LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102021\
Data File : BG050656.D
Acq On : 20 Oct 2021 22:17
Operator : CG/JU
Sample : M4288-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SA-01-101921

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-BG100621.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG050656.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.791	384	392	403	rBV	669580	1168380	60.57%	4.761%
2	7.389	655	664	676	rBV	697555	1249431	64.77%	5.092%
3	8.247	803	810	823	rVB	180522	322184	16.70%	1.313%
4	9.422	1002	1010	1020	rBV	425902	792537	41.09%	3.230%
5	10.815	1241	1247	1256	rBV	128892	226050	11.72%	0.921%
6	11.073	1284	1291	1298	rBV	239139	445871	23.11%	1.817%
7	13.494	1695	1703	1710	rVB	968149	1563781	81.07%	6.373%
8	13.670	1726	1733	1741	rBV5	13448	30793	1.60%	0.125%
9	14.593	1886	1890	1897	rBV	16059	30551	1.58%	0.124%
10	14.869	1923	1937	1943	rBV	355141	589311	30.55%	2.401%
11	14.927	1943	1947	1953	rVV3	14434	23521	1.22%	0.096%
12	15.257	1999	2003	2011	rVB3	13072	26335	1.37%	0.107%
13	15.680	2072	2075	2079	rVB2	16556	21153	1.10%	0.086%
14	15.909	2108	2114	2120	rBV2	23257	38199	1.98%	0.156%
15	16.349	2183	2189	2197	rVB	676573	1050129	54.44%	4.279%
16	16.537	2217	2221	2223	rBV2	23414	33051	1.71%	0.135%
17	17.131	2316	2322	2325	rBV8	10339	19939	1.03%	0.081%
18	17.331	2352	2356	2361	rVV	35311	49049	2.54%	0.200%
19	17.383	2361	2365	2369	rVV4	21933	36819	1.91%	0.150%
20	17.425	2369	2372	2379	rVB	19122	34497	1.79%	0.141%
21	17.613	2397	2404	2408	rBV	410515	674268	34.96%	2.748%
22	17.654	2408	2411	2417	rVB	271889	375940	19.49%	1.532%
23	17.742	2422	2426	2431	rVB	59874	90245	4.68%	0.368%
24	18.006	2467	2471	2477	rBV	22895	37204	1.93%	0.152%
25	18.071	2479	2482	2487	rVB	27505	33299	1.73%	0.136%
26	18.124	2487	2491	2496	rBV3	21778	31997	1.66%	0.130%
27	18.188	2499	2502	2507	rVV	14694	20579	1.07%	0.084%
28	18.247	2507	2512	2518	rVB	236500	296561	15.37%	1.209%
29	18.341	2523	2528	2533	rBV6	11459	20965	1.09%	0.085%
30	18.470	2545	2550	2556	rBV2	121177	226588	11.75%	0.923%
31	18.529	2556	2560	2566	rVB2	78915	123890	6.42%	0.505%
32	18.611	2570	2574	2578	rVB2	31977	47493	2.46%	0.194%
33	18.676	2579	2585	2590	rBV4	105603	218805	11.34%	0.892%
34	18.758	2596	2599	2603	rBV2	22312	28821	1.49%	0.117%
35	18.970	2631	2635	2639	rBV	55946	87909	4.56%	0.358%
36	19.017	2639	2643	2651	rVB2	42826	75687	3.92%	0.308%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102021\
Data File : BG050656.D
Acq On : 20 Oct 2021 22:17
Operator : CG/JU
Sample : M4288-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SA-01-101921

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-BG100621.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	19.211	2674	2676	2681	rVB3	29246	37021	1.92%	0.151%
38	19.264	2682	2685	2688	rVV2	22065	28592	1.48%	0.117%
39	19.375	2699	2704	2706	rBV2	630253	759066	39.35%	3.093%
40	19.405	2706	2709	2713	rVB	711071	840939	43.60%	3.427%
41	19.534	2725	2731	2736	rBV	207359	297238	15.41%	1.211%
42	19.651	2746	2751	2757	rVB	1098569	1586542	82.25%	6.465%
43	19.980	2800	2807	2808	rBV2	158698	257609	13.36%	1.050%
44	20.016	2808	2813	2819	rVB	974671	1479276	76.69%	6.028%
45	20.080	2820	2824	2829	rVB2	54714	80249	4.16%	0.327%
46	20.198	2838	2844	2849	rVB	1429288	1928932	100.00%	7.861%
47	20.333	2861	2867	2873	rBV3	81899	211578	10.97%	0.862%
48	20.597	2908	2912	2915	rBV	125405	165915	8.60%	0.676%
49	20.656	2920	2922	2926	rVB2	92514	114890	5.96%	0.468%
50	20.797	2943	2946	2949	rBV	70769	84174	4.36%	0.343%
51	20.979	2974	2977	2981	rVB	128117	155730	8.07%	0.635%
52	21.279	3024	3028	3036	rVB	91212	152399	7.90%	0.621%
53	21.508	3063	3067	3073	rVB2	278830	388790	20.16%	1.584%
54	21.896	3126	3133	3140	rBV2	612608	1670777	86.62%	6.809%
55	21.960	3140	3144	3155	rVB	434112	829043	42.98%	3.378%
56	22.078	3161	3164	3168	rBV2	102348	125246	6.49%	0.510%
57	24.234	3524	3531	3539	rBV2	333102	1144730	59.35%	4.665%
58	25.010	3657	3663	3675	rVB	217393	631560	32.74%	2.574%
59	25.163	3682	3689	3700	rVB2	303364	881327	45.69%	3.591%
60	25.327	3710	3717	3725	rVB3	214326	545923	28.30%	2.225%

Sum of corrected areas: 24539378

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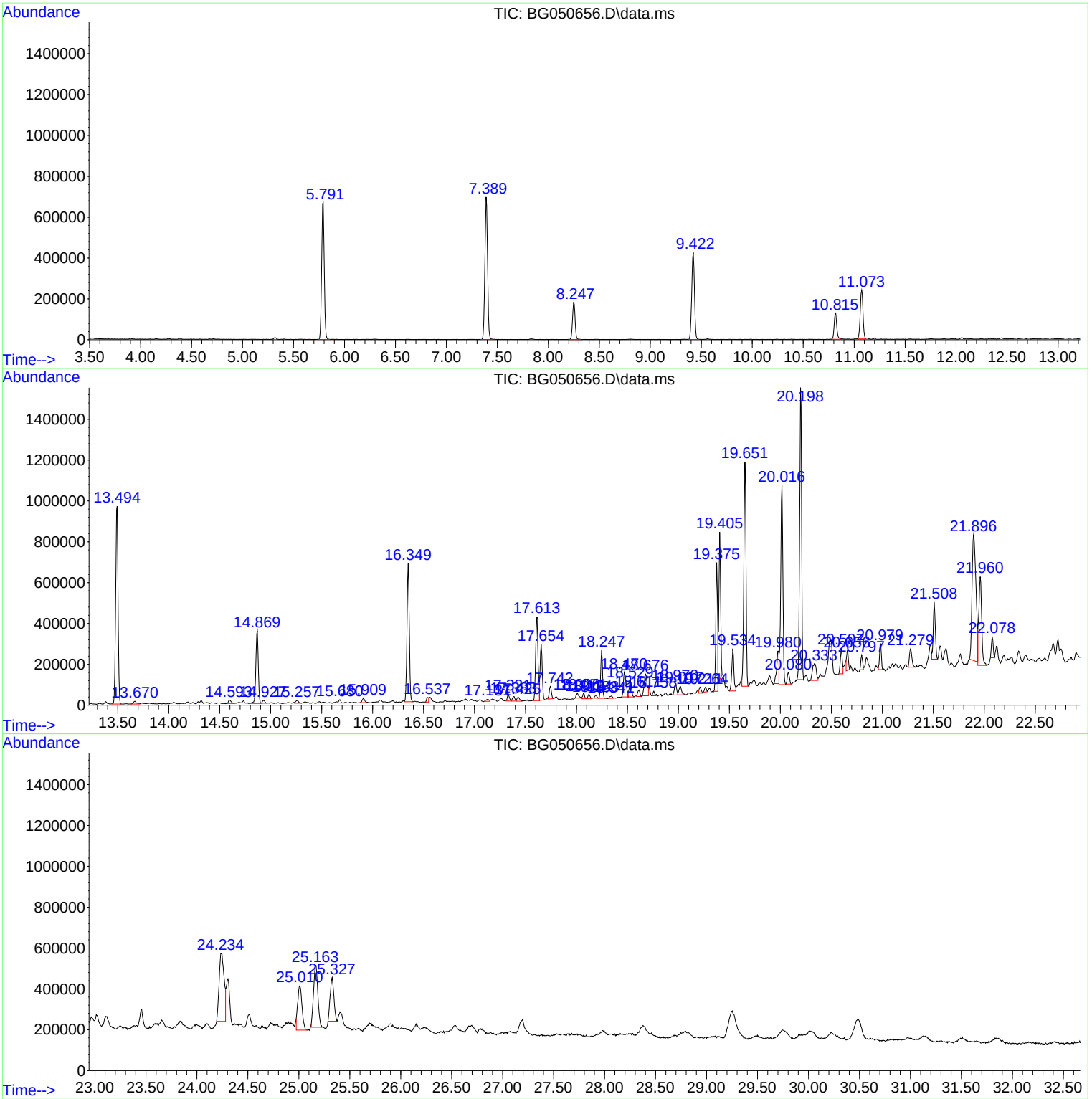
Sample : M4288-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SA-01-101921

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102021\
Data File : BG050656.D
Acq On : 20 Oct 2021 22:17
Operator : CG/JU
Sample : M4288-01
Misc :
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Instrument : BNA_G
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

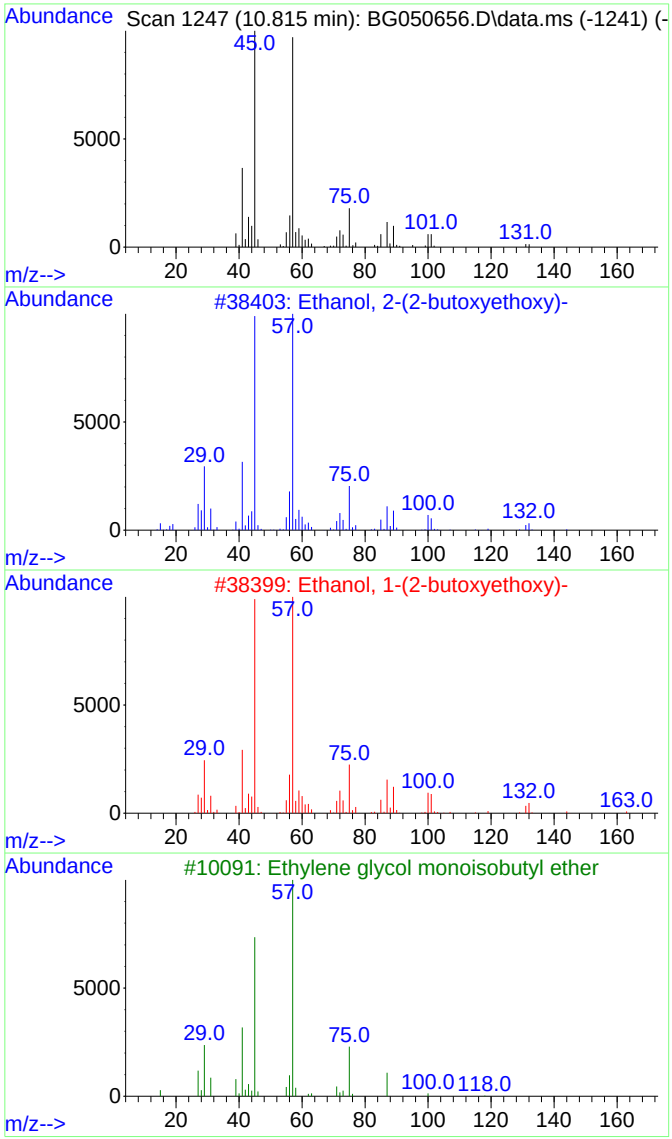
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

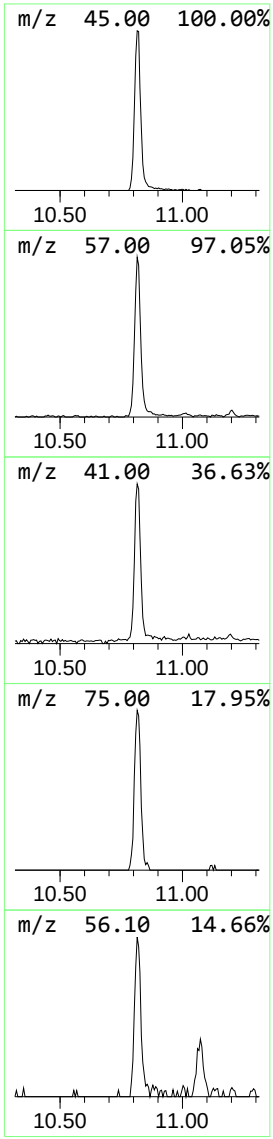
****

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.815	10.14 ng	226050	Naphthalene-d8	11.073

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
3	Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
4	3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	59
5	Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102021\
Data File : BG050656.D
Acq On : 20 Oct 2021 22:17
Operator : CG/JU
Sample : M4288-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SA-01-101921

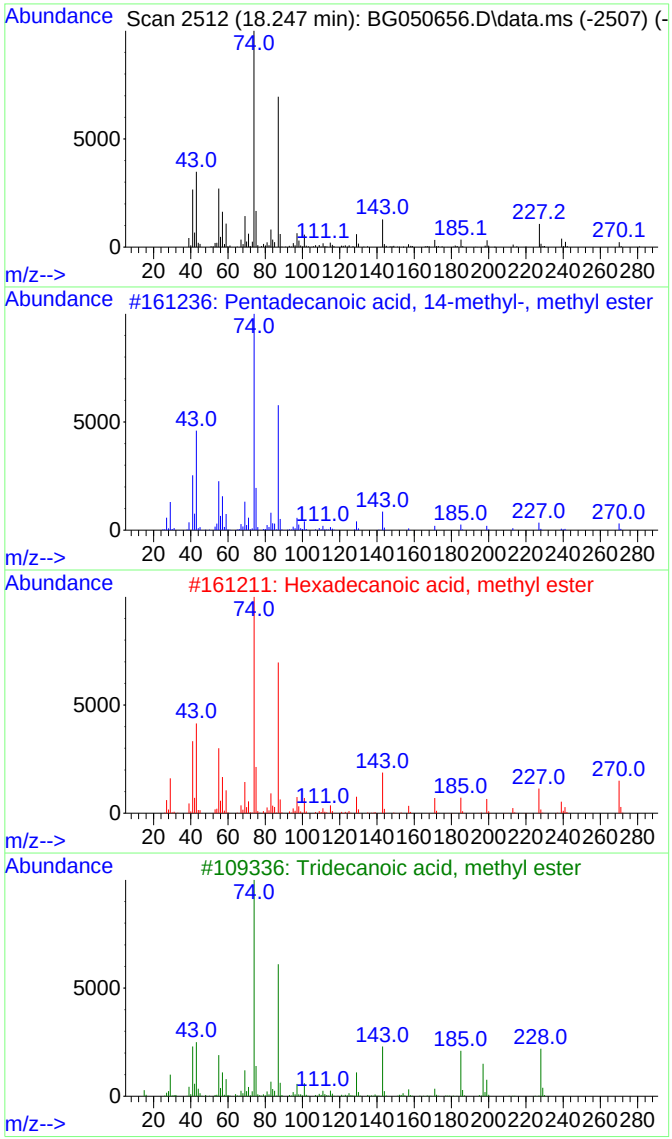
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

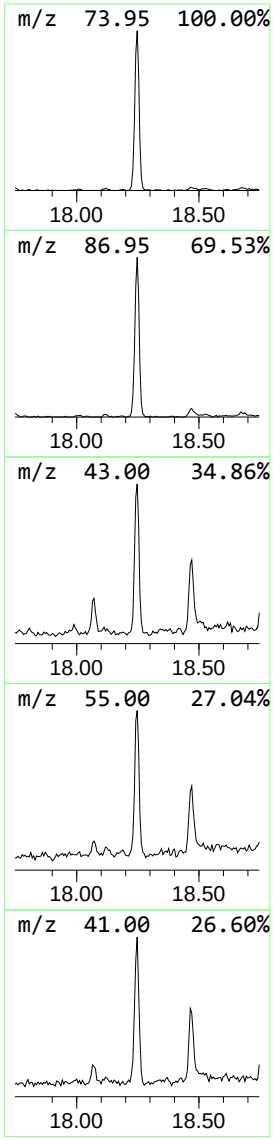
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Pentadecanoic acid, 14-meth... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.247	8.80 ng	296561	Phenanthrene-d10	17.613

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	96
2	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	91
3	Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	87
4	Methyl tetradecanoate	242	C15H30O2	000124-10-7	86
5	Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102021\
Data File : BG050656.D
Acq On : 20 Oct 2021 22:17
Operator : CG/JU
Sample : M4288-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SA-01-101921

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

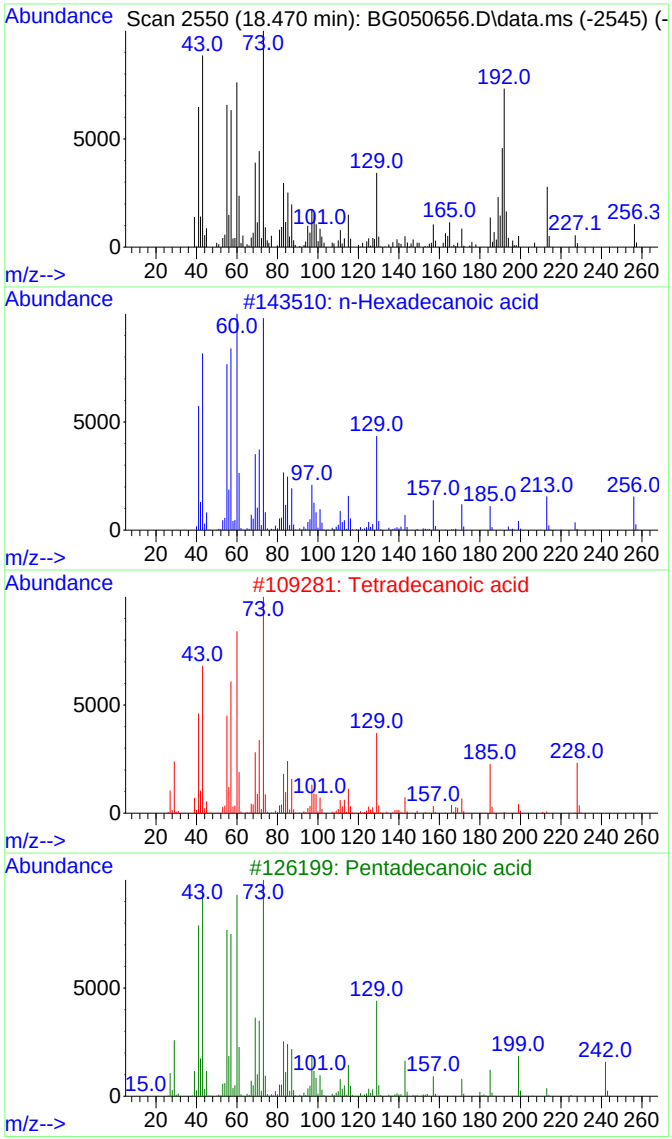
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

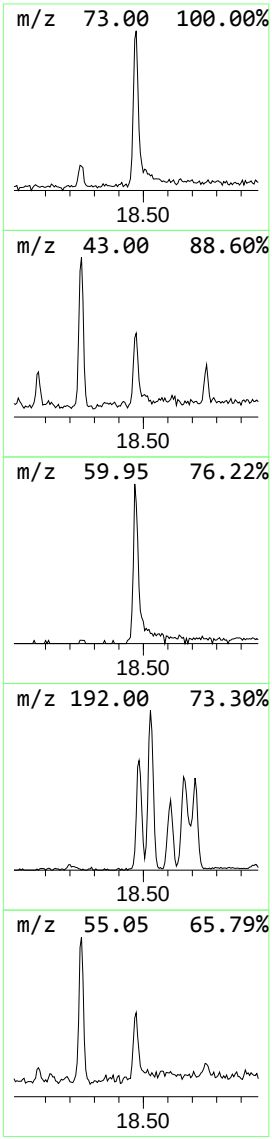
****

Peak Number 3 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.470	6.72 ng	226588	Phenanthrene-d10	17.613

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	70
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	50
4	Tridecanoic acid	214	C13H26O2	000638-53-9	50
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102021\
Data File : BG050656.D
Acq On : 20 Oct 2021 22:17
Operator : CG/JU
Sample : M4288-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SA-01-101921

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

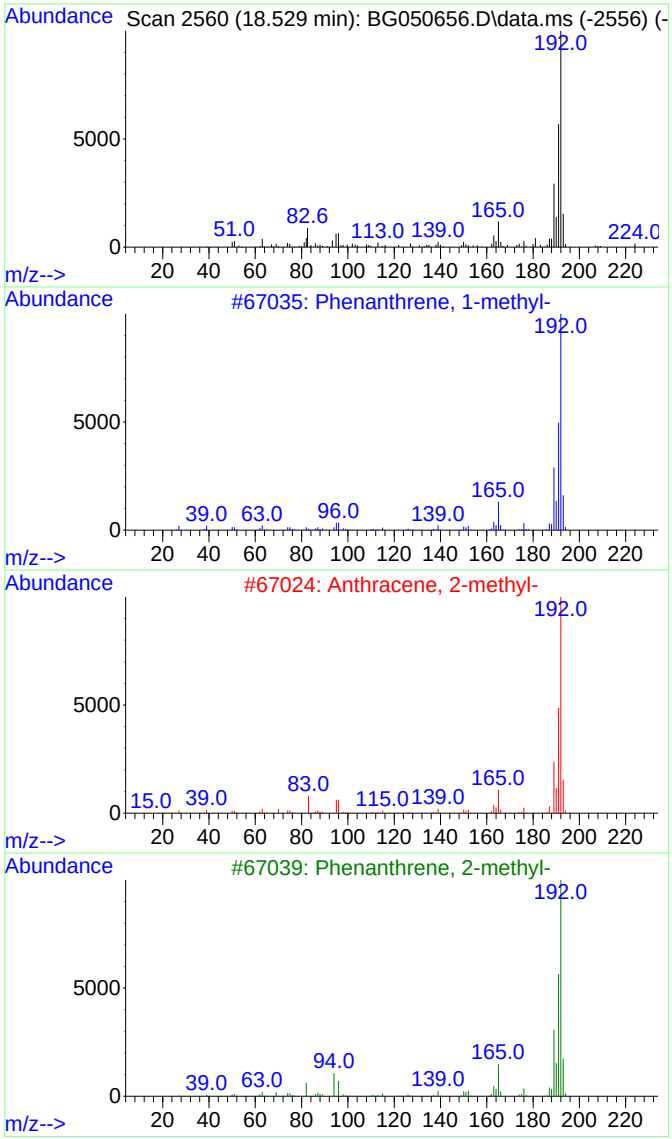
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

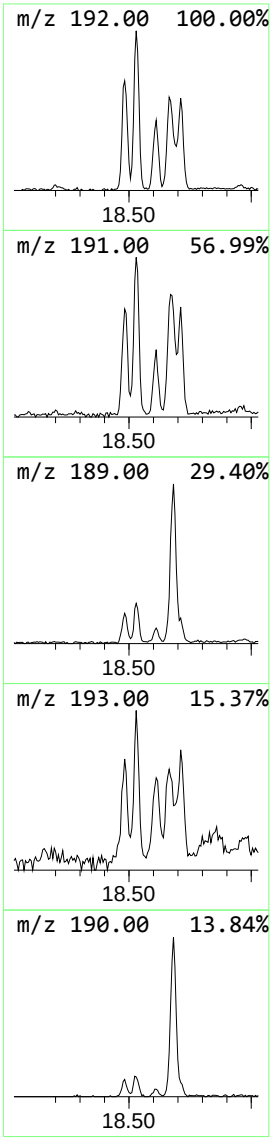
****

Peak Number 4 Phenanthrene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.529	3.67 ng	123890	Phenanthrene-d10	17.613

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	95
5	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102021\
Data File : BG050656.D
Acq On : 20 Oct 2021 22:17
Operator : CG/JU
Sample : M4288-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SA-01-101921

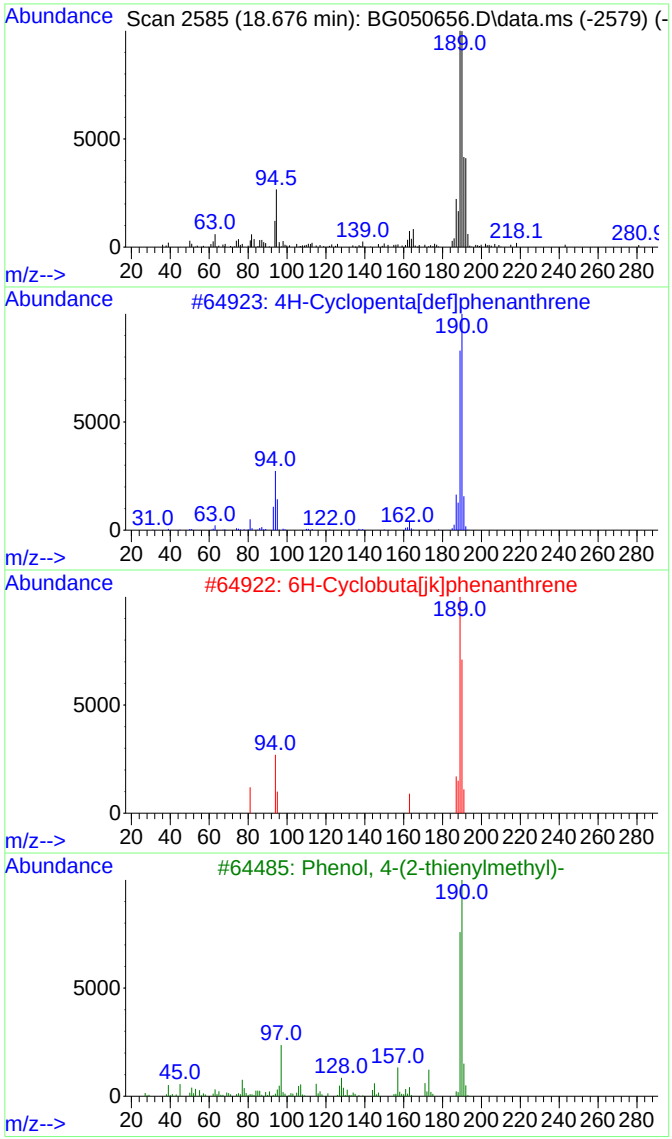
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

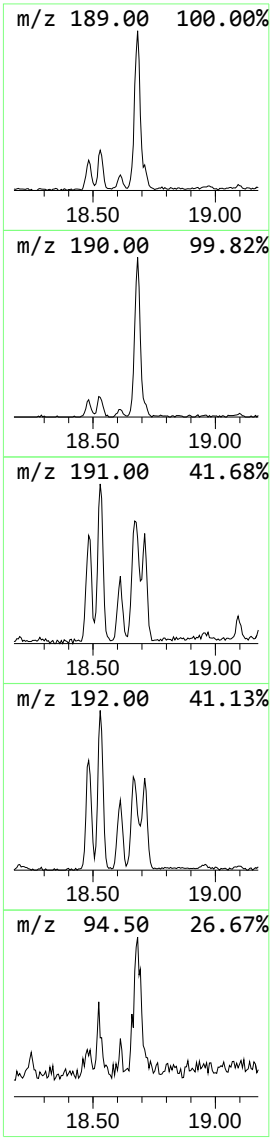
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.676	6.49 ng	218805	Phenanthrene-d10	17.613

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	53
3	Phenol, 4-(2-thienylmethyl)-	190	C11H10OS	091680-55-6	25
4	1,1'-Biphenyl, 4,4'-difluoro-	190	C12H8F2	000398-23-2	23
5	4-Bromothiophene-2-aldehyde	190	C5H3BrOS	018791-75-8	17





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102021\
Data File : BG050656.D
Acq On : 20 Oct 2021 22:17
Operator : CG/JU
Sample : M4288-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SA-01-101921

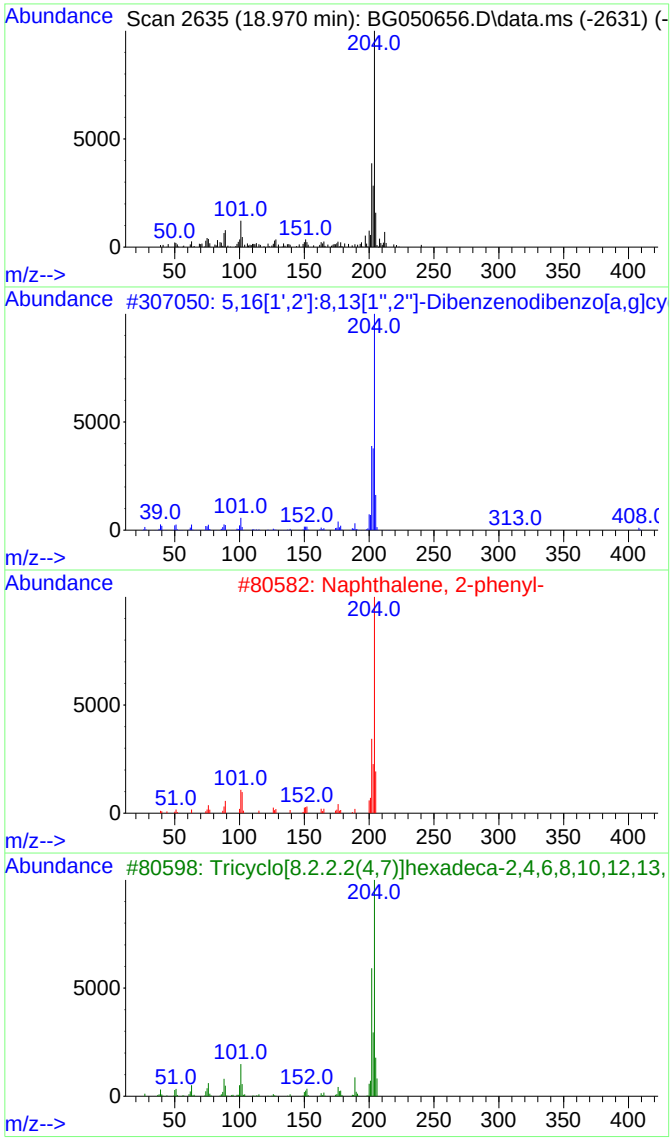
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

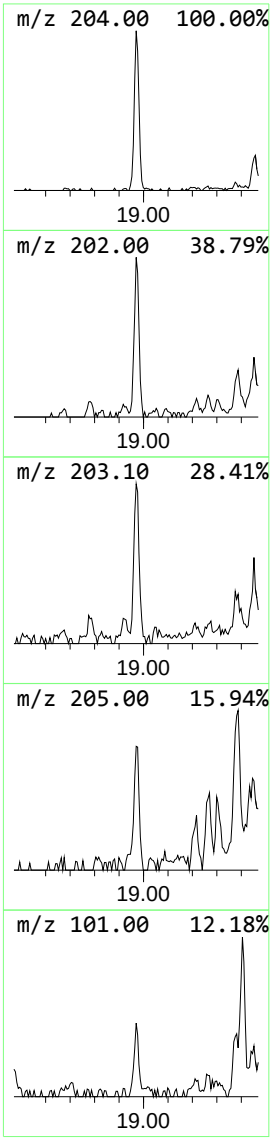
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.970	2.61 ng	87909	Phenanthrene-d10	17.613

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	86
2	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	78
3	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	64
4	Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	55
5	[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	49





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102021\
Data File : BG050656.D
Acq On : 20 Oct 2021 22:17
Operator : CG/JU
Sample : M4288-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SA-01-101921

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

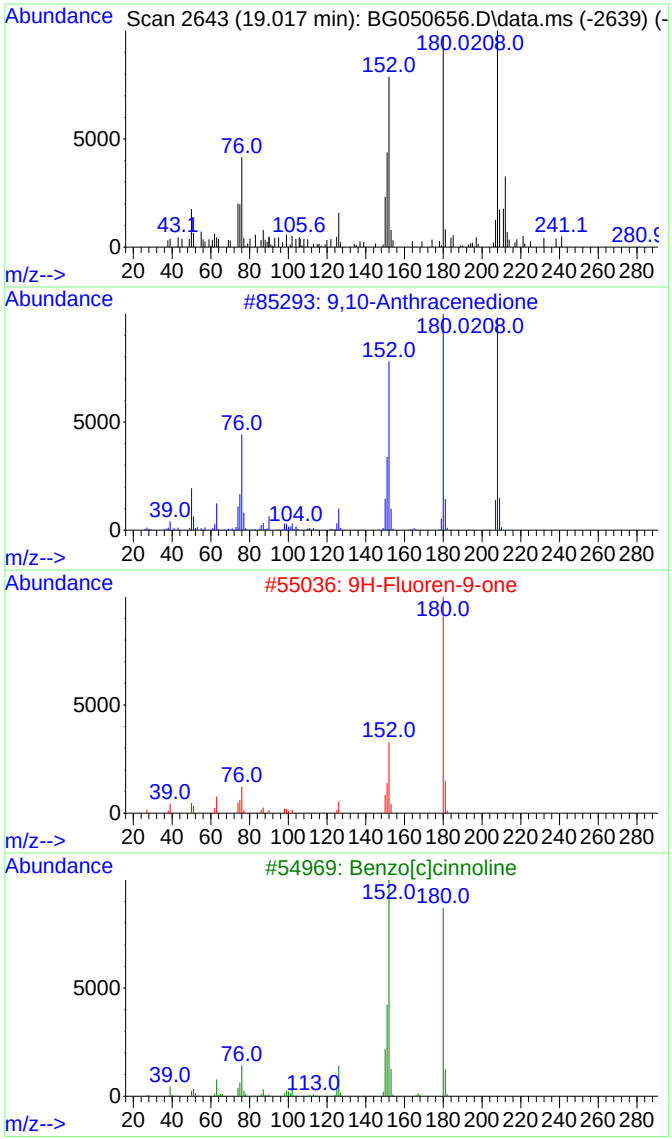
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

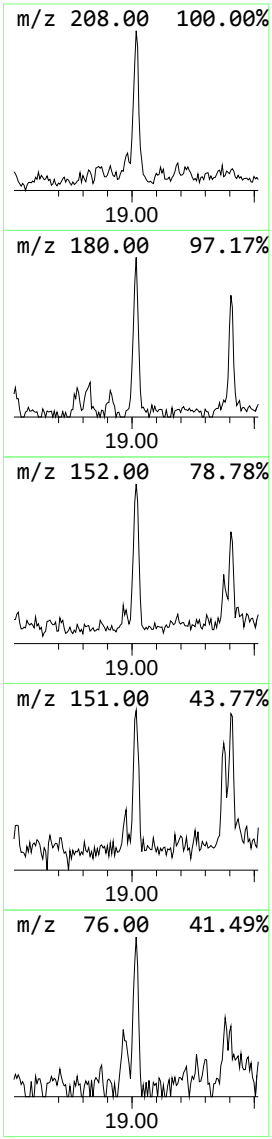
****

Peak Number 7 9,10-Anthracenedione Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.017	2.25 ng	75687	Phenanthrene-d10	17.613

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	93
2	9H-Fluoren-9-one	180	C13H8O	000486-25-9	86
3	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	70
4	Diphenic anhydride	224	C14H8O3	006050-13-1	49
5	Benzo[h]cinnoline	180	C12H8N2	000230-31-9	46





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102021\
Data File : BG050656.D
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Sample : M4288-01
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SA-01-101921

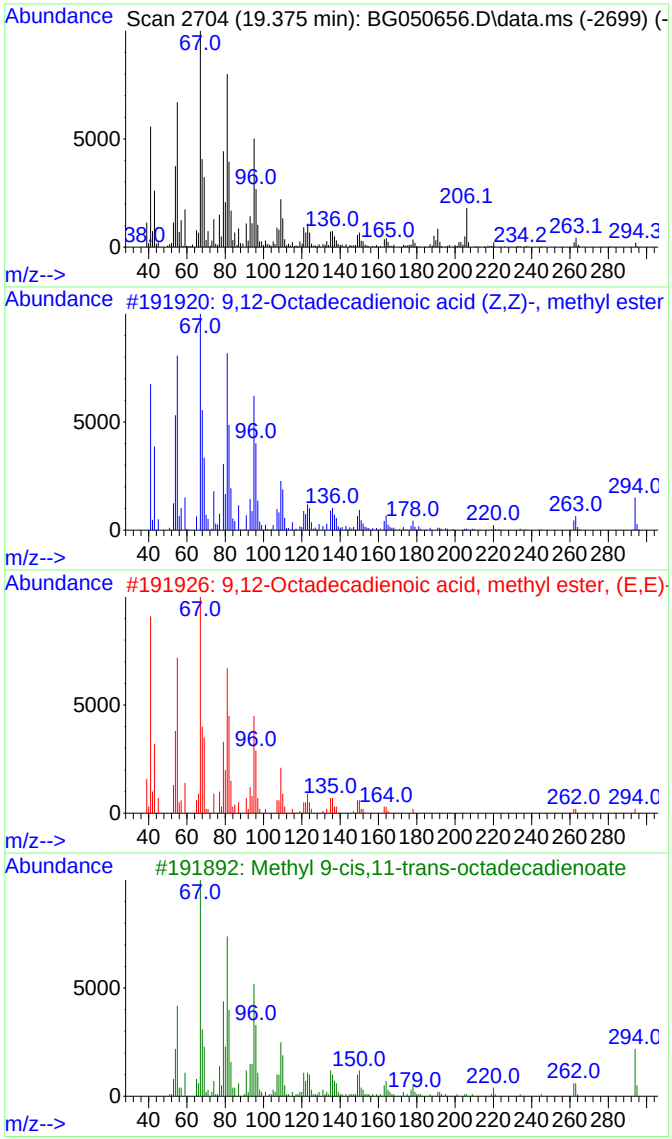
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

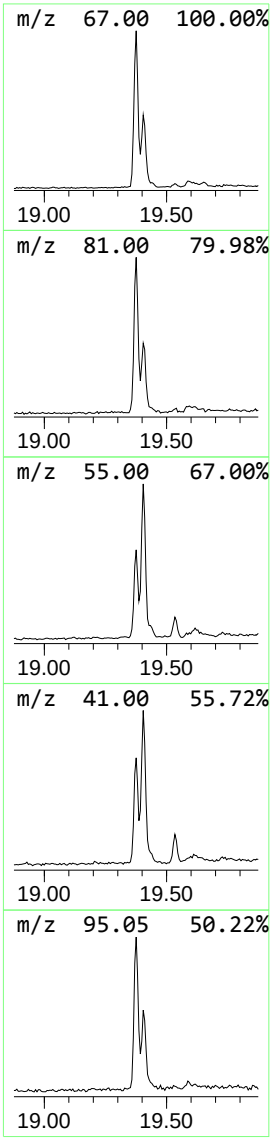
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 9,12-Octadecadienoic acid (... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.375	22.52 ng	759066	Phenanthrene-d10	17.613

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
2	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
3	Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	013058-52-1	99
4	9,11-Octadecadienoic acid, methy...	294	C19H34O2	013038-47-6	98
5	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	98





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102021\
Data File : BG050656.D
Acq On : 20 Oct 2021 22:17
Operator : CG/JU
Sample : M4288-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SA-01-101921

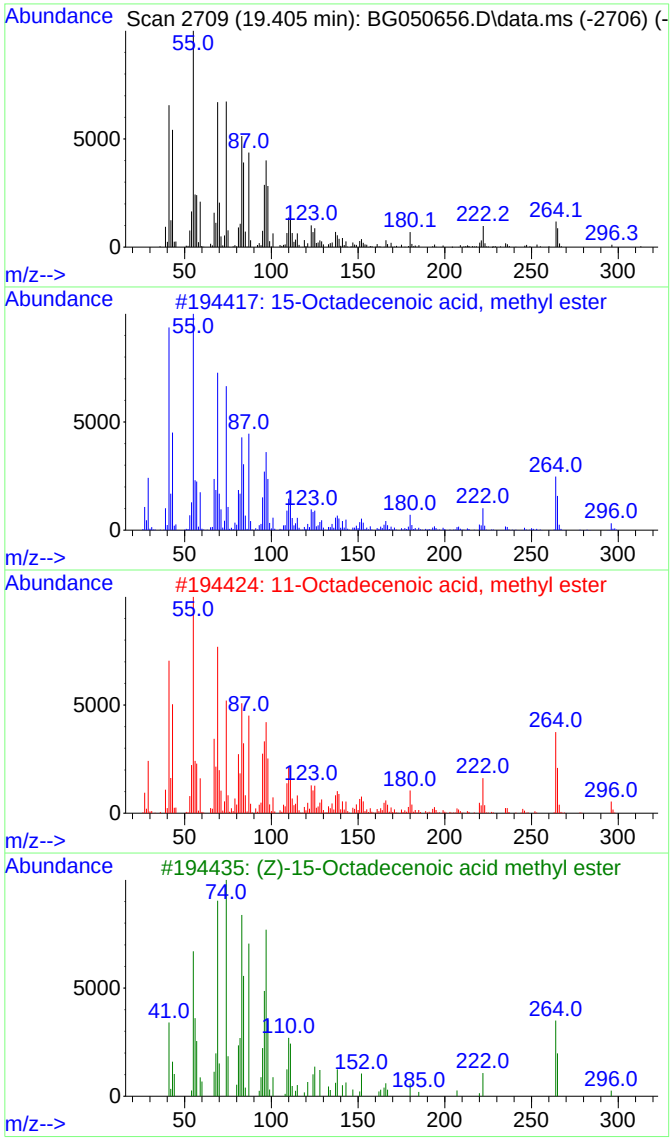
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

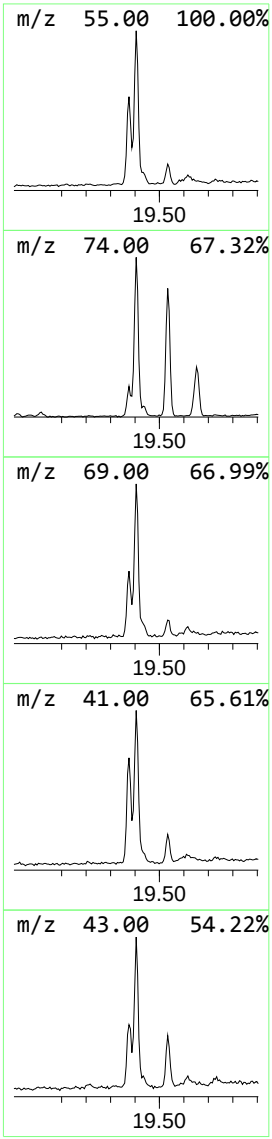
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 15-Octadecenoic acid, methy... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.405	24.94 ng	840939	Phenanthrene-d10	17.613

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	96
2	11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	94
3	(Z)-15-Octadecenoic acid methyl ...	296	C19H36O2	1000427-69-3	94
4	11-Octadecenoic acid, methyl est...	296	C19H36O2	001937-63-9	94
5	7-Octadecenoic acid, methyl ester	296	C19H36O2	057396-98-2	93





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Operator : CG/JU
Sample : M4288-01
Misc :
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Instrument : BNA_G
ClientSampleId : SA-01-101921

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
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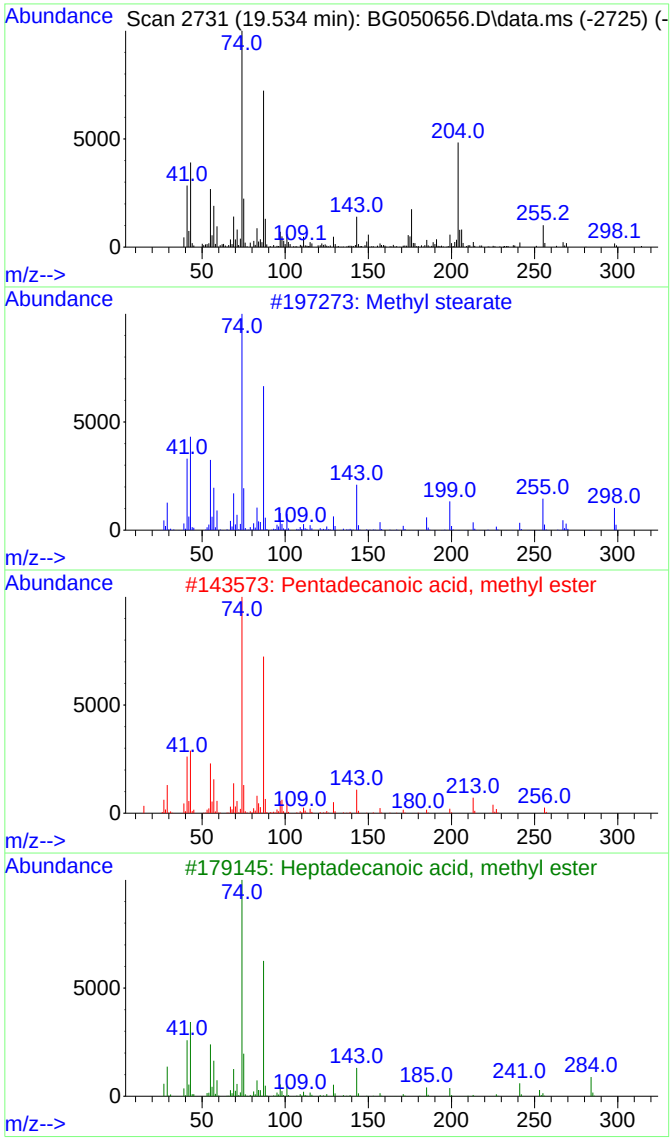
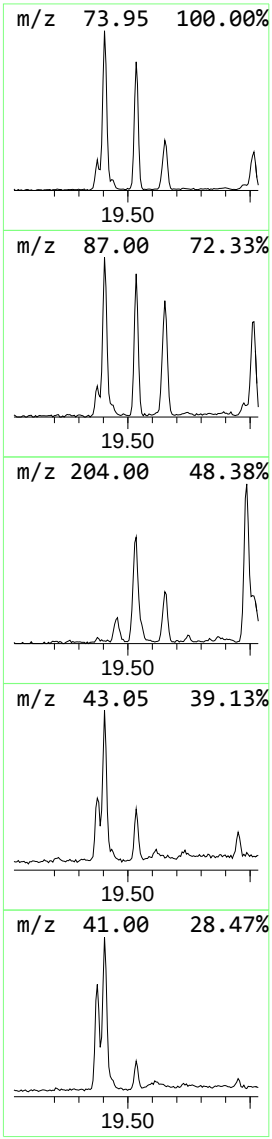
****
Peak Number 10 Methyl stearate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.534	8.82 ng	297238	Phenanthrene-d10	17.613

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl stearate	298	C19H38O2	000112-61-8	86
2	Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	64
3	Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	62
4	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	58
5	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	58





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Sample : M4288-01
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Instrument : BNA_G
ClientSampleId : SA-01-101921

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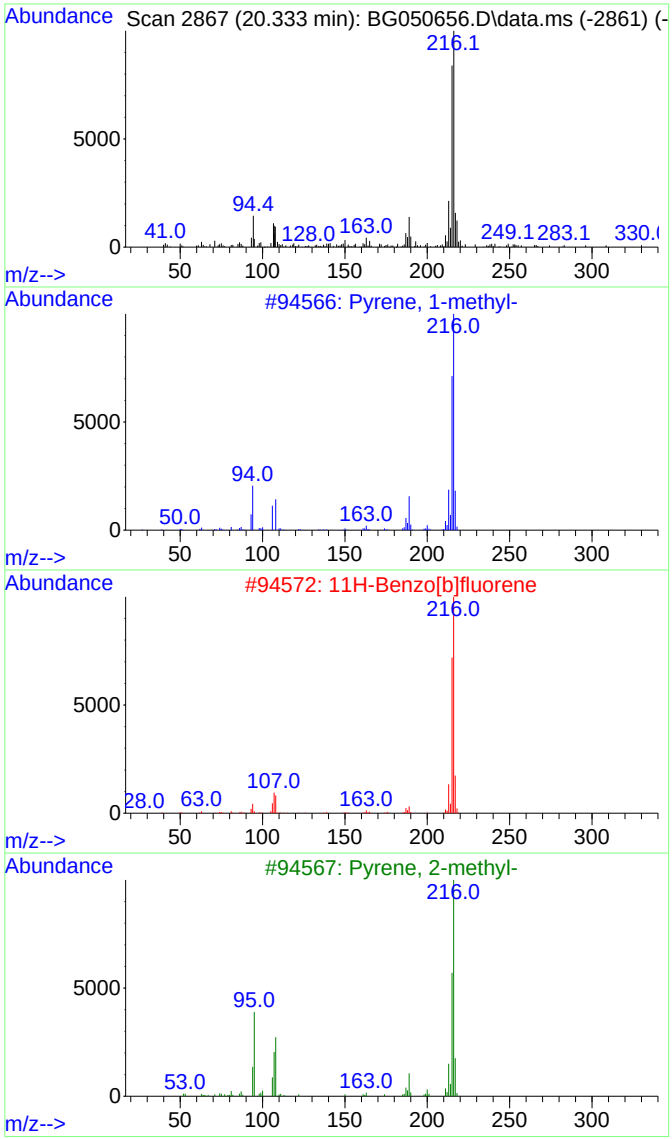
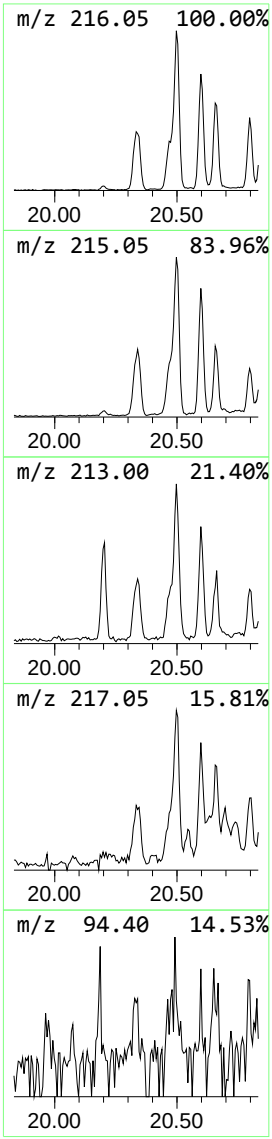
Peak Number 11 Pyrene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.333	2.53 ng	211578	Chrysene-d12	21.913

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	95
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
3	Pyrene, 2-methyl-	216	C17H12	003442-78-2	91
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	90





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Instrument : BNA_G
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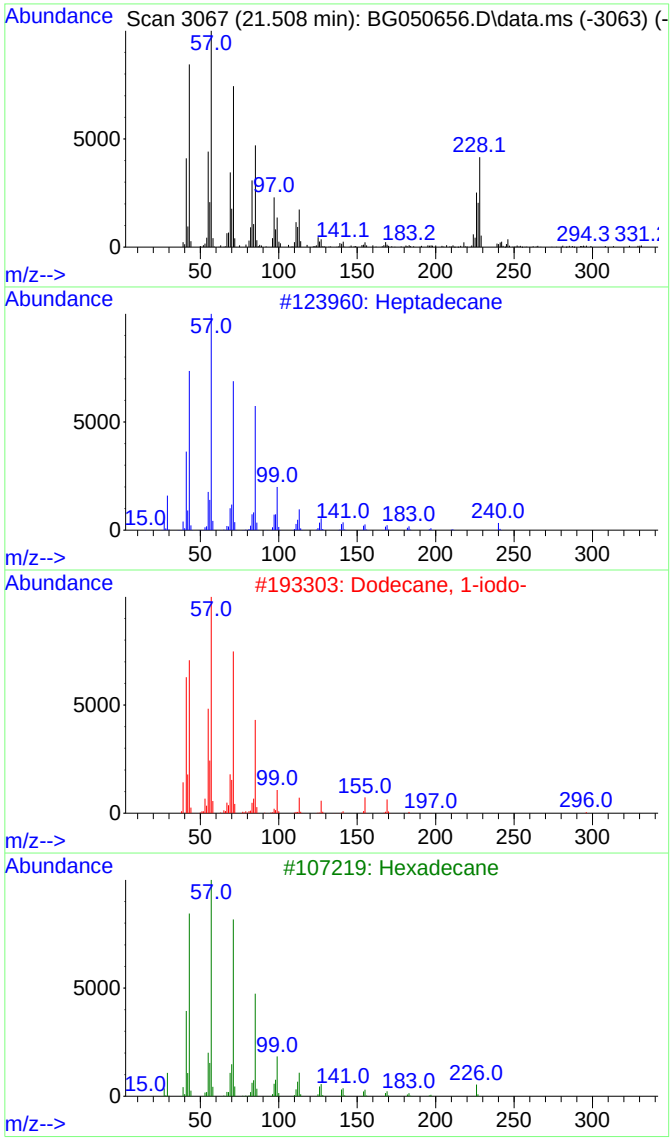
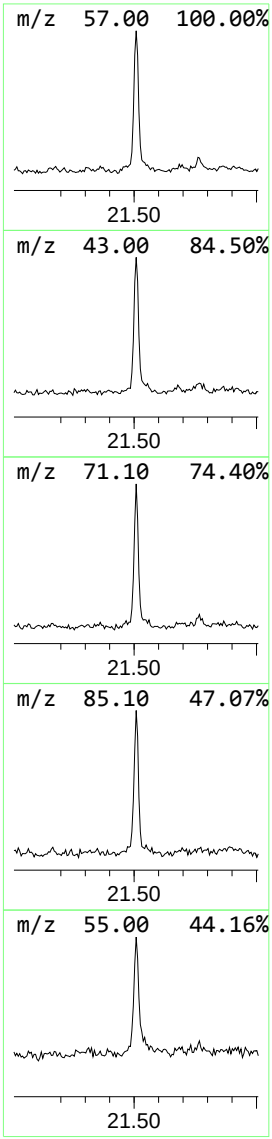
Peak Number 12 Heptadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.508	4.65 ng	388790	Chrysene-d12	21.913

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	86
2	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	64
3	Hexadecane	226	C16H34	000544-76-3	64
4	Carbonic acid, hexadecyl prop-1-...	326	C20H38O3	1000382-90-3	60
5	Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	55





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Instrument : BNA_G
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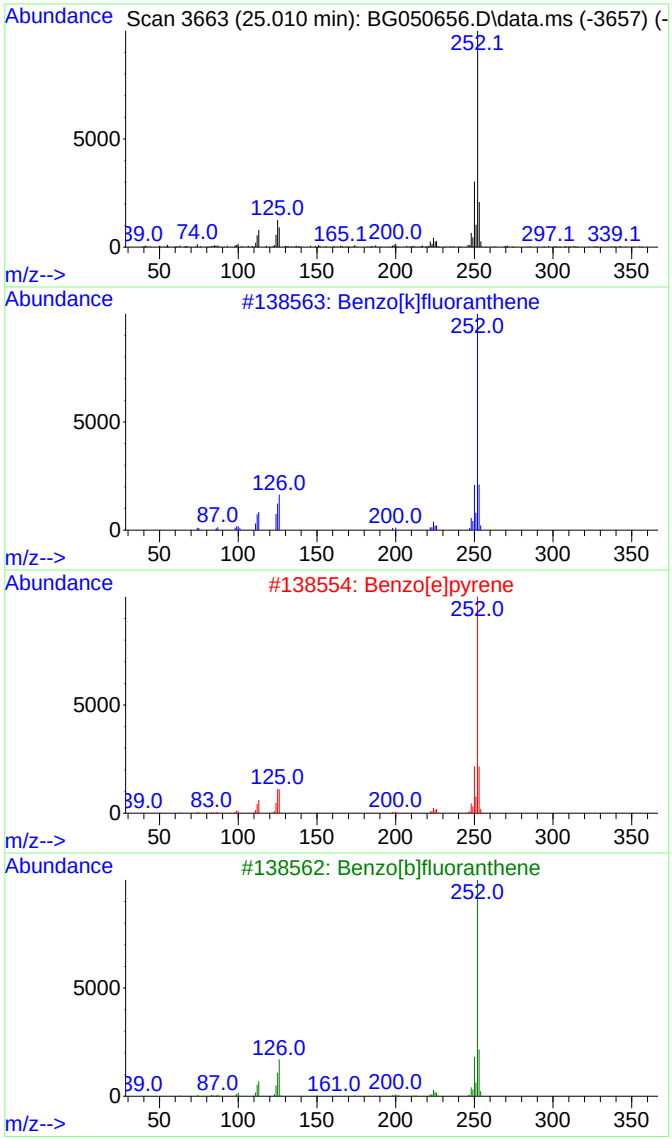
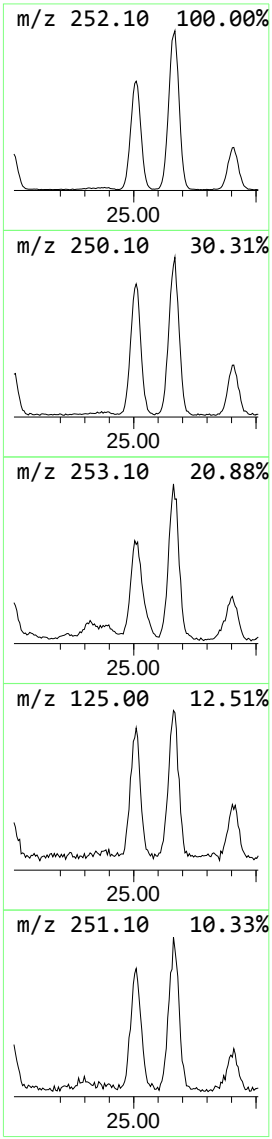
****

Peak Number 13 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.010	23.14 ng	631560	Perylene-d12	25.327

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
2	Benzo[e]pyrene	252	C20H12	000192-97-2	97
3	Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
4	Perylene	252	C20H12	000198-55-0	93
5	Benzo[a]pyrene	252	C20H12	000050-32-8	93





Tentatively Identified Compound (LSC) summary

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TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Ethanol, 2-(2-b...	10.815	10.1	ng	226050	2	11.073	445871	20.0
Pentadecanoic a...	18.247	8.8	ng	296561	4	17.613	674268	20.0
n-Hexadecanoic ...	18.470	6.7	ng	226588	4	17.613	674268	20.0
Phenanthrene, 1...	18.529	3.7	ng	123890	4	17.613	674268	20.0
4H-Cyclopenta[d...	18.676	6.5	ng	218805	4	17.613	674268	20.0
5,16[1',2']:8,1...	18.970	2.6	ng	87909	4	17.613	674268	20.0
9,10-Anthracene...	19.017	2.3	ng	75687	4	17.613	674268	20.0
9,12-Octadecadi...	19.375	22.5	ng	759066	4	17.613	674268	20.0
15-Octadecenoic...	19.405	24.9	ng	840939	4	17.613	674268	20.0
Methyl stearate	19.534	8.8	ng	297238	4	17.613	674268	20.0
Pyrene, 1-methyl-	20.333	2.5	ng	211578	5	21.913	1670780	20.0
Heptadecane	21.508	4.7	ng	388790	5	21.913	1670780	20.0
Benzo[e]pyrene	25.010	23.1	ng	631560	6	25.327	545923	20.0