

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111020\
 Data File : BP003908.D
 Acq On : 10 Nov 2020 18:26
 Operator : CG/JU
 Sample : L4680-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PAV-B7-110520-0-10

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_P\METHODS\8270-BP110320.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP003908.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.934	3	8	27	rVB	5659293	7373882	13.38%	5.073%
2	5.805	487	496	507	rBV	1998694	2939368	5.33%	2.022%
3	7.446	767	775	789	rBV	1949115	3294446	5.98%	2.267%
4	8.281	909	917	923	rBV	664087	1063682	1.93%	0.732%
5	9.440	1103	1114	1121	rBV	1156098	1905330	3.46%	1.311%
6	11.110	1388	1398	1404	rBV	697080	1206883	2.19%	0.830%
7	13.522	1801	1808	1819	rVB	2448076	3506434	6.36%	2.412%
8	14.898	2036	2042	2048	rVB2	981762	1422490	2.58%	0.979%
9	16.375	2287	2293	2300	rBV	1802021	2615495	4.75%	1.799%
10	17.628	2500	2506	2510	rBV	1007397	1451516	2.63%	0.999%
11	17.669	2510	2513	2520	rVB	634373	831504	1.51%	0.572%
12	18.128	2587	2591	2597	rVB	674967	808960	1.47%	0.557%
13	18.251	2607	2612	2617	rBV	10107874	11784126	21.39%	8.108%
14	19.339	2791	2797	2799	rBV	14968441	21166328	38.42%	14.563%
15	19.381	2799	2804	2812	rVB	32812767	55099068	100.00%	37.909%
16	19.486	2817	2822	2829	rVB	5515127	6099269	11.07%	4.196%
17	19.604	2837	2842	2847	rVB	1084106	1396423	2.53%	0.961%
18	19.951	2897	2901	2909	rVB	1316011	1786498	3.24%	1.229%
19	20.128	2926	2931	2936	rVB	3447120	4122840	7.48%	2.837%
20	20.422	2974	2981	2991	rVB	1514558	2142254	3.89%	1.474%
21	20.528	2995	2999	3002	rBV2	846528	957033	1.74%	0.658%
22	20.563	3002	3005	3011	rVV2	1049445	1750858	3.18%	1.205%
23	20.616	3011	3014	3019	rVB	1006753	1223202	2.22%	0.842%
24	20.816	3044	3048	3053	rVB2	441625	589652	1.07%	0.406%
25	21.110	3095	3098	3110	rVB4	212519	596529	1.08%	0.410%
26	21.439	3151	3154	3161	rVB2	497970	600249	1.09%	0.413%
27	21.657	3184	3191	3195	rBV2	1373878	2774382	5.04%	1.909%
28	21.698	3195	3198	3205	rVB	621284	805184	1.46%	0.554%
29	23.380	3479	3484	3489	rBV2	421020	965354	1.75%	0.664%
30	23.916	3571	3575	3585	rVB	379364	702053	1.27%	0.483%
31	24.021	3588	3593	3600	rVB	466468	925666	1.68%	0.637%
32	24.133	3606	3612	3618	rBV	733034	1439466	2.61%	0.990%

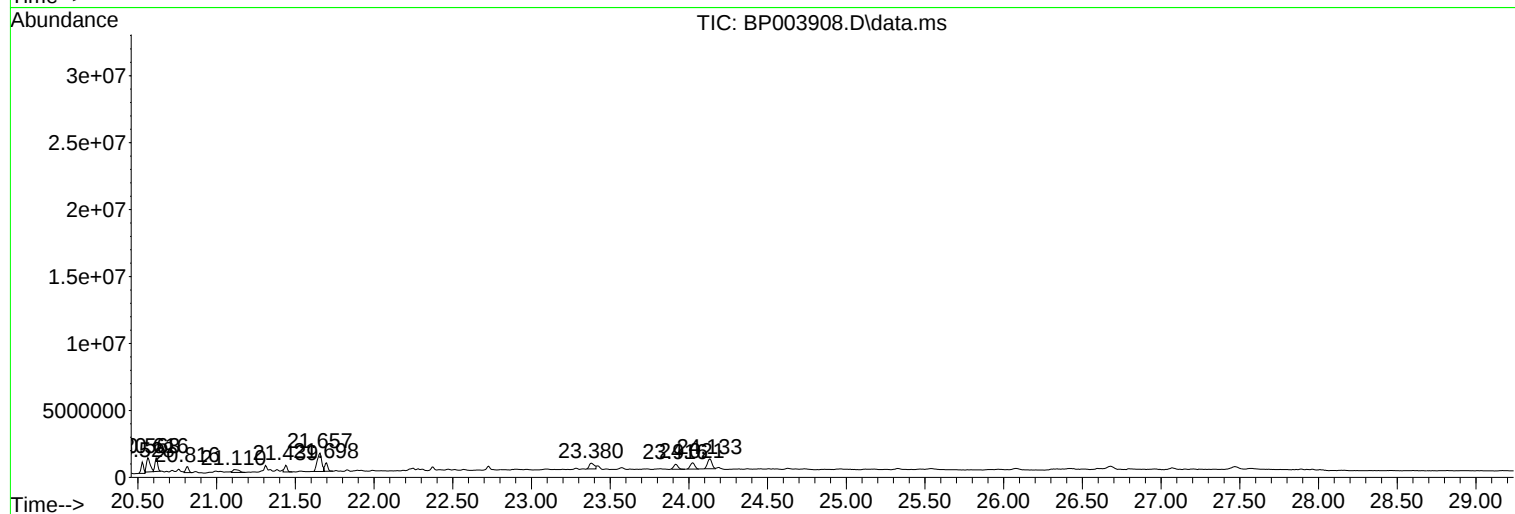
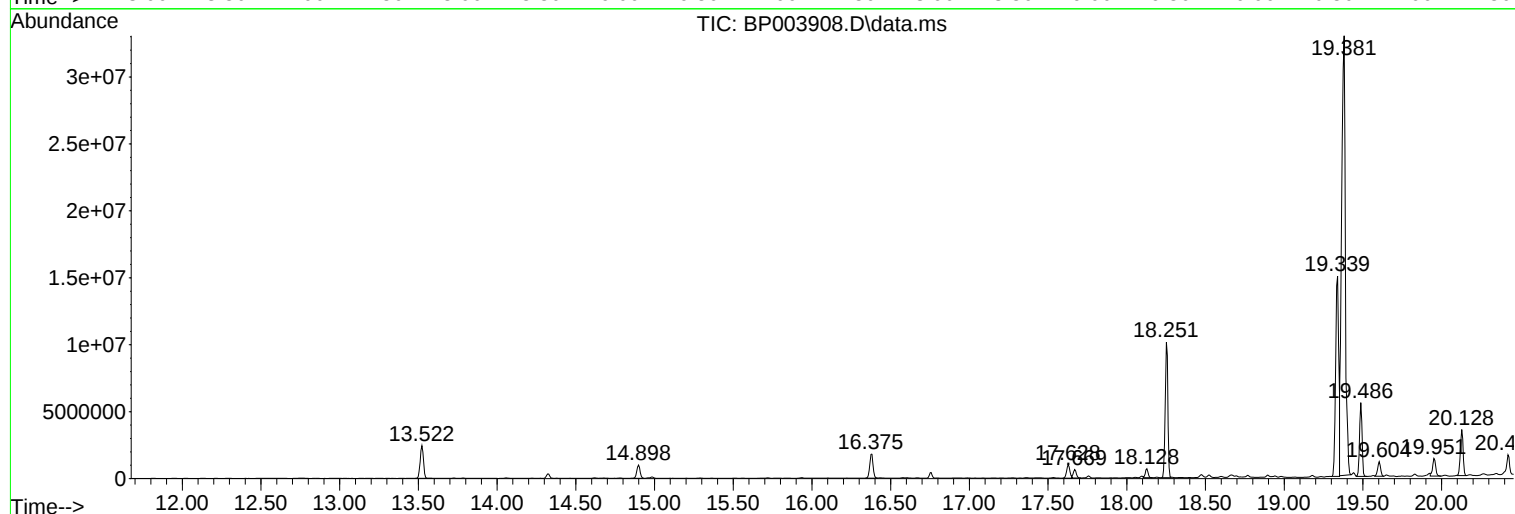
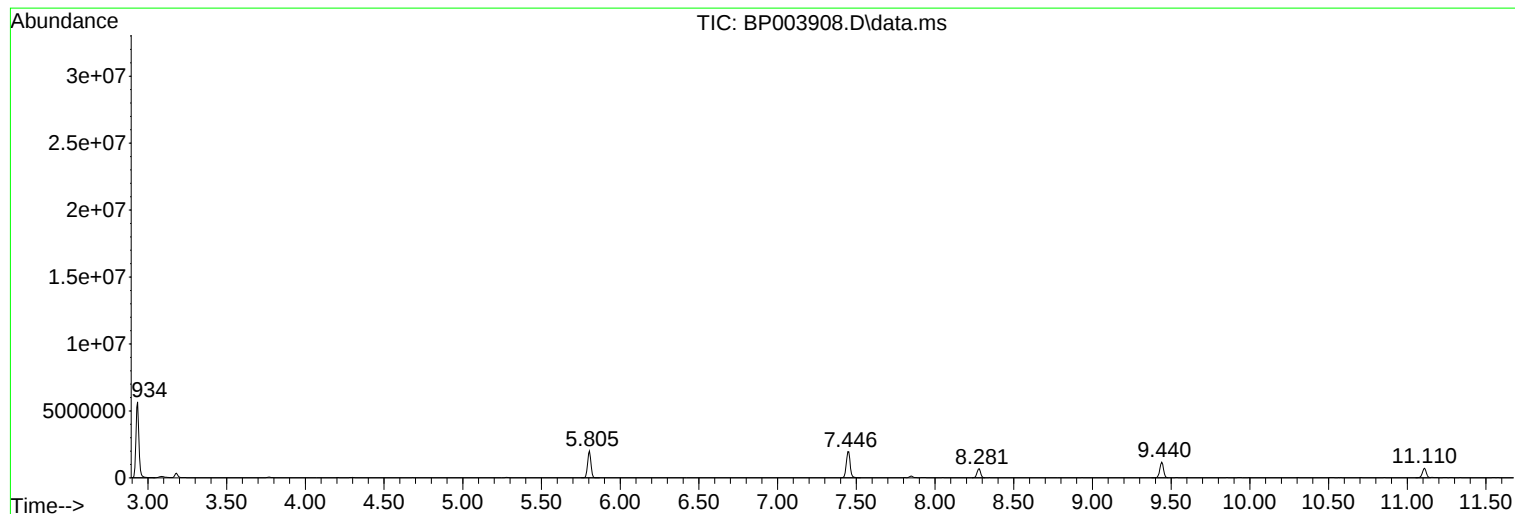
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Instrument :
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ClientSampleId :
PAV-B7-110520-0-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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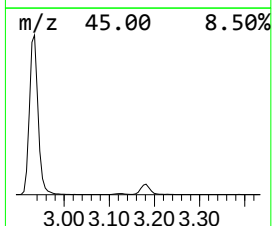
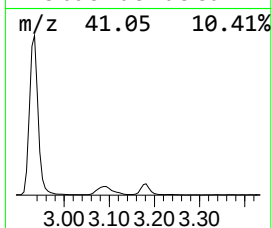
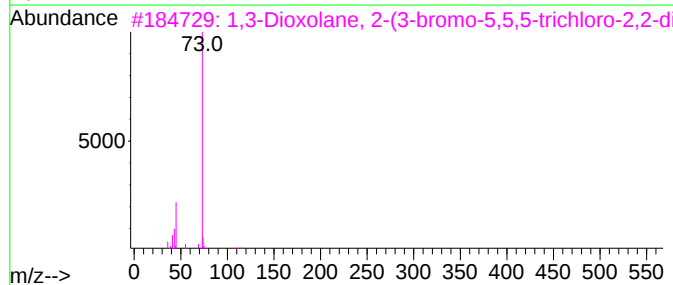
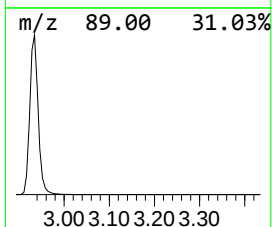
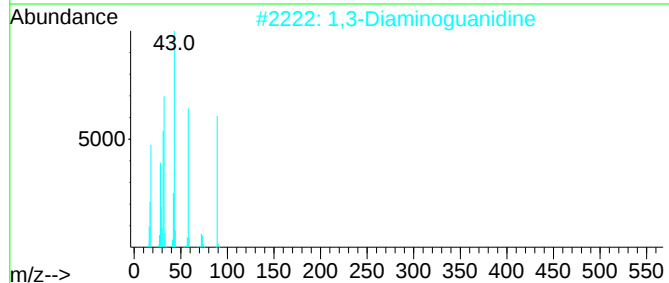
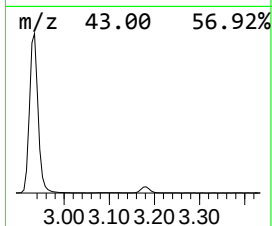
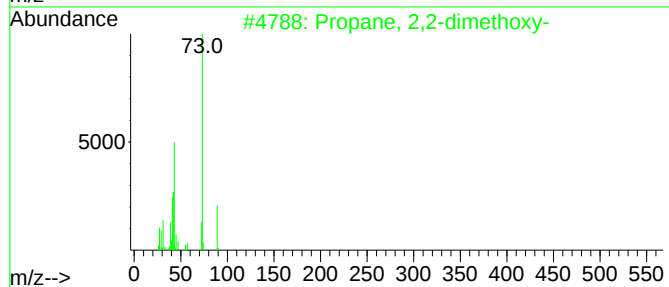
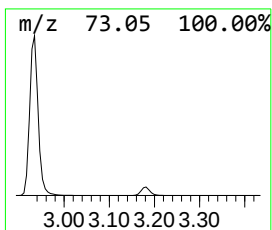
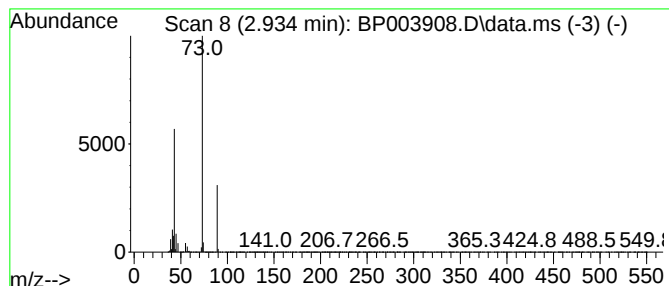
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown2.934 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.934	138.65 ng	7373880	1,4-Dichlorobenzene-d4	8.281

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	38
2			1,3-Diaminoguanidine	89	CH7N5	004364-78-7	9
3			1,3-Dioxolane, 2-(3-bromo-5,5,5-...	352	C10H16BrCl3O2	1000115-31-4	9
4			2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9
5			Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9



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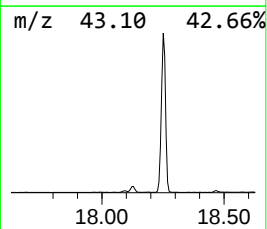
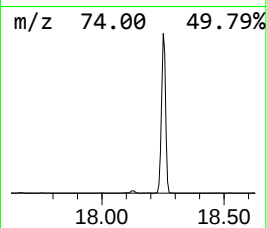
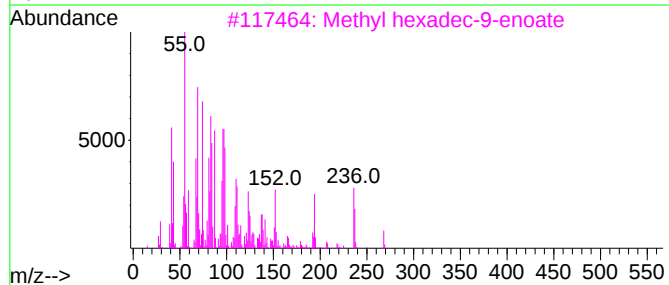
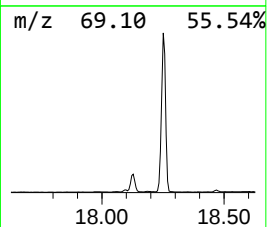
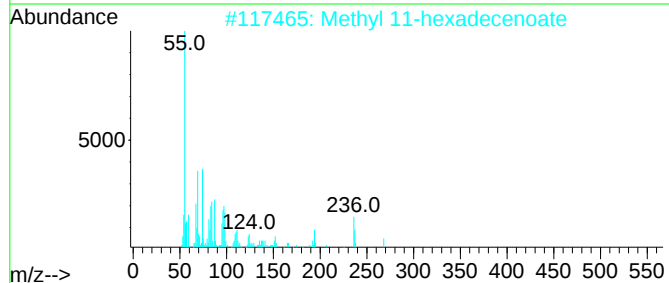
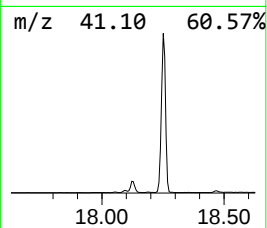
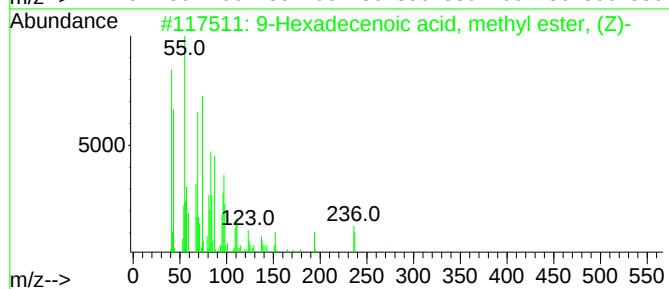
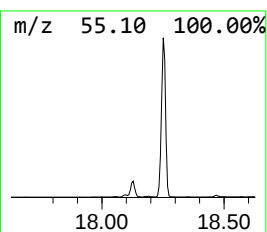
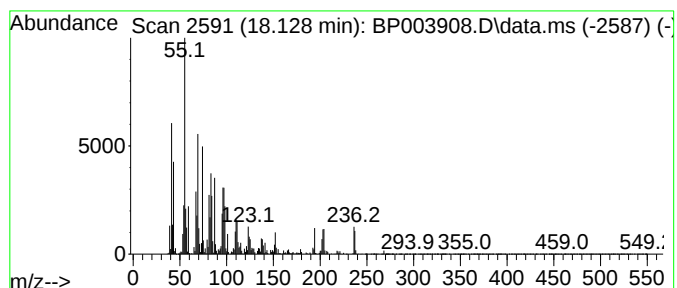
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 9-Hexadecenoic acid, methyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.128	11.15 ng	808960	Phenanthrene-d10	17.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Hexadecenoic acid, methyl este...	268	C17H32O2	001120-25-8	99
2			Methyl 11-hexadecenoate	268	C17H32O2	1000336-35-5	91
3			Methyl hexadec-9-enoate	268	C17H32O2	010030-74-7	90
4			Methyl myristoleate	240	C15H28O2	056219-06-8	87
5			9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	83



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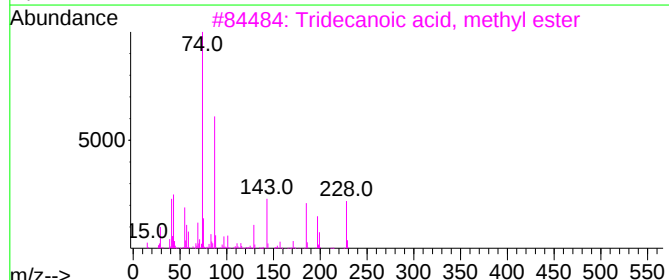
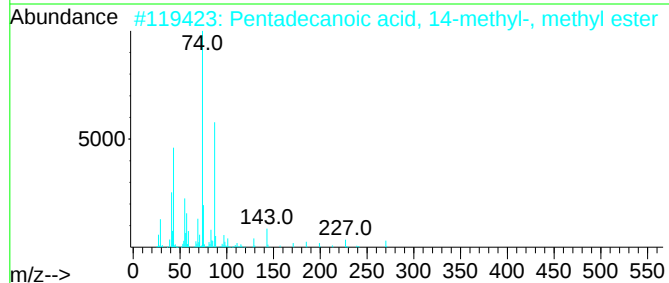
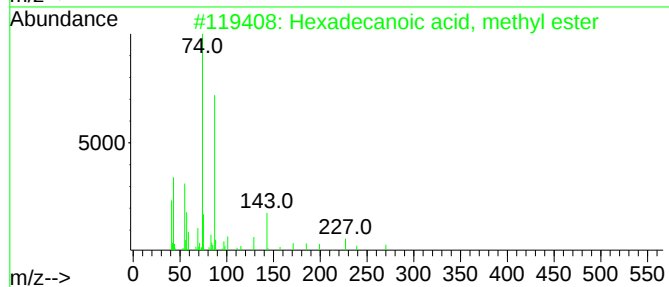
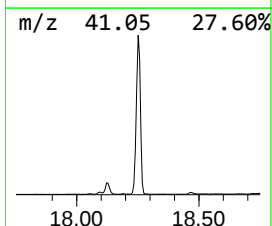
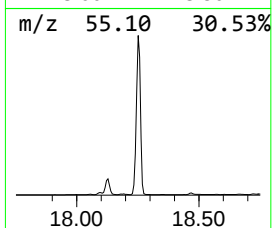
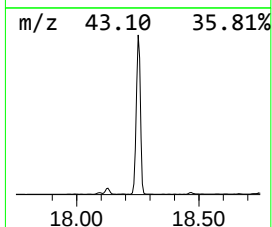
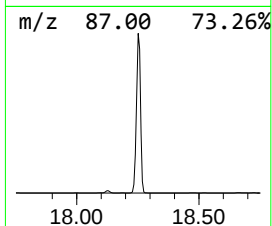
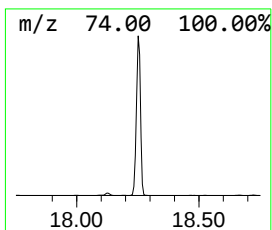
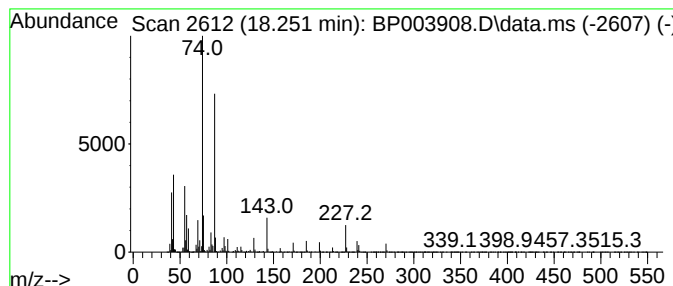
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.251	162.37 ng	11784100	Phenanthrene-d10	17.628

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	96
3			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	91
4			Methyl tetradecanoate	242	C15H30O2	000124-10-7	91
5			Tridecanoic acid, 12-methyl-, me...	242	C15H30O2	005129-58-8	87



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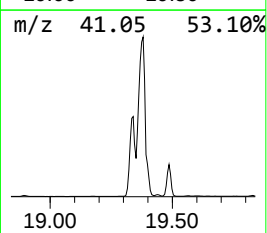
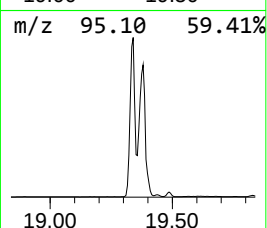
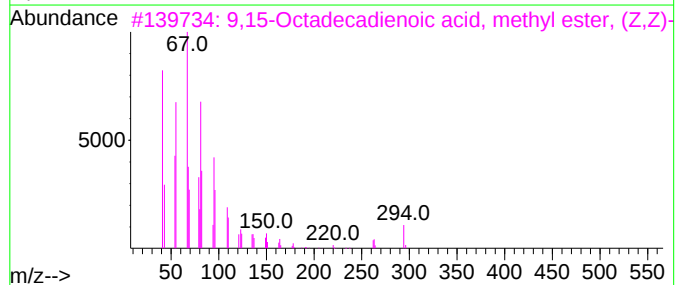
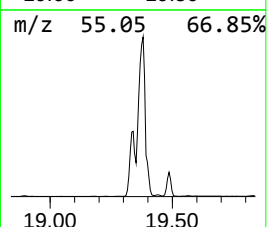
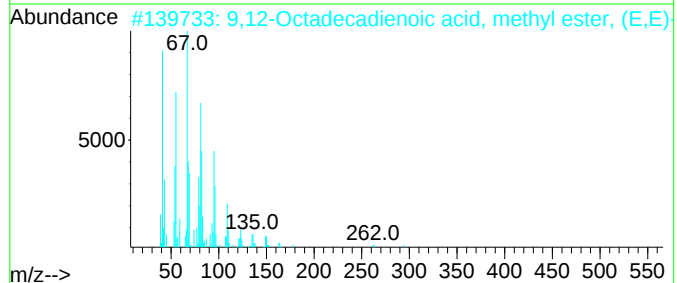
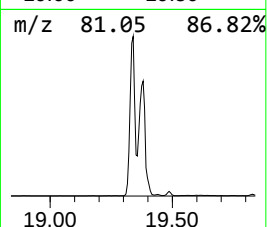
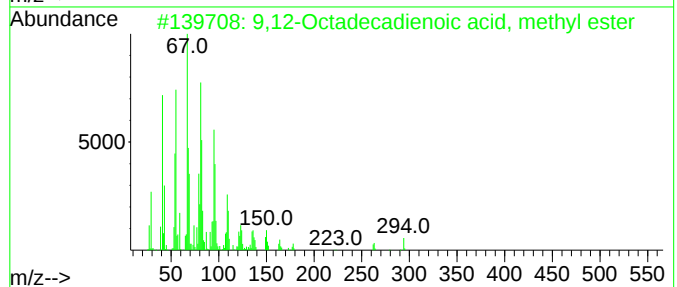
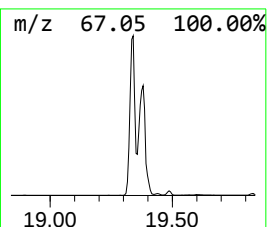
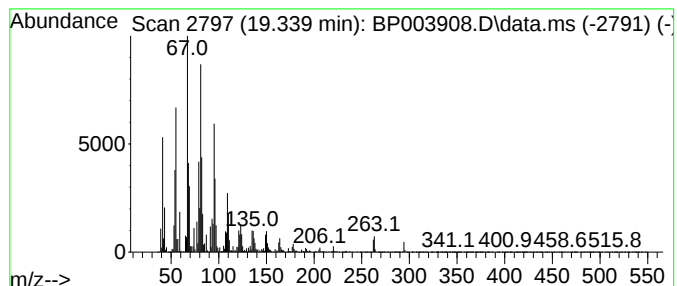
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 9,12-Octadecadienoic acid, ... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.339	291.64 ng	21166300	Phenanthrene-d10	17.628	
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	10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99
5	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99



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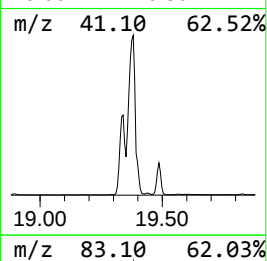
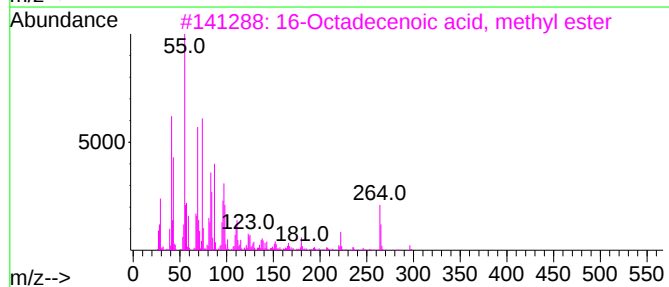
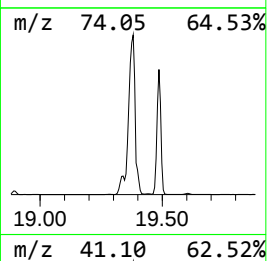
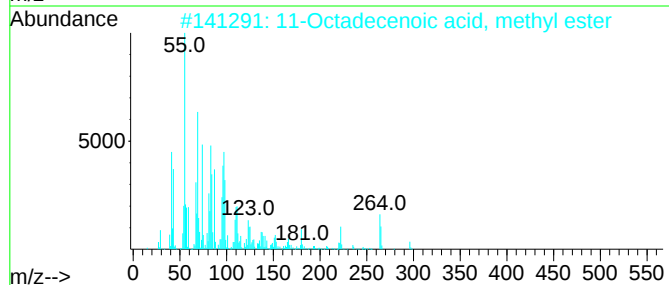
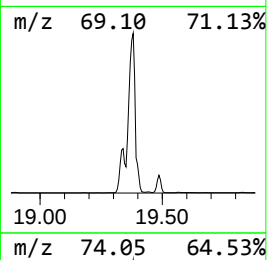
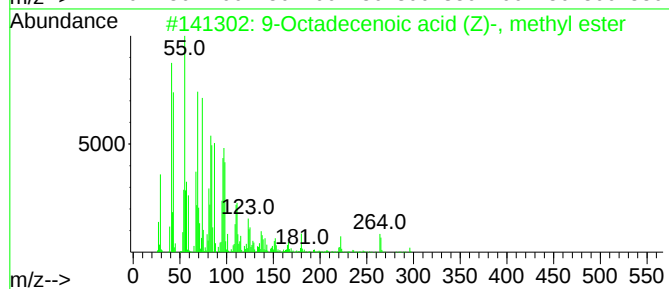
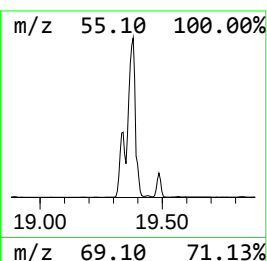
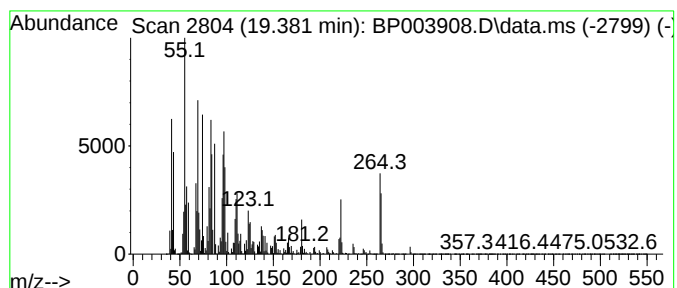
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Peak Number 5 9-Octadecenoic acid (Z)-, m... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.381	759.19 ng	55099100	Phenanthrene-d10	17.628	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2	11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	99
3	16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	99
4	15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	99
5	14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	99



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ClientSampleId :
PAV-B7-110520-0-10

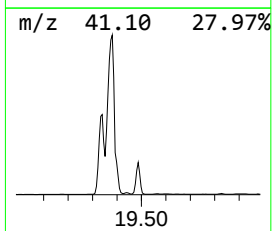
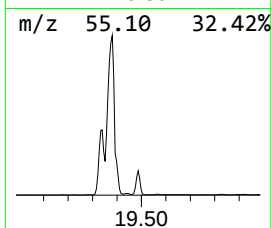
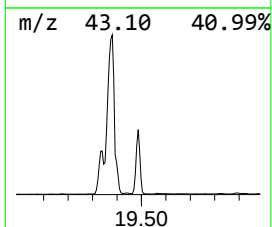
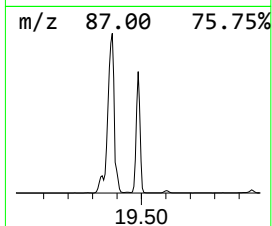
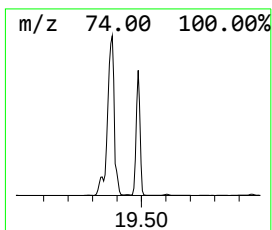
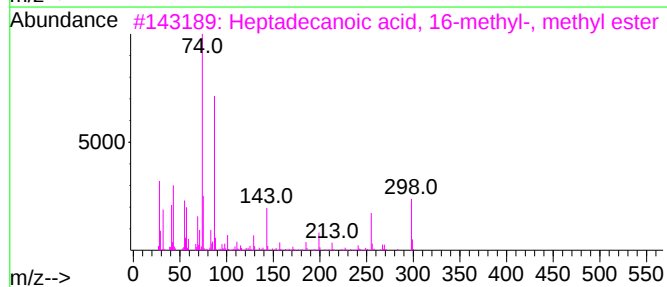
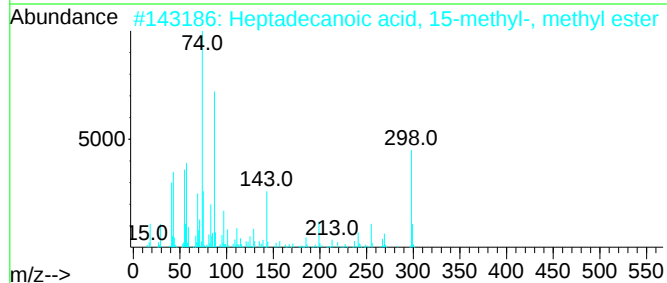
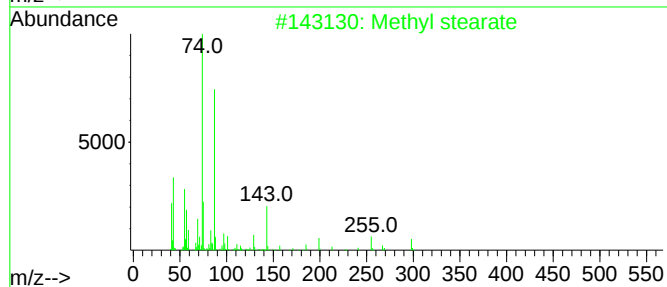
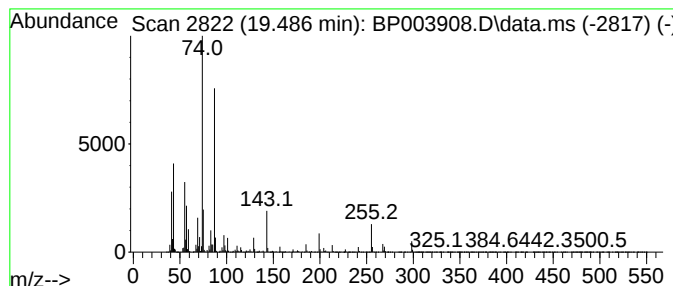
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Methyl stearate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.486	84.04 ng	6099270	Phenanthrene-d10	17.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl stearate	298	C19H38O2	000112-61-8	98
2			Heptadecanoic acid, 15-methyl-, ...	298	C19H38O2	054833-55-5	94
3			Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	92
4			Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	90
5			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111020\
Data File : BP003908.D
Acq On : 10 Nov 2020 18:26
Operator : CG/JU
Sample : L4680-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PAV-B7-110520-0-10

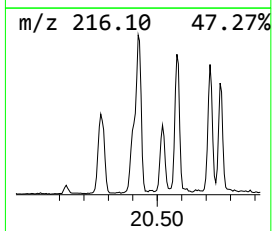
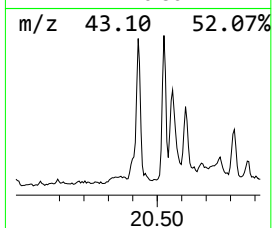
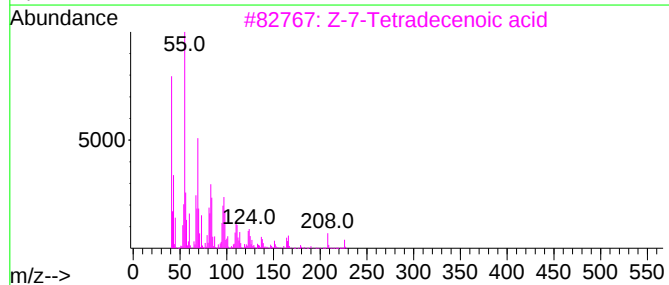
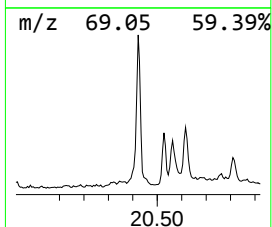
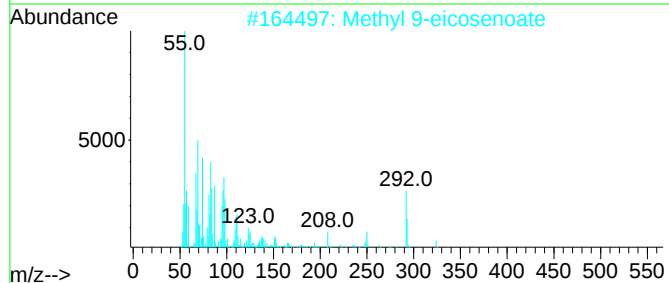
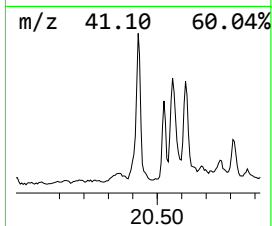
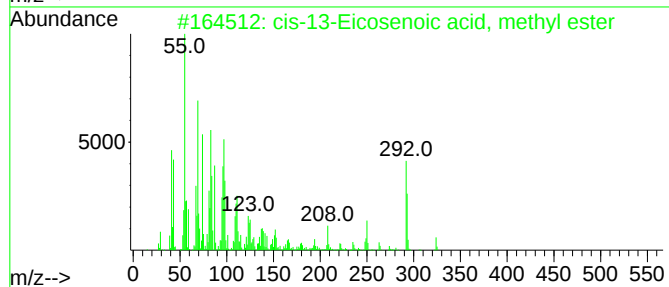
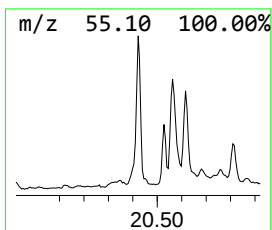
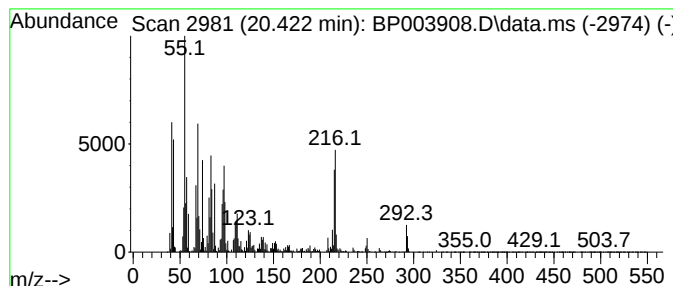
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 cis-13-Eicosenoic acid, met... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.422	15.44 ng	2142250	Chrysene-d12	21.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-13-Eicosenoic acid, methyl e...	324	C21H40O2	1000333-52-1	99
2			Methyl 9-eicosenoate	324	C21H40O2	1000336-50-5	99
3			Z-7-Tetradecenoic acid	226	C14H26O2	1000130-98-4	70
4			cis-11-Eicosenoic acid, methyl e...	324	C21H40O2	1000333-63-8	70
5			11-Eicosenoic acid, methyl ester	324	C21H40O2	003946-08-5	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111020\
Data File : BP003908.D
Acq On : 10 Nov 2020 18:26
Operator : CG/JU
Sample : L4680-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

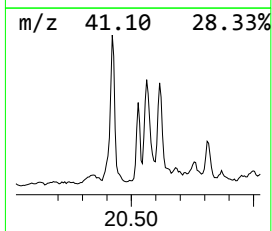
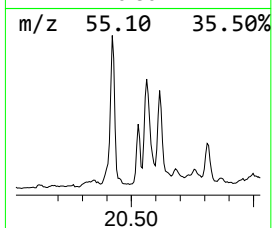
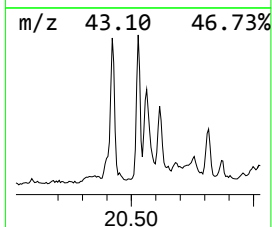
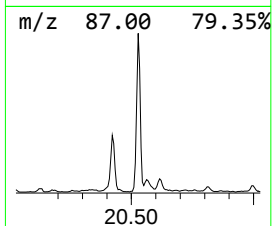
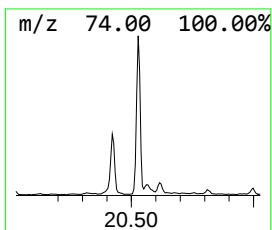
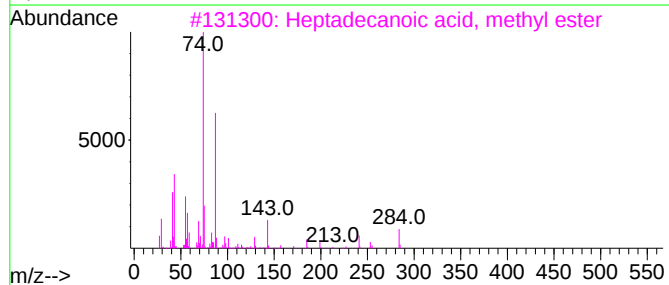
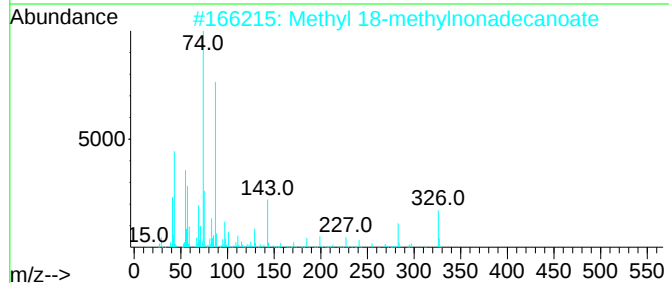
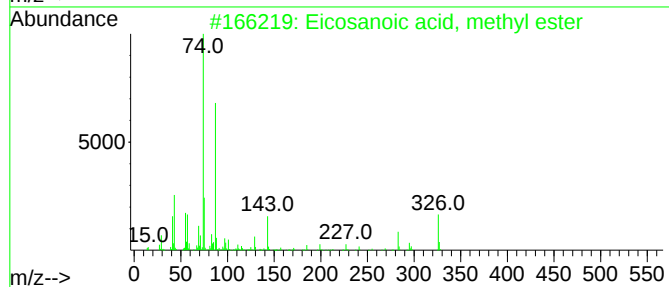
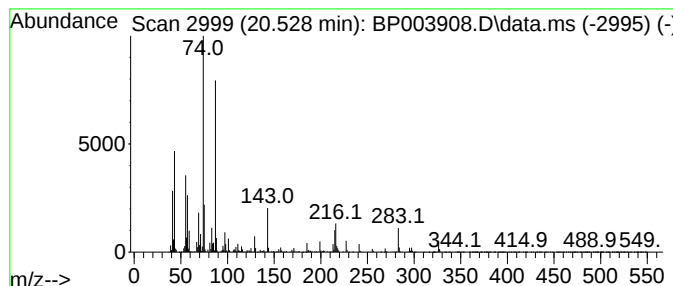
Instrument :
BNA_P
ClientSampleId :
PAV-B7-110520-0-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Eicosanoic acid, methyl ester Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.528	6.90 ng	957033	Chrysene-d12		21.657	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Eicosanoic acid, methyl ester		326	C21H42O2	001120-28-1	98
2	Methyl 18-methylnonadecanoate		326	C21H42O2	1000352-20-6	97
3	Heptadecanoic acid, methyl ester		284	C18H36O2	001731-92-6	94
4	Methyl stearate		298	C19H38O2	000112-61-8	94
5	Hexadecanoic acid, 15-methyl-, m...		284	C18H36O2	006929-04-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111020\
Data File : BP003908.D
Acq On : 10 Nov 2020 18:26
Operator : CG/JU
Sample : L4680-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PAV-B7-110520-0-10

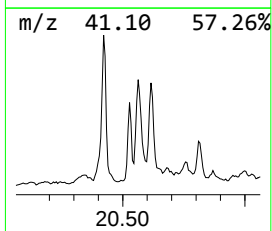
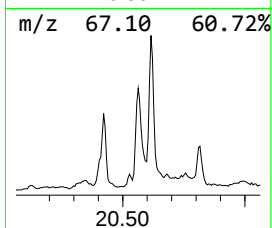
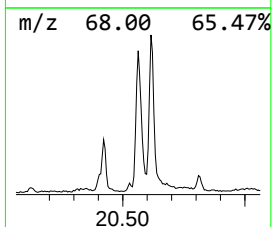
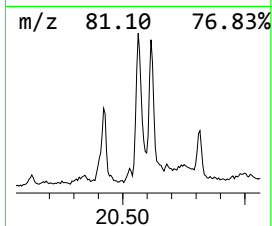
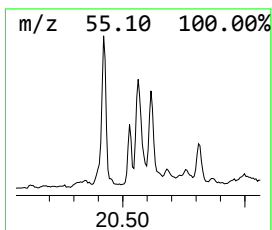
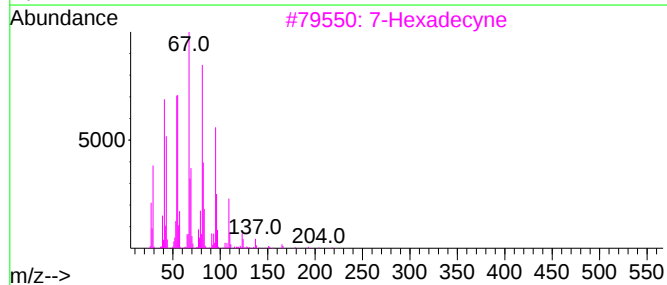
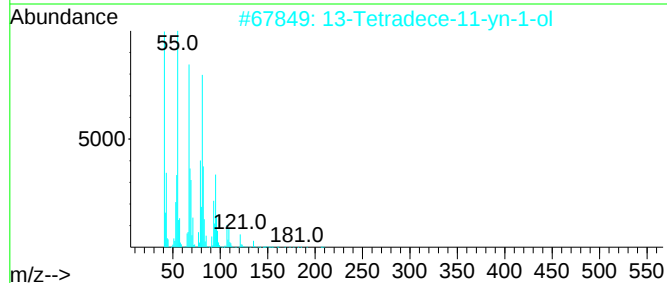
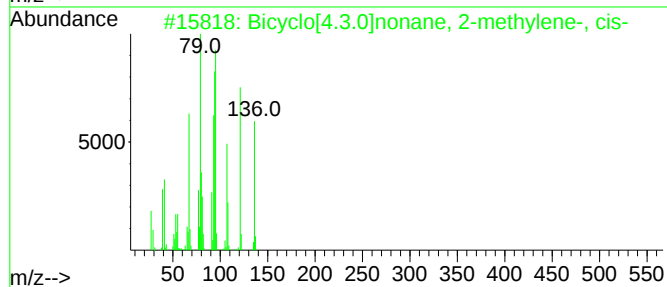
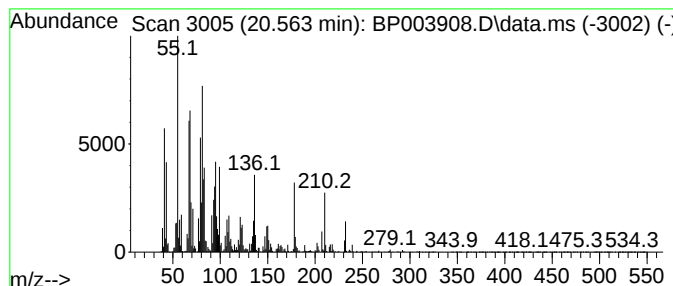
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown20.563 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.563	12.62 ng	1750860	Chrysene-d12	21.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[4.3.0]nonane, 2-methylen...	136	C10H16	040954-37-8	47
2			13-Tetradecene-11-yn-1-ol	208	C14H24O	1000131-00-4	46
3			7-Hexadecyne	222	C16H30	074685-28-2	42
4			9-Oxabicyclo[3.3.1]non-6-en-2-ol...	140	C8H12O2	029946-76-7	38
5			1,E-8,Z-10-Dodecatriene	164	C12H20	1000098-05-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111020\
Data File : BP003908.D
Acq On : 10 Nov 2020 18:26
Operator : CG/JU
Sample : L4680-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

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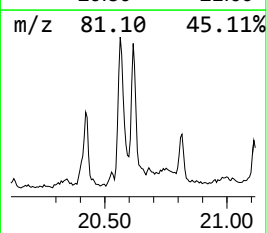
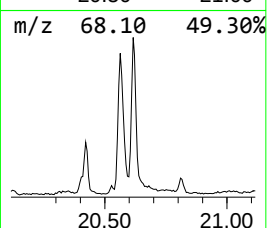
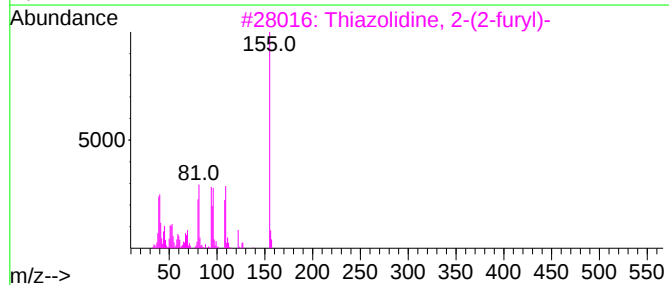
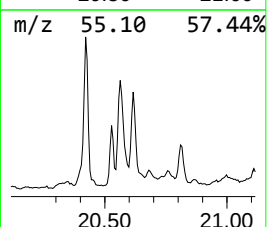
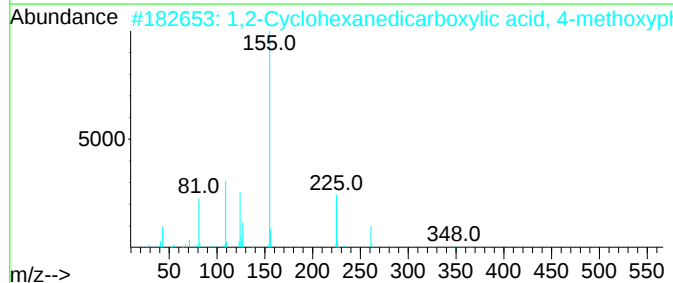
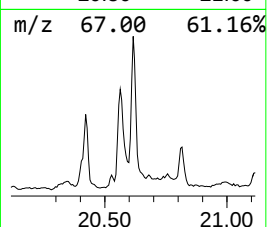
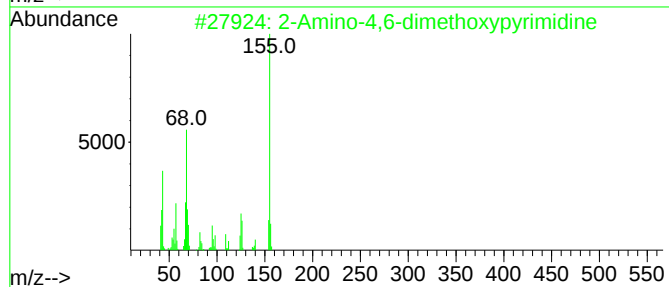
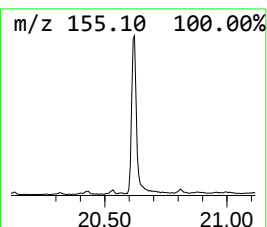
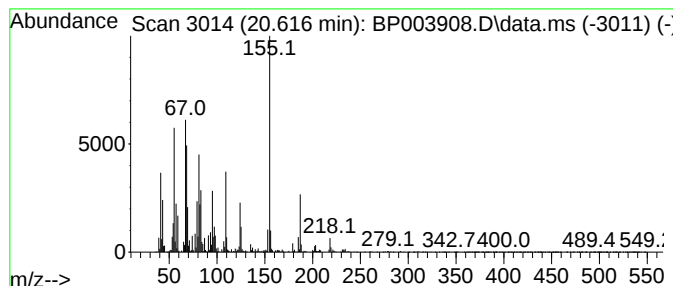
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown20.616 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.616	8.82 ng	1223200	Chrysene-d12	21.657

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Amino-4,6-dimethoxypyrimidine	155	C6H9N3O2	036315-01-2	43
2		1,2-Cyclohexanedicarboxylic acid...	348	C20H28O5	1000339-66-9	38
3		Thiazolidine, 2-(2-furyl)-	155	C7H9NOS	051859-60-0	38
4		1,2-Cyclohexanedicarboxylic acid...	404	C24H36O5	1000339-67-4	35
5		2,6-Difluoro-3-methylbenzoic aci...	384	C14H6Cl4F2O2	1000357-30-0	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111020\
Data File : BP003908.D
Acq On : 10 Nov 2020 18:26
Operator : CG/JU
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Misc :
ALS Vial : 13 Sample Multiplier: 1

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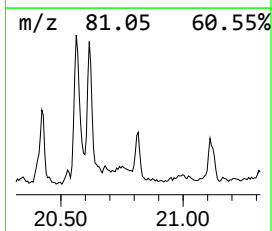
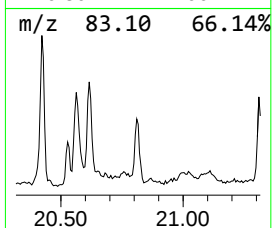
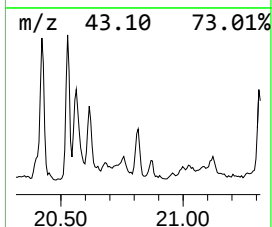
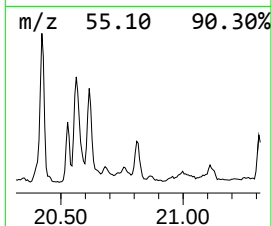
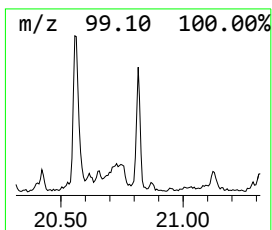
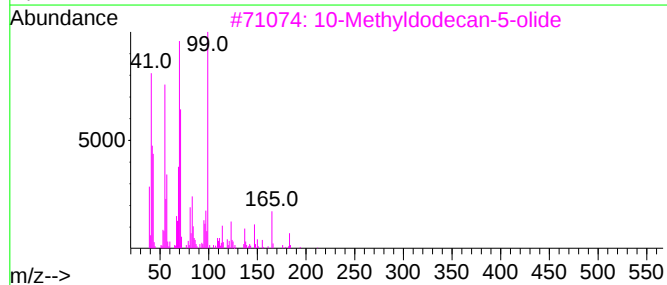
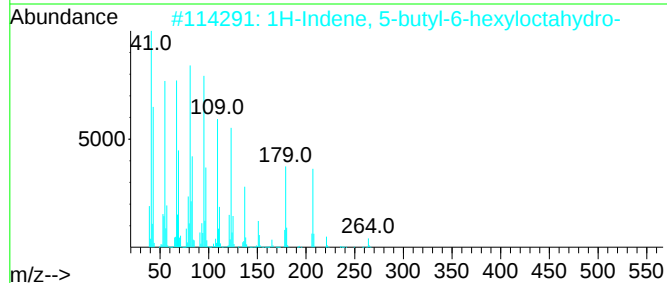
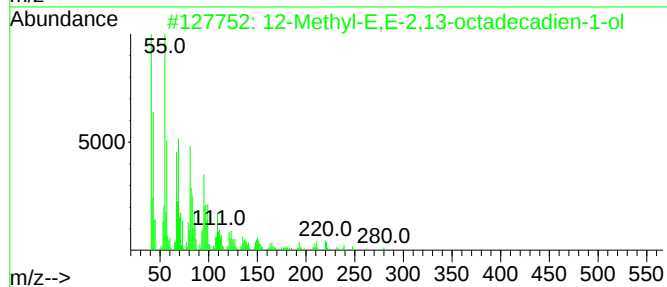
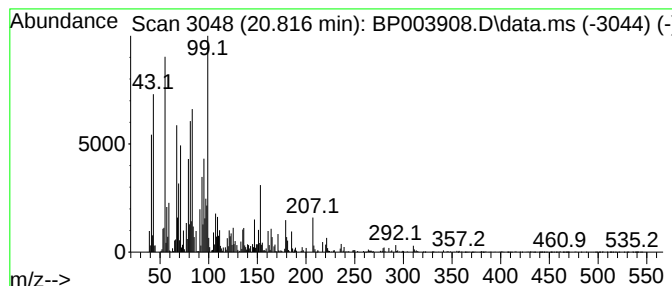
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 12-Methyl-E,E-2,13-octadeca... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.816	4.25 ng	589652	Chrysene-d12	21.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			12-Methyl-E,E-2,13-octadecadien-...	280	C19H360	1000130-90-4	51
2			1H-Indene, 5-butyl-6-hexyloctahy...	264	C19H36	055044-36-5	50
3			10-Methyldodecan-5-olide	212	C13H24O2	1000370-40-8	30
4			Thiazole, 5-methyl-	99	C4H5NS	003581-89-3	30
5			Cyclohexanol, 1-(1,2-propadienyl)-	138	C9H14O	034761-56-3	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111020\
Data File : BP003908.D
Acq On : 10 Nov 2020 18:26
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ALS Vial : 13 Sample Multiplier: 1

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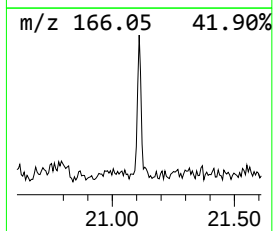
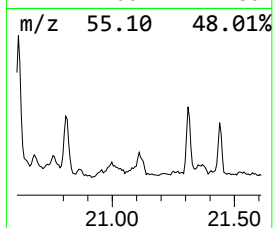
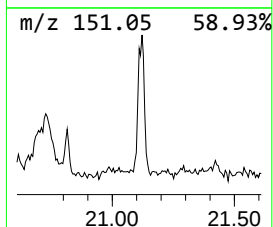
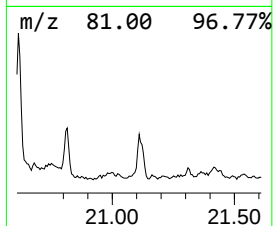
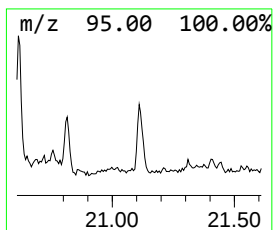
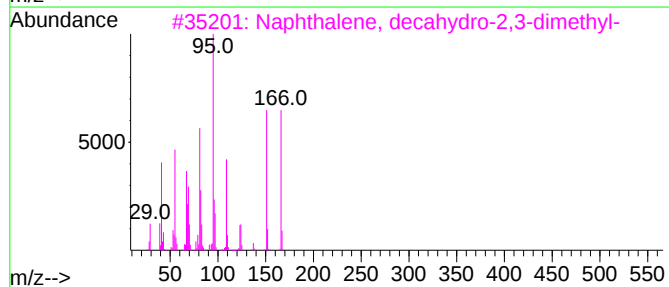
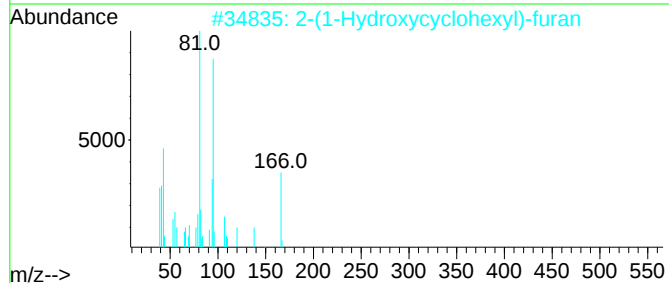
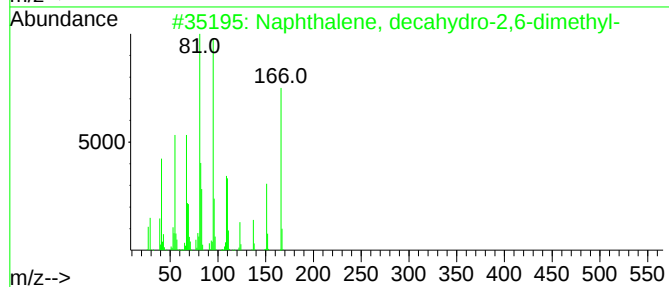
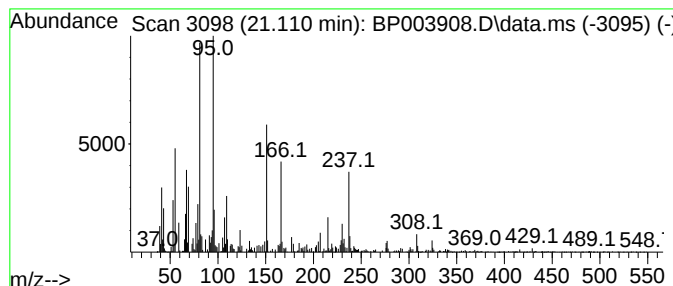
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Naphthalene, decahydro-2,6-... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.110	4.30 ng	596529	Chrysene-d12	21.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-2,6-dimet...	166	C12H22	001618-22-0	74
2			2-(1-Hydroxycyclohexyl)-furan	166	C10H14O2	036169-67-2	52
3			Naphthalene, decahydro-2,3-dimet...	166	C12H22	001008-80-6	30
4			2,4-Hexadiene, 3-methyl-	96	C7H12	028823-42-9	20
5			2-Hexyne, 4-methyl-	96	C7H12	020198-49-6	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111020\
Data File : BP003908.D
Acq On : 10 Nov 2020 18:26
Operator : CG/JU
Sample : L4680-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

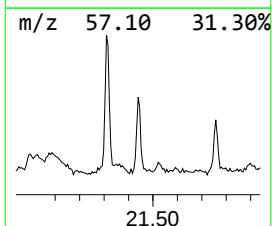
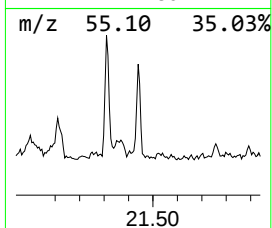
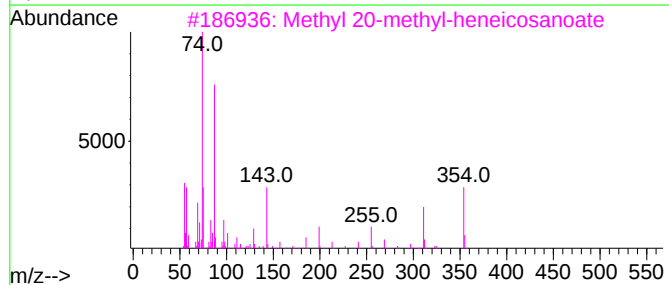
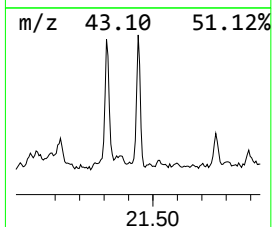
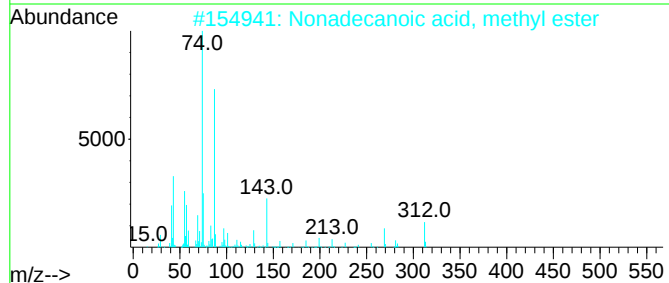
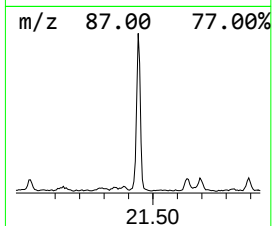
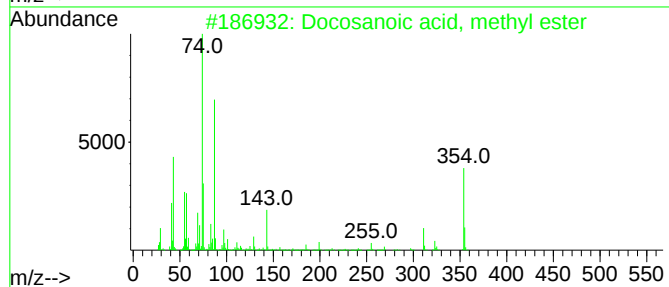
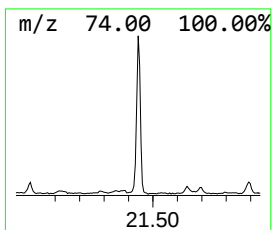
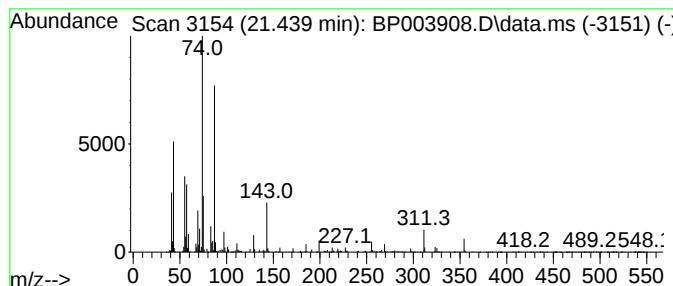
Instrument :
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ClientSampleId :
PAV-B7-110520-0-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Docosanoic acid, methyl ester Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.439	4.33 ng	600249	Chrysene-d12	21.657	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosanoic acid, methyl ester	354	C23H46O2	000929-77-1	95
2	Nonadecanoic acid, methyl ester	312	C20H40O2	001731-94-8	93
3	Methyl 20-methyl-heneicosanoate	354	C23H46O2	1000336-47-4	90
4	Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	87
5	Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111020\
Data File : BP003908.D
Acq On : 10 Nov 2020 18:26
Operator : CG/JU
Sample : L4680-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PAV-B7-110520-0-10

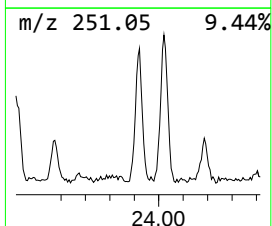
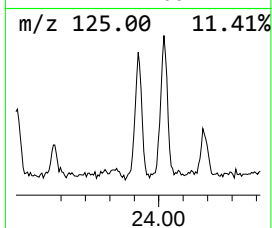
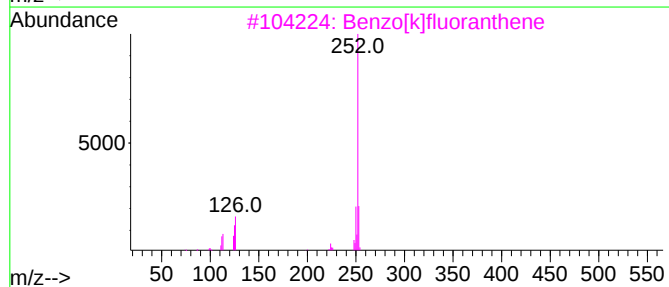
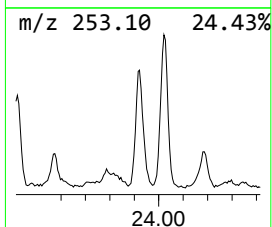
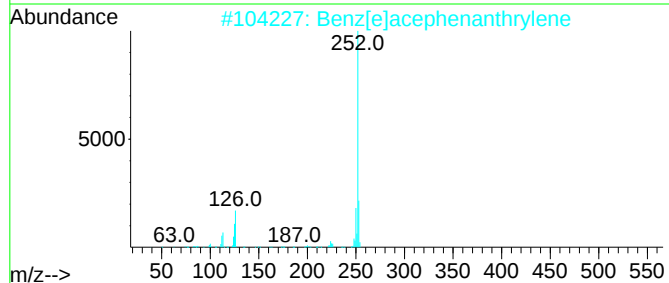
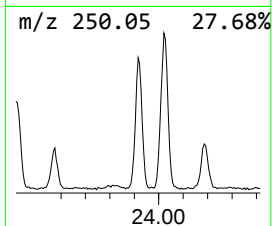
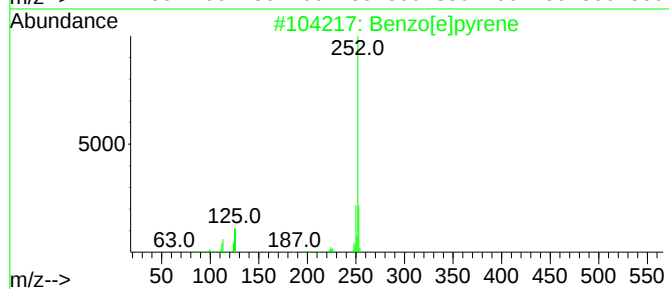
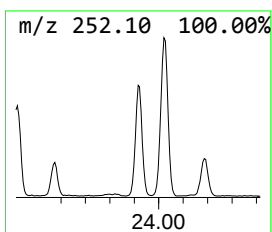
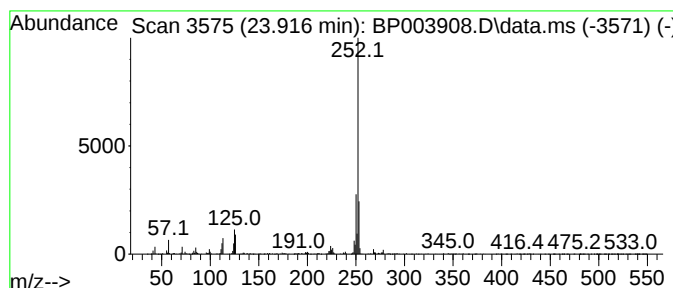
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Benzo[e]pyrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.916	9.75 ng	702053	Perylene-d12	24.133

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[e]pyrene	252	C20H12	000192-97-2	97
2		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	93
3		Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
4		Benzo[a]pyrene	252	C20H12	000050-32-8	93
5		Perylene	252	C20H12	000198-55-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111020\
Data File : BP003908.D
Acq On : 10 Nov 2020 18:26
Operator : CG/JU
Sample : L4680-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PAV-B7-110520-0-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown2.934	2.934	138.7	ng	7373880	1	8.281	1063680	20.0
9-Hexadecenoic ...	18.128	11.2	ng	808960	4	17.628	1451520	20.0
Hexadecanoic ac...	18.251	162.4	ng	11784100	4	17.628	1451520	20.0
9,12-Octadecadi...	19.339	291.6	ng	21166300	4	17.628	1451520	20.0
9-Octadecenoic ...	19.381	759.2	ng	55099100	4	17.628	1451520	20.0
Methyl stearate	19.486	84.0	ng	6099270	4	17.628	1451520	20.0
cis-13-Eicoseno...	20.422	15.4	ng	2142250	5	21.657	2774380	20.0
Eicosanoic acid...	20.528	6.9	ng	957033	5	21.657	2774380	20.0
unknown20.563	20.563	12.6	ng	1750860	5	21.657	2774380	20.0
unknown20.616	20.616	8.8	ng	1223200	5	21.657	2774380	20.0
12-Methyl-E,E-2...	20.816	4.3	ng	589652	5	21.657	2774380	20.0
Naphthalene, de...	21.110	4.3	ng	596529	5	21.657	2774380	20.0
Docosanoic acid...	21.439	4.3	ng	600249	5	21.657	2774380	20.0
Benzo[e]pyrene	23.916	9.8	ng	702053	6	24.133	1439470	20.0