LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
Data File : BG062100.D
Acq On : 16 Jul 2024 18:59
Operator : MA/JU
Sample : P3177-14DL 5X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : A4BX8DL

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-BG070224.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BG062100.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.890	97	102	108	rBV4	4241	7572	0.40%	0.056%
2	4.653	227	232	235	rVB2	1937	3183	0.17%	0.023%
3	5.029	290	296	300	rBV2	1770	4448	0.23%	0.033%
4	5.106	307	309	311	rBV2	3272	3257	0.17%	0.024%
5	5.223	328	329	333	rVB2	2883	2630	0.14%	0.019%
6	5.323	345	346	353	rVB	2458	3920	0.20%	0.029%
7	5.799	425	427	432	rBV2	2278	3551	0.19%	0.026%
8	7.227	665	670	679	rVB3	31202	55534	2.90%	0.408%
9	7.380	688	696	702	rBV2	17094	31199	1.63%	0.229%
10	7.585	723	731	738	rBV3	28961	51828	2.70%	0.381%
11	7.673	743	746	749	rVB3	2149	2796	0.15%	0.021%
12	7.732	753	756	759	rVB	1951	2564	0.13%	0.019%
13	7.767	759	762	766	rBV	2984	3429	0.18%	0.025%
14	7.920	785	788	791	rVB2	1813	2582	0.13%	0.019%
15	8.055	802	811	818	rBV	164311	270998	14.14%	1.993%
16	8.508	883	888	893	rBV	3600	6910	0.36%	0.051%
17	8.778	926	934	939	rBV2	39978	77250	4.03%	0.568%
18	8.837	941	944	950	rVB3	4084	7506	0.39%	0.055%
19	9.219	1003	1009	1016	rBV2	31734	60215	3.14%	0.443%
20	9.371	1031	1035	1038	rVB3	1910	2560	0.13%	0.019%
21	9.941	1125	1132	1140	rBV3	29015	55595	2.90%	0.409%
22	10.353	1199	1202	1204	rBV2	3167	3095	0.16%	0.023%
23	10.406	1210	1211	1214	rBV2	3117	3537	0.18%	0.026%
24	10.500	1221	1227	1235	rVB4	38023	68024	3.55%	0.500%
25	10.870	1283	1290	1294	rBV	230553	413376	21.57%	3.040%
26	10.923	1294	1299	1305	rVB	183599	321138	16.76%	2.362%
27	11.005	1308	1313	1319	rBV3	21449	42338	2.21%	0.311%
28	11.234	1349	1352	1355	rVB	2040	2757	0.14%	0.020%
29	11.604	1410	1415	1420	rBV2	3532	6239	0.33%	0.046%
30	11.698	1426	1431	1436	rBV4	12805	19962	1.04%	0.147%
31	11.775	1438	1444	1448	rVB4	2228	4107	0.21%	0.030%
32	12.027	1484	1487	1490	rVB	2884	3457	0.18%	0.025%
33	12.057	1490	1492	1496	rBV	2362	3070	0.16%	0.023%
34	12.245	1521	1524	1529	rBV	3187	5300	0.28%	0.039%
35	12.521	1563	1571	1578	rBV2	70902	127671	6.66%	0.939%
36	12.626	1588	1589	1591	rBV2	3618	3453	0.18%	0.025%

LSC Area Percent Report

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Sample : P3177-14DL 5X
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Instrument :
BNA_G
ClientSampleId :
A4BX8DL

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-BG070224.MA.M
Title : SVOA CALIBRATION

37	12.738	1601	1608	1615	rVB2	38961	68583	3.58%	0.504%
38	13.032	1655	1658	1661	rVB	2763	3827	0.20%	0.028%
39	13.238	1687	1693	1697	rVB2	4345	7923	0.41%	0.058%
40	13.490	1732	1736	1737	rBV2	2399	2774	0.14%	0.020%
41	13.525	1737	1742	1746	rBV2	21290	34522	1.80%	0.254%
42	13.678	1761	1768	1770	rBV3	4662	9902	0.52%	0.073%
43	13.719	1770	1775	1778	rBV3	11953	14861	0.78%	0.109%
44	14.007	1818	1824	1829	rBV4	24151	41673	2.17%	0.307%
45	14.084	1829	1837	1843	rVB2	90079	150165	7.84%	1.104%
46	14.136	1843	1846	1848	rBV2	2288	2569	0.13%	0.019%
47	14.248	1861	1865	1867	rBV2	11650	13415	0.70%	0.099%
48	14.313	1873	1876	1877	rBV	4009	3263	0.17%	0.024%
49	14.383	1881	1888	1898	rBV	84455	145691	7.60%	1.072%
50	14.689	1930	1940	1946	rBV	393465	629380	32.85%	4.629%
51	14.748	1946	1950	1957	rVB	142219	219565	11.46%	1.615%
52	14.824	1957	1963	1968	rBV	11194	15429	0.81%	0.113%
53	14.871	1968	1971	1972	rBV3	5218	4842	0.25%	0.036%
54	14.912	1974	1978	1982	rVB5	27964	41620	2.17%	0.306%
55	15.082	2001	2007	2014	rVV	173484	274289	14.32%	2.017%
56	15.153	2017	2019	2022	rVV4	9642	9550	0.50%	0.070%
57	15.206	2026	2028	2029	rBV	3451	2811	0.15%	0.021%
58	15.235	2032	2033	2036	rVB2	4378	3757	0.20%	0.028%
59	15.282	2038	2041	2042	rVV2	6533	6111	0.32%	0.045%
60	15.300	2042	2044	2048	rVV4	10686	10854	0.57%	0.080%
61	15.353	2048	2053	2057	rVB5	6715	10378	0.54%	0.076%
62	15.505	2077	2079	2084	rVB4	7125	6309	0.33%	0.046%
63	15.599	2094	2095	2096	rBV	4735	3125	0.16%	0.023%
64	15.676	2099	2108	2112	rVV2	104246	165121	8.62%	1.214%
65	15.735	2112	2118	2125	rVB	167800	268261	14.00%	1.973%
66	15.799	2126	2129	2131	rBV2	10101	11292	0.59%	0.083%
67	15.893	2142	2145	2150	rVB2	17595	24131	1.26%	0.177%
68	15.975	2156	2159	2163	rBV3	13430	22195	1.16%	0.163%
69	16.022	2163	2167	2173	rVB2	36333	54941	2.87%	0.404%
70	16.169	2187	2192	2198	rBV2	44200	81178	4.24%	0.597%
71	16.293	2211	2213	2215	rBV2	2606	2715	0.14%	0.020%
72	16.334	2218	2220	2223	rBV2	3484	2921	0.15%	0.021%
73	16.363	2223	2225	2227	rBV2	6728	5205	0.27%	0.038%
74	16.575	2257	2261	2262	rBV4	5204	5412	0.28%	0.040%
75	16.733	2279	2288	2292	rVB4	31457	62347	3.25%	0.459%
76	16.780	2293	2296	2298	rBV3	8801	9932	0.52%	0.073%

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Operator : MA/JU
Sample : P3177-14DL 5X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : A4BX8DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Title : SVOA CALIBRATION

77	16.963	2322	2327	2330	rBV7	12853	19845	1.04%	0.146%
78	17.086	2343	2348	2355	rVB	85564	132788	6.93%	0.977%
79	17.156	2357	2360	2367	rVB7	21943	36508	1.91%	0.269%
80	17.250	2372	2376	2385	rVB	77533	116129	6.06%	0.854%
81	17.433	2401	2407	2410	rBV	456763	696209	36.34%	5.121%
82	17.480	2410	2415	2420	rVV	1301625	1916032	100.00%	14.093%
83	17.532	2420	2424	2426	rVV2	141872	206160	10.76%	1.516%
84	17.568	2426	2430	2435	rVV	231242	348513	18.19%	2.563%
85	17.626	2436	2440	2445	rVB2	37916	58035	3.03%	0.427%
86	17.838	2468	2476	2482	rBV3	142655	249713	13.03%	1.837%
87	18.308	2551	2556	2560	rBV	86236	131457	6.86%	0.967%
88	18.355	2560	2564	2569	rVB2	139988	199040	10.39%	1.464%
89	18.431	2573	2577	2582	rBV	42571	61406	3.20%	0.452%
90	18.502	2583	2589	2593	rBV2	139105	255494	13.33%	1.879%
91	18.802	2636	2640	2643	rBV	81481	104619	5.46%	0.769%
92	18.843	2644	2647	2652	rVB	147785	195247	10.19%	1.436%
93	19.483	2751	2756	2762	rVB	1206529	1625032	84.81%	11.952%
94	19.812	2807	2812	2814	rBV2	293143	448445	23.40%	3.298%
95	19.842	2814	2817	2825	rVB	520058	751057	39.20%	5.524%
96	21.686	3126	3131	3137	rBV4	501824	1201934	62.73%	8.840%
97	21.745	3137	3141	3153	rVB	376285	702470	36.66%	5.167%

Sum of corrected areas: 13596023

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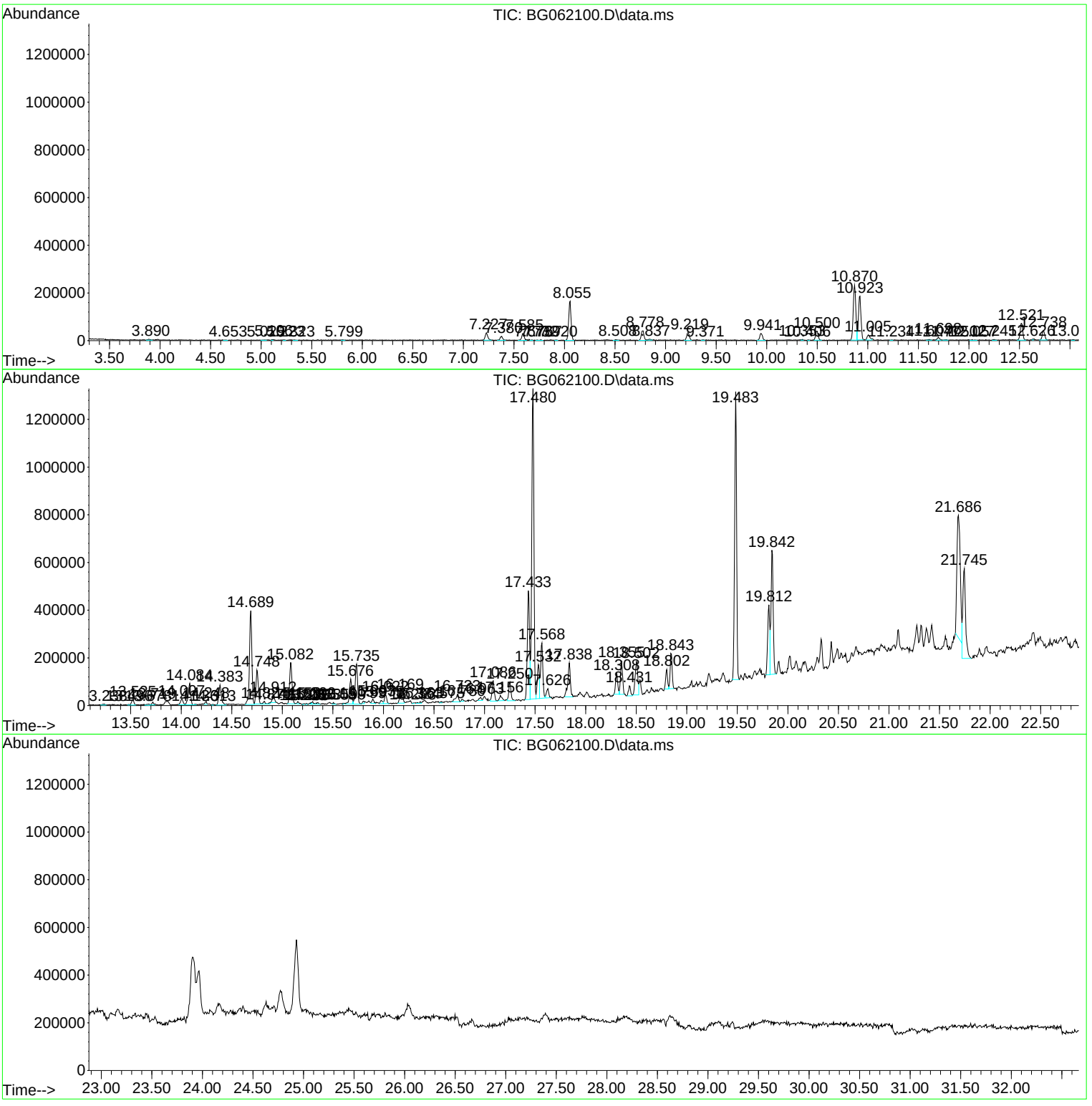
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4BX8DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
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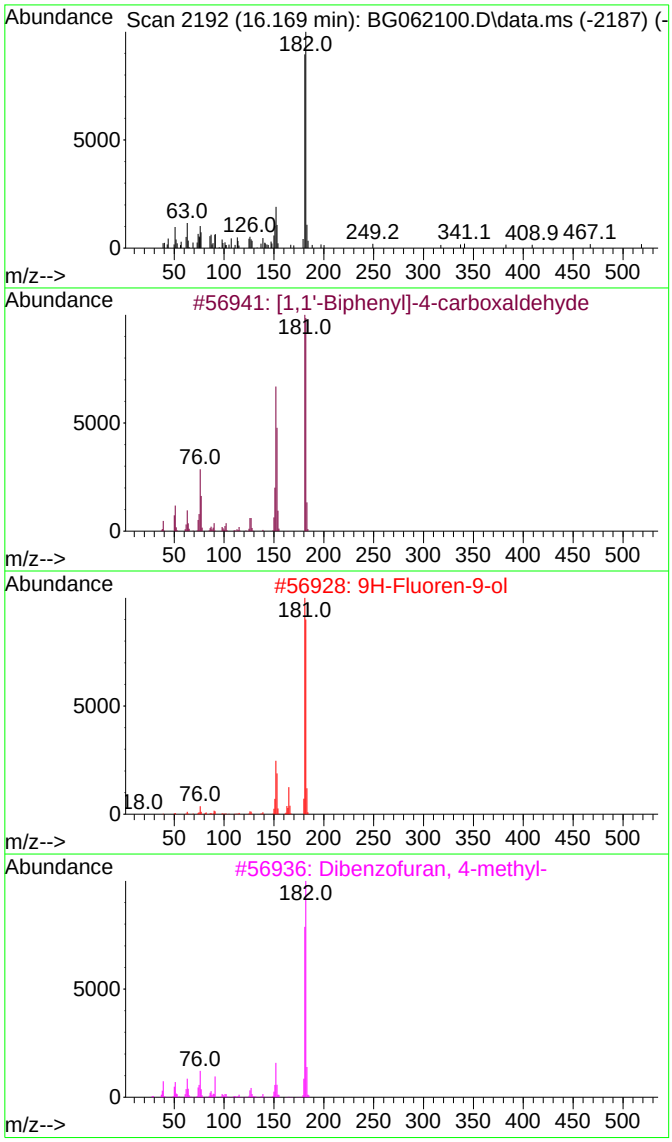
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

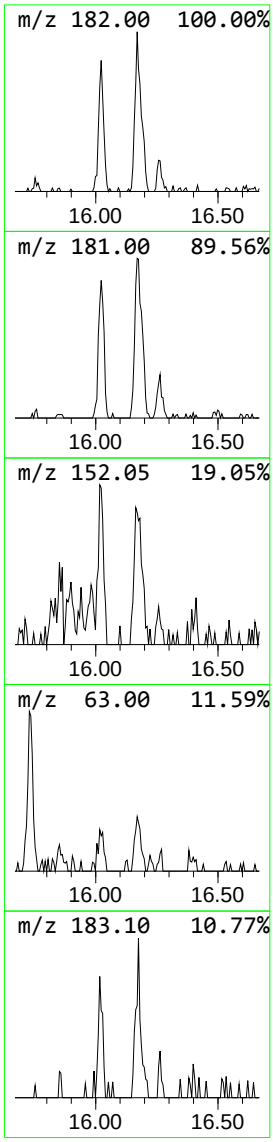
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 [1,1'-Biphenyl]-4-carboxald... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.169	2.33 ng/ul	81178	Phenanthrene-d10	17.433

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	87
2	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	86
3	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	83
4	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	74
5	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	68





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
Data File : BG062100.D
Acq On : 16 Jul 2024 18:59
Operator : MA/JU
Sample : P3177-14DL 5X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4BX8DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

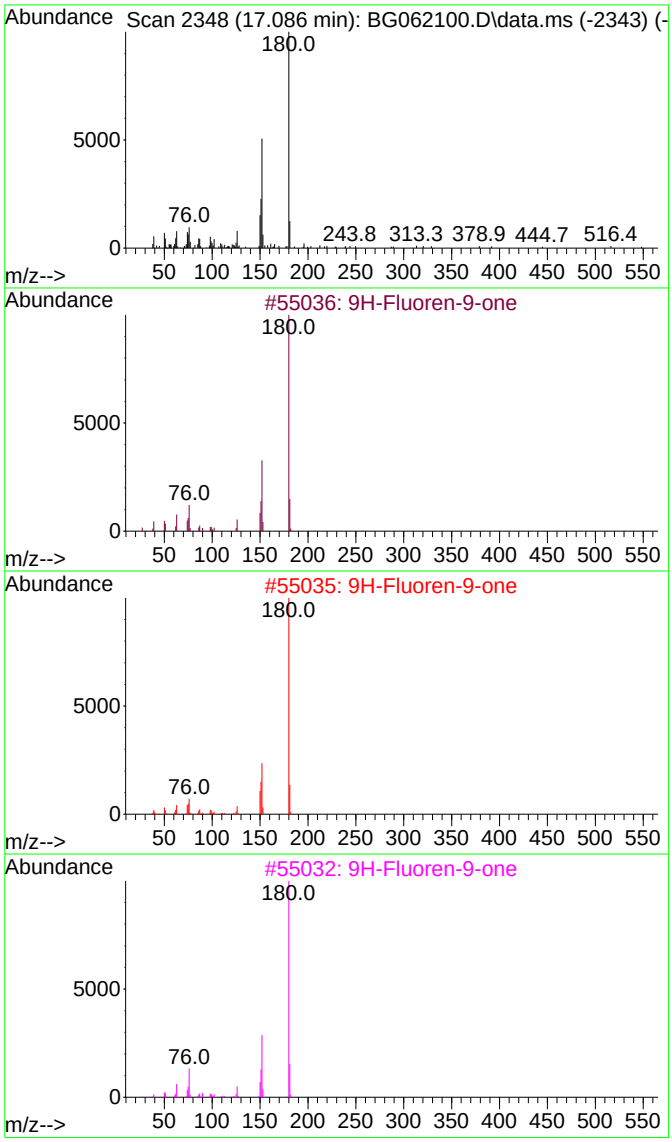
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

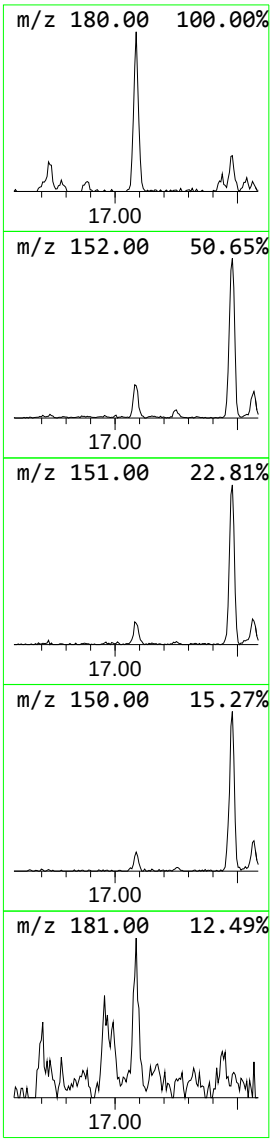
****

Peak Number 2 9H-Fluoren-9-one Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.086	3.81 ng/ul	132788	Phenanthrene-d10	17.433

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluoren-9-one	180	C13H8O	000486-25-9	95
2	9H-Fluoren-9-one	180	C13H8O	000486-25-9	94
3	9H-Fluoren-9-one	180	C13H8O	000486-25-9	91
4	9H-Fluoren-9-one	180	C13H8O	000486-25-9	91
5	9H-Fluoren-9-one	180	C13H8O	000486-25-9	91





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Instrument :
BNA_G
ClientSampleId :
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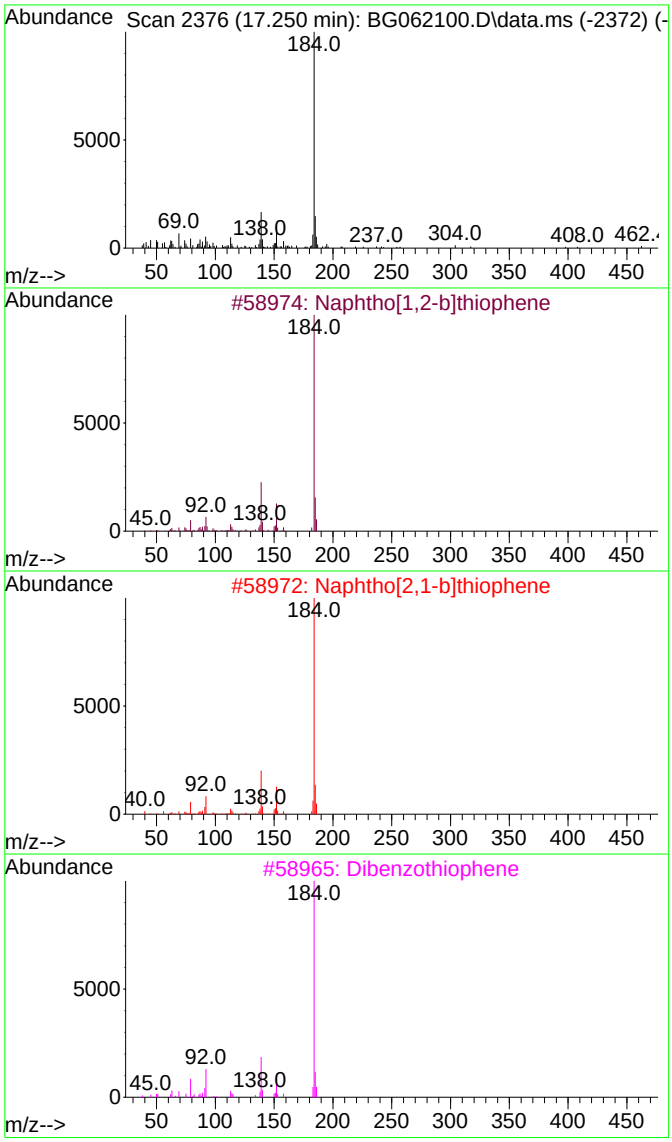
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

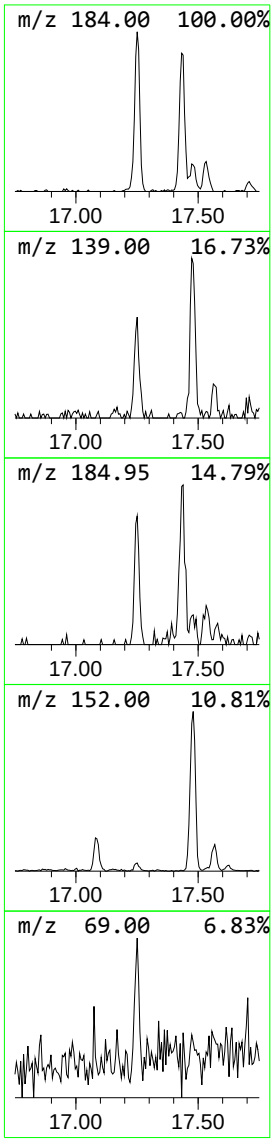
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Naphtho[1,2-b]thiophene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.250	3.34 ng/ul	116129	Phenanthrene-d10	17.433

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	95
2	Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	95
3	Dibenzothiophene	184	C12H8S	000132-65-0	94
4	Dibenzothiophene	184	C12H8S	000132-65-0	94
5	Dibenzothiophene	184	C12H8S	000132-65-0	94





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Instrument :
BNA_G
ClientSampleId :
A4BX8DL

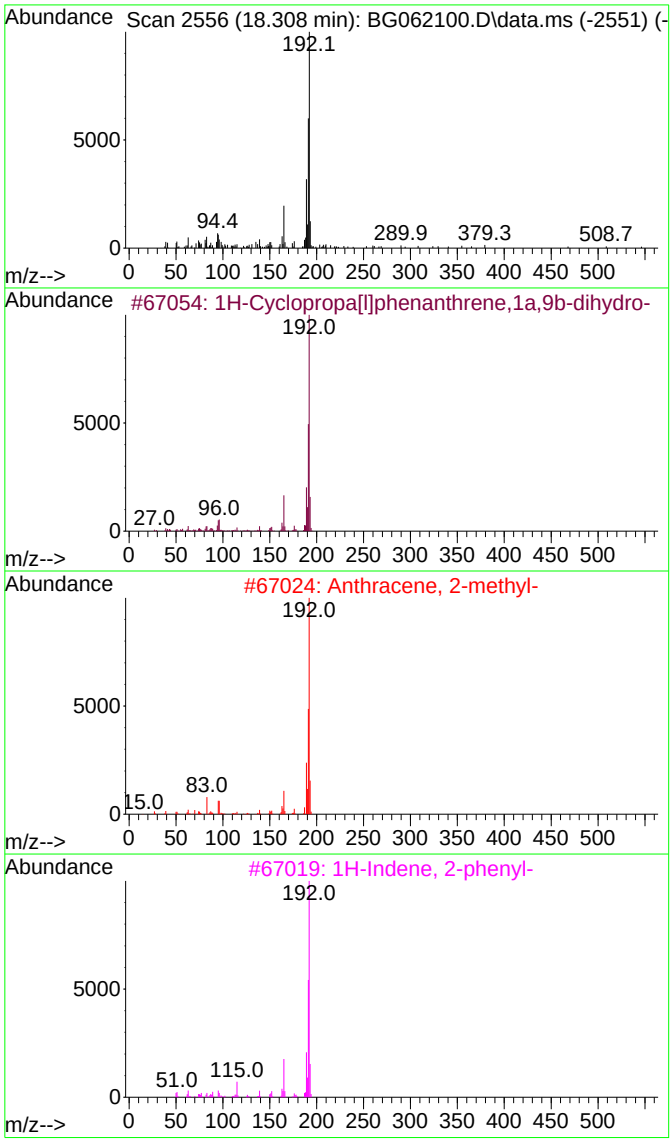
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

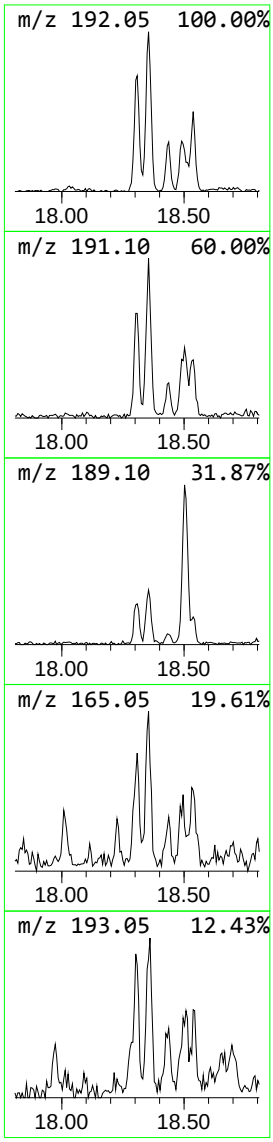
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 1H-Cyclopropa[l]phenanthren... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.308	3.78 ng/ul	131457	Phenanthrene-d10	17.433

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	90
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
3	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	87
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	83
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	81





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Data File : BG062100.D
Acq On : 16 Jul 2024 18:59
Operator : MA/JU
Sample : P3177-14DL 5X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4BX8DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

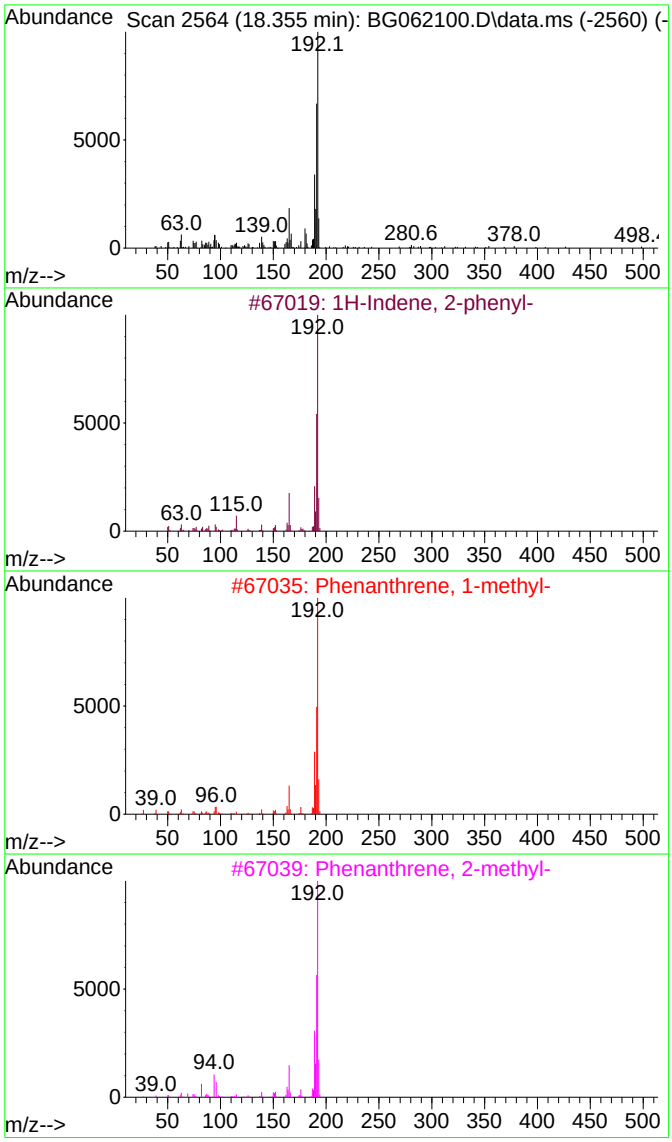
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

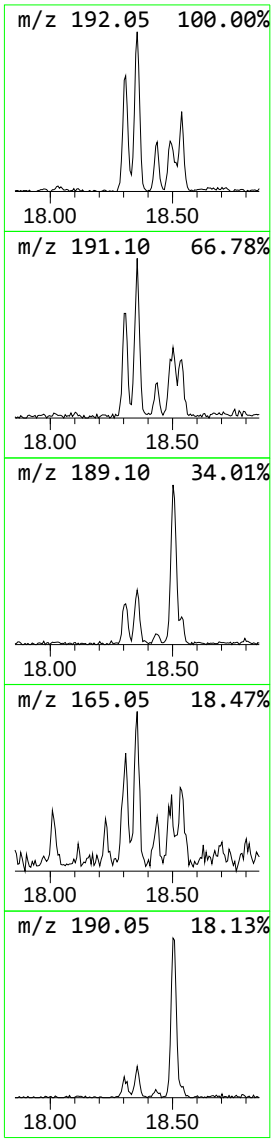
****

Peak Number 5 1H-Indene, 2-phenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.355	5.72 ng/ul	199040	Phenanthrene-d10	17.433

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	89
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	83
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	81
4	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	76
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	76





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
Data File : BG062100.D
Acq On : 16 Jul 2024 18:59
Operator : MA/JU
Sample : P3177-14DL 5X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4BX8DL

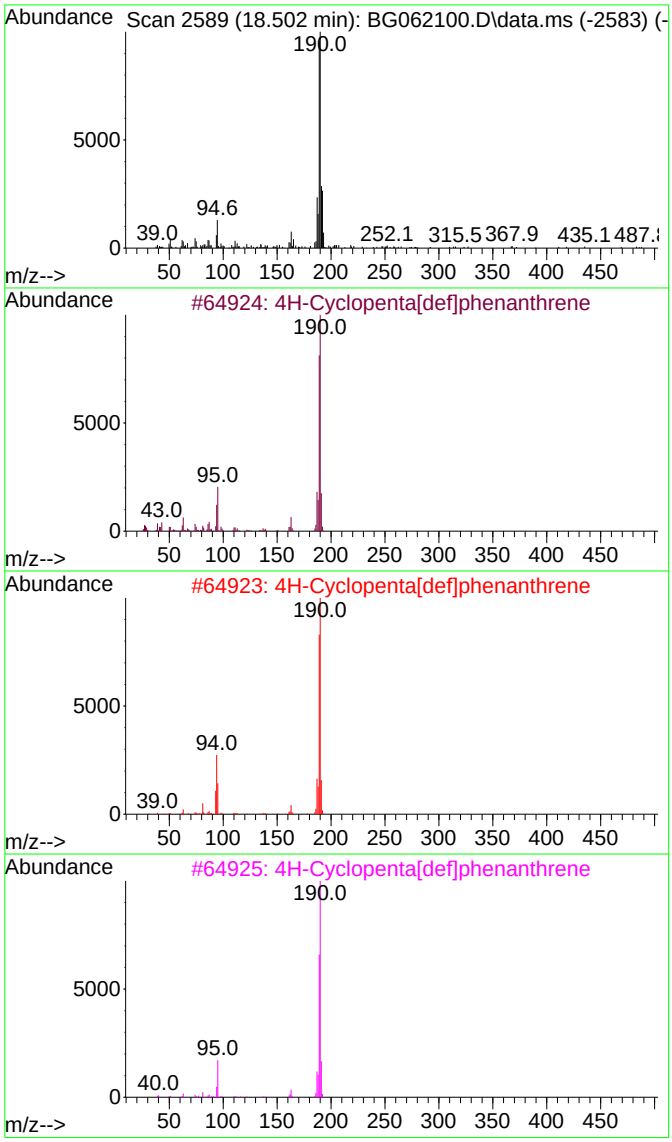
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

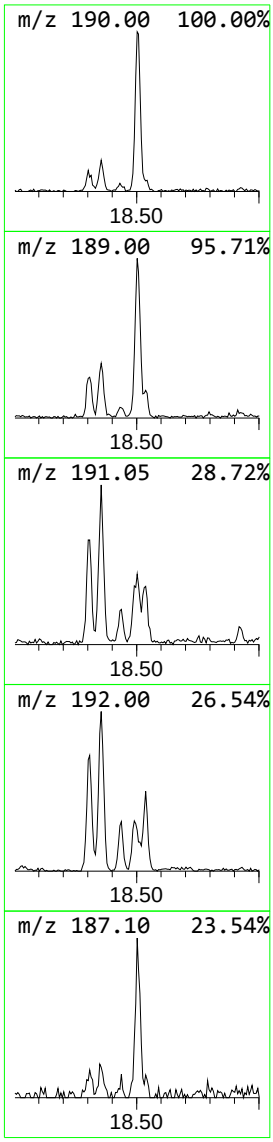
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.502	7.34 ng/ul	255494	Phenanthrene-d10	17.433

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	83
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
4	2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	50
5	Methyl diselenide	190	C2H6Se2	007101-31-7	45





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
Data File : BG062100.D
Acq On : 16 Jul 2024 18:59
Operator : MA/JU
Sample : P3177-14DL 5X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4BX8DL

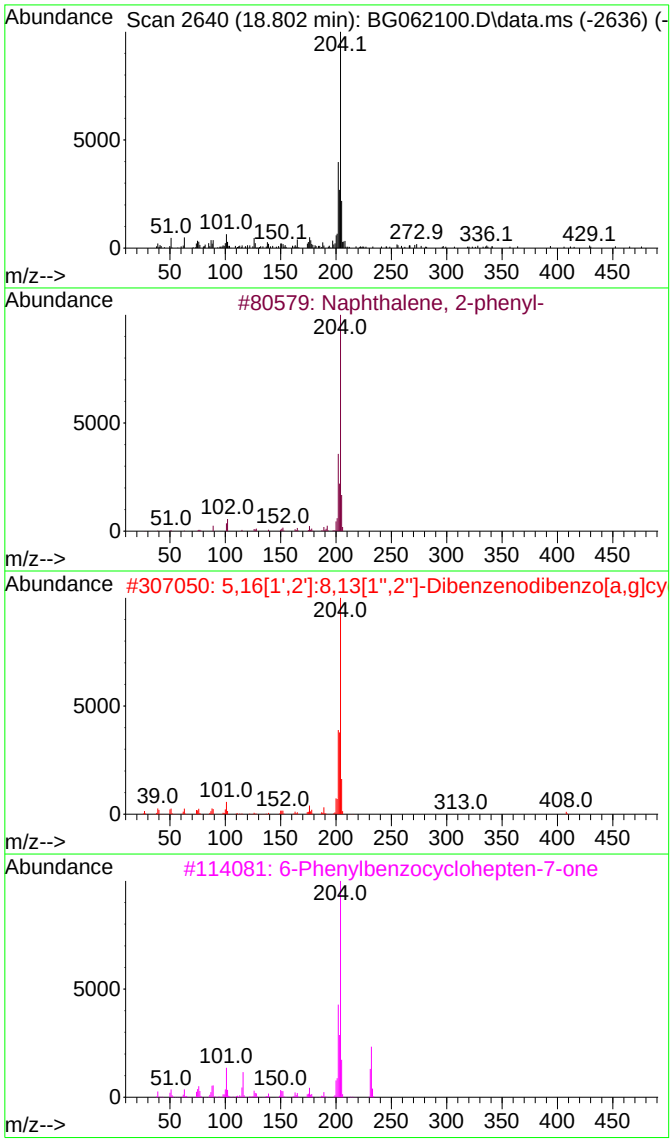
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

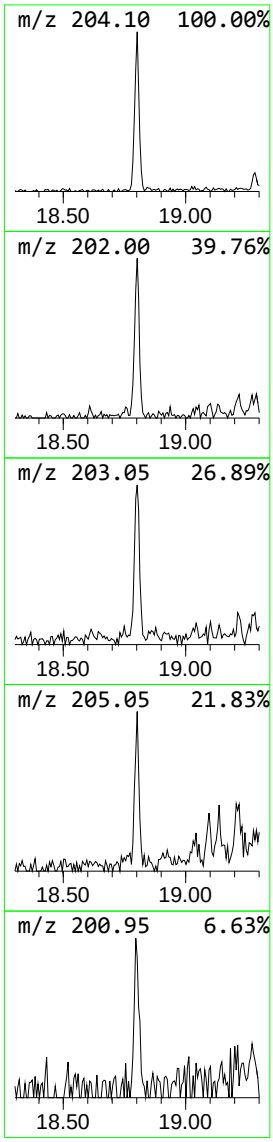
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Naphthalene, 2-phenyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.802	3.01 ng/ul	104619	Phenanthrene-d10	17.433

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	91
2	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	80
3	6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	72
4	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	64
5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	60





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
Data File : BG062100.D
Acq On : 16 Jul 2024 18:59
Operator : MA/JU
Sample : P3177-14DL 5X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4BX8DL

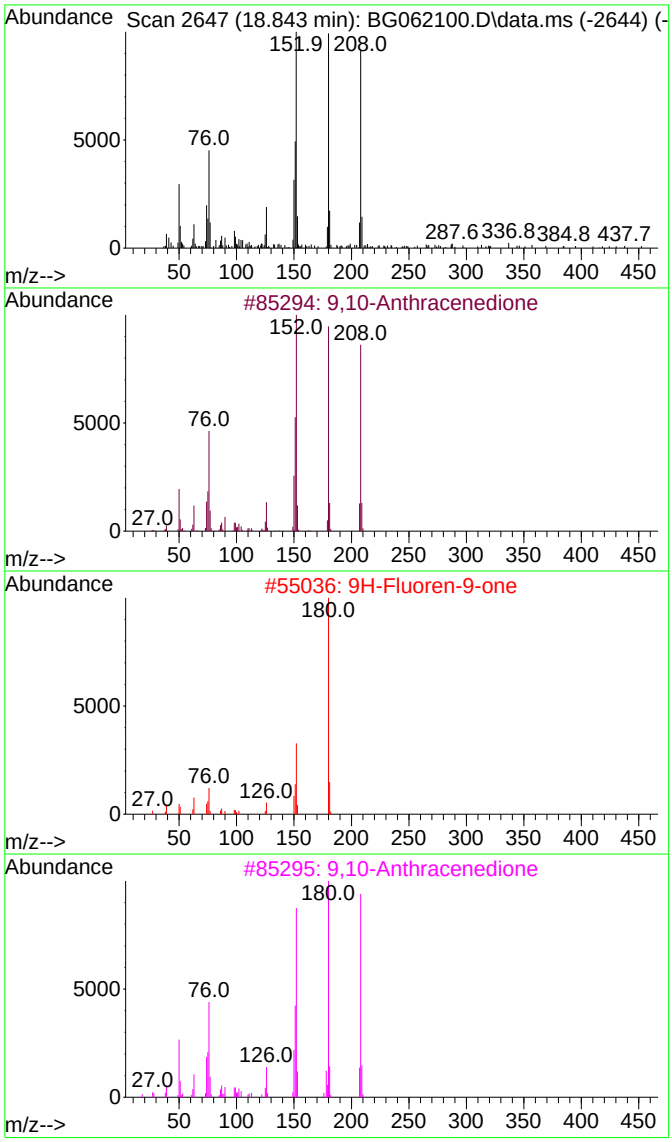
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

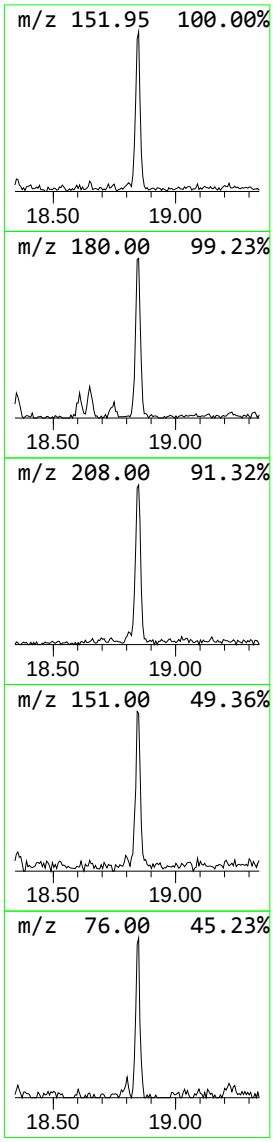
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 9,10-Anthracenedione Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.843	5.61 ng/ul	195247	Phenanthrene-d10	17.433

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	98
2	9H-Fluoren-9-one	180	C13H8O	000486-25-9	97
3	9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
4	9H-Fluoren-9-one	180	C13H8O	000486-25-9	93
5	9,10-Anthracenedione	208	C14H8O2	000084-65-1	93





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
Data File : BG062100.D
Acq On : 16 Jul 2024 18:59
Operator : MA/JU
Sample : P3177-14DL 5X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4BX8DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
[1,1'-Biphenyl]...	16.169	2.3	ng/ul	81178	4	17.433	696209	20.0
9H-Fluoren-9-one	17.086	3.8	ng/ul	132788	4	17.433	696209	20.0
Naphtho[1,2-b]t...	17.250	3.3	ng/ul	116129	4	17.433	696209	20.0
1H-Cyclopropa[l...	18.308	3.8	ng/ul	131457	4	17.433	696209	20.0
1H-Indene, 2-ph...	18.355	5.7	ng/ul	199040	4	17.433	696209	20.0
4H-Cyclopenta[d...	18.502	7.3	ng/ul	255494	4	17.433	696209	20.0
Naphthalene, 2-...	18.802	3.0	ng/ul	104619	4	17.433	696209	20.0
9,10-Anthracene...	18.843	5.6	ng/ul	195247	4	17.433	696209	20.0