

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP060120\
 Data File : BP002337.D
 Acq On : 01 Jun 2020 14:02
 Operator : CG/JU
 Sample : L2773-17
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 NM33-COMP-2020-0-6

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-BP052720.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	4.558	274	284	301	rBV	140705	710583	4.80%	0.508%
2	5.228	391	398	407	rBV	5227651	8036822	54.28%	5.751%
3	6.799	653	665	680	rBV	5438444	8902840	60.13%	6.370%
4	7.605	795	802	812	rBV	1549107	2445692	16.52%	1.750%
5	7.928	849	857	865	rBV	4729105	7713627	52.10%	5.519%
6	8.752	986	997	1012	rBV	3744800	6302811	42.57%	4.510%
7	10.381	1266	1274	1281	rBV	2035400	3428013	23.15%	2.453%
8	12.875	1690	1698	1716	rBV	7601350	12571657	84.91%	8.996%
9	13.716	1834	1841	1850	rBV	730336	1087300	7.34%	0.778%
10	14.251	1926	1932	1939	rBV2	2940298	4490419	30.33%	3.213%
11	15.751	2176	2187	2199	rBV	5889913	8877560	59.96%	6.352%
12	16.404	2288	2298	2303	rBV2	400730	670644	4.53%	0.480%
13	16.945	2385	2390	2395	rBV	143131	239722	1.62%	0.172%
14	17.010	2395	2401	2405	rVV	3560990	5223085	35.28%	3.737%
15	17.051	2405	2408	2413	rVV	822726	1105904	7.47%	0.791%
16	17.104	2413	2417	2420	rVV3	119566	183875	1.24%	0.132%
17	17.139	2420	2423	2429	rVB	280001	380719	2.57%	0.272%
18	17.410	2465	2469	2476	rVB	219784	318927	2.15%	0.228%
19	17.822	2534	2539	2544	rBV2	138101	246289	1.66%	0.176%
20	17.886	2546	2550	2554	rVV2	179959	265160	1.79%	0.190%
21	17.939	2554	2559	2567	rVV2	726921	1093198	7.38%	0.782%
22	18.075	2577	2582	2586	rBV2	167096	274758	1.86%	0.197%
23	18.233	2605	2609	2614	rBV2	122562	190055	1.28%	0.136%
24	18.416	2636	2640	2647	rVB4	121346	209373	1.41%	0.150%
25	18.780	2698	2702	2707	rBV2	104742	161360	1.09%	0.115%
26	18.904	2719	2723	2727	rBV2	334532	465841	3.15%	0.333%
27	19.004	2733	2740	2743	rBV	3313815	4132722	27.91%	2.957%
28	19.033	2743	2745	2750	rVB	1637853	1823578	12.32%	1.305%
29	19.386	2797	2805	2809	rBV	1525248	2302026	15.55%	1.647%
30	19.598	2833	2841	2853	rVB	10297914	14806146	100.00%	10.594%
31	19.875	2885	2888	2890	rBV2	142074	184146	1.24%	0.132%
32	20.027	2911	2914	2918	rVB2	143968	188204	1.27%	0.135%
33	20.245	2947	2951	2954	rBV2	317005	382823	2.59%	0.274%
34	20.280	2954	2957	2962	rVB	779324	885416	5.98%	0.634%

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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_P\METHODS\8270-BP052720.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	20.580	3005	3008	3010	rBV2	150459	173058	1.17%	0.124%
36	20.857	3051	3055	3066	rVB3	1577247	2190364	14.79%	1.567%
37	21.116	3092	3099	3103	rBV2	4106927	6475930	43.74%	4.634%
38	21.151	3103	3105	3111	rVB	709359	767134	5.18%	0.549%
39	21.469	3157	3159	3163	rVB2	151726	158636	1.07%	0.114%
40	21.739	3201	3205	3214	rVB2	1003235	1497539	10.11%	1.072%
41	22.416	3317	3320	3326	rVB	434051	564395	3.81%	0.404%
42	22.663	3357	3362	3365	rBV	673480	1287977	8.70%	0.922%
43	22.704	3365	3369	3372	rVV2	2005731	3104147	20.97%	2.221%
44	22.739	3372	3375	3386	rVB	1307380	2082415	14.06%	1.490%
45	23.133	3438	3442	3448	rVB	432717	757881	5.12%	0.542%
46	23.227	3453	3458	3462	rBV	539639	965620	6.52%	0.691%
47	23.327	3469	3475	3481	rBV	2568684	4507116	30.44%	3.225%
48	23.580	3513	3518	3525	rVB	895661	1529890	10.33%	1.095%
49	23.927	3571	3577	3583	rBV	2798442	5330994	36.01%	3.815%
50	23.998	3584	3589	3602	rVB	1299081	2576776	17.40%	1.844%
51	25.551	3847	3853	3864	rVB2	725162	1984538	13.40%	1.420%
52	25.686	3870	3876	3884	rBV	701151	1667256	11.26%	1.193%
53	26.392	3991	3996	4013	rVB4	611638	1830998	12.37%	1.310%

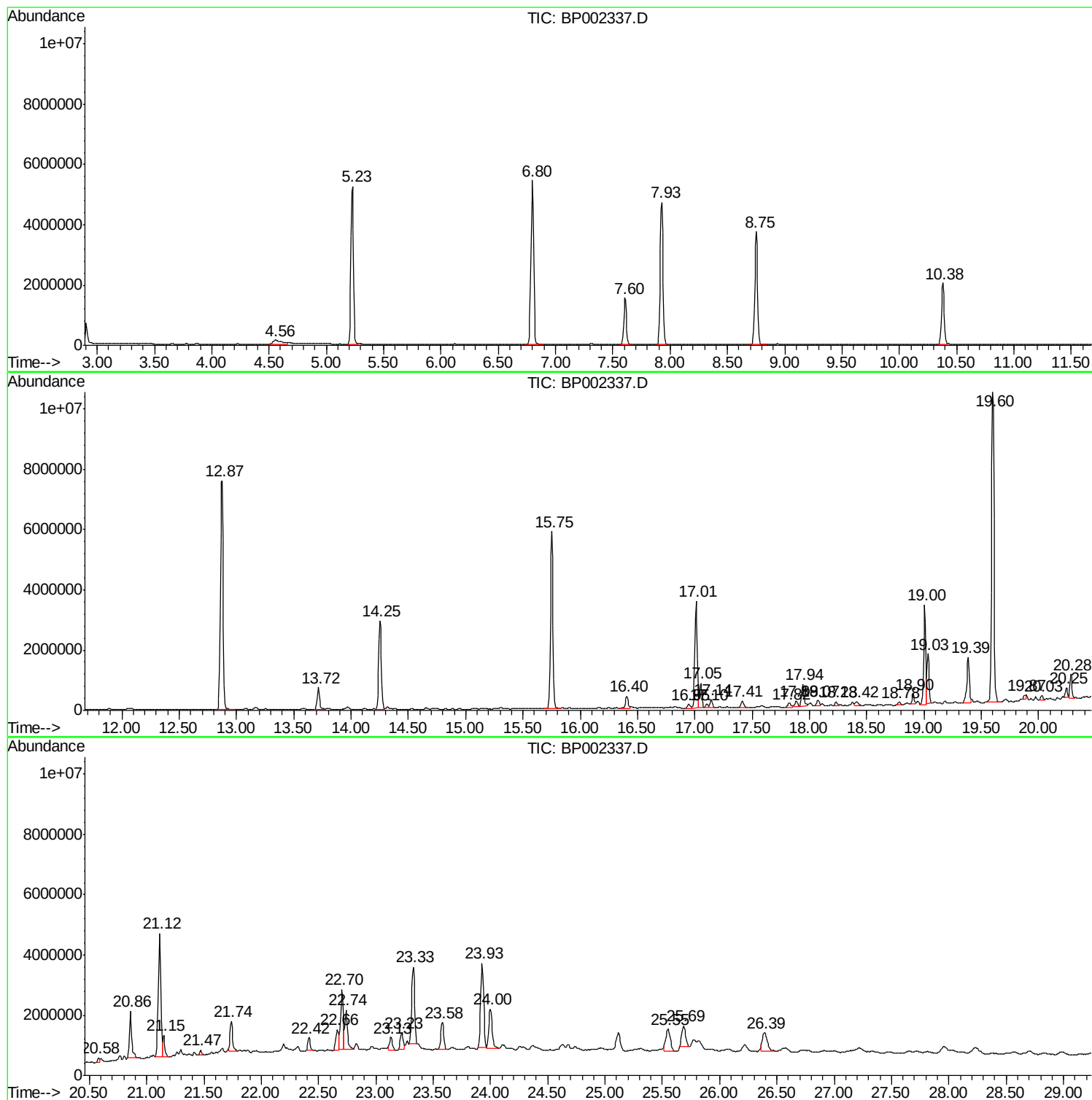
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP052720.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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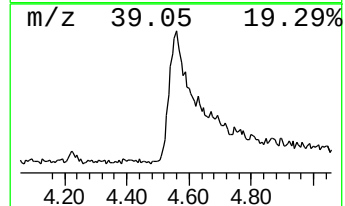
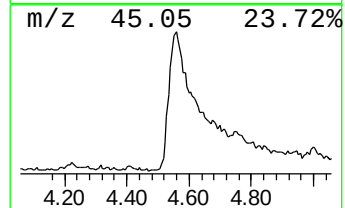
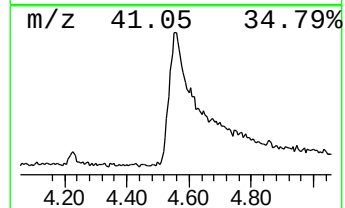
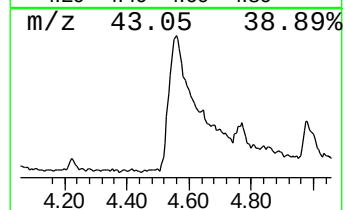
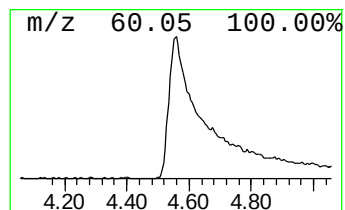
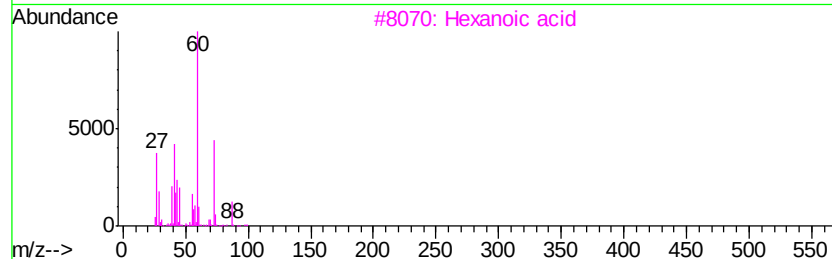
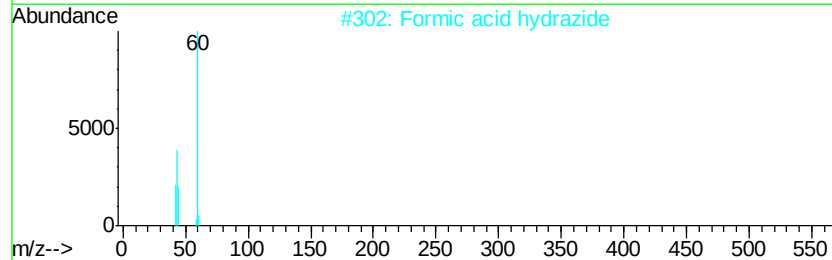
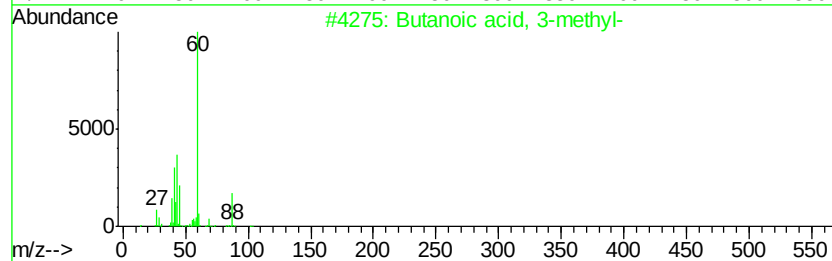
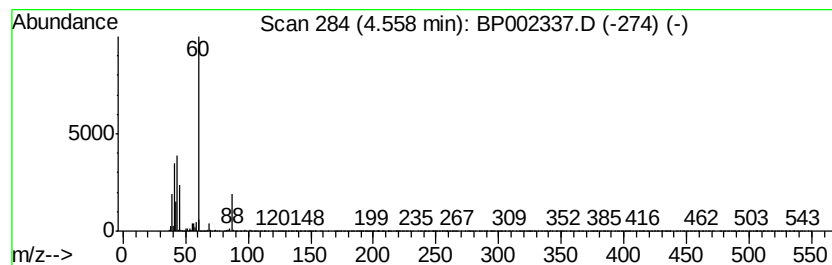
Instrument :
BNA_P
ClientSampleId :
NM33-COMP-2020-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP052720.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butanoic acid, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
4.56	5.81 ng	710583	1,4-Dichlorobenzene-d4	7.60	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	90
2	Formic acid hydrazide	60	CH4N2O	000624-84-0	50
3	Hexanoic acid	116	C6H12O2	000142-62-1	38
4	Pentanoic acid	102	C5H10O2	000109-52-4	33
5	Hydrazine, 1,1-dimethyl-	60	C2H8N2	000057-14-7	9



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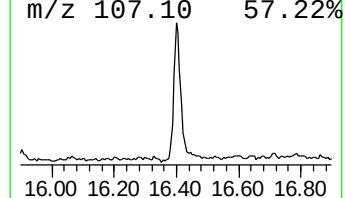
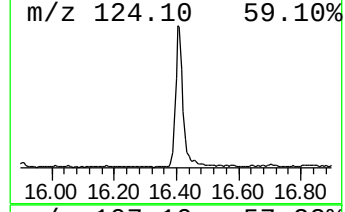
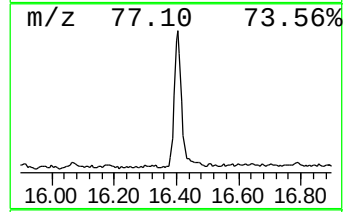
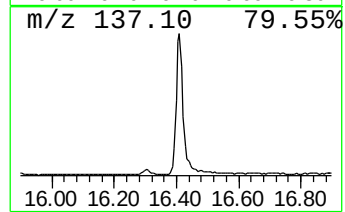
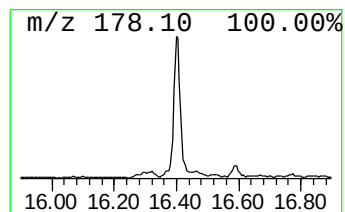
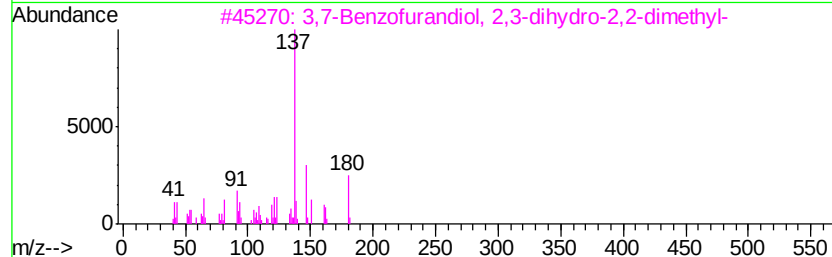
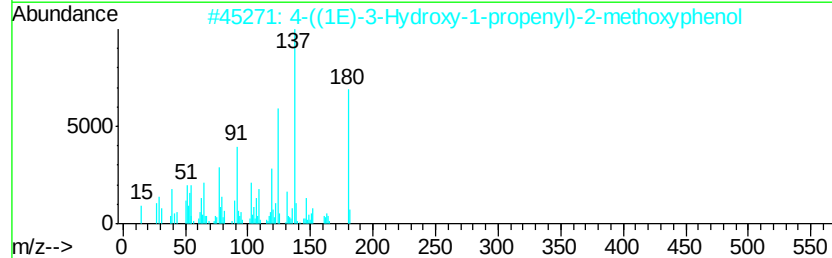
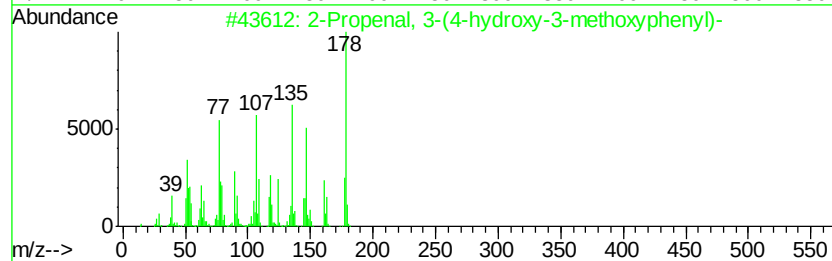
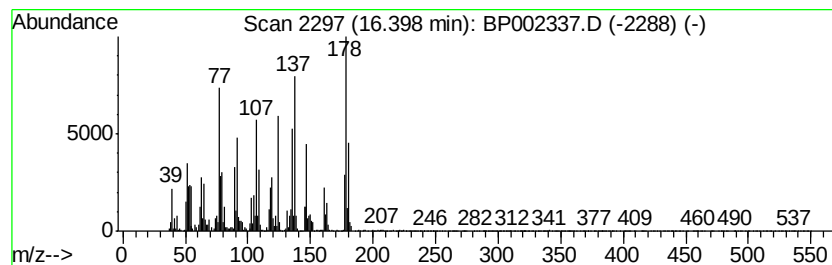
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Propenal, 3-(4-hydroxy-3-... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.40	2.57 ng	670644	Phenanthrene-d10	17.01		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propenal, 3-(4-hydroxy-3-metho...	178	C10H10O3	000458-36-6	95	
2	4-((1E)-3-Hydroxy-1-propenyl)-2-...	180	C10H12O3	1000297-95-5	83	
3	3,7-Benzofurandiol, 2,3-dihydro-...	180	C10H12O3	017781-15-6	50	
4	Benzaldehyde, 2-hydroxy-, oxime	137	C7H7NO2	000094-67-7	35	
5	4-Hydroxy-2-methoxycinnamaldehyde	178	C10H10O3	127321-19-1	25	



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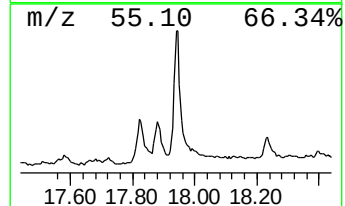
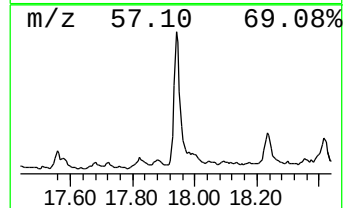
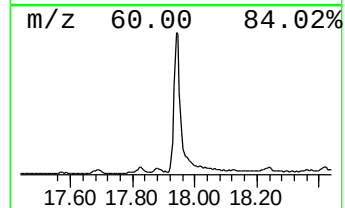
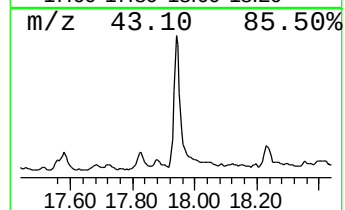
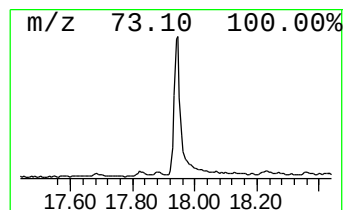
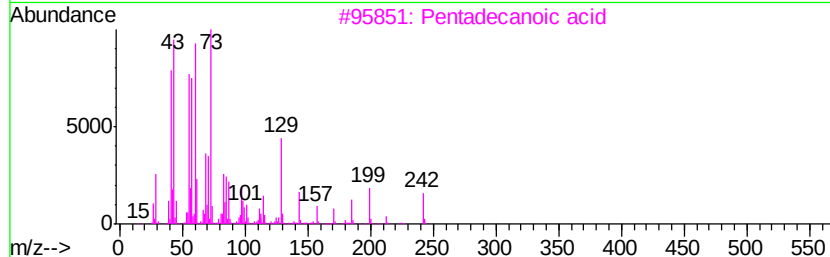
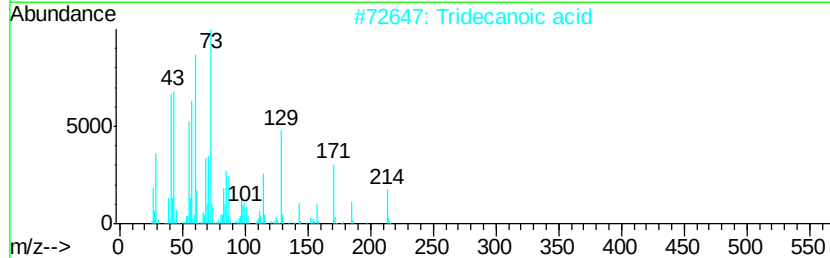
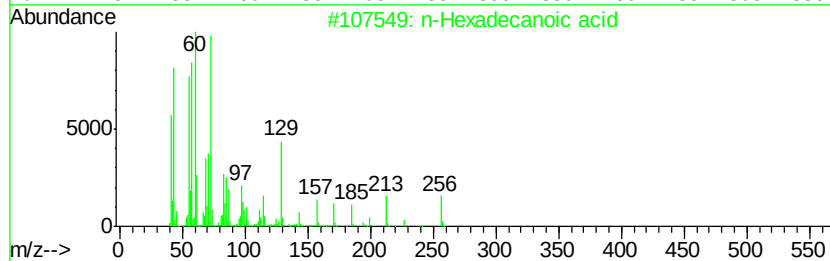
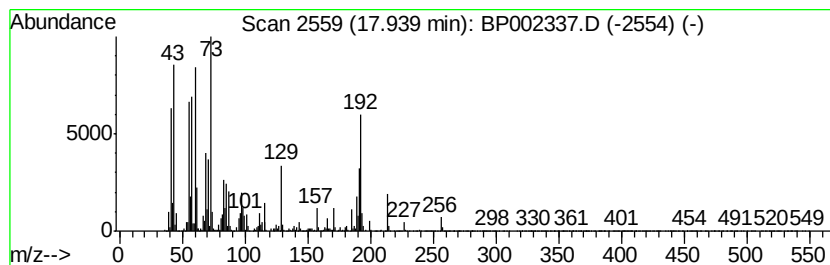
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.94	4.19 ng	1093200	Phenanthrene-d10	17.01

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	95
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	70
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	64
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	64



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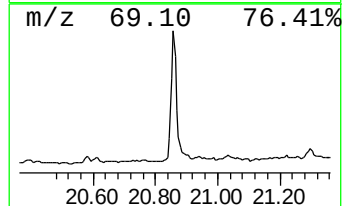
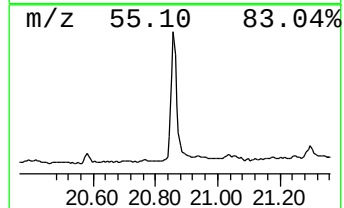
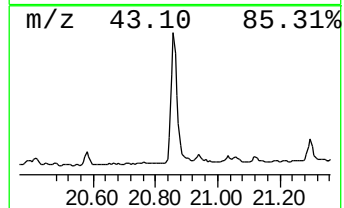
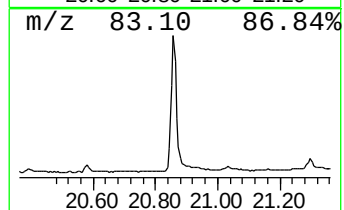
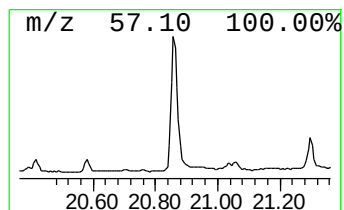
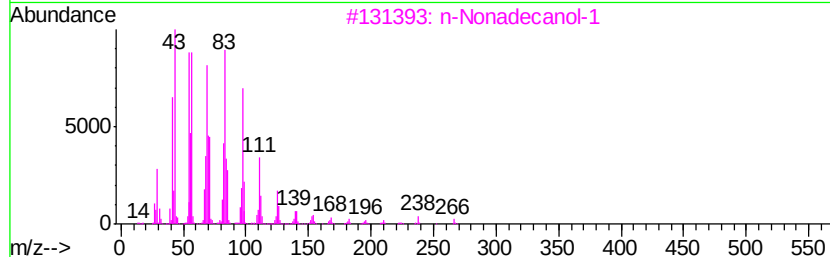
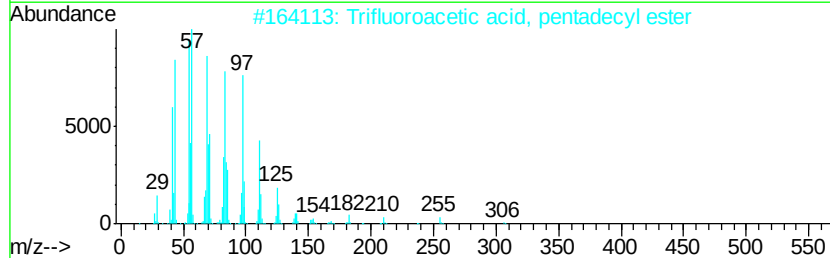
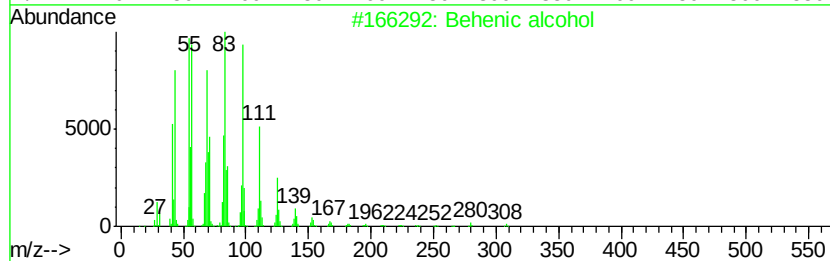
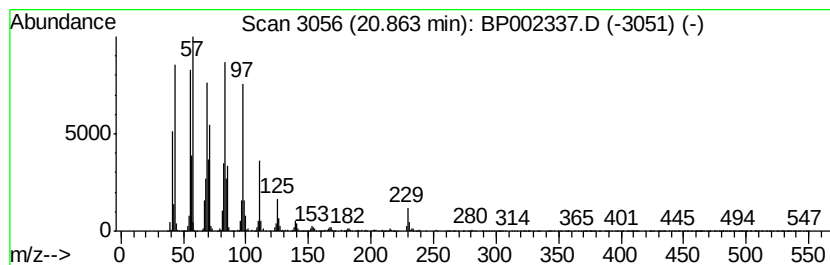
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Trifluoroacetic acid, penta... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.86	6.76 ng	2190360	Chrysene-d12	21.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Behenic alcohol	326	C22H46O	000661-19-8	94
2		Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	94
3		n-Nonadecanol-1	284	C19H40O	001454-84-8	94
4		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
5		1-Heneicosanol	312	C21H44O	015594-90-8	94



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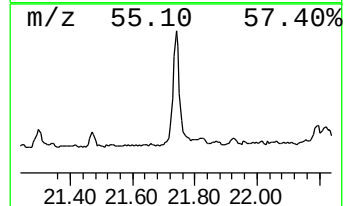
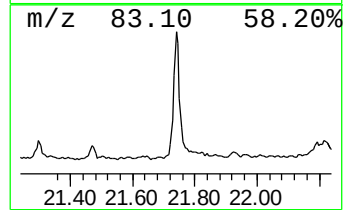
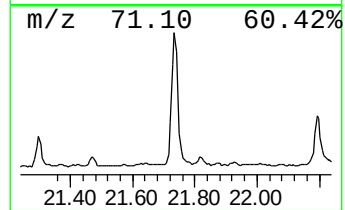
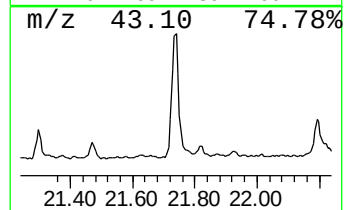
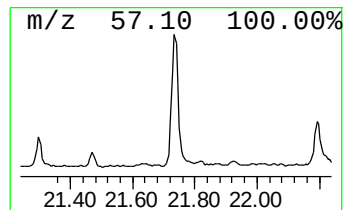
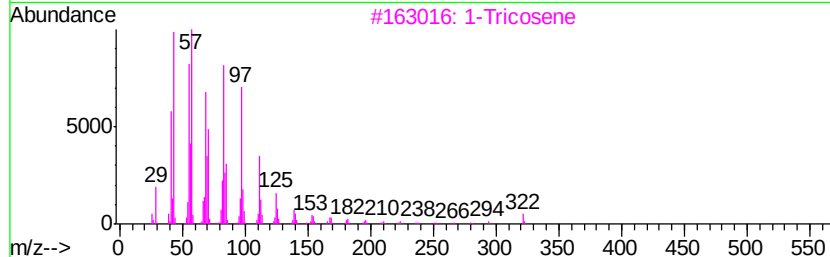
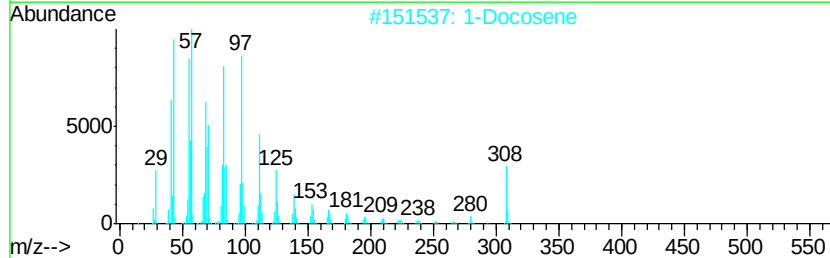
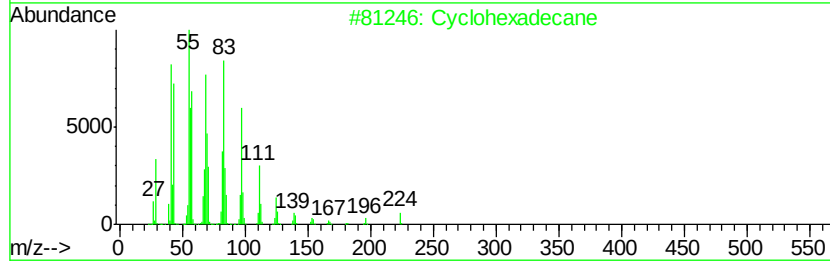
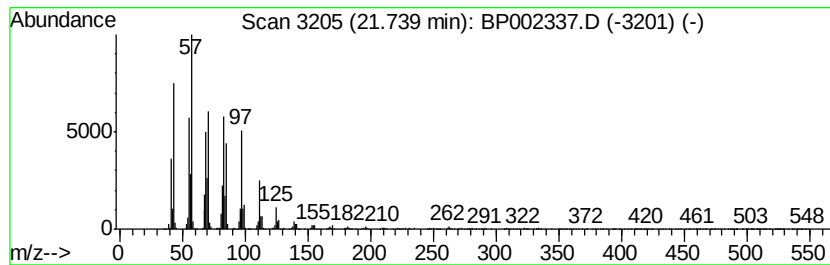
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Cyclohexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.74	4.62 ng	1497540	Chrysene-d12	21.12
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexadecane	224 C16H32	000295-65-8 96
2		1-Docosene	308 C22H44	001599-67-3 95
3		1-Tricosene	322 C23H46	018835-32-0 95
4		17-Pentatriacontene	491 C35H70	006971-40-0 93
5		1-Heneicosanol	312 C21H44O	015594-90-8 92



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP060120\
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Acq On : 01 Jun 2020 14:02
Operator : CG/JU
Sample : L2773-17
Misc :
ALS Vial : 7 Sample Multiplier: 1

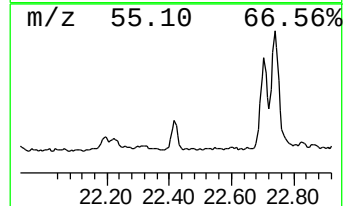
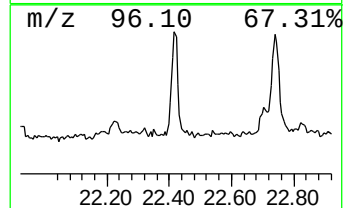
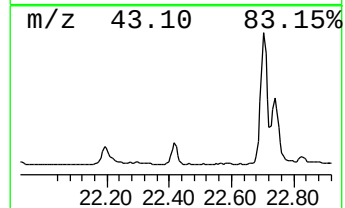
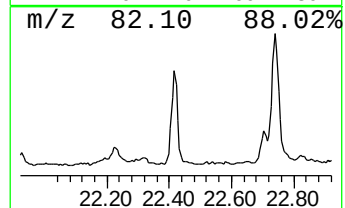
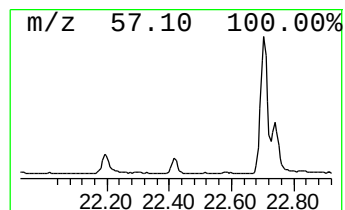
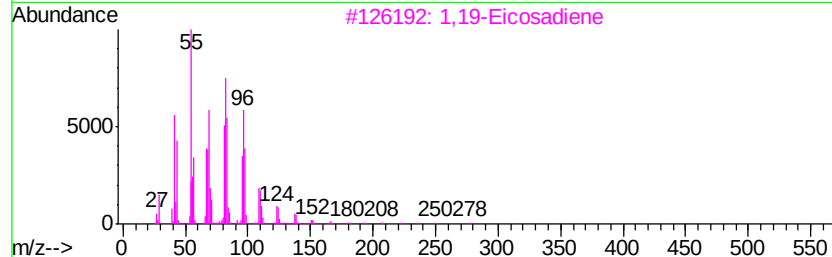
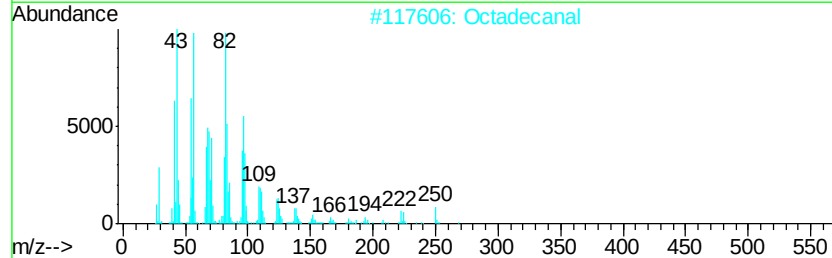
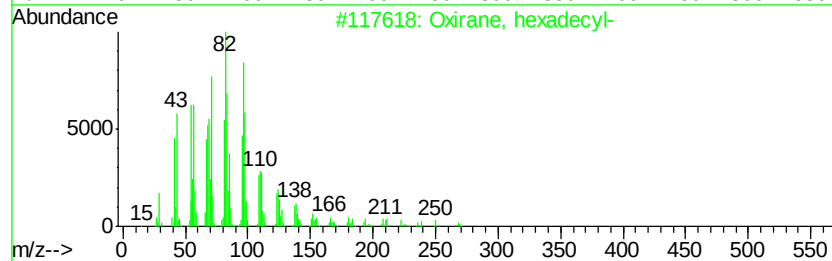
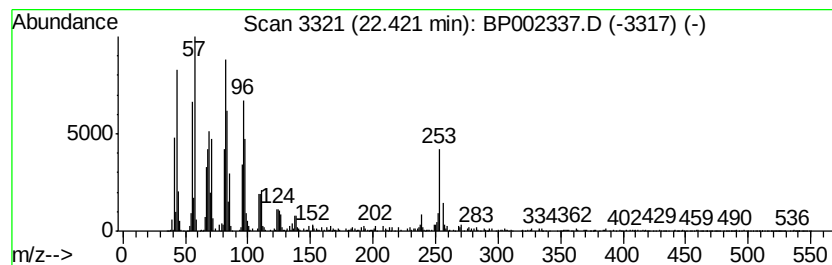
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Oxirane, hexadecyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.42	2.50 ng	564395	Perylene-d12	23.33	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	95
2	Octadecanal	268	C18H36O	000638-66-4	95
3	1,19-Eicosadiene	278	C20H38	014811-95-1	95
4	16-Heptadecenal	252	C17H32O	1000144-57-9	93
5	14-Heptadecenal	252	C17H32O	1000144-58-0	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP060120\
Data File : BP002337.D
Acq On : 01 Jun 2020 14:02
Operator : CG/JU
Sample : L2773-17
Misc :
ALS Vial : 7 Sample Multiplier: 1

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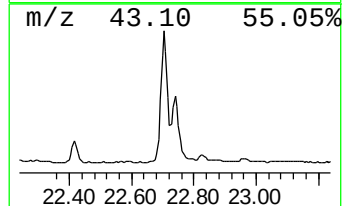
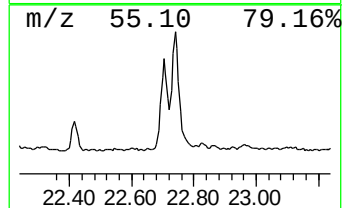
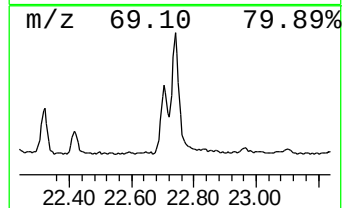
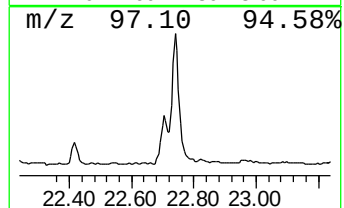
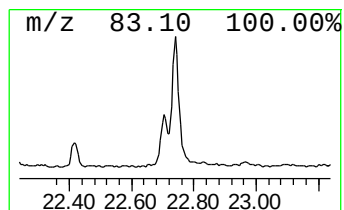
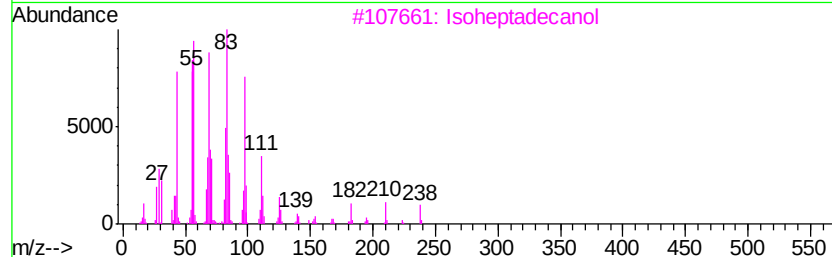
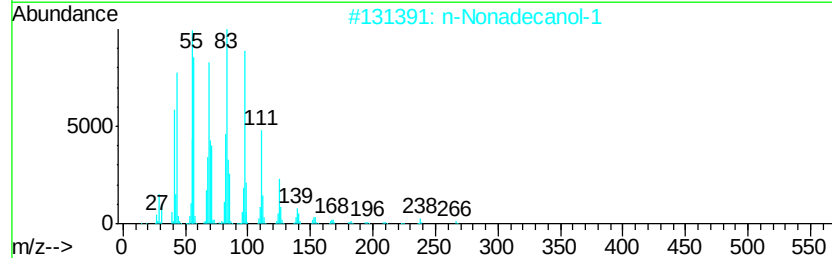
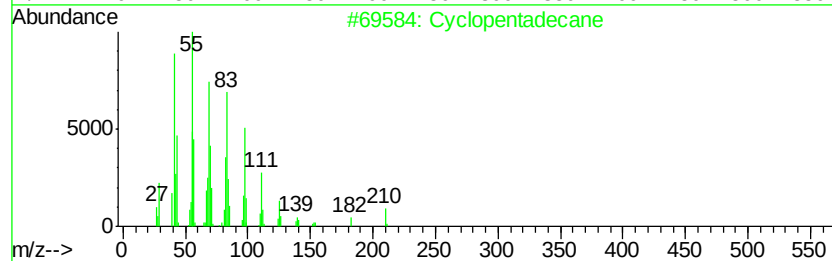
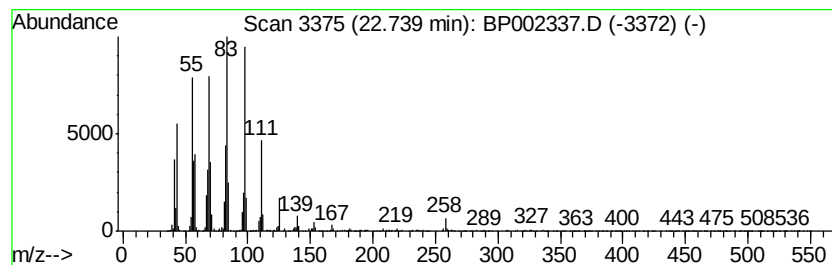
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Cyclopentadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.74	9.24 ng	2082420	Perylene-d12	23.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentadecane	210	C15H30	000295-48-7	91
2		n-Nonadecanol-1	284	C19H40O	001454-84-8	90
3		Isoheptadecanol	256	C17H36O	057289-07-3	80
4		n-Heptadecanol-1	256	C17H36O	001454-85-9	72
5		1-Heneicosanol	312	C21H44O	015594-90-8	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP060120\
Data File : BP002337.D
Acq On : 01 Jun 2020 14:02
Operator : CG/JU
Sample : L2773-17
Misc :
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Instrument :
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ClientSampleId :
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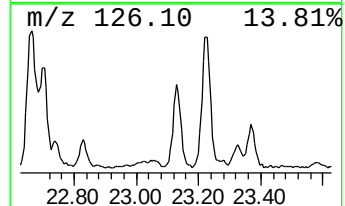
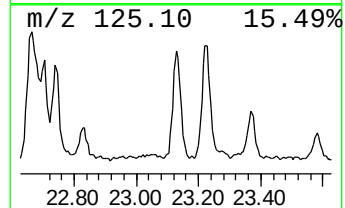
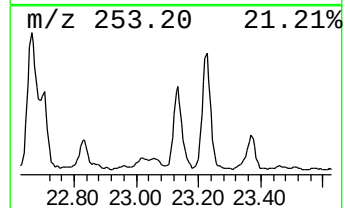
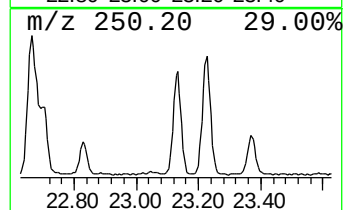
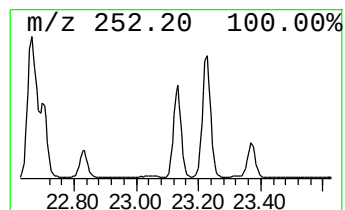
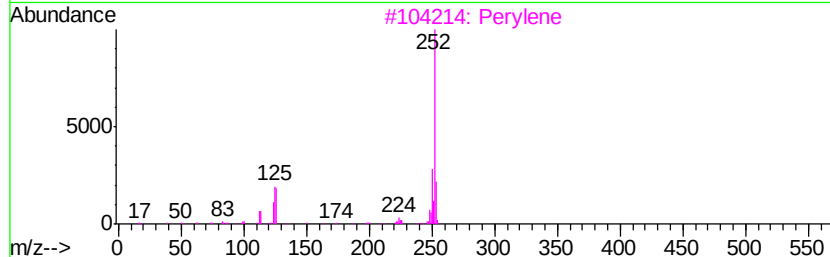
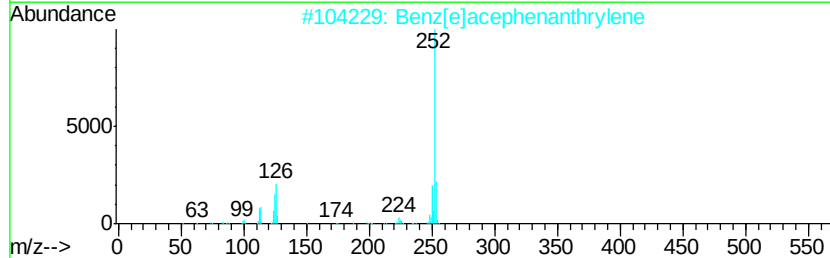
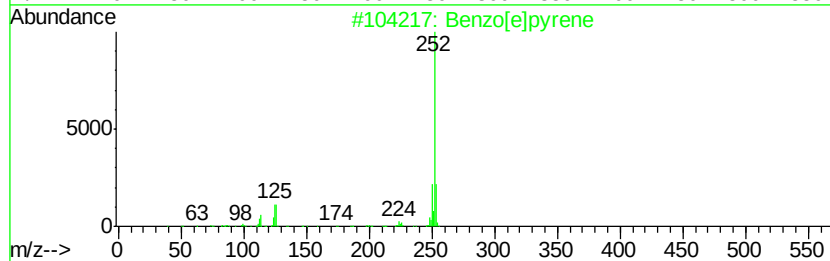
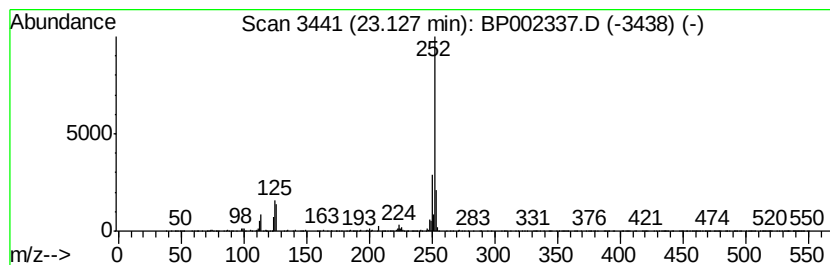
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzo[e]pyrene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.13	3.36 ng	757881	Perylene-d12	23.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	99
2		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	95
3		Perylene	252	C20H12	000198-55-0	94
4		Benzo[k]fluoranthene	252	C20H12	000207-08-9	91
5		Benzo[a]pyrene	252	C20H12	000050-32-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP060120\
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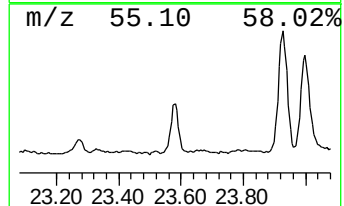
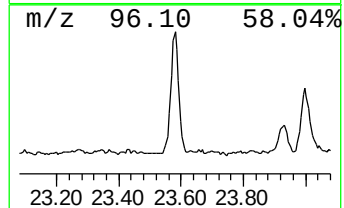
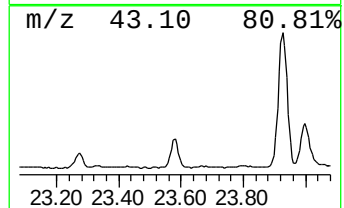
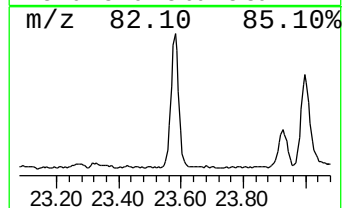
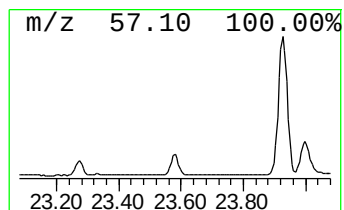
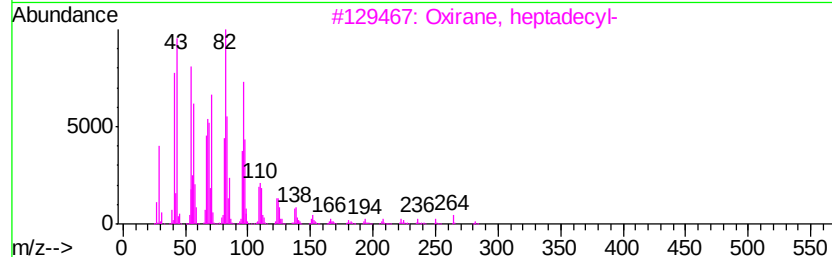
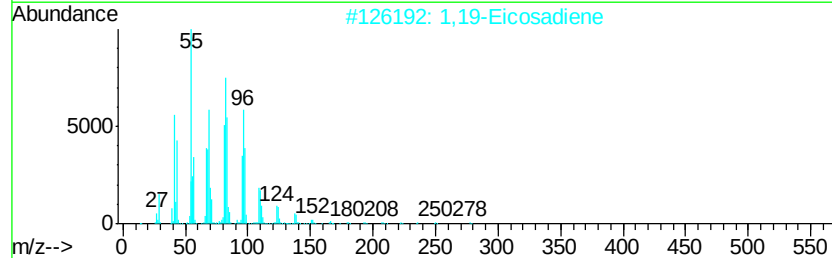
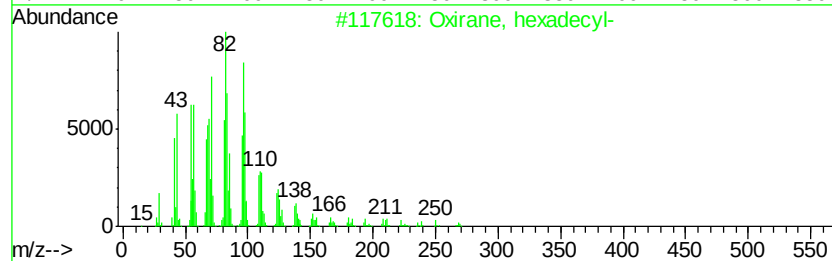
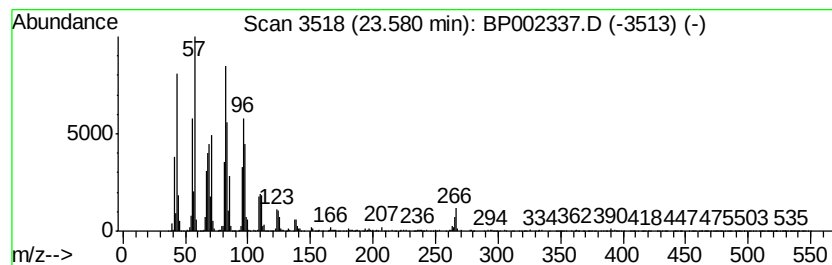
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 10 1,19-Eicosadiene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
23.58	6.79 ng	1529890	Perylene-d12		23.33	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
2		1,19-Eicosadiene	278	C20H38	014811-95-1	90
3		Oxirane, heptadecyl-	282	C19H38O	067860-04-2	87
4		Octadecanal	268	C18H36O	000638-66-4	87
5		17-Octadecenal	266	C18H34O	056554-86-0	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP060120\
Data File : BP002337.D
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Sample : L2773-17
Misc :
ALS Vial : 7 Sample Multiplier: 1

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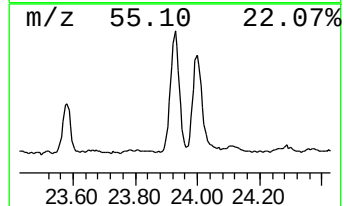
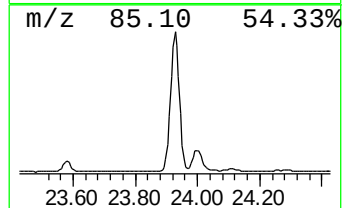
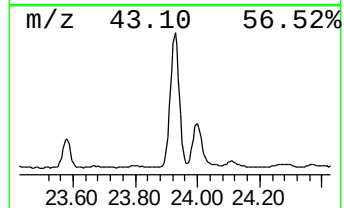
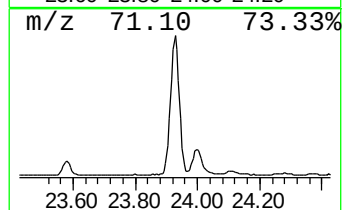
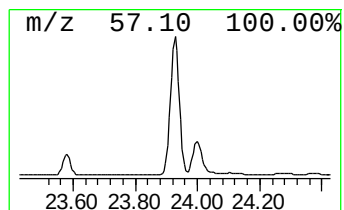
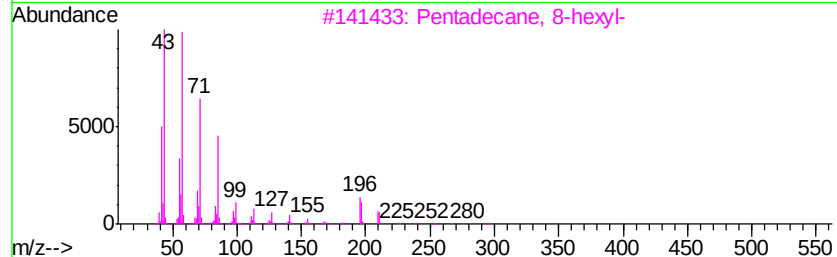
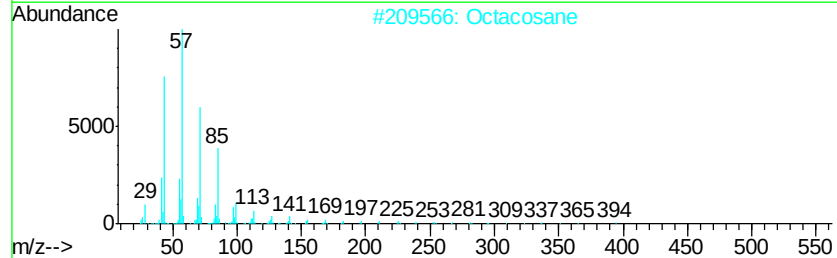
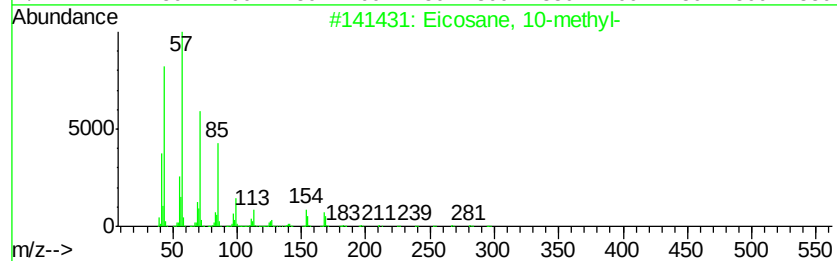
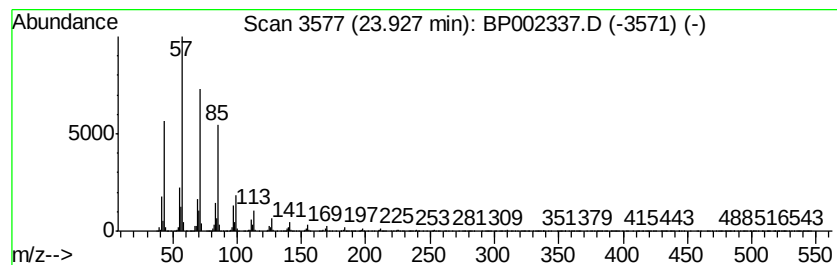
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane, 10-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.93	23.66 ng	5330990	Perylene-d12	23.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	94
2		Octacosane	394	C28H58	000630-02-4	91
3		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91
4		Hexacosane	366	C26H54	000630-01-3	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP060120\
 Data File : BP002337.D
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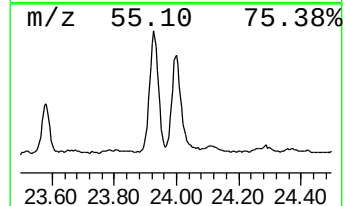
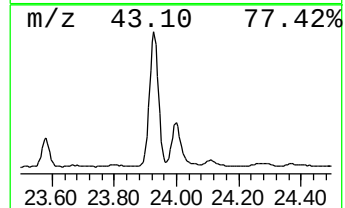
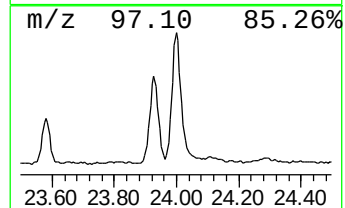
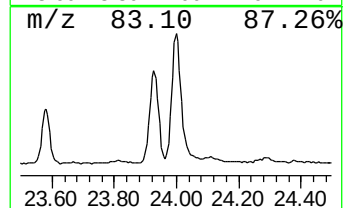
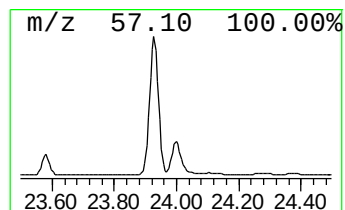
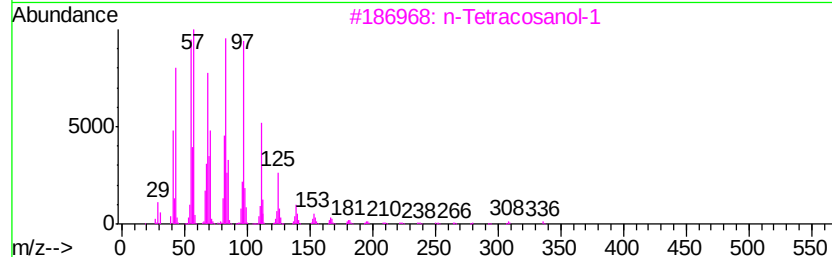
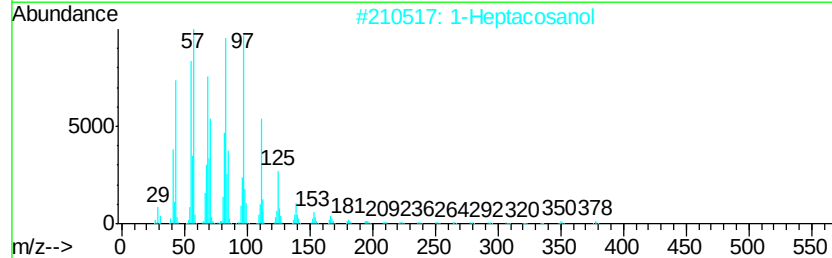
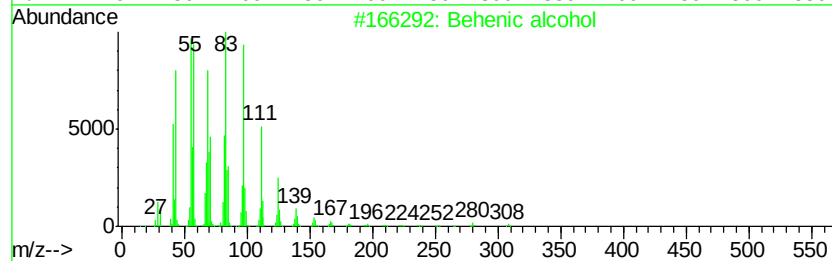
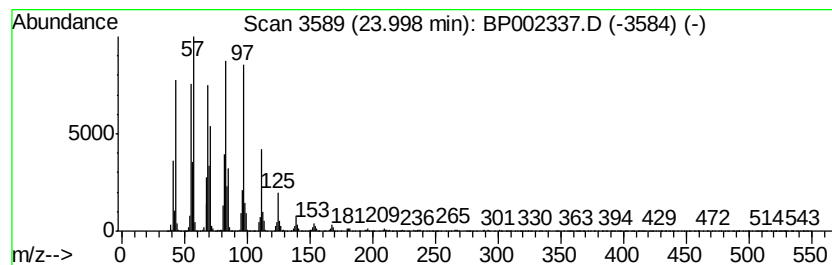
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 12 Behenic alcohol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.00	11.43 ng	2576780	Perylene-d12	23.33

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Behenic alcohol	326	C22H46O	000661-19-8	95
2		1-Heptacosanol	396	C27H56O	002004-39-9	95
3		n-Tetracosanol-1	354	C24H50O	000506-51-4	95
4		Cyclopentadecane	210	C15H30	000295-48-7	90
5		Octacosyl acetate	452	C30H60O2	018206-97-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP060120\
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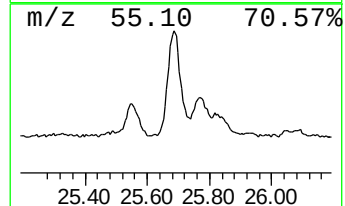
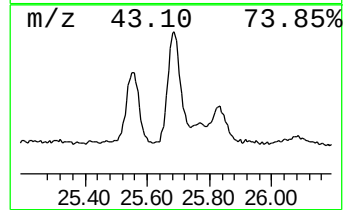
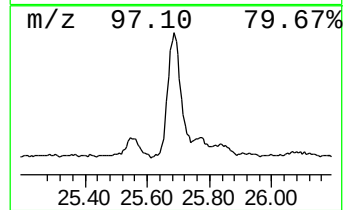
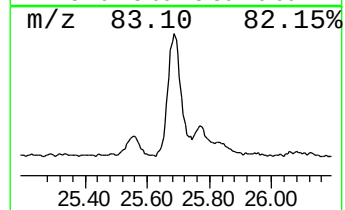
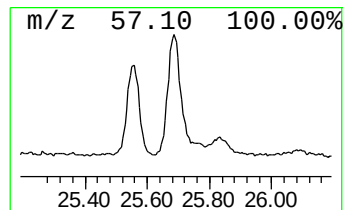
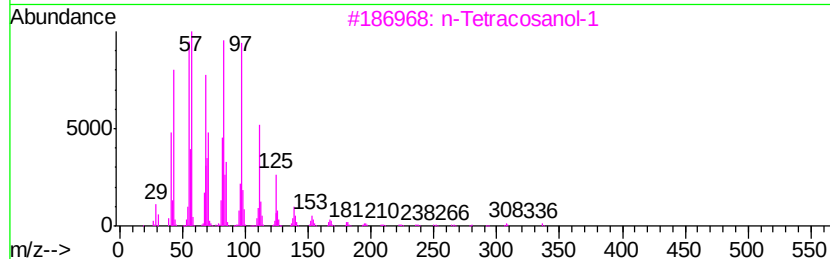
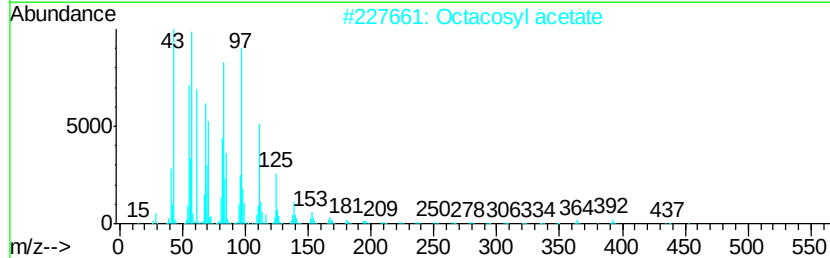
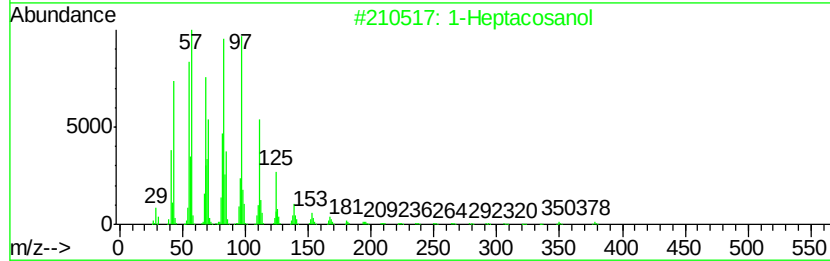
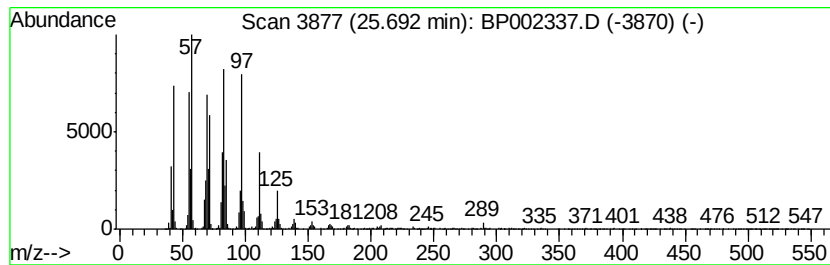
Instrument :
BNA_P
ClientSampleId :
NM33-COMP-2020-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP052720.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1-Heptacosanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
25.69	7.40 ng	1667260	Perylene-d12	23.33	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptacosanol	396	C27H56O	002004-39-9	94
2	Octacosyl acetate	452	C30H60O2	018206-97-8	94
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	93
4	1-Hexacosanol	382	C26H54O	000506-52-5	91
5	1-Triacontanol	438	C30H62O	000593-50-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP060120\
Data File : BP002337.D
Acq On : 01 Jun 2020 14:02
Operator : CG/JU
Sample : L2773-17
Misc :
ALS Vial : 7 Sample Multiplier: 1

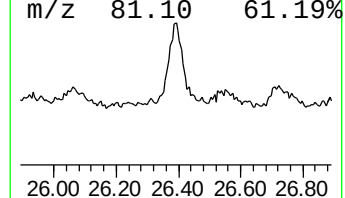
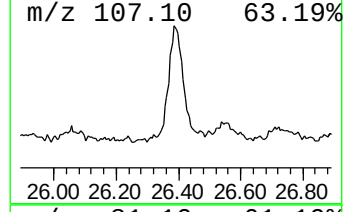
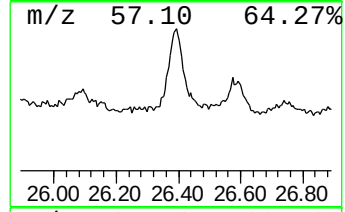
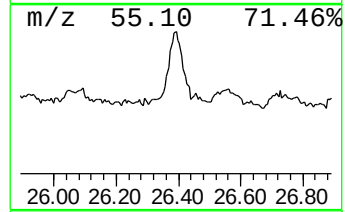
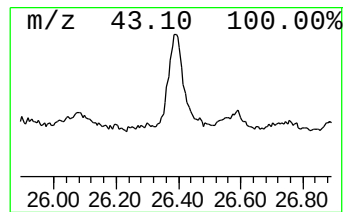
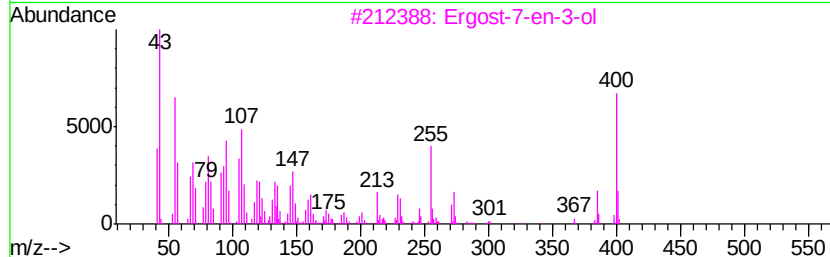
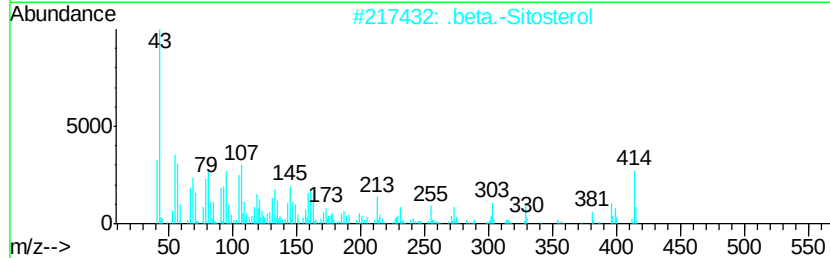
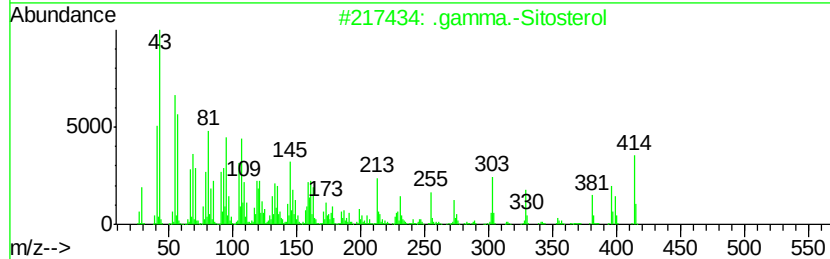
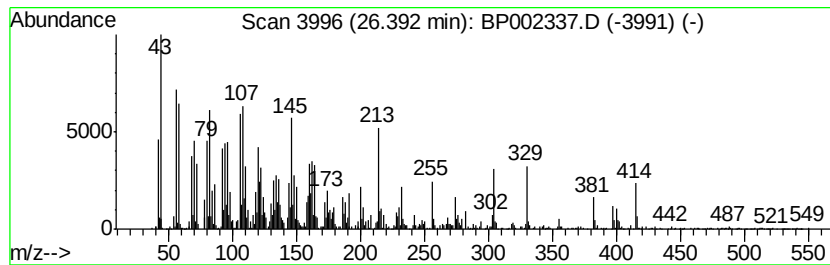
Instrument :
BNA_P
ClientSampleId :
NM33-COMP-2020-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP052720.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 .gamma.-Sitosterol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
26.39	8.12 ng	1831000	Perylene-d12	23.33	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.gamma.-Sitosterol	414	C29H50O	000083-47-6	99
2	.beta.-Sitosterol	414	C29H50O	000083-46-5	93
3	Ergost-7-en-3-ol	400	C28H48O	116179-22-7	35
4	Pyridine-2-carboxylic acid, (2-a...	213	C12H11N3O	1000317-79-6	25
5	Isocholesteryl methyl ether	400	C28H48O	029944-53-4	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP060120\
 Data File : BP002337.D
 Acq On : 01 Jun 2020 14:02
 Operator : CG/JU
 Sample : L2773-17
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 NM33-COMP-2020-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP052720.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
Butanoic acid, 3-...	4.56	5.8	ng	710583	1	7.60	2445690	20.0
2-Propenal, 3-(4-...	16.40	2.6	ng	670644	4	17.01	5223090	20.0
n-Hexadecanoic acid	17.94	4.2	ng	1093200	4	17.01	5223090	20.0
Trifluoroacetic a...	20.86	6.8	ng	2190360	5	21.12	6475930	20.0
Cyclohexadecane	21.74	4.6	ng	1497540	5	21.12	6475930	20.0
Oxirane, hexadecyl-	22.42	2.5	ng	564395	6	23.33	4507120	20.0
Cyclopentadecane	22.74	9.2	ng	2082420	6	23.33	4507120	20.0
Benzo[e]pyrene	23.13	3.4	ng	757881	6	23.33	4507120	20.0
1,19-Eicosadiene	23.58	6.8	ng	1529890	6	23.33	4507120	20.0
Eicosane, 10-methyl-	23.93	23.7	ng	5330990	6	23.33	4507120	20.0
Behenic alcohol	24.00	11.4	ng	2576780	6	23.33	4507120	20.0
1-Heptacosanol	25.69	7.4	ng	1667260	6	23.33	4507120	20.0
.gamma.-Sitosterol	26.39	8.1	ng	1831000	6	23.33	4507120	20.0