

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP063020\
 Data File : BP002579.D
 Acq On : 30 Jun 2020 17:28
 Operator : CG/JU
 Sample : L3153-31
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SP2-H

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

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-BP062920.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.193	385	392	409	rBV	1427042	2285566	27.92%	4.401%
2	6.764	653	659	680	rVB	1742557	2846384	34.78%	5.481%
3	7.569	789	796	806	rBV	414978	649233	7.93%	1.250%
4	7.887	843	850	861	rBV	1428318	2306733	28.18%	4.442%
5	8.716	982	991	1012	rBV	1195051	2054536	25.10%	3.956%
6	10.340	1260	1267	1285	rBV	585374	1019668	12.46%	1.963%
7	12.081	1550	1563	1579	rBV5	28430	88212	1.08%	0.170%
8	12.834	1684	1691	1708	rBV	3183750	4887426	59.71%	9.411%
9	13.681	1830	1835	1848	rVB	336434	483962	5.91%	0.932%
10	14.222	1918	1927	1936	rVB	1024083	1664609	20.34%	3.205%
11	14.510	1967	1976	1981	rBV3	36336	84239	1.03%	0.162%
12	15.722	2174	2182	2194	rBV	2972877	4369061	53.38%	8.413%
13	16.981	2390	2396	2407	rVB	1451282	2070136	25.29%	3.986%
14	19.486	2817	2822	2828	rVB	546572	667455	8.15%	1.285%
15	19.575	2831	2837	2849	rVB	6138652	8185017	100.00%	15.761%
16	19.975	2902	2905	2910	rVB	126925	144240	1.76%	0.278%
17	20.404	2975	2978	2983	rVB	126712	143679	1.76%	0.277%
18	20.457	2984	2987	2991	rBV	67320	86407	1.06%	0.166%
19	20.657	3017	3021	3026	rBV	278772	323786	3.96%	0.623%
20	20.828	3046	3050	3061	rBV	891713	1198018	14.64%	2.307%
21	21.086	3089	3094	3099	rBV2	2085533	2621612	32.03%	5.048%
22	21.263	3121	3124	3134	rVB	164118	205422	2.51%	0.396%
23	21.704	3193	3199	3210	rBV2	727366	1289506	15.75%	2.483%
24	22.151	3271	3275	3290	rVB	282691	474868	5.80%	0.914%
25	22.375	3309	3313	3320	rVB	154372	209736	2.56%	0.404%
26	22.657	3355	3361	3364	rBV	1156927	1665708	20.35%	3.207%
27	22.692	3364	3367	3377	rVB	414141	681136	8.32%	1.312%
28	23.216	3451	3456	3461	rBV	277312	429351	5.25%	0.827%
29	23.280	3461	3467	3476	rVB2	1354874	2562703	31.31%	4.935%
30	23.522	3503	3508	3516	rVB	273296	461487	5.64%	0.889%
31	23.857	3559	3565	3573	rVB	1015625	1983149	24.23%	3.819%
32	23.933	3573	3578	3591	rVB	333693	699296	8.54%	1.347%
33	25.033	3759	3765	3775	rVB	241805	560343	6.85%	1.079%
34	25.463	3832	3838	3852	rVB2	292784	812181	9.92%	1.564%

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SP2-H

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

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35	25.592	3856	3860	3871	rvB2	133379	355945	4.35%	0.685%
36	26.310	3974	3982	3999	rvB2	421811	1361089	16.63%	2.621%

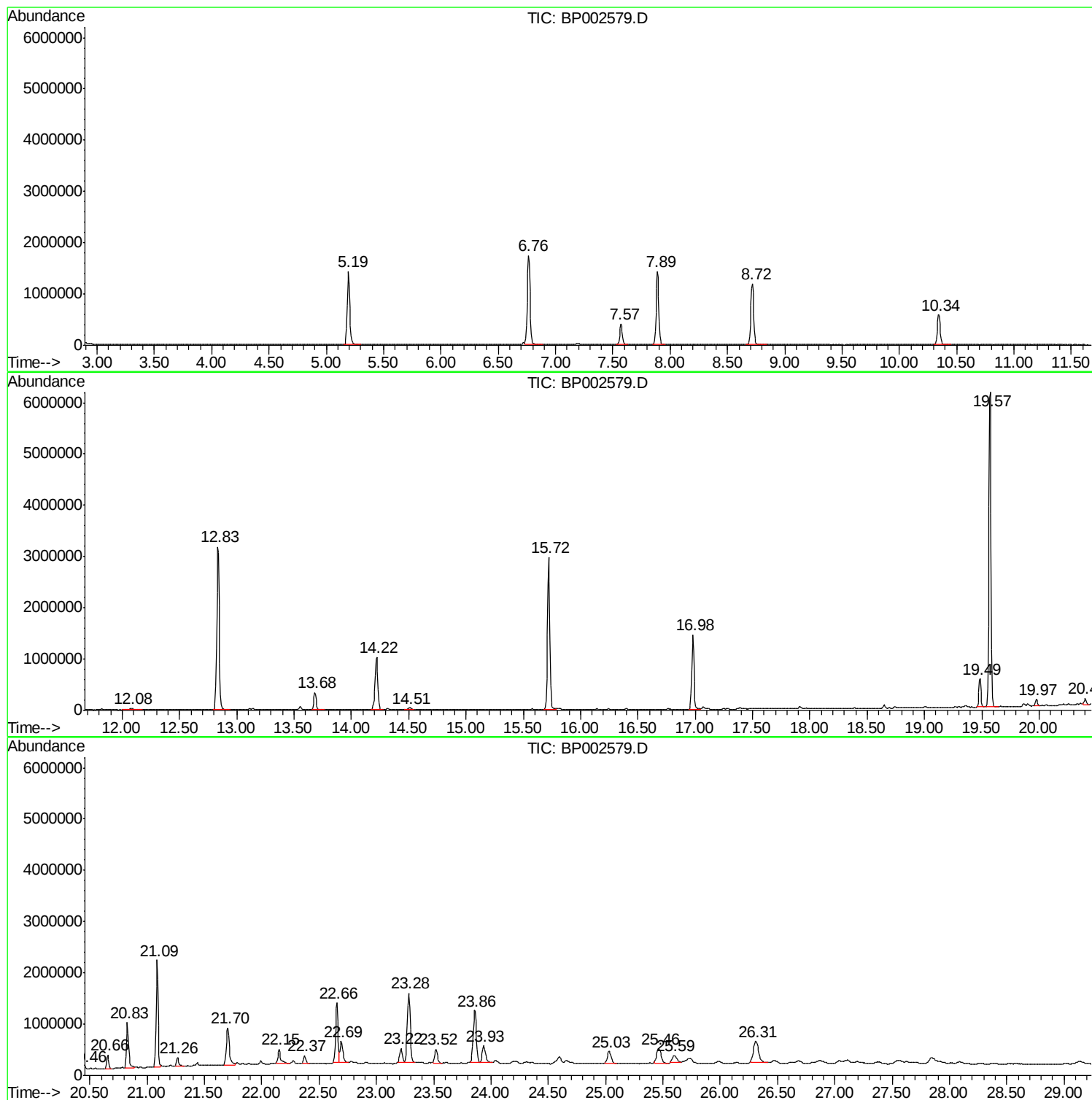
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BNA_P
ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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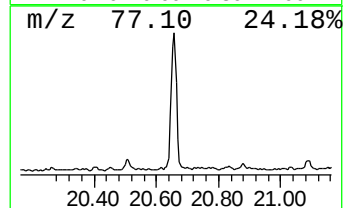
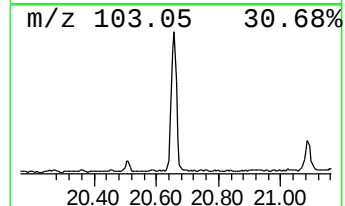
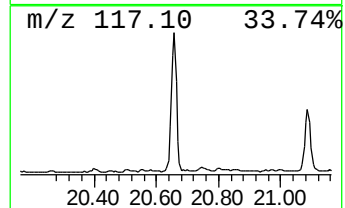
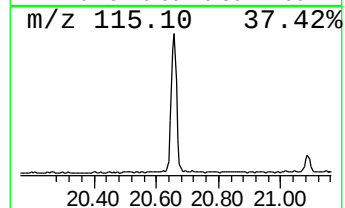
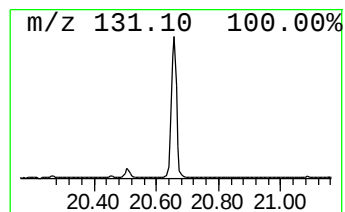
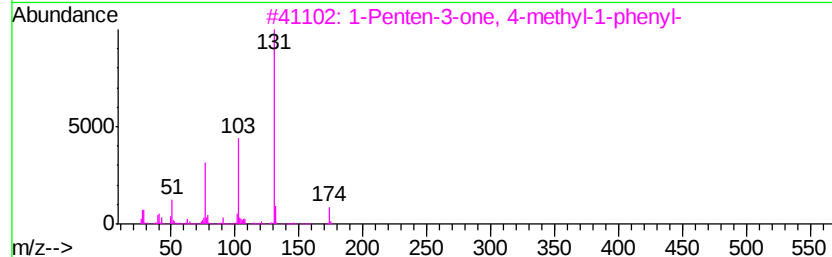
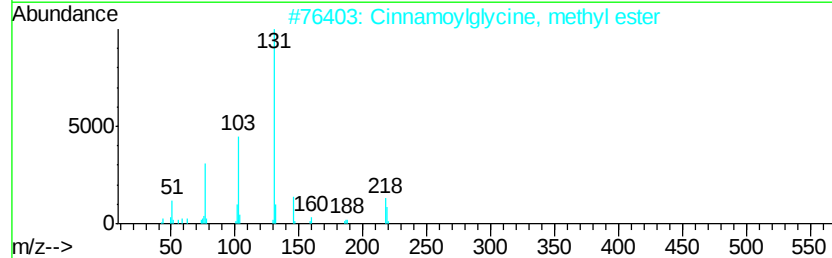
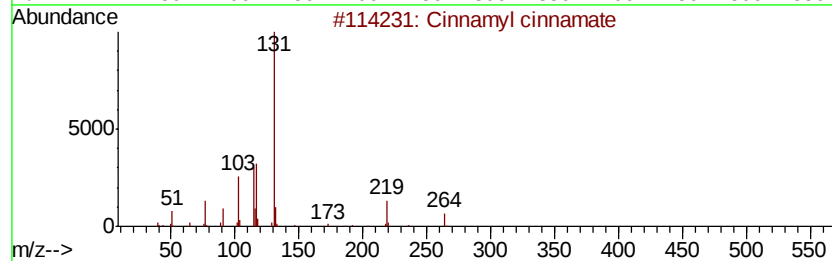
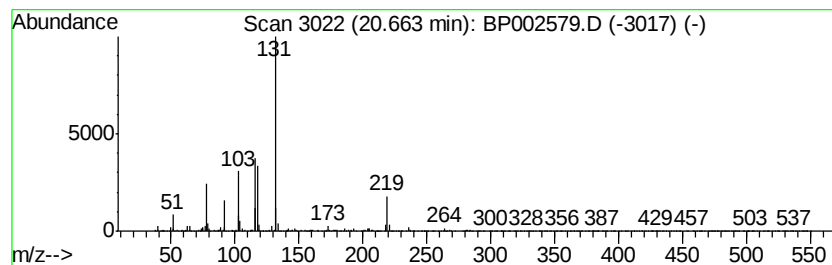
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Cinnamyl cinnamate Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.66	2.47 ng	323786	Chrysene-d12	21.09

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cinnamyl cinnamate	264	C18H16O2	000122-69-0	87	
2	Cinnamoylglycine, methyl ester	219	C12H13NO3	040778-04-9	53	
3	1-Penten-3-one, 4-methyl-1-phenyl-	174	C12H14O	003160-32-5	50	
4	2-Iodo-benzoic acid N'-(3-phenyl...	392	C16H13IN2O2	315672-62-9	50	
5	1-Penten-3-one, 4,4-dimethyl-1-p...	188	C13H16O	000538-44-3	50	



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ClientSampleId :
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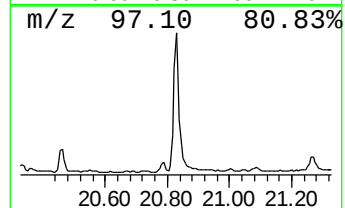
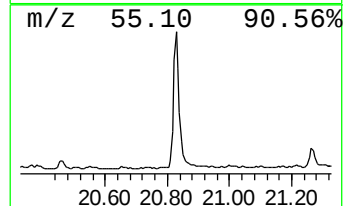
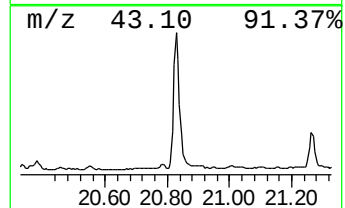
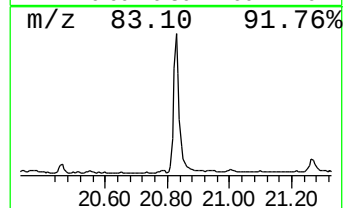
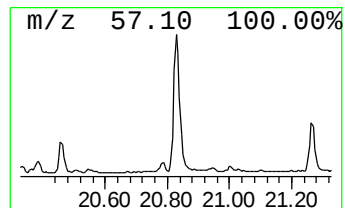
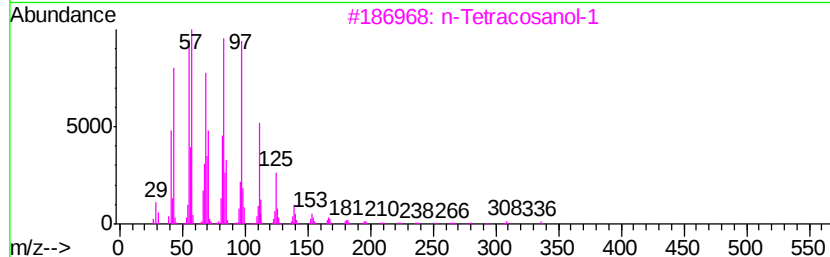
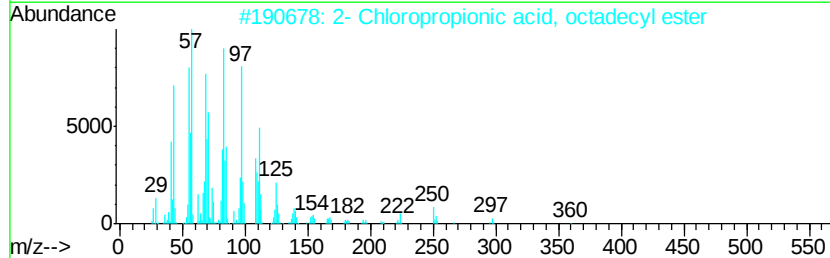
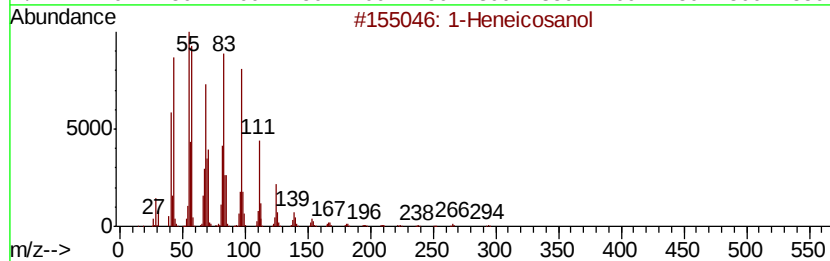
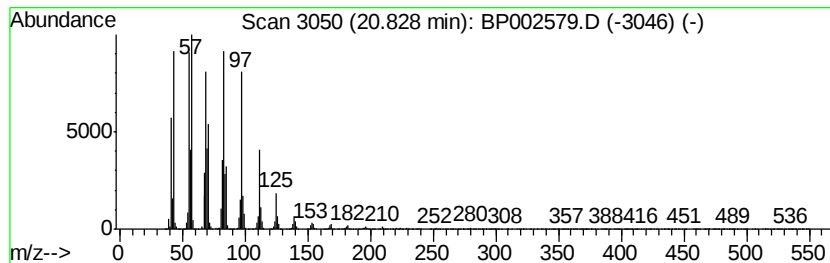
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1-Heneicosanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.83	9.14 ng	1198020	Chrysene-d12	21.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Heneicosanol	312	C21H44O	015594-90-8	95
2		2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94
3		n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4		Behenic alcohol	326	C22H46O	000661-19-8	94
5		1-Heptacosanol	396	C27H56O	002004-39-9	93



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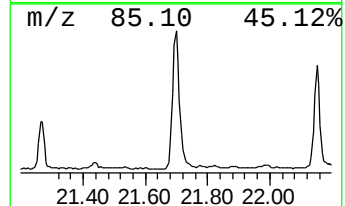
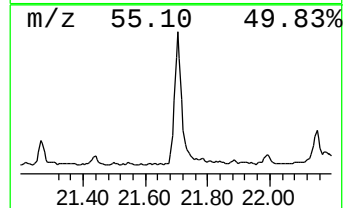
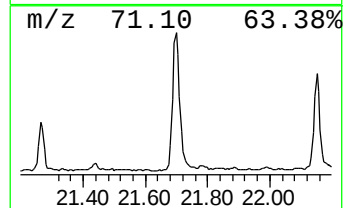
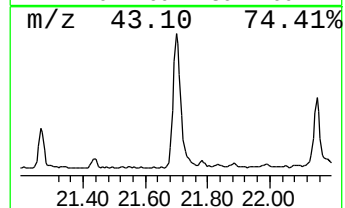
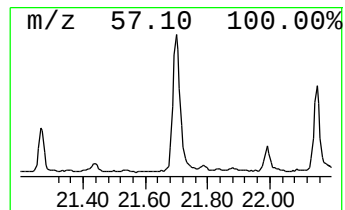
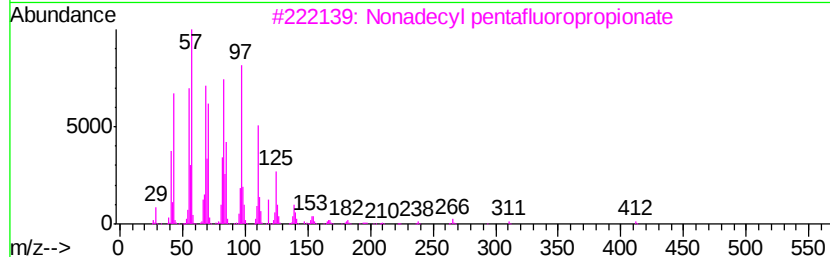
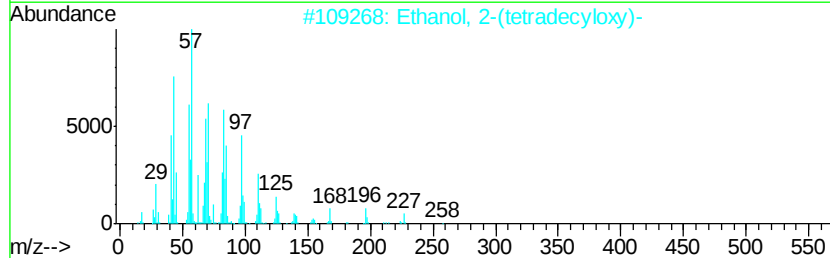
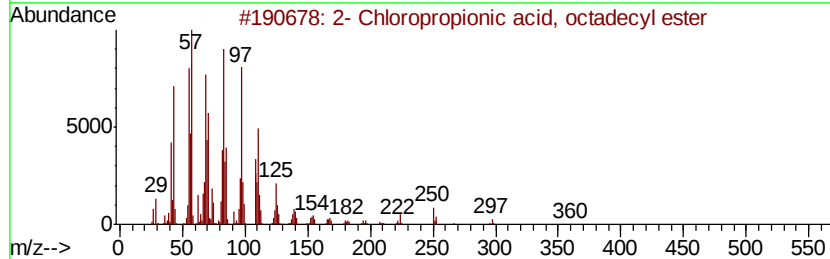
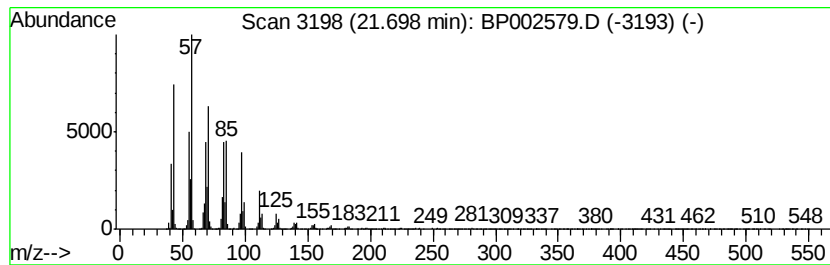
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2- Chloropropionic acid, oc... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.70	9.84 ng	1289510	Chrysene-d12	21.09

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	93
2		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
3		Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
4		Eicosyl heptafluorobutyrte	494	C24H41F7O2	1000351-83-5	91
5		Eicosyl pentafluoropropionate	444	C23H41F5O2	1000351-80-8	91



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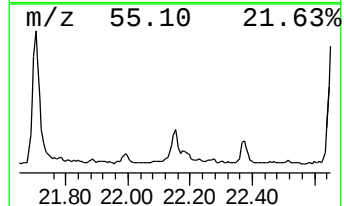
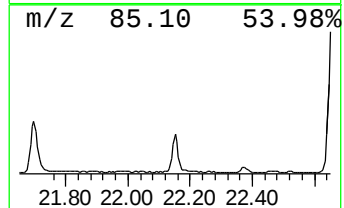
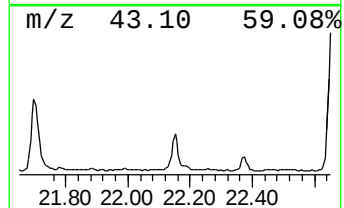
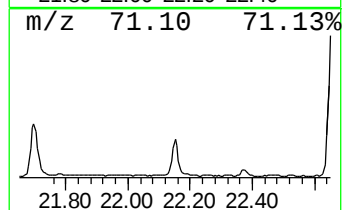
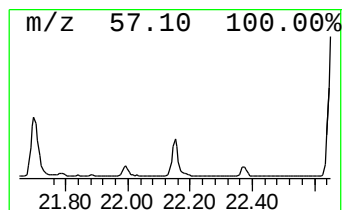
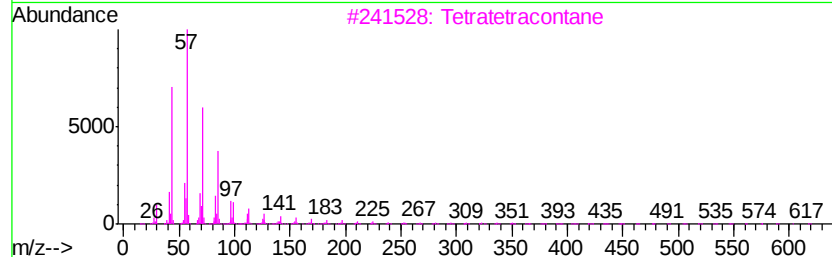
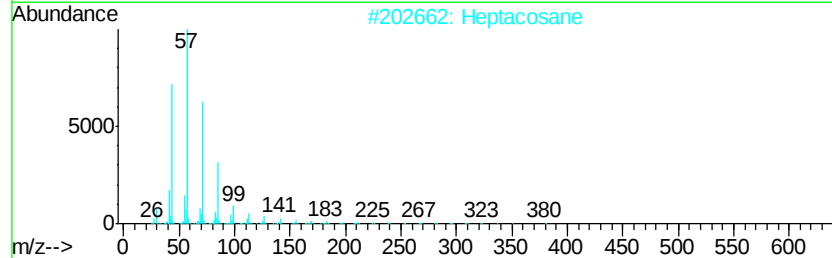
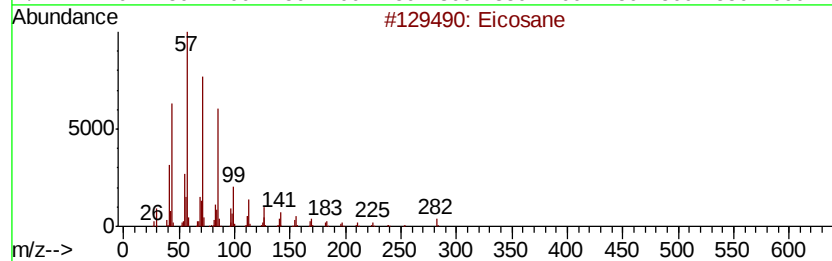
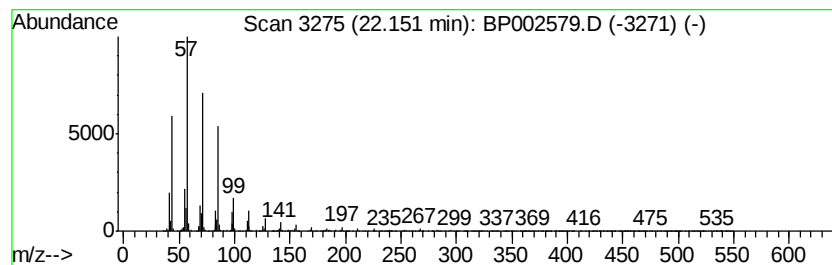
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Peak Number 6 Eicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.15	3.62 ng	474868	Chrysene-d12	21.09

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	94
2	Heptacosane		380	C27H56	000593-49-7	94
3	Tetratetracontane		619	C44H90	007098-22-8	87
4	Octadecane, 2-methyl-		268	C19H40	001560-88-9	81
5	triacontane		422	C30H62	000638-68-6	76



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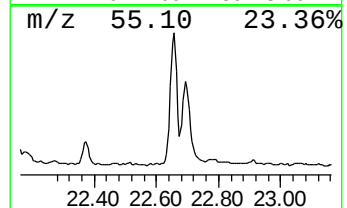
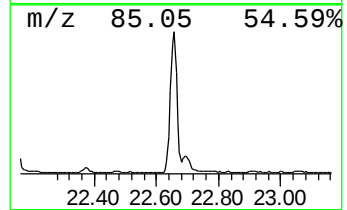
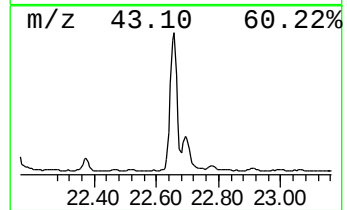
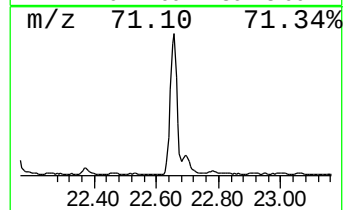
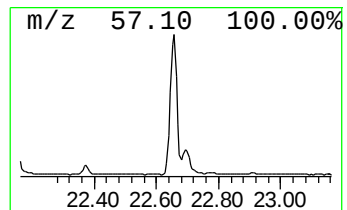
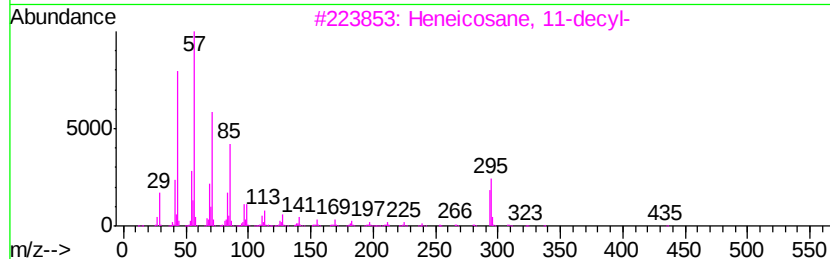
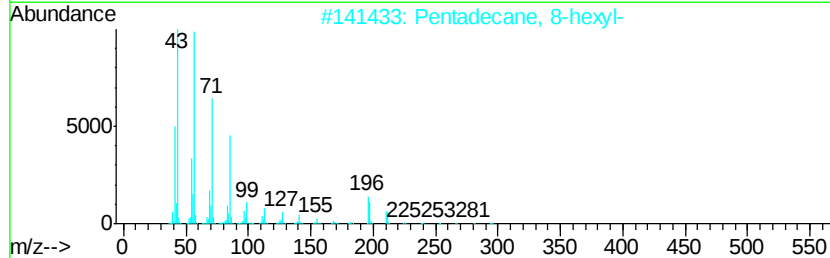
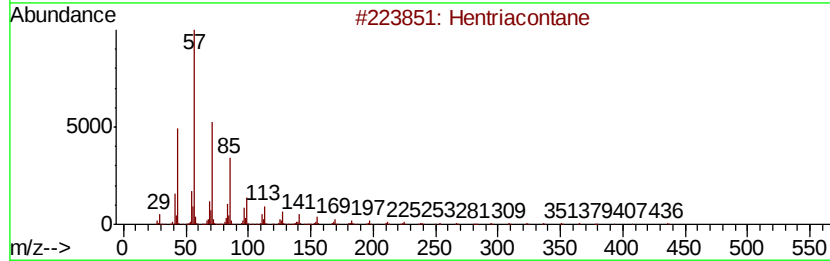
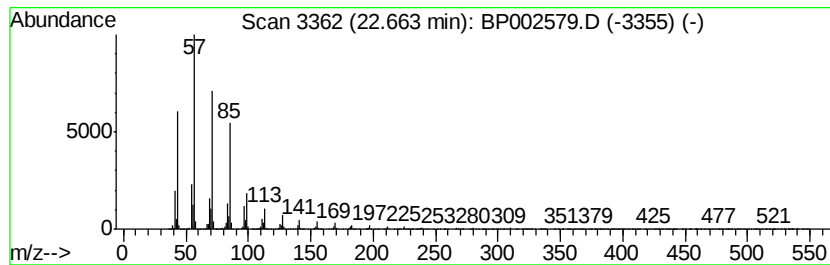
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Peak Number 7 Hentriacontane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.66	13.00 ng	1665710	Perylene-d12	23.28

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hentriacontane		437	C31H64	000630-04-6	96
2	Pentadecane, 8-hexyl-		296	C21H44	013475-75-7	94
3	Heneicosane, 11-decyl-		437	C31H64	055320-06-4	93
4	Eicosane, 10-methyl-		296	C21H44	054833-23-7	93
5	Nonacosane		408	C29H60	000630-03-5	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\
Data File : BP002579.D
Acq On : 30 Jun 2020 17:28
Operator : CG/JU
Sample : L3153-31
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SP2-H

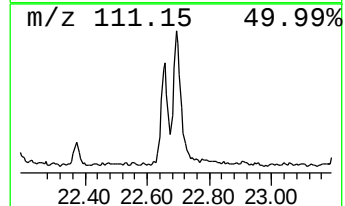
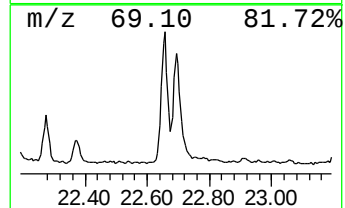
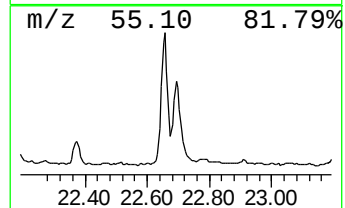
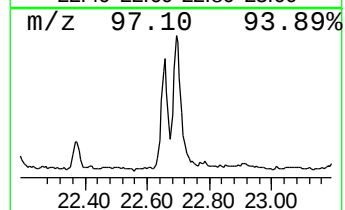
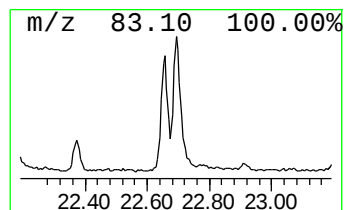
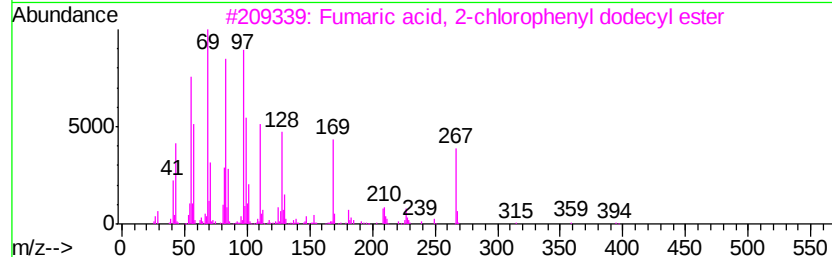
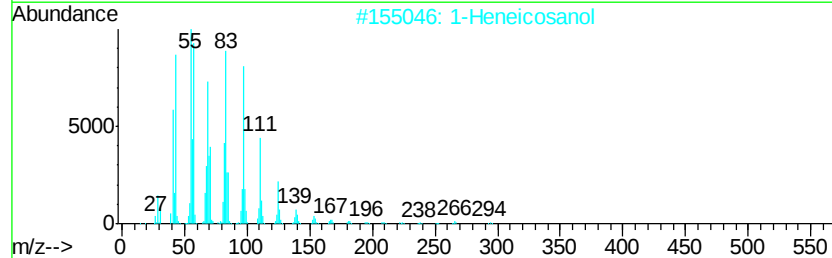
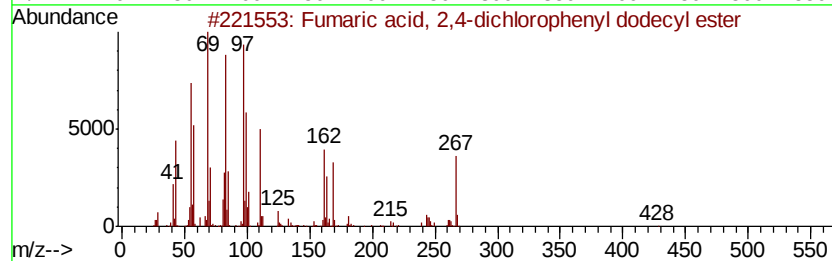
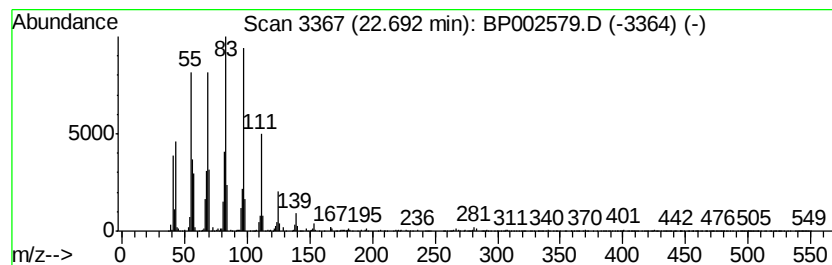
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Fumaric acid, 2,4-dichlorop... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.69	5.32 ng	681136	Perylene-d12	23.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Fumaric acid, 2,4-dichlorophenvl...	428	C22H30Cl2O4	1000345-34-4	94
2		1-Heneicosanol	312	C21H44O	015594-90-8	83
3		Fumaric acid, 2-chlorophenvl dod...	394	C22H31ClO4	1000344-83-4	76
4		Chloroacetic acid, tridecyl ester	276	C15H29ClO2	018277-85-5	72
5		Fumaric acid, 2-methylallyl tetr...	366	C22H38O4	1000330-55-3	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\
Data File : BP002579.D
Acq On : 30 Jun 2020 17:28
Operator : CG/JU
Sample : L3153-31
Misc :
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Instrument :
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ClientSampleId :
SP2-H

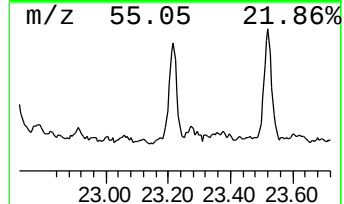
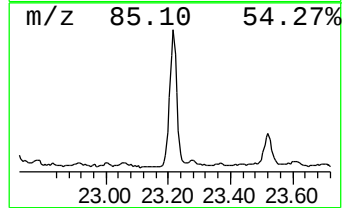
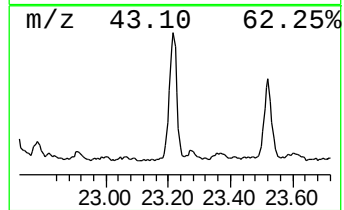
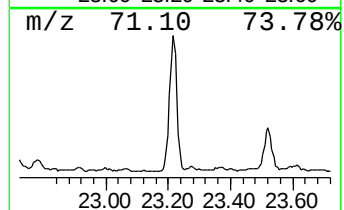
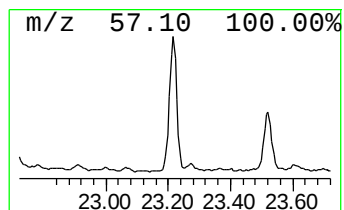
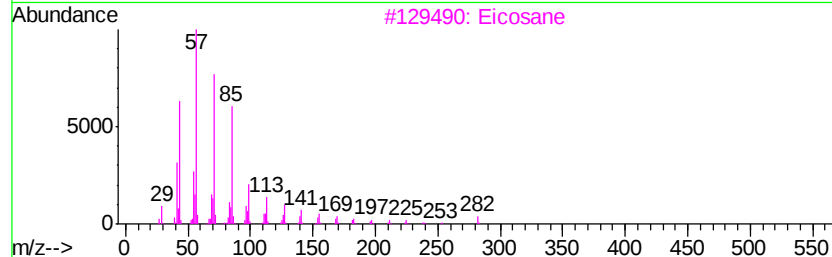
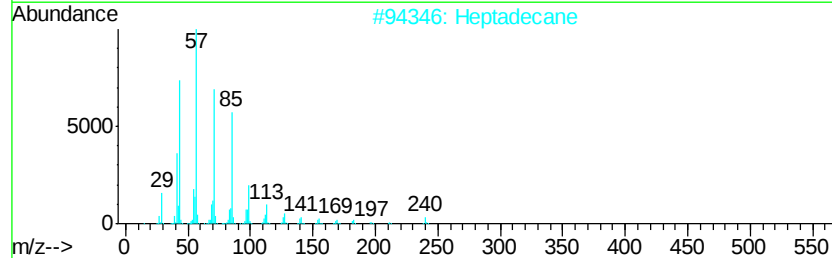
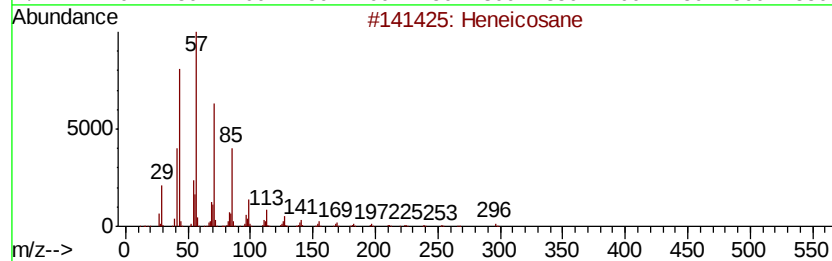
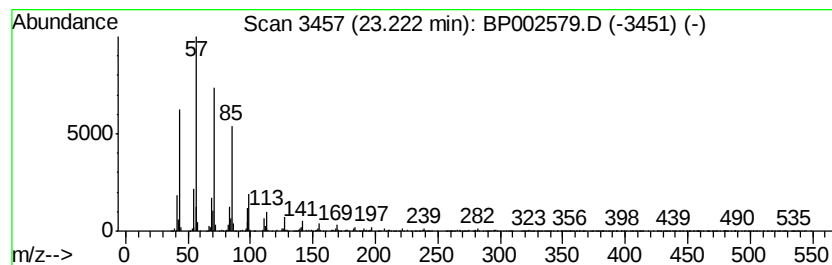
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heneicosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.22	3.35 ng	429351	Perylene-d12	23.28

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	90
2	Heptadecane		240	C17H36	000629-78-7	90
3	Eicosane		282	C20H42	000112-95-8	87
4	Tetratetracontane		619	C44H90	007098-22-8	87
5	Octadecane		254	C18H38	000593-45-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\
Data File : BP002579.D
Acq On : 30 Jun 2020 17:28
Operator : CG/JU
Sample : L3153-31
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SP2-H

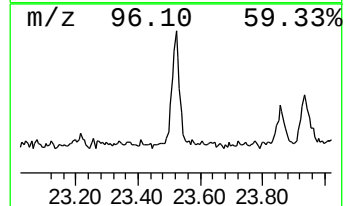
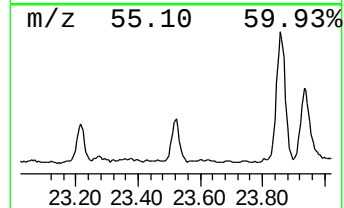
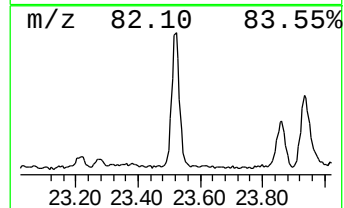
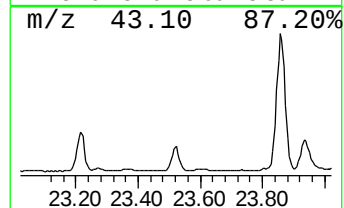
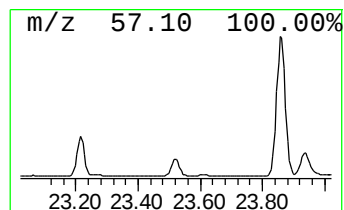
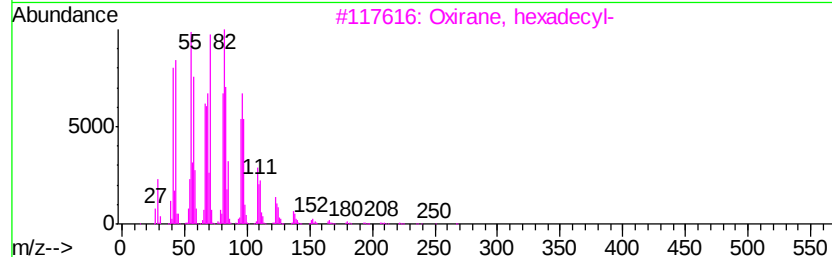
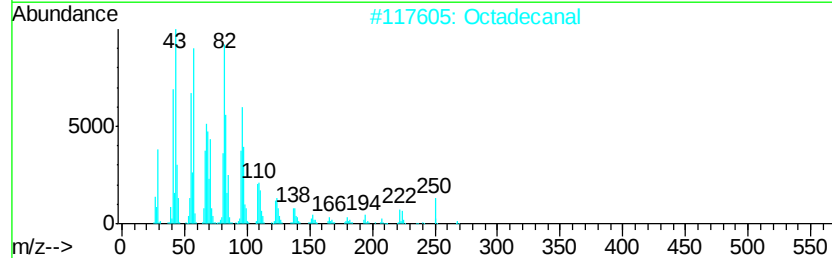
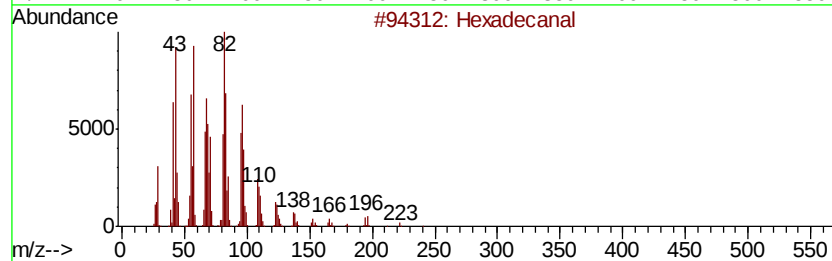
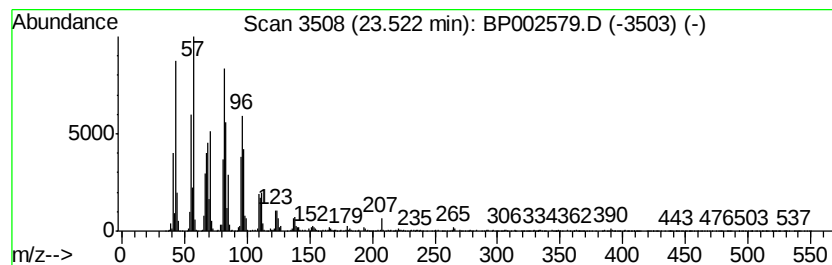
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexadecanal Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.52	3.60 ng	461487	Perylene-d12	23.28

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanal	240	C16H32O	000629-80-1	93
2			Octadecanal	268	C18H36O	000638-66-4	91
3			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
4			Tetracosanal	352	C24H48O	057866-08-7	91
5			1,19-Eicosadiene	278	C20H38	014811-95-1	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\
Data File : BP002579.D
Acq On : 30 Jun 2020 17:28
Operator : CG/JU
Sample : L3153-31
Misc :
ALS Vial : 9 Sample Multiplier: 1

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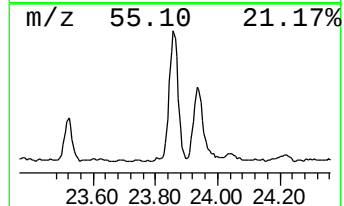
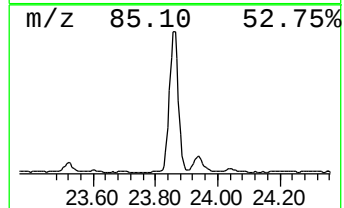
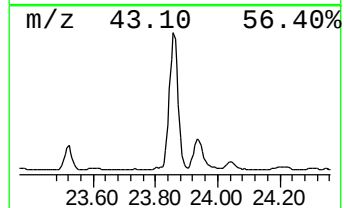
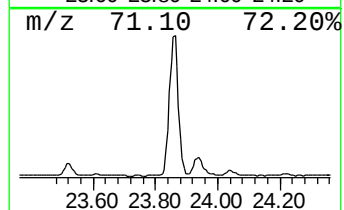
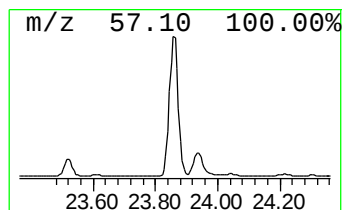
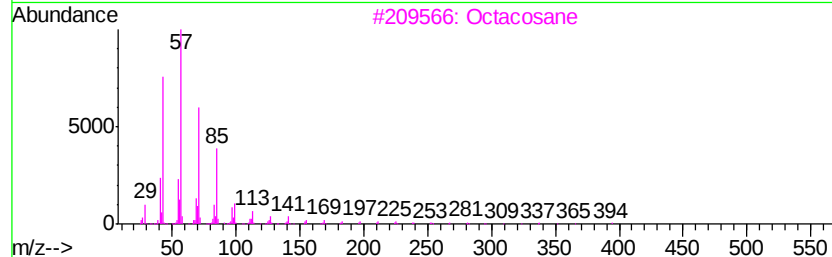
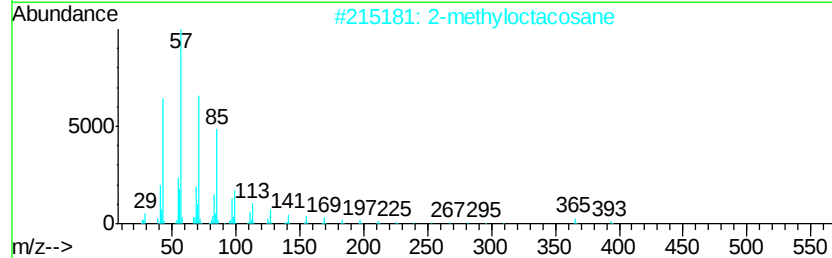
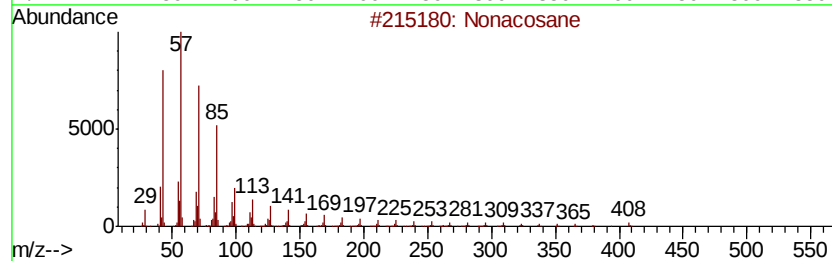
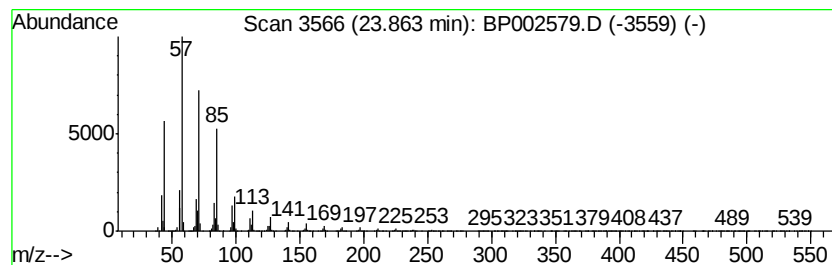
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Nonacosane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.86	15.48 ng	1983150	Perylene-d12	23.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonacosane	408	C29H60	000630-03-5	95
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Tetratetracontane	619	C44H90	007098-22-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\
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Acq On : 30 Jun 2020 17:28
Operator : CG/JU
Sample : L3153-31
Misc :
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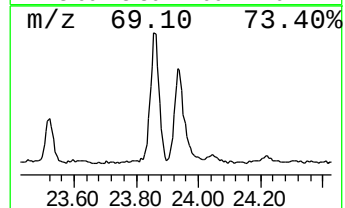
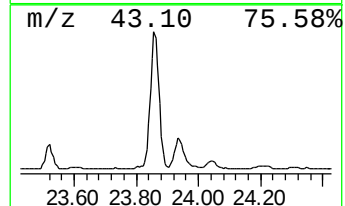
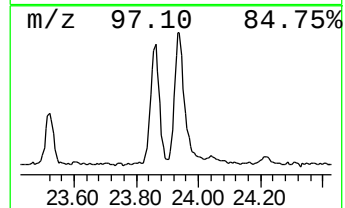
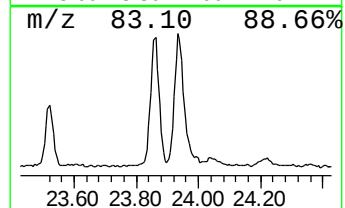
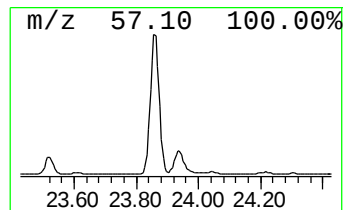
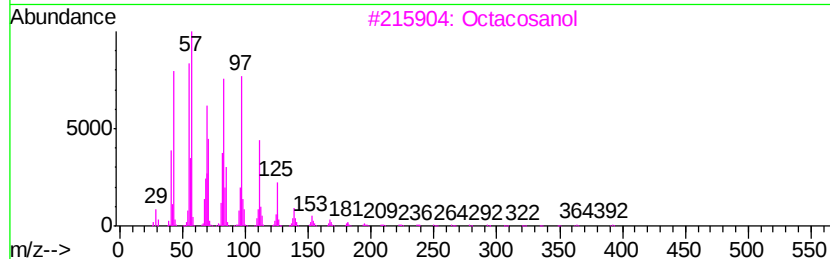
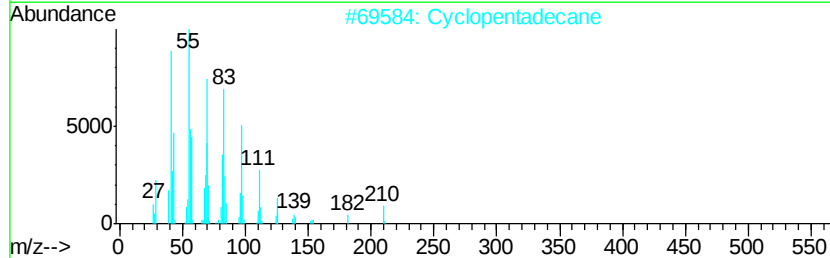
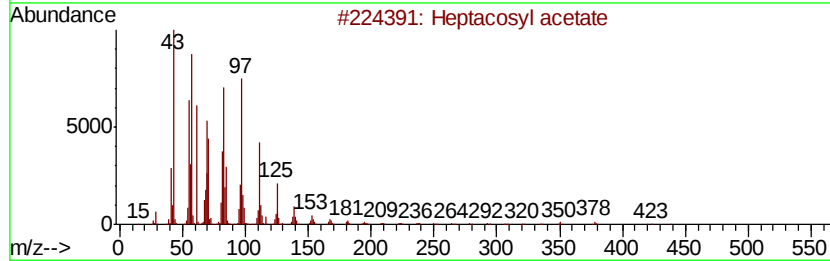
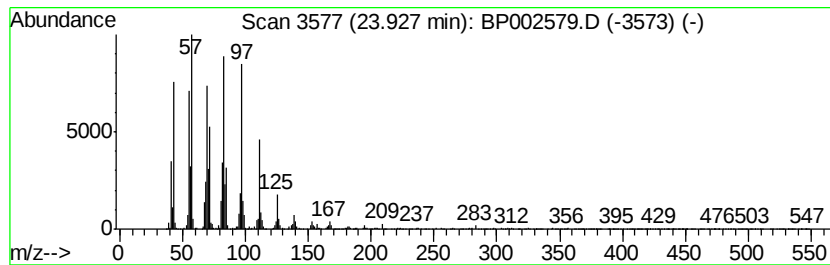
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heptacosyl acetate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.93	5.46 ng	699296	Perylene-d12	23.28	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosyl acetate	438	C29H58O2	1000351-78-2	99
2	Cyclopentadecane	210	C15H30	000295-48-7	93
3	Octacosanol	410	C28H58O	000557-61-9	93
4	1-Heptacosanol	396	C27H56O	002004-39-9	93
5	n-Tetracosanol-1	354	C24H50O	000506-51-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\
Data File : BP002579.D
Acq On : 30 Jun 2020 17:28
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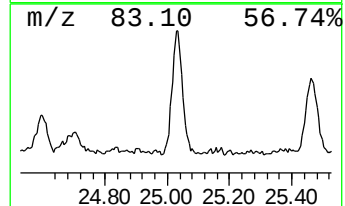
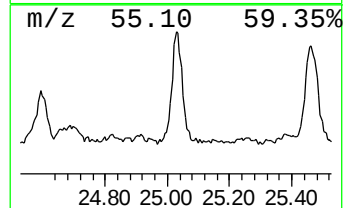
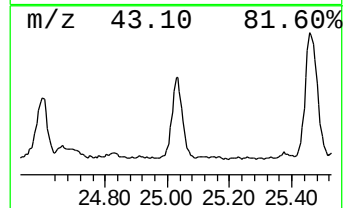
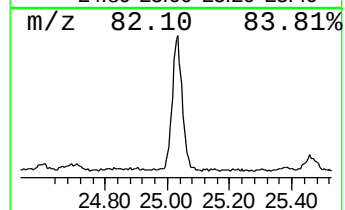
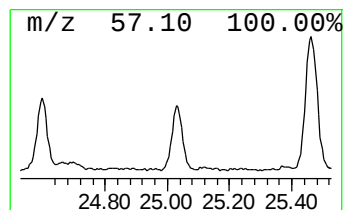
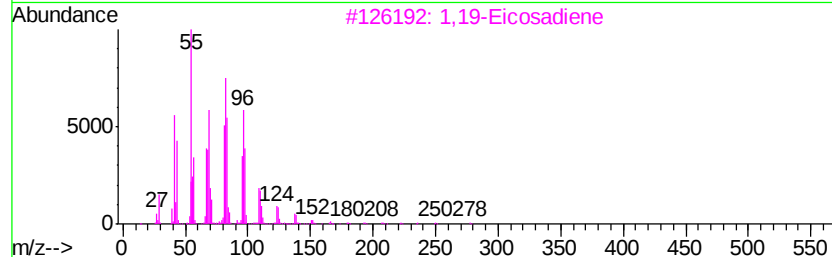
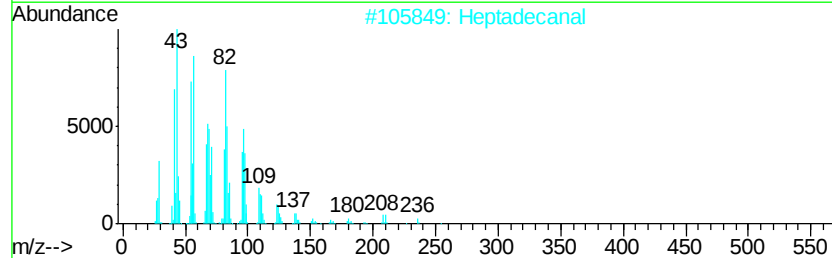
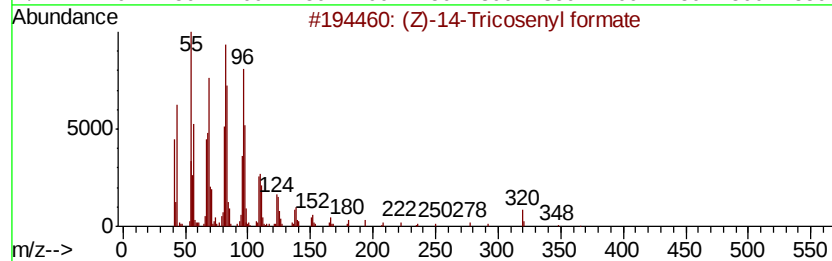
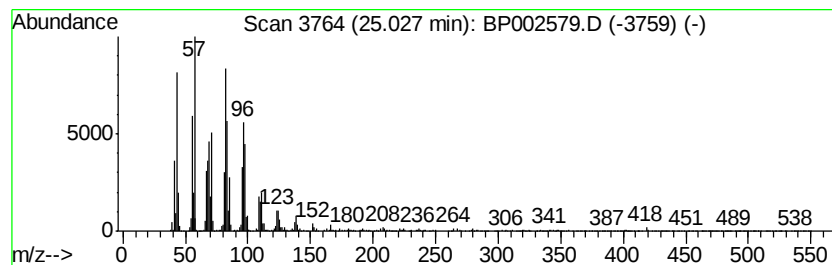
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 (Z)-14-Tricosenyl formate Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.03	4.37 ng	560343	Perylene-d12	23.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(Z)-14-Tricosenyl formate	366	C24H46O2	077899-10-6	92
2		Heptadecanal	254	C17H34O	1000376-70-0	91
3		1,19-Eicosadiene	278	C20H38	014811-95-1	90
4		Oxirane, heptadecyl-	282	C19H38O	067860-04-2	87
5		Octadecanal	268	C18H36O	000638-66-4	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\
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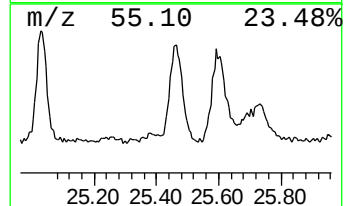
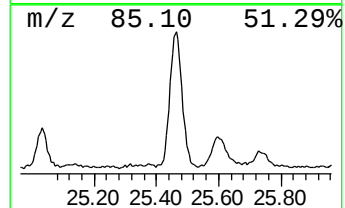
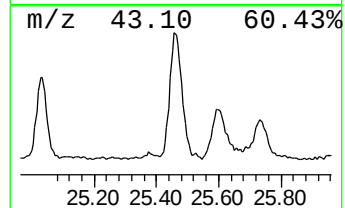
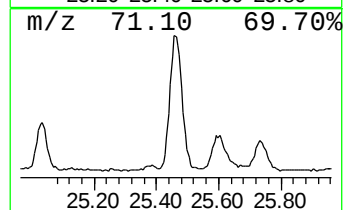
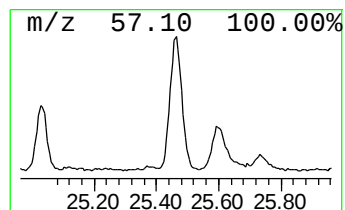
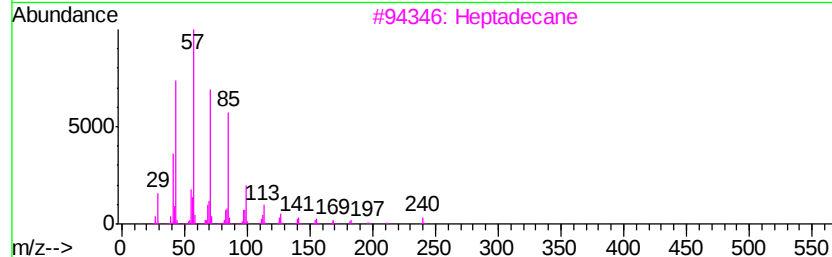
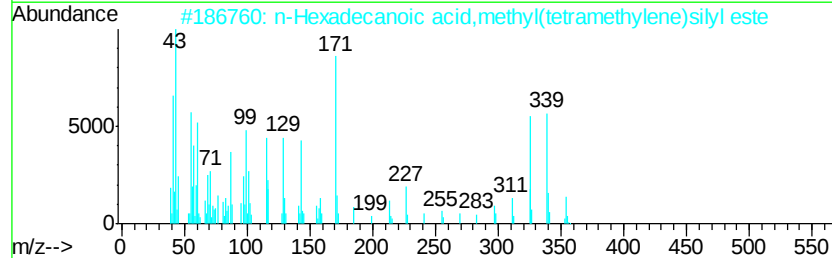
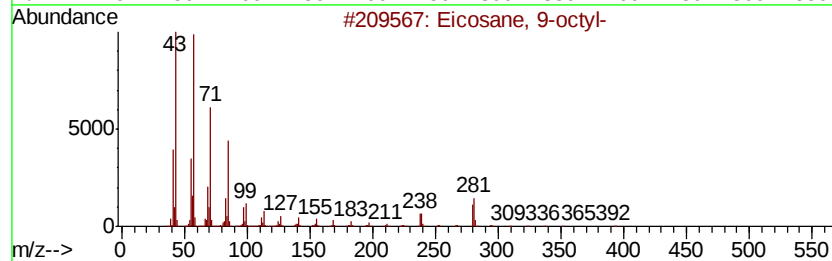
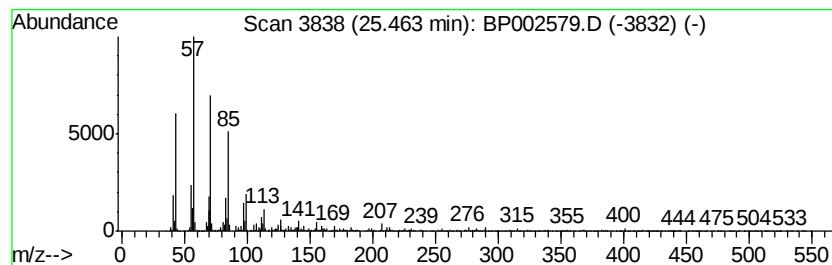
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Eicosane, 9-octyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.46	6.34 ng	812181	Perylene-d12	23.28
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane, 9-octyl-	394 C28H58	013475-77-9	91
2	n-Hexadecanoic acid,methyl(tetra...	354 C21H42O2Si	1000217-04-1	90
3	Heptadecane	240 C17H36	000629-78-7	87
4	Hexacosane	366 C26H54	000630-01-3	87
5	Tetracosane	338 C24H50	000646-31-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\
Data File : BP002579.D
Acq On : 30 Jun 2020 17:28
Operator : CG/JU
Sample : L3153-31
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SP2-H

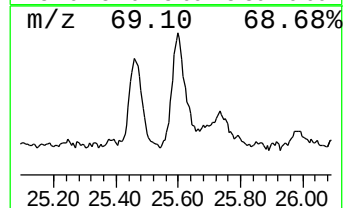
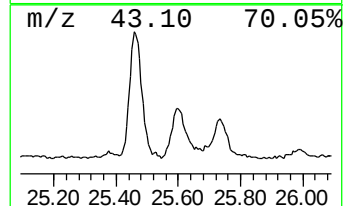
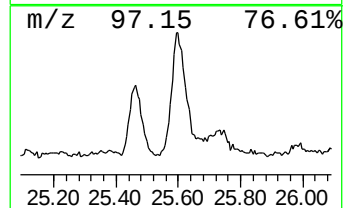
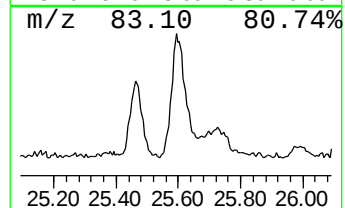
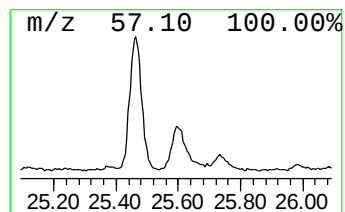
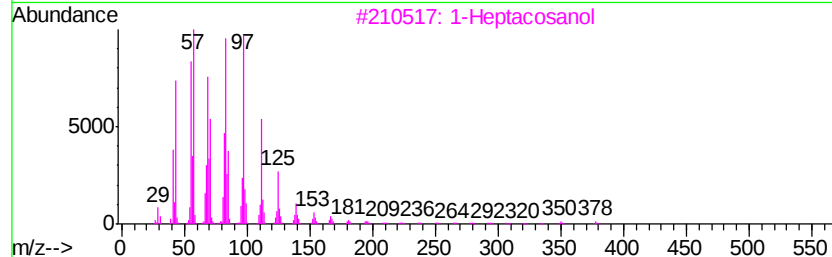
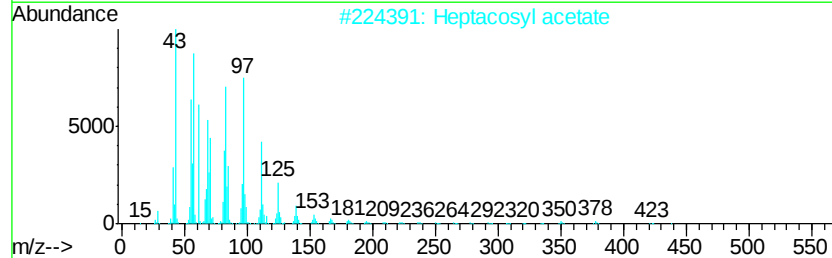
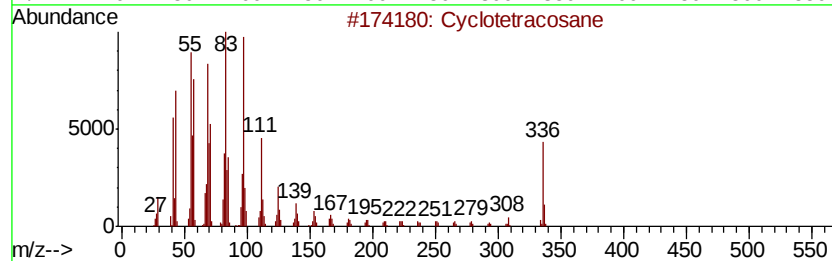
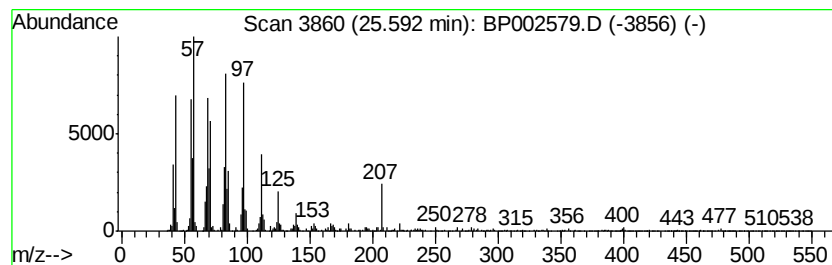
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Cyclotetracosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.59	2.78 ng	355945	Perylene-d12	23.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclotetracosane	336	C24H48	000297-03-0	96
2		Heptacosyl acetate	438	C29H58O2	1000351-78-2	93
3		1-Heptacosanol	396	C27H56O	002004-39-9	91
4		1-Triacontanol	438	C30H62O	000593-50-0	90
5		Octacosanol	410	C28H58O	000557-61-9	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\
Data File : BP002579.D
Acq On : 30 Jun 2020 17:28
Operator : CG/JU
Sample : L3153-31
Misc :
ALS Vial : 9 Sample Multiplier: 1

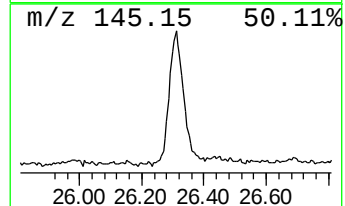
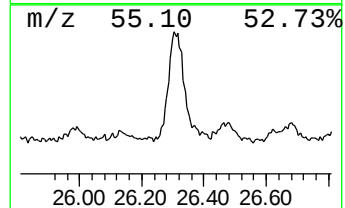
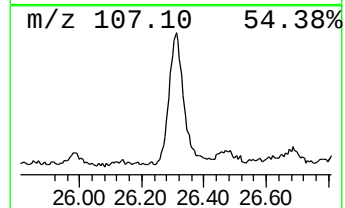
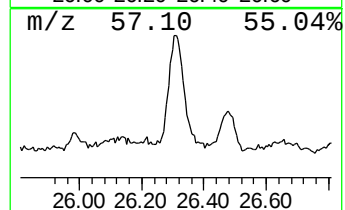
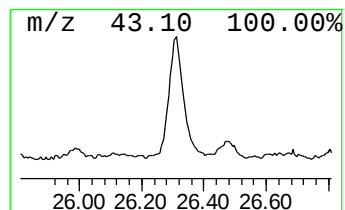
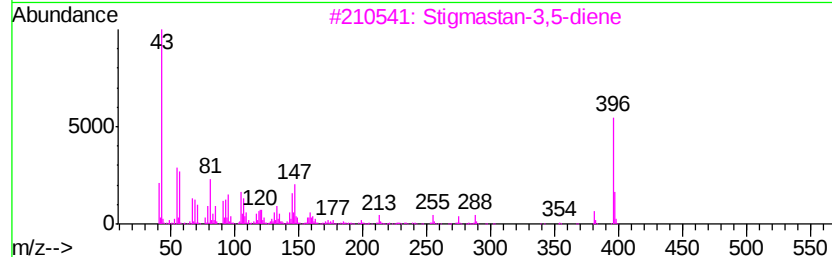
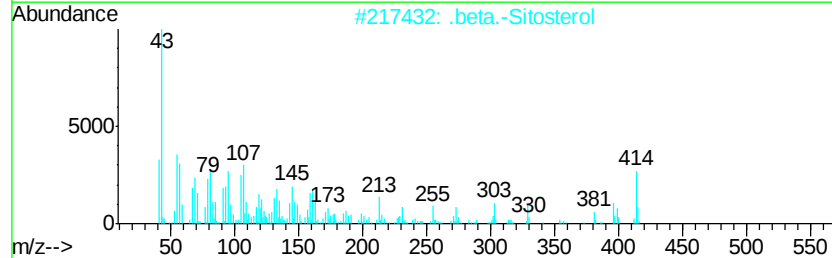
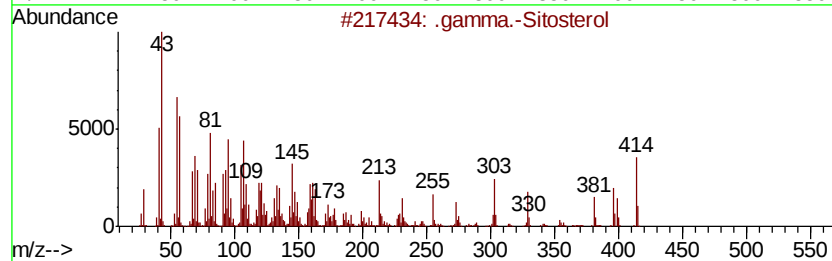
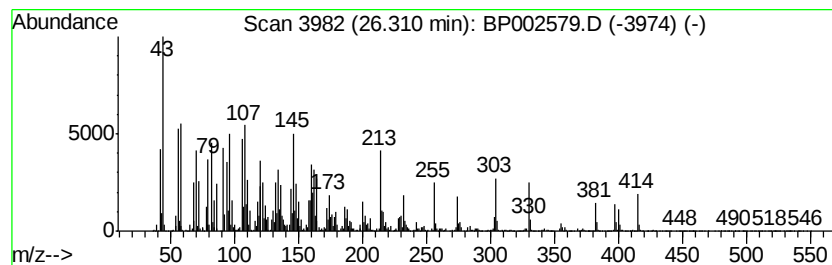
Instrument :
BNA_P
ClientSampleId :
SP2-H

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 .gamma.-Sitosterol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
26.31	10.62 ng	1361090	Perylene-d12	23.28	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.gamma.-Sitosterol	414	C29H50O	000083-47-6	96
2	.beta.-Sitosterol	414	C29H50O	000083-46-5	89
3	Stigmastan-3,5-diene	396	C29H48	1000214-16-4	25
4	Pregn-5-en-3-ol, 21-bromo-20-met...	394	C22H35BrO	055103-80-5	18
5	17-(1,5-Dimethylhexyl)-10,13-dim...	414	C29H50O	1000210-86-9	16



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP063020\
 Data File : BP002579.D
 Acq On : 30 Jun 2020 17:28
 Operator : CG/JU
 Sample : L3153-31
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SP2-H

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP062920.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
Cinnamyl cinnamate	20.66	2.5 ng		323786	5	21.09	2621610	20.0
1-Heneicosanol	20.83	9.1 ng		1198020	5	21.09	2621610	20.0
2-Chloropropioni...	21.70	9.8 ng		1289510	5	21.09	2621610	20.0
Eicosane	22.15	3.6 ng		474868	5	21.09	2621610	20.0
Hentriacontane	22.66	13.0 ng		1665710	6	23.28	2562700	20.0
Fumaric acid, 2,4...	22.69	5.3 ng		681136	6	23.28	2562700	20.0
Heneicosane	23.22	3.4 ng		429351	6	23.28	2562700	20.0
Hexadecanal	23.52	3.6 ng		461487	6	23.28	2562700	20.0
Nonacosane	23.86	15.5 ng		1983150	6	23.28	2562700	20.0
Heptacosyl acetate	23.93	5.5 ng		699296	6	23.28	2562700	20.0
(Z)-14-Tricosenyl...	25.03	4.4 ng		560343	6	23.28	2562700	20.0
Eicosane, 9-octyl-	25.46	6.3 ng		812181	6	23.28	2562700	20.0
Cyclotetracosane	25.59	2.8 ng		355945	6	23.28	2562700	20.0
.gamma.-Sitosterol	26.31	10.6 ng		1361090	6	23.28	2562700	20.0