LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040423\
Data File : BG056977.D
Acq On : 4 Apr 2023 23:04
Operator : CG/JU
Sample : 02176-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TR-04-040323

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-BG033023.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG056977.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.441	317	324	335	rVB	131377	208656	2.35%	0.837%
2	5.923	398	406	415	rBV	326038	536518	6.04%	2.153%
3	7.544	674	682	691	rBV	355545	613767	6.91%	2.463%
4	8.414	823	830	837	rVB	63943	110135	1.24%	0.442%
5	9.588	1022	1030	1042	rVB	204929	385986	4.34%	1.549%
6	11.251	1306	1313	1319	rVB	79860	145095	1.63%	0.582%
7	13.654	1715	1722	1731	rVB	534182	842963	9.48%	3.382%
8	15.022	1950	1955	1962	rVB	125705	205447	2.31%	0.824%
9	15.821	2083	2091	2095	rBV	72566	101970	1.15%	0.409%
10	16.503	2201	2207	2216	rBV	424184	637041	7.17%	2.556%
11	16.679	2232	2237	2240	rBV2	71483	99399	1.12%	0.399%
12	17.766	2415	2422	2425	rBV	153114	248495	2.80%	0.997%
13	17.807	2425	2429	2433	rVB	108704	162720	1.83%	0.653%
14	18.382	2522	2527	2533	rVB	1043305	1340142	15.08%	5.377%
15	18.612	2560	2566	2573	rBV2	169350	291230	3.28%	1.169%
16	18.829	2596	2603	2606	rBV3	53962	111365	1.25%	0.447%
17	19.522	2713	2721	2723	rBV	6731465	8888071	100.00%	35.662%
18	19.546	2723	2725	2734	rVB2	3331883	4075354	45.85%	16.352%
19	19.663	2740	2745	2750	rBV	559576	669726	7.54%	2.687%
20	19.722	2750	2755	2758	rBV	245121	340957	3.84%	1.368%
21	19.792	2763	2767	2773	rVV	334801	530495	5.97%	2.129%
22	19.857	2773	2778	2786	rVB6	65653	124867	1.40%	0.501%
23	20.115	2818	2822	2825	rBV2	61407	97721	1.10%	0.392%
24	20.157	2825	2829	2835	rVB	324410	446423	5.02%	1.791%
25	20.333	2853	2859	2865	rBV	886162	1169351	13.16%	4.692%
26	20.638	2907	2911	2918	rVB	115650	181936	2.05%	0.730%
27	20.738	2923	2928	2932	rBV2	103834	139340	1.57%	0.559%
28	22.066	3148	3154	3162	rBV3	231647	655746	7.38%	2.631%
29	22.136	3162	3166	3176	rVB	181952	330700	3.72%	1.327%
30	24.498	3561	3568	3576	rBV2	114430	399589	4.50%	1.603%
31	25.297	3698	3704	3719	rVB2	78474	246172	2.77%	0.988%
32	25.461	3724	3732	3745	rVB2	115448	365803	4.12%	1.468%
33	25.631	3753	3761	3769	rBV2	77341	220070	2.48%	0.883%

Sum of corrected areas: 24923250

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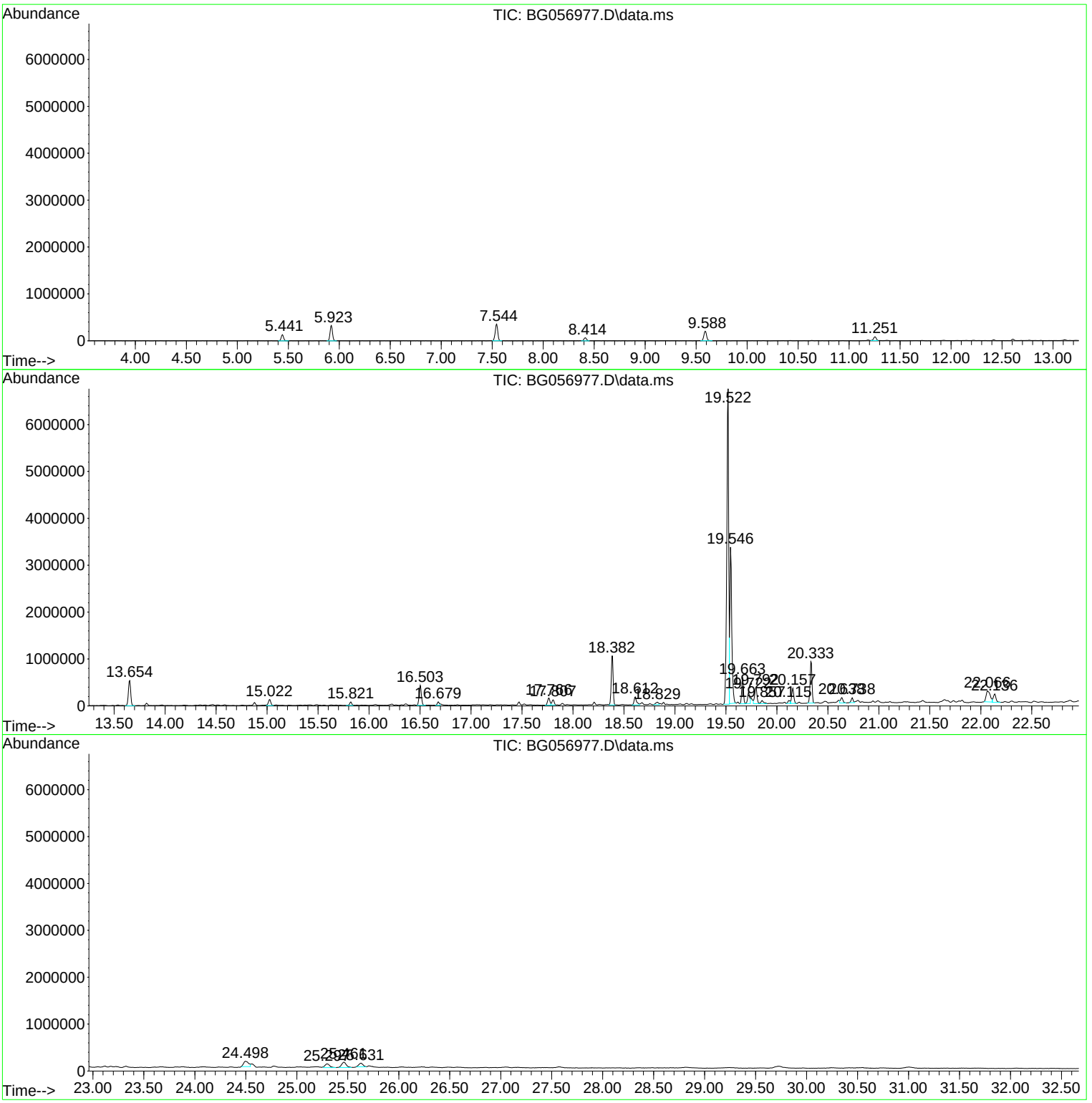
Instrument :
BNA_G
ClientSampleId :
TR-04-040323

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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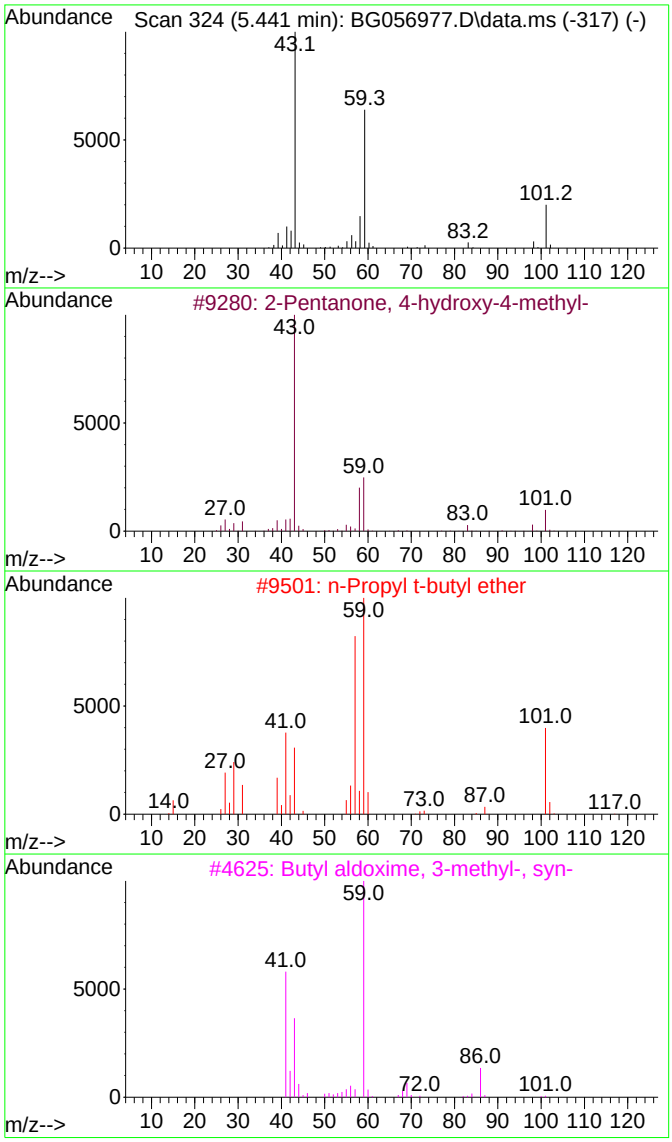
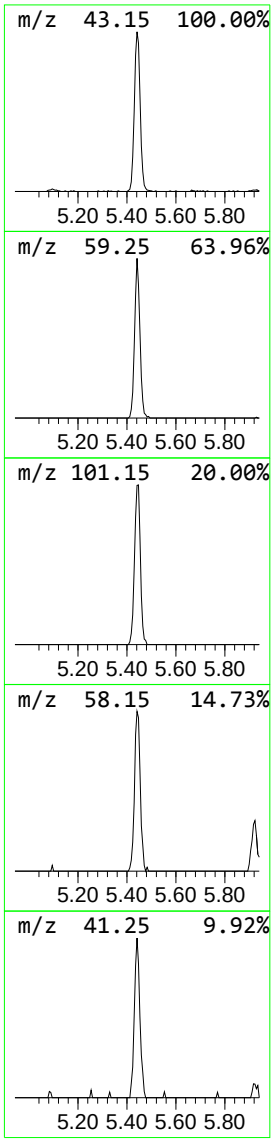
****
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.441	37.89 ng	208656	1,4-Dichlorobenzene-d4	8.414

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	33
3	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
5	4-Methoxy-1-pentene	100	C6H12O	098386-09-5	17





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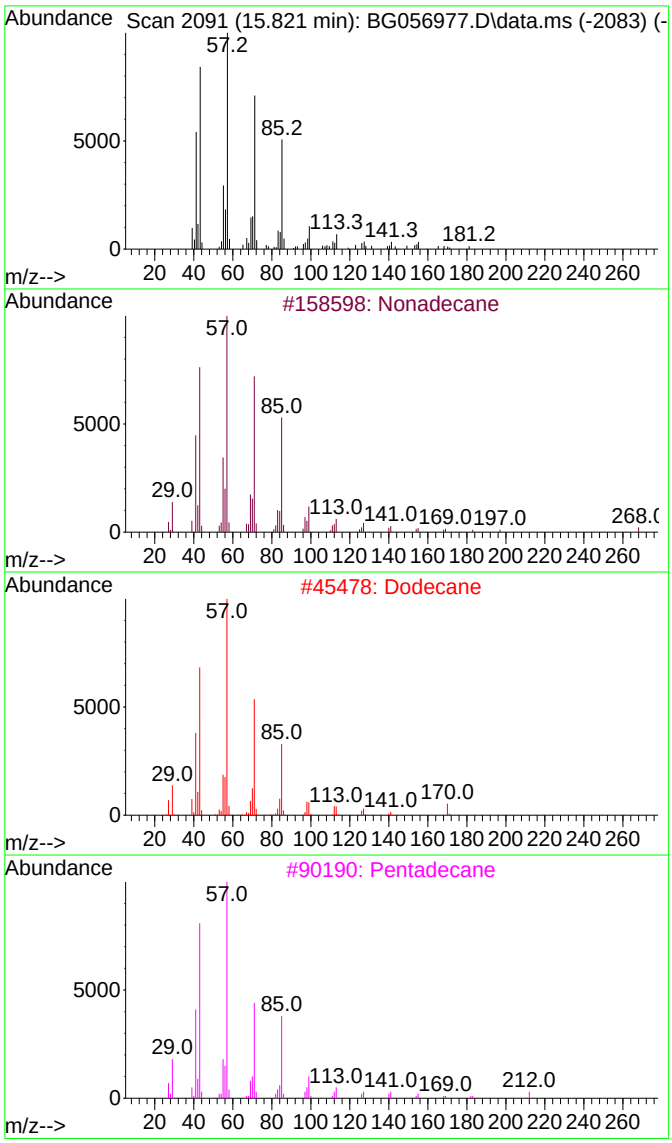
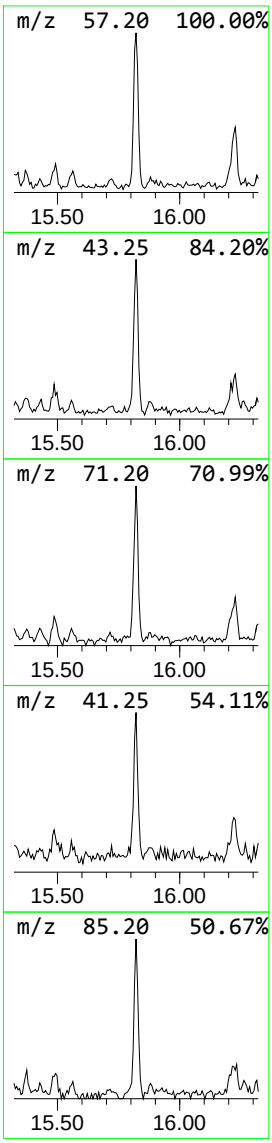
****

Peak Number 2 Nonadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.821	9.93 ng	101970	Acenaphthene-d10	15.028

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	90
2	Dodecane	170	C12H26	000112-40-3	87
3	Pentadecane	212	C15H32	000629-62-9	86
4	Heptadecane	240	C17H36	000629-78-7	86
5	Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	86





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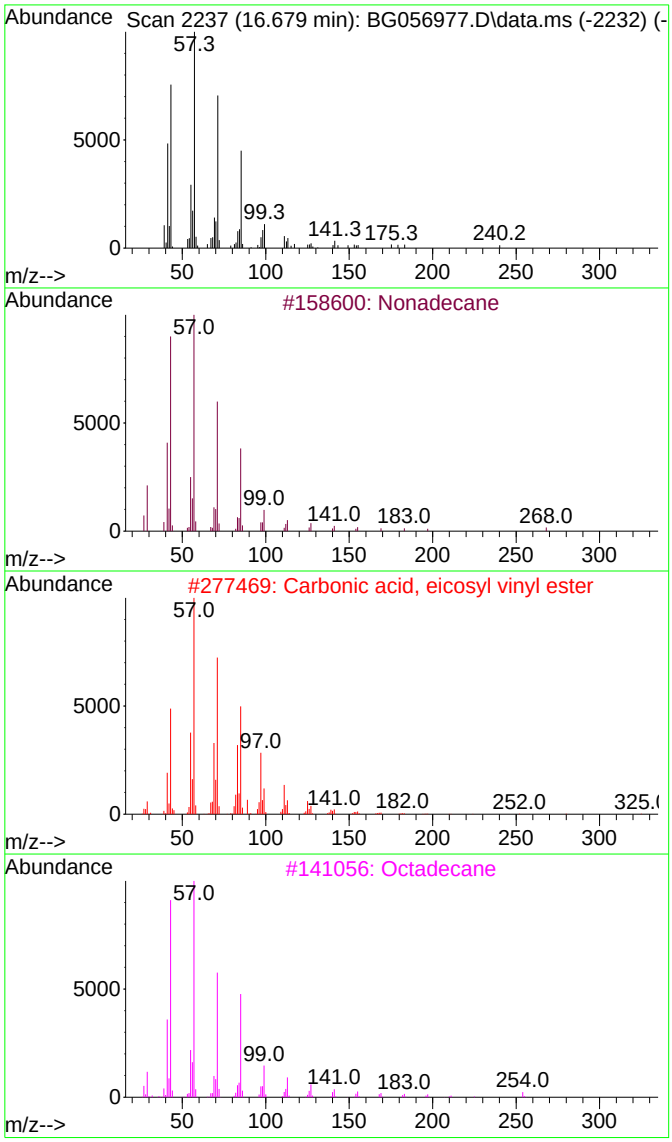
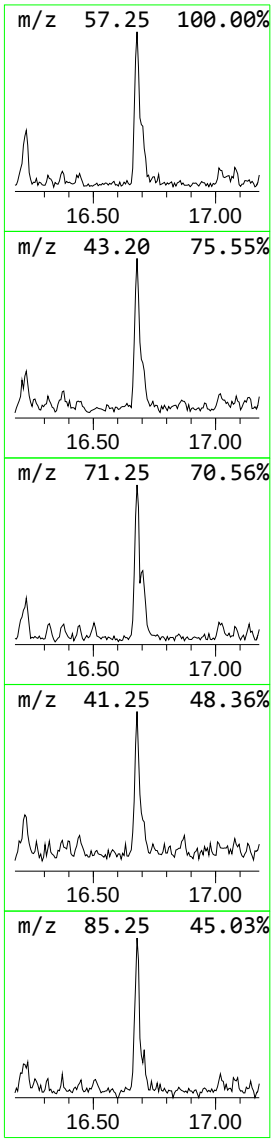
****
Peak Number 3 Carbonic acid, eicosyl viny... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.679	8.00 ng	99399	Phenanthrene-d10	17.760

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	90
2	Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	86
3	Octadecane	254	C18H38	000593-45-3	86
4	Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	86
5	Tetratetracontane	619	C44H90	007098-22-8	86





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Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
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TIC Library : C:\Database\NIST20.L
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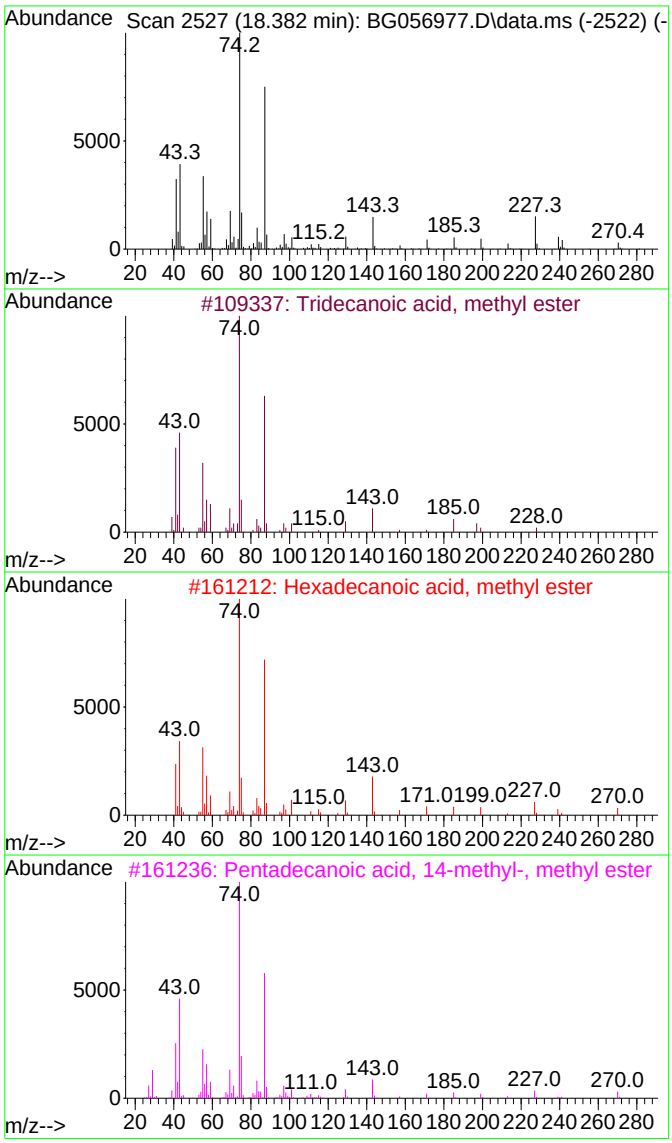
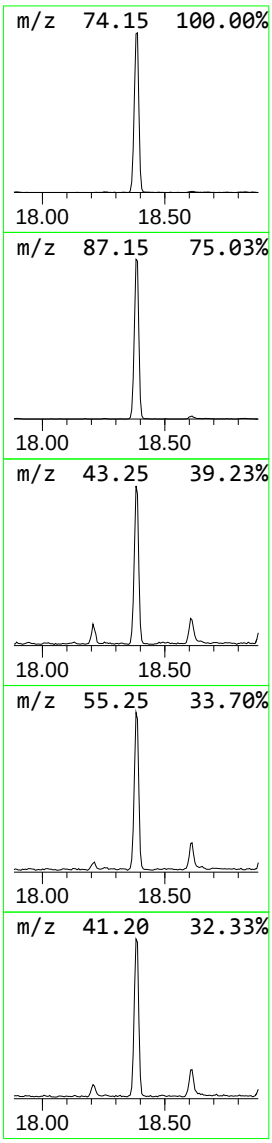
****
Peak Number 4 Tridecanoic acid, methyl ester Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.382	107.86 ng	1340140	Phenanthrene-d10	17.760

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	97
2	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	97
3	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	95
4	Dodecanoic acid, 10-methyl-, met...	228	C14H28O2	005129-65-7	86
5	Hexadecanoic acid, 2-methyl-	270	C17H34O2	027147-71-3	83





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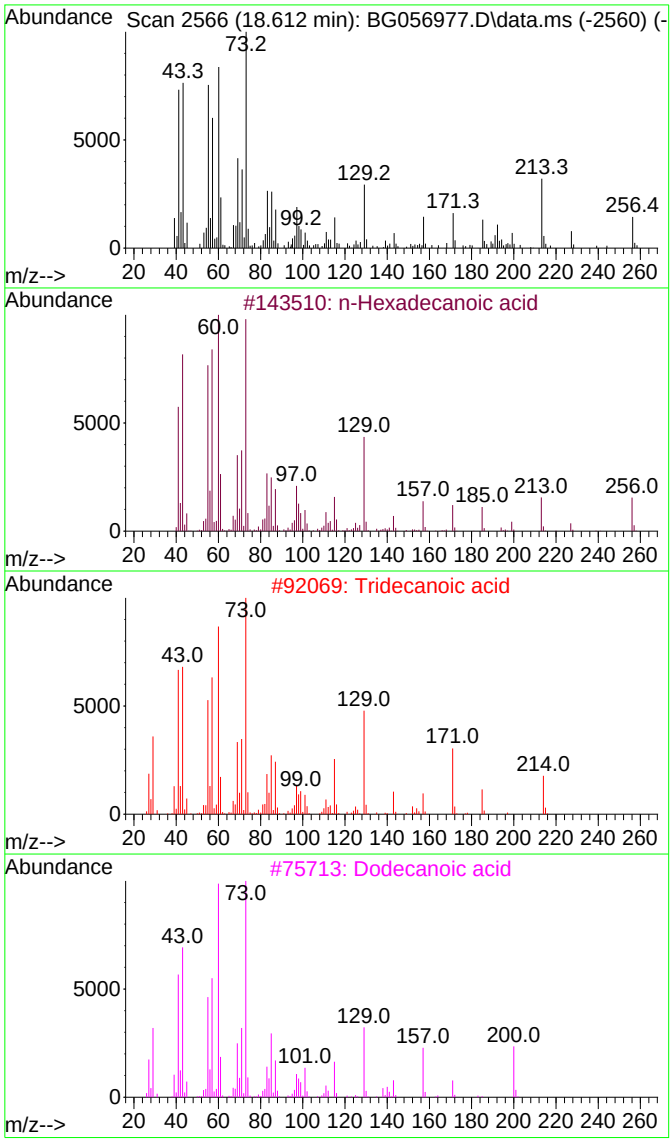
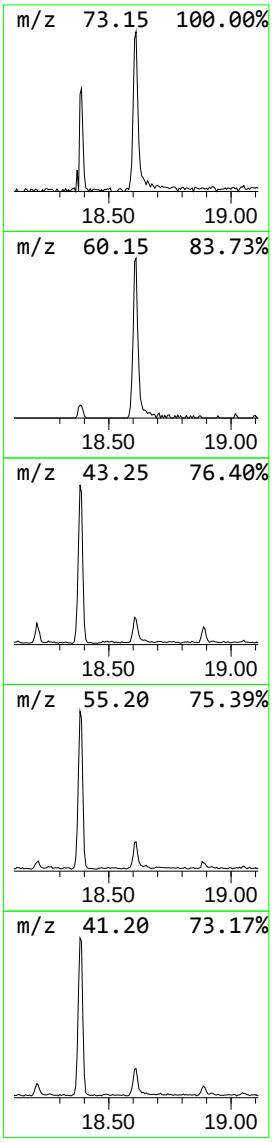
****

Peak Number 5 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.611	23.44 ng	291230	Phenanthrene-d10	17.760

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tridecanoic acid	214	C13H26O2	000638-53-9	93
3	Dodecanoic acid	200	C12H24O2	000143-07-7	70
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
5	n-Decanoic acid	172	C10H20O2	000334-48-5	53





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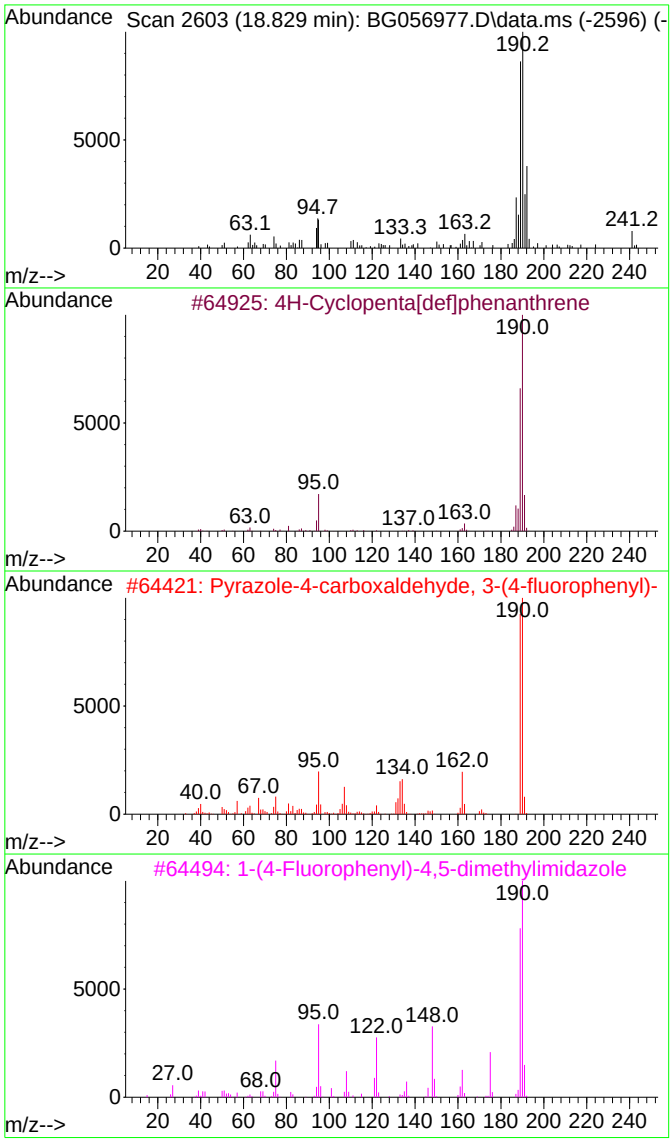
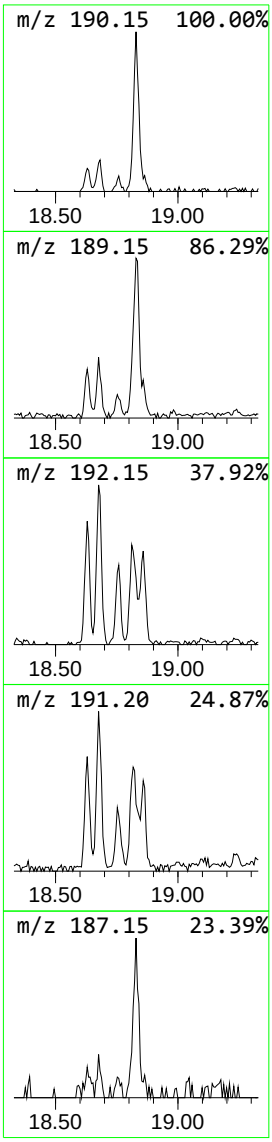
****
Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.829	8.96 ng	111365	Phenanthrene-d10	17.760

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	83
2	Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	50
3	1-(4-Fluorophenyl)-4,5-dimethyli...	190	C11H11FN2	1000443-18-7	42
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	40
5	Pyrimidine, 2-chloro-5-phenyl-	190	C10H7ClN2	022536-62-5	30





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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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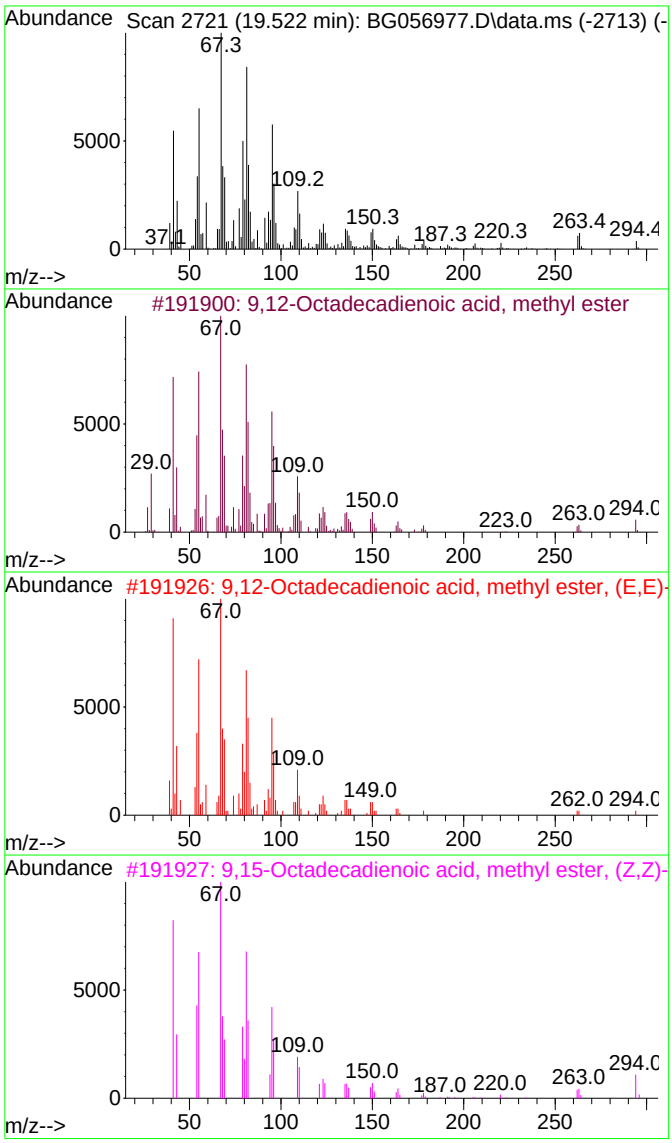
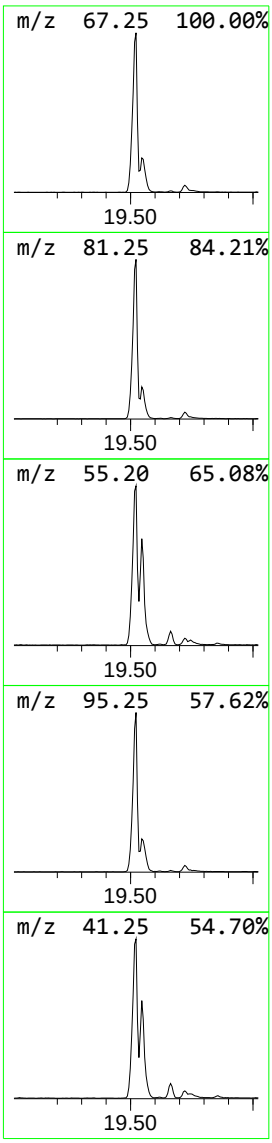
****
Peak Number 7 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.522	715.35 ng	8888070	Phenanthrene-d10	17.760

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99
2	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
3	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
4	Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	013058-52-1	98
5	10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	96





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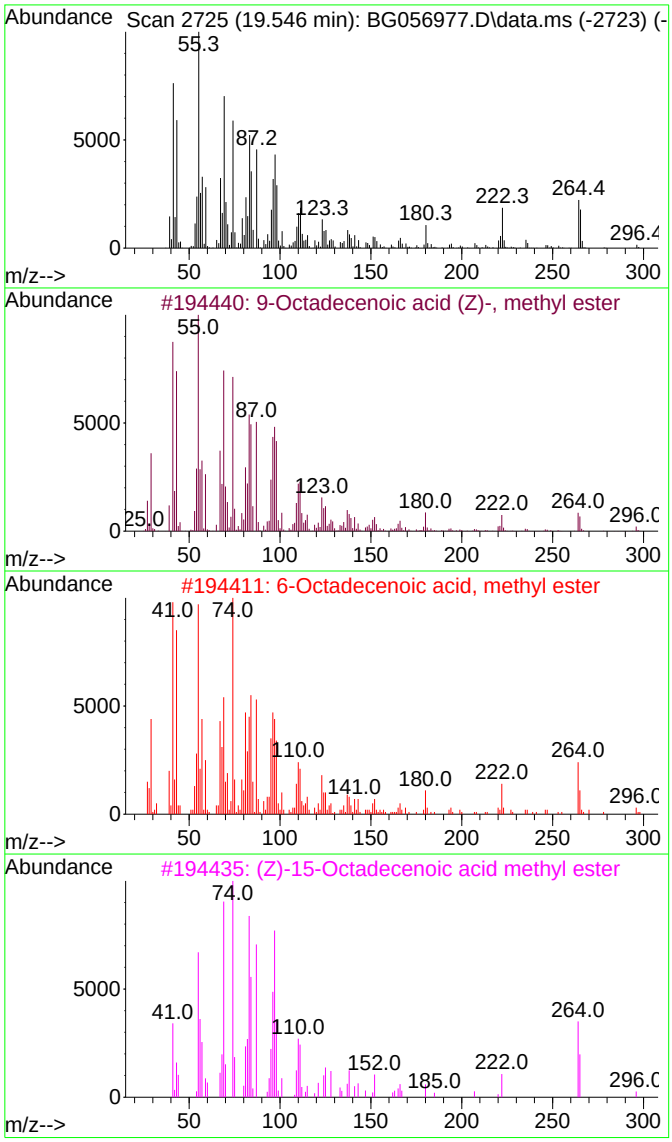
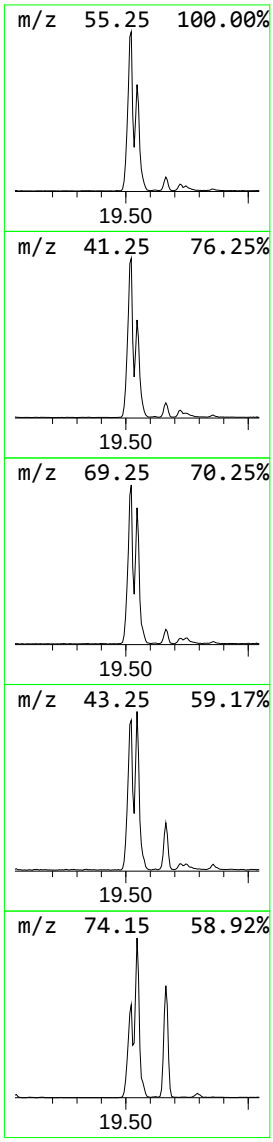
****
Peak Number 8 9-Octadecenoic acid (Z)-, m... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.546	328.00 ng	4075350	Phenanthrene-d10	17.760

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2	6-Octadecenoic acid, methyl ester	296	C19H36O2	052355-31-4	96
3	(Z)-15-Octadecenoic acid methyl ...	296	C19H36O2	1000427-69-3	95
4	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	94
5	10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	94





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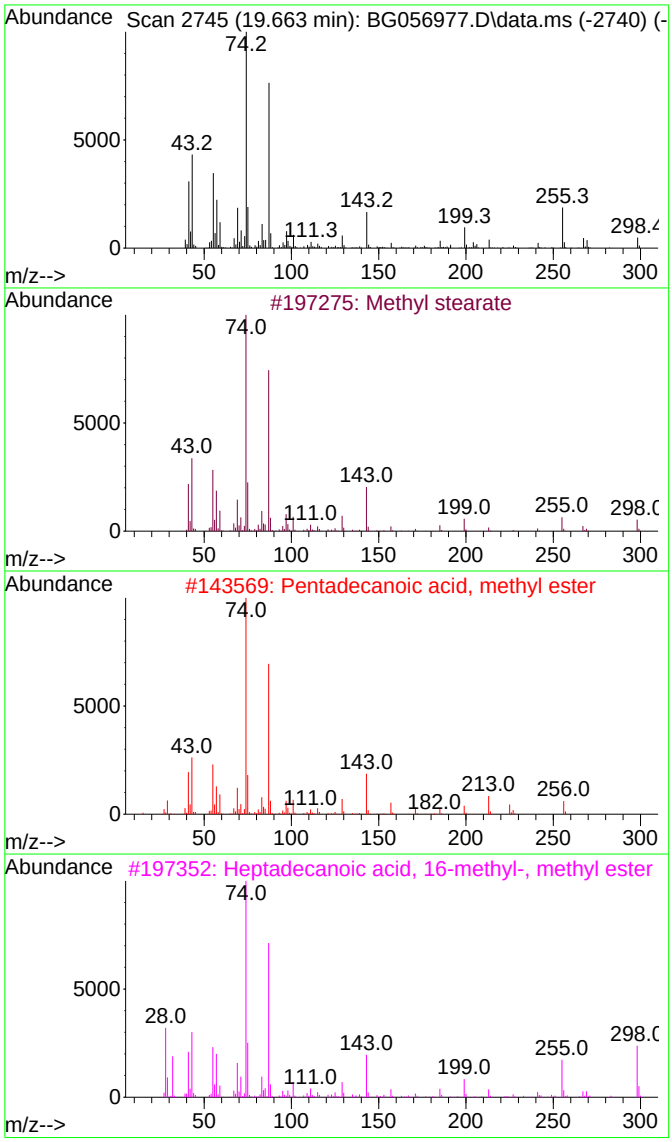
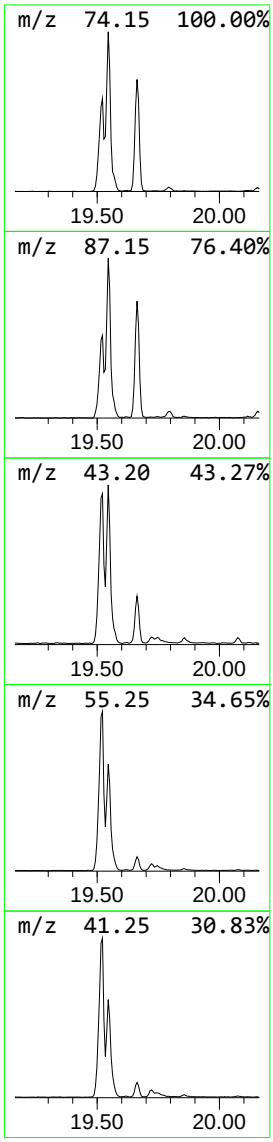
****

Peak Number 9 Methyl stearate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.663	53.90 ng	669726	Phenanthrene-d10	17.760

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl stearate	298	C19H38O2	000112-61-8	97
2	Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	91
3	Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	91
4	2-Methylpentadecanoic acid	256	C16H32O2	025354-92-1	83
5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	81





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Data File : BG056977.D
Acq On : 4 Apr 2023 23:04
Operator : CG/JU
Sample : 02176-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TR-04-040323

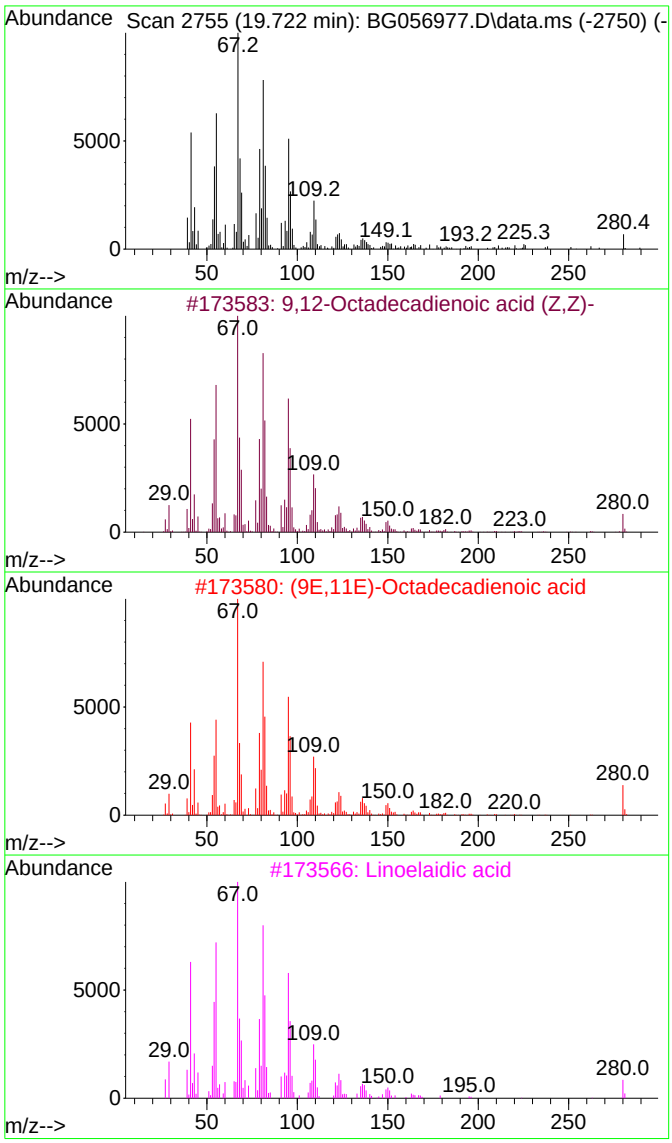
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

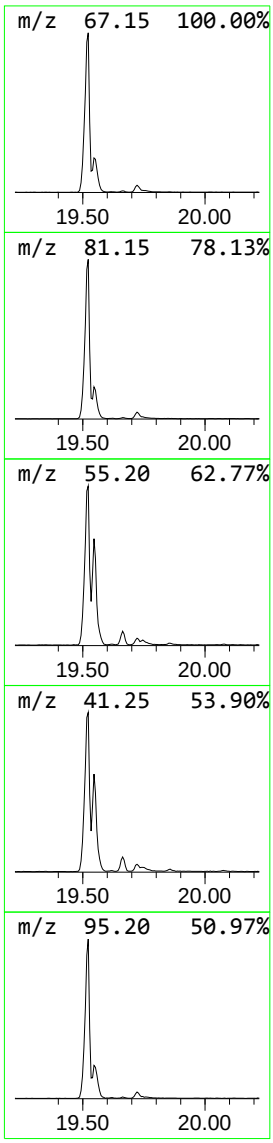
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 9,12-Octadecadienoic acid (... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.722	27.44 ng	340957	Phenanthrene-d10	17.760

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	96
2	(9E,11E)-Octadecadienoic acid	280	C18H32O2	000544-71-8	94
3	Linoelaidic acid	280	C18H32O2	000506-21-8	93
4	10E,12Z-Octadecadienoic acid	280	C18H32O2	002420-56-6	93
5	Methyl 9,12-heptadecadienoate	280	C18H32O2	1000336-36-2	90





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Operator : CG/JU
Sample : 02176-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TR-04-040323

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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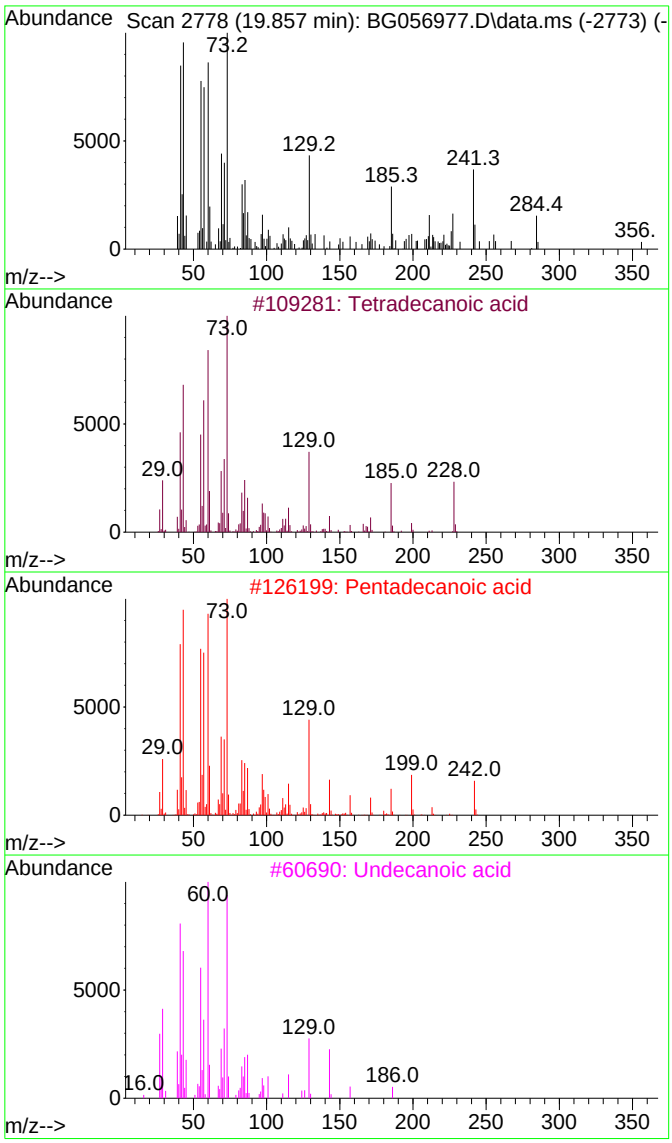
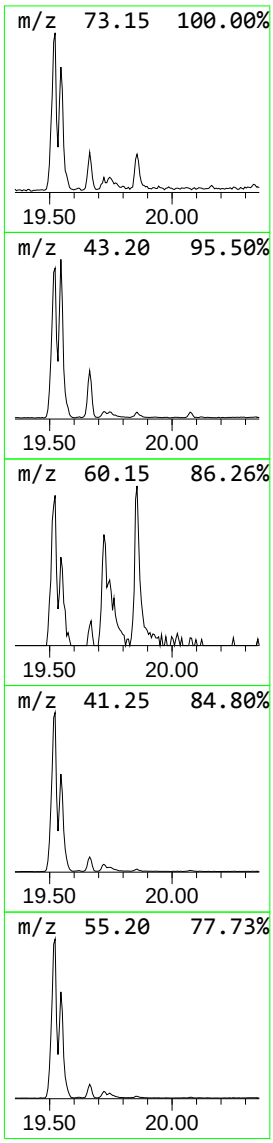
Peak Number 11 Tetradecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.857	10.05 ng	124867	Phenanthrene-d10	17.760

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecanoic acid	228	C14H28O2	000544-63-8	87
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	86
3	Undecanoic acid	186	C11H22O2	000112-37-8	83
4	Tridecanoic acid	214	C13H26O2	000638-53-9	81
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	74





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Operator : CG/JU
Sample : 02176-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TR-04-040323

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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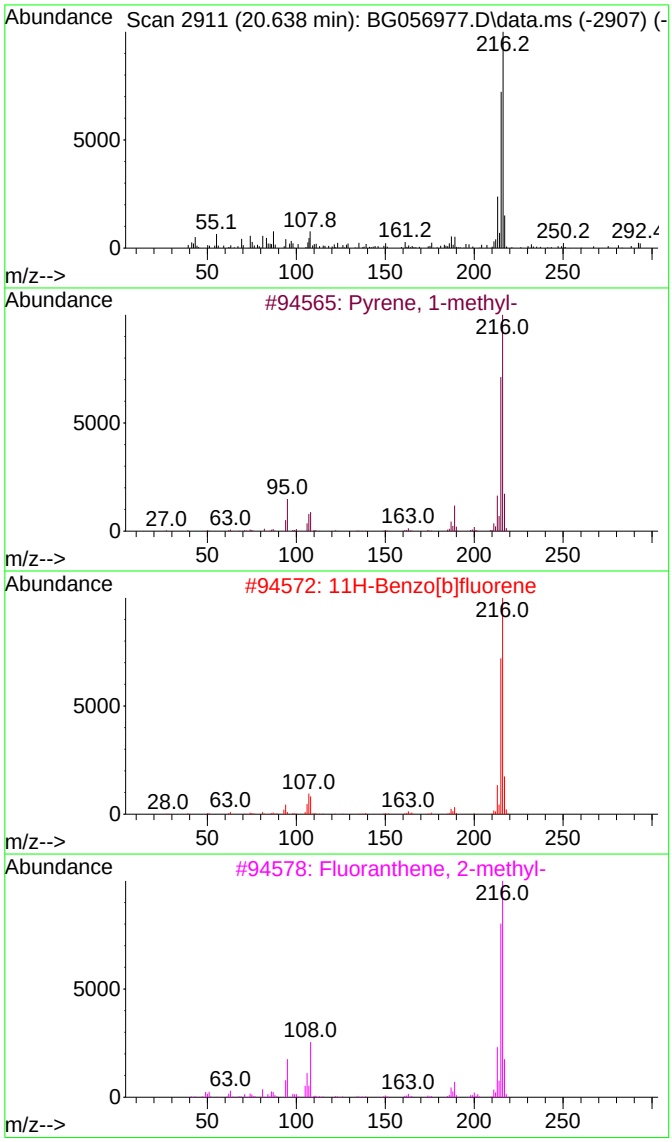
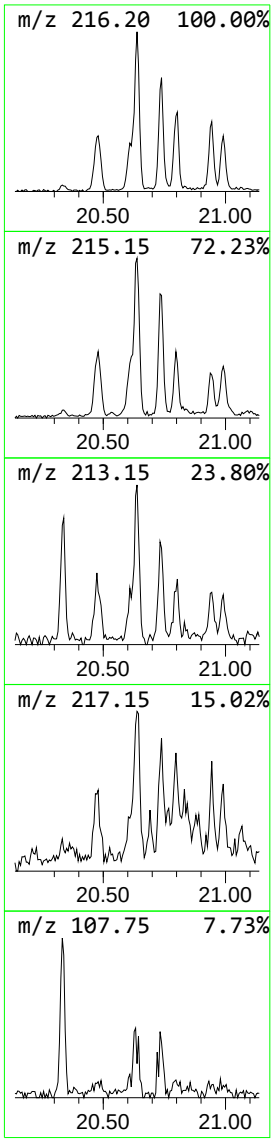
Peak Number 12 Pyrene, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.638	5.55 ng	181936	Chrysene-d12	22.083

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	80
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	74
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	74
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	74
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	59





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Acq On : 4 Apr 2023 23:04
Operator : CG/JU
Sample : 02176-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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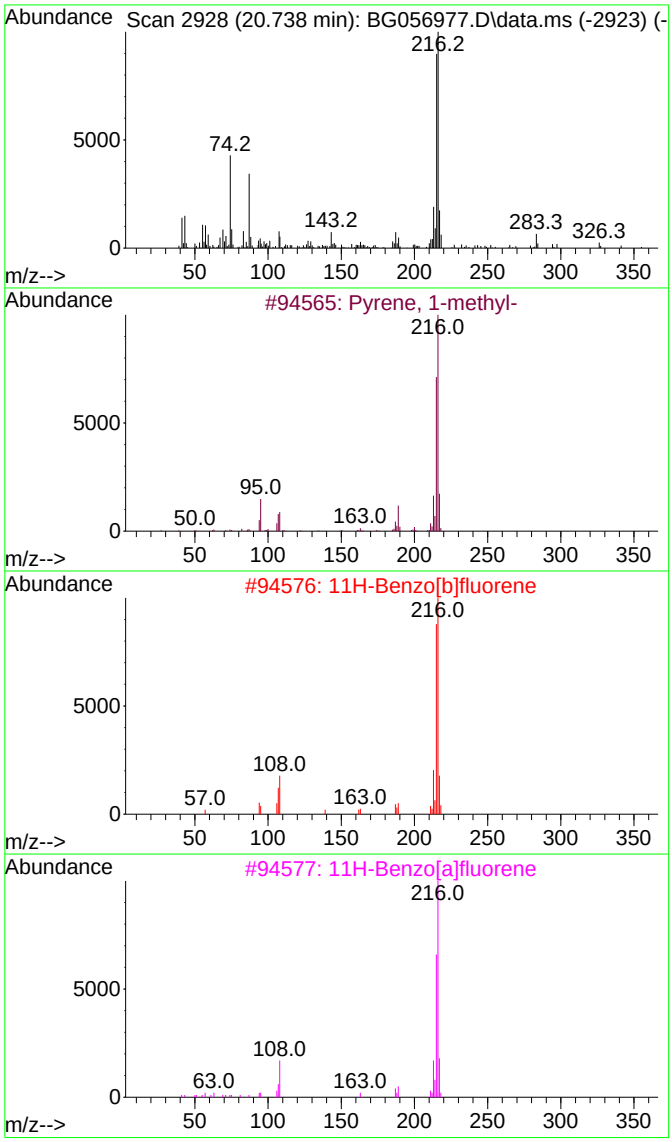
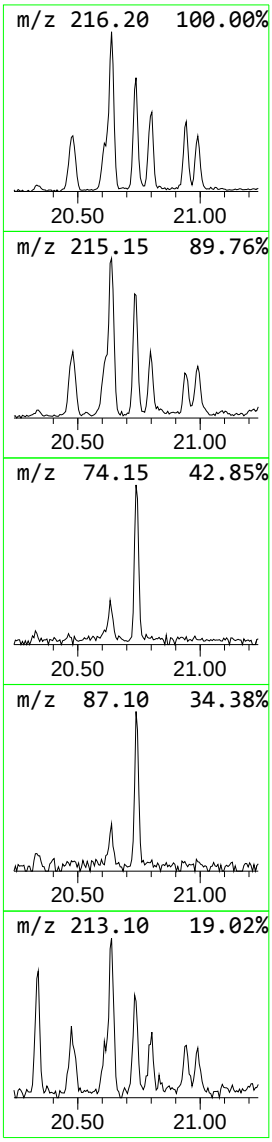
Peak Number 13 11H-Benzo[b]fluorene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.738	4.25 ng	139340	Chrysene-d12	22.083

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	76
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	70
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	70
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	64
5	7H-Benzanthrene	216	C17H12	000199-94-0	53





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Operator : CG/JU
Sample : 02176-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TR-04-040323

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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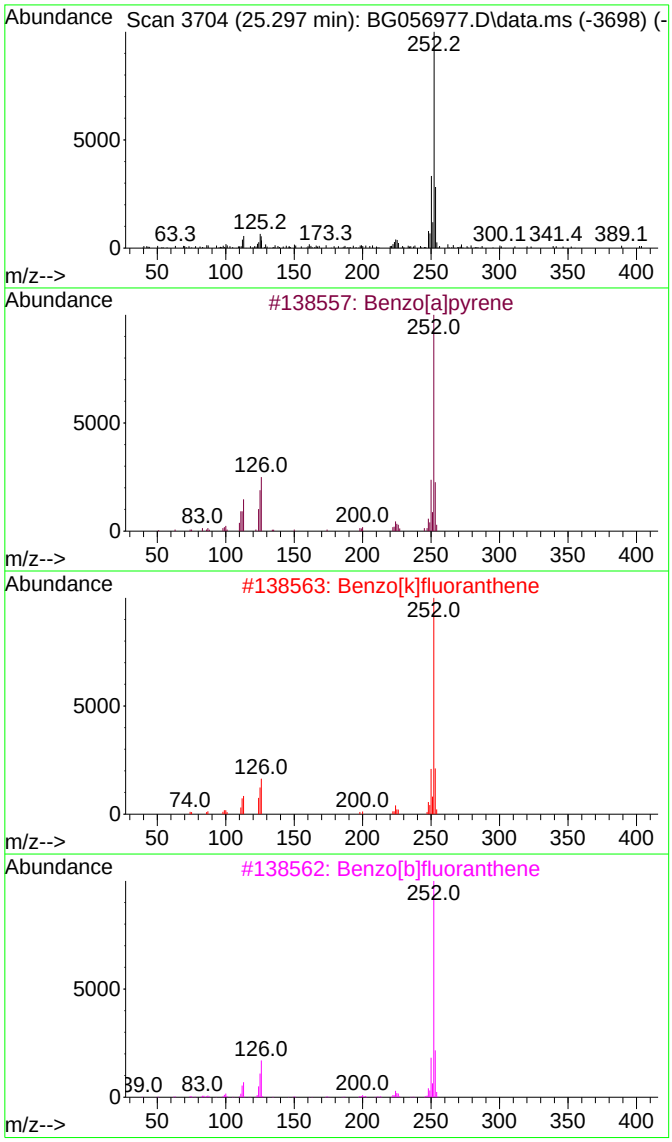
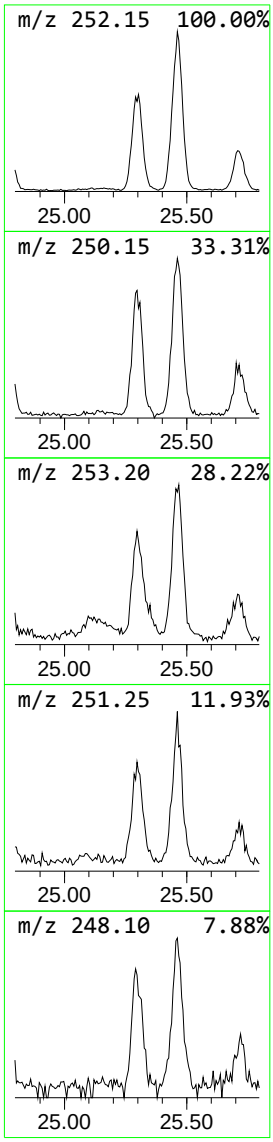
Peak Number 14 Benzo[j]fluoranthene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.297	22.37 ng	246172	Perylene-d12	25.631

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[a]pyrene	252	C20H12	000050-32-8	89
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	89
3	Benzo[b]fluoranthene	252	C20H12	000205-99-2	81
4	Benzo[j]fluoranthene	252	C20H12	000205-82-3	76
5	Perylene	252	C20H12	000198-55-0	76





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040423\
Data File : BG056977.D
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Operator : CG/JU
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Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2-Pentanone, 4-...	5.441	37.9	ng	208656	1	8.414	110135	20.0
Nonadecane	15.821	9.9	ng	101970	3	15.028	205447	20.0
Carbonic acid, ...	16.679	8.0	ng	99399	4	17.760	248495	20.0
Tridecanoic aci...	18.382	107.9	ng	1340140	4	17.760	248495	20.0
n-Hexadecanoic ...	18.611	23.4	ng	291230	4	17.760	248495	20.0
4H-Cyclopenta[d...	18.829	9.0	ng	111365	4	17.760	248495	20.0
9,12-Octadecadi...	19.522	715.4	ng	8888070	4	17.760	248495	20.0
9-Octadecenoic ...	19.546	328.0	ng	4075350	4	17.760	248495	20.0
Methyl stearate	19.663	53.9	ng	669726	4	17.760	248495	20.0
9,12-Octadecadi...	19.722	27.4	ng	340957	4	17.760	248495	20.0
Tetradecanoic acid	19.857	10.1	ng	124867	4	17.760	248495	20.0
Pyrene, 1-methyl-	20.638	5.5	ng	181936	5	22.083	655746	20.0
11H-Benzo[b]flu...	20.738	4.3	ng	139340	5	22.083	655746	20.0
Benzo[j]fluoran...	25.297	22.4	ng	246172	6	25.631	220070	20.0