

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041423\
 Data File : BG057109.D
 Acq On : 14 Apr 2023 21:20
 Operator : CG/JU
 Sample : 02348-01 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TR-05-041323

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: BG057109.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.912	398	404	413	rVB	44460	70187	1.90%	0.565%
2	7.533	673	680	690	rBB	46119	80017	2.16%	0.644%
3	8.402	821	828	835	rBB	73367	126091	3.41%	1.015%
4	9.577	1020	1028	1037	rBB2	27034	48537	1.31%	0.391%
5	11.240	1303	1311	1318	rBV	97867	176824	4.78%	1.423%
6	13.637	1712	1719	1725	rBV	78635	115383	3.12%	0.929%
7	15.017	1943	1954	1960	rBV	157824	253258	6.84%	2.038%
8	16.051	2125	2130	2138	rBV2	30071	53013	1.43%	0.427%
9	16.492	2200	2205	2213	rVB2	63265	98026	2.65%	0.789%
10	17.755	2414	2420	2423	rBV	203089	296724	8.02%	2.388%
11	17.796	2423	2427	2433	rVB	207286	312180	8.43%	2.512%
12	17.884	2437	2442	2447	rBV	49178	74821	2.02%	0.602%
13	18.371	2520	2525	2530	rVB	472260	603466	16.30%	4.857%
14	18.595	2558	2563	2565	rBV	46184	66866	1.81%	0.538%
15	18.618	2565	2567	2572	rVV	46781	67446	1.82%	0.543%
16	18.665	2572	2575	2582	rVB	47823	67492	1.82%	0.543%
17	18.818	2594	2601	2604	rBV3	58508	116891	3.16%	0.941%
18	19.100	2645	2649	2653	rBV	27570	37909	1.02%	0.305%
19	19.499	2710	2717	2719	rBV	2895026	3701228	100.00%	29.787%
20	19.529	2719	2722	2731	rVB2	1535640	2102686	56.81%	16.922%
21	19.646	2738	2742	2748	rBV	299284	357810	9.67%	2.880%
22	19.705	2748	2752	2755	rBV	120486	168454	4.55%	1.356%
23	19.787	2760	2766	2771	rVB	371297	540359	14.60%	4.349%
24	20.110	2816	2821	2823	rBV3	56731	91630	2.48%	0.737%
25	20.145	2823	2827	2833	rVV	331356	470172	12.70%	3.784%
26	20.322	2852	2857	2861	rBV2	141486	188894	5.10%	1.520%
27	20.463	2876	2881	2886	rVB5	23294	49065	1.33%	0.395%
28	20.627	2905	2909	2916	rVB2	78885	127973	3.46%	1.030%
29	20.727	2921	2926	2929	rBV	74617	102523	2.77%	0.825%
30	20.980	2966	2969	2976	rVB3	31064	55088	1.49%	0.443%
31	21.626	3074	3079	3082	rBV3	27106	43290	1.17%	0.348%
32	21.720	3091	3095	3100	rBV2	28530	44517	1.20%	0.358%
33	22.066	3145	3154	3160	rBV2	198477	582638	15.74%	4.689%
34	22.119	3160	3163	3170	rVB	121688	200159	5.41%	1.611%
35	22.290	3187	3192	3196	rBV	24732	43500	1.18%	0.350%
36	24.475	3557	3564	3585	rVB	79155	382681	10.34%	3.080%

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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

37	25.274	3694	3700	3710	rVB2	44172	127118	3.43%	1.023%
38	25.438	3720	3728	3738	rVB2	61710	187624	5.07%	1.510%
39	25.609	3752	3757	3765	rVB2	72361	193230	5.22%	1.555%

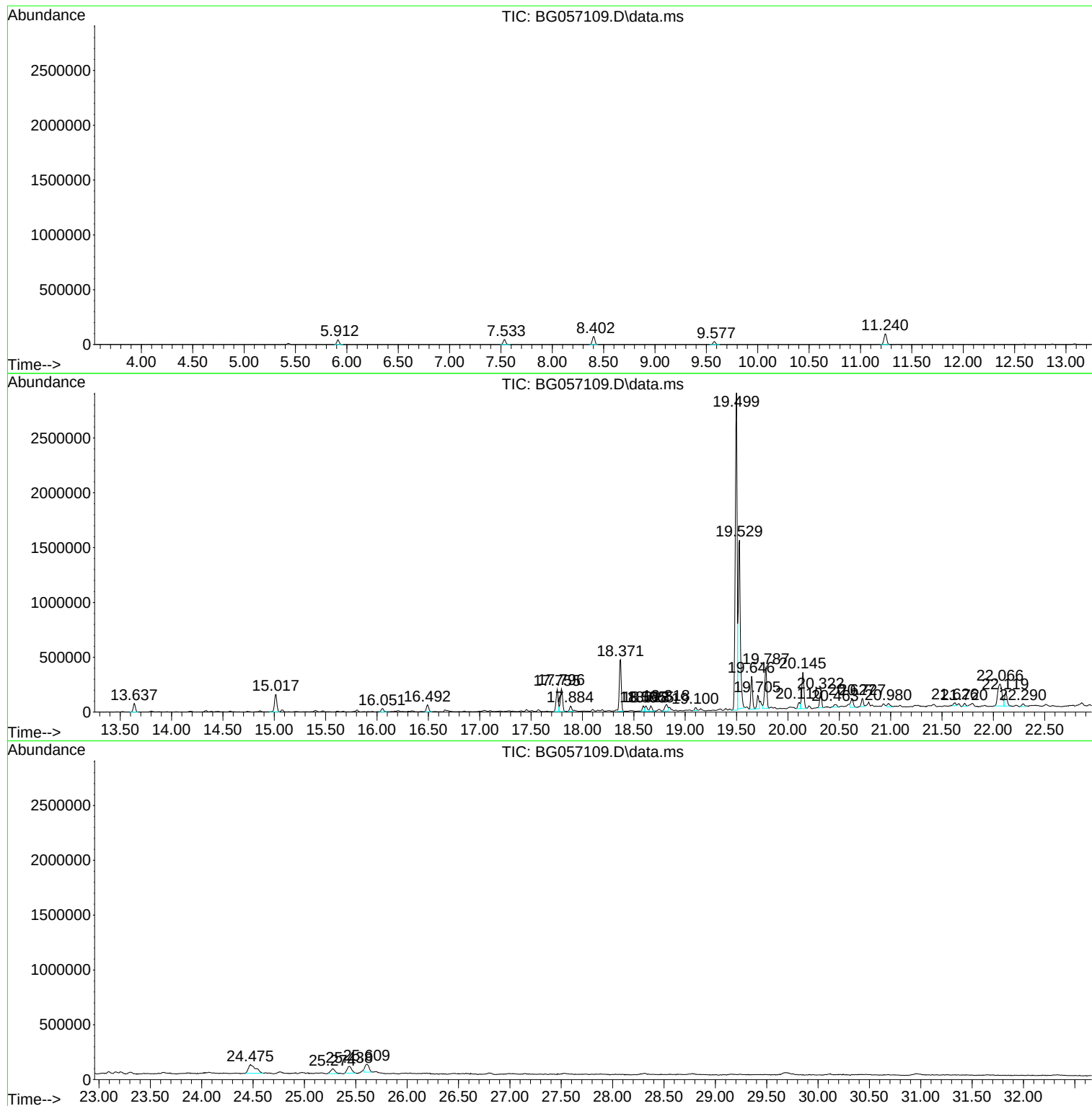
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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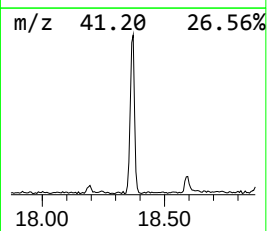
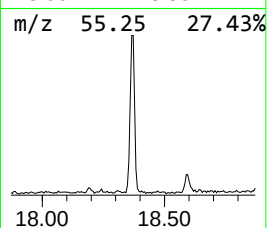
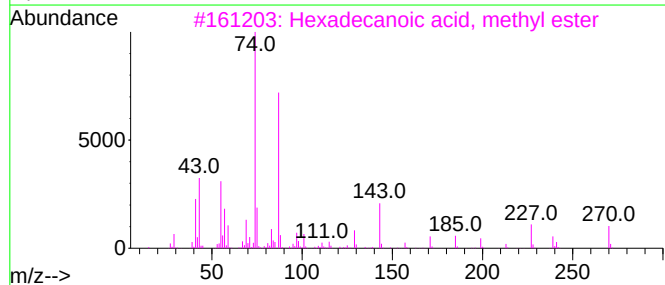
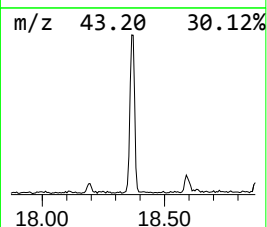
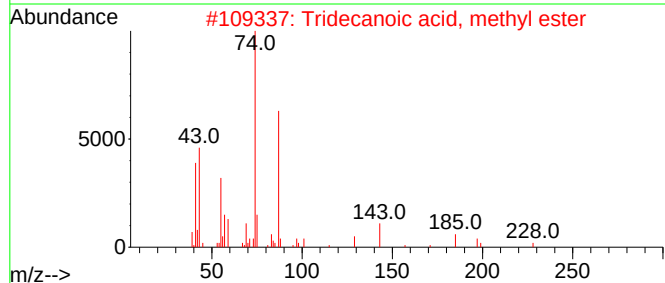
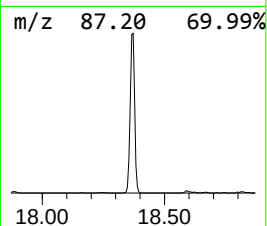
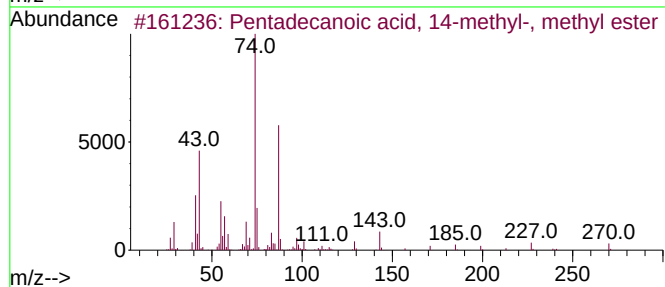
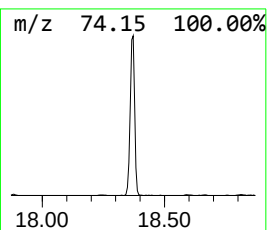
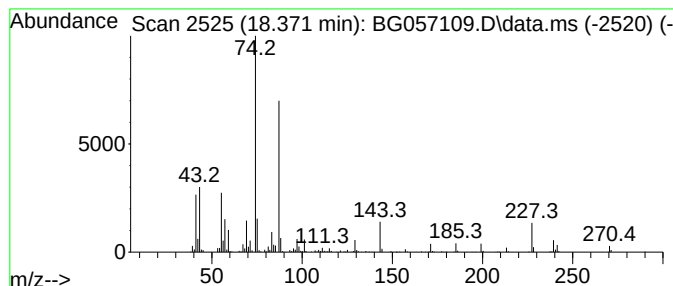
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Pentadecanoic acid, 14-meth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.371	40.68 ng	603466	Phenanthrene-d10	17.755

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	95
2			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	93
3			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	91
4			Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86
5			Methyl 11-methyl-dodecanoate	228	C14H28O2	1000336-45-1	80



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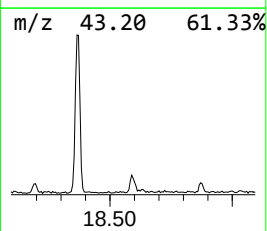
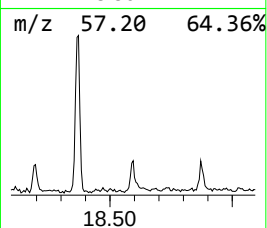
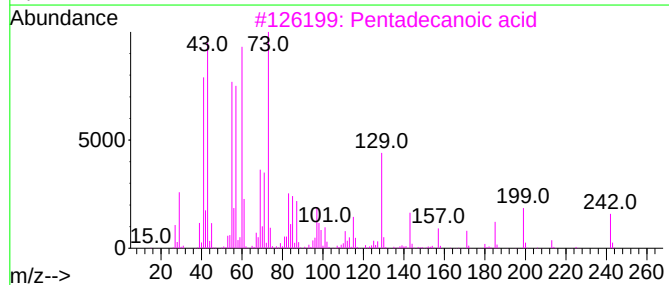
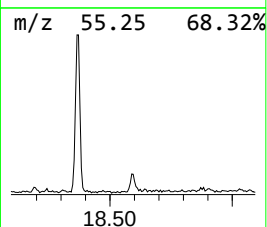
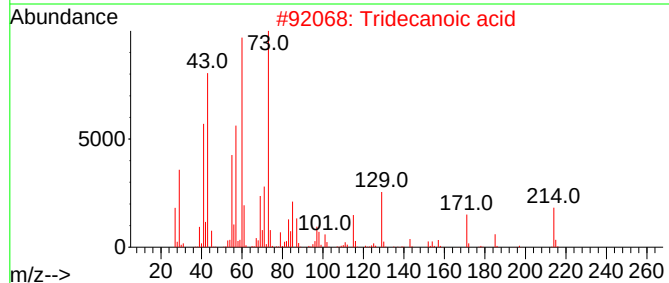
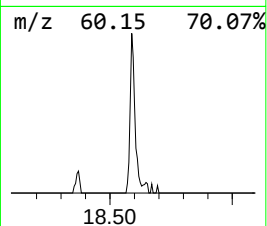
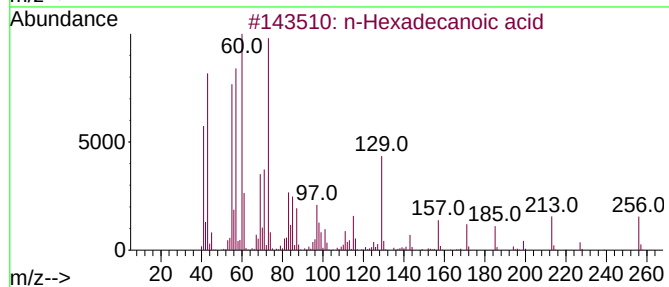
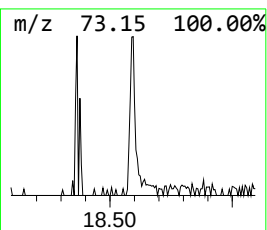
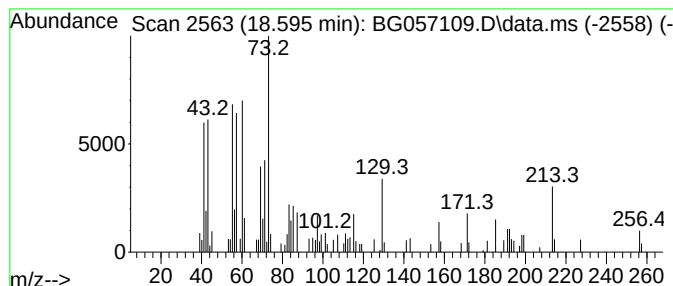
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.595	4.51 ng	66866	Phenanthrene-d10	17.755

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
2			Tridecanoic acid	214	C13H26O2	000638-53-9	83
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	47
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	43
5			n-Decanoic acid	172	C10H20O2	000334-48-5	41



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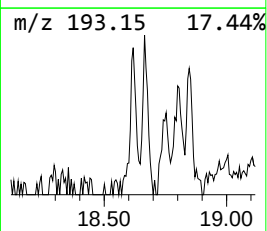
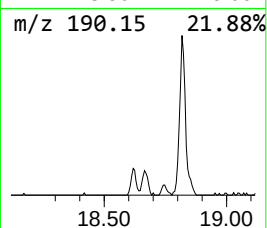
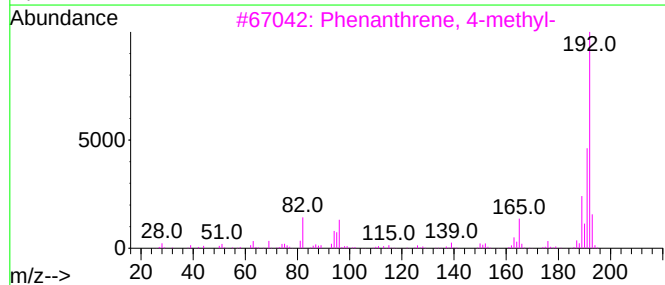
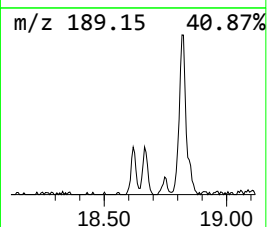
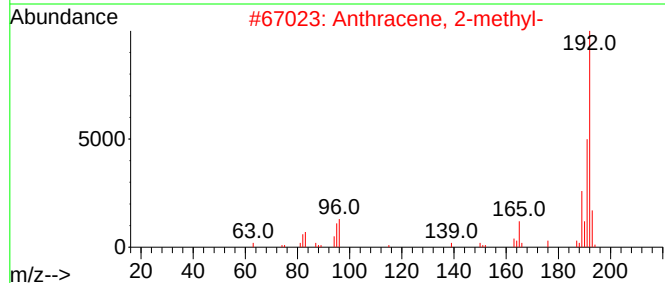
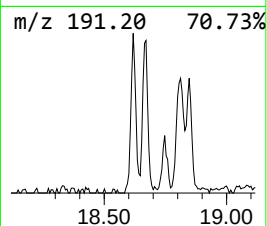
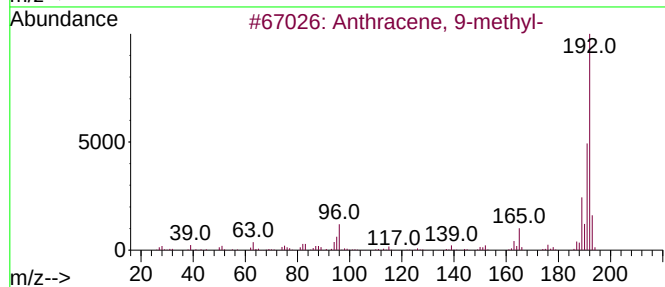
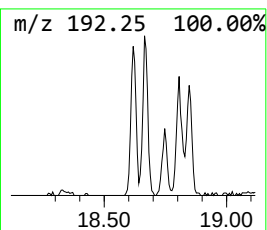
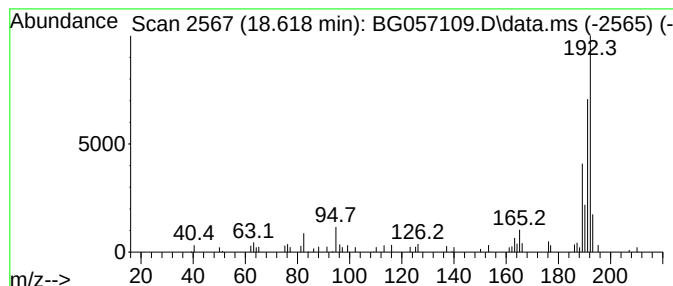
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Anthracene, 9-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.618	4.55 ng	67446	Phenanthrene-d10	17.755

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 9-methyl-	192	C15H12	000779-02-2	70
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	62
3			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	62
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	60
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	58



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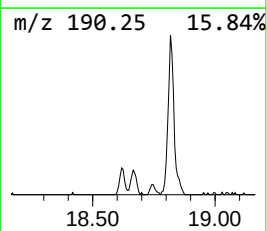
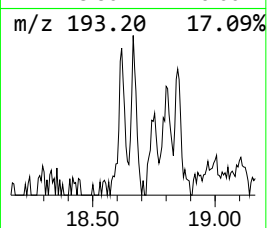
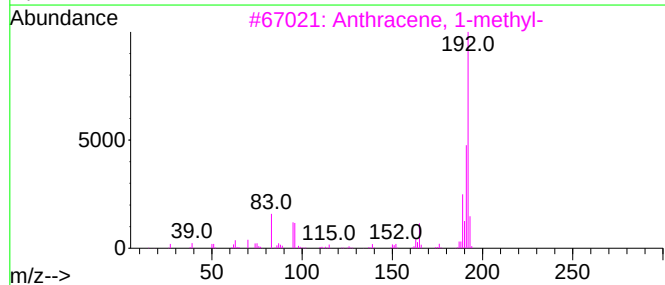
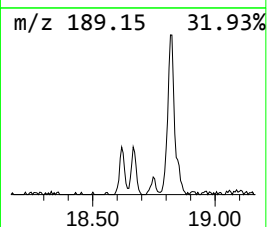
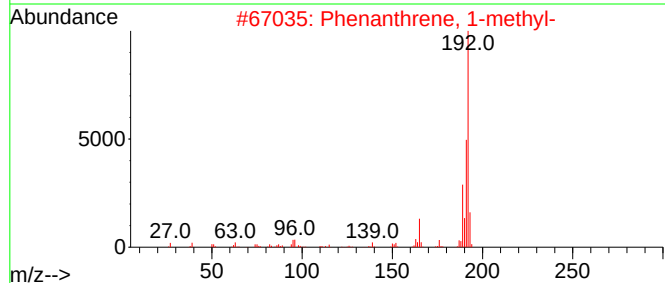
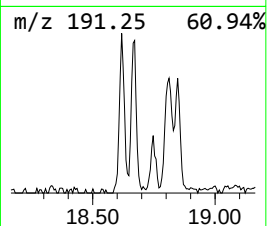
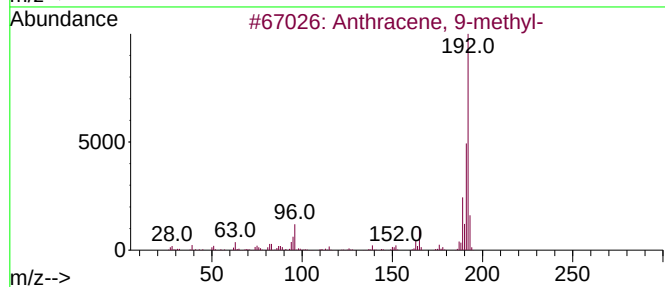
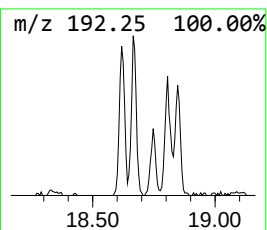
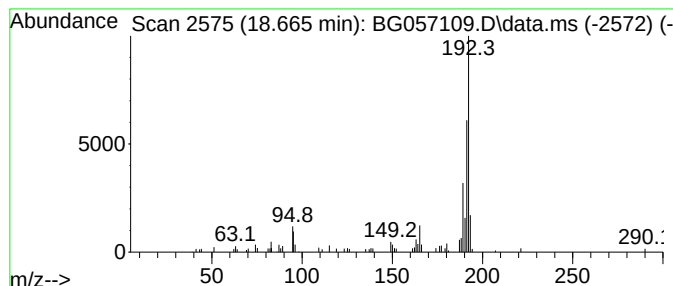
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.665	4.55 ng	67492	Phenanthrene-d10	17.755

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3			Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	89
5			Anthracene, 2-methyl-	192	C15H12	000613-12-7	81



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ALS Vial : 17 Sample Multiplier: 1

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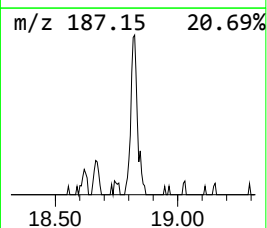
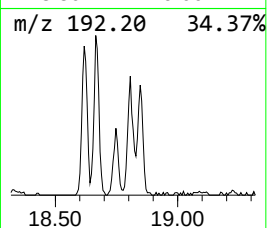
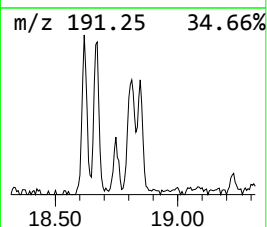
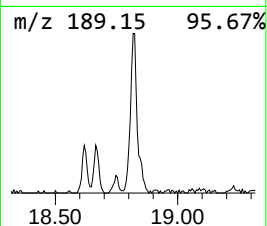
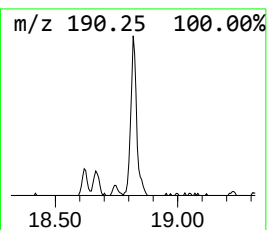
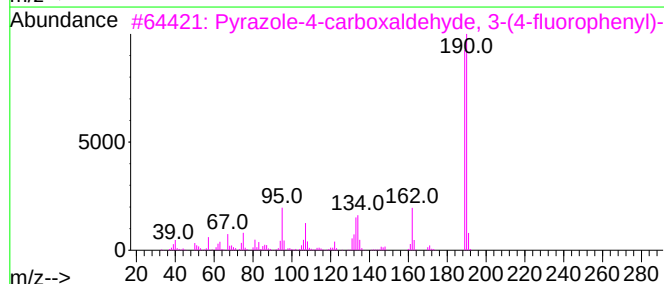
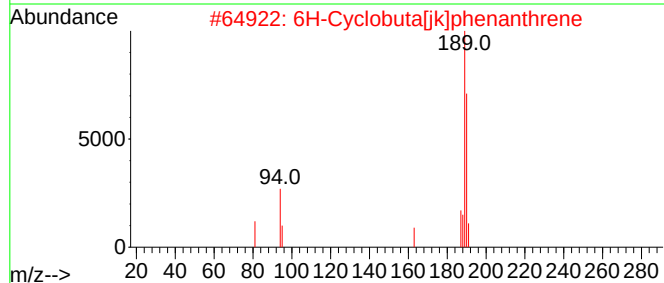
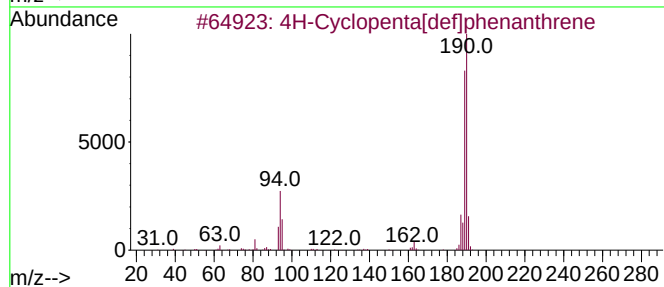
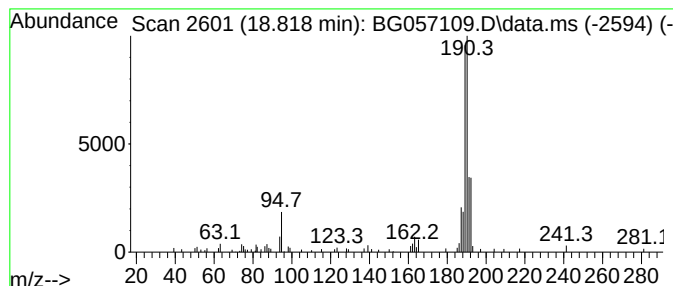
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.818	7.88 ng	116891	Phenanthrene-d10	17.755

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	83
2			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	64
3			Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	58
4			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	53
5			Cyclohexanone, 2-(2-furanylmethy...	190	C12H14O2	056053-07-7	27



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Data File : BG057109.D
Acq On : 14 Apr 2023 21:20
Operator : CG/JU
Sample : 02348-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TR-05-041323

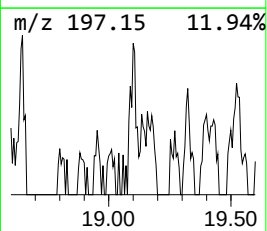
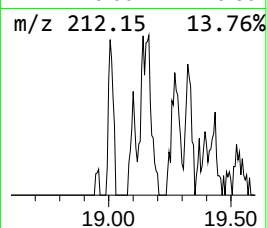
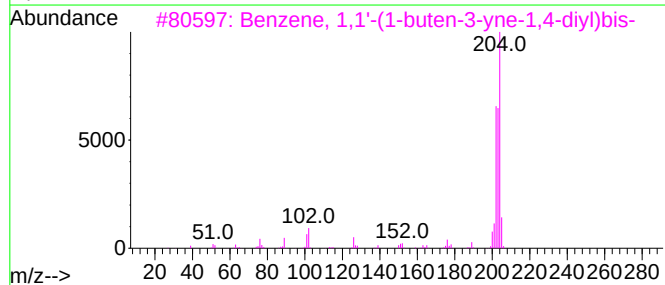
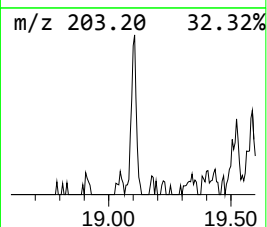
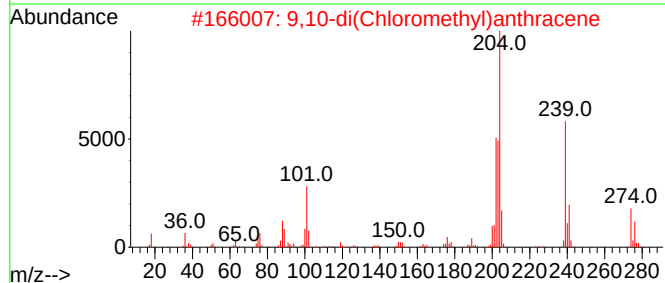
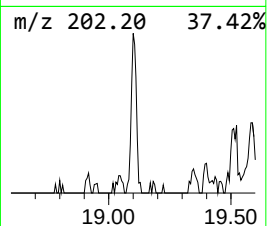
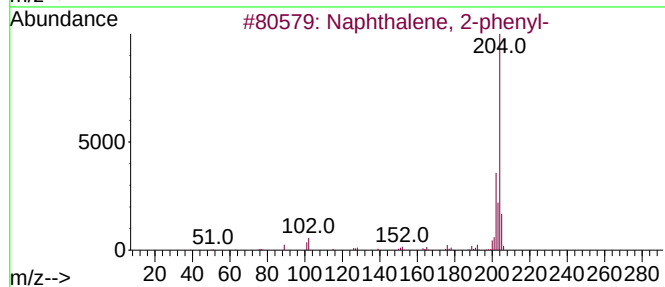
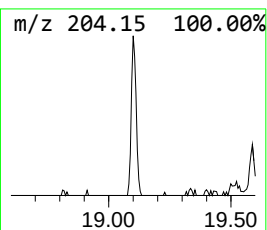
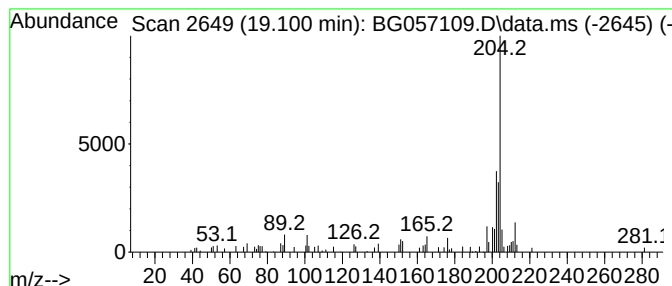
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Naphthalene, 2-phenyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.100	2.56 ng	37909	Phenanthrene-d10	17.755

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	58
2			9,10-di(Chloromethyl)anthracene	274	C16H12Cl2	010387-13-0	47
3			Benzene, 1,1'-(1-buten-3-yne-1,4...	204	C16H12	013141-45-2	43
4			Tetracyano-p-quinodimethane	204	C12H4N4	001518-16-7	43
5			1,4-Naphthalenedione, 2-hydroxy-...	204	C11H8O4	005257-83-0	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041423\
Data File : BG057109.D
Acq On : 14 Apr 2023 21:20
Operator : CG/JU
Sample : 02348-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

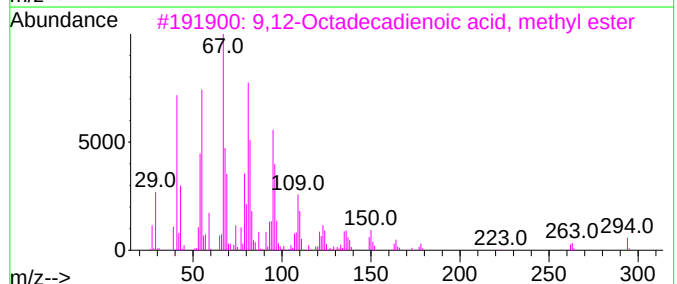
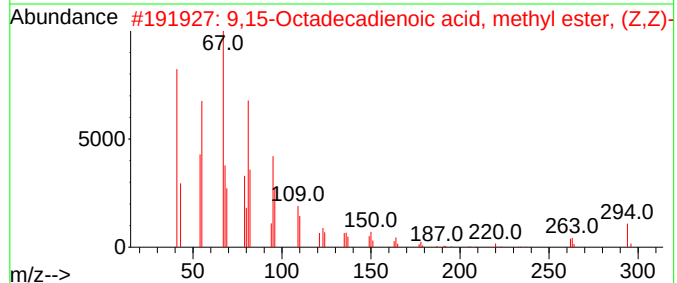
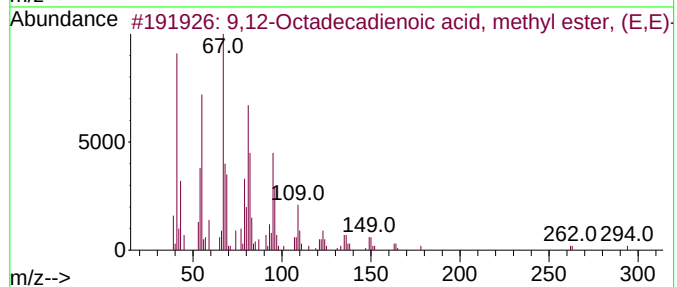
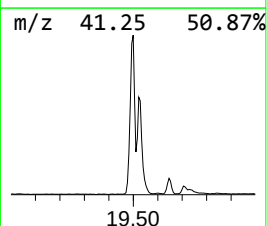
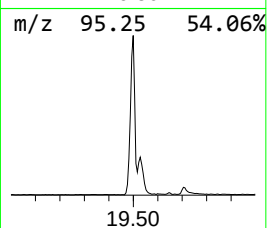
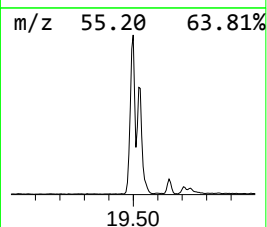
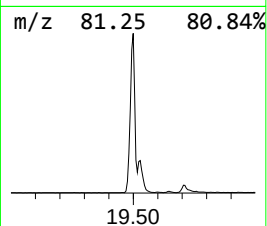
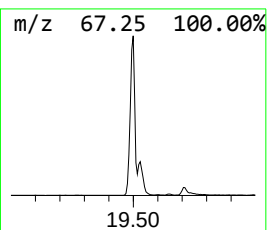
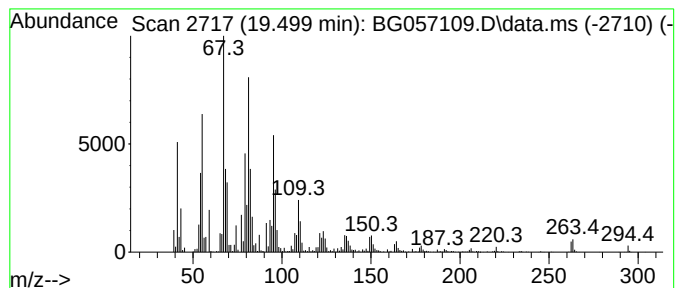
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.499	249.47 ng	3701230	Phenanthrene-d10	17.755
Hit#	of 5	Tentative ID	MW MolForm	CAS# Qual
1		9,12-Octadecadienoic acid, methy...	294 C19H34O2	002566-97-4 99
2		9,15-Octadecadienoic acid, methy...	294 C19H34O2	017309-05-6 98
3		9,12-Octadecadienoic acid, methy...	294 C19H34O2	002462-85-3 98
4		10,13-Octadecadienoic acid, meth...	294 C19H34O2	056554-62-2 97
5		9,12-Octadecadienoic acid (Z,Z)-...	294 C19H34O2	000112-63-0 96



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041423\
Data File : BG057109.D
Acq On : 14 Apr 2023 21:20
Operator : CG/JU
Sample : 02348-01 5X
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ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TR-05-041323

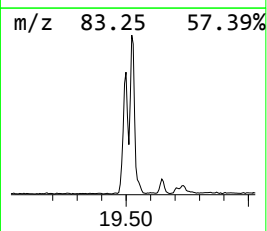
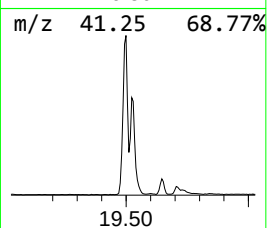
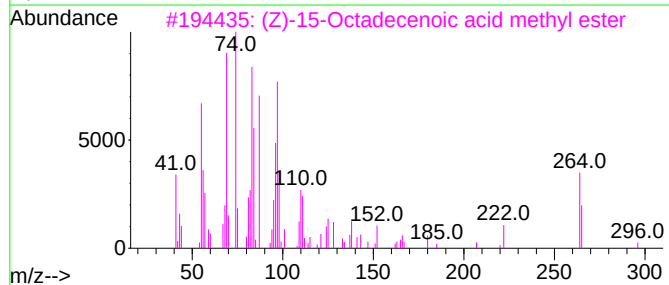
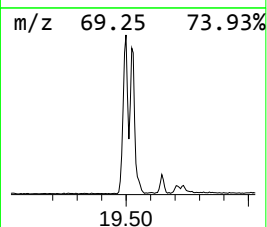
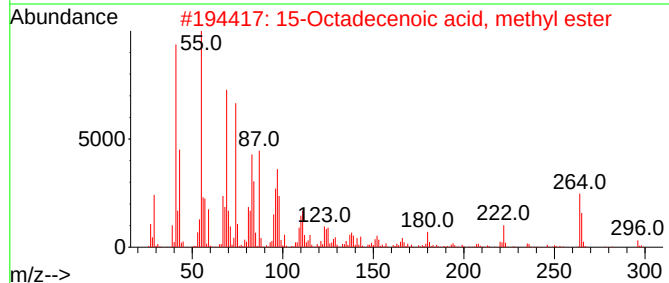
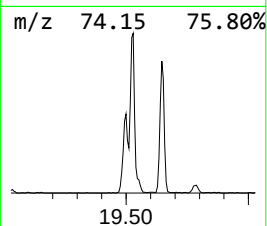
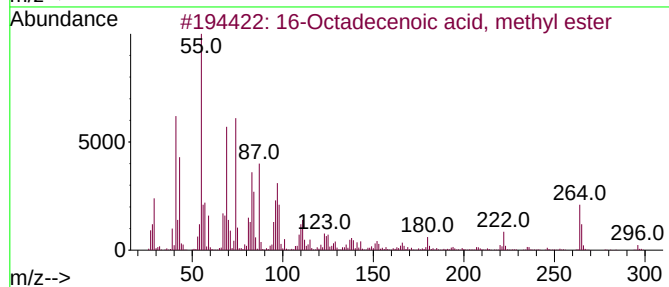
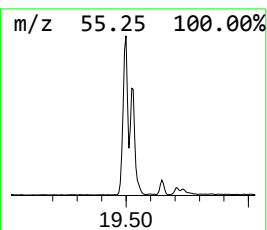
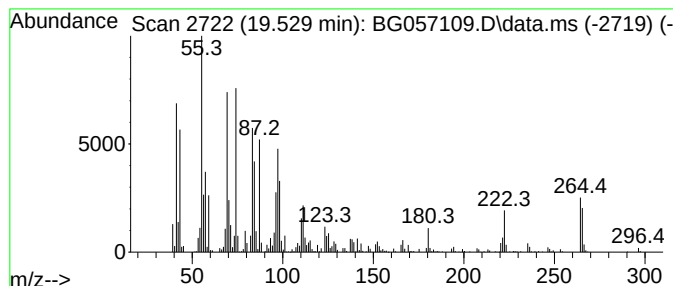
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 16-Octadecenoic acid, methy... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.529	141.73 ng	2102690	Phenanthrene-d10	17.755

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	93
2			15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	89
3			(Z)-15-Octadecenoic acid methyl ...	296	C19H36O2	1000427-69-3	87
4			9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	86
5			13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041423\
Data File : BG057109.D
Acq On : 14 Apr 2023 21:20
Operator : CG/JU
Sample : 02348-01 5X
Misc :
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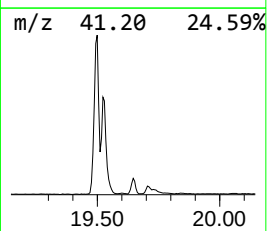
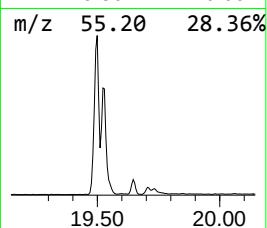
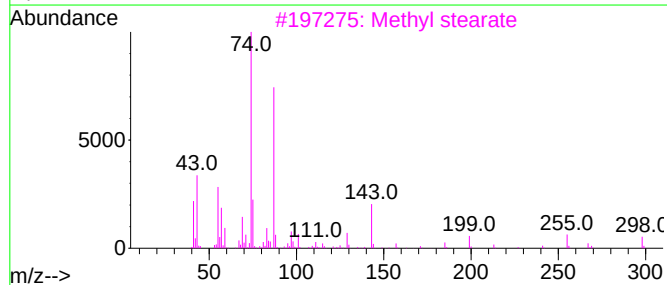
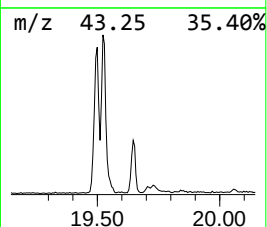
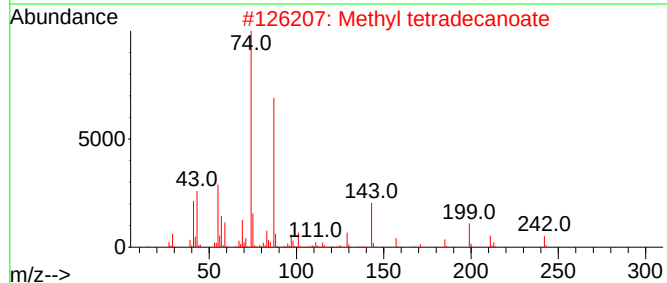
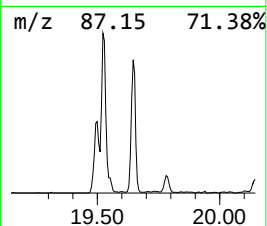
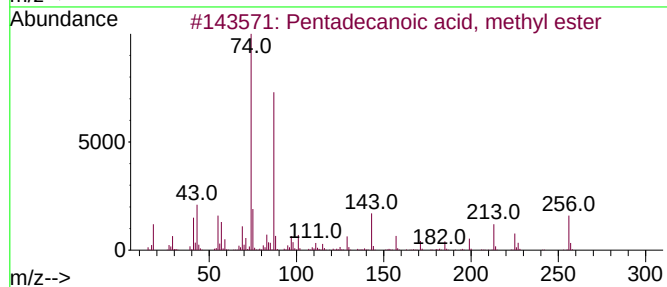
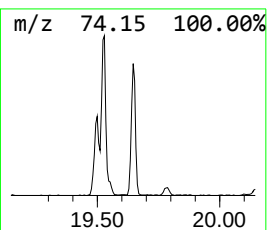
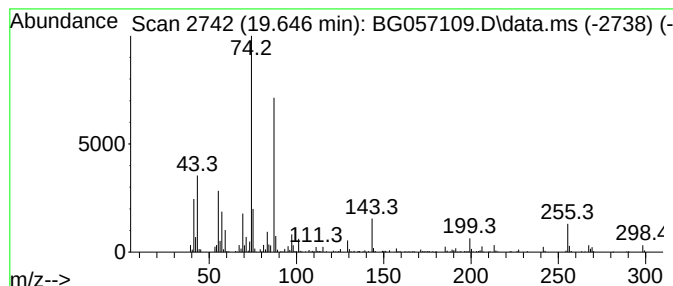
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Pentadecanoic acid, methyl ... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.646	24.12 ng	357810	Phenanthrene-d10		17.755	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pentadecanoic acid, methyl ester		256	C16H32O2	007132-64-1	93
2	Methyl tetradecanoate		242	C15H30O2	000124-10-7	93
3	Methyl stearate		298	C19H38O2	000112-61-8	92
4	Pentadecanoic acid, 14-methyl-, ...		270	C17H34O2	005129-60-2	87
5	Heptadecanoic acid, methyl ester		284	C18H36O2	001731-92-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041423\
Data File : BG057109.D
Acq On : 14 Apr 2023 21:20
Operator : CG/JU
Sample : 02348-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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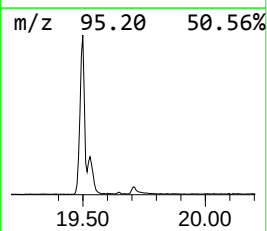
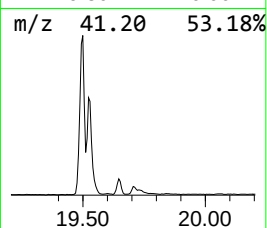
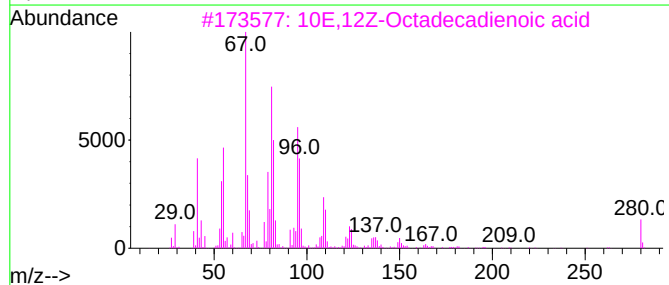
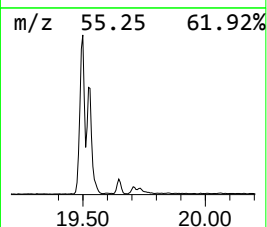
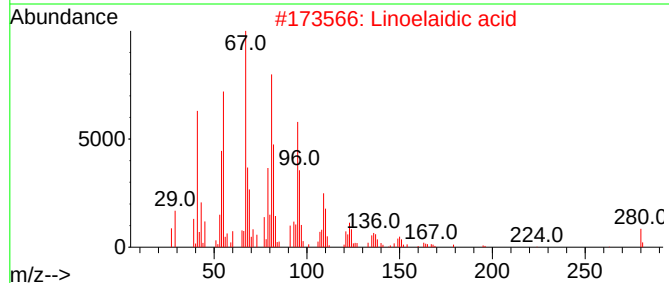
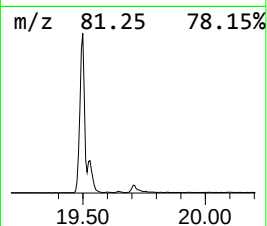
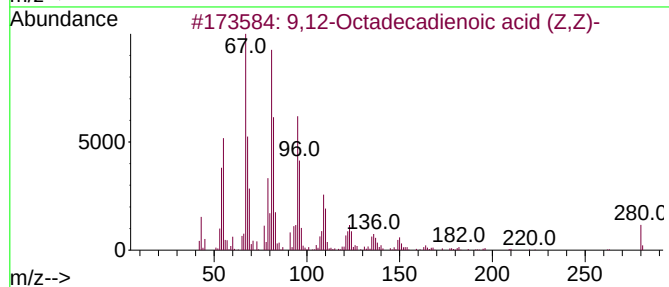
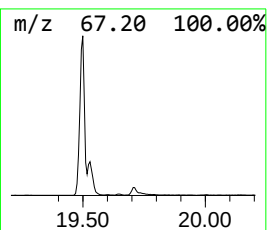
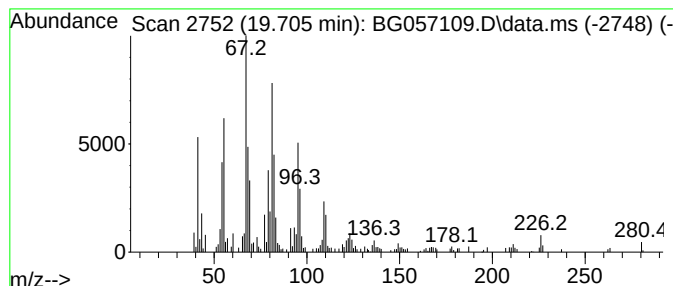
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
19.705	11.35 ng	168454	Phenanthrene-d10	17.755

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	97
2		Linoelaidic acid	280	C18H32O2	000506-21-8	94
3		10E,12Z-Octadecadienoic acid	280	C18H32O2	002420-56-6	93
4		Cyclodecene	138	C10H18	003618-12-0	91
5		(9E,11E)-Octadecadienoic acid	280	C18H32O2	000544-71-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041423\
Data File : BG057109.D
Acq On : 14 Apr 2023 21:20
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TR-05-041323

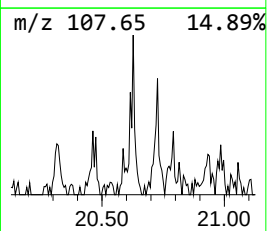
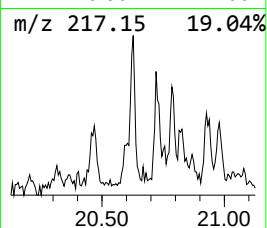
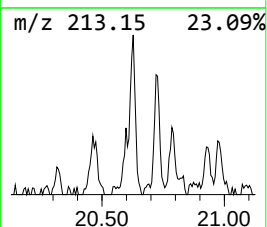
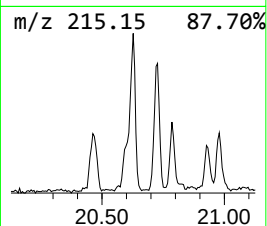
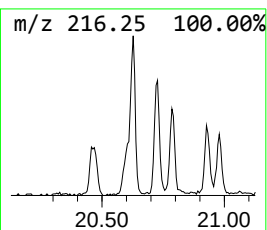
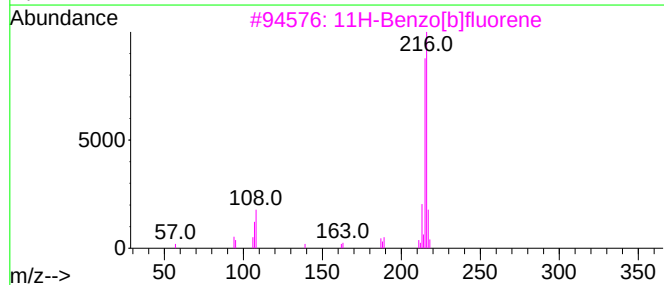
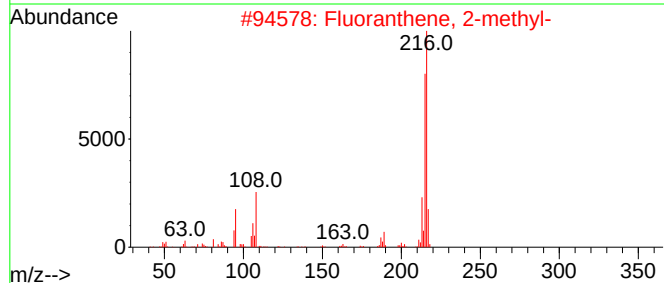
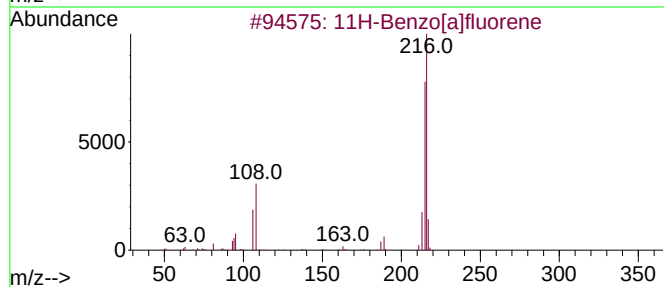
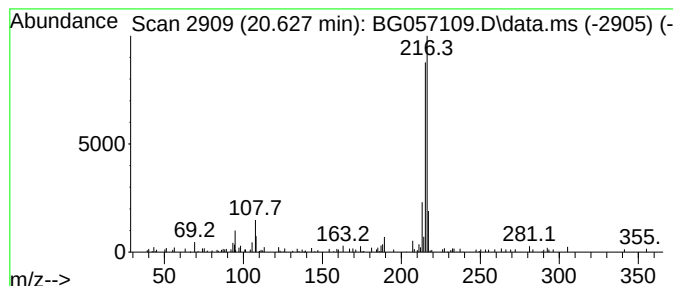
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 11H-Benzo[a]fluorene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.627	4.39 ng	127973	Chrysene-d12	22.072

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
2			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87
3			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
4			Pyrene, 1-methyl-	216	C17H12	002381-21-7	80
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041423\
Data File : BG057109.D
Acq On : 14 Apr 2023 21:20
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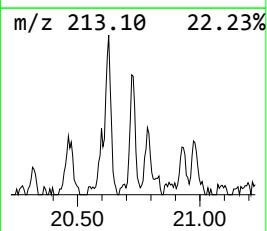
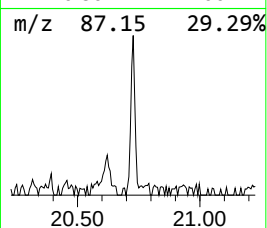
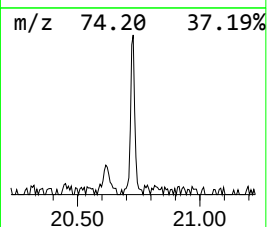
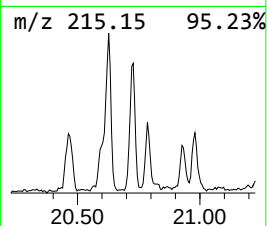
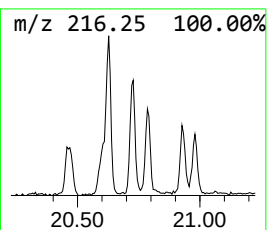
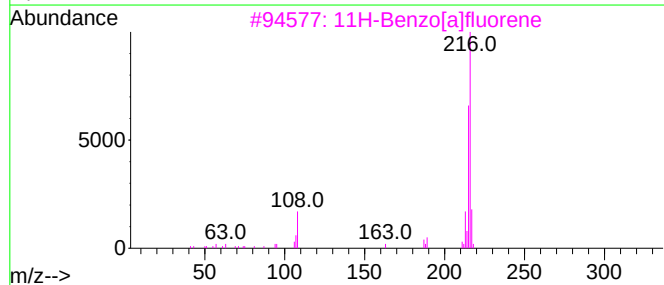
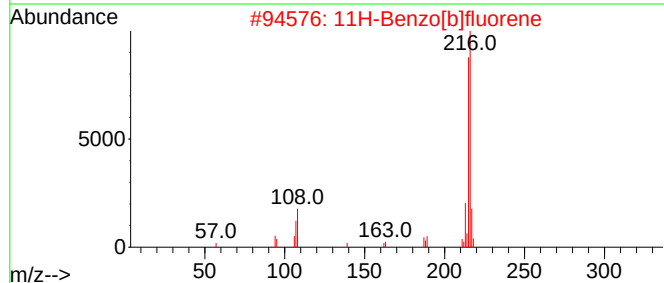
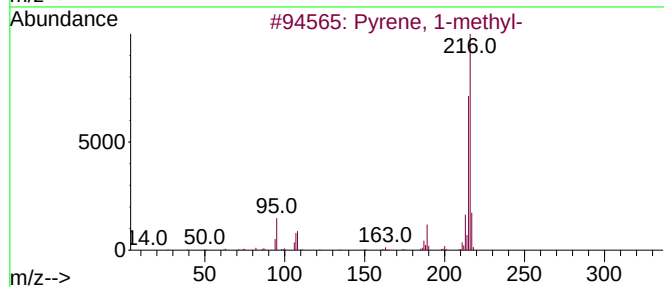
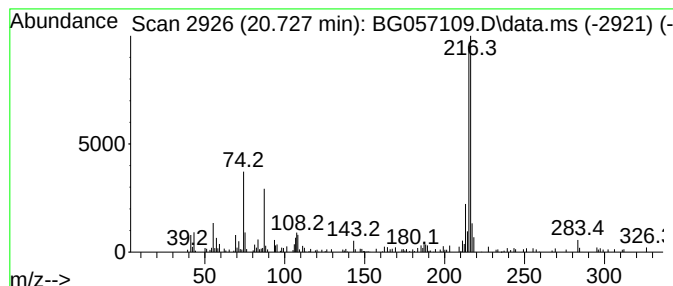
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.727	3.52 ng	102523	Chrysene-d12	22.072

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	62
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	58
5		7H-Benzo[c]fluorene	216	C17H12	000205-12-9	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041423\
Data File : BG057109.D
Acq On : 14 Apr 2023 21:20
Operator : CG/JU
Sample : 02348-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TR-05-041323

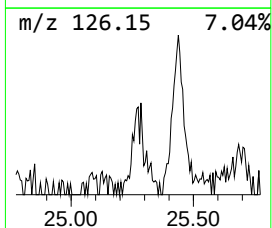
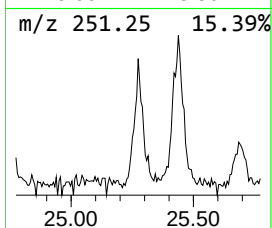
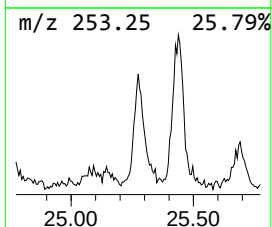
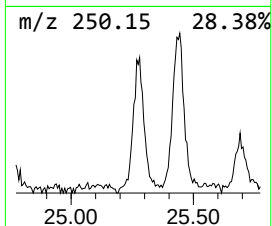
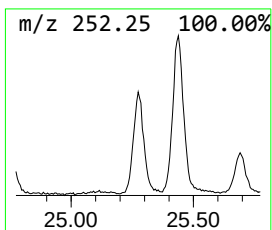
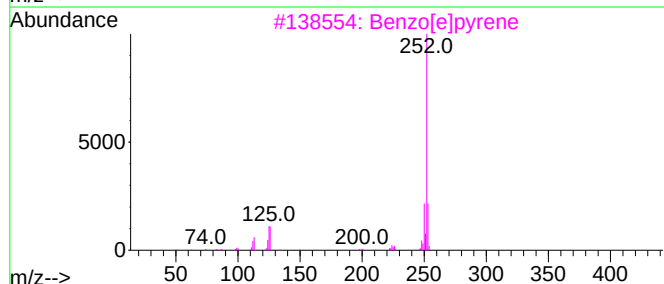
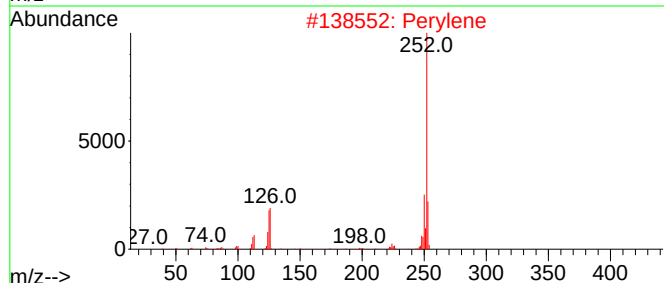
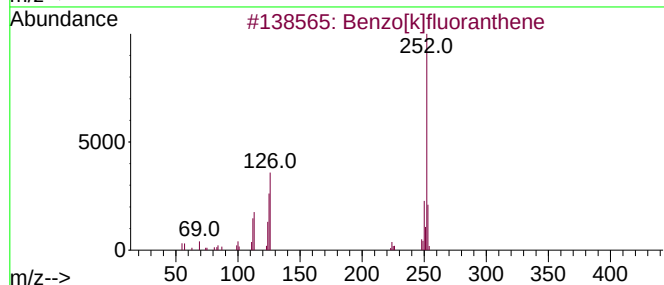
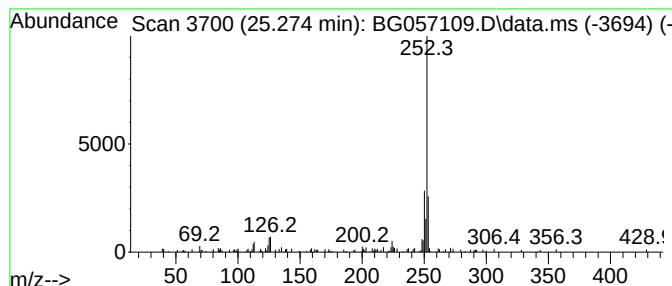
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Perylene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.274	13.16 ng	127118	Perylene-d12	25.609

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[k]fluoranthene	252	C20H12	000207-08-9	90
2		Perylene	252	C20H12	000198-55-0	76
3		Benzo[e]pyrene	252	C20H12	000192-97-2	76
4		Benzo[j]fluoranthene	252	C20H12	000205-82-3	76
5		Benzo[a]pyrene	252	C20H12	000050-32-8	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041423\
Data File : BG057109.D
Acq On : 14 Apr 2023 21:20
Operator : CG/JU
Sample : 02348-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TR-05-041323

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

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Pentadecanoic a...	18.371	40.7	ng	603466	4	17.755	296724	20.0
n-Hexadecanoic ...	18.595	4.5	ng	66866	4	17.755	296724	20.0
Anthracene, 9-m...	18.618	4.5	ng	67446	4	17.755	296724	20.0
Phenanthrene, 1...	18.665	4.5	ng	67492	4	17.755	296724	20.0
4H-Cyclopenta[d...	18.818	7.9	ng	116891	4	17.755	296724	20.0
Naphthalene, 2-...	19.100	2.6	ng	37909	4	17.755	296724	20.0
9,12-Octadecadi...	19.499	249.5	ng	3701230	4	17.755	296724	20.0
16-Octadecenoic...	19.529	141.7	ng	2102690	4	17.755	296724	20.0
Pentadecanoic a...	19.646	24.1	ng	357810	4	17.755	296724	20.0
9,12-Octadecadi...	19.705	11.4	ng	168454	4	17.755	296724	20.0
11H-Benzo[a]flu...	20.627	4.4	ng	127973	5	22.072	582638	20.0
Pyrene, 1-methyl-	20.727	3.5	ng	102523	5	22.072	582638	20.0
Perylene	25.274	13.2	ng	127118	6	25.609	193230	20.0