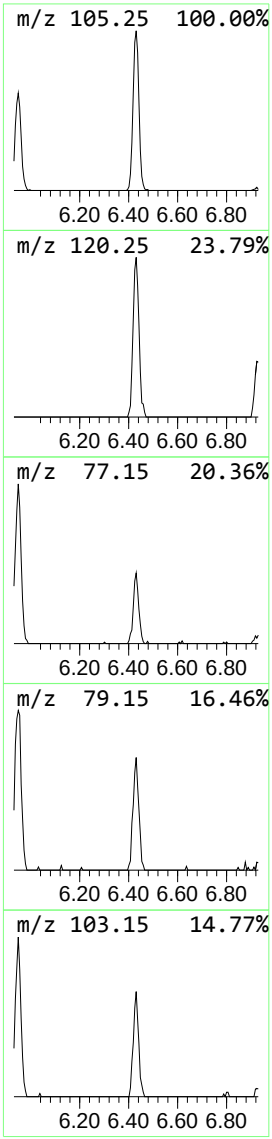
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, (1-methylethyl)- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.430	17.64 ng/ul	176744	1,4-Dichlorobenzene-d4	7.964

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	87
2	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	87
3	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	87
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	80
5	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
Data File : BG050067.D
Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YAS53

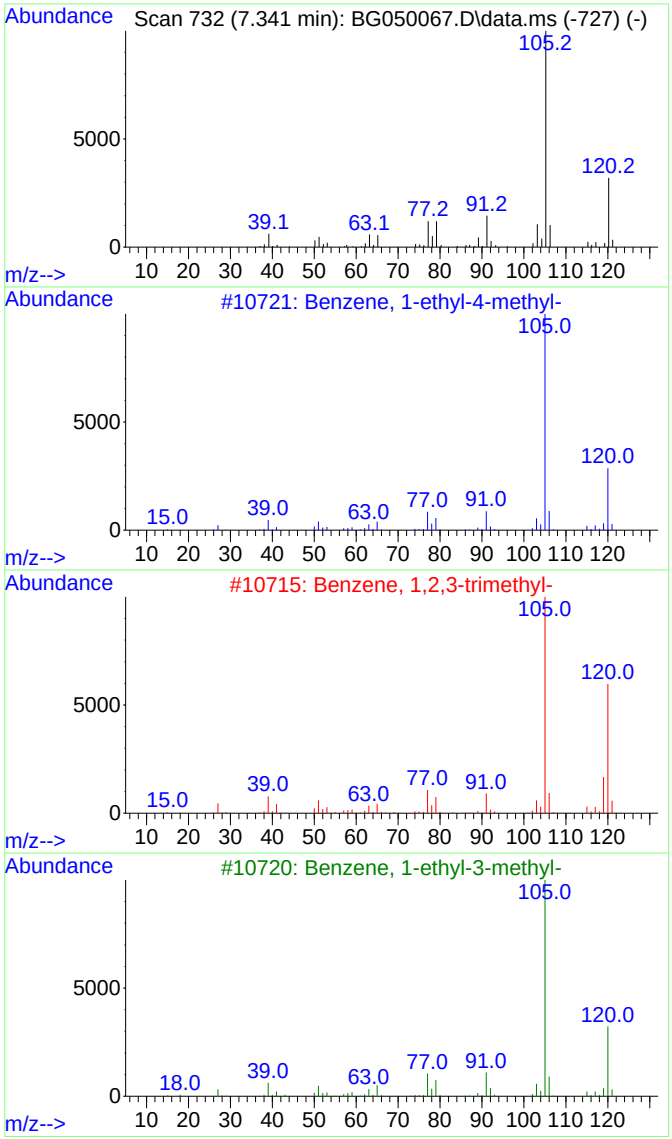
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

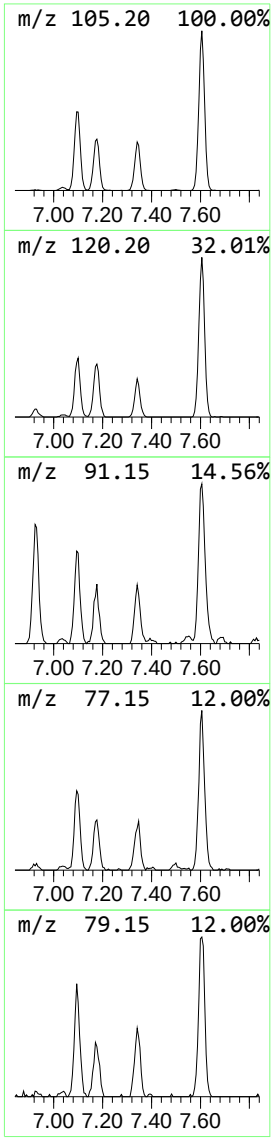
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1-ethyl-4-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.341	21.31 ng/ul	213516	1,4-Dichlorobenzene-d4	7.964

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
Data File : BG050067.D
Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YAS53

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

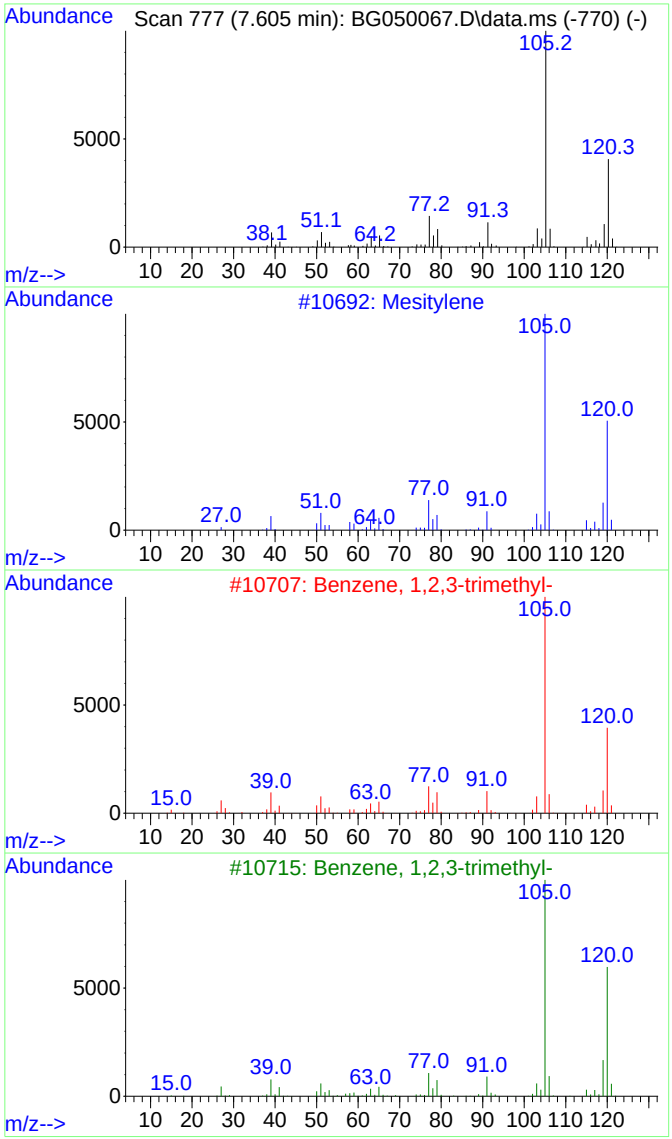
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

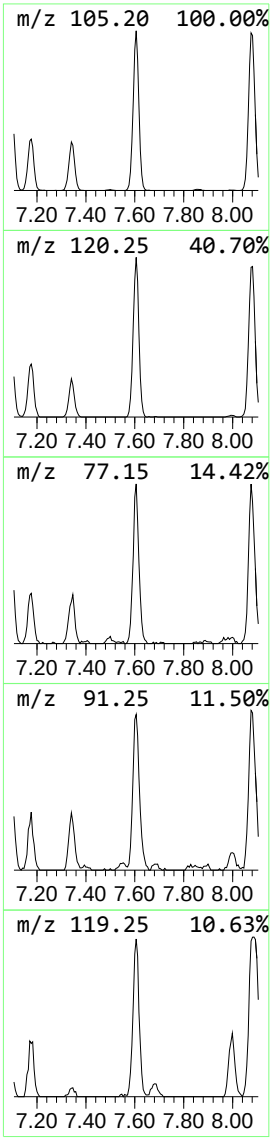
****

Peak Number 6 Mesitylene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.605	80.64 ng/ul	807854	1,4-Dichlorobenzene-d4	7.964

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	95
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	Mesitylene	120	C9H12	000108-67-8	91
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
Data File : BG050067.D
Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YAS53

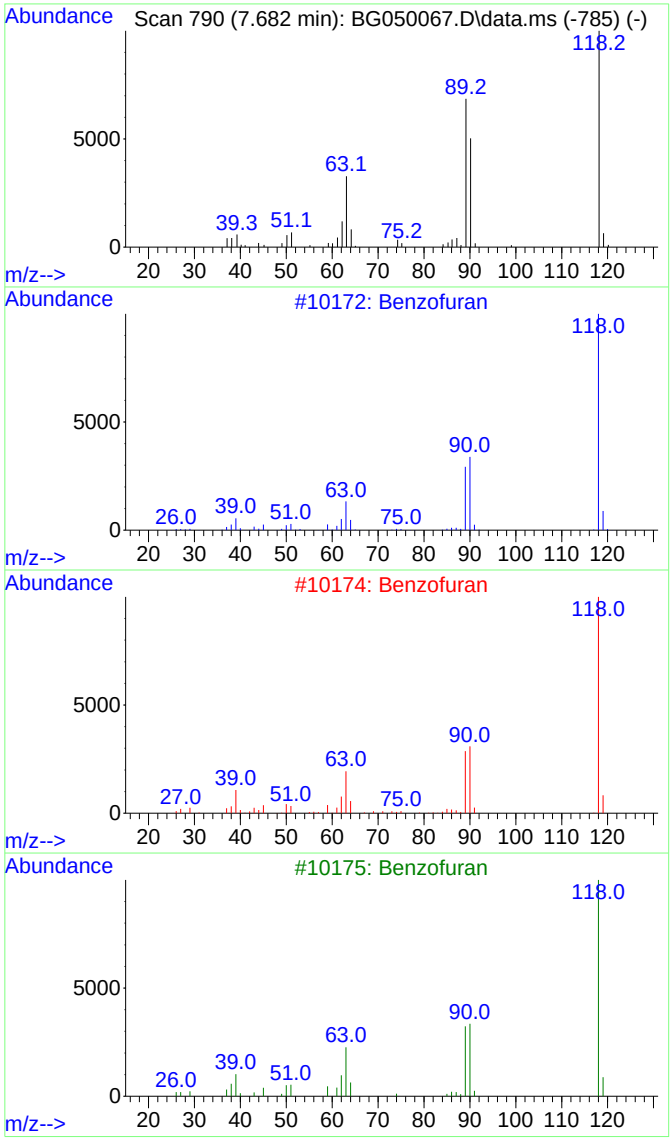
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

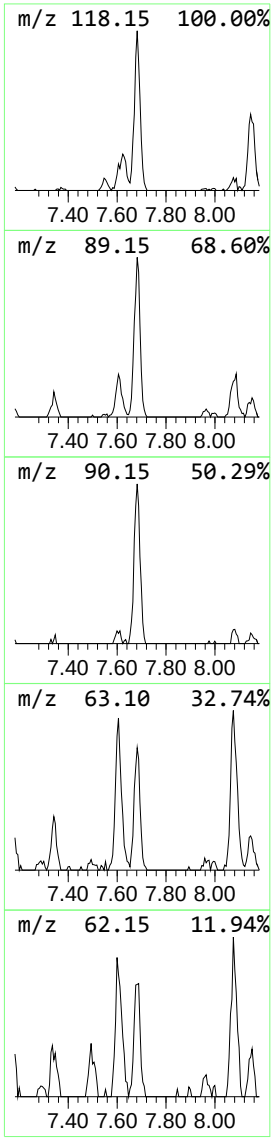
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzofuran Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.682	11.95 ng/ul	119687	1,4-Dichlorobenzene-d4	7.964

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	87
4	Benzocyclobuten-1(2H)-one	118	C8H6O	003469-06-5	59
5	2H-Cyclopenta[d]pyridazine	118	C7H6N2	000270-64-4	53





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Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YAS53

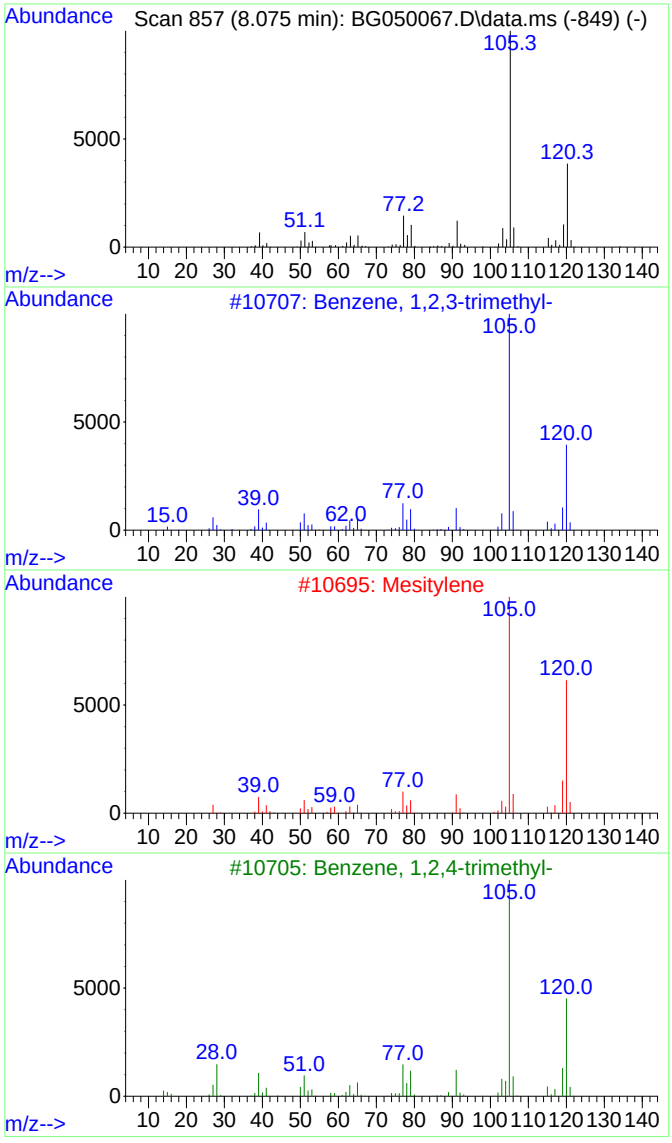
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

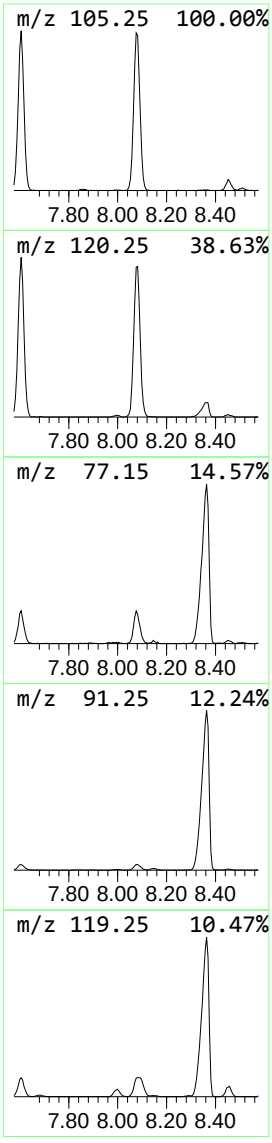
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.075	87.01 ng/ul	871641	1,4-Dichlorobenzene-d4	7.964

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





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Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YAS53

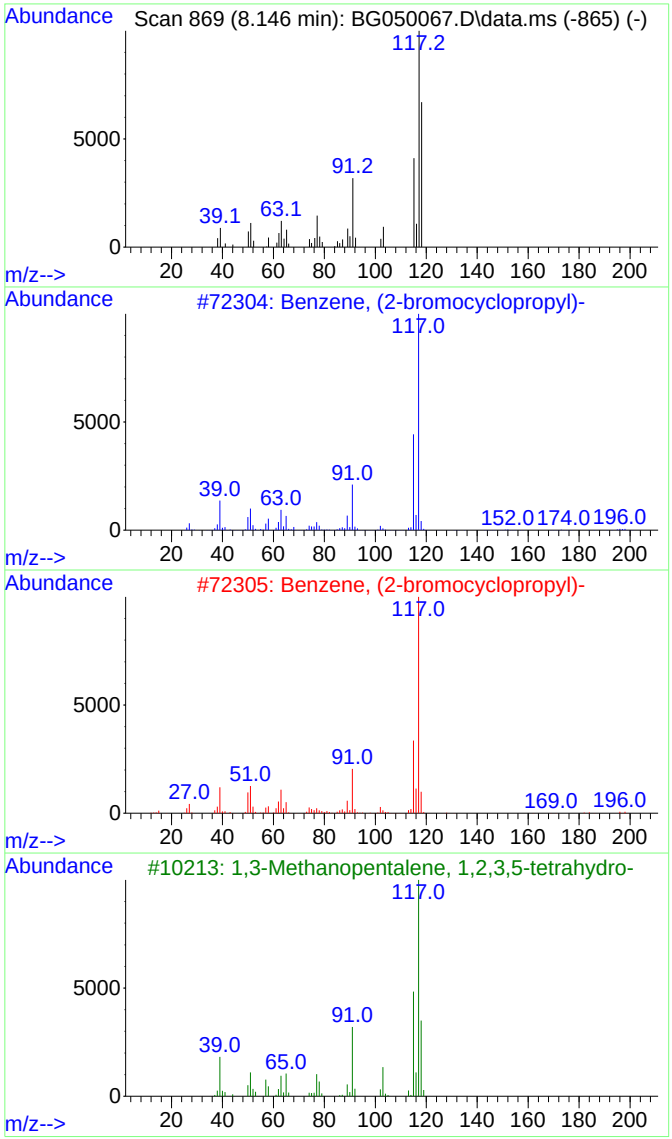
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

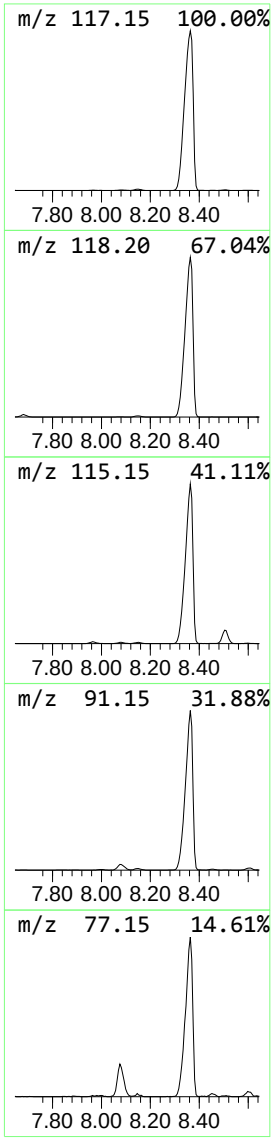
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, (2-bromocyclopropyl)- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.146	11.78 ng/ul	118041	1,4-Dichlorobenzene-d4	7.964

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	59
2	Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	59
3	1,3-Methanopentalene, 1,2,3,5-te...	118	C9H10	128600-88-4	50
4	Deltacyclene	118	C9H10	007785-10-6	50
5	trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	42





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Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Title : SVOA CALIBRATION

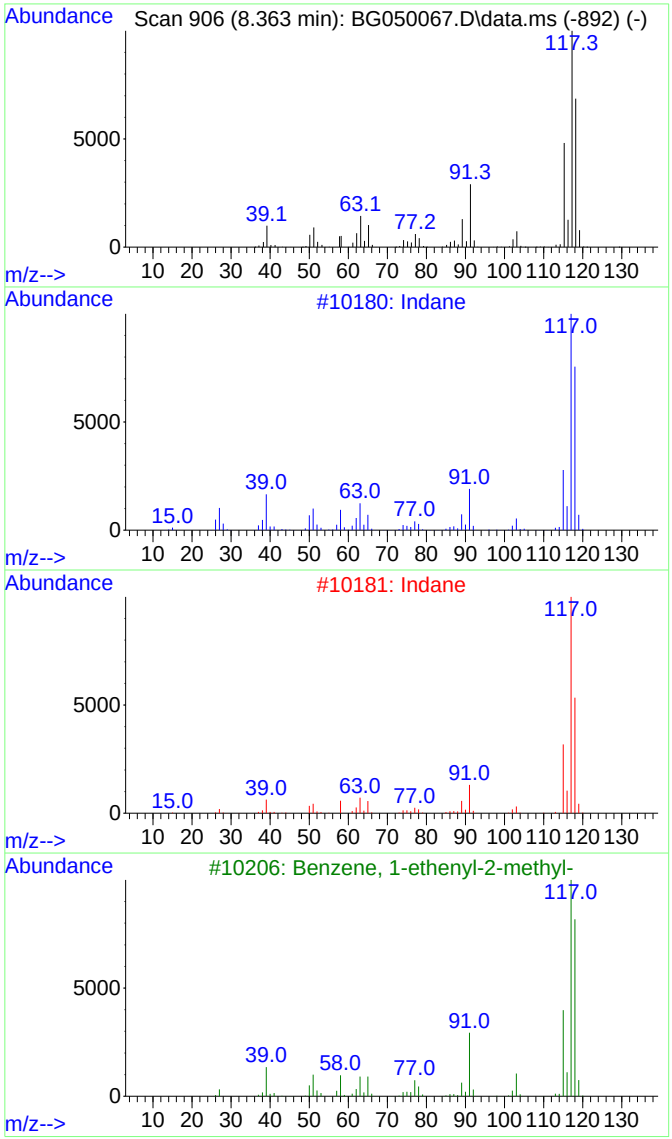
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

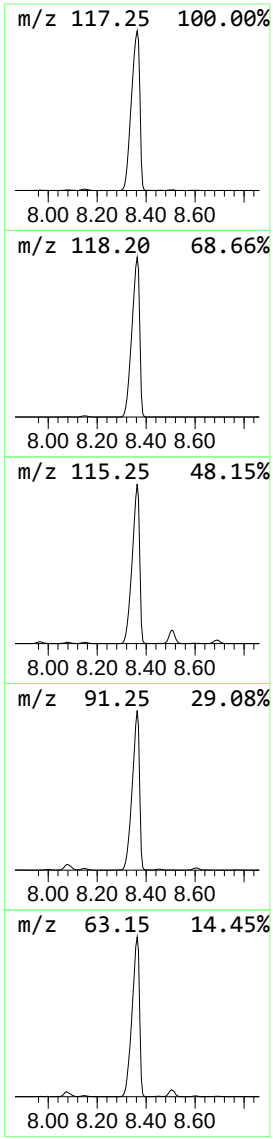
****

Peak Number 10 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.363	1660.89 ng/ul	16639000	1,4-Dichlorobenzene-d4	7.964

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	95
2	Indane	118	C9H10	000496-11-7	94
3	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	94
4	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	93
5	Indane	118	C9H10	000496-11-7	93





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Operator : CG/JU
Sample : M3494-01
Misc :
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Instrument : BNA_G
ClientSampleId : YAS53

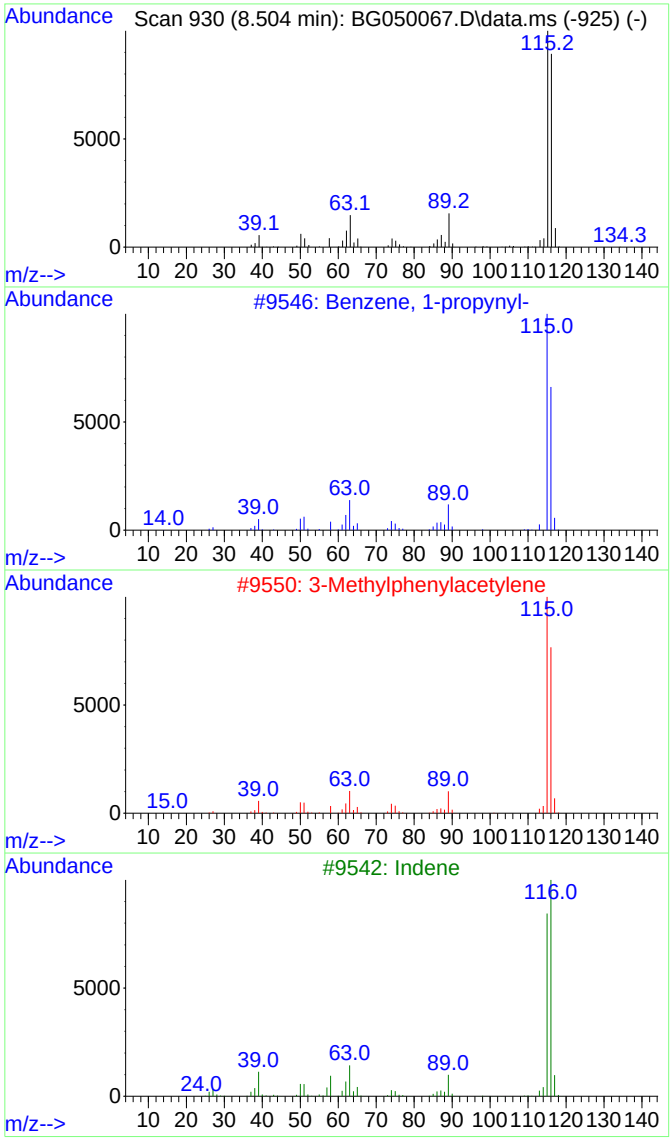
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

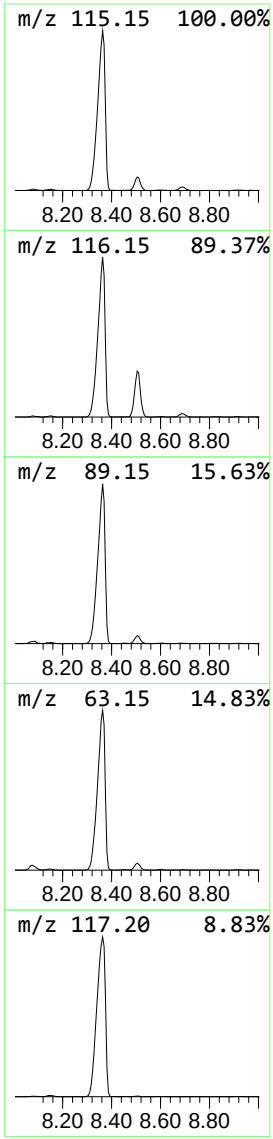
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, 1-propynyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.504	40.01 ng/ul	400800	1,4-Dichlorobenzene-d4	7.964

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-propynyl-	116	C9H8	000673-32-5	97
2	3-Methylphenylacetylene	116	C9H8	000766-82-5	94
3	Indene	116	C9H8	000095-13-6	94
4	Indene	116	C9H8	000095-13-6	91
5	2-Methylphenylacetylene	116	C9H8	000766-47-2	91





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Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

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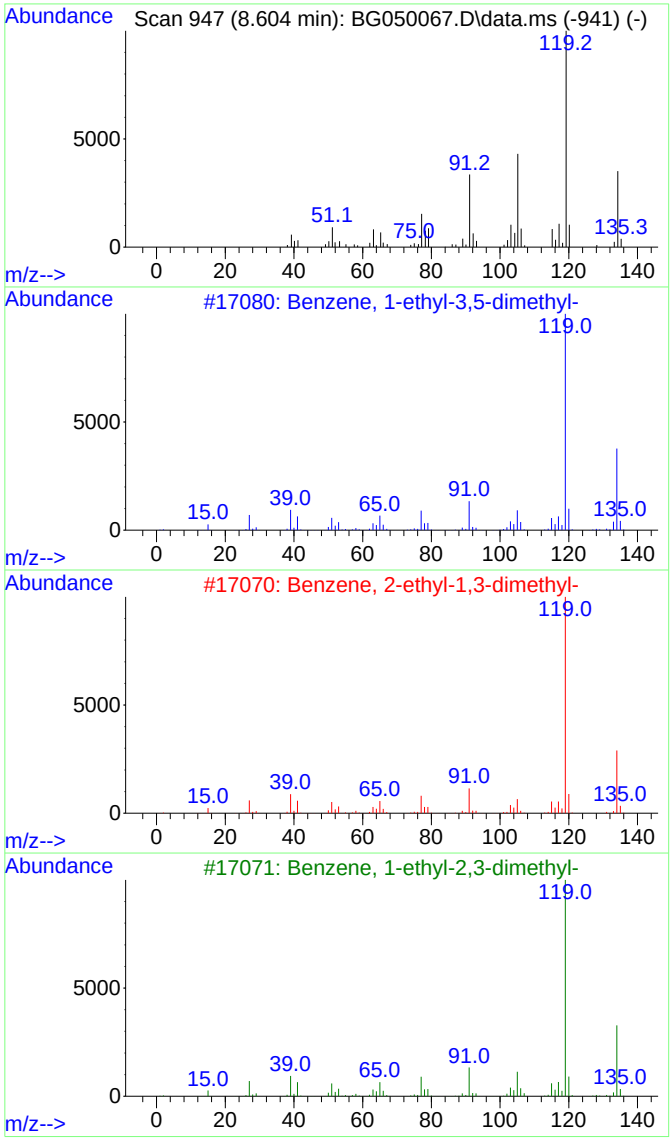
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

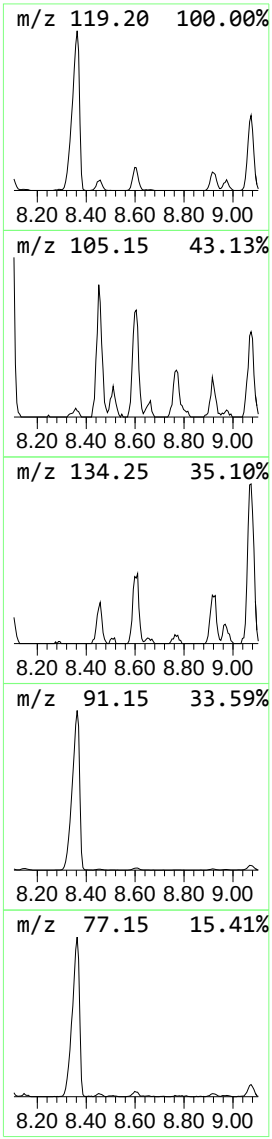
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.604	13.75 ng/ul	137799	1,4-Dichlorobenzene-d4	7.964

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	95
2	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
4	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	93





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Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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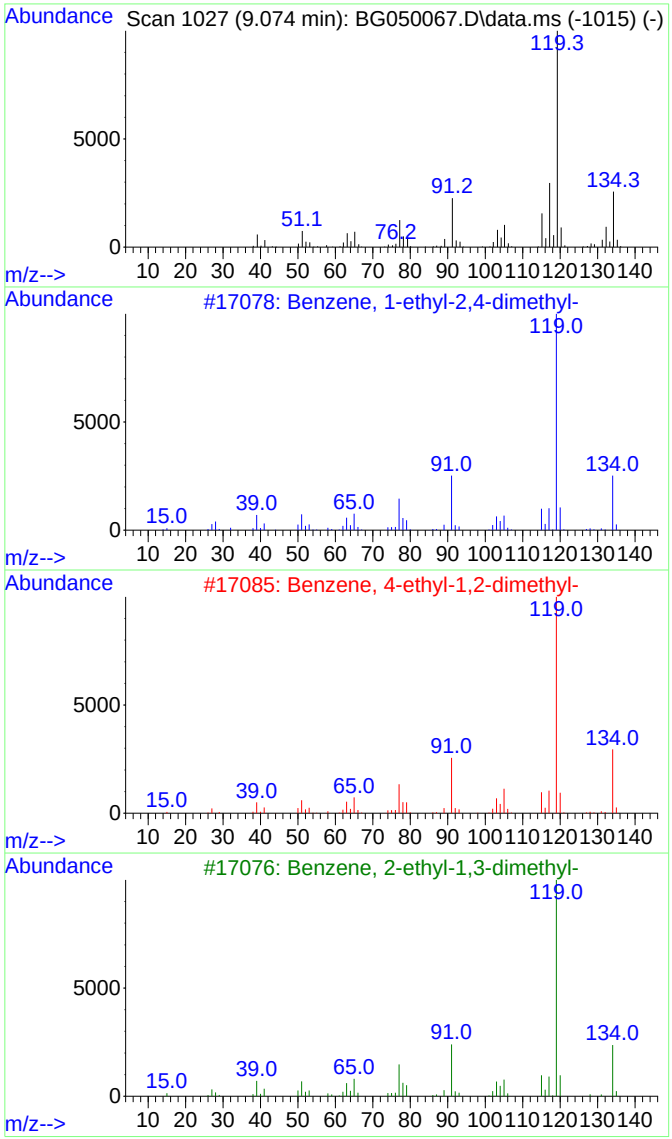
Quant Title : SVOA CALIBRATION

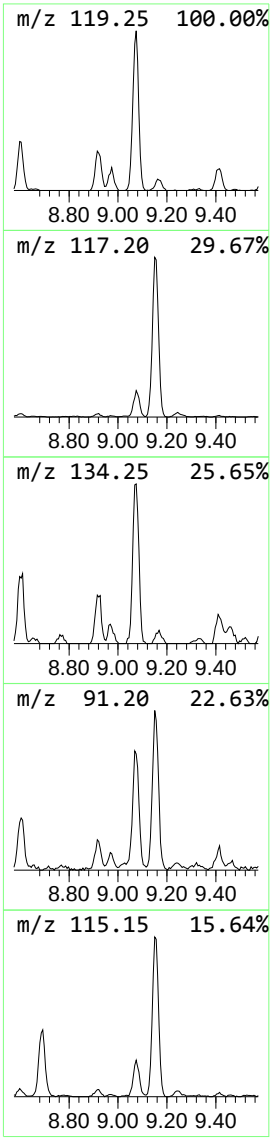
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.074	42.24 ng/ul	423143	1,4-Dichlorobenzene-d4	7.964

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	81





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Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YAS53

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

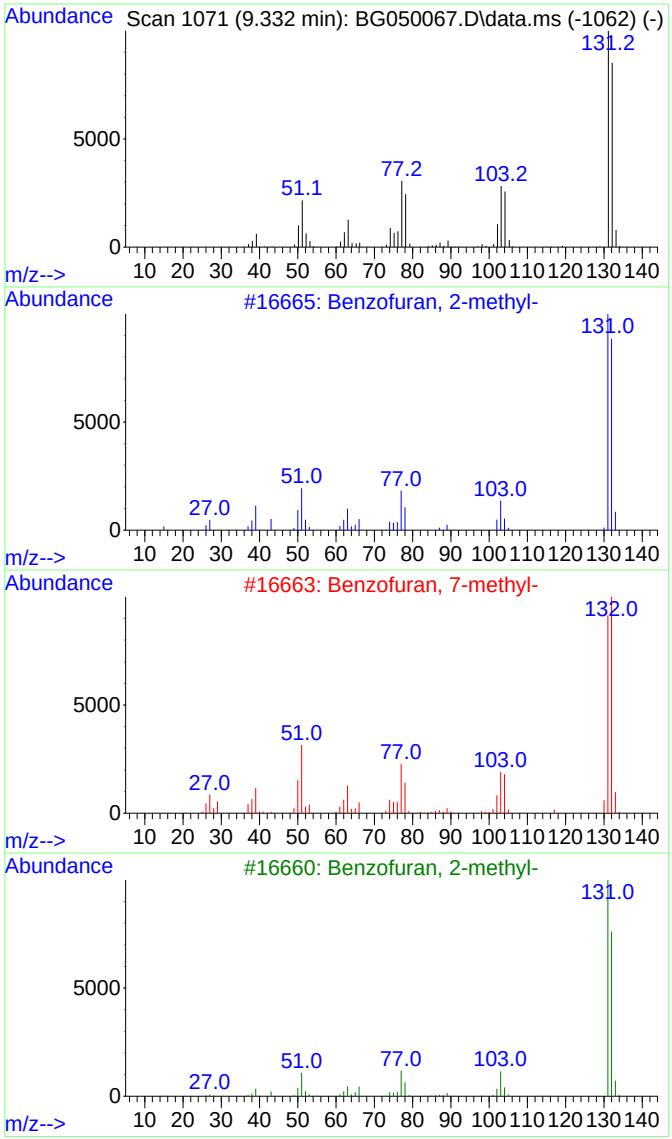
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

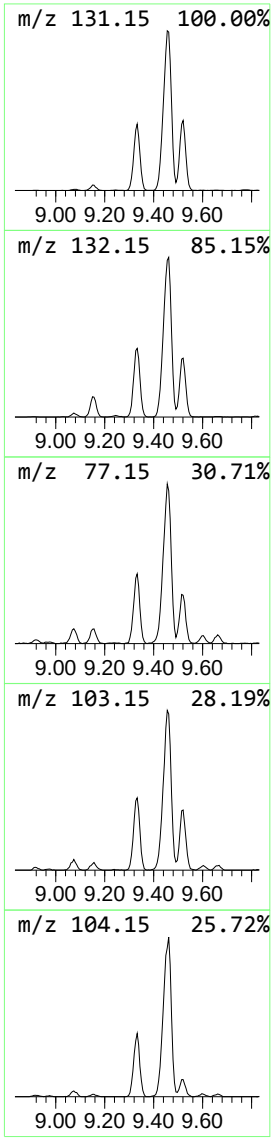
****

Peak Number 14 Benzofuran, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.332	104.72 ng/ul	1049110	1,4-Dichlorobenzene-d4	7.964

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
2	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	94
3	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
4	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
5	Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	93





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Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YAS53

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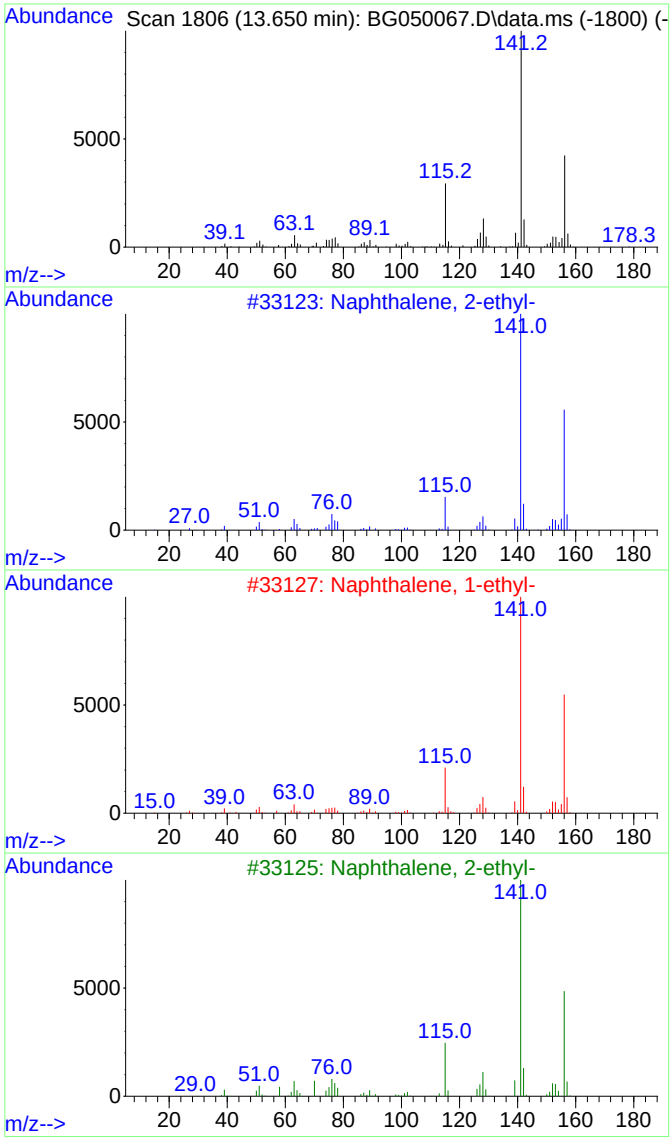
Quant Title : SVOA CALIBRATION

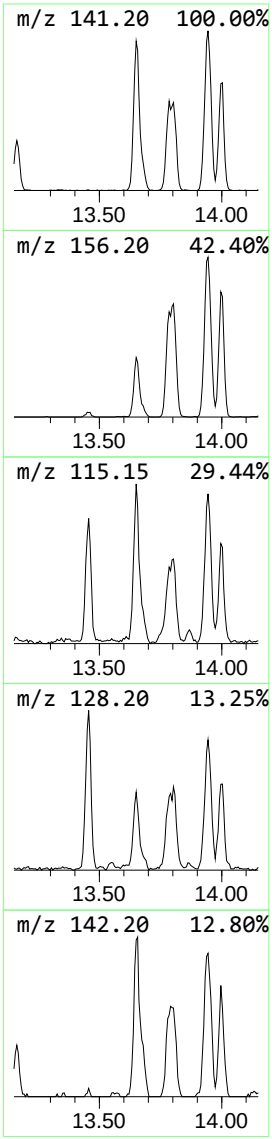
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Naphthalene, 2-ethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.650	33.62 ng/ul	548157	Acenaphthene-d10	14.625

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	94
2	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	94
3	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	93
4	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	93
5	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	91





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Sample : M3494-01
Misc :
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Instrument : BNA_G
ClientSampleId : YAS53

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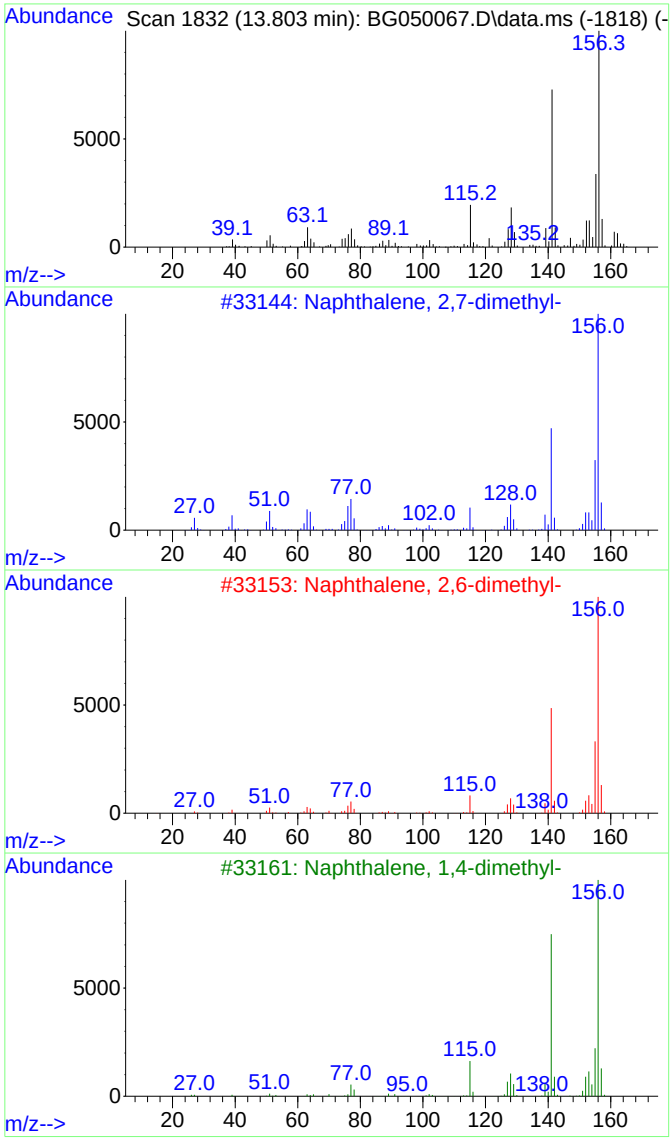
Quant Title : SVOA CALIBRATION

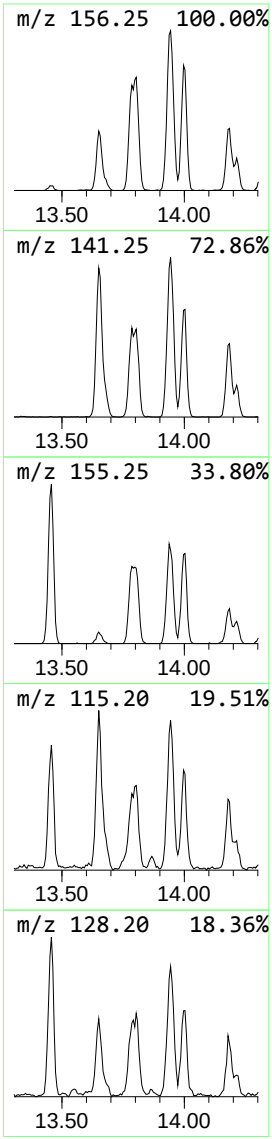
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Naphthalene, 2,7-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.803	52.76 ng/ul	860170	Acenaphthene-d10	14.625

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
2	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
3	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95





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Sample : M3494-01
Misc :
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Instrument : BNA_G
ClientSampleId : YAS53

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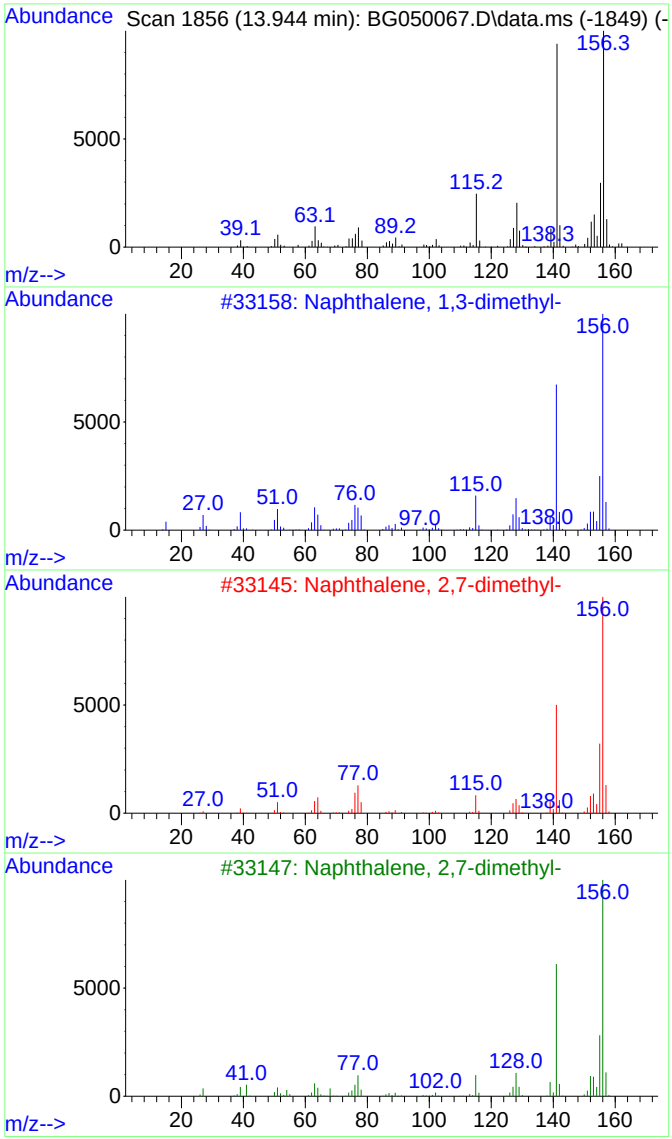
Quant Title : SVOA CALIBRATION

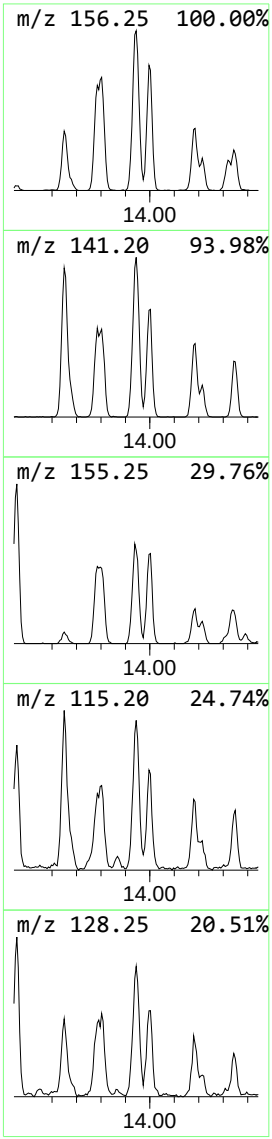
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Naphthalene, 1,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.944	52.30 ng/ul	852812	Acenaphthene-d10	14.625

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
2	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
4	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95
5	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95





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ClientSampleId :
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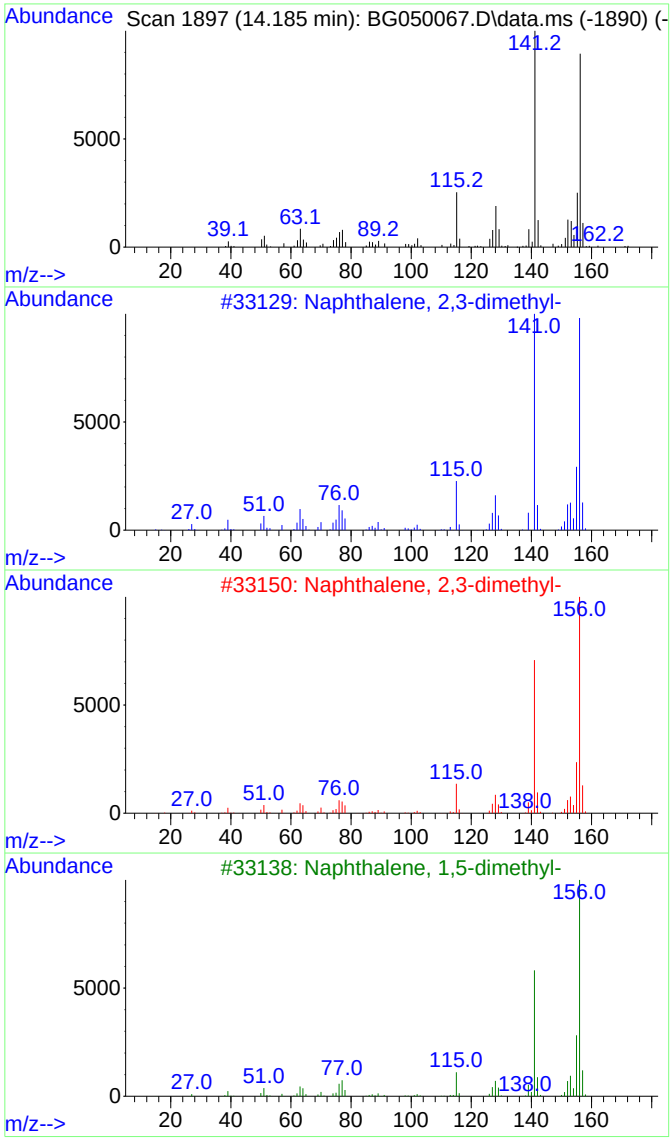
Quant Title : SVOA CALIBRATION

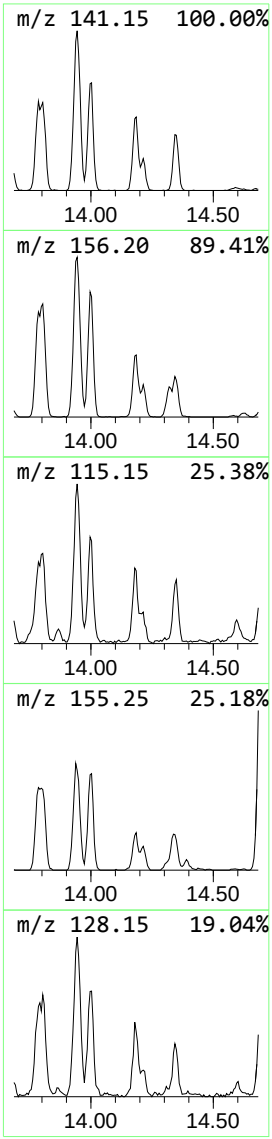
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Naphthalene, 2,3-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.185	18.88 ng/ul	307877	Acenaphthene-d10	14.625

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	96
4	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96





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Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

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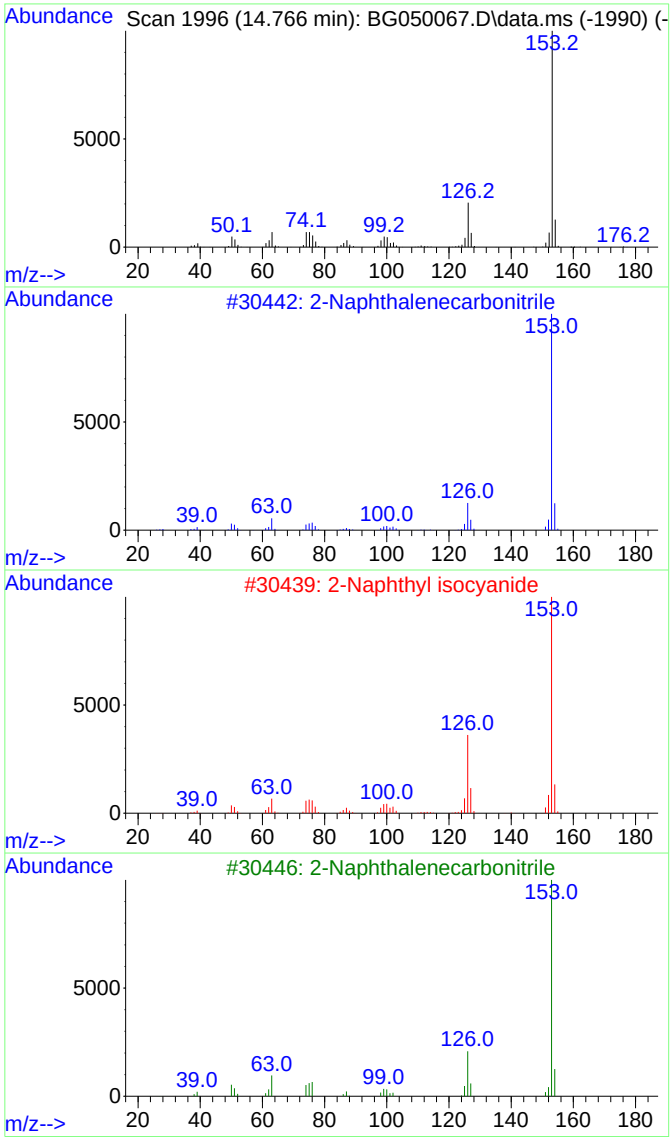
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

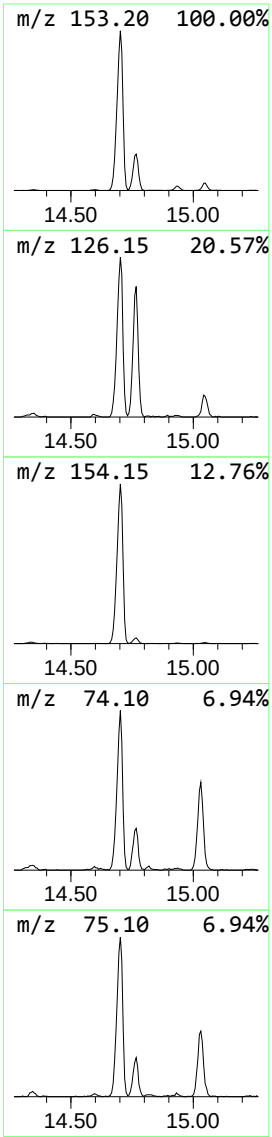
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 2-Naphthalenecarbonitrile Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.766	51.66 ng/ul	842257	Acenaphthene-d10	14.625

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





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Data File : BG050067.D
Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YAS53

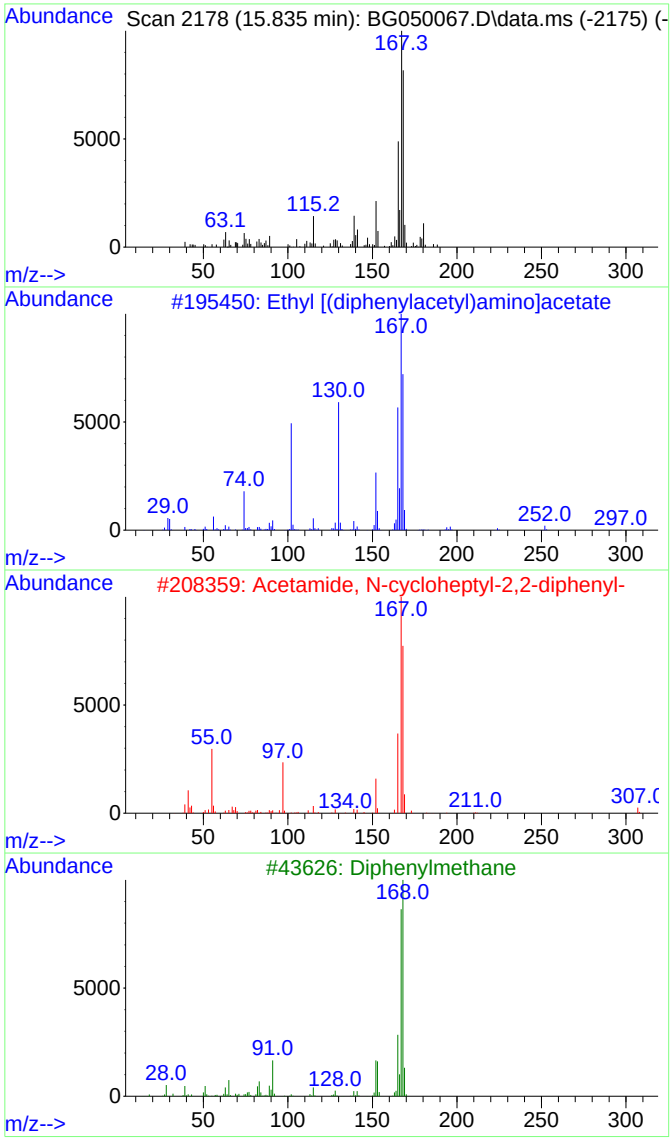
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

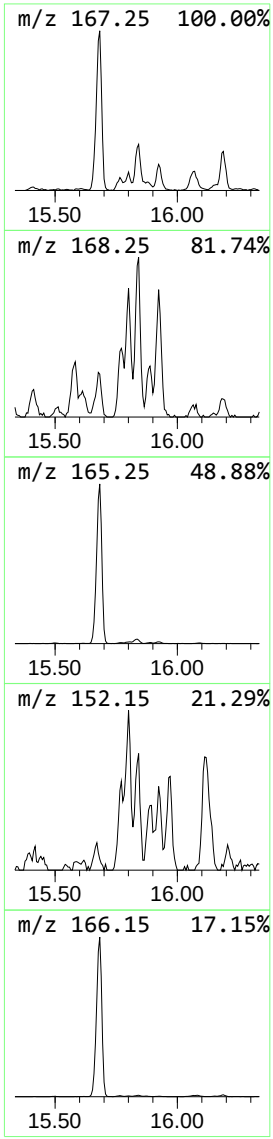
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Ethyl [(diphenylacetyl)amin... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.835	14.51 ng/ul	236536	Acenaphthene-d10	14.625

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethyl [(diphenylacetyl)amino]ace...	297	C18H19NO3	006325-31-1	72
2	Acetamide, N-cycloheptyl-2,2-dip...	307	C21H25NO	351062-01-6	72
3	Diphenylmethane	168	C13H12	000101-81-5	58
4	Fluorene, 1,4-dihydro-	168	C13H12	041593-21-9	53
5	Fluorene, 2,4a-dihydro-	168	C13H12	059247-36-8	50





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Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YAS53

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

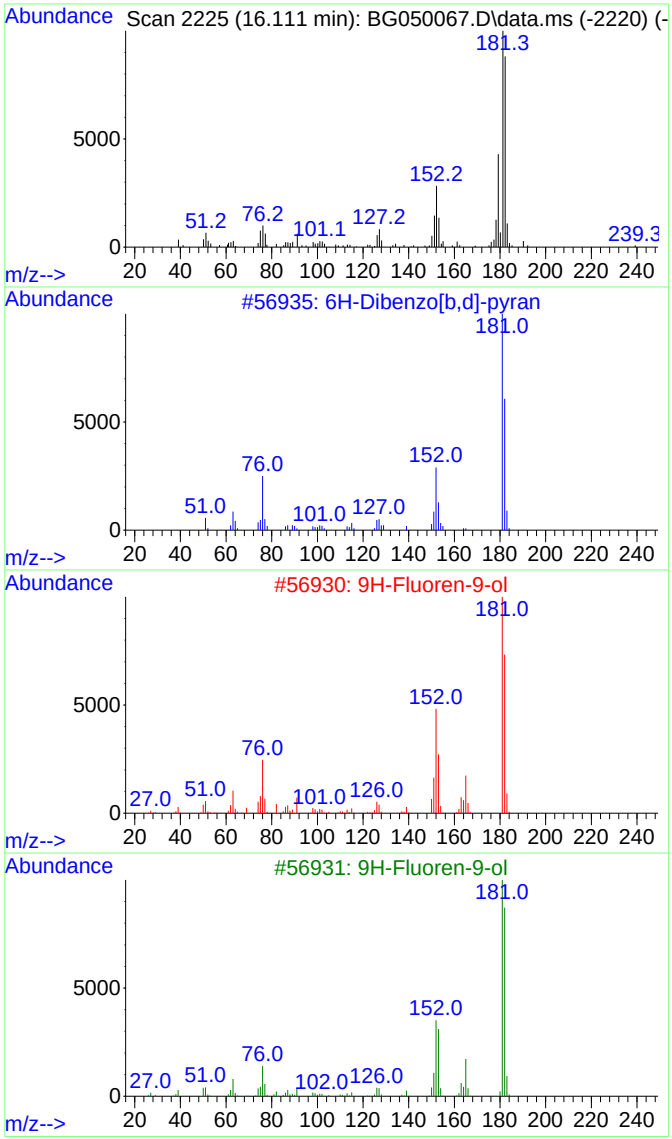
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

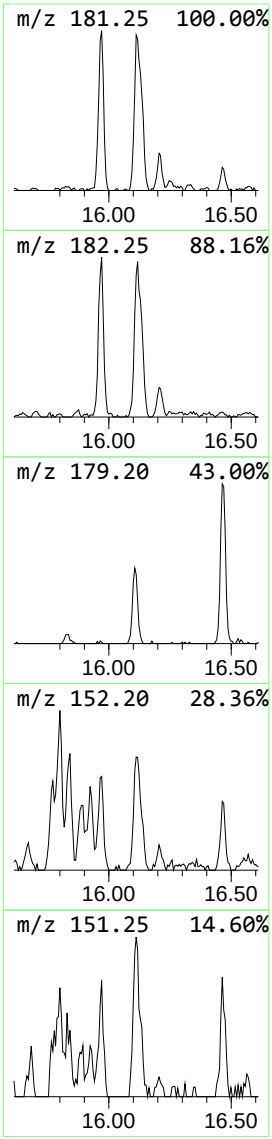
****

Peak Number 26 6H-Dibenzo[b,d]-pyran Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.111	10.71 ng/ul	237826	Phenanthrene-d10	17.380

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6H-Dibenzo[b,d]-pyran	182	C13H10O	000229-95-8	90
2	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	72
3	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	68
4	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	64
5	2-Hydroxyfluorene	182	C13H10O	002443-58-5	58





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Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YAS53

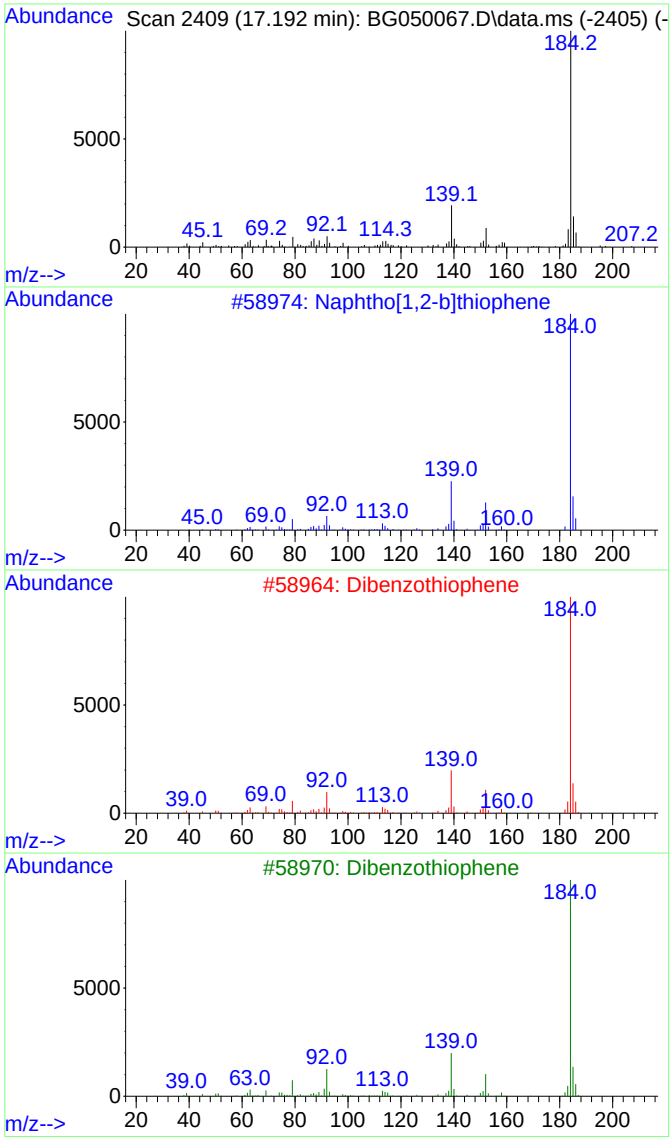
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

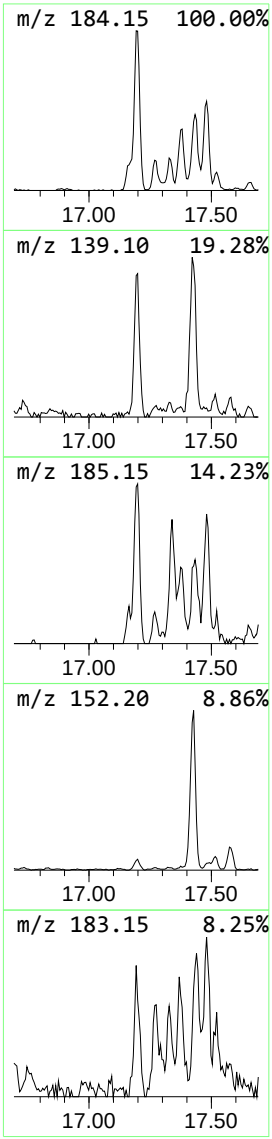
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 32 Naphtho[1,2-b]thiophene Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.192	9.53 ng/ul	211723	Phenanthrene-d10	17.380

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	96
2	Dibenzothiophene	184	C12H8S	000132-65-0	95
3	Dibenzothiophene	184	C12H8S	000132-65-0	95
4	Dibenzothiophene	184	C12H8S	000132-65-0	95
5	Dibenzothiophene	184	C12H8S	000132-65-0	94





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Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YAS53

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

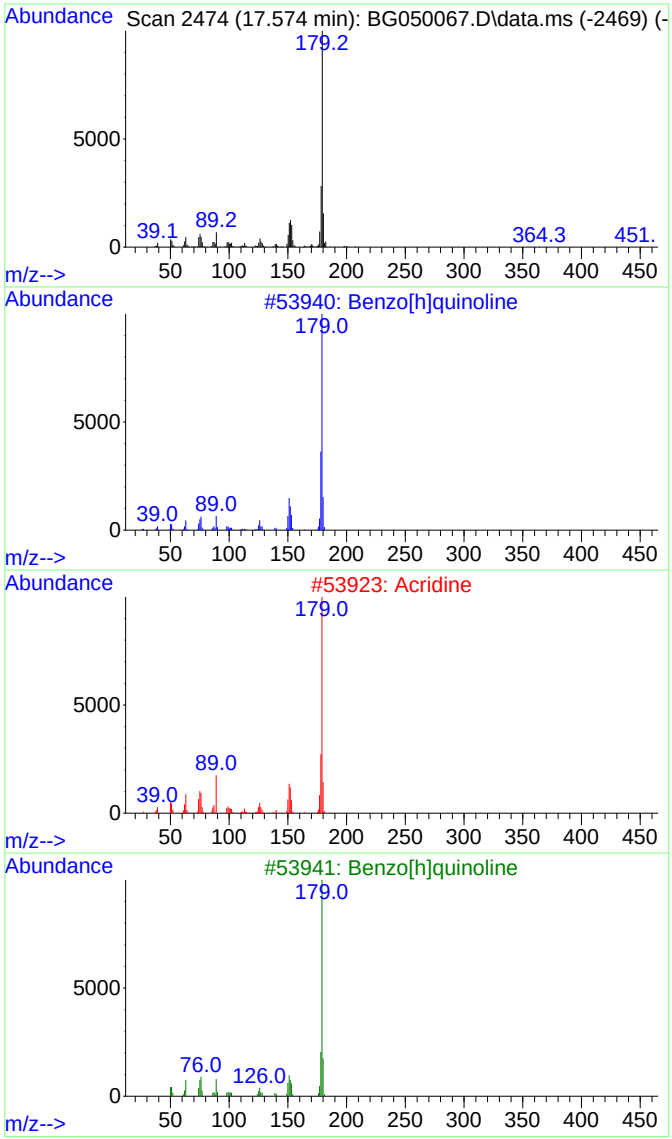
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

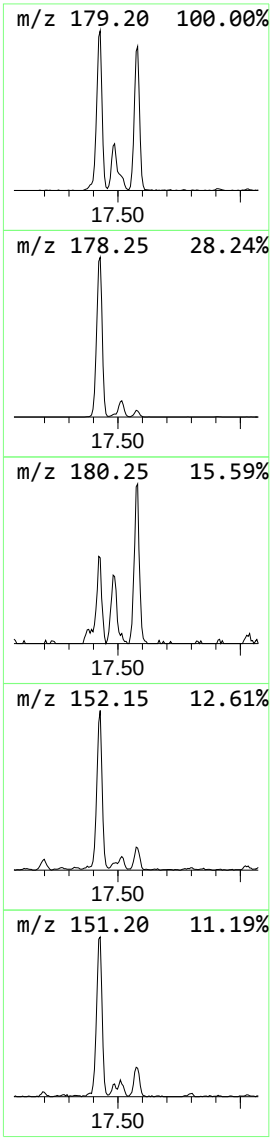
****

Peak Number 33 Benzo[h]quinoline Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.574	16.40 ng/ul	364247	Phenanthrene-d10	17.380

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[h]quinoline	179	C13H9N	000230-27-3	96
2	Acridine	179	C13H9N	000260-94-6	94
3	Benzo[h]quinoline	179	C13H9N	000230-27-3	93
4	Phenanthridine	179	C13H9N	000229-87-8	91
5	Acridine	179	C13H9N	000260-94-6	90





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Acq On : 31 Aug 2021 16:24
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Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YAS53

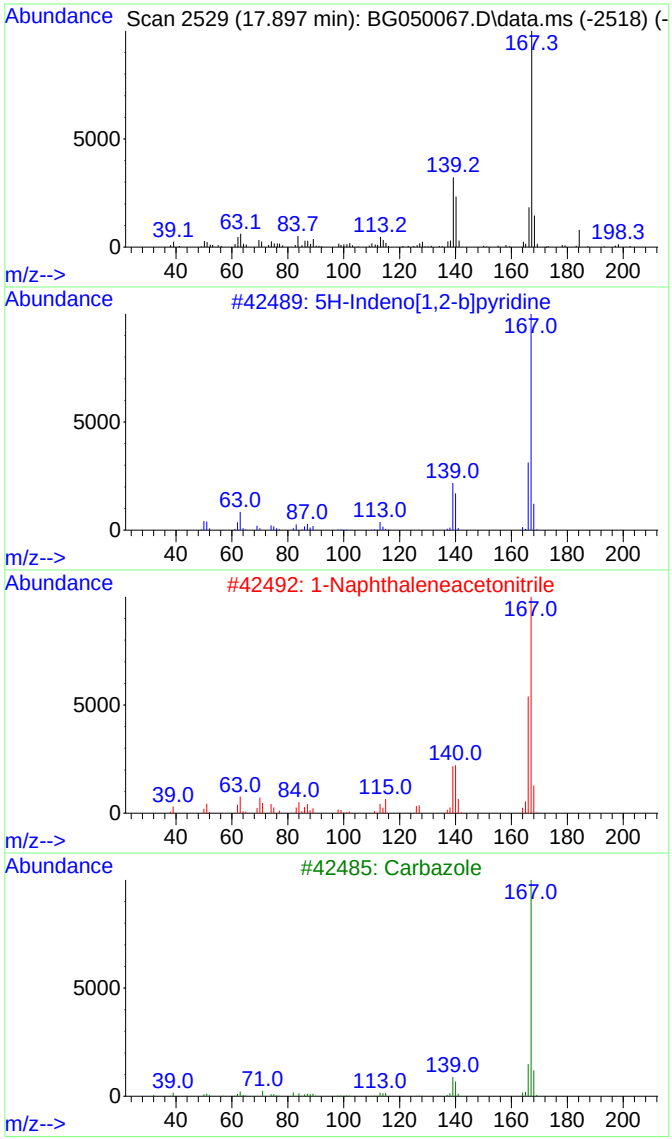
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

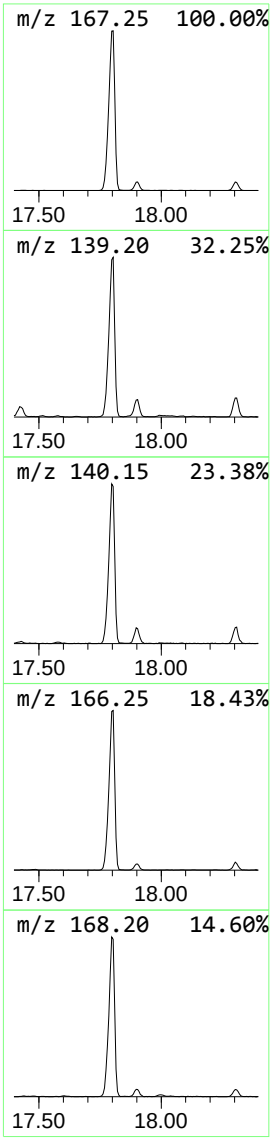
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 34 5H-Indeno[1,2-b]pyridine Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.897	14.16 ng/ul	314566	Phenanthrene-d10	17.380

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	90
2	1-Naphthaleneacetonitrile	167	C12H9N	000132-75-2	87
3	Carbazole	167	C12H9N	000086-74-8	83
4	5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	81
5	Carbazole	167	C12H9N	000086-74-8	81





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Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YAS53

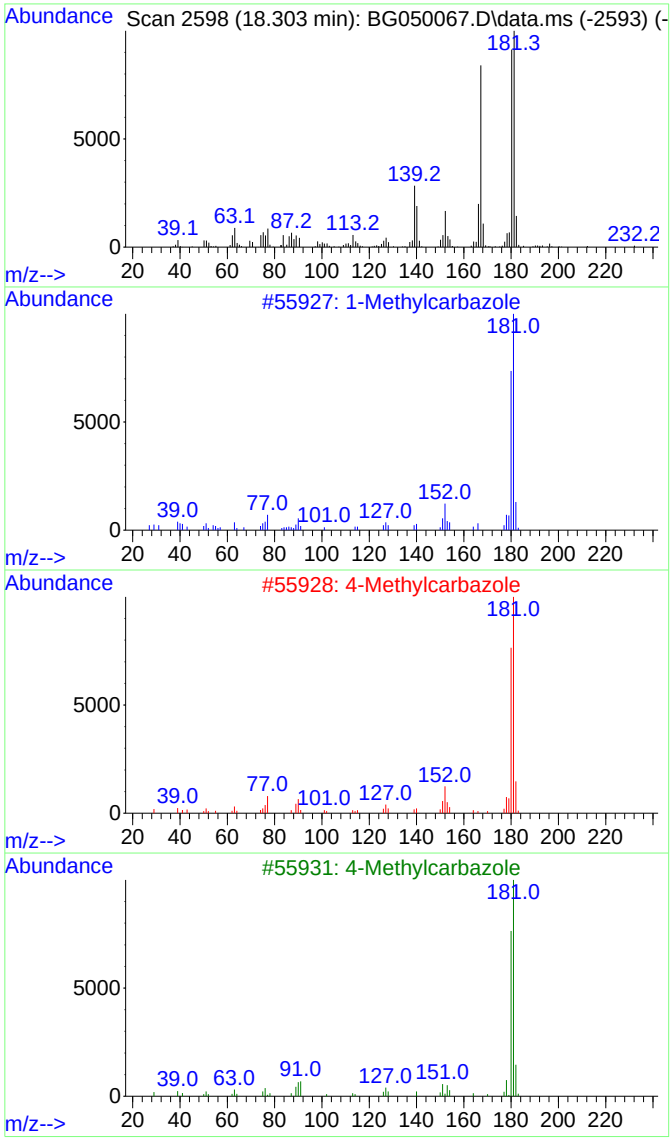
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

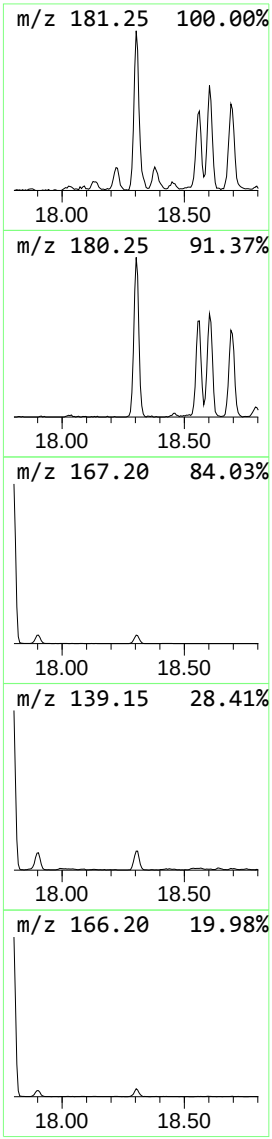
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 38 1-Methylcarbazole Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.303	25.10 ng/ul	557481	Phenanthrene-d10	17.380

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Methylcarbazole	181	C13H11N	006510-65-2	86
2	4-Methylcarbazole	181	C13H11N	003770-48-7	70
3	4-Methylcarbazole	181	C13H11N	003770-48-7	70
4	9H-Carbazole, 9-methyl-	181	C13H11N	001484-12-4	70
5	5H-Indeno[1,2-b]pyridine, 4-methyl-	181	C13H11N	064292-01-9	64





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Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YAS53

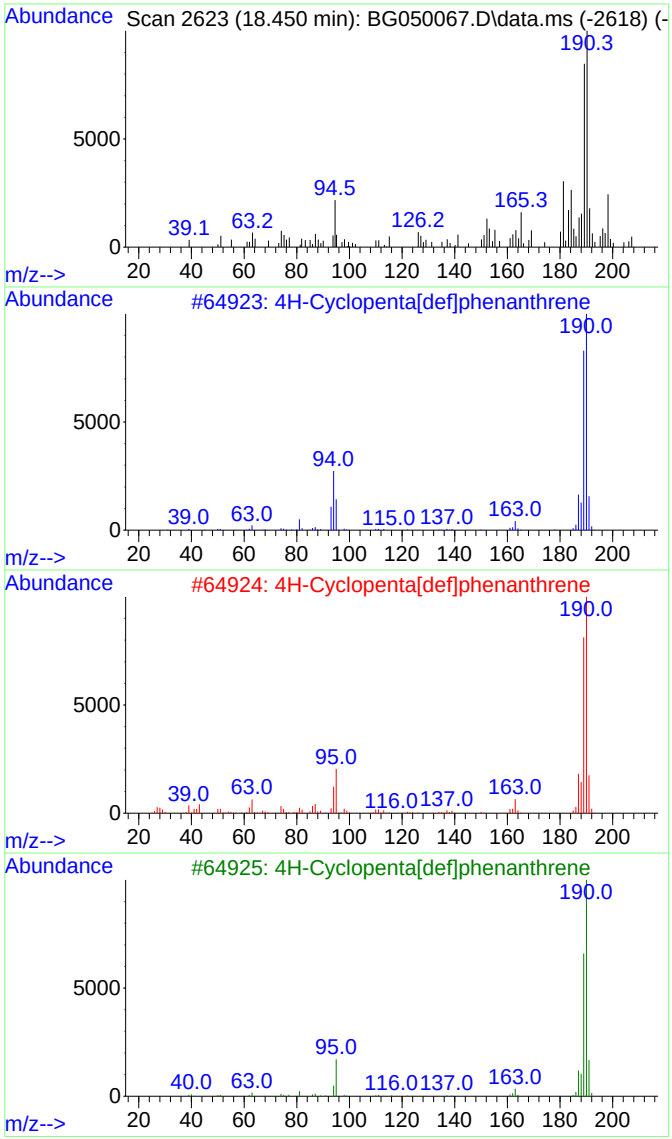
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

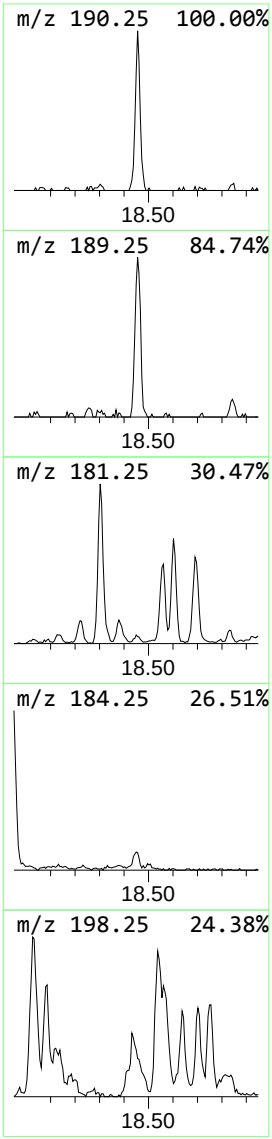
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 39 4H-Cyclopenta[def]phenanthrene Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.450	5.24 ng/ul	116460	Phenanthrene-d10	17.380

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	62
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	53
4	Phenol, 4-(2-thienylmethyl)-	190	C11H10OS	091680-55-6	50
5	2-Propynyl 1H-purin-6-yl sulfide	190	C8H6N4S	1000486-13-9	50





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Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

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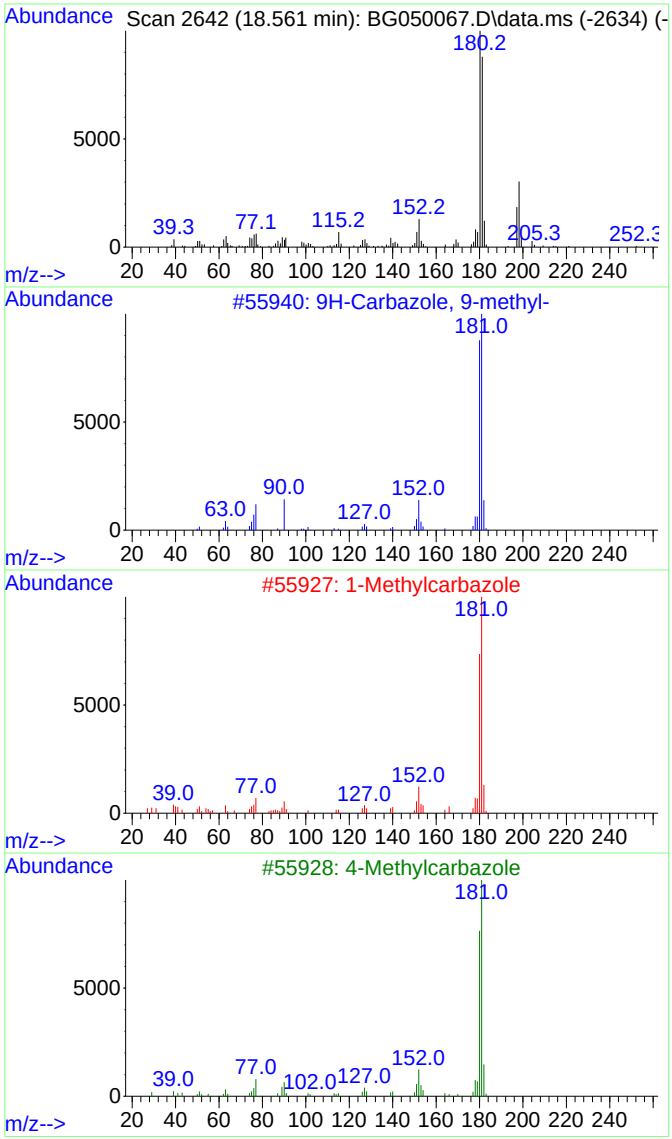
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

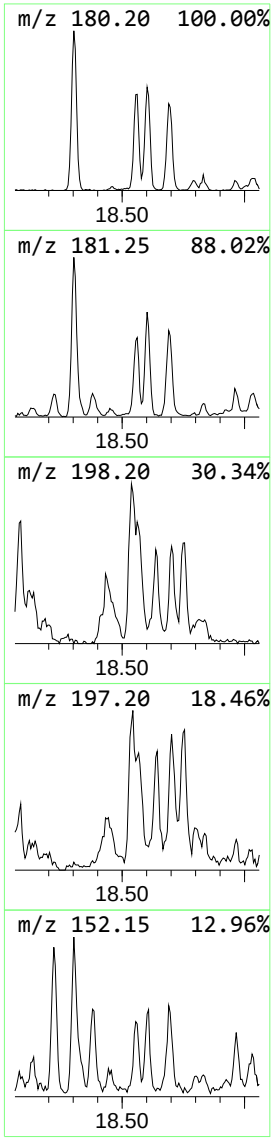
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 40 9H-Carbazole, 9-methyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.561	14.12 ng/ul	313633	Phenanthrene-d10	17.380

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Carbazole, 9-methyl-	181	C13H11N	001484-12-4	90
2	1-Methylcarbazole	181	C13H11N	006510-65-2	83
3	4-Methylcarbazole	181	C13H11N	003770-48-7	83
4	2-Fluorenamine	181	C13H11N	000153-78-6	72
5	Acridine, 9,10-dihydro-	181	C13H11N	000092-81-9	72





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Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YAS53

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

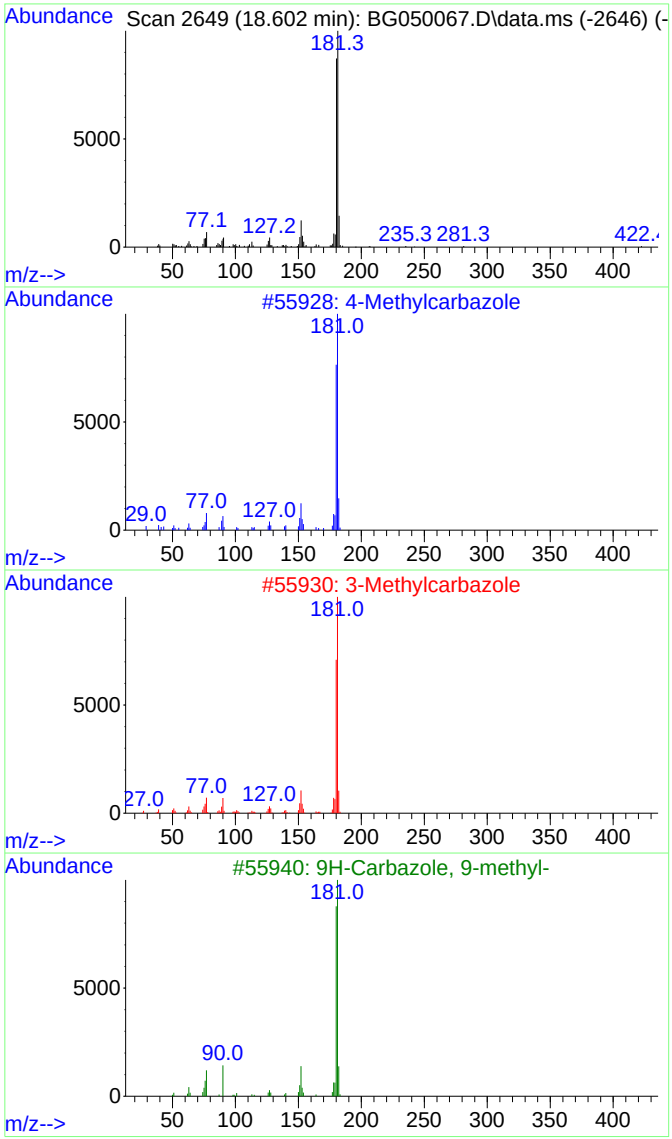
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

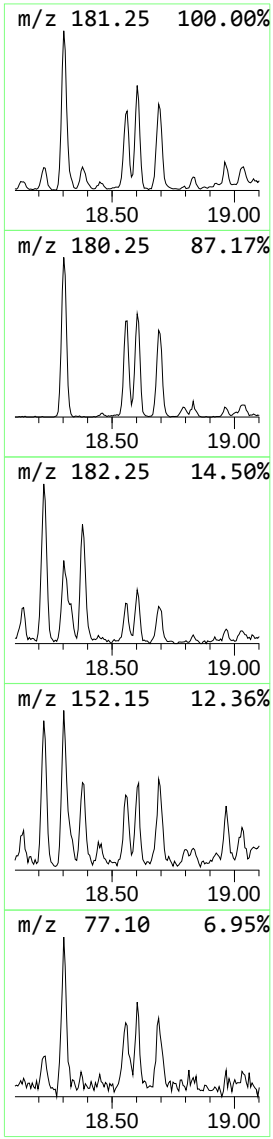
****

Peak Number 41 4-Methylcarbazole Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.602	8.23 ng/ul	182754	Phenanthrene-d10	17.380

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Methylcarbazole	181	C13H11N	003770-48-7	96
2	3-Methylcarbazole	181	C13H11N	004630-20-0	94
3	9H-Carbazole, 9-methyl-	181	C13H11N	001484-12-4	94
4	1-Methylcarbazole	181	C13H11N	006510-65-2	94
5	4aH-carbazole, 6-methyl-	181	C13H11N	1000404-86-1	91





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Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YAS53

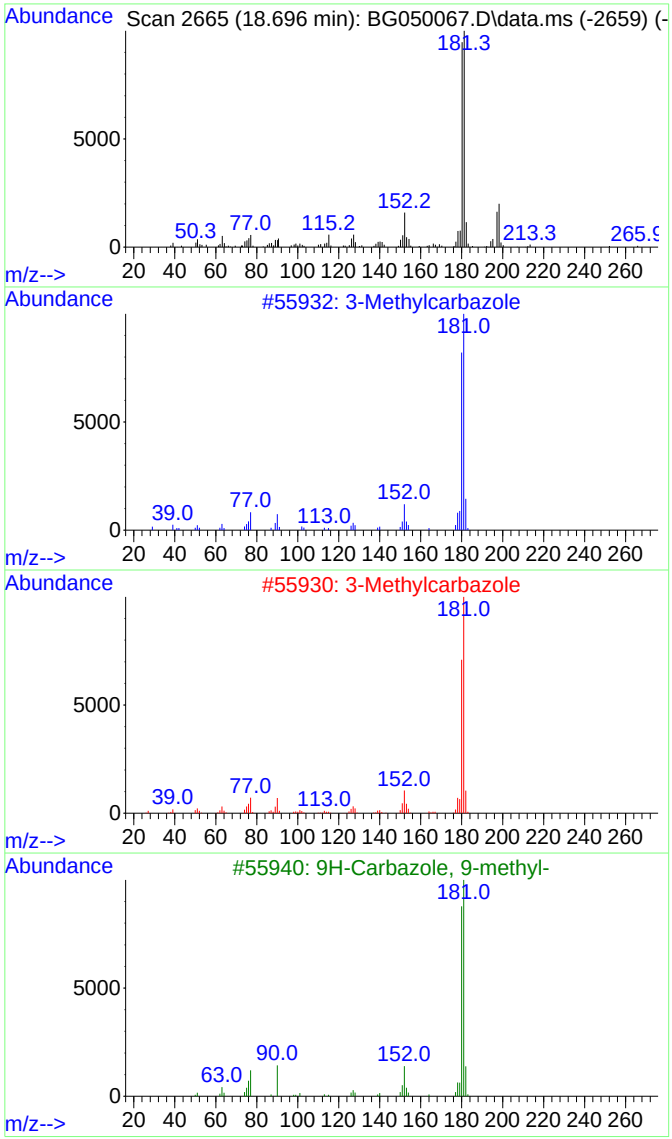
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

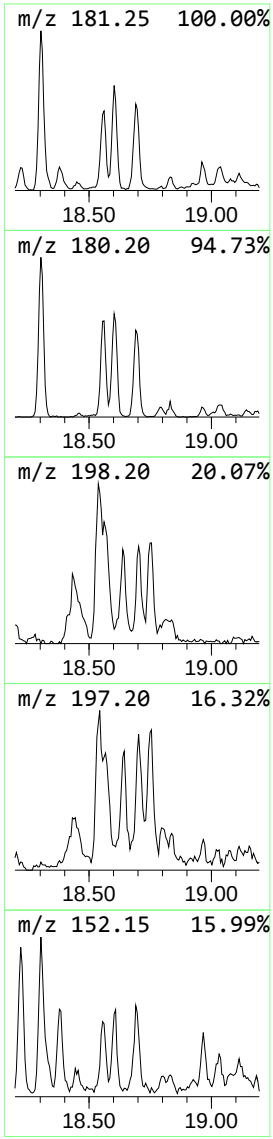
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 42 3-Methylcarbazole Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.696	9.82 ng/ul	218118	Phenanthrene-d10	17.380

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Methylcarbazole	181	C13H11N	004630-20-0	95
2	3-Methylcarbazole	181	C13H11N	004630-20-0	94
3	9H-Carbazole, 9-methyl-	181	C13H11N	001484-12-4	93
4	1-Methylcarbazole	181	C13H11N	006510-65-2	93
5	2-Fluorenamine	181	C13H11N	000153-78-6	93





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Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YAS53

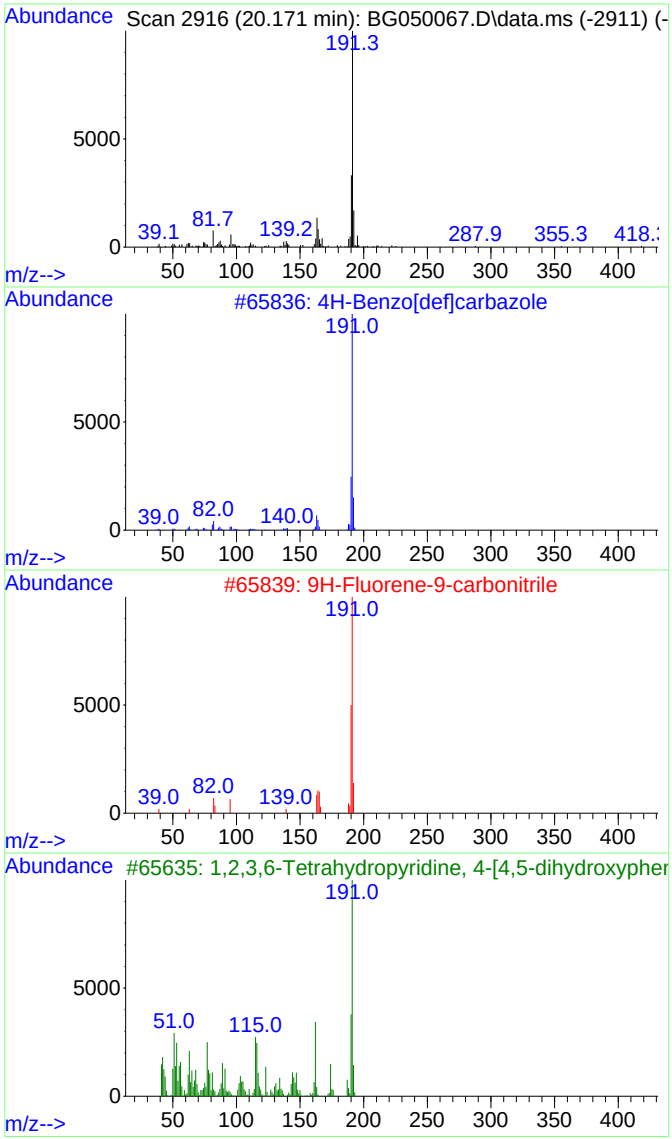
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

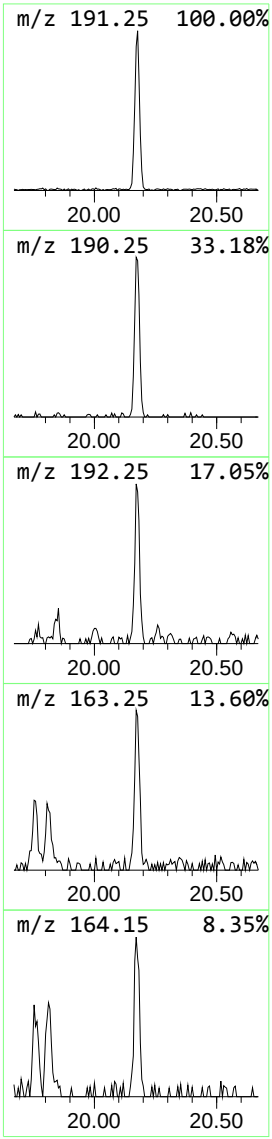
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 43 4H-Benzo[def]carbazole Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.171	7.81 ng/ul	149007	Chrysene-d12	21.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Benzo[def]carbazole	191	C14H9N	000203-65-6	93
2	9H-Fluorene-9-carbonitrile	191	C14H9N	001529-40-4	59
3	1,2,3,6-Tetrahydropyridine, 4-[4...	191	C11H13NO2	090684-17-6	59
4	9H-Fluorene-2-carbonitrile	191	C14H9N	002523-48-0	53
5	9H-Fluorene-9-carbonitrile	191	C14H9N	001529-40-4	53





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
Data File : BG050067.D
Acq On : 31 Aug 2021 16:24
Operator : CG/JU
Sample : M3494-01
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : YAS53

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...	5.949	65.0	ng/ul	651342	1	7.964	200363	20.0
Benzene, (1-met...	6.430	17.6	ng/ul	176744	1	7.964	200363	20.0
Benzene, 1-ethy...	7.341	21.3	ng/ul	213516	1	7.964	200363	20.0
Mesitylene	7.605	80.6	ng/ul	807854	1	7.964	200363	20.0
Benzofuran	7.682	11.9	ng/ul	119687	1	7.964	200363	20.0
Benzene, 1,2,3-...	8.075	87.0	ng/ul	871641	1	7.964	200363	20.0
Benzene, (2-bro...	8.146	11.8	ng/ul	118041	1	7.964	200363	20.0
Indane	8.363	1660.9	ng/ul	16639000	1	7.964	200363	20.0
Benzene, 1-prop...	8.504	40.0	ng/ul	400800	1	7.964	200363	20.0
Benzene, 1-ethy...	8.604	13.8	ng/ul	137799	1	7.964	200363	20.0
Benzene, 1-ethy...	9.074	42.2	ng/ul	423143	1	7.964	200363	20.0
Benzofuran, 2-m...	9.332	104.7	ng/ul	1049110	1	7.964	200363	20.0
Naphthalene, 2-...	13.650	33.6	ng/ul	548157	3	14.625	326093	20.0
Naphthalene, 2,...	13.803	52.8	ng/ul	860170	3	14.625	326093	20.0
Naphthalene, 1,...	13.944	52.3	ng/ul	852812	3	14.625	326093	20.0
Naphthalene, 2,...	14.185	18.9	ng/ul	307877	3	14.625	326093	20.0
2-Naphthaleneca...	14.766	51.7	ng/ul	842257	3	14.625	326093	20.0
Ethyl [(dipheny...	15.835	14.5	ng/ul	236536	3	14.625	326093	20.0
6H-Dibenzo[b,d]...	16.111	10.7	ng/ul	237826	4	17.380	444212	20.0
Naphtho[1,2-b]t...	17.192	9.5	ng/ul	211723	4	17.380	444212	20.0
Benzo[h]quinoline	17.574	16.4	ng/ul	364247	4	17.380	444212	20.0
5H-Indeno[1,2-b...	17.897	14.2	ng/ul	314566	4	17.380	444212	20.0
1-Methylcarbazole	18.303	25.1	ng/ul	557481	4	17.380	444212	20.0
4H-Cyclopenta[d...	18.450	5.2	ng/ul	116460	4	17.380	444212	20.0
9H-Carbazole, 9...	18.561	14.1	ng/ul	313633	4	17.380	444212	20.0
4-Methylcarbazole	18.602	8.2	ng/ul	182754	4	17.380	444212	20.0
3-Methylcarbazole	18.696	9.8	ng/ul	218118	4	17.380	444212	20.0
4H-Benzo[def]ca...	20.171	7.8	ng/ul	149007	5	21.651	381368	20.0