LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112120\
Data File : BG047400.D
Acq On : 21 Nov 2020 18:28
Operator : CG/JU
Sample : L4711-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0B36

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\SOM-EPA-BG111920MA.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.613	41	47	52	rBV5	8017	14939	0.93%	0.074%
2	3.889	90	94	100	rBV2	9391	14912	0.93%	0.074%
3	4.606	212	216	219	rBV3	6834	9817	0.61%	0.049%
4	5.029	285	288	292	rBV3	5851	8577	0.53%	0.043%
5	7.227	657	662	665	rBV3	5358	10737	0.67%	0.053%
6	7.409	686	693	701	rBV2	210972	377872	23.54%	1.876%
7	7.591	715	724	732	rBV	119689	207281	12.91%	1.029%
8	7.802	752	760	766	rBV	196605	357005	22.24%	1.773%
9	7.849	766	768	774	rVB2	24501	30840	1.92%	0.153%
10	8.178	821	824	831	rVB2	3546	6174	0.38%	0.031%
11	8.278	831	841	847	rBV	190081	325939	20.31%	1.618%
12	8.325	847	849	854	rVB3	3867	4369	0.27%	0.022%
13	8.819	930	933	937	rBV3	2869	4409	0.27%	0.022%
14	8.966	949	958	966	rBV2	247314	425930	26.54%	2.115%
15	9.089	977	979	983	rBV4	3432	3973	0.25%	0.020%
16	9.442	1030	1039	1047	rVB	183856	343735	21.42%	1.707%
17	10.170	1156	1163	1170	rBV2	185090	338737	21.10%	1.682%
18	10.711	1247	1255	1262	rVV	275372	506805	31.58%	2.516%
19	11.104	1315	1322	1329	rVV	242217	442051	27.54%	2.195%
20	11.228	1333	1343	1354	rVB	218607	413699	25.77%	2.054%
21	11.404	1369	1373	1376	rVB	3630	4966	0.31%	0.025%
22	11.774	1430	1436	1441	rVB2	15662	28340	1.77%	0.141%
23	12.103	1488	1492	1498	rVB2	2822	4950	0.31%	0.025%
24	12.391	1538	1541	1546	rVB4	8347	11615	0.72%	0.058%
25	12.591	1569	1575	1580	rBV2	18607	29969	1.87%	0.149%
26	12.661	1585	1587	1592	rVB2	4555	7309	0.46%	0.036%
27	12.955	1633	1637	1639	rBV3	5365	5968	0.37%	0.030%
28	13.625	1748	1751	1753	rBV	3358	3943	0.25%	0.020%
29	13.719	1762	1767	1770	rBV4	5830	8863	0.55%	0.044%
30	13.772	1770	1776	1785	rVB	185251	310373	19.34%	1.541%
31	14.207	1846	1850	1852	rBV	3745	5024	0.31%	0.025%
32	14.283	1857	1863	1867	rVV	465001	682611	42.53%	3.389%
33	14.324	1867	1870	1877	rVV	178669	262774	16.37%	1.305%
34	14.506	1898	1901	1904	rBV2	4424	6131	0.38%	0.030%

LSC Area Percent Report

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Data File : BG047400.D
Acq On : 21 Nov 2020 18:28
Operator : CG/JU
Sample : L4711-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0B36

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\SOM-EPA-BG111920MA.M
Title : SVOA CALIBRATION

35	14.583	1908	1914	1923	rVB	480248	774541	48.26%	3.846%
36	14.765	1943	1945	1951	rVB4	3883	5268	0.33%	0.026%
37	14.888	1958	1966	1972	rVV	391370	623306	38.83%	3.095%
38	14.947	1972	1976	1982	rVB2	34362	54098	3.37%	0.269%
39	15.041	1982	1992	2000	rBV	214661	359911	22.42%	1.787%
40	15.200	2013	2019	2023	rVB6	6410	12074	0.75%	0.060%
41	15.282	2028	2033	2038	rVV2	24674	38494	2.40%	0.191%
42	15.352	2040	2045	2047	rBV3	4016	4837	0.30%	0.024%
43	15.605	2083	2088	2092	rVV3	7988	8747	0.54%	0.043%
44	15.664	2095	2098	2099	rVV2	5306	5871	0.37%	0.029%
45	15.705	2103	2105	2112	rVB3	5757	9176	0.57%	0.046%
46	15.869	2127	2133	2139	rBV	591206	912027	56.82%	4.529%
47	15.928	2139	2143	2146	rVV	41152	56490	3.52%	0.280%
48	15.975	2146	2151	2161	rVB2	208889	329185	20.51%	1.635%
49	16.104	2171	2173	2180	rVB7	11345	15805	0.98%	0.078%
50	16.169	2181	2184	2185	rVV2	4524	3965	0.25%	0.020%
51	16.216	2188	2192	2198	rVB2	14355	25592	1.59%	0.127%
52	16.369	2209	2218	2224	rBV5	17357	40739	2.54%	0.202%
53	16.545	2242	2248	2250	rBV6	6398	9998	0.62%	0.050%
54	16.592	2250	2256	2262	rVV6	9460	20396	1.27%	0.101%
55	16.651	2262	2266	2272	rVV3	13505	20223	1.26%	0.100%
56	16.727	2276	2279	2284	rBV5	5829	8750	0.55%	0.043%
57	16.851	2296	2300	2304	rBV5	8570	14375	0.90%	0.071%
58	16.886	2304	2306	2308	rVV3	4440	4923	0.31%	0.024%
59	16.921	2310	2312	2316	rVB4	10423	12946	0.81%	0.064%
60	17.068	2334	2337	2344	rBV6	7340	13045	0.81%	0.065%
61	17.150	2347	2351	2354	rVB4	11974	15842	0.99%	0.079%
62	17.185	2354	2357	2361	rBV4	9652	15398	0.96%	0.076%
63	17.221	2361	2363	2366	rVB4	6113	4447	0.28%	0.022%
64	17.268	2366	2371	2380	rVB	33250	59846	3.73%	0.297%
65	17.350	2380	2385	2387	rBV3	15941	24629	1.53%	0.122%
66	17.444	2395	2401	2406	rVB	28125	47473	2.96%	0.236%
67	17.526	2413	2415	2419	rVB4	6211	7416	0.46%	0.037%
68	17.620	2423	2431	2435	rBV	490820	800323	49.86%	3.974%
69	17.661	2435	2438	2443	rVB	587899	830802	51.76%	4.125%
70	17.720	2443	2448	2453	rBV2	580202	875824	54.57%	4.349%
71	17.802	2458	2462	2468	rVB3	22084	37238	2.32%	0.185%

LSC Area Percent Report

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Operator : CG/JU
Sample : L4711-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0B36

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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Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Title : SVOA CALIBRATION

72	17.879	2473	2475	2482	rBV3	10976	18250	1.14%	0.091%
73	17.932	2482	2484	2486	rVB2	7350	4415	0.28%	0.022%
74	17.979	2488	2492	2494	rBV5	15056	23122	1.44%	0.115%
75	18.008	2494	2497	2503	rVB2	50193	77735	4.84%	0.386%
76	18.131	2515	2518	2525	rBV6	17799	26724	1.66%	0.133%
77	18.196	2525	2529	2532	rBV4	21672	28118	1.75%	0.140%
78	18.390	2561	2562	2563	rBV	7590	3849	0.24%	0.019%
79	18.408	2563	2565	2569	rVV5	9661	11002	0.69%	0.055%
80	18.490	2575	2579	2583	rBV2	71363	100484	6.26%	0.499%
81	18.537	2583	2587	2596	rVB	95752	149579	9.32%	0.743%
82	18.619	2596	2601	2605	rVB	23654	36227	2.26%	0.180%
83	18.684	2605	2612	2616	rBV3	104396	209760	13.07%	1.042%
84	18.977	2658	2662	2665	rBV2	63813	92139	5.74%	0.458%
85	19.019	2665	2669	2679	rVB	151640	227494	14.17%	1.130%
86	19.260	2706	2710	2715	rBV2	37326	52121	3.25%	0.259%
87	19.389	2729	2732	2735	rBV2	39189	53662	3.34%	0.266%
88	19.530	2752	2756	2762	rBV3	54810	82234	5.12%	0.408%
89	19.653	2771	2777	2783	rVB	1147854	1605047	100.00%	7.970%
90	19.982	2827	2833	2836	rBV2	815785	1374077	85.61%	6.823%
91	20.017	2836	2839	2845	rVV	800138	1070164	66.67%	5.314%
92	20.076	2846	2849	2854	rVB	50449	68099	4.24%	0.338%
93	20.182	2864	2867	2872	rVB	61564	88298	5.50%	0.438%
94	20.493	2918	2920	2925	rVB	143282	170024	10.59%	0.844%
95	20.593	2935	2937	2941	rVB	68790	71264	4.44%	0.354%
96	21.269	3049	3052	3057	rVB2	90786	131089	8.17%	0.651%
97	21.510	3090	3093	3098	rVB2	210878	279905	17.44%	1.390%
98	21.892	3152	3158	3164	rBV2	490338	1330506	82.90%	6.607%
99	21.950	3165	3168	3178	rVB	402732	669192	41.69%	3.323%
100	24.224	3549	3555	3564	rBV2	241724	760060	47.35%	3.774%

Sum of corrected areas: 20139297

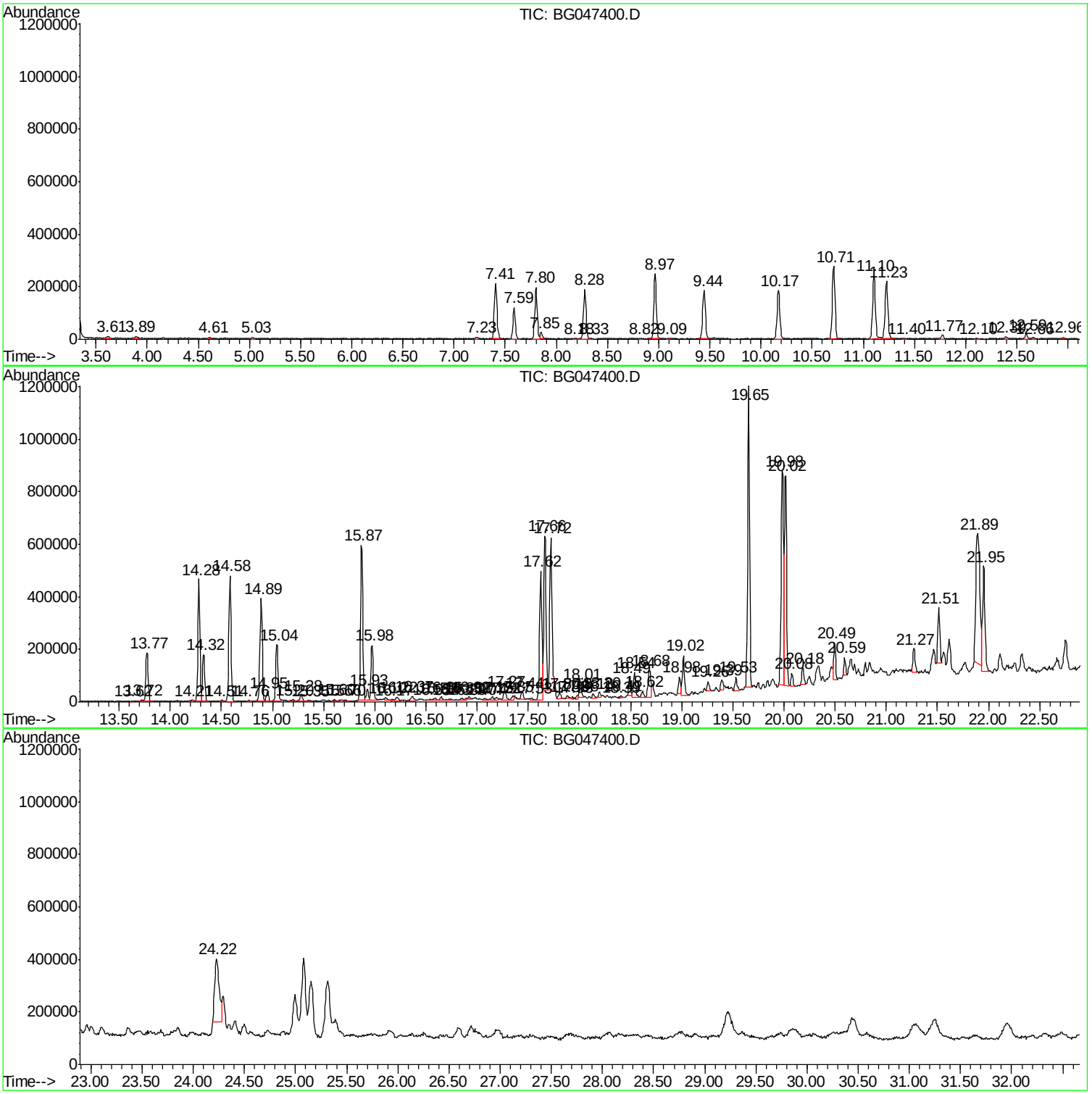
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112120\
Data File : BG047400.D
Acq On : 21 Nov 2020 18:28
Operator : CG/JU
Sample : L4711-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0B36

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112120\
Data File : BG047400.D
Acq On : 21 Nov 2020 18:28
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Instrument : BNA_G
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

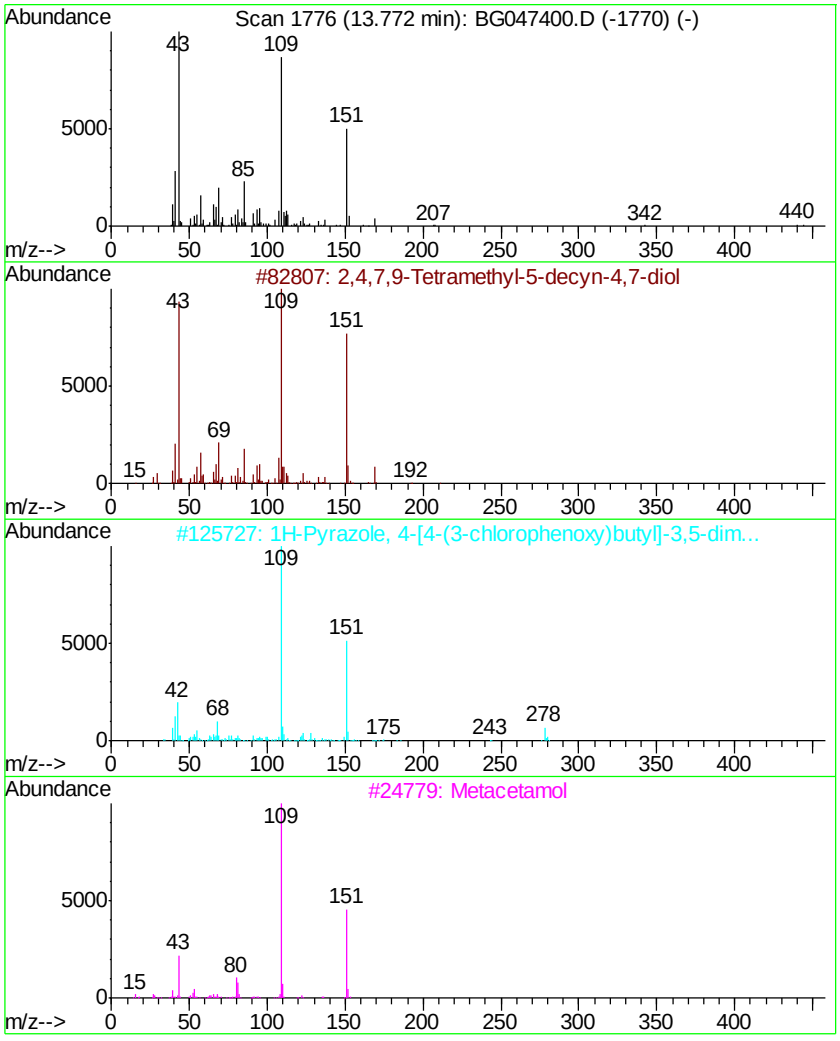
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

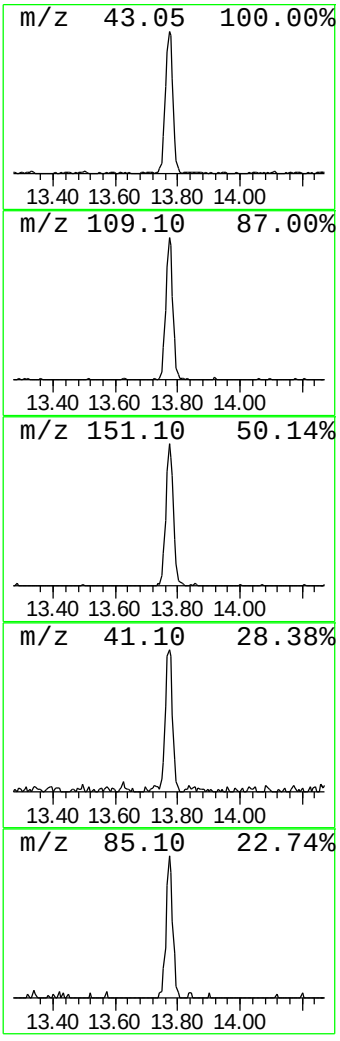
****

Peak Number 1 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.77	9.96 ng/ul	310373	Acenaphthene-d10	14.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	86
2	1H-Pyrazole, 4-[4-(3-chloropheno...	278	C15H19ClN2O	1000302-33-9	59
3	Metacetamol	151	C8H9NO2	000621-42-1	58
4	Acetaminophen	151	C8H9NO2	000103-90-2	58
5	Tiocarlide	400	C23H32N2O2S	000910-86-1	53





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112120\
Data File : BG047400.D
Acq On : 21 Nov 2020 18:28
Operator : CG/JU
Sample : L4711-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0B36

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

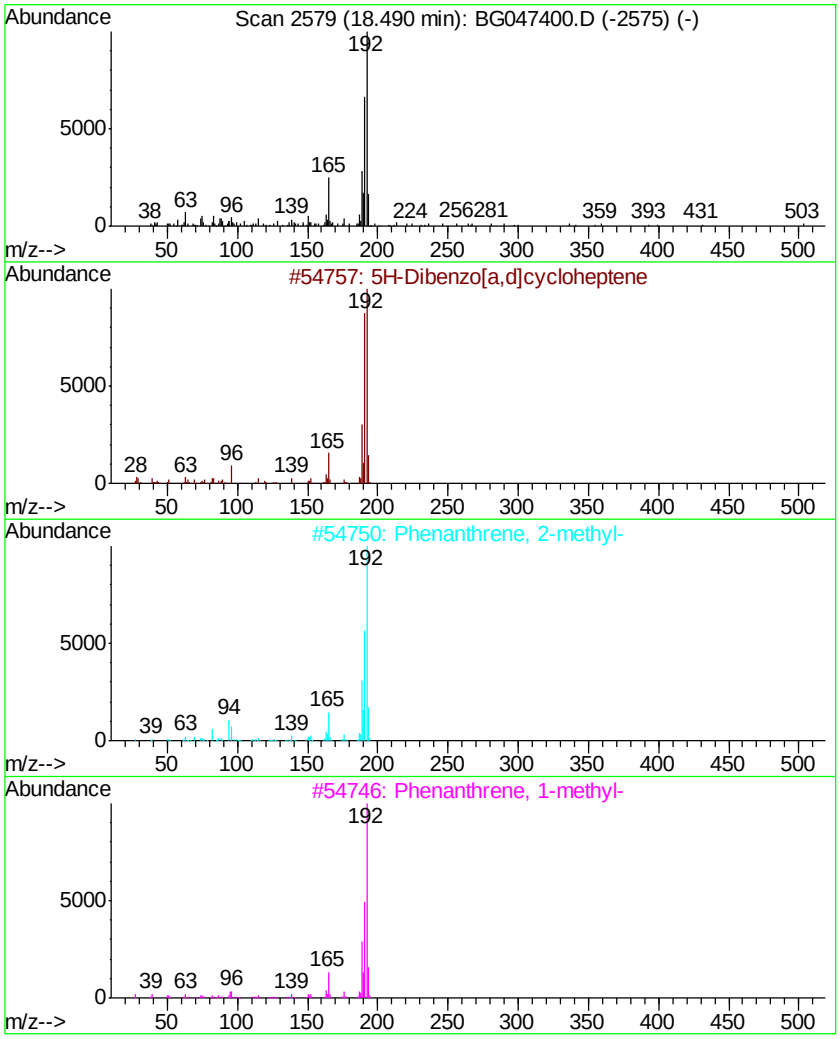
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

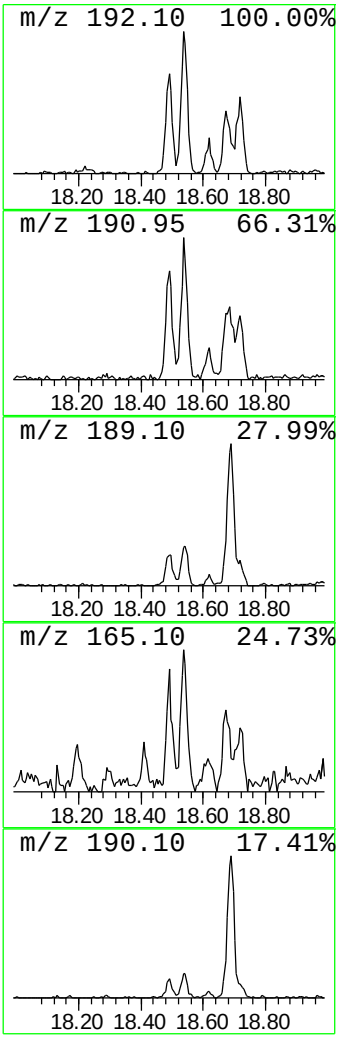
****

Peak Number 2 5H-Dibenzo[a,d]cycloheptene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.49	2.51 ng/ul	100484	Phenanthrene-d10	17.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	91
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	89
4	Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	87
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	83





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112120\
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Sample : L4711-01
Misc :
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Instrument : BNA_G
ClientSampleId : C0B36

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

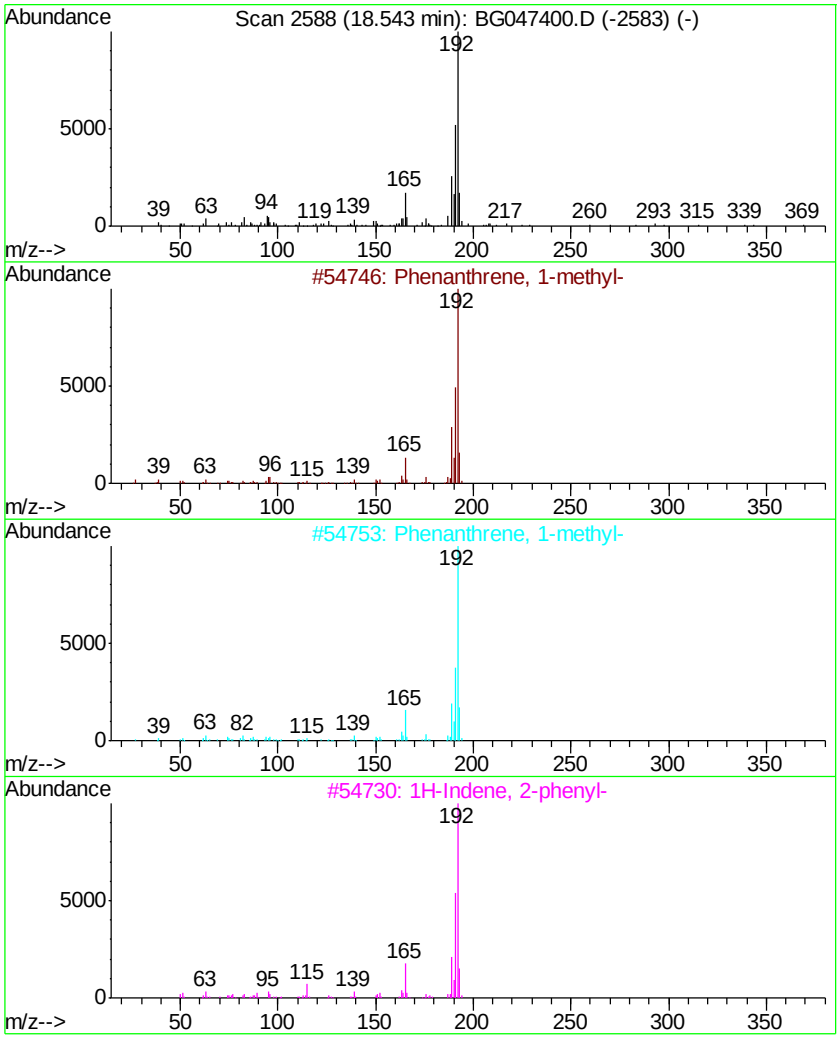
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

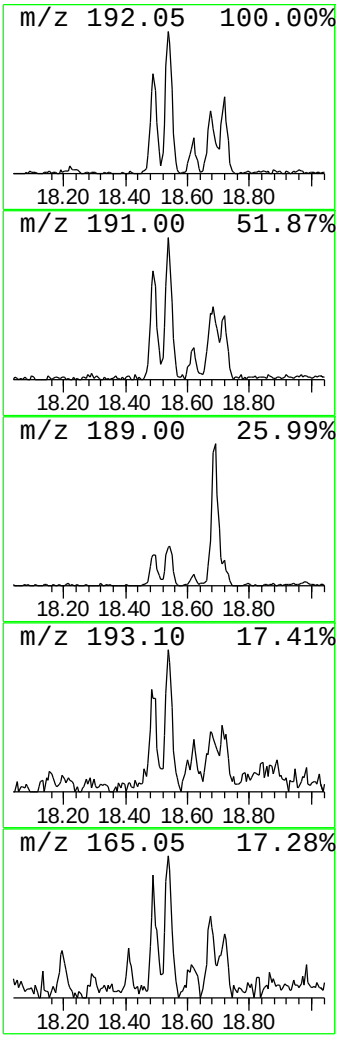
****

Peak Number 3 Phenanthrene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.54	3.74 ng/ul	149579	Phenanthrene-d10	17.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
3	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	95
4	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	94
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	94





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112120\
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Misc :
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Instrument : BNA_G
ClientSampleId : C0B36

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

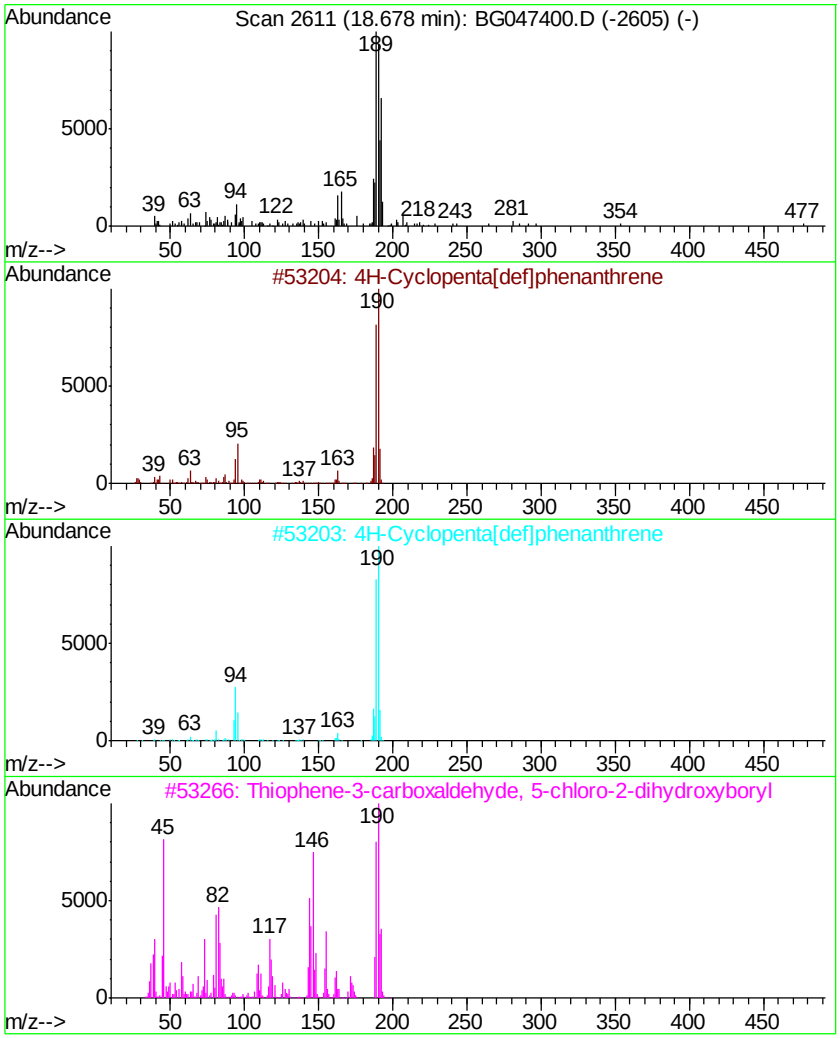
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

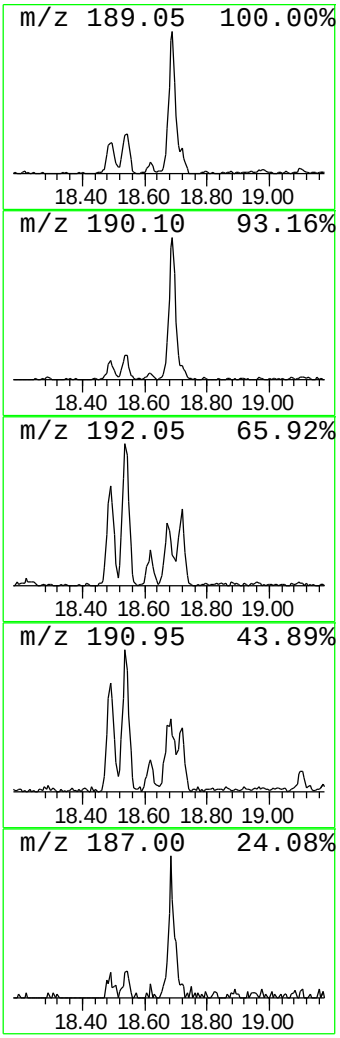
****

Peak Number 4 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.68	5.24 ng/ul	209760	Phenanthrene-d10	17.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	89
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
3	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	53
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	50
5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	38





Library Search Compound Report

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Data File : BG047400.D
Acq On : 21 Nov 2020 18:28
Operator : CG/JU
Sample : L4711-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0B36

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

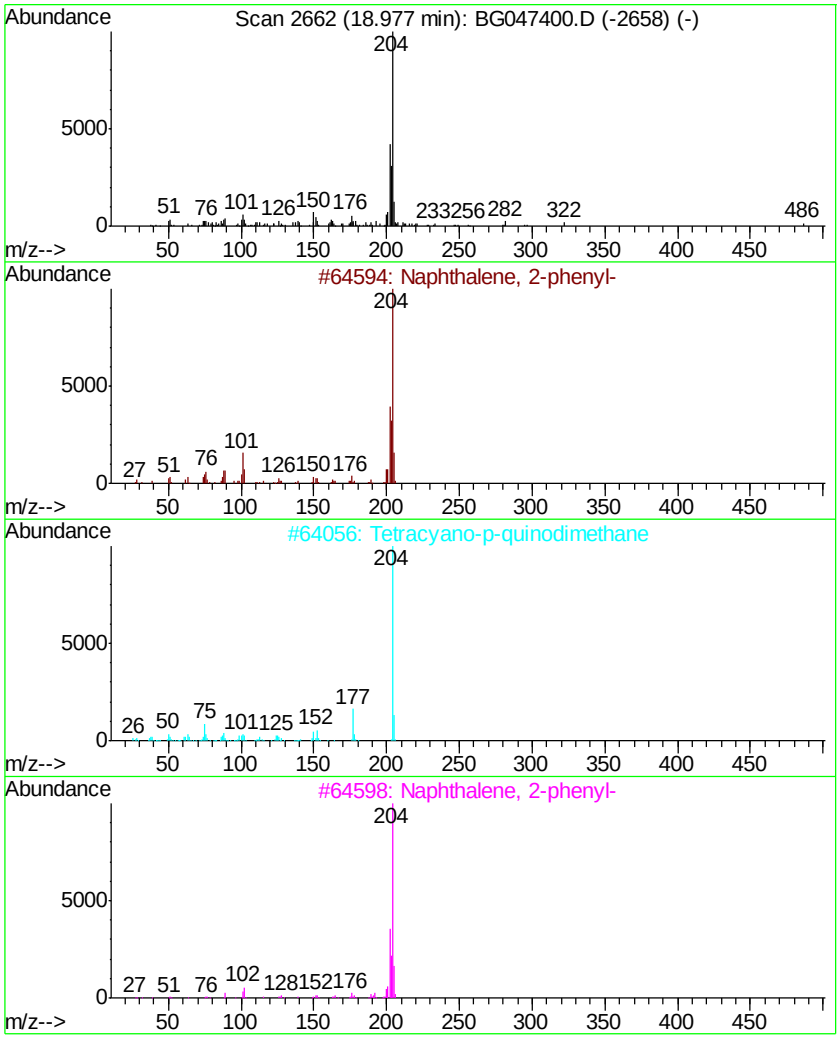
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

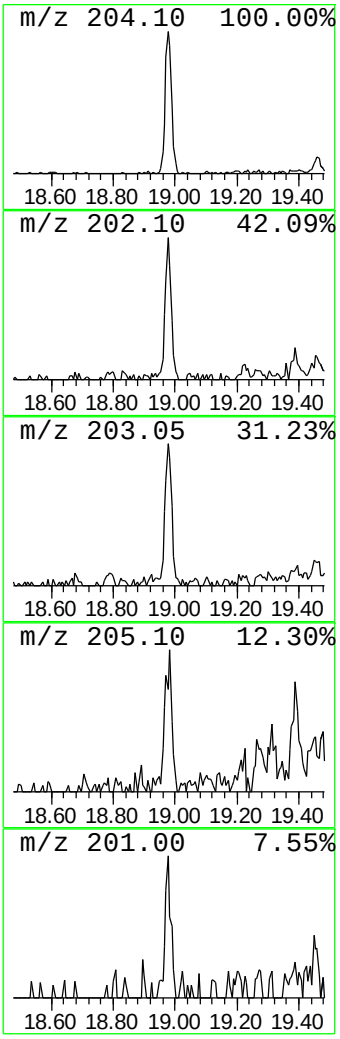
****

Peak Number 5 Naphthalene, 2-phenyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.98	2.30 ng/ul	92139	Phenanthrene-d10	17.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	93
2	Tetracyano-p-quinodimethane	204	C12H4N4	001518-16-7	55
3	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	53
4	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	53
5	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	53





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Acq On : 21 Nov 2020 18:28
Operator : CG/JU
Sample : L4711-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0B36

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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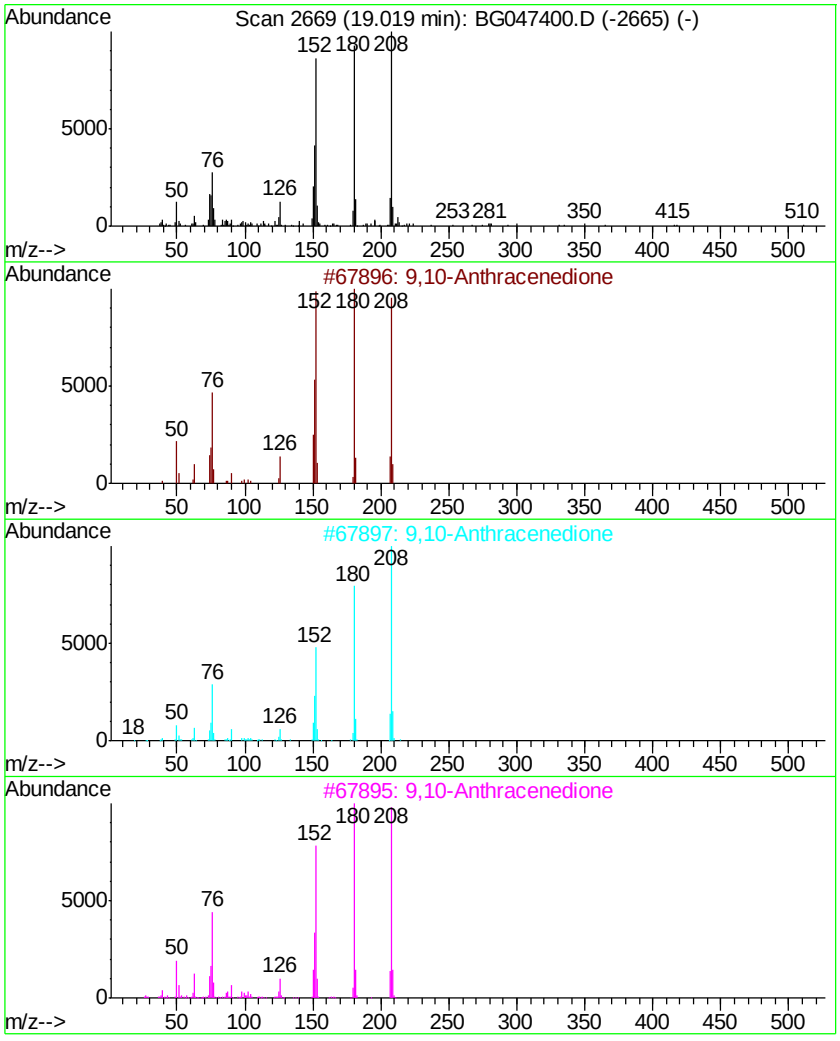
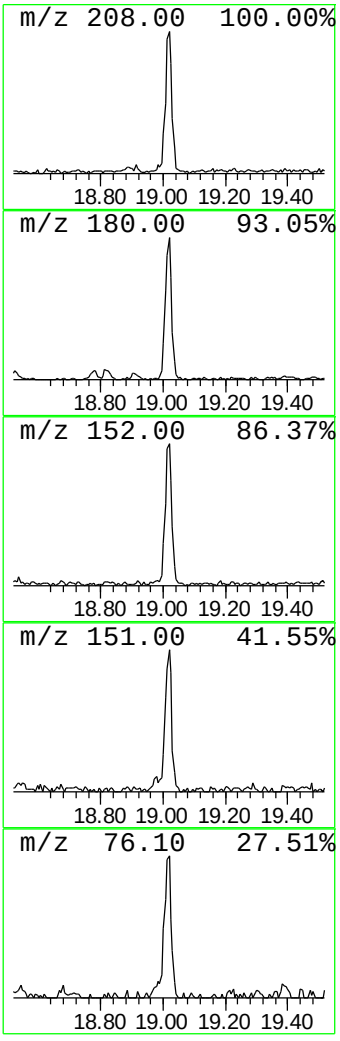
Peak Number 6 9,10-Anthracenedione Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.02	5.69 ng/ul	227494	Phenanthrene-d10	17.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	97
2	9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
3	9,10-Anthracenedione	208	C14H8O2	000084-65-1	93
4	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	70
5	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	64





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Data File : BG047400.D
Acq On : 21 Nov 2020 18:28
Operator : CG/JU
Sample : L4711-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0B36

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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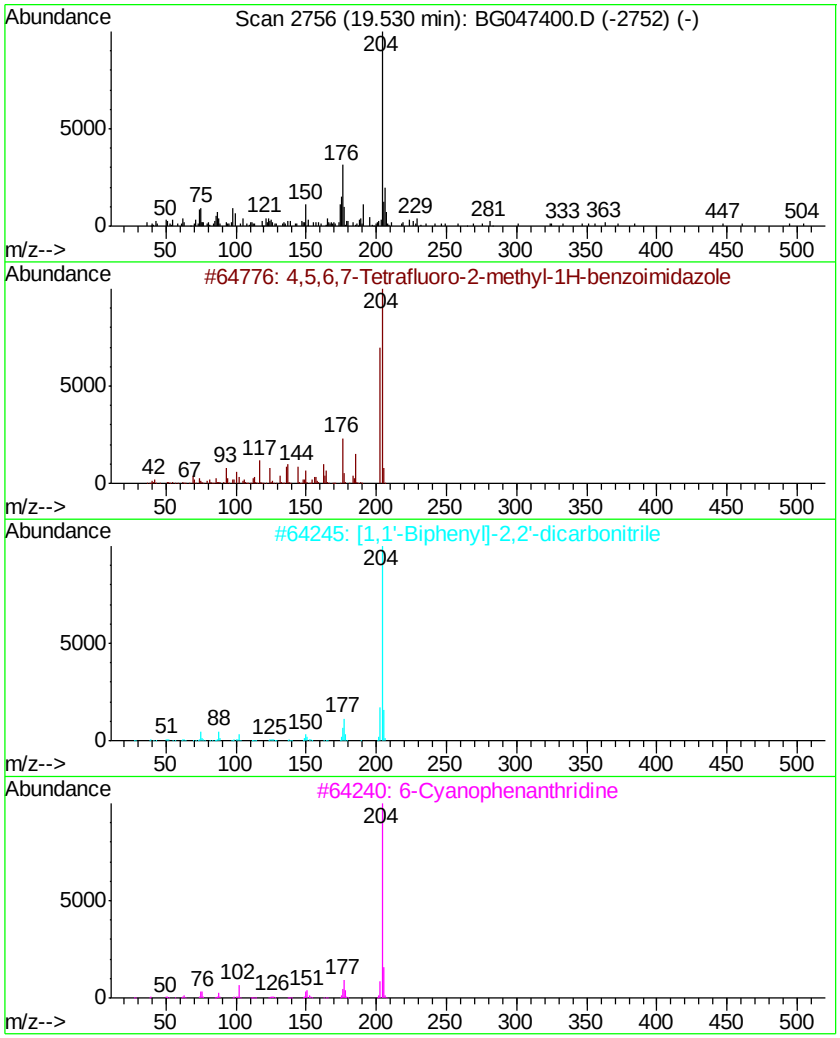
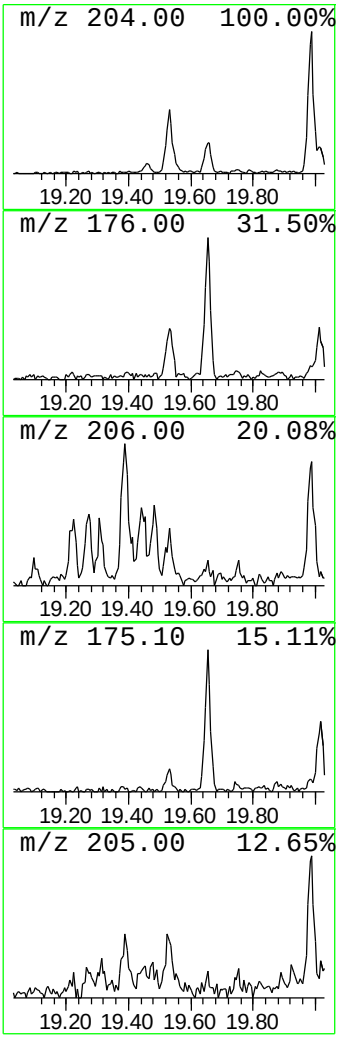
Peak Number 7 4,5,6,7-Tetrafluoro-2-methy... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.53	2.06 ng/ul	82234	Phenanthrene-d10	17.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4,5,6,7-Tetrafluoro-2-methyl-1H-...	204	C8H4F4N2	081430-75-3	59
2	[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	52
3	6-Cyanophenanthridine	204	C14H8N2	002176-28-5	52
4	[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	50
5	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	50





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Data File : BG047400.D
Acq On : 21 Nov 2020 18:28
Operator : CG/JU
Sample : L4711-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0B36

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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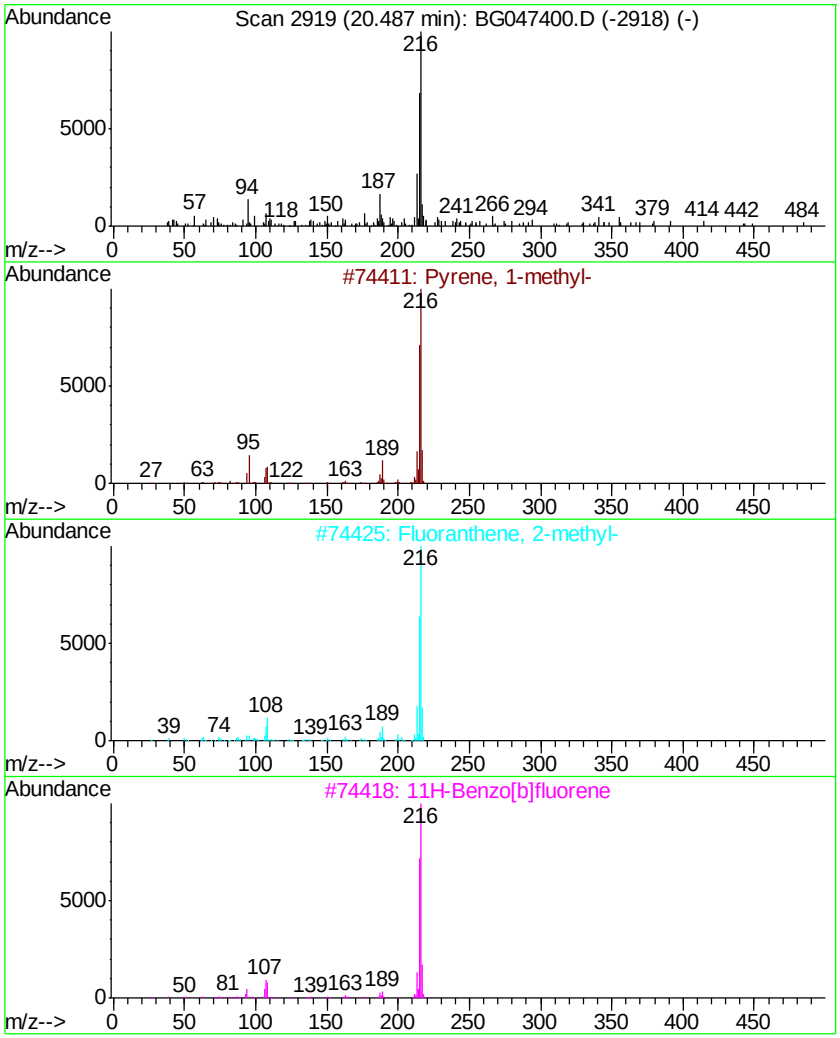
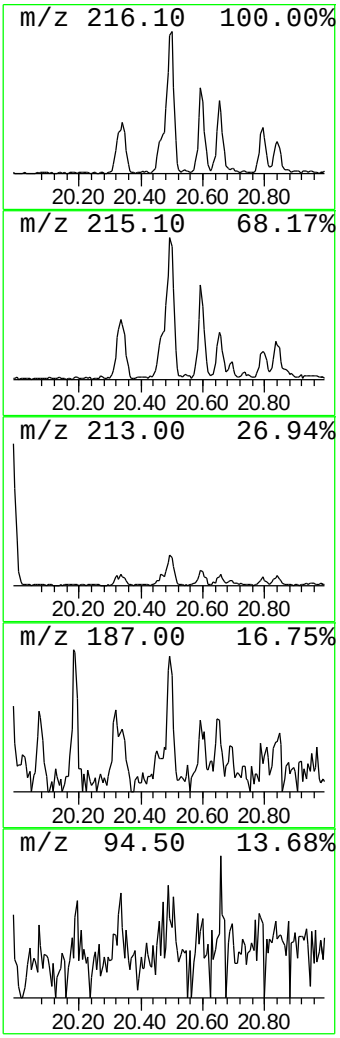
Peak Number 8 Pyrene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.49	2.56 ng/ul	170024	Chrysene-d12	21.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
3	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
4	Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	87





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Data File : BG047400.D
Acq On : 21 Nov 2020 18:28
Operator : CG/JU
Sample : L4711-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0B36

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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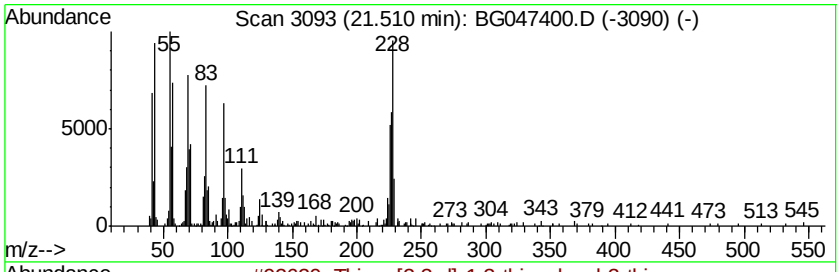
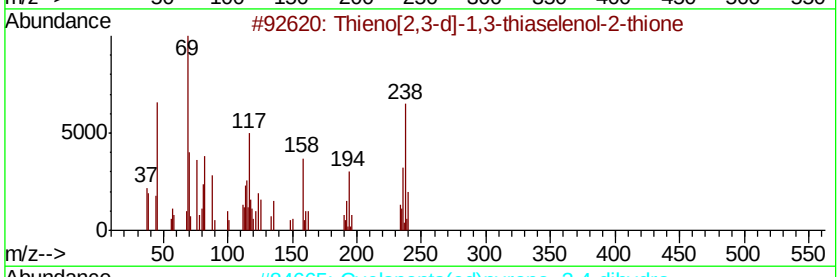
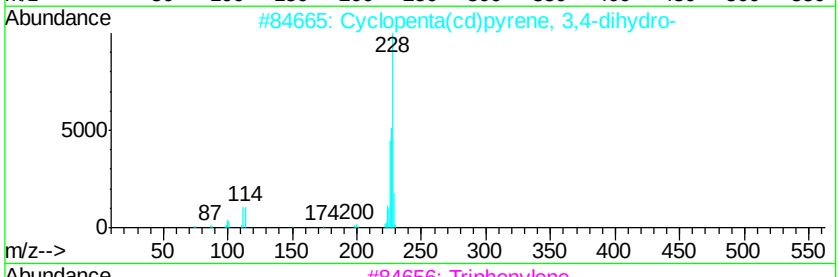
Peak Number 9 Thieno[2,3-d]-1,3-thiaselen... Concentration Rank 4

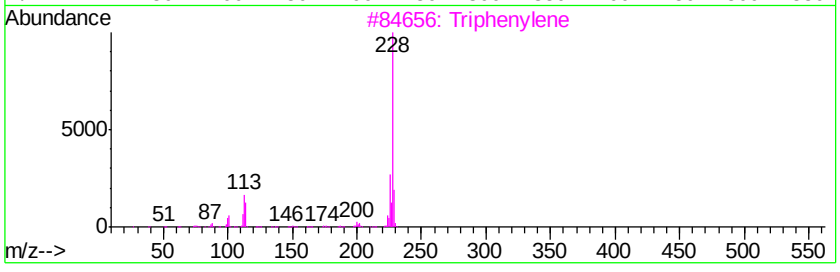
R.T.	EstConc	Area	Relative to ISTD	R.T.
21.51	4.21 ng/ul	279905	Chrysene-d12	21.90

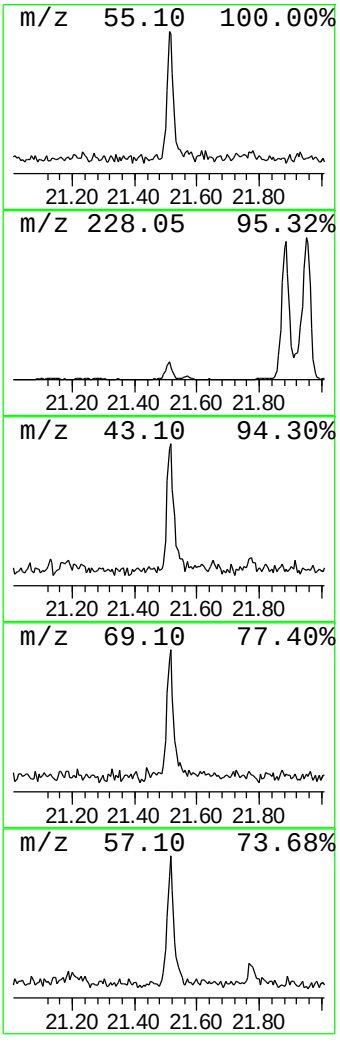
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Thieno[2,3-d]-1,3-thiaselenol-2-...	238	C5H2S3Se	081803-09-0	53
2	Cyclopenta(cd)pyrene, 3,4-dihydro-	228	C18H12	025732-74-5	43
3	Triphenylene	228	C18H12	000217-59-4	38
4	Naphthacene	228	C18H12	000092-24-0	38
5	Thiazolo[5,4-f]quinoline, 2,7,9-...	228	C13H12N2S	038463-47-7	35











Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG112120\
Data File : BG047400.D
Acq On : 21 Nov 2020 18:28
Operator : CG/JU
Sample : L4711-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0B36

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2,4,7,9-Tetrameth...	13.77	10.0	ng/ul	310373	3	14.89	623306	20.0
5H-Dibenzo[a,d]cy...	18.49	2.5	ng/ul	100484	4	17.62	800323	20.0
Phenanthrene, 1-m...	18.54	3.7	ng/ul	149579	4	17.62	800323	20.0
4H-Cyclopenta[def...	18.68	5.2	ng/ul	209760	4	17.62	800323	20.0
Naphthalene, 2-ph...	18.98	2.3	ng/ul	92139	4	17.62	800323	20.0
9,10-Anthracenedione	19.02	5.7	ng/ul	227494	4	17.62	800323	20.0
4,5,6,7-Tetrafluo...	19.53	2.1	ng/ul	82234	4	17.62	800323	20.0
Pyrene, 1-methyl-	20.49	2.6	ng/ul	170024	5	21.90	1330510	20.0
Thieno[2,3-d]-1,3...	21.51	4.2	ng/ul	279905	5	21.90	1330510	20.0