LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG110515\
Data File : BG019494.D
Acq On : 5 Nov 2015 16:06
Operator : UM/NP
Sample : G4273-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AP4

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:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102815.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	2.893	6	9	15	rVB5	11650	17899	1.02%	0.073%
2	3.128	43	49	57	rBV	93503	184319	10.45%	0.755%
3	3.205	57	62	74	rVB	642389	933729	52.95%	3.824%
4	3.481	107	109	113	rBV4	5647	7077	0.40%	0.029%
5	4.033	199	203	209	rBV3	3814	6164	0.35%	0.025%
6	4.609	296	301	307	rVB5	3583	6594	0.37%	0.027%
7	5.220	401	405	411	rVB3	6590	9579	0.54%	0.039%
8	7.323	752	763	773	rBV	137803	248093	14.07%	1.016%
9	7.494	785	792	800	rVB	101832	163048	9.25%	0.668%
10	7.705	821	828	837	rBV2	125035	222689	12.63%	0.912%
11	8.181	897	909	915	rBV	153778	261354	14.82%	1.070%
12	8.557	970	973	980	rVB5	6264	9936	0.56%	0.041%
13	8.880	1022	1028	1036	rVB	173994	304289	17.26%	1.246%
14	9.350	1101	1108	1115	rBV	136474	238545	13.53%	0.977%
15	10.079	1225	1232	1241	rVB2	125677	224848	12.75%	0.921%
16	10.619	1317	1324	1332	rBV2	184733	324700	18.41%	1.330%
17	11.007	1383	1390	1396	rBV	216957	395315	22.42%	1.619%
18	11.060	1396	1399	1404	rVB2	16517	21049	1.19%	0.086%
19	11.137	1405	1412	1420	rVV	170291	299535	16.99%	1.227%
20	12.576	1653	1657	1661	rVB2	4222	6993	0.40%	0.029%
21	12.652	1663	1670	1677	rBV	55978	93710	5.31%	0.384%
22	12.864	1701	1706	1714	rVB	43528	73805	4.19%	0.302%
23	13.175	1753	1759	1767	rVV7	4337	9794	0.56%	0.040%
24	13.410	1794	1799	1804	rVB4	6919	9917	0.56%	0.041%
25	13.651	1834	1840	1842	rVV2	6113	8741	0.50%	0.036%
26	13.839	1867	1872	1882	rVB	14618	27704	1.57%	0.113%
27	13.968	1888	1894	1895	rBV2	21163	26182	1.48%	0.107%
28	14.133	1915	1922	1926	rVV4	32699	63139	3.58%	0.259%
29	14.204	1926	1934	1938	rVV	383087	577477	32.75%	2.365%
30	14.251	1938	1942	1949	rVB	170083	242841	13.77%	0.995%
31	14.368	1956	1962	1966	rBV3	18299	30592	1.73%	0.125%
32	14.503	1979	1985	1994	rBV	372385	599985	34.03%	2.457%
33	14.809	2026	2037	2043	rBV2	408986	677819	38.44%	2.776%
34	14.873	2043	2048	2056	rVB	234333	364988	20.70%	1.495%

LSC Area Percent Report

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Data File : BG019494.D
Acq On : 5 Nov 2015 16:06
Operator : UM/NP
Sample : G4273-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AP4

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:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102815.M
Title : SVOA CALIBRATION

35	14.991	2062	2068	2074	rBV	198184	318408	18.06%	1.304%
36	15.114	2081	2089	2093	rBV4	14808	26391	1.50%	0.108%
37	15.202	2098	2104	2111	rVB	160061	240500	13.64%	0.985%
38	15.273	2111	2116	2120	rBV2	15192	21303	1.21%	0.087%
39	15.414	2134	2140	2145	rBV6	13264	23890	1.35%	0.098%
40	15.467	2145	2149	2153	rVB3	11333	16320	0.93%	0.067%
41	15.590	2161	2170	2172	rBV6	13152	27404	1.55%	0.112%
42	15.625	2173	2176	2179	rVB2	13946	16284	0.92%	0.067%
43	15.796	2197	2205	2210	rVV	519144	790576	44.84%	3.238%
44	15.849	2210	2214	2219	rVV	303803	441747	25.05%	1.809%
45	15.902	2219	2223	2233	rVB	180720	266044	15.09%	1.090%
46	15.978	2233	2236	2238	rBV3	9506	10457	0.59%	0.043%
47	16.013	2238	2242	2248	rVV4	35433	58613	3.32%	0.240%
48	16.101	2253	2257	2260	rVV2	28345	47646	2.70%	0.195%
49	16.142	2260	2264	2269	rVB	54443	83536	4.74%	0.342%
50	16.289	2278	2289	2296	rBV2	66719	158504	8.99%	0.649%
51	16.383	2301	2305	2309	rVB5	14413	21533	1.22%	0.088%
52	16.489	2319	2323	2333	rVV6	37365	83123	4.71%	0.340%
53	16.642	2345	2349	2355	rVB3	29359	48962	2.78%	0.201%
54	16.847	2374	2384	2388	rBV2	51602	104684	5.94%	0.429%
55	16.894	2389	2392	2398	rVB5	18695	28134	1.60%	0.115%
56	17.000	2406	2410	2415	rVB2	20908	33382	1.89%	0.137%
57	17.071	2415	2422	2426	rBV7	21938	54548	3.09%	0.223%
58	17.276	2452	2457	2459	rBV3	31057	49086	2.78%	0.201%
59	17.370	2468	2473	2479	rVB	68432	110984	6.29%	0.455%
60	17.553	2497	2504	2507	rBV	604462	978212	55.48%	4.006%
61	17.594	2507	2511	2516	rVV	955123	1384928	78.54%	5.672%
62	17.647	2516	2520	2523	rVV	597432	906838	51.43%	3.714%
63	17.682	2523	2526	2531	rVB	591455	805152	45.66%	3.298%
64	17.817	2546	2549	2554	rVV3	11427	18776	1.06%	0.077%
65	17.946	2566	2571	2580	rBV	85748	139847	7.93%	0.573%
66	18.011	2580	2582	2587	rVB5	12441	14485	0.82%	0.059%
67	18.058	2587	2590	2595	rBV3	19772	30595	1.74%	0.125%
68	18.263	2623	2625	2630	rVB4	11326	12036	0.68%	0.049%
69	18.369	2642	2643	2646	rBV3	7723	7992	0.45%	0.033%
70	18.416	2646	2651	2656	rBV	97346	150213	8.52%	0.615%
71	18.469	2656	2660	2667	rVB	125475	182482	10.35%	0.747%

LSC Area Percent Report

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Acq On : 5 Nov 2015 16:06
Operator : UM/NP
Sample : G4273-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AP4

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:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102815.M
Title : SVOA CALIBRATION

72	18.545	2669	2673	2678	rVB	88398	124651	7.07%	0.511%
73	18.622	2679	2686	2689	rBV	176913	305472	17.32%	1.251%
74	18.757	2706	2709	2712	rVB4	12458	15534	0.88%	0.064%
75	18.910	2730	2735	2741	rBV	82361	114545	6.50%	0.469%
76	19.151	2773	2776	2782	rVB6	23813	31464	1.78%	0.129%
77	19.245	2788	2792	2797	rVB3	16880	22668	1.29%	0.093%
78	19.321	2800	2805	2811	rBV3	43542	79677	4.52%	0.326%
79	19.456	2825	2828	2831	rBV5	17948	27183	1.54%	0.111%
80	19.591	2845	2851	2856	rVB	843037	1173403	66.55%	4.806%
81	19.685	2863	2867	2869	rVB4	19155	24020	1.36%	0.098%
82	19.738	2872	2876	2881	rVB	55880	76943	4.36%	0.315%
83	19.832	2886	2892	2900	rVV3	28767	67759	3.84%	0.278%
84	19.920	2901	2907	2910	rVV2	892612	1408368	79.87%	5.768%
85	19.950	2910	2912	2920	rVV	571927	753420	42.73%	3.086%
86	20.014	2920	2923	2929	rVB3	46754	72716	4.12%	0.298%
87	20.126	2937	2942	2947	rVB	63584	104050	5.90%	0.426%
88	20.261	2960	2965	2972	rBV4	57412	143669	8.15%	0.588%
89	20.437	2991	2995	3000	rVB	189618	265000	15.03%	1.085%
90	20.531	3007	3011	3016	rBV	186952	267422	15.17%	1.095%
91	20.737	3042	3046	3049	rBV	36138	55415	3.14%	0.227%
92	21.395	3155	3158	3161	rBV3	31335	41767	2.37%	0.171%
93	21.436	3162	3165	3170	rVB	43776	70209	3.98%	0.288%
94	21.689	3205	3208	3213	rVB2	57938	80218	4.55%	0.329%
95	21.824	3223	3231	3236	rBV2	807846	1763297	100.00%	7.222%
96	21.877	3236	3240	3248	rVB	229787	397106	22.52%	1.626%
97	24.110	3613	3620	3626	rBV	104970	310363	17.60%	1.271%
98	24.850	3744	3746	3753	rBV3	25202	47593	2.70%	0.195%
99	24.944	3753	3762	3769	rBV	373646	1013612	57.48%	4.151%
100	25.173	3792	3801	3811	rBV	393068	1152196	65.34%	4.719%

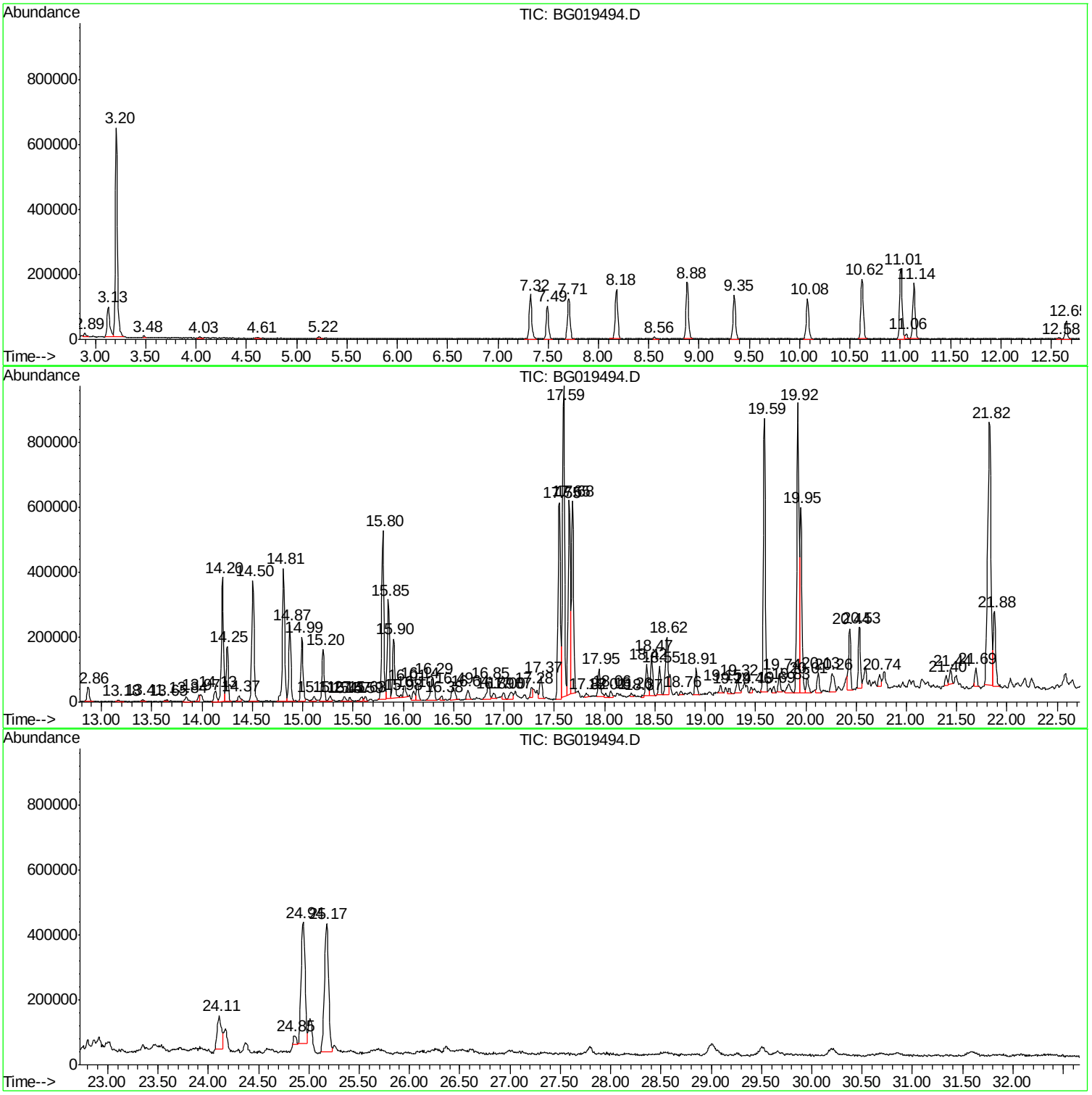
Sum of corrected areas: 24415804

Data Path : Z:\HPCHEM1\BNA G\DATA\BG110515\
Data File : BG019494.D
Acq On : 5 Nov 2015 16:06
Operator : UM/NP
Sample : G4273-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AP4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102815.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG110515\
Data File : BG019494.D
Acq On : 5 Nov 2015 16:06
Operator : UM/NP
Sample : G4273-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102815.M
Quant Title : SVOA CALIBRATION

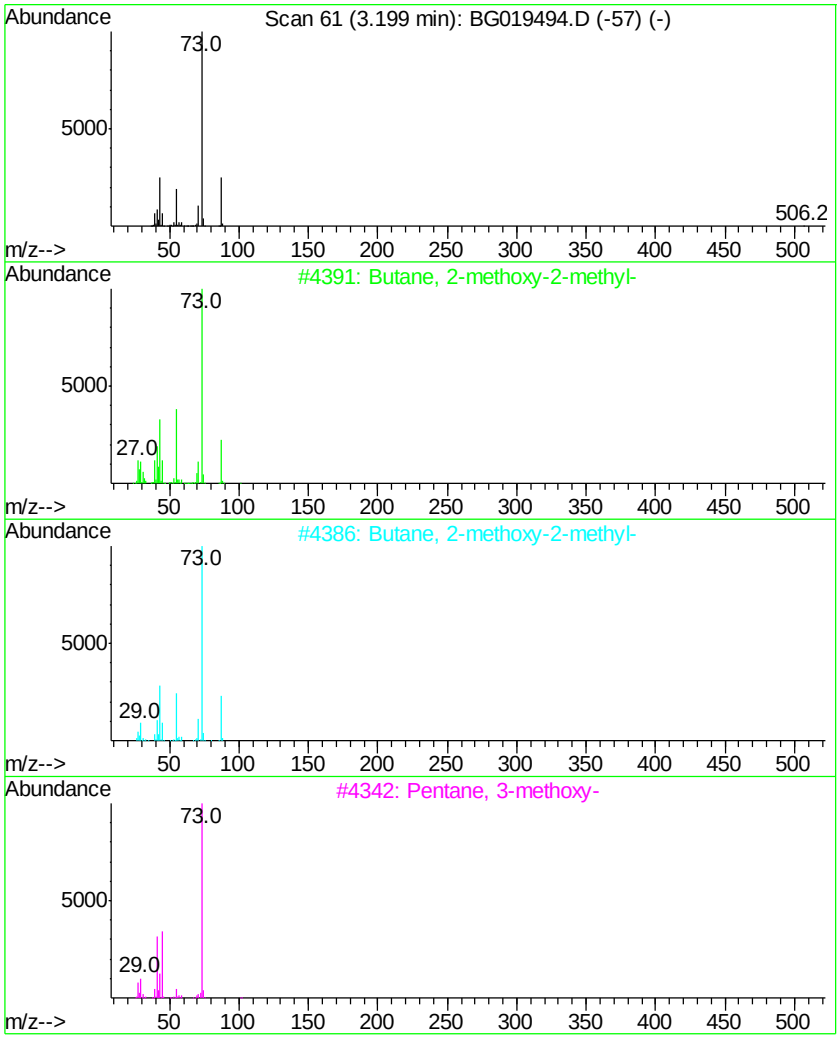
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

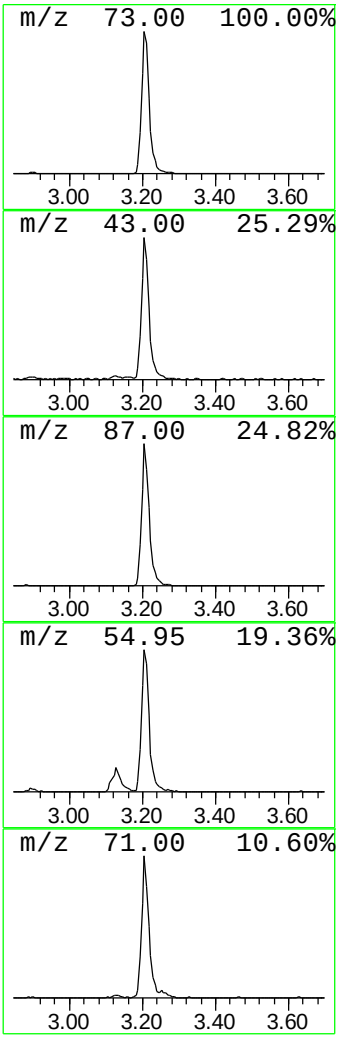
****

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.20	71.45 ng/ul	933729	1,4-Dichlorobenzene-d4	8.18

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9





Data Path : Z:\HPCHEM1\BNA G\DATA\BG110515\
Data File : BG019494.D
Acq On : 5 Nov 2015 16:06
Operator : UM/NP
Sample : G4273-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AP4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102815.M
Quant Title : SVOA CALIBRATION

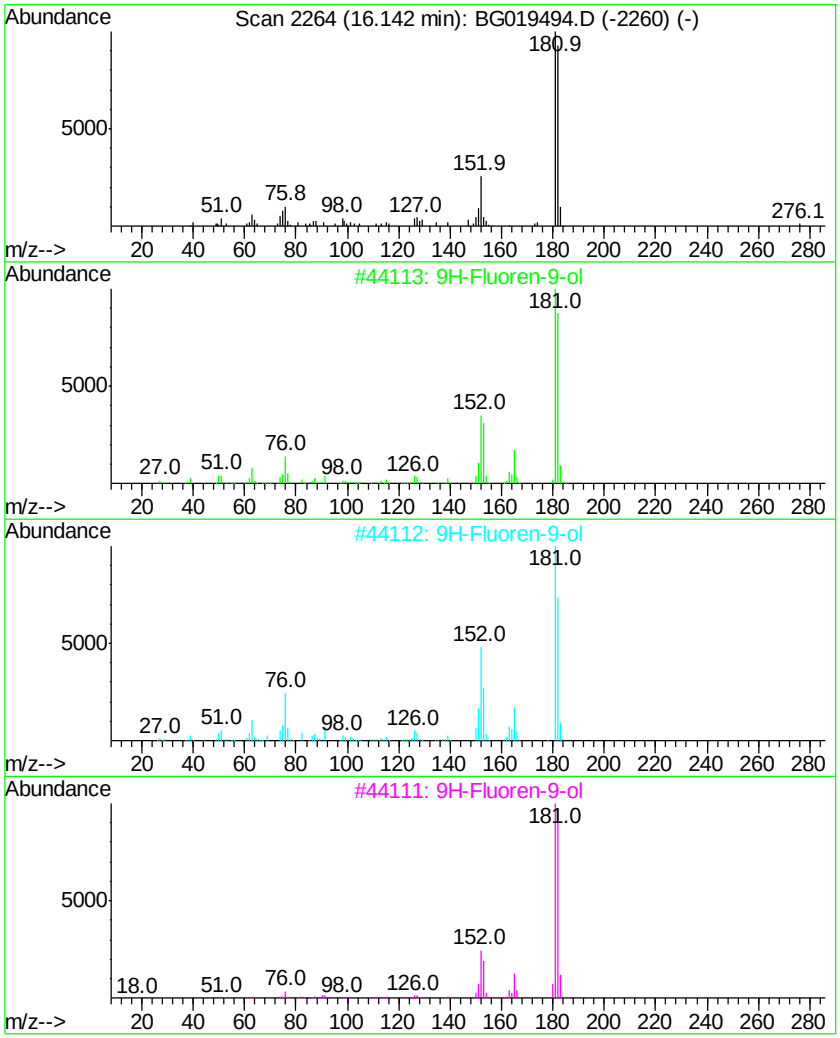
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

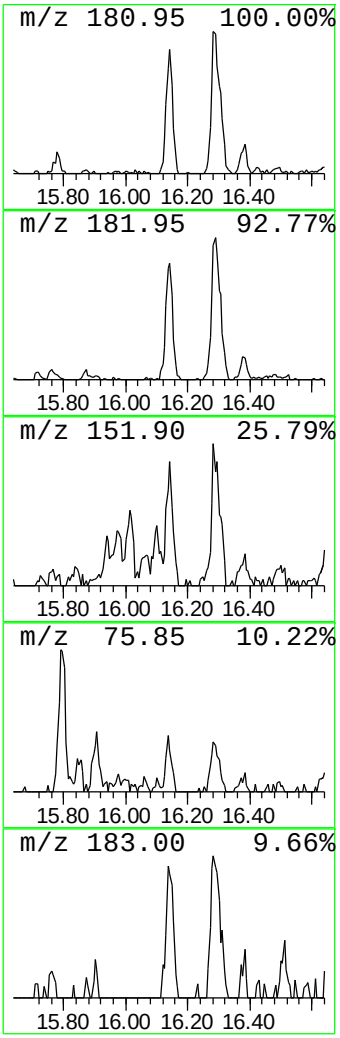
****

Peak Number 3 9H-Fluoren-9-ol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.14	2.46 ng/ul	83536	Acenaphthene-d10	14.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	94
2	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	83
3	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	78
4	3-(1-Naphthyl)acrylaldehyde	182	C13H10O	001504-73-0	78
5	2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	72





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG110515\
Data File : BG019494.D
Acq On : 5 Nov 2015 16:06
Operator : UM/NP
Sample : G4273-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AP4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102815.M
Quant Title : SVOA CALIBRATION

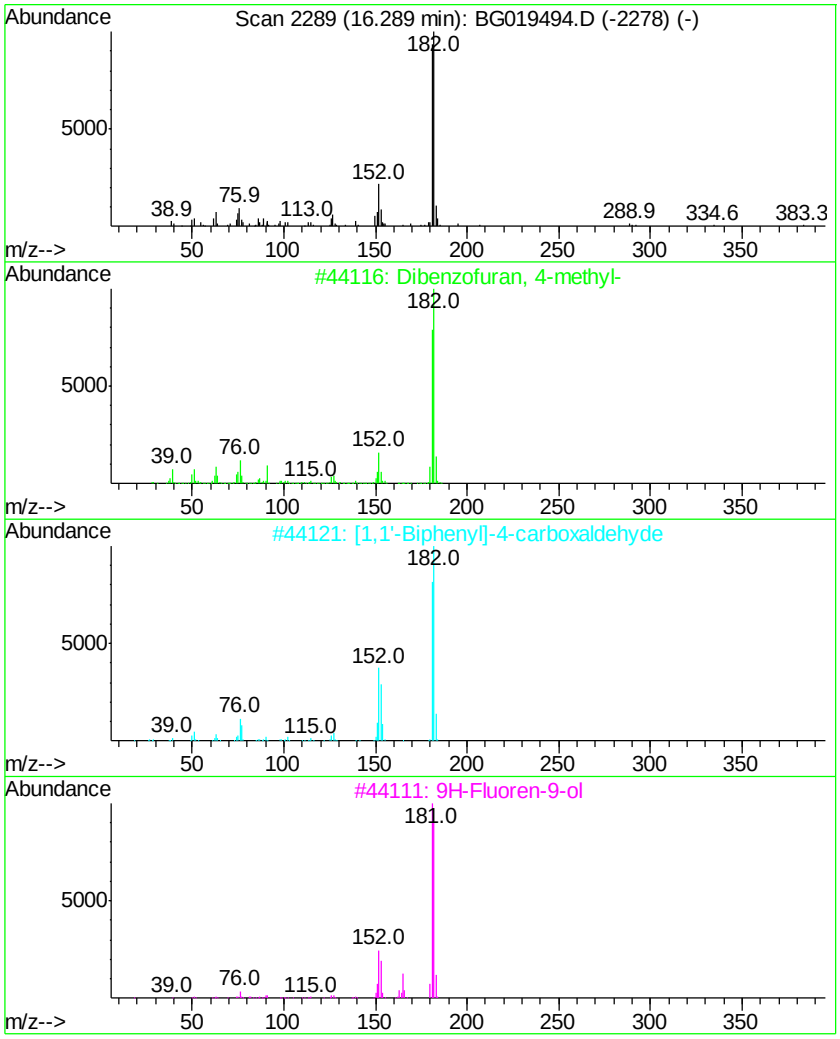
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

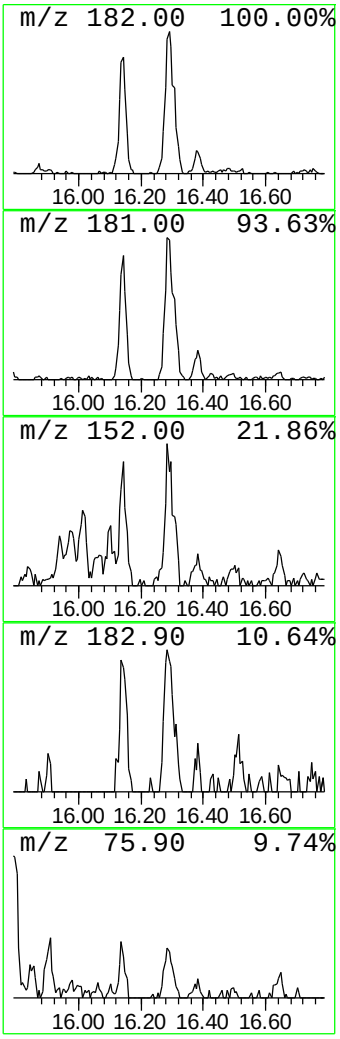
****

Peak Number 4 Dibenzofuran, 4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.29	3.24 ng/ul	158504	Phenanthrene-d10	17.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	76
2	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	74
3	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	72
4	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	59
5	2-Fluorenamine	181	C13H11N	000153-78-6	55





Data Path : Z:\HPCHEM1\BNA G\DATA\BG110515\
Data File : BG019494.D
Acq On : 5 Nov 2015 16:06
Operator : UM/NP
Sample : G4273-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AP4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102815.M
Quant Title : SVOA CALIBRATION

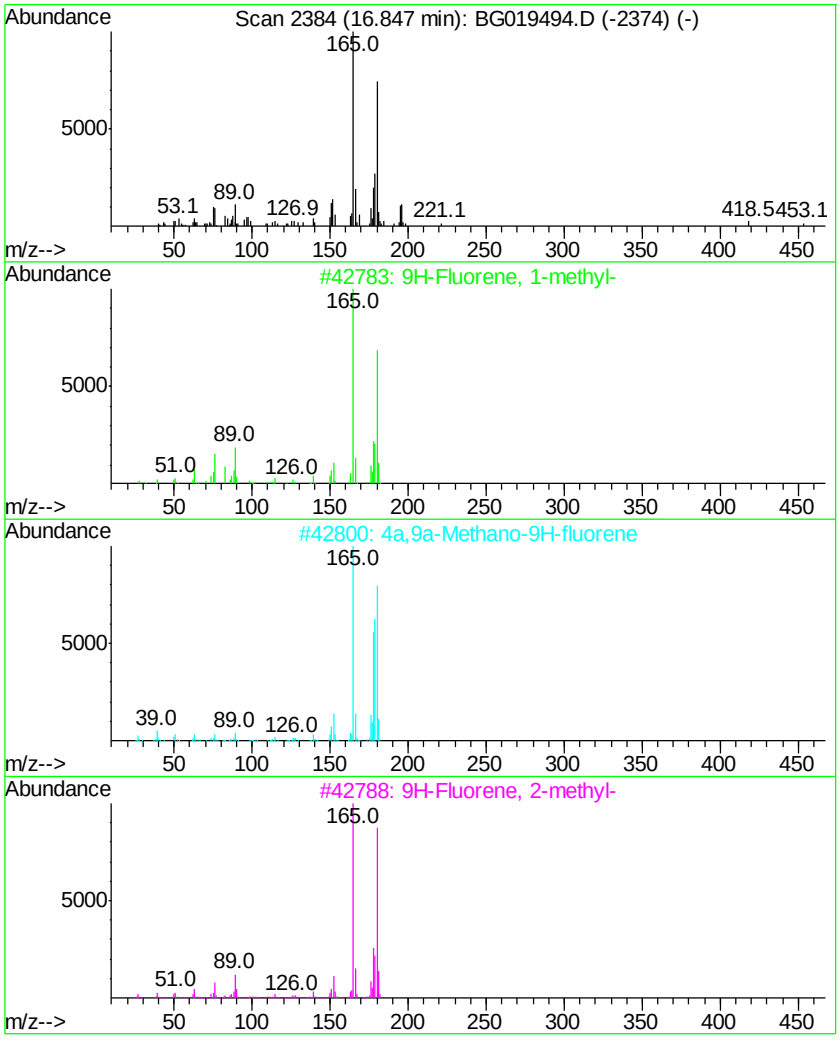
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

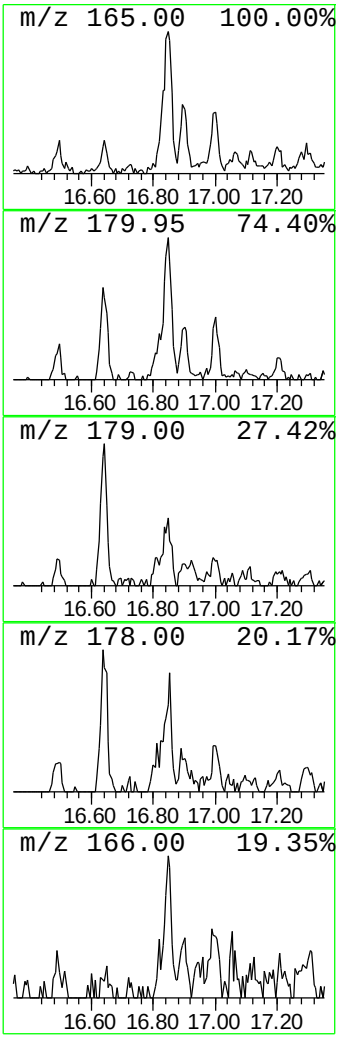
****

Peak Number 5 9H-Fluorene, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.85	2.14 ng/ul	104684	Phenanthrene-d10	17.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	89
2	4a,9a-Methano-9H-fluorene	180	C14H12	019540-84-2	86
3	9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	76
4	9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	76
5	9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	76





Data Path : Z:\HPCHEM1\BNA G\DATA\BG110515\
Data File : BG019494.D
Acq On : 5 Nov 2015 16:06
Operator : UM/NP
Sample : G4273-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AP4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102815.M
Quant Title : SVOA CALIBRATION

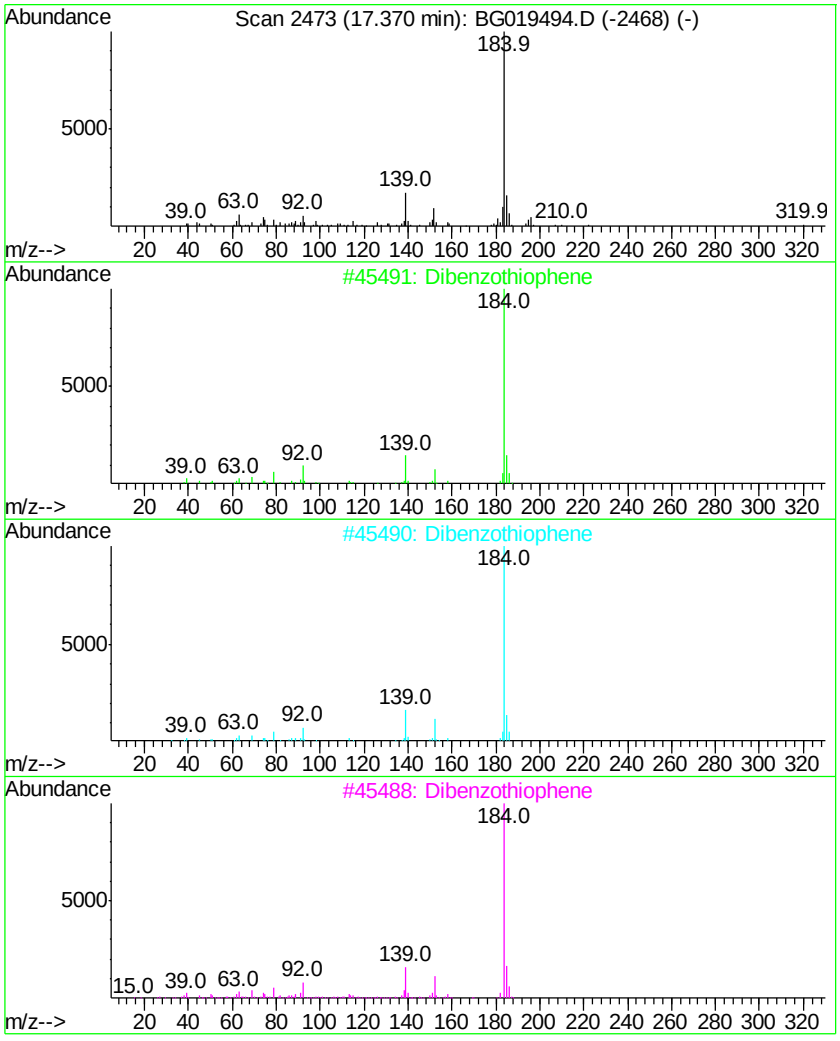
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

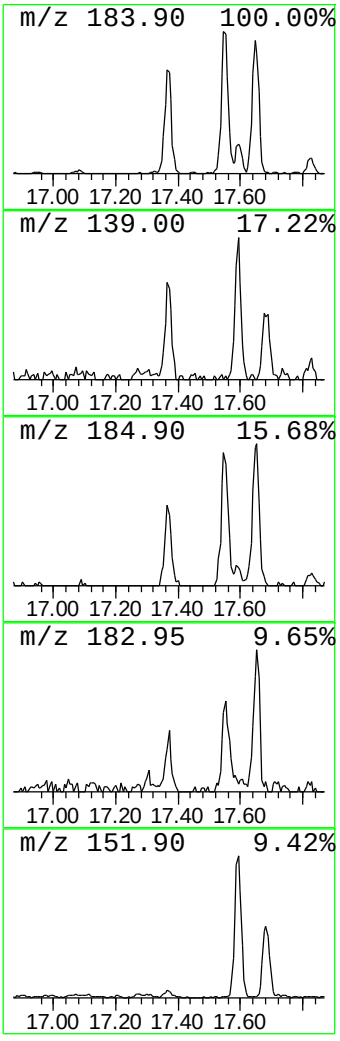
****

Peak Number 6 Dibenzothiophene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.37	2.27 ng/ul	110984	Phenanthrene-d10	17.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzothiophene	184	C12H8S	000132-65-0	91
2	Dibenzothiophene	184	C12H8S	000132-65-0	90
3	Dibenzothiophene	184	C12H8S	000132-65-0	90
4	Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	90
5	Dibenzothiophene	184	C12H8S	000132-65-0	87





Data Path : Z:\HPCHEM1\BNA G\DATA\BG110515\
Data File : BG019494.D
Acq On : 5 Nov 2015 16:06
Operator : UM/NP
Sample : G4273-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AP4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102815.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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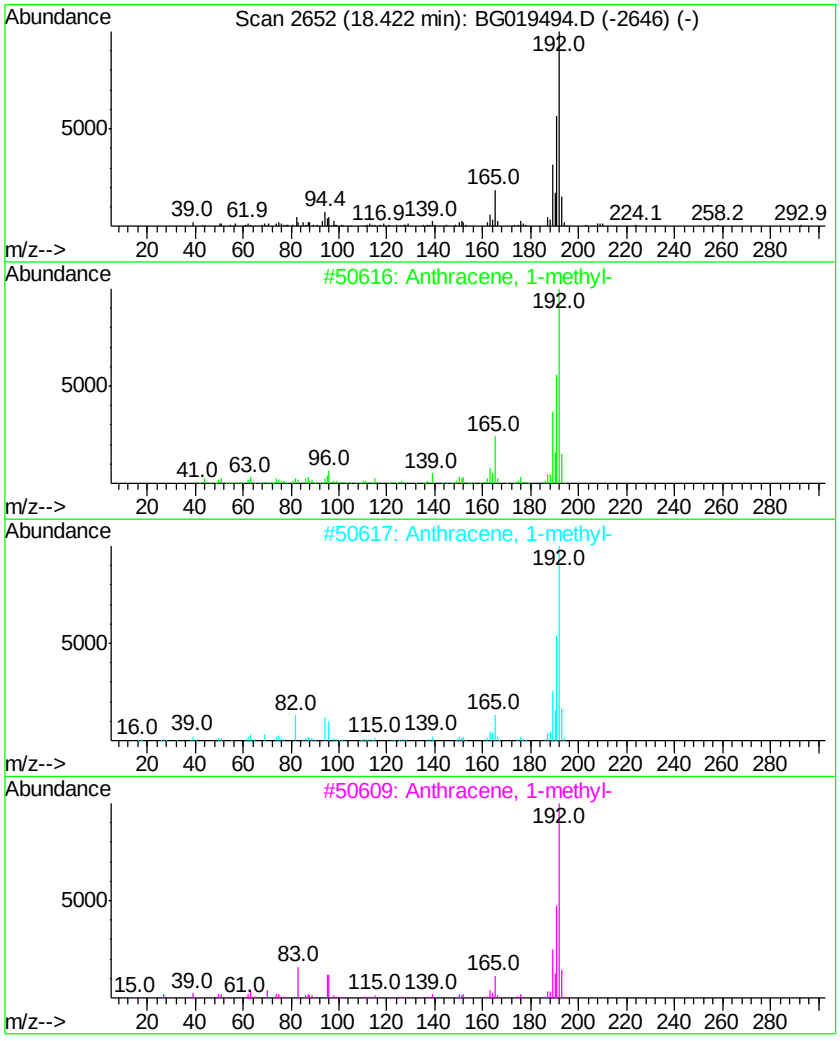
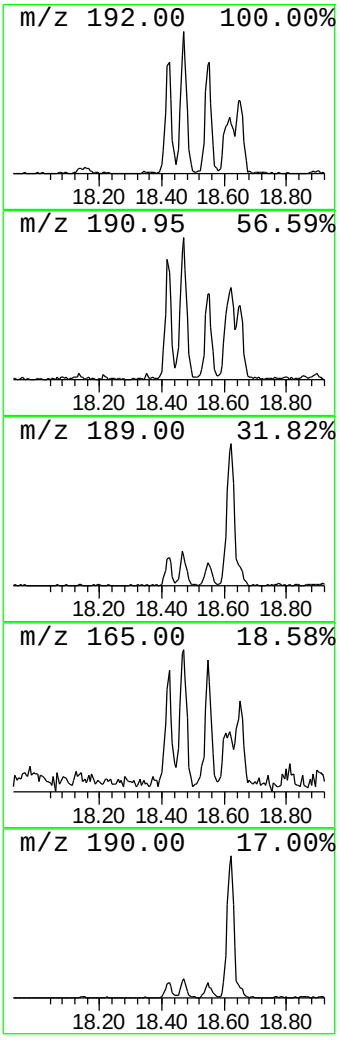
Peak Number 7 Anthracene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.42	3.07 ng/ul	150213	Phenanthrene-d10	17.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
2	Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	87
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	87





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Data File : BG019494.D
Acq On : 5 Nov 2015 16:06
Operator : UM/NP
Sample : G4273-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AP4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102815.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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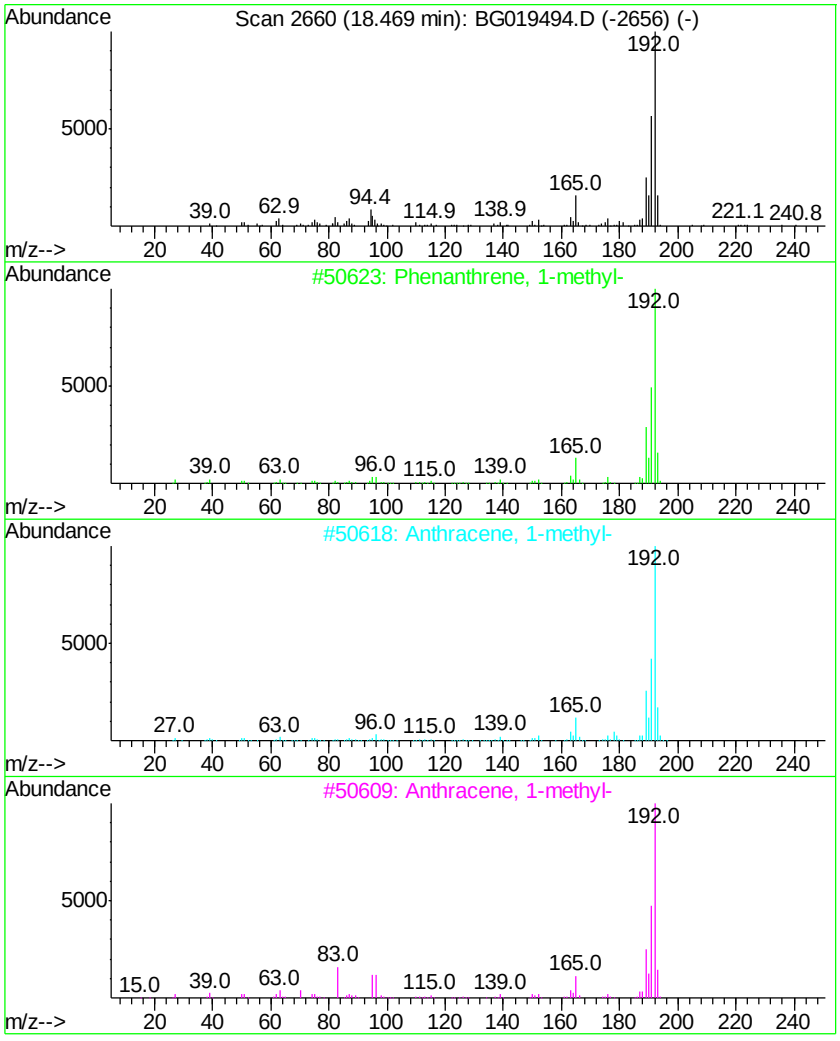
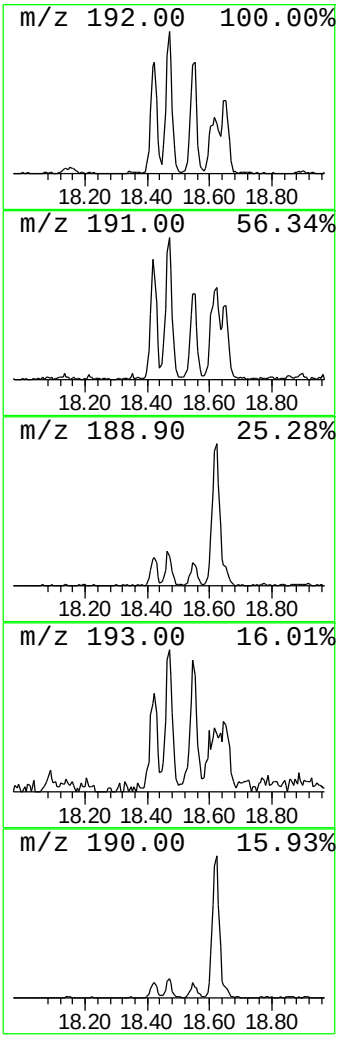
Peak Number 8 Phenanthrene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.47	3.73 ng/ul	182482	Phenanthrene-d10	17.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2	Anthracene, 1-methyl-	192	C15H12	000610-48-0	95
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	94
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93





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Data File : BG019494.D
Acq On : 5 Nov 2015 16:06
Operator : UM/NP
Sample : G4273-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AP4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102815.M
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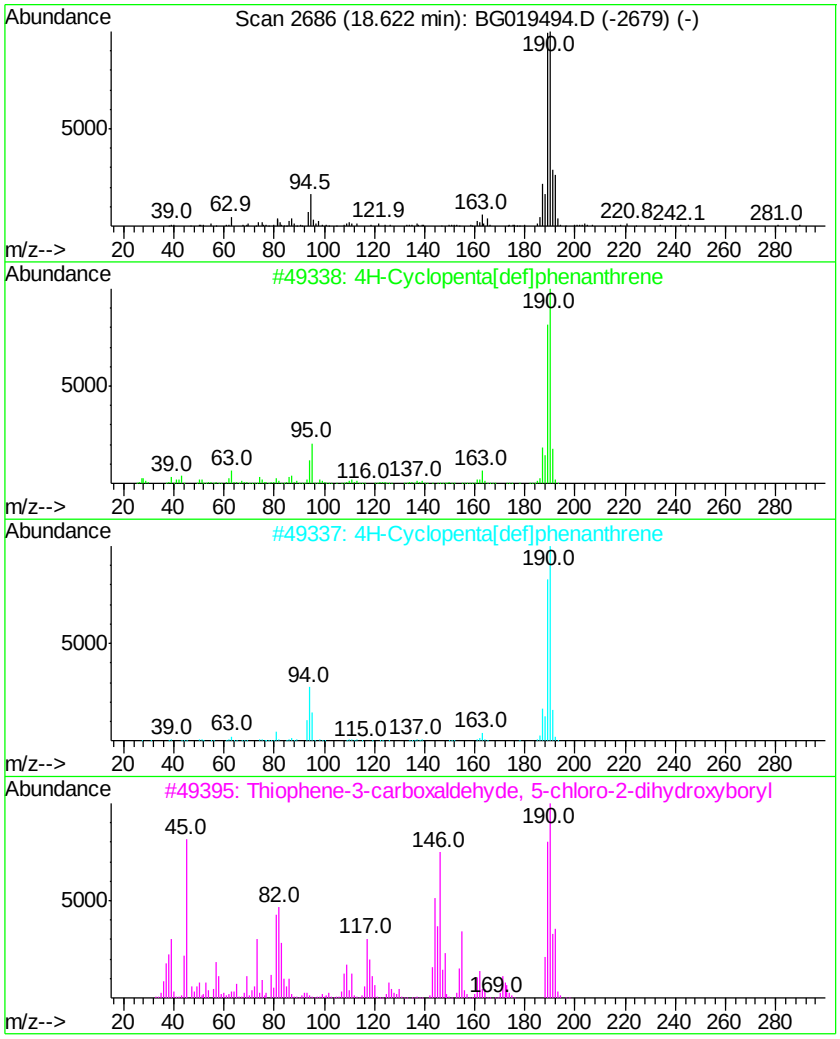
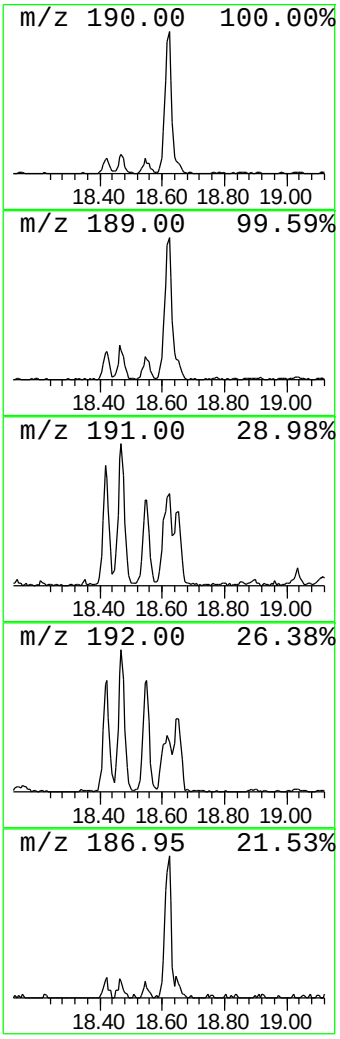
Peak Number 10 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.62	6.25 ng/ul	305472	Phenanthrene-d10	17.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	94
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	74
3	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	52
4	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	37
5	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	25





Library Search Compound Report

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Operator : UM/NP
Sample : G4273-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AP4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102815.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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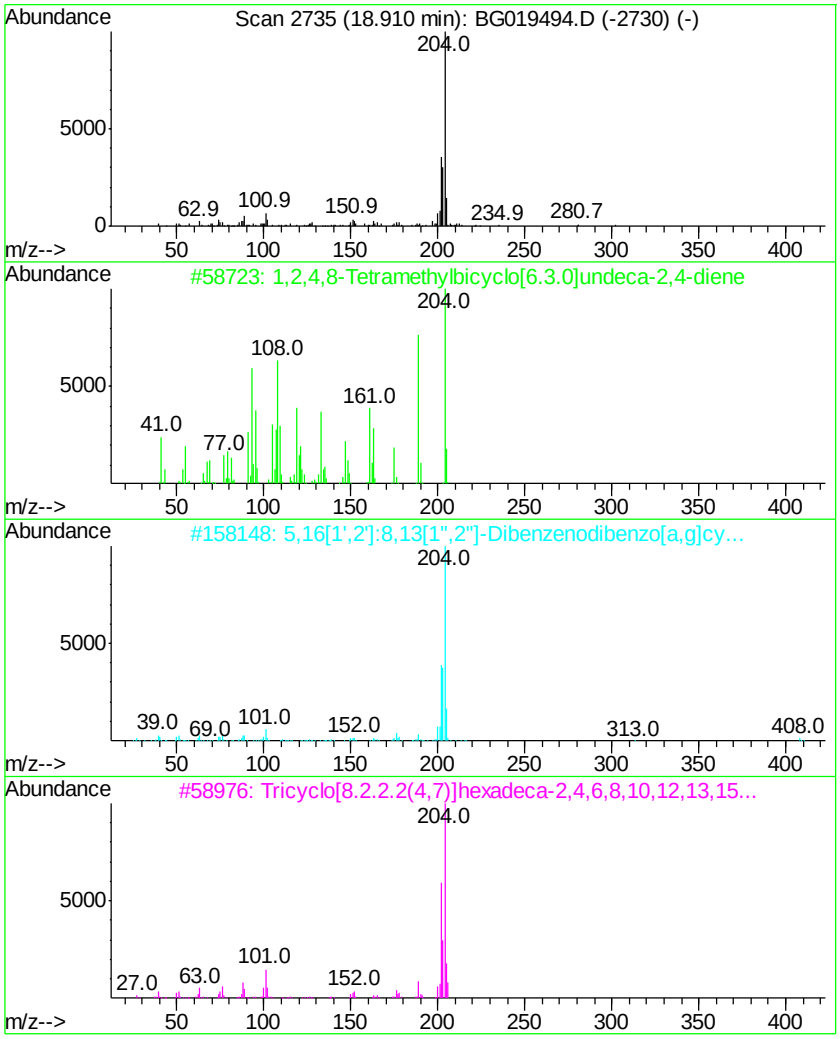
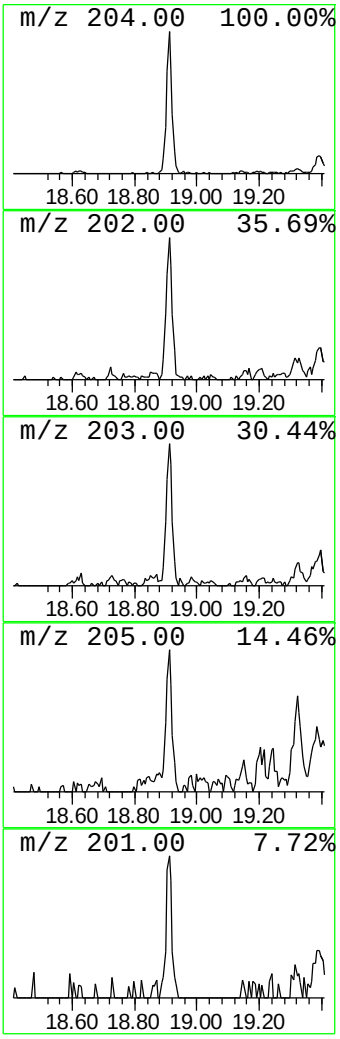
Peak Number 11 1,2,4,8-Tetramethylbicyclo[... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.91	2.34 ng/ul	114545	Phenanthrene-d10	17.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	90
2	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	86
3	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	72
4	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	68
5	2-Phenylnaphthalene	204	C16H12	035465-71-5	68





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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AP4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102815.M
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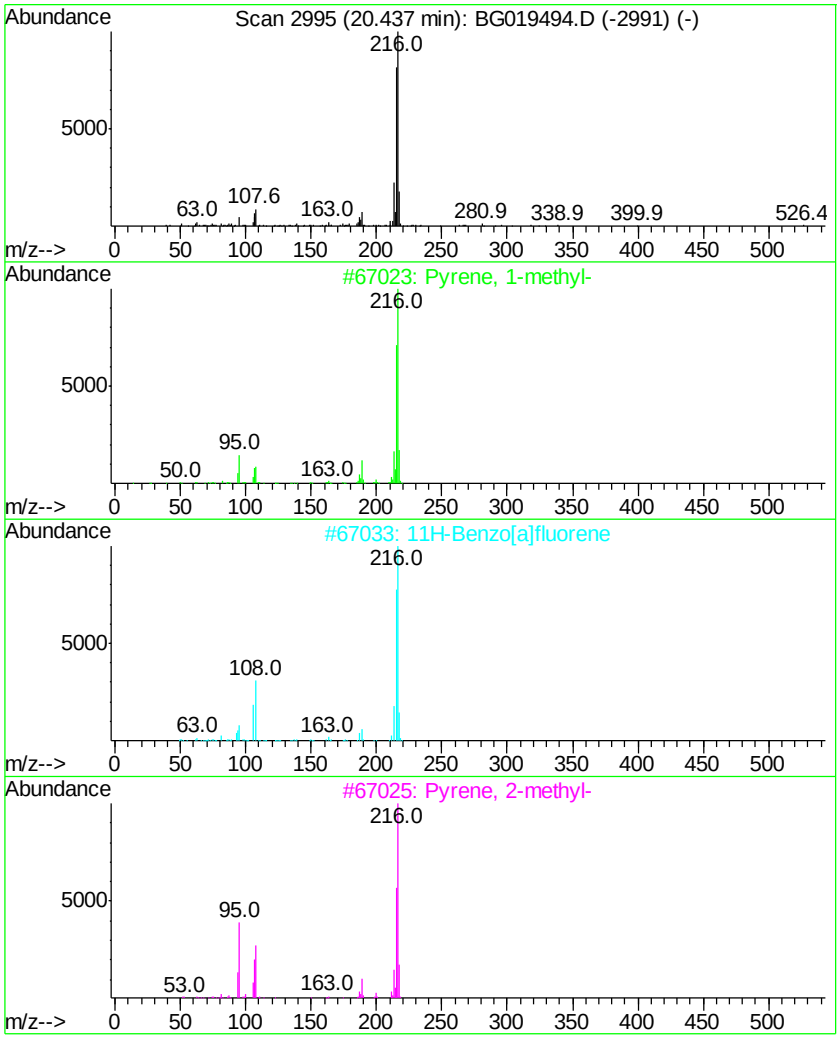
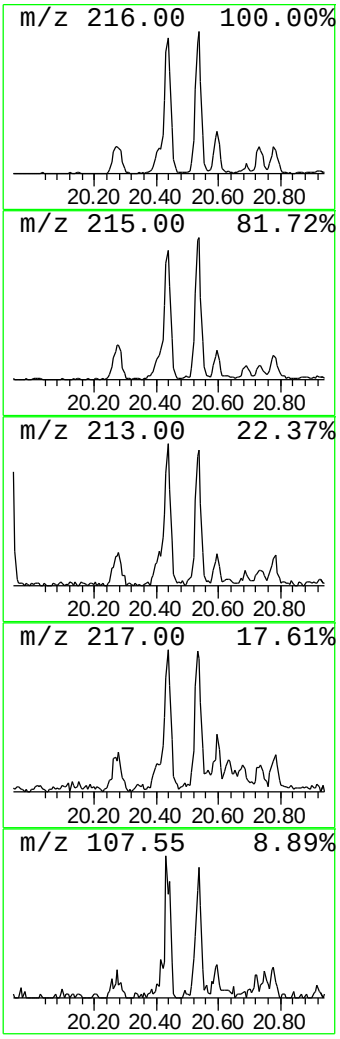
Peak Number 12 Pyrene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.44	3.01 ng/ul	265000	Chrysene-d12	21.83

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
3	Pyrene, 2-methyl-	216	C17H12	003442-78-2	91
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91





Tentatively Identified Compound (LSC) summary

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ClientSampleId : C0AP4

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-methoxy...	3.20	71.5	ng/ul	933729	1	8.18	261354	20.0
9H-Fluoren-9-ol	16.14	2.5	ng/ul	83536	3	14.81	677819	20.0
Dibenzofuran, 4-m...	16.29	3.2	ng/ul	158504	4	17.55	978212	20.0
9H-Fluorene, 1-me...	16.85	2.1	ng/ul	104684	4	17.55	978212	20.0
Dibenzothiophene	17.37	2.3	ng/ul	110984	4	17.55	978212	20.0
Anthracene, 1-met...	18.42	3.1	ng/ul	150213	4	17.55	978212	20.0
Phenanthrene, 1-m...	18.47	3.7	ng/ul	182482	4	17.55	978212	20.0
4H-Cyclopenta[def...	18.62	6.3	ng/ul	305472	4	17.55	978212	20.0
1,2,4,8-Tetrameth...	18.91	2.3	ng/ul	114545	4	17.55	978212	20.0
Pyrene, 1-methyl-	20.44	3.0	ng/ul	265000	5	21.83	1763300	20.0