LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011620\
Data File : BG044070.D
Acq On : 16 Jan 2020 16:02
Operator : JU
Sample : L1126-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BW-1-3.0

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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\8270-BG123019.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.129	3	7	26	rVB	297774	696661	18.47%	1.554%
2	5.050	328	334	346	rVB	296794	491957	13.04%	1.098%
3	5.526	406	415	432	rBV	1309050	2196238	58.22%	4.901%
4	7.113	677	685	704	rBV	1395139	2509387	66.52%	5.599%
5	7.477	739	747	759	rBV	1393067	2448670	64.91%	5.464%
6	7.941	819	826	835	rBV	409143	703241	18.64%	1.569%
7	8.264	871	881	898	rBV	1055846	1890425	50.11%	4.218%
8	9.093	1014	1022	1034	rBV	785649	1467116	38.89%	3.274%
9	10.744	1295	1303	1310	rBV	527183	986674	26.16%	2.202%
10	13.200	1713	1721	1733	rBV	1935786	3182230	84.36%	7.101%
11	13.476	1762	1768	1773	rBV	43173	68478	1.82%	0.153%
12	14.022	1855	1861	1867	rBV	172969	265836	7.05%	0.593%
13	14.222	1890	1895	1902	rBV	63562	114724	3.04%	0.256%
14	14.298	1903	1908	1923	rVB	43167	96319	2.55%	0.215%
15	14.575	1947	1955	1962	rBV	805586	1291435	34.23%	2.882%
16	16.061	2201	2208	2217	rBV	1488603	2340987	62.06%	5.223%
17	16.296	2242	2248	2260	rVB4	33108	73875	1.96%	0.165%
18	16.631	2296	2305	2309	rBV8	15482	41282	1.09%	0.092%
19	17.142	2387	2392	2400	rVB3	23093	47143	1.25%	0.105%
20	17.313	2414	2421	2426	rBV	889782	1427866	37.85%	3.186%
21	17.354	2426	2428	2435	rVB	231848	317739	8.42%	0.709%
22	17.448	2440	2444	2450	rVB	75412	118832	3.15%	0.265%
23	17.506	2450	2454	2460	rBV4	21867	42717	1.13%	0.095%
24	17.718	2485	2490	2493	rBV4	23457	45863	1.22%	0.102%
25	17.830	2506	2509	2514	rVB3	28676	44398	1.18%	0.099%
26	18.194	2566	2571	2575	rBV	84746	162950	4.32%	0.364%
27	18.241	2575	2579	2587	rVB	126924	197964	5.25%	0.442%
28	18.323	2588	2593	2597	rBV	54630	84593	2.24%	0.189%
29	18.382	2597	2603	2607	rBV2	150873	280962	7.45%	0.627%
30	18.687	2651	2655	2659	rBV	76034	111895	2.97%	0.250%
31	18.934	2695	2697	2701	rVB	44797	47116	1.25%	0.105%
32	18.987	2703	2706	2709	rBV	41188	45131	1.20%	0.101%
33	19.022	2709	2712	2716	rVB2	45954	62986	1.67%	0.141%
34	19.075	2718	2721	2722	rBV2	50843	51825	1.37%	0.116%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011620\
Data File : BG044070.D
Acq On : 16 Jan 2020 16:02
Operator : JU
Sample : L1126-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BW-1-3.0

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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\8270-BG123019.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	19.105	2723	2726	2730	rVB	111914	164446	4.36%	0.367%
36	19.163	2731	2736	2740	rBV3	60153	118479	3.14%	0.264%
37	19.240	2745	2749	2758	rBV4	75343	162630	4.31%	0.363%
38	19.369	2765	2771	2777	rBV	1207694	1727493	45.79%	3.855%
39	19.516	2793	2796	2799	rBV	56312	74191	1.97%	0.166%
40	19.727	2822	2832	2841	rBV	1289815	2191403	58.09%	4.890%
41	19.798	2842	2844	2851	rVB3	68494	125408	3.32%	0.280%
42	19.933	2861	2867	2871	rVB	2871198	3772284	100.00%	8.417%
43	20.051	2882	2887	2894	rBV5	143551	353701	9.38%	0.789%
44	20.221	2913	2916	2919	rVB	226814	267027	7.08%	0.596%
45	20.321	2930	2933	2937	rVB	130695	153404	4.07%	0.342%
46	20.380	2939	2943	2947	rVB2	184925	267711	7.10%	0.597%
47	20.521	2963	2967	2970	rBV	128626	160163	4.25%	0.357%
48	20.562	2971	2974	2978	rVB2	139230	178160	4.72%	0.398%
49	21.196	3079	3082	3085	rBV	106571	144010	3.82%	0.321%
50	21.231	3085	3088	3095	rVV3	423266	680722	18.05%	1.519%
51	21.561	3136	3144	3149	rBV2	1209123	3137415	83.17%	7.001%
52	21.608	3149	3152	3161	rVB	744347	1209987	32.08%	2.700%
53	23.229	3424	3428	3434	rVB3	160859	281790	7.47%	0.629%
54	23.699	3501	3508	3515	rBV	573123	1834772	48.64%	4.094%
55	24.392	3621	3626	3638	rVB	391073	1087956	28.84%	2.428%
56	24.533	3643	3650	3660	rVB	535955	1393785	36.95%	3.110%
57	24.680	3668	3675	3683	rBV2	510165	1374023	36.42%	3.066%

Sum of corrected areas: 44816475

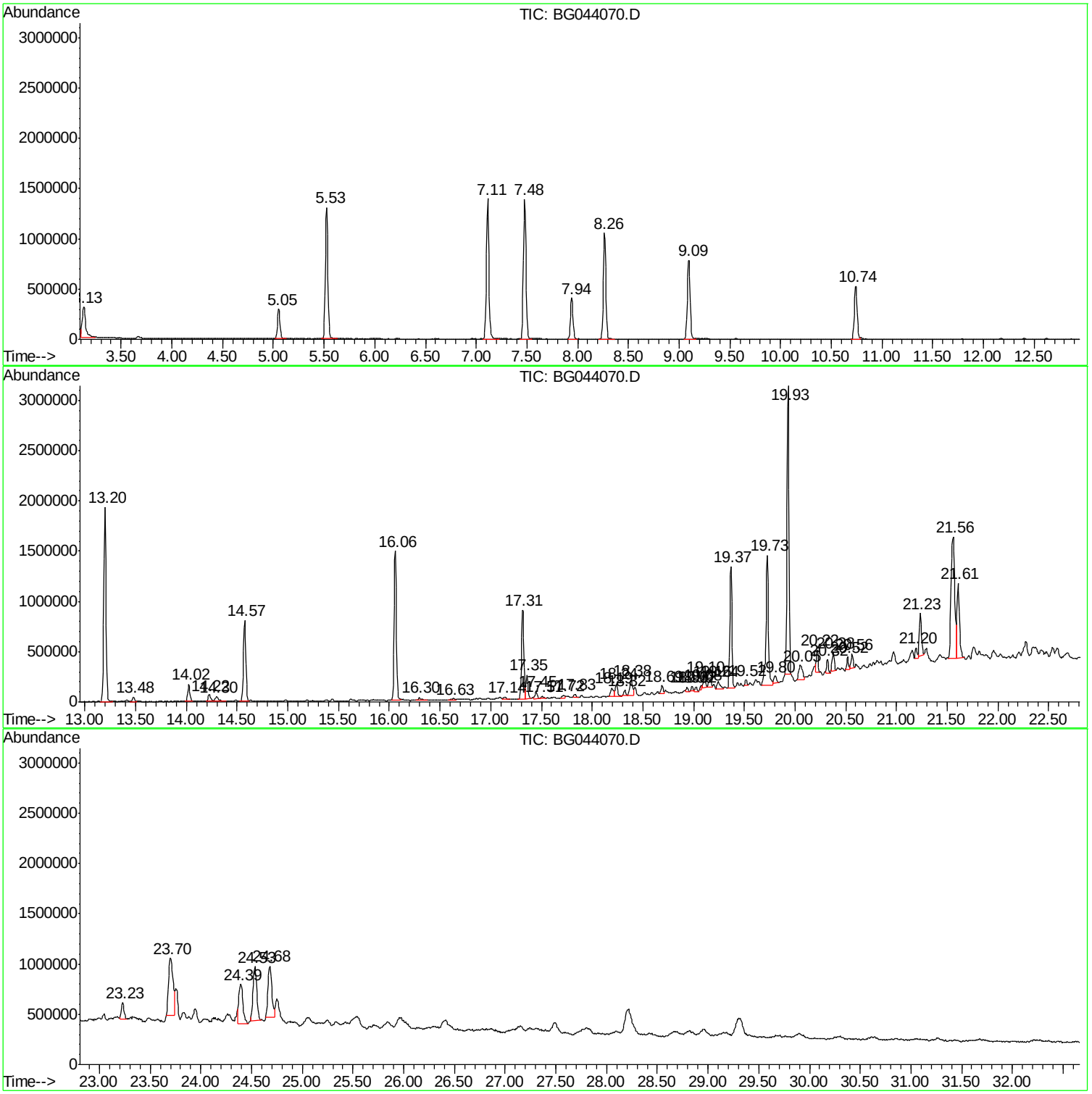
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011620\
Data File : BG044070.D
Acq On : 16 Jan 2020 16:02
Operator : JU
Sample : L1126-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BW-1-3.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011620\
Data File : BG044070.D
Acq On : 16 Jan 2020 16:02
Operator : JU
Sample : L1126-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BW-1-3.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

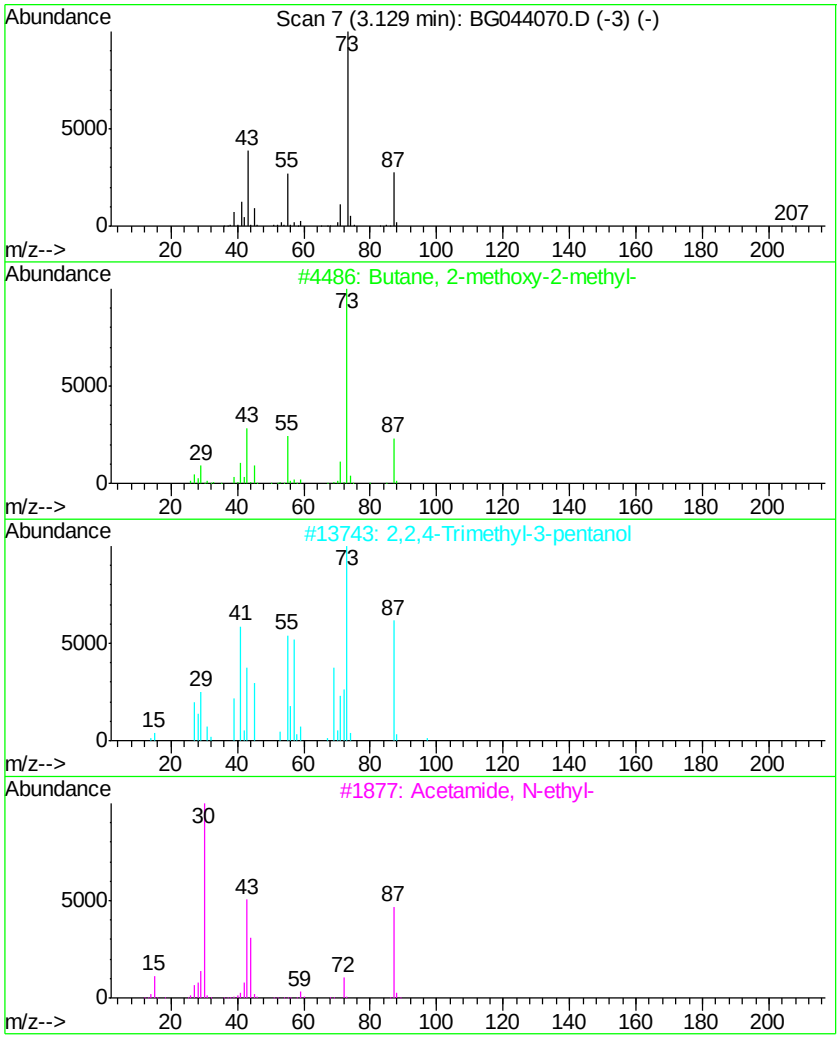
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

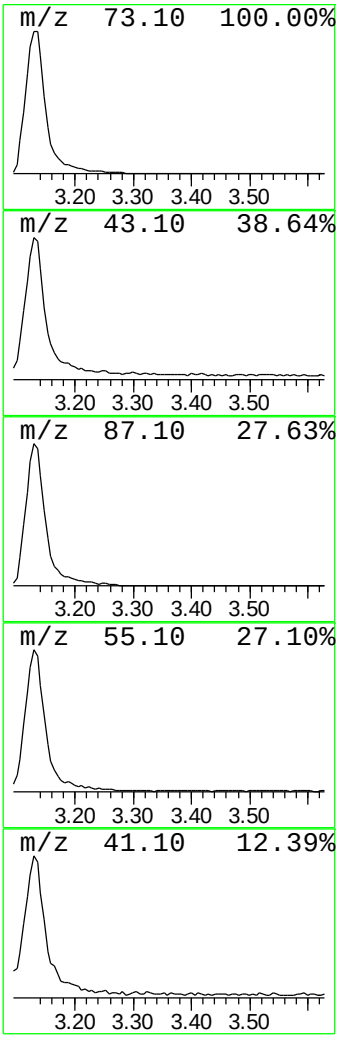
****

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.13	19.81 ng	696661	1,4-Dichlorobenzene-d4	7.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	35
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011620\
Data File : BG044070.D
Acq On : 16 Jan 2020 16:02
Operator : JU
Sample : L1126-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BW-1-3.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

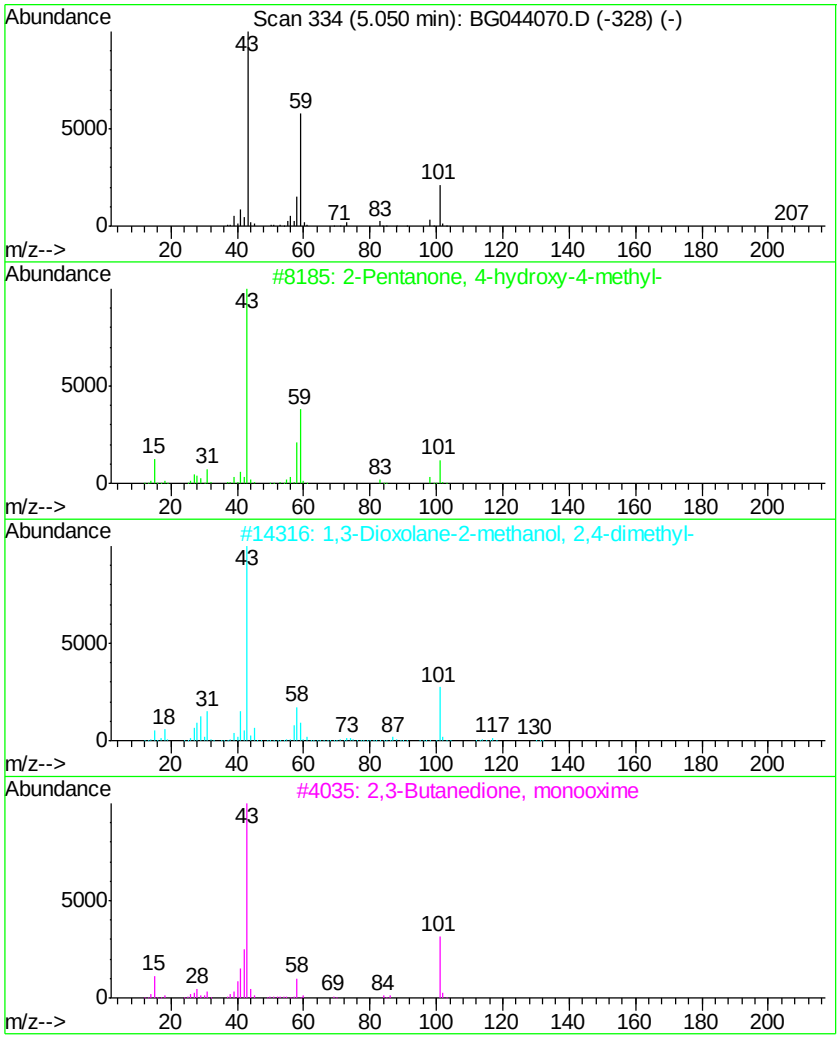
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

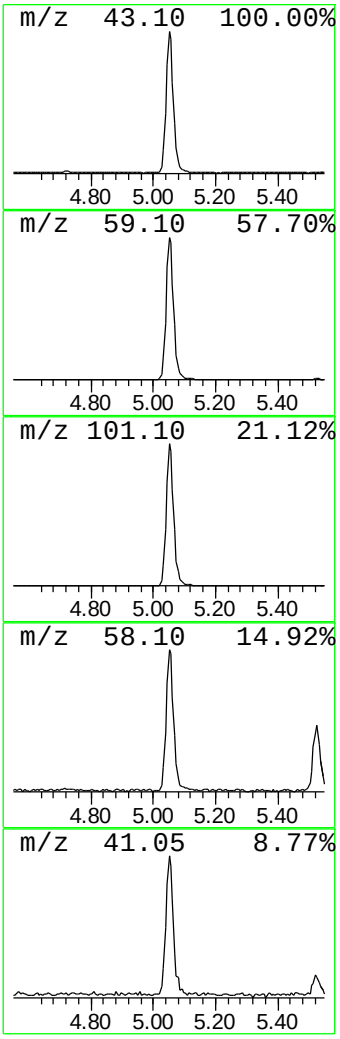
****

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	13.99 ng	491957	1,4-Dichlorobenzene-d4	7.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	1,3-Dioxolane-2-methanol, 2,4-di...	132	C6H12O3	053951-43-2	28
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5	Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	17





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011620\
Data File : BG044070.D
Acq On : 16 Jan 2020 16:02
Operator : JU
Sample : L1126-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BW-1-3.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

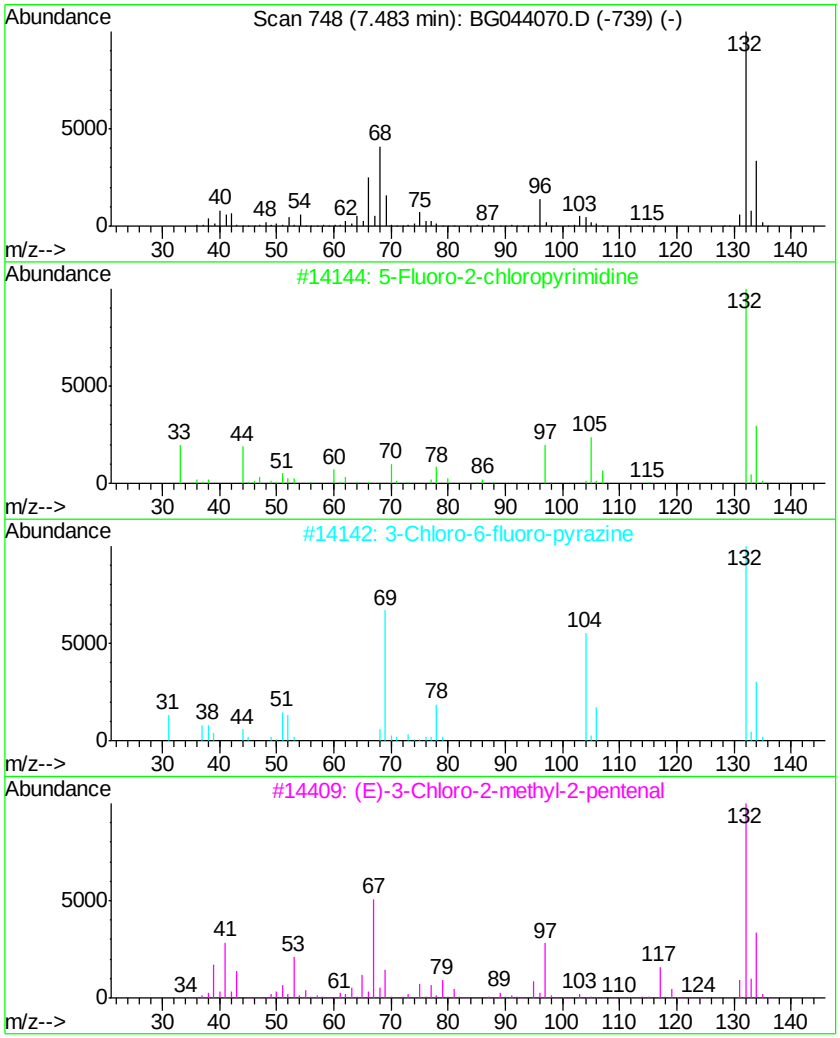
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

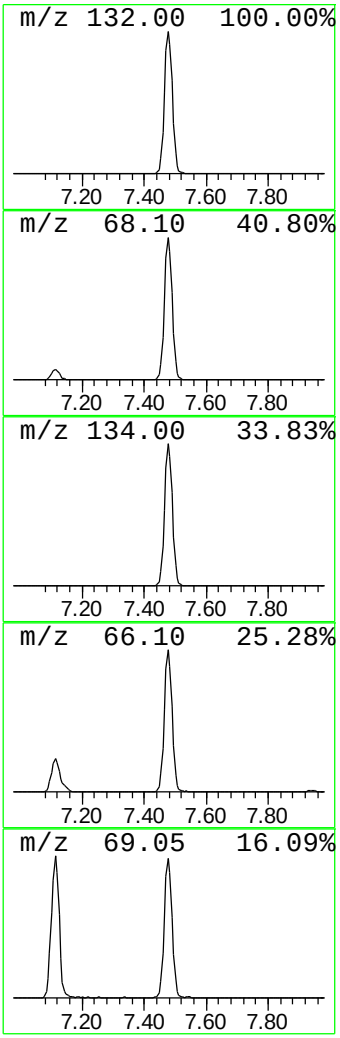
****

Peak Number 3 unknown7.48 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.48	69.64 ng	2448670	1,4-Dichlorobenzene-d4	7.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
4	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5	Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011620\
Data File : BG044070.D
Acq On : 16 Jan 2020 16:02
Operator : JU
Sample : L1126-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BW-1-3.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

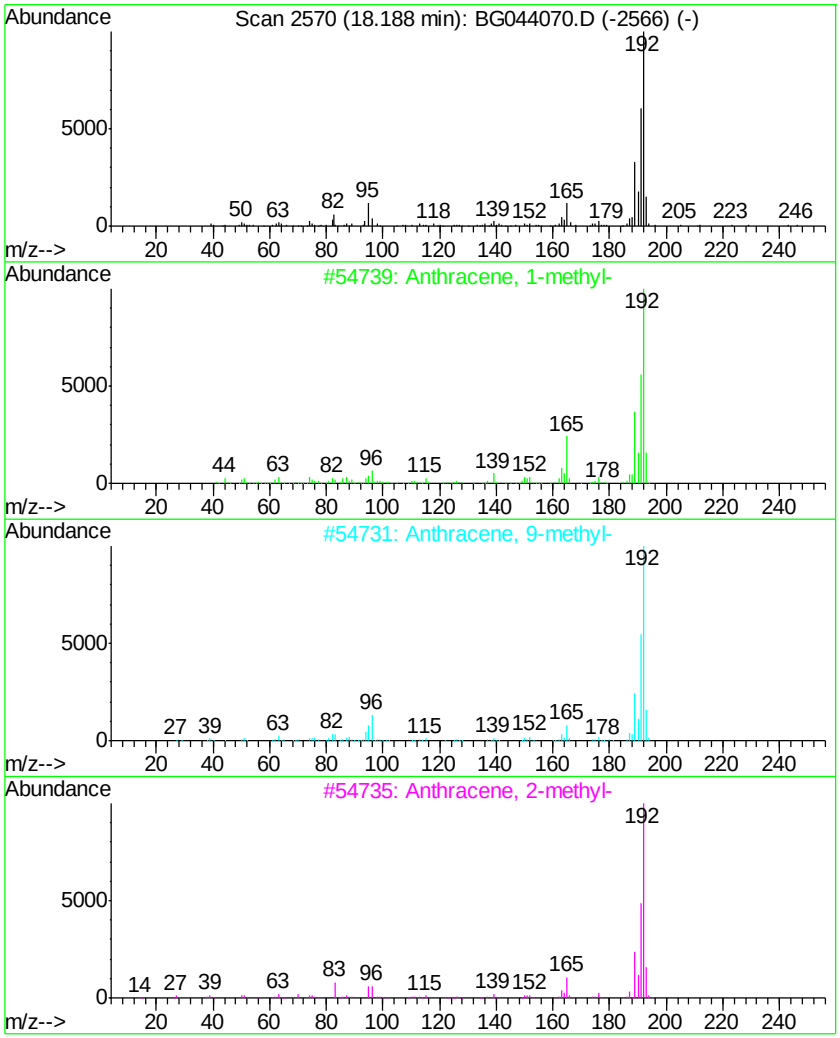
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

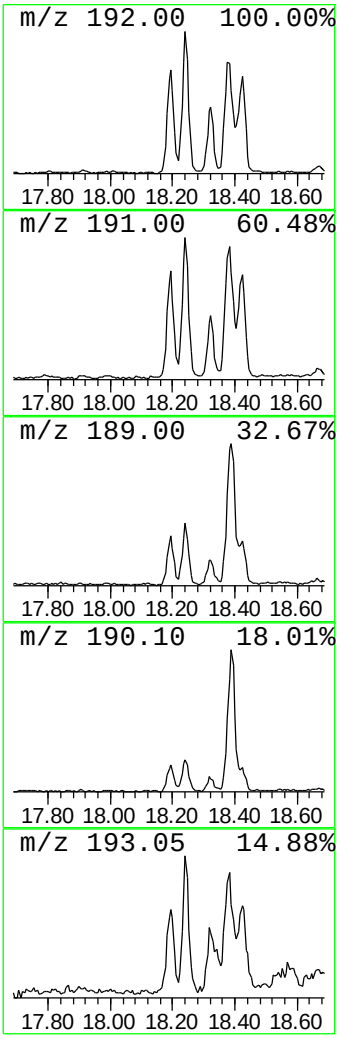
****

Peak Number 5 Anthracene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.19	2.28 ng	162950	Phenanthrene-d10	17.31

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
2	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	91
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	89





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011620\
Data File : BG044070.D
Acq On : 16 Jan 2020 16:02
Operator : JU
Sample : L1126-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BW-1-3.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

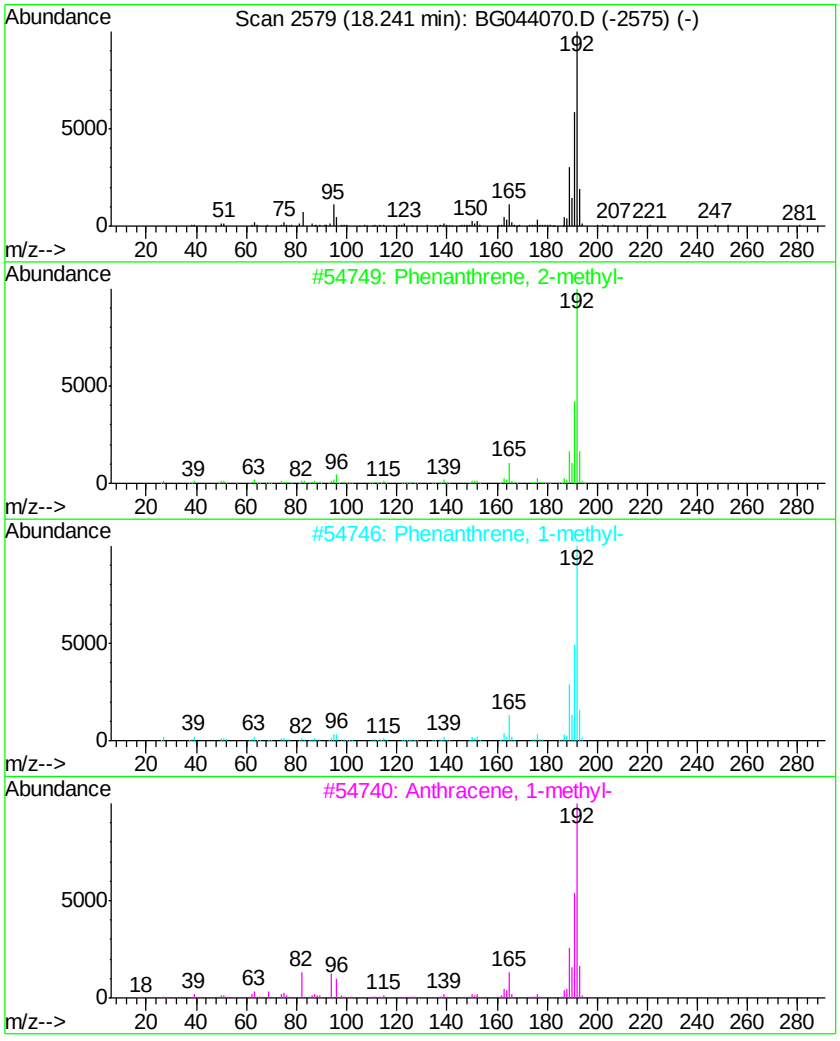
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

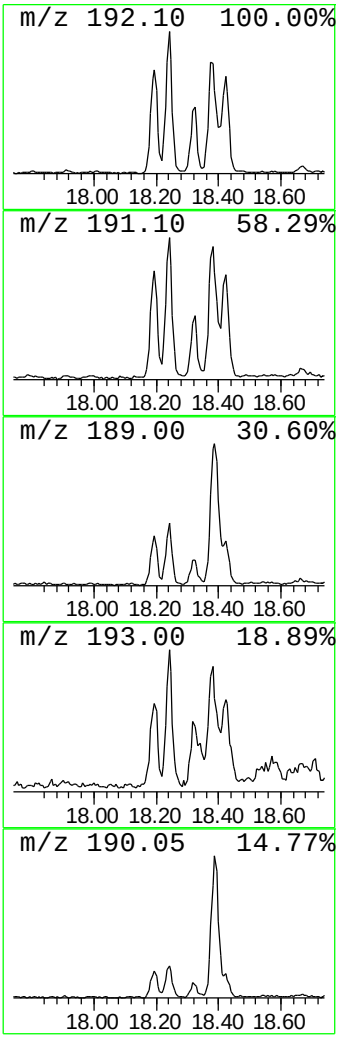
****

Peak Number 6 Phenanthrene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.24	2.77 ng	197964	Phenanthrene-d10	17.31

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
4	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	93
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011620\
Data File : BG044070.D
Acq On : 16 Jan 2020 16:02
Operator : JU
Sample : L1126-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BW-1-3.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

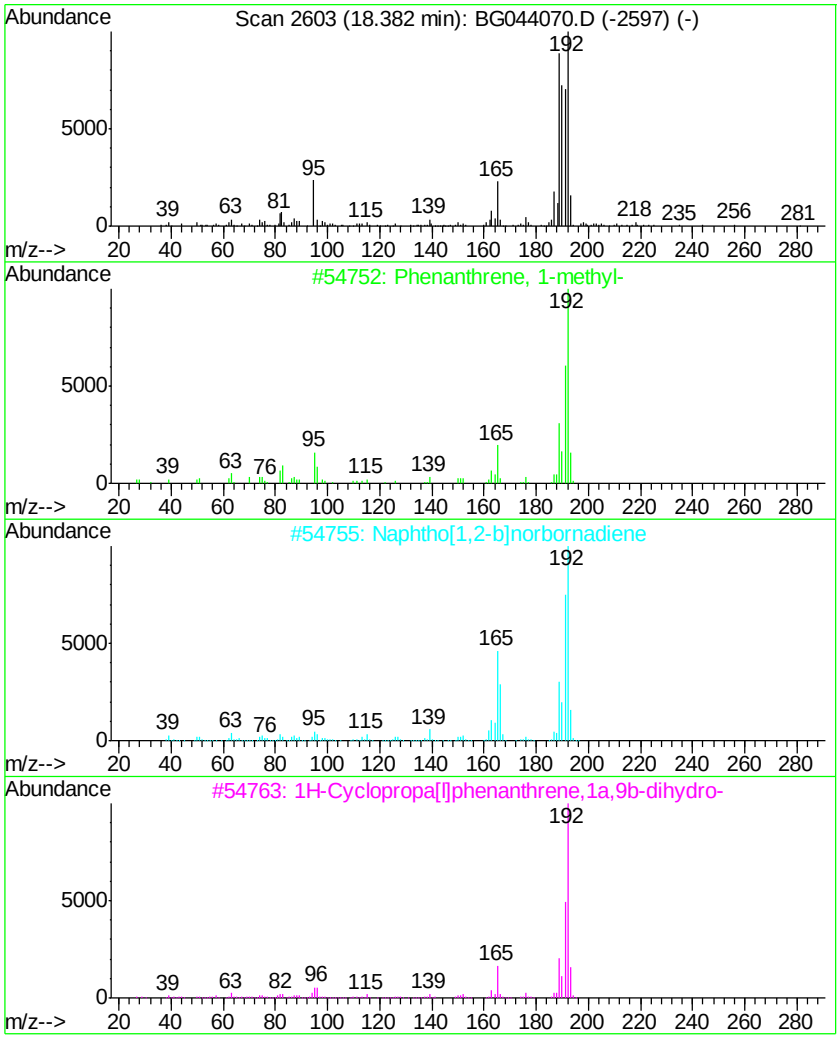
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

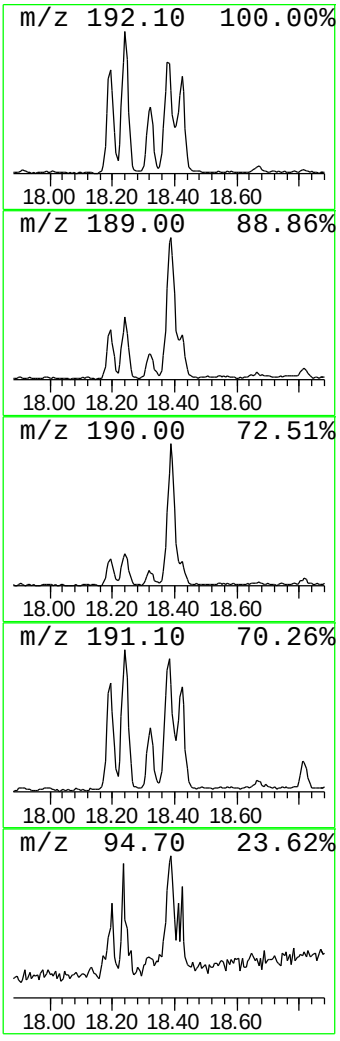
****

Peak Number 7 Phenanthrene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.38	3.94 ng	280962	Phenanthrene-d10	17.31

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	55
2	Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	55
3	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	49
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	46
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	46





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011620\
Data File : BG044070.D
Acq On : 16 Jan 2020 16:02
Operator : JU
Sample : L1126-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BW-1-3.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

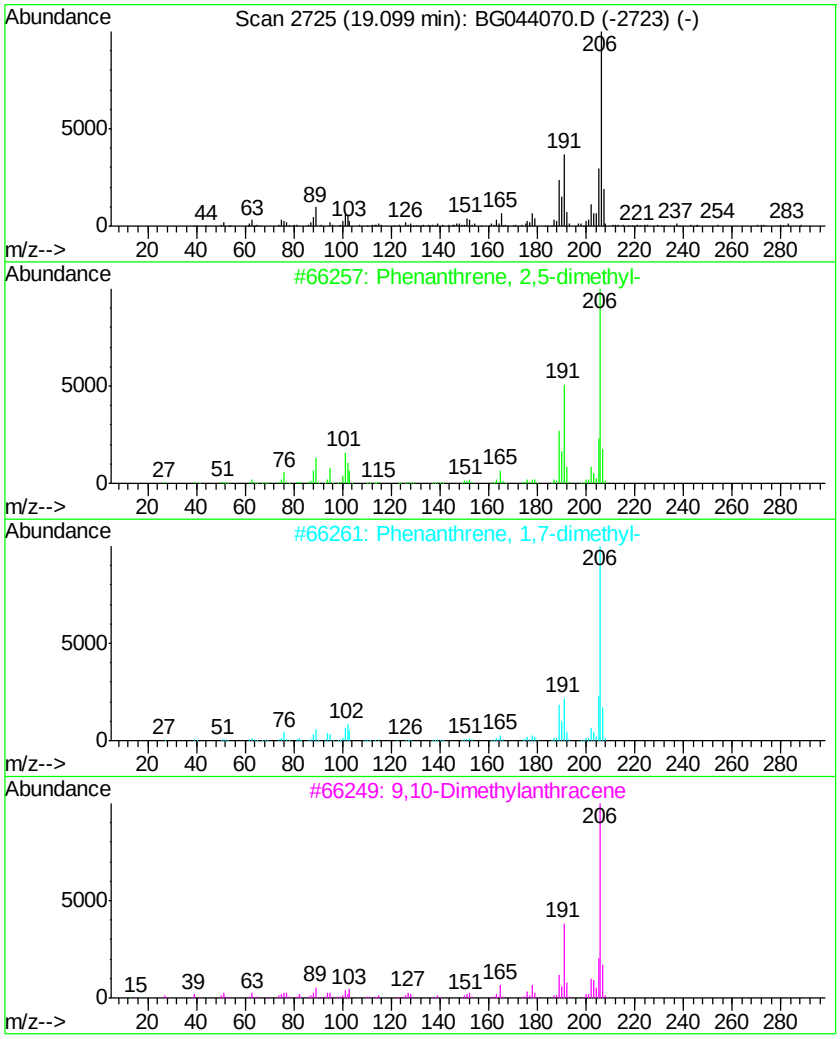
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

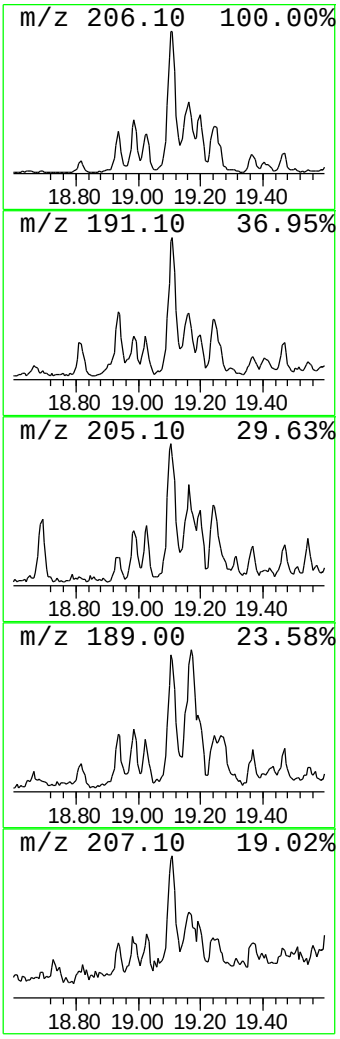
****

Peak Number 8 Phenanthrene, 2,5-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.10	2.30 ng	164446	Phenanthrene-d10	17.31

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	91
3	9,10-Dimethylanthracene	206	C16H14	000781-43-1	91
4	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	90
5	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	90





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011620\
Data File : BG044070.D
Acq On : 16 Jan 2020 16:02
Operator : JU
Sample : L1126-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BW-1-3.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

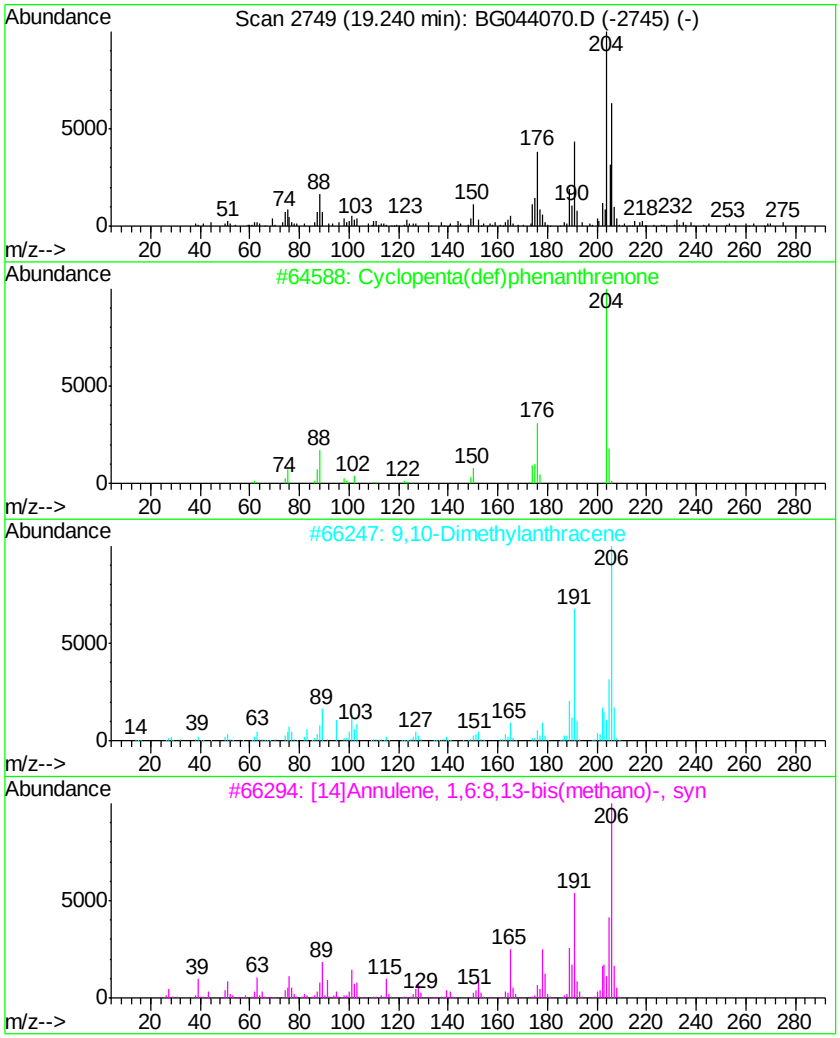
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

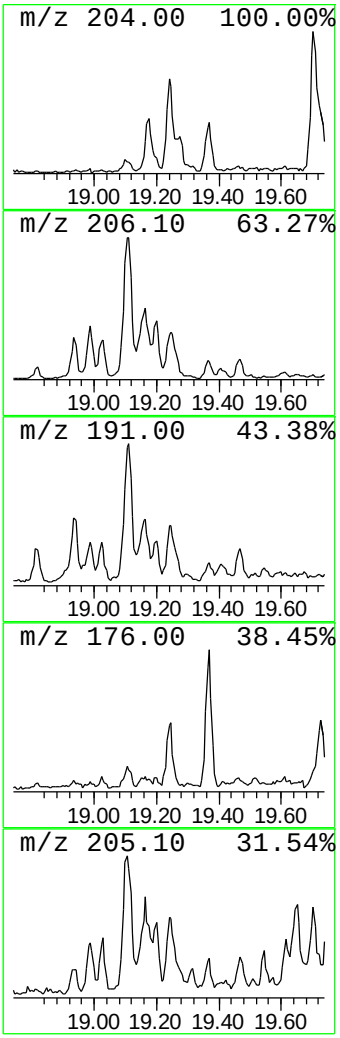
****

Peak Number 9 Cyclopenta(def)phenanthrenone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.24	2.28 ng	162630	Phenanthrene-d10	17.31

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	83
2	9,10-Dimethylanthracene	206	C16H14	000781-43-1	46
3	[14]Annulene, 1,6:8,13-bis(metha...	206	C16H14	085385-68-8	45
4	1,3-Diphenyl-3-methylcyclopropene	206	C16H14	086544-79-8	42
5	1H-Indene, 2-methyl-3-phenyl-	206	C16H14	035099-60-6	42





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011620\
Data File : BG044070.D
Acq On : 16 Jan 2020 16:02
Operator : JU
Sample : L1126-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BW-1-3.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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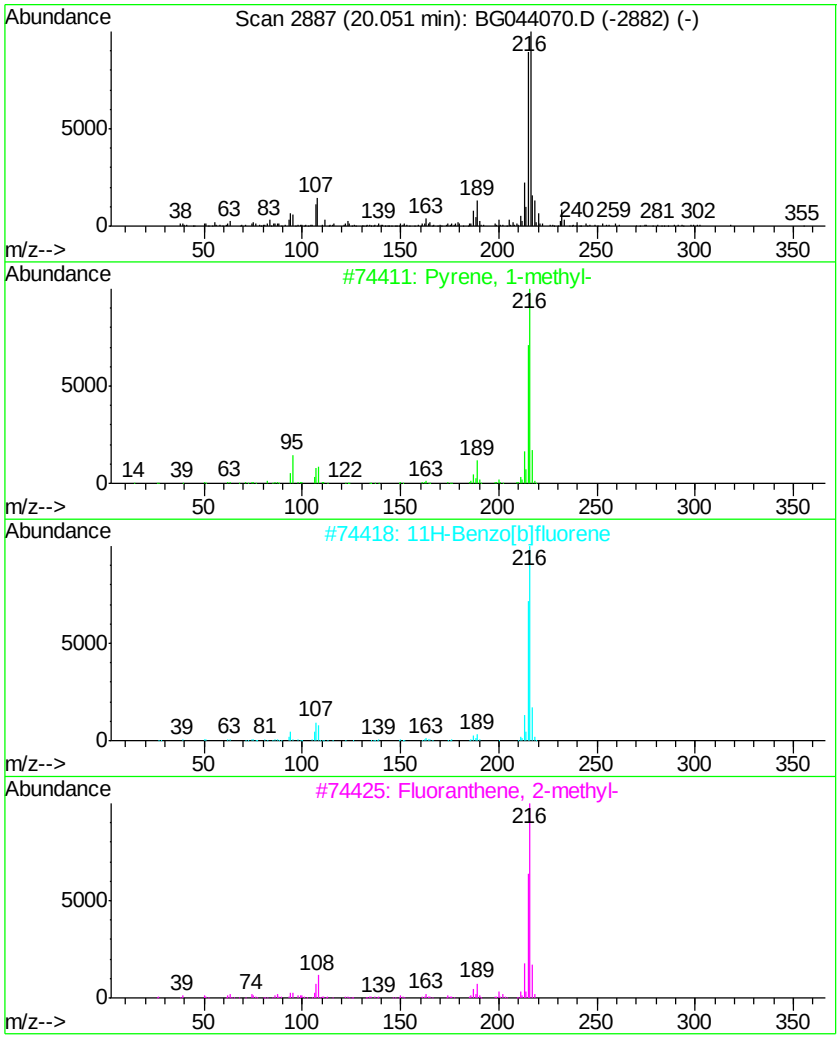
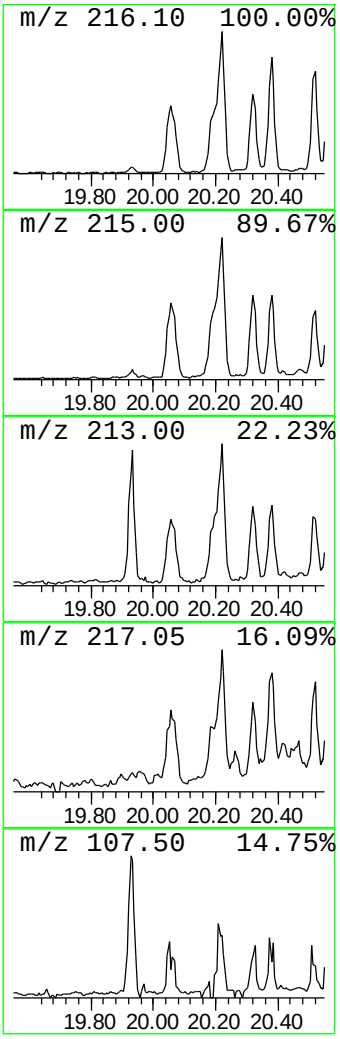
Peak Number 10 Pyrene, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.05	2.25 ng	353701	Chrysene-d12	21.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	83
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	76
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	74
5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	70





Library Search Compound Report

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Data File : BG044070.D
Acq On : 16 Jan 2020 16:02
Operator : JU
Sample : L1126-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BW-1-3.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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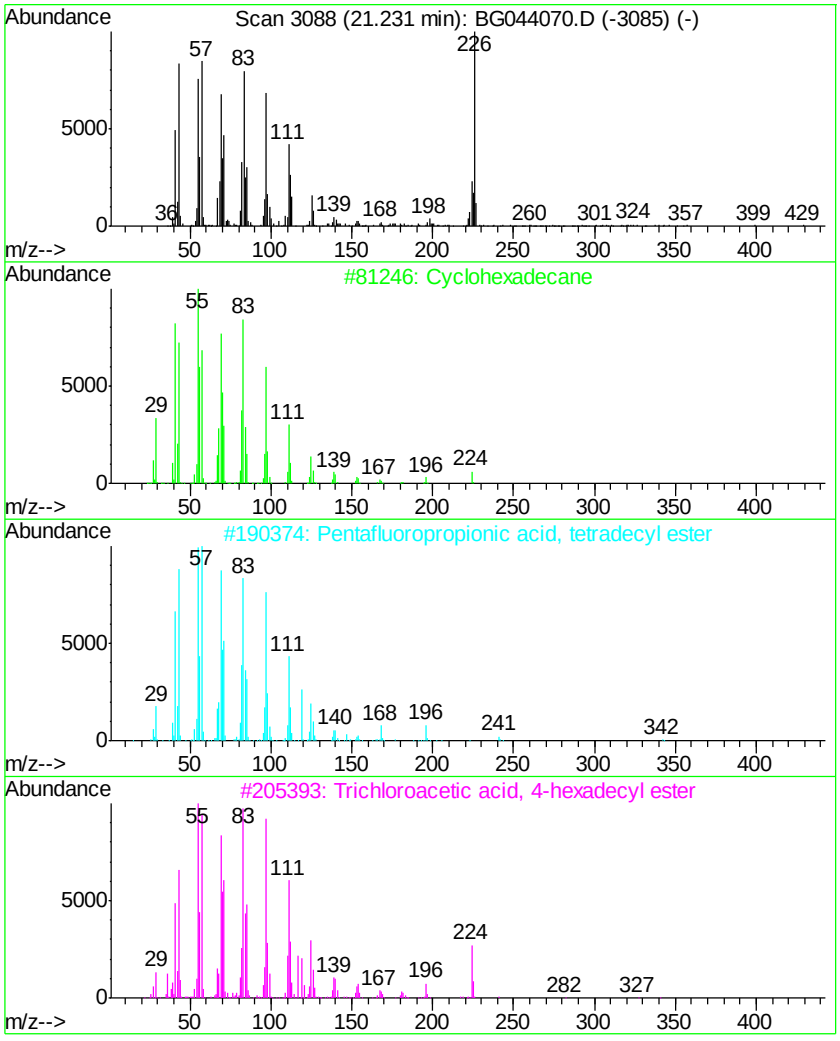
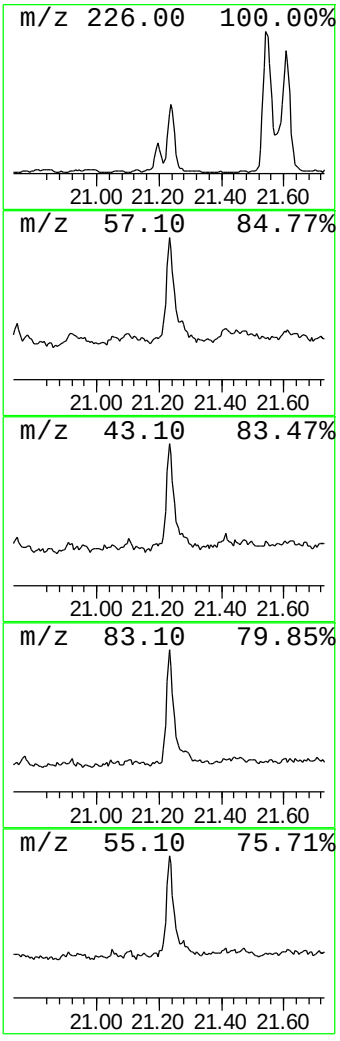
Peak Number 11 Cyclohexadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.23	4.34 ng	680722	Chrysene-d12	21.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane	224	C16H32	000295-65-8	98
2	Pentafluoropropionic acid, tetra...	360	C17H29F5O2	006222-06-6	83
3	Trichloroacetic acid, 4-hexadecy...	386	C18H33Cl3O2	1000282-92-4	70
4	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	64
5	3-Eicosene, (E)-	280	C20H40	074685-33-9	64





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011620\
Data File : BG044070.D
Acq On : 16 Jan 2020 16:02
Operator : JU
Sample : L1126-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BW-1-3.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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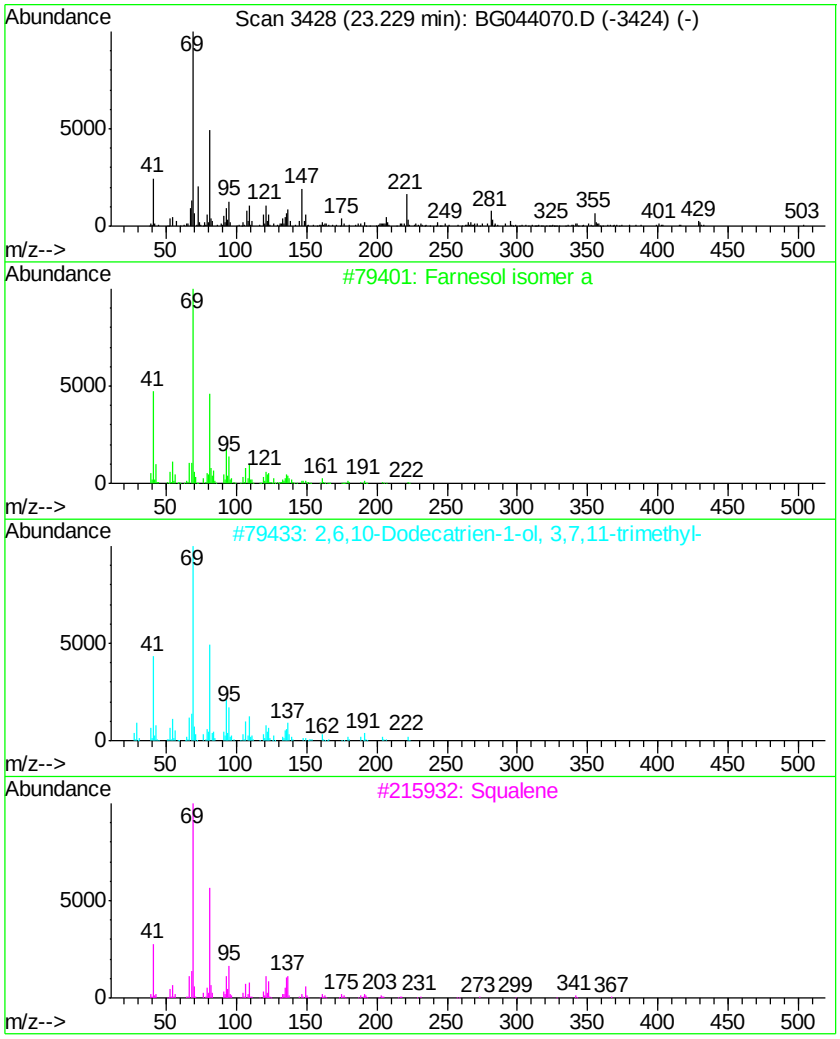
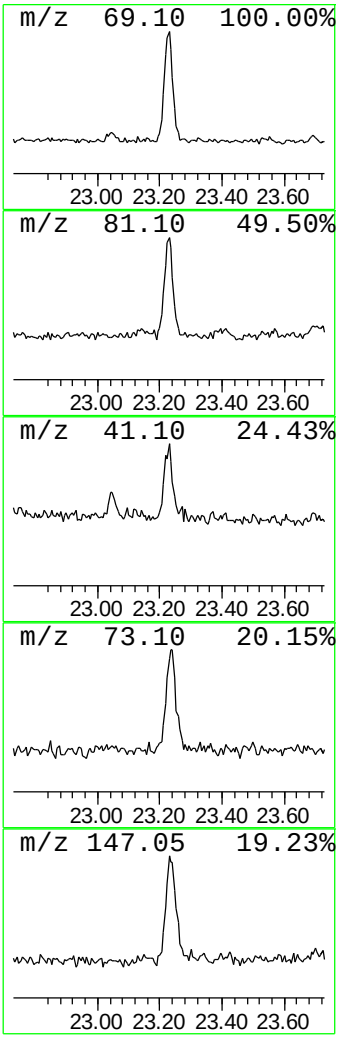
Peak Number 12 Farnesol isomer a Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.23	4.10 ng	281790	Perylene-d12	24.68

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Farnesol isomer a	222	C15H26O	1000108-92-4	55
2	2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	55
3	Squalene	410	C30H50	000111-02-4	49
4	Supraene	410	C30H50	007683-64-9	46
5	4-Fluoroaniline, N,N-di(trifluor...	303	C10H4F7NO2	1000328-30-6	43





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011620\
Data File : BG044070.D
Acq On : 16 Jan 2020 16:02
Operator : JU
Sample : L1126-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BW-1-3.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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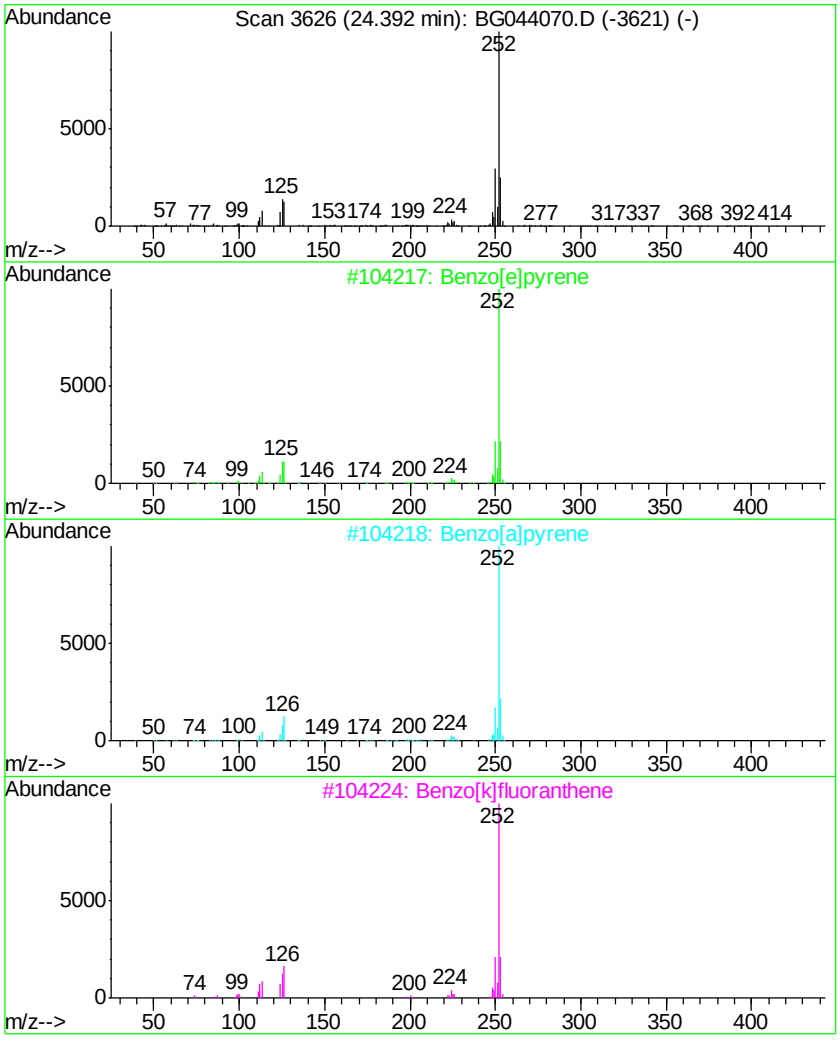
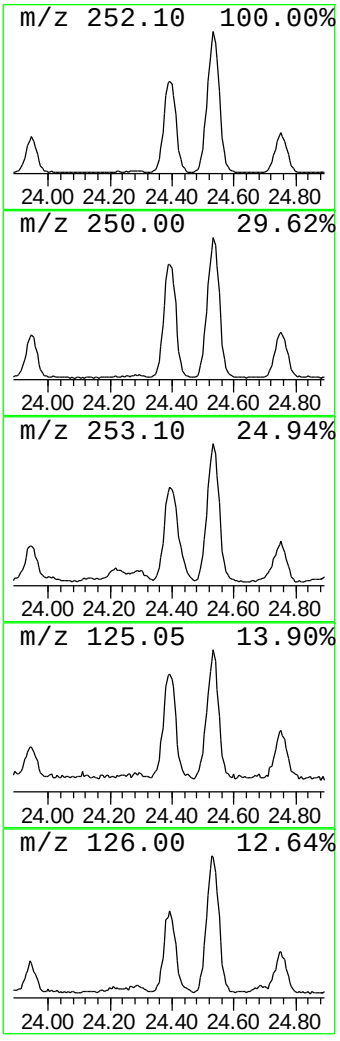
Peak Number 13 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.39	15.84 ng	1087960	Perylene-d12	24.68

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2	Benzo[a]pyrene	252	C20H12	000050-32-8	97
3	Benzo[k]fluoranthene	252	C20H12	000207-08-9	96
4	Benzo[j]fluoranthene	252	C20H12	000205-82-3	95
5	Perylene	252	C20H12	000198-55-0	94





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG011620\
Data File : BG044070.D
Acq On : 16 Jan 2020 16:02
Operator : JU
Sample : L1126-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BW-1-3.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
Butane, 2-methoxy...	3.13	19.8	ng	696661	1	7.94	703241	20.0
2-Pentanone, 4-hy...	5.05	14.0	ng	491957	1	7.94	703241	20.0
unknown7.48	7.48	69.6	ng	2448670	1	7.94	703241	20.0
Anthracene, 1-met...	18.19	2.3	ng	162950	4	17.31	1427870	20.0
Phenanthrene, 2-m...	18.24	2.8	ng	197964	4	17.31	1427870	20.0
Phenanthrene, 1-m...	18.38	3.9	ng	280962	4	17.31	1427870	20.0
Phenanthrene, 2,5...	19.10	2.3	ng	164446	4	17.31	1427870	20.0
Cyclopenta(def)ph...	19.24	2.3	ng	162630	4	17.31	1427870	20.0
Pyrene, 1-methyl-	20.05	2.3	ng	353701	5	21.57	3137420	20.0
Cyclohexadecane	21.23	4.3	ng	680722	5	21.57	3137420	20.0
Farnesol isomer a	23.23	4.1	ng	281790	6	24.68	1374020	20.0
Benzo[e]pyrene	24.39	15.8	ng	1087960	6	24.68	1374020	20.0