

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111222\
 Data File : BG055596.D
 Acq On : 12 Nov 2022 19:26
 Operator : CG/JU
 Sample : N5580-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SU-02-111022

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-BG110922.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG055596.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.323	3	6	11	rVB	113624	149139	1.87%	0.390%
2	5.321	340	346	361	rVB	212805	354332	4.44%	0.926%
3	5.797	419	427	441	rBV	752934	1276337	16.01%	3.334%
4	7.401	692	700	713	rBV	773158	1364750	17.11%	3.565%
5	8.264	840	847	860	rVB	205924	367447	4.61%	0.960%
6	9.440	1037	1047	1054	rBV	455339	828400	10.39%	2.164%
7	10.838	1279	1285	1294	rBV	47856	86361	1.08%	0.226%
8	11.096	1322	1329	1343	rVB	297418	543757	6.82%	1.421%
9	13.511	1733	1740	1748	rVB	1190658	1838353	23.05%	4.803%
10	14.886	1966	1974	1981	rBV	461209	718445	9.01%	1.877%
11	16.367	2217	2226	2234	rBV	913681	1396434	17.51%	3.648%
12	17.624	2434	2440	2445	rBV	513698	816375	10.24%	2.133%
13	17.665	2445	2447	2456	rVB	161647	233391	2.93%	0.610%
14	18.270	2545	2550	2557	rVB	1253531	1533484	19.23%	4.006%
15	18.494	2582	2588	2594	rBV2	466074	713952	8.95%	1.865%
16	18.547	2594	2597	2603	rVB	67144	96463	1.21%	0.252%
17	18.699	2616	2623	2626	rBV4	83469	176838	2.22%	0.462%
18	19.404	2736	2743	2745	rBV	5984732	7974431	100.00%	20.834%
19	19.434	2745	2748	2757	rVB2	4491162	5995636	75.19%	15.664%
20	19.557	2764	2769	2774	rBV	982310	1122668	14.08%	2.933%
21	19.616	2774	2779	2781	rVV	741624	1022583	12.82%	2.672%
22	19.639	2781	2783	2785	rVV	620080	768496	9.64%	2.008%
23	19.663	2785	2787	2792	rVV	611613	803388	10.07%	2.099%
24	19.710	2793	2795	2799	rVV2	89206	126188	1.58%	0.330%
25	19.751	2799	2802	2809	rVB2	195215	254240	3.19%	0.664%
26	19.904	2824	2828	2837	rVB5	70492	136899	1.72%	0.358%
27	19.986	2838	2842	2844	rVV2	70261	119406	1.50%	0.312%
28	20.027	2844	2849	2856	rVB	485691	719192	9.02%	1.879%
29	20.215	2876	2881	2886	rVB	1683981	2198620	27.57%	5.744%
30	20.509	2927	2931	2940	rVB3	126923	277224	3.48%	0.724%
31	20.609	2945	2948	2951	rBV	73401	92234	1.16%	0.241%
32	20.644	2951	2954	2956	rVV2	92303	106651	1.34%	0.279%
33	20.679	2957	2960	2967	rVB4	80847	150601	1.89%	0.393%
34	21.531	3102	3105	3110	rVB2	99674	136489	1.71%	0.357%
35	21.696	3131	3133	3138	rVB	84396	100860	1.26%	0.264%
36	21.925	3164	3172	3177	rBV2	559345	1384372	17.36%	3.617%

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ClientSampleId :
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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-BG110922.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	21.972	3177	3180	3189	rVB	241134	421251	5.28%	1.101%
38	24.258	3563	3569	3577	rBV	208963	631867	7.92%	1.651%
39	25.186	3721	3727	3741	rVB	180469	536408	6.73%	1.401%
40	25.356	3749	3756	3767	rBV2	232883	702801	8.81%	1.836%

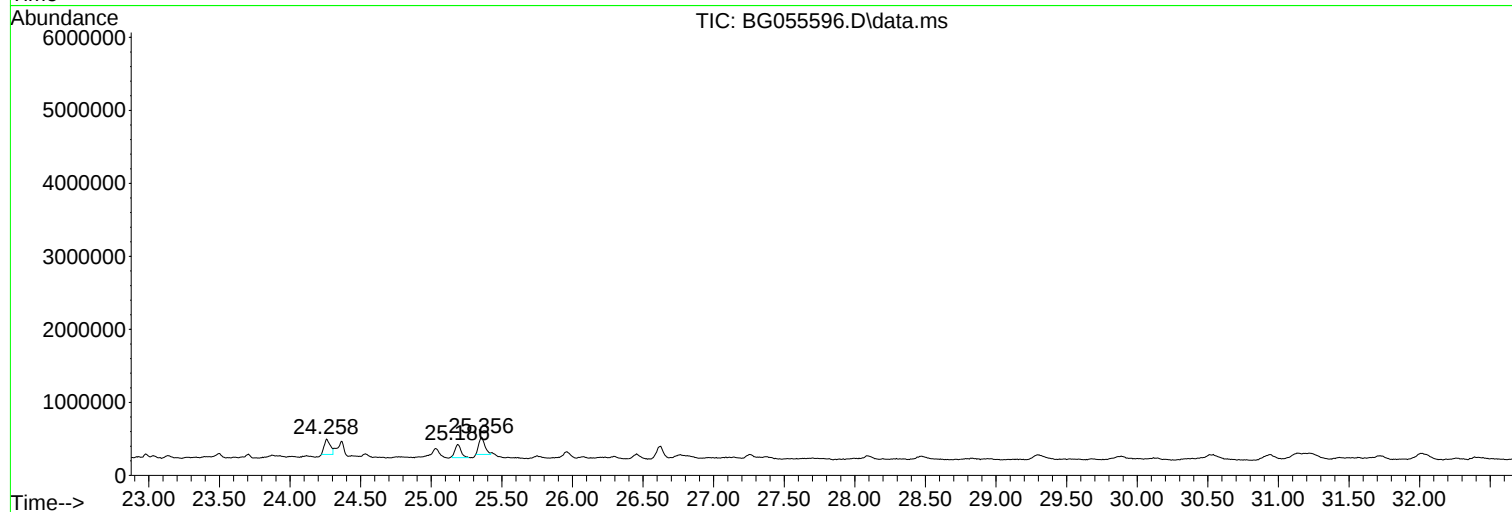
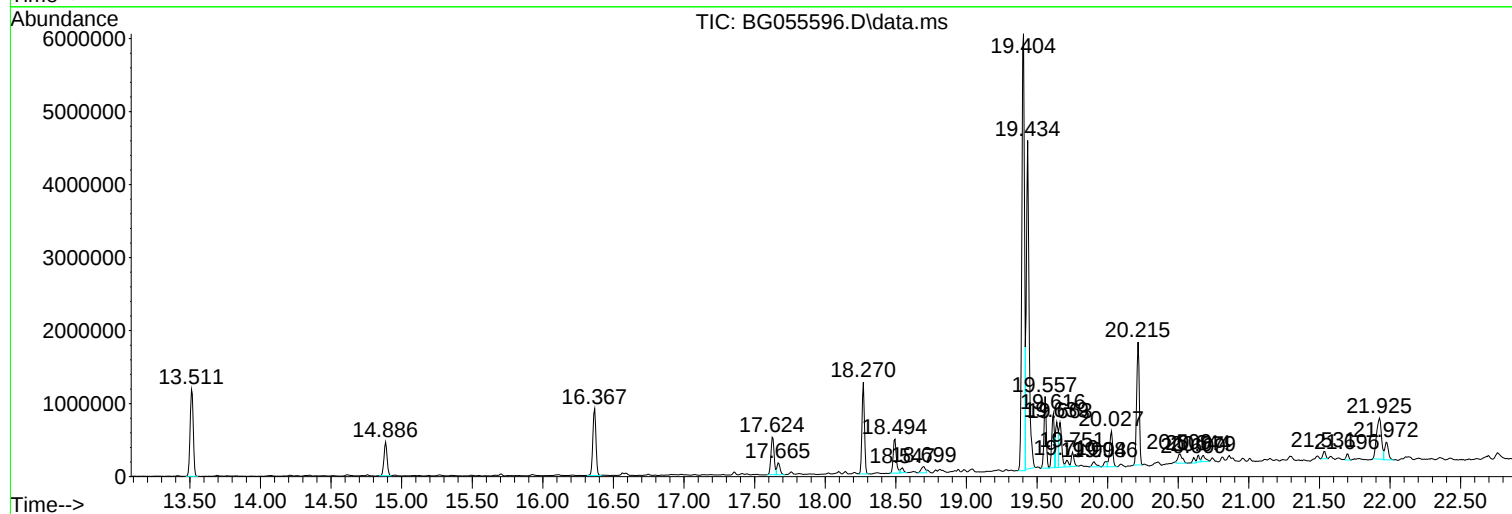
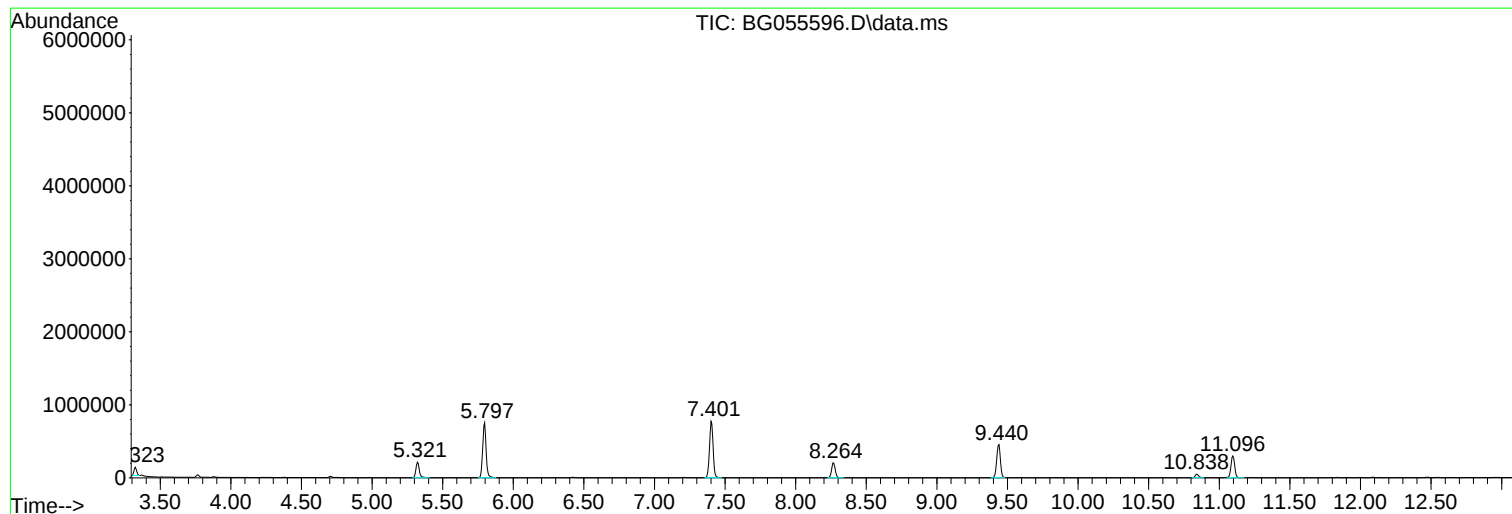
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :
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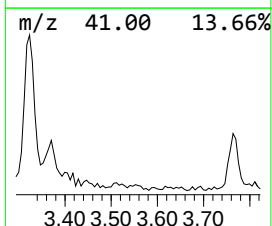
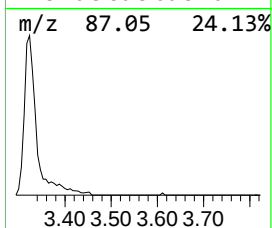
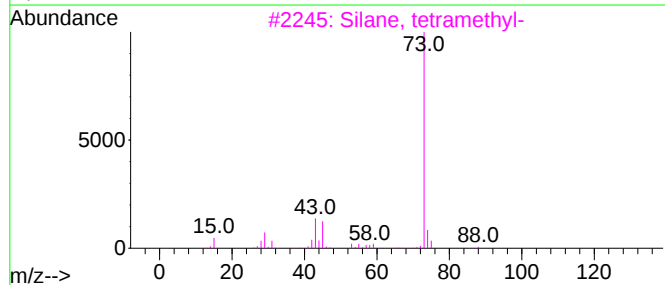
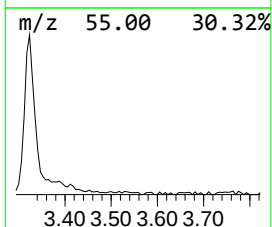
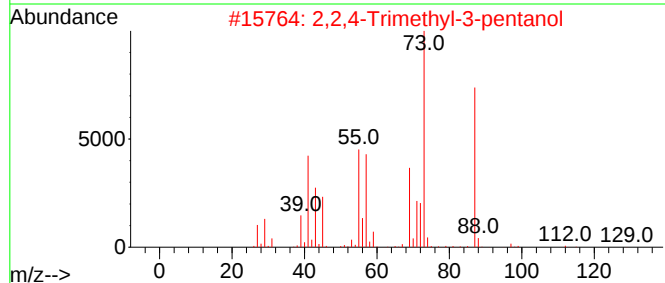
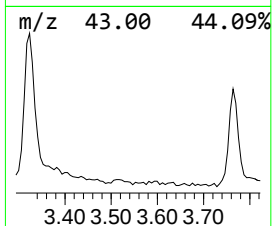
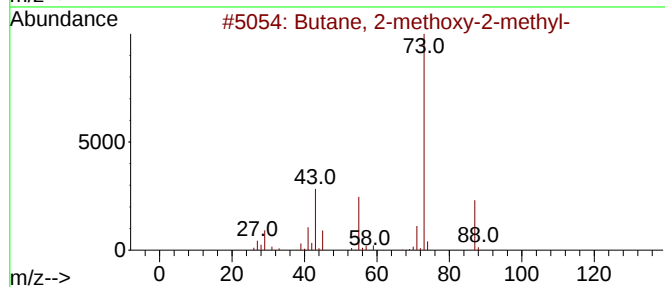
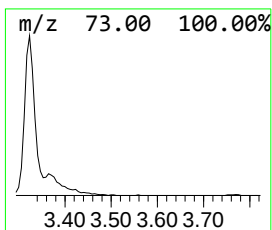
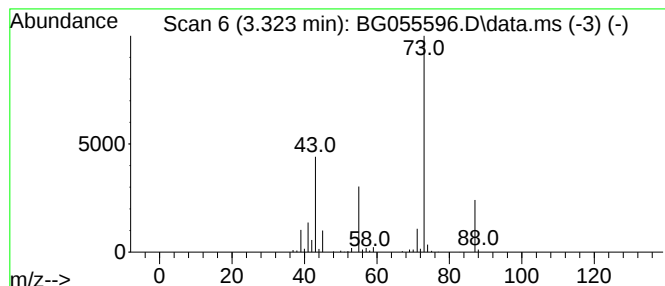
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.323	8.12 ng	149139	1,4-Dichlorobenzene-d4	8.270

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	40
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
4			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	37
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	32



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Misc :
ALS Vial : 12 Sample Multiplier: 1

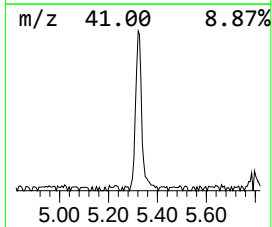
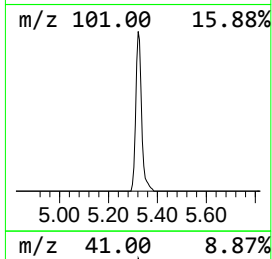
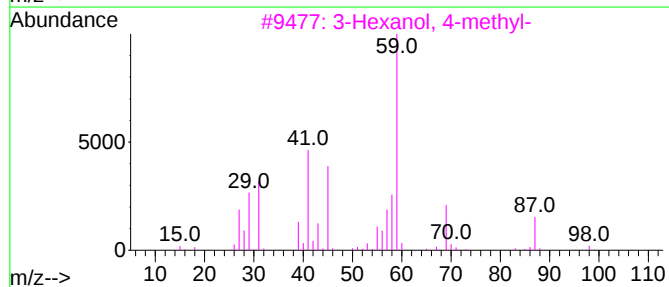
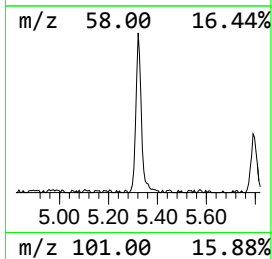
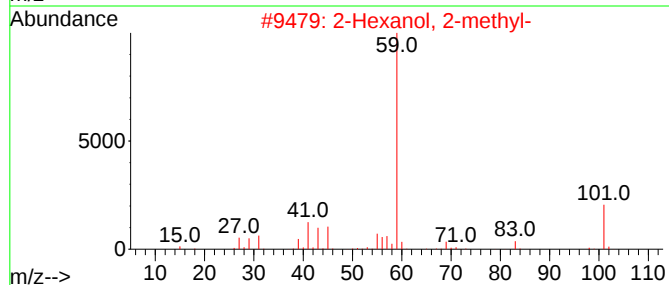
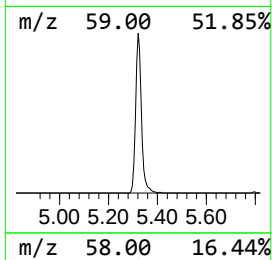
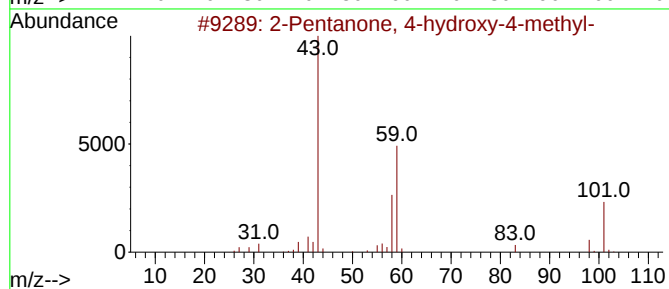
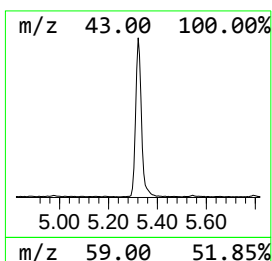
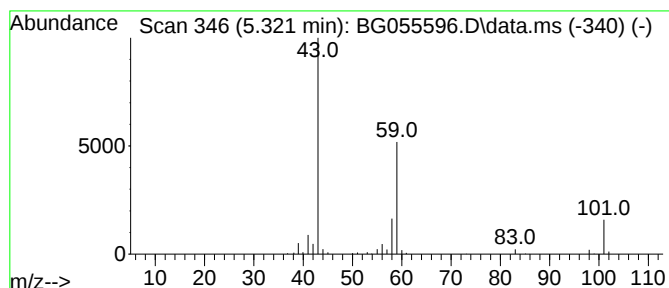
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.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.
5.321	19.29 ng	354332	1,4-Dichlorobenzene-d4			8.270
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	72
2	2-Hexanol, 2-methyl-		116	C7H16O	000625-23-0	33
3	3-Hexanol, 4-methyl-		116	C7H16O	000615-29-2	33
4	Acetic acid, cyano-, 1,1-dimethy...		141	C7H11NO2	001116-98-9	25
5	1-Propen-2-ol, acetate		100	C5H8O2	000108-22-5	10



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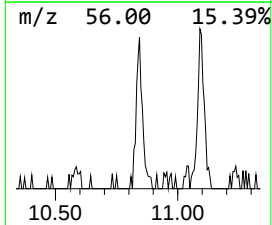
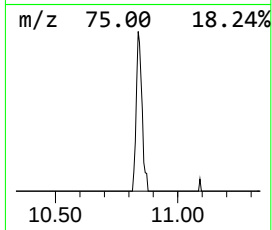
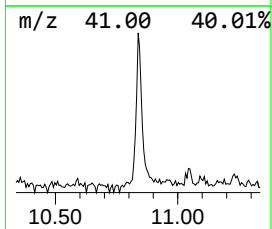
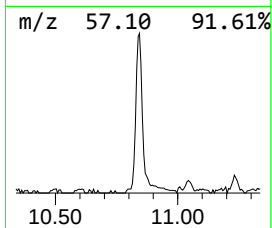
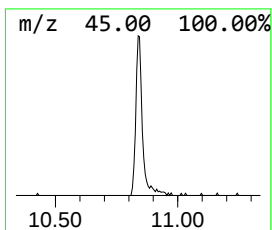
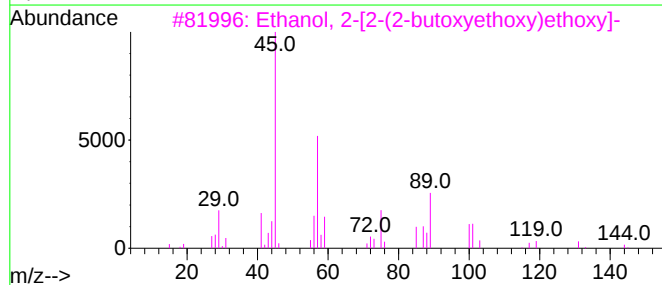
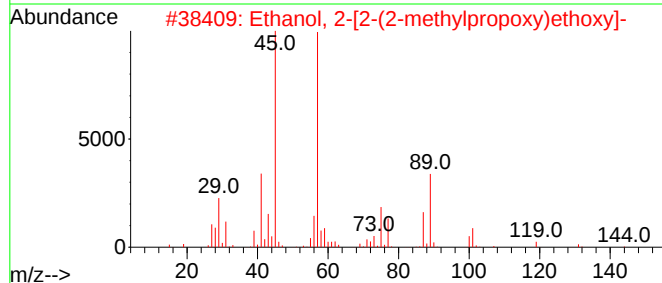
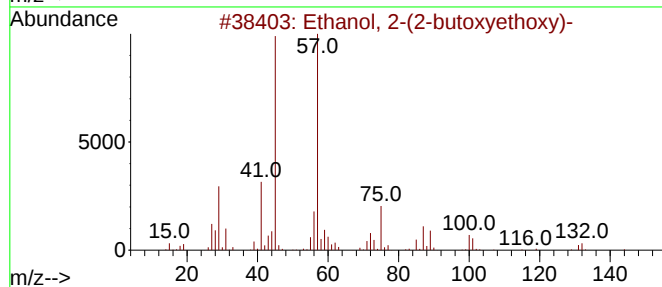
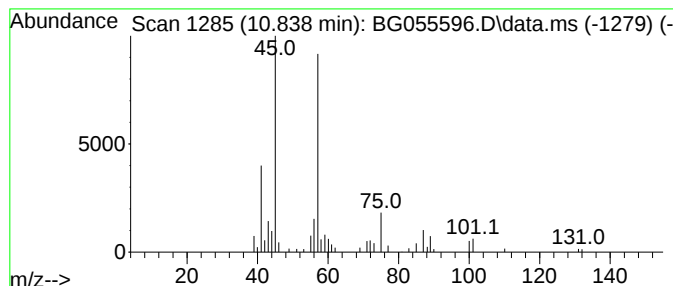
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.838	3.18 ng	86361	Naphthalene-d8	11.097	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
2	Ethanol, 2-[2-(2-methylpropoxy)et...	162	C8H18O3	018912-80-6	56
3	Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	50
4	Propanoic acid, 2-(methoxymethoxy)-	134	C5H10O4	081327-29-9	47
5	Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	45



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ClientSampleId :
SU-02-111022

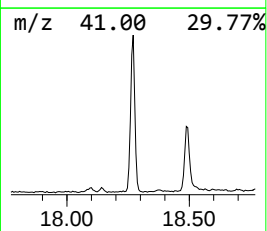
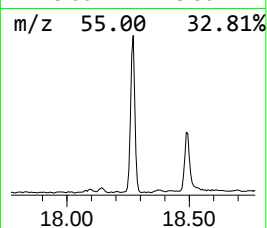
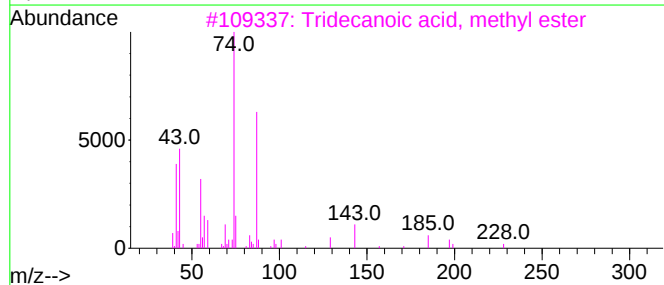
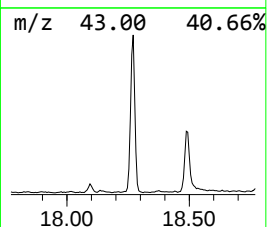
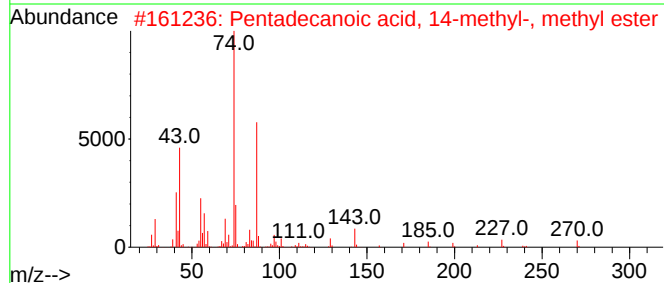
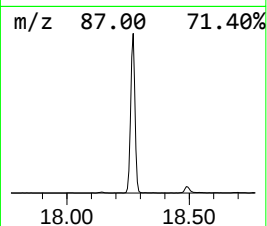
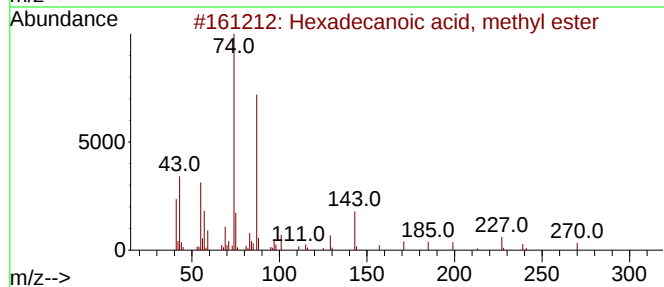
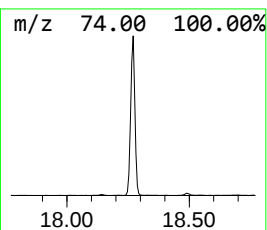
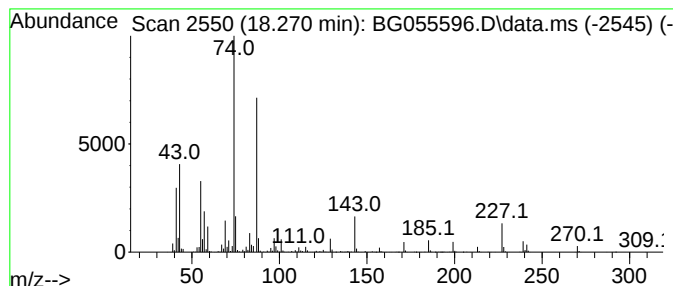
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.270	37.57 ng	1533480	Phenanthrene-d10	17.624

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	95
3			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	93
4			Methyl tetradecanoate	242	C15H30O2	000124-10-7	87
5			Dodecanoic acid, 10-methyl-, met...	228	C14H28O2	005129-65-7	83



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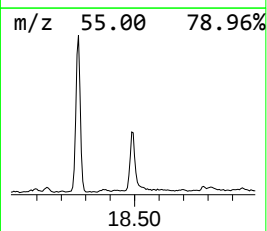
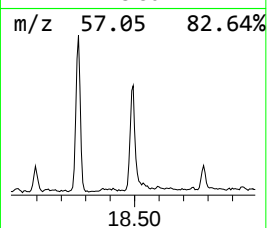
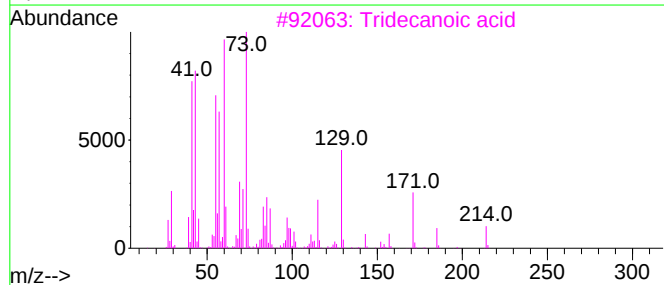
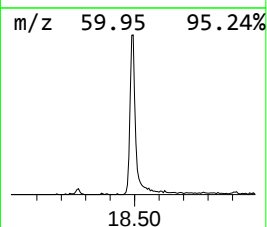
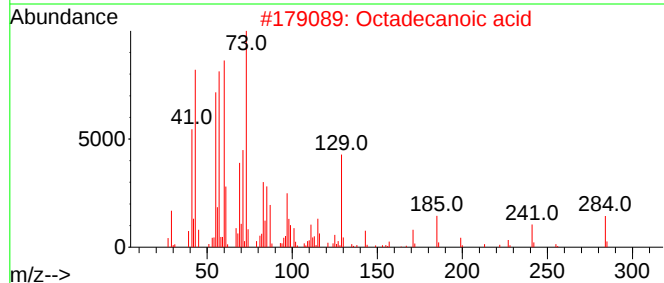
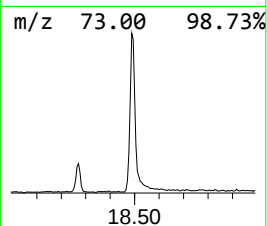
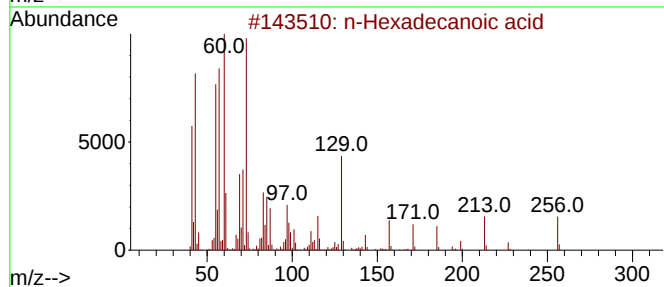
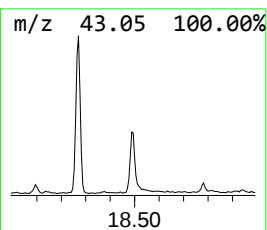
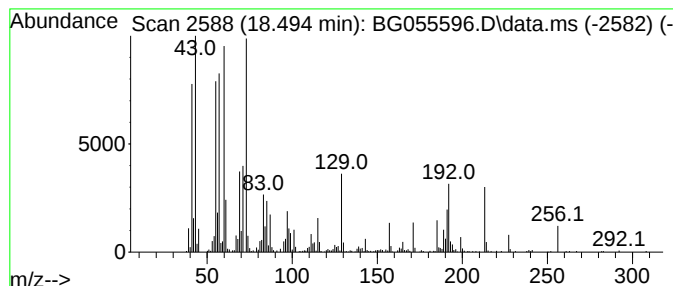
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.494	17.49 ng	713952	Phenanthrene-d10	17.624

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	93
3			Tridecanoic acid	214	C13H26O2	000638-53-9	91
4			Dodecanoic acid	200	C12H24O2	000143-07-7	78
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111222\
Data File : BG055596.D
Acq On : 12 Nov 2022 19:26
Operator : CG/JU
Sample : N5580-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

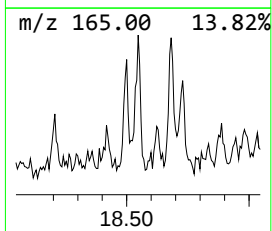
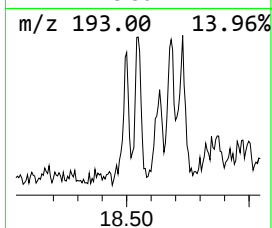
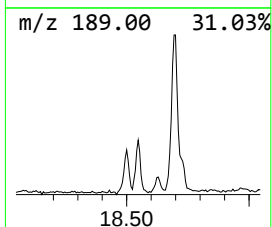
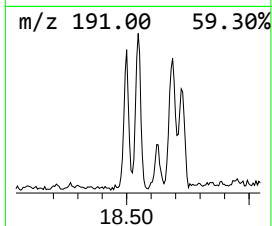
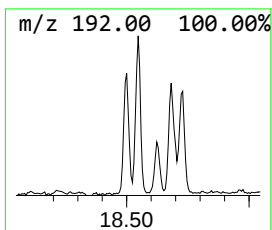
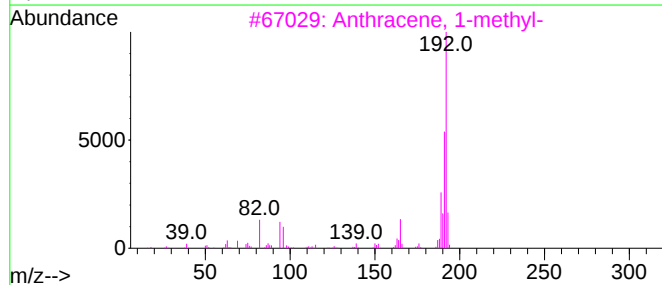
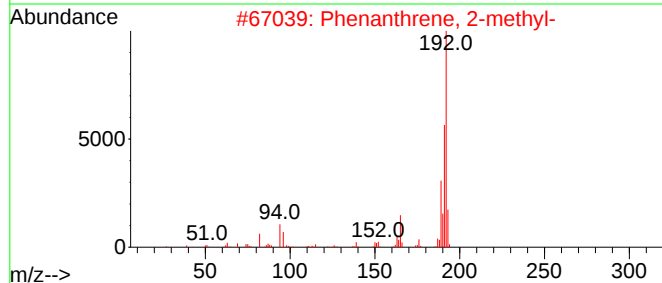
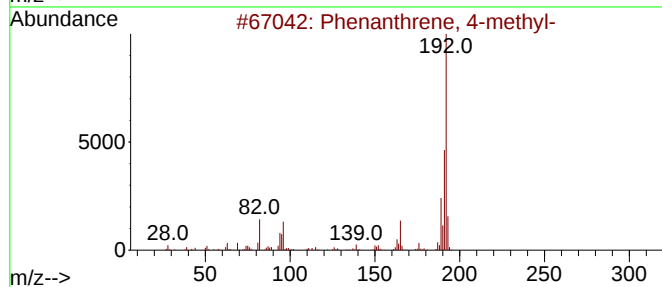
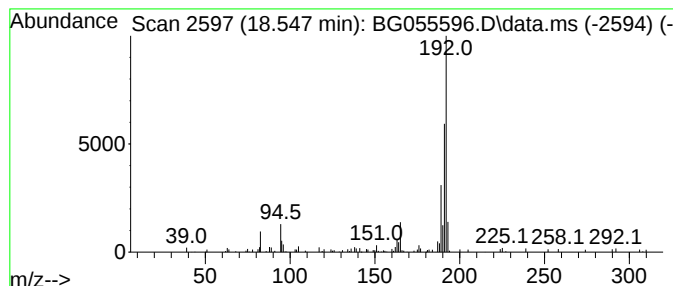
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ClientSampleId :
SU-02-111022

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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 Phenanthrene, 4-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.547	2.36 ng	96463	Phenanthrene-d10	17.624	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	91
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	83
5	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111222\
Data File : BG055596.D
Acq On : 12 Nov 2022 19:26
Operator : CG/JU
Sample : N5580-01
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ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SU-02-111022

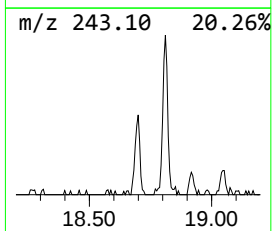
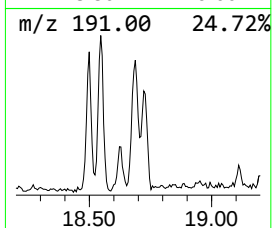
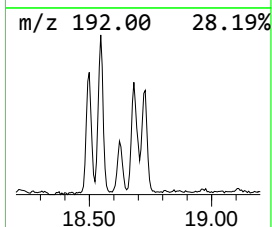
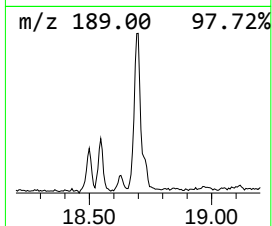
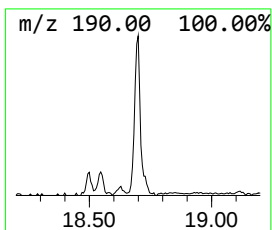
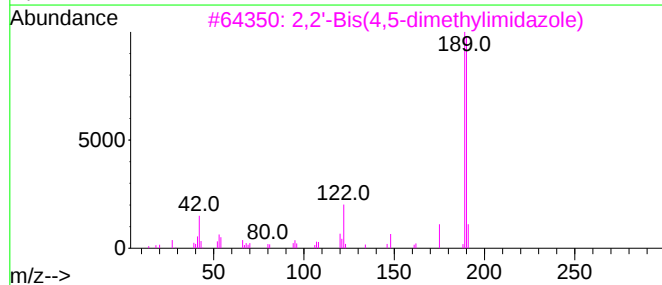
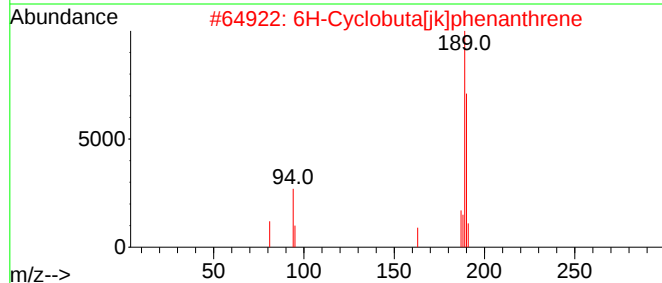
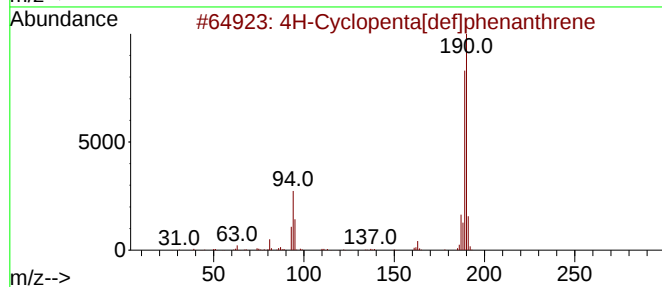
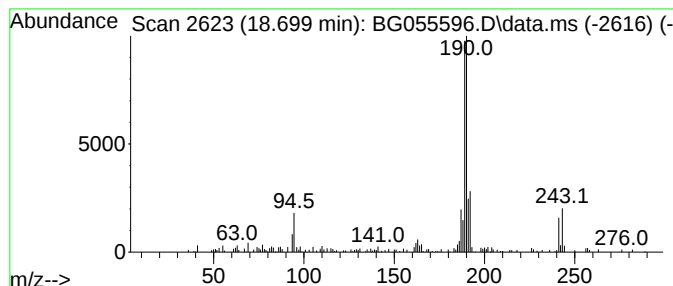
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.699	4.33 ng	176838	Phenanthrene-d10	17.624

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	68
2			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	59
3			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	38
4			2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	35
5			Methyl diselenide	190	C2H6Se2	007101-31-7	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111222\
Data File : BG055596.D
Acq On : 12 Nov 2022 19:26
Operator : CG/JU
Sample : N5580-01
Misc :
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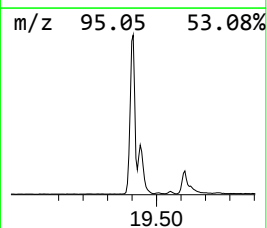
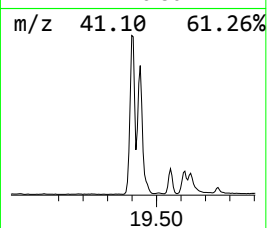
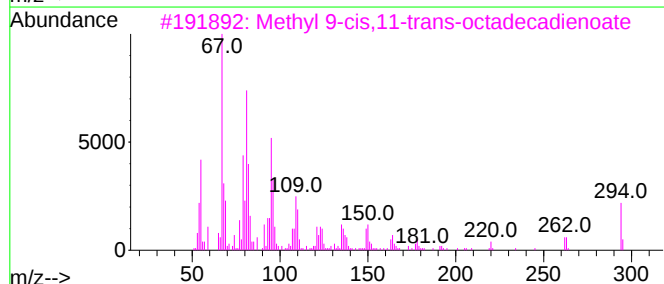
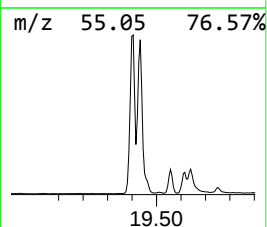
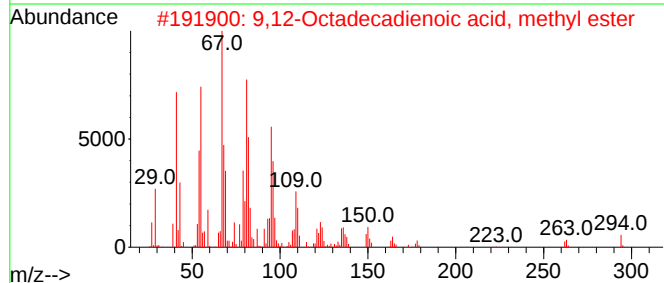
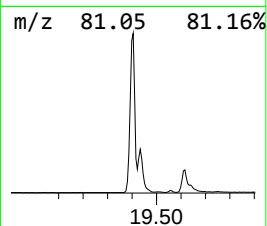
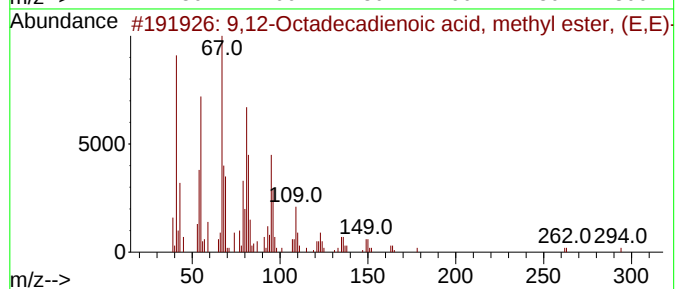
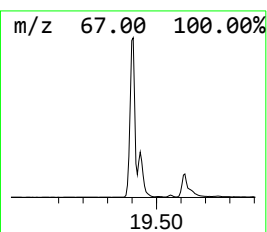
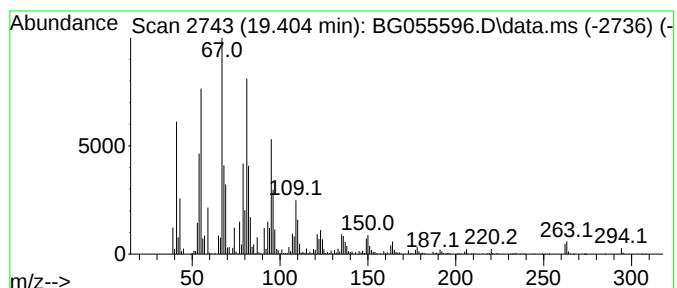
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.404	195.36 ng	7974430	Phenanthrene-d10	17.624
Hit#	of 5	Tentative ID	MW MolForm	CAS# Qual
1		9,12-Octadecadienoic acid, methy...	294 C19H34O2	002566-97-4 99
2		9,12-Octadecadienoic acid, methy...	294 C19H34O2	002462-85-3 99
3		Methyl 9-cis,11-trans-octadecadi...	294 C19H34O2	013058-52-1 99
4		8,11-Octadecadienoic acid, methy...	294 C19H34O2	056599-58-7 97
5		9,12-Octadecadienoic acid (Z,Z)-...	294 C19H34O2	000112-63-0 95



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111222\
Data File : BG055596.D
Acq On : 12 Nov 2022 19:26
Operator : CG/JU
Sample : N5580-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

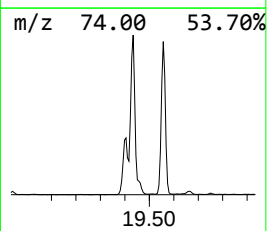
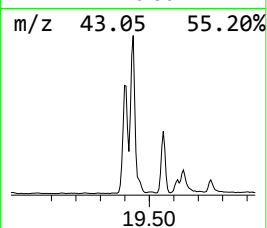
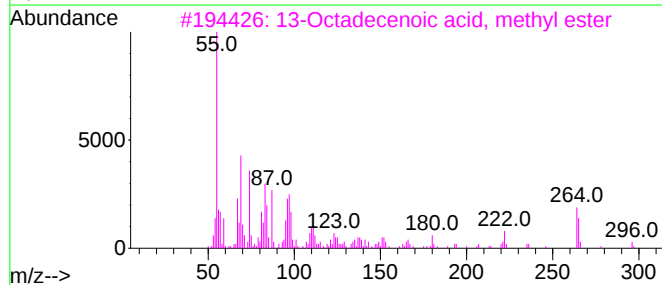
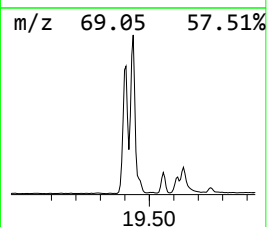
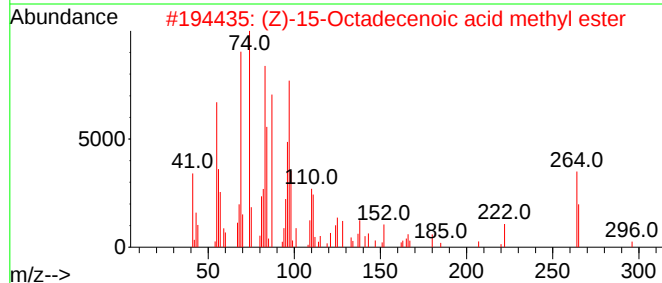
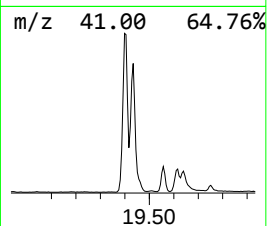
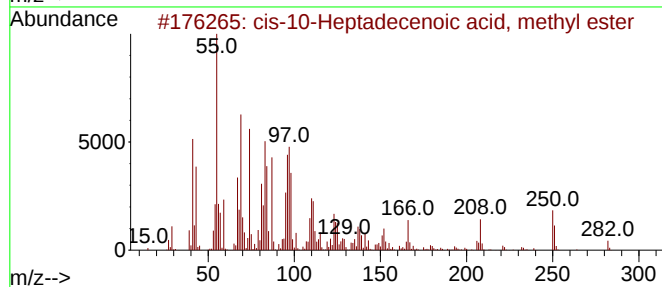
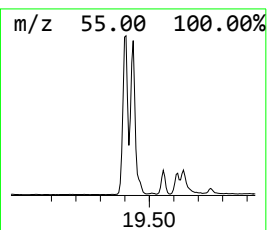
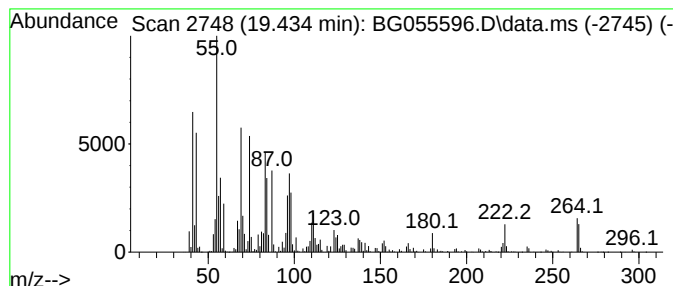
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 cis-10-Heptadecenoic acid, ... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.434	146.88 ng	5995640	Phenanthrene-d10	17.624	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	cis-10-Heptadecenoic acid, methy...	282	C18H34O2	1000333-62-1	96
2	(Z)-15-Octadecenoic acid methyl ...	296	C19H36O2	1000427-69-3	93
3	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	93
4	9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	91
5	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111222\
Data File : BG055596.D
Acq On : 12 Nov 2022 19:26
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ALS Vial : 12 Sample Multiplier: 1

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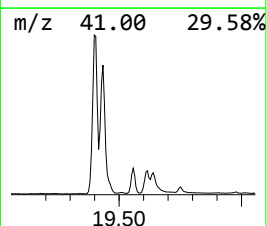
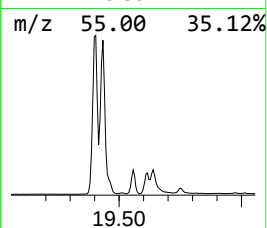
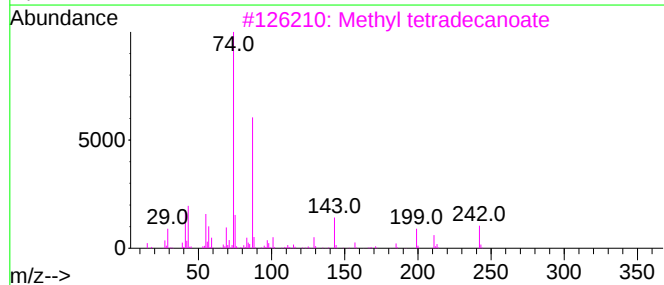
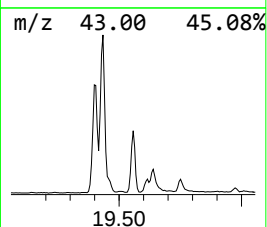
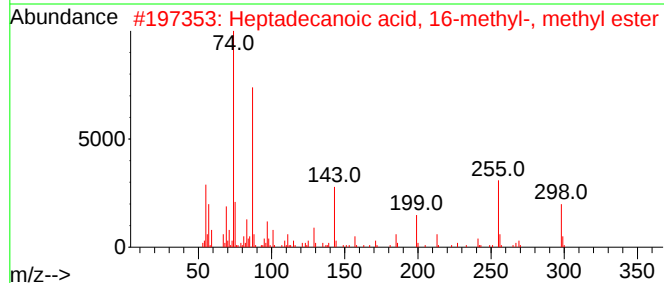
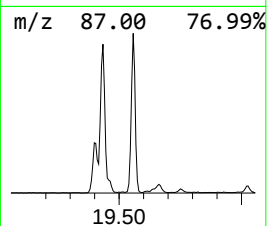
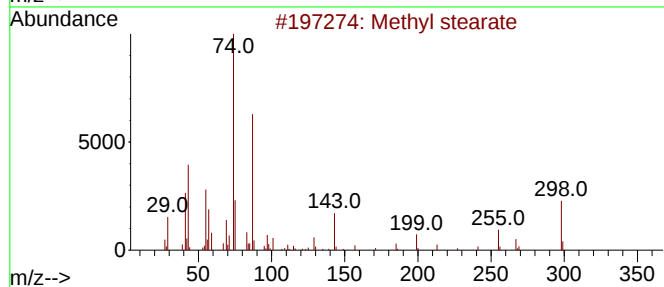
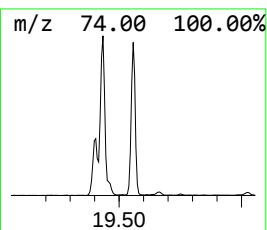
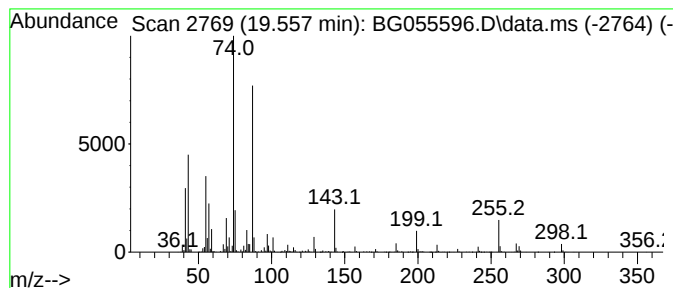
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Methyl stearate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.557	27.50 ng	1122670	Phenanthrene-d10	17.624

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl stearate	298	C19H38O2	000112-61-8	97
2			Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	95
3			Methyl tetradecanoate	242	C15H30O2	000124-10-7	90
4			Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	87
5			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111222\
Data File : BG055596.D
Acq On : 12 Nov 2022 19:26
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ALS Vial : 12 Sample Multiplier: 1

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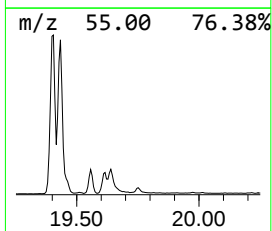
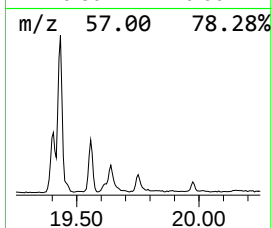
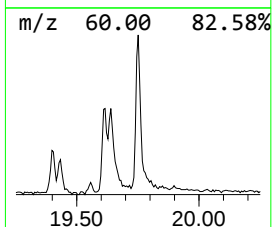
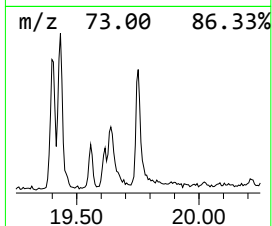
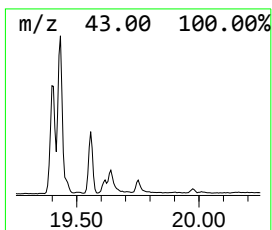
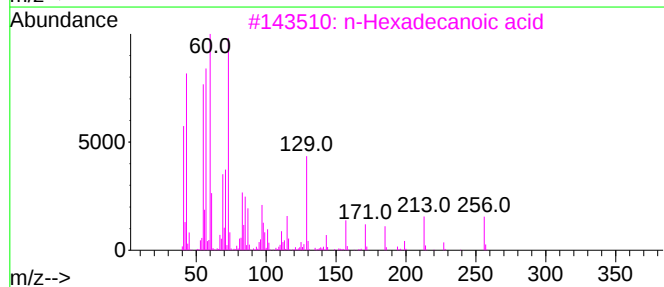
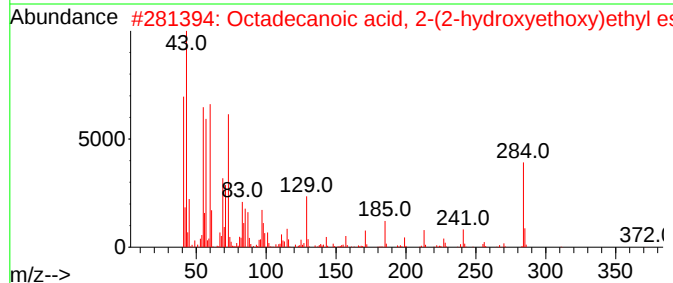
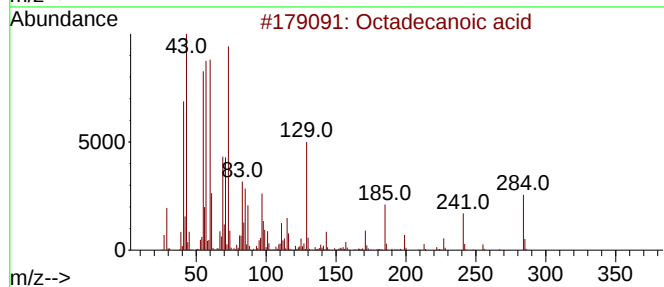
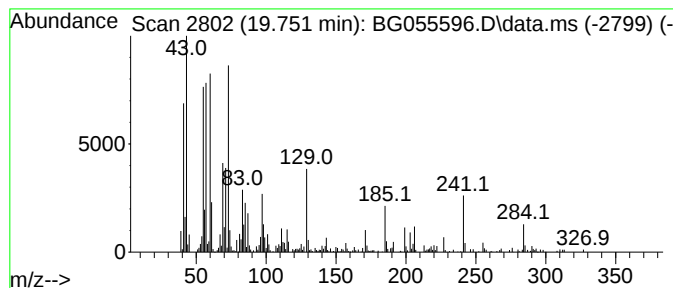
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.751	6.23 ng	254240	Phenanthrene-d10	17.624

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	86
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	76
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5			n-Decanoic acid	172	C10H20O2	000334-48-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111222\
Data File : BG055596.D
Acq On : 12 Nov 2022 19:26
Operator : CG/JU
Sample : N5580-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SU-02-111022

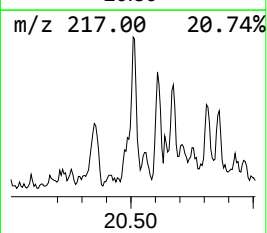
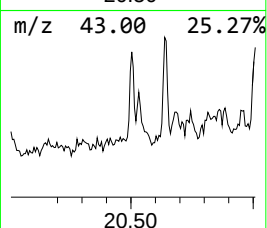
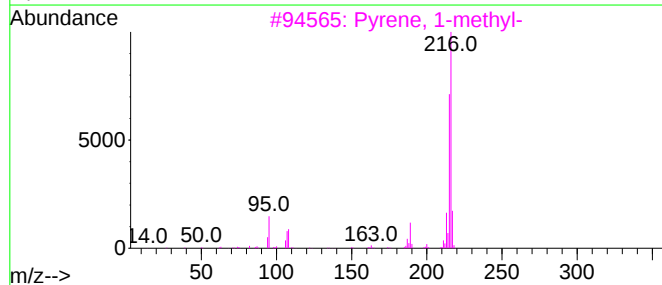
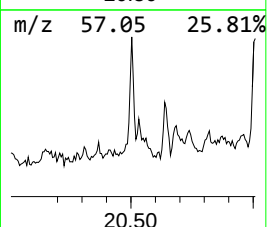
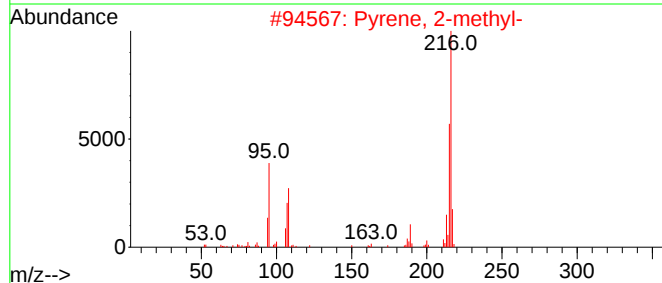
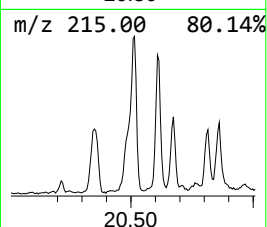
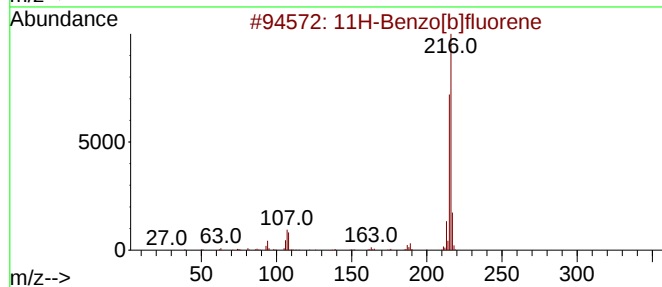
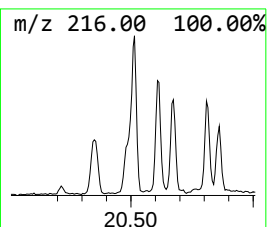
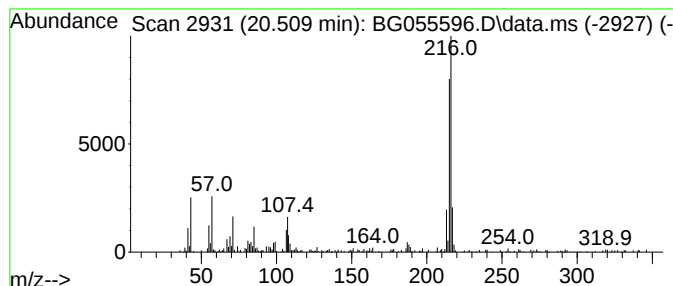
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 11H-Benzo[b]fluorene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.509	4.01 ng	277224	Chrysene-d12	21.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2			Pyrene, 2-methyl-	216	C17H12	003442-78-2	90
3			Pyrene, 1-methyl-	216	C17H12	002381-21-7	86
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	74
5			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111222\
Data File : BG055596.D
Acq On : 12 Nov 2022 19:26
Operator : CG/JU
Sample : N5580-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

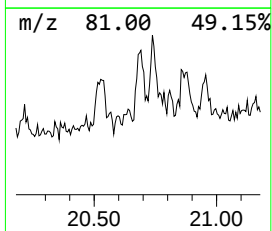
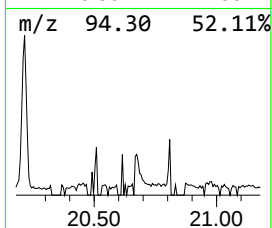
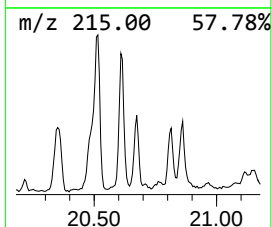
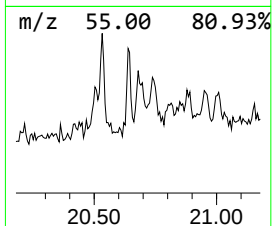
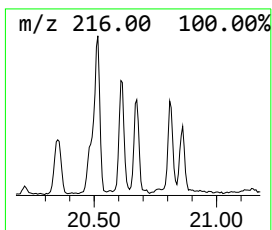
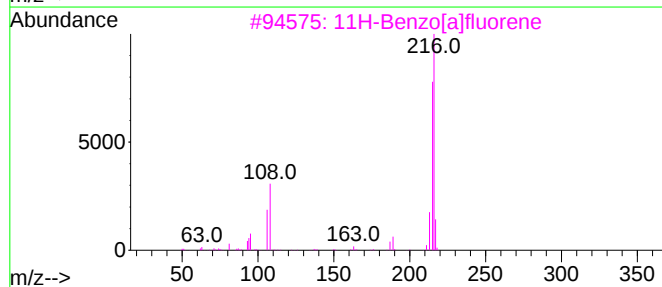
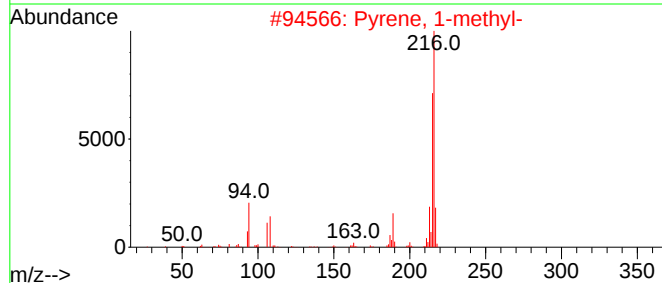
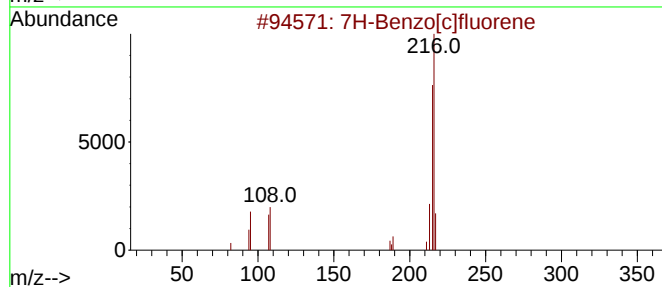
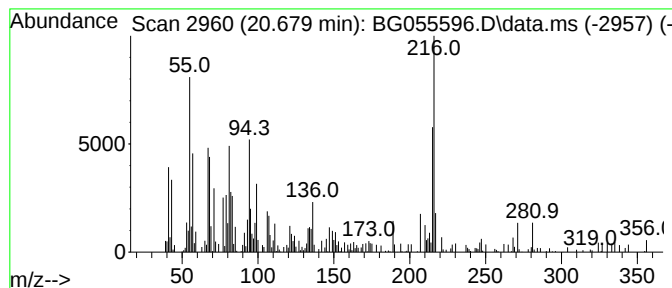
Instrument :
BNA_G
ClientSampleId :
SU-02-111022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown20.679 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.679	2.18 ng	150601	Chrysene-d12	21.925	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	49
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	43
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	38
4	Hydrazidiamantane	216	C14H20N2	1000138-39-6	35
5	1,3-Difluoro-fluoren-9-one	216	C13H6F2O	1000443-10-1	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111222\
Data File : BG055596.D
Acq On : 12 Nov 2022 19:26
Operator : CG/JU
Sample : N5580-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SU-02-111022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110922.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
Butane, 2-metho...	3.323	8.1	ng	149139	1	8.270	367447	20.0
2-Pentanone, 4-...	5.321	19.3	ng	354332	1	8.270	367447	20.0
Ethanol, 2-(2-b...	10.838	3.2	ng	86361	2	11.097	543757	20.0
Hexadecanoic ac...	18.270	37.6	ng	1533480	4	17.624	816375	20.0
n-Hexadecanoic ...	18.494	17.5	ng	713952	4	17.624	816375	20.0
Phenanthrene, 4...	18.547	2.4	ng	96463	4	17.624	816375	20.0
4H-Cyclopenta[d...	18.699	4.3	ng	176838	4	17.624	816375	20.0
9,12-Octadecadi...	19.404	195.4	ng	7974430	4	17.624	816375	20.0
cis-10-Heptadec...	19.434	146.9	ng	5995640	4	17.624	816375	20.0
Methyl stearate	19.557	27.5	ng	1122670	4	17.624	816375	20.0
Octadecanoic acid	19.751	6.2	ng	254240	4	17.624	816375	20.0
11H-Benzo[b]flu...	20.509	4.0	ng	277224	5	21.925	1384370	20.0
unknown20.679	20.679	2.2	ng	150601	5	21.925	1384370	20.0