

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP082820\
 Data File : BP003156.D
 Acq On : 28 Aug 2020 18:15
 Operator : CG/JU
 Sample : L3837-13
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 789UMR-TP11-0006-01

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-BP082420.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	2.928	3	7	26	rVB	697303	996999	11.85%	1.722%
2	5.269	399	405	422	rBV	1681999	2574424	30.59%	4.447%
3	6.846	665	673	690	rBV	1581100	2592072	30.80%	4.478%
4	7.657	805	811	821	rBV	643495	1060463	12.60%	1.832%
5	8.804	996	1006	1021	rBV	1095701	1829249	21.73%	3.160%
6	10.439	1275	1284	1298	rBV	885873	1534878	18.24%	2.651%
7	12.922	1698	1706	1729	rBV	3350605	5121507	60.85%	8.847%
8	13.763	1842	1849	1861	rBV	125555	215566	2.56%	0.372%
9	14.304	1934	1941	1948	rBV	1388762	2139495	25.42%	3.696%
10	15.792	2188	2194	2219	rBV	2362116	3778158	44.89%	6.526%
11	17.051	2402	2408	2413	rBV	1787075	2702579	32.11%	4.668%
12	17.092	2413	2415	2421	rVB	290154	369458	4.39%	0.638%
13	17.974	2560	2565	2573	rBV2	320312	489068	5.81%	0.845%
14	18.116	2584	2589	2593	rBV2	56379	87183	1.04%	0.151%
15	19.068	2746	2751	2761	rBV	752632	1076092	12.79%	1.859%
16	19.210	2771	2775	2780	rBV3	71741	100968	1.20%	0.174%
17	19.421	2803	2811	2820	rBV	656953	1029878	12.24%	1.779%
18	19.627	2838	2846	2861	rBV	6515963	8416770	100.00%	14.539%
19	19.910	2890	2894	2899	rVB	91496	119100	1.42%	0.206%
20	20.633	3014	3017	3025	rVB	57559	91224	1.08%	0.158%
21	20.880	3055	3059	3065	rBV4	746046	1092492	12.98%	1.887%
22	21.145	3097	3104	3108	rBV2	2957982	4312718	51.24%	7.450%
23	21.180	3108	3110	3119	rVB	480957	551171	6.55%	0.952%
24	21.298	3127	3130	3132	rBV	74177	101100	1.20%	0.175%
25	21.751	3202	3207	3220	rBV4	422868	958333	11.39%	1.655%
26	22.198	3279	3283	3296	rBV4	181308	367478	4.37%	0.635%
27	22.339	3304	3307	3315	rVB	175484	242684	2.88%	0.419%
28	22.439	3320	3324	3333	rVB2	347012	497414	5.91%	0.859%
29	22.721	3364	3372	3377	rBV2	993717	2307711	27.42%	3.986%
30	22.762	3377	3379	3392	rVB2	490604	885824	10.52%	1.530%
31	23.186	3446	3451	3459	rVB	280796	523653	6.22%	0.905%
32	23.280	3462	3467	3476	rVB	379386	795722	9.45%	1.375%
33	23.380	3476	3484	3491	rBV	2052693	3959647	47.04%	6.840%
34	23.615	3519	3524	3534	rVB	421508	734735	8.73%	1.269%

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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

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35	23.962	3577	3583	3589	rVB	489106	943975	11.22%	1.631%
36	24.045	3591	3597	3610	rVV	268091	615341	7.31%	1.063%
37	25.168	3783	3788	3798	rVB3	280421	669699	7.96%	1.157%
38	25.627	3859	3866	3880	rVB3	314808	1228172	14.59%	2.122%
39	26.492	4007	4013	4033	rVB9	217357	777499	9.24%	1.343%

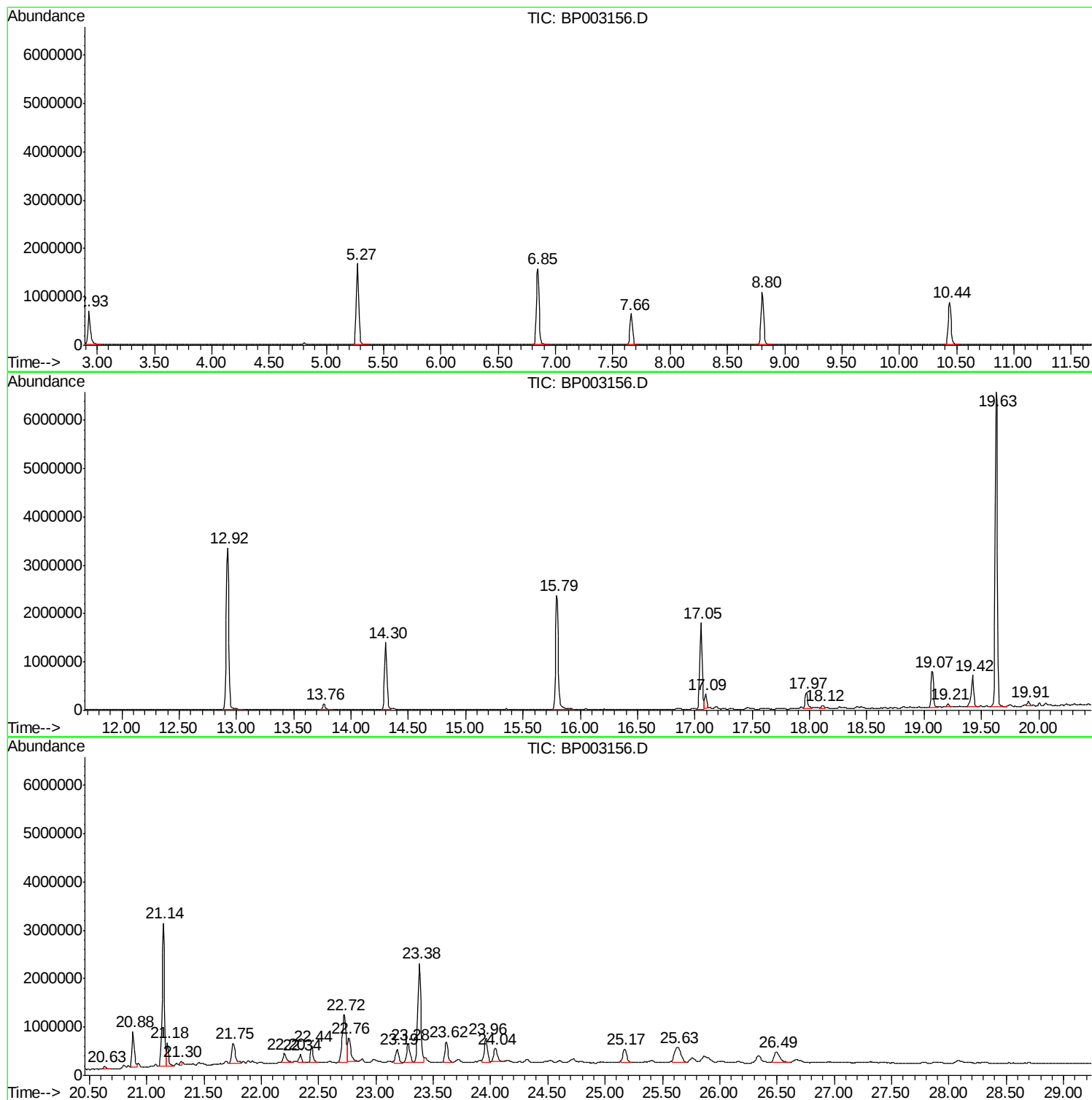
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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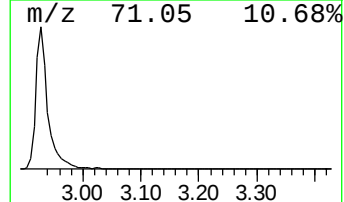
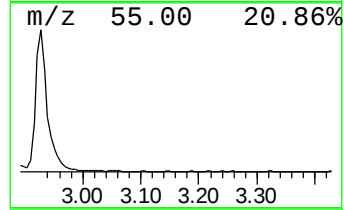
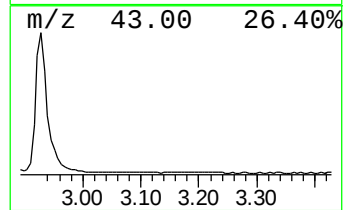
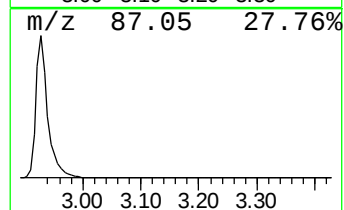
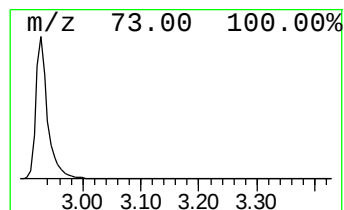
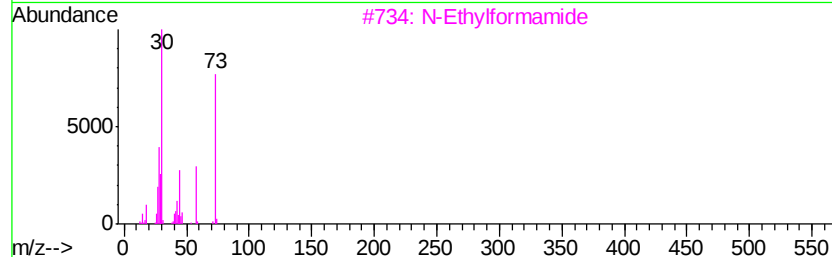
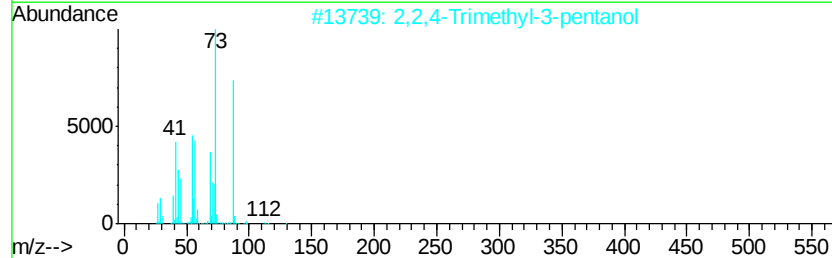
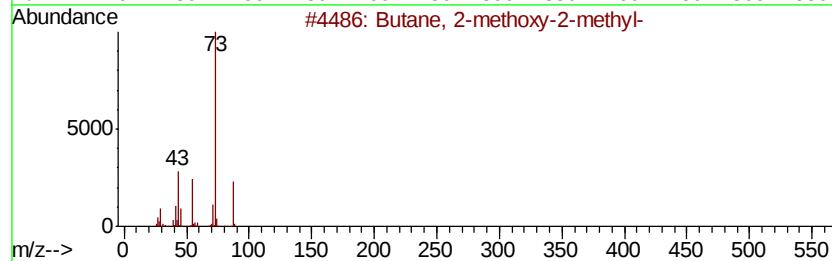
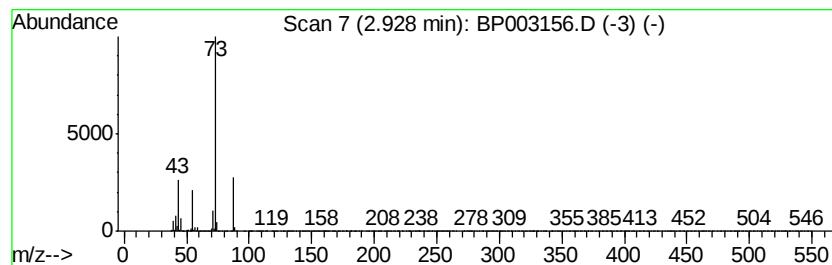
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.93	18.80 ng	996999	1,4-Dichlorobenzene-d4	7.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	43
3			N-Ethylformamide	73	C3H7NO	000627-45-2	9
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9
5			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	7



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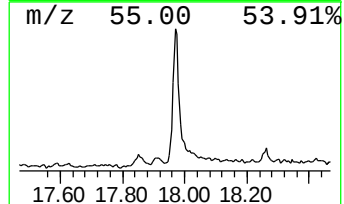
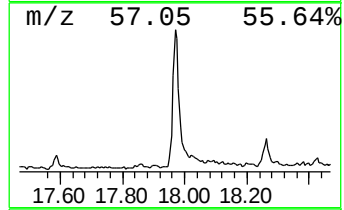
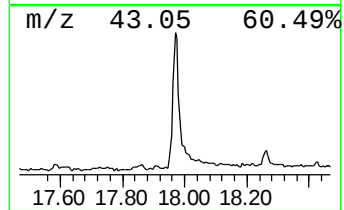
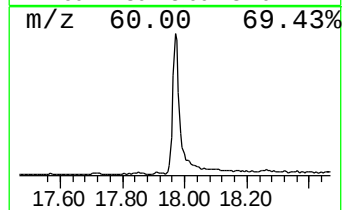
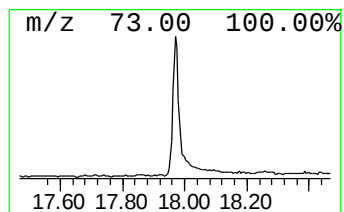
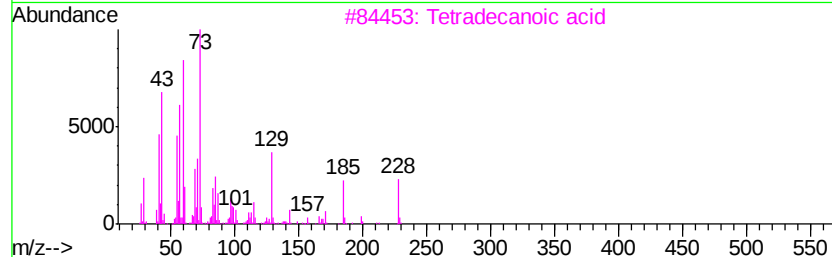
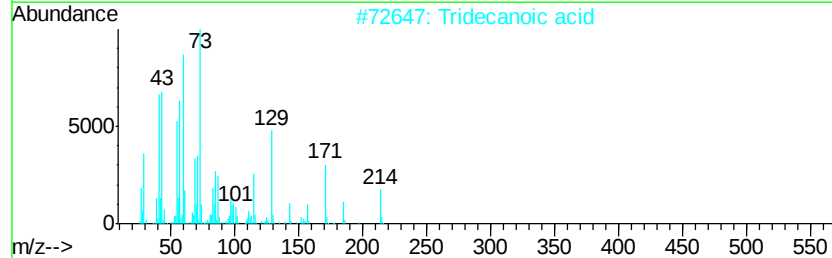
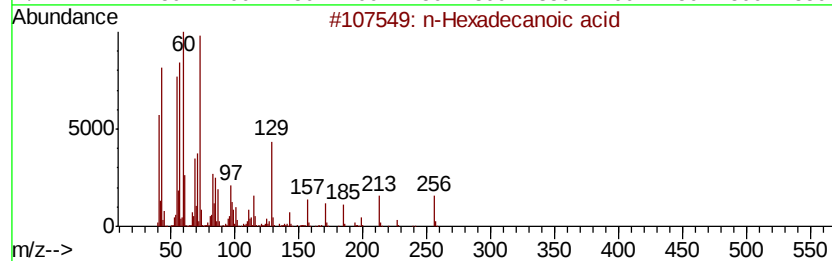
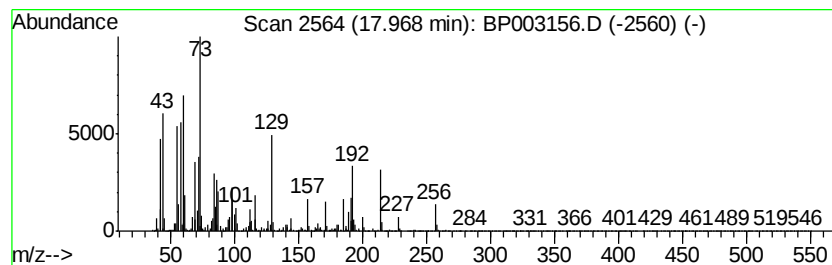
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
17.97	3.62 ng	489068	Phenanthrene-d10		17.05
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	78
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
4	Octadecanoic acid	284	C18H36O2	000057-11-4	50
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	43



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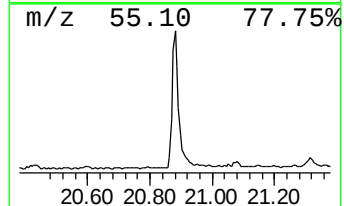
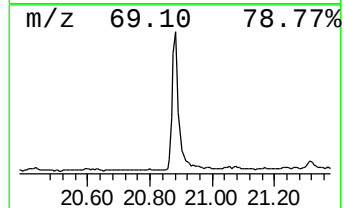
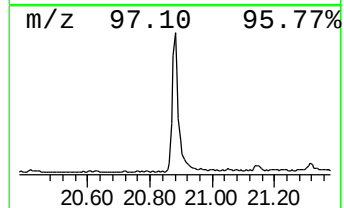
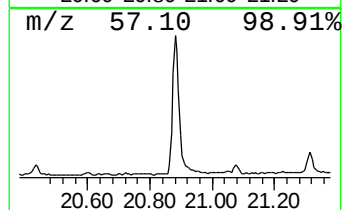
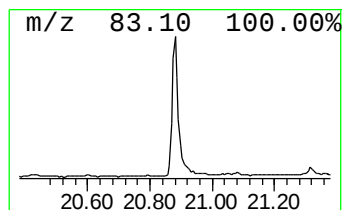
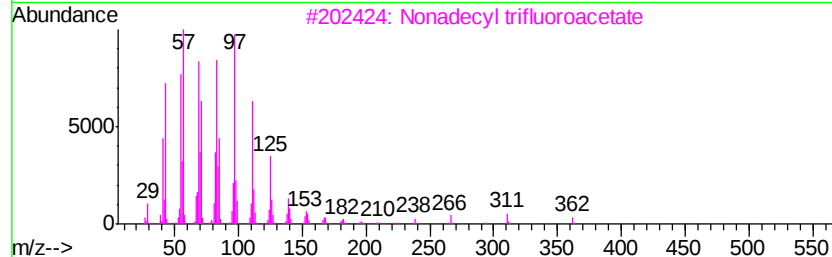
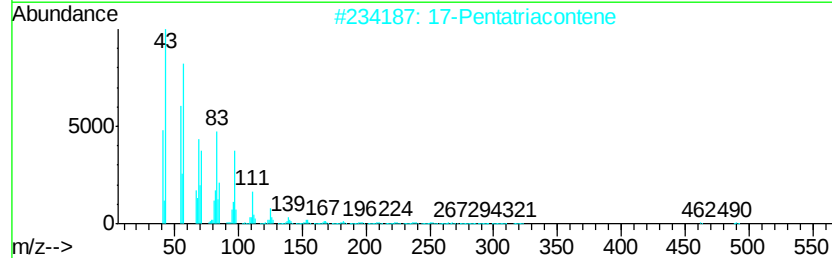
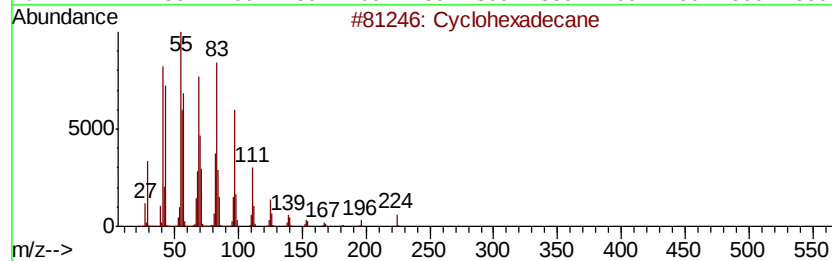
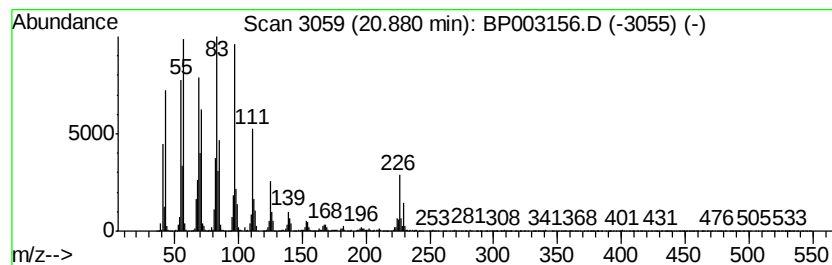
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Cyclohexadecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.88	5.07 ng	1092490	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexadecane	224	C16H32	000295-65-8	94
2		17-Pentatriacontene	491	C35H70	006971-40-0	94
3		Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	93
4		E-15-Heptadecenal	252	C17H32O	1000130-97-9	93
5		n-Tetracosanol-1	354	C24H50O	000506-51-4	93



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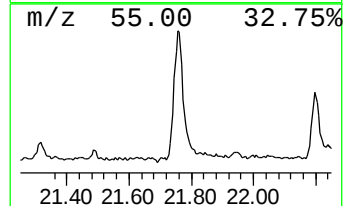
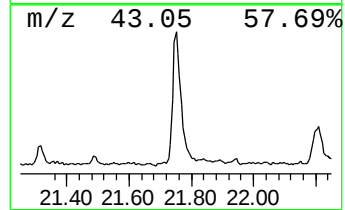
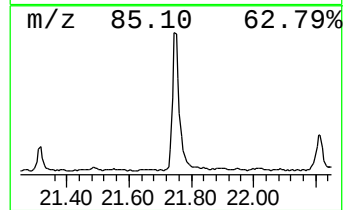
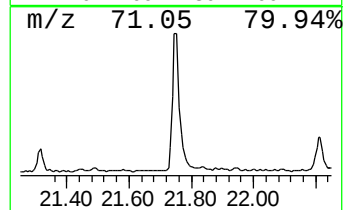
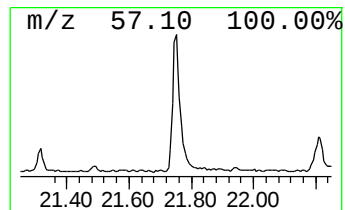
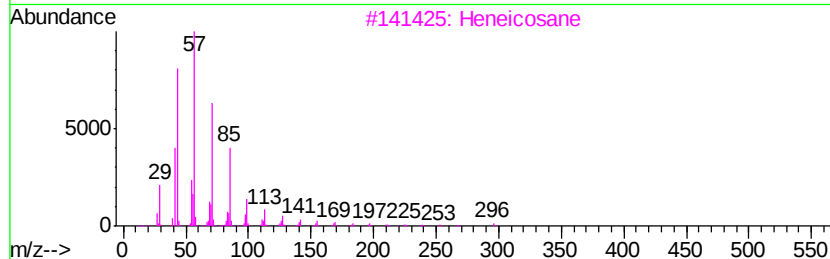
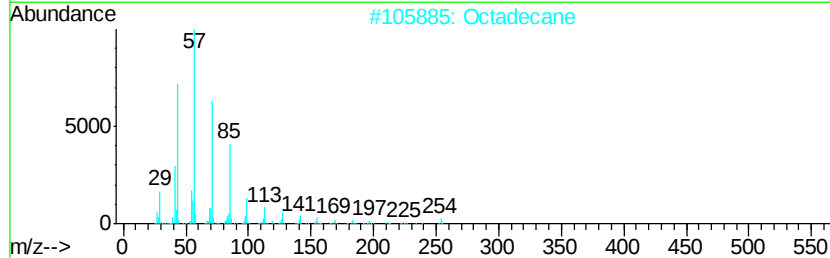
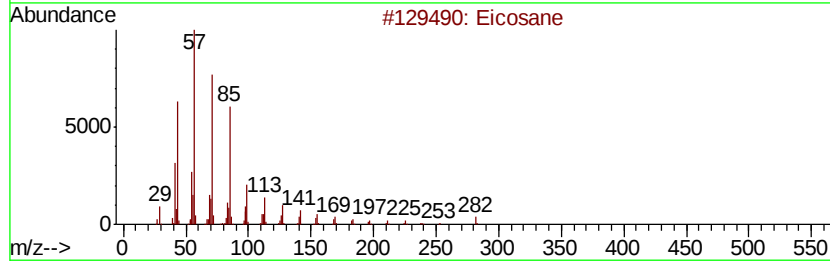
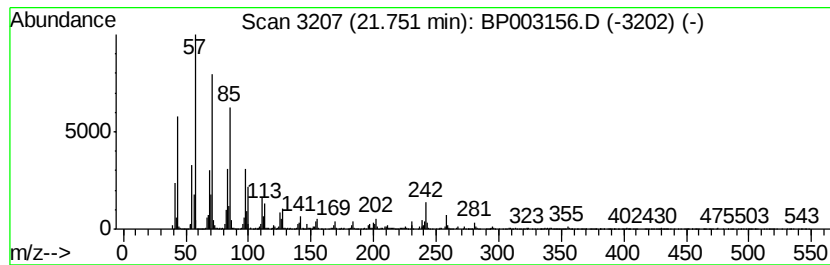
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Eicosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.		
21.75	4.44 ng	958333	Chrysene-d12		21.14		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	97
2			Octadecane	254	C18H38	000593-45-3	95
3			Heneicosane	296	C21H44	000629-94-7	95
4			Heptadecane	240	C17H36	000629-78-7	95
5			triacontane, 1-bromo-	500	C30H61Br	004209-22-7	95



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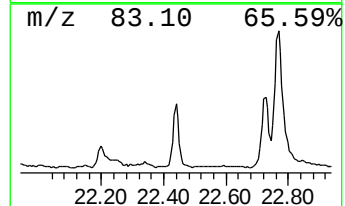
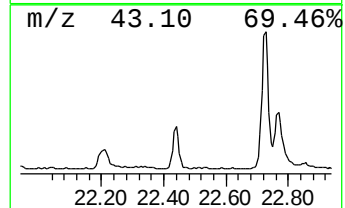
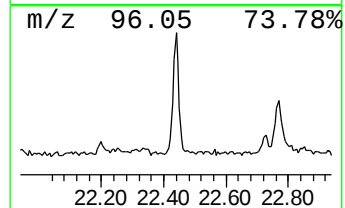
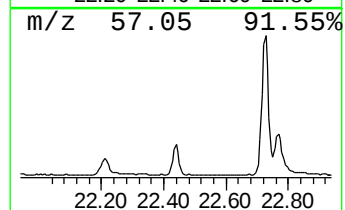
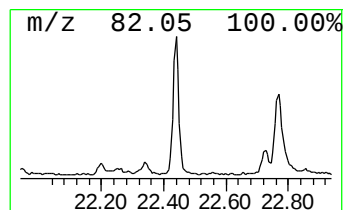
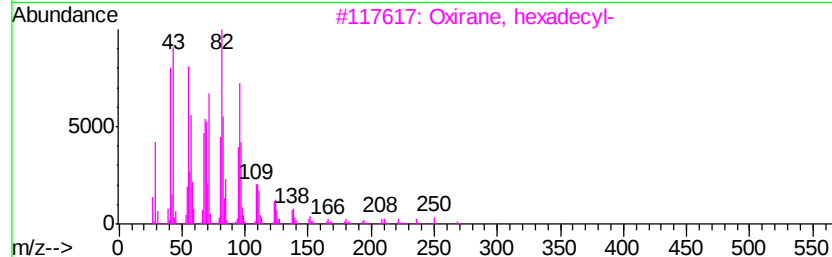
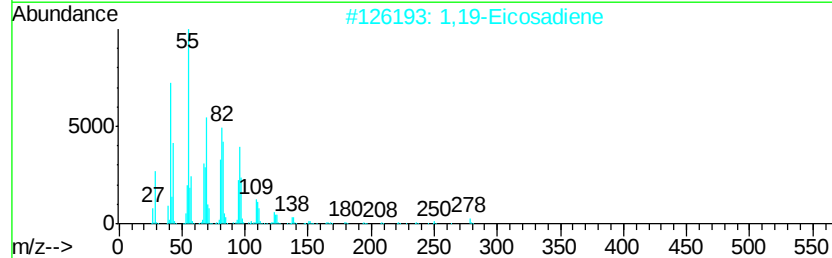
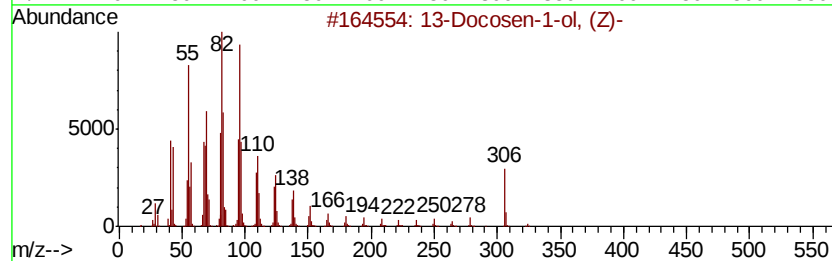
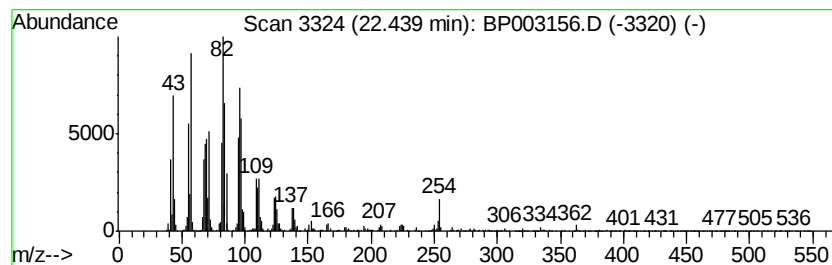
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 13-Docosen-1-ol, (Z)- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.44	2.51 ng	497414	Perylene-d12	23.38	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	13-Docosen-1-ol, (Z)-	324	C22H44O	000629-98-1	96
2	1,19-Eicosadiene	278	C20H38	014811-95-1	96
3	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	96
4	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	96
5	Octadecanal	268	C18H36O	000638-66-4	94



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Operator : CG/JU
Sample : L3837-13
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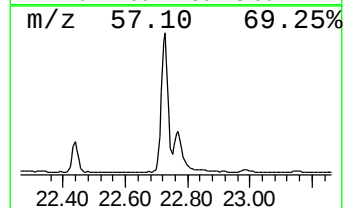
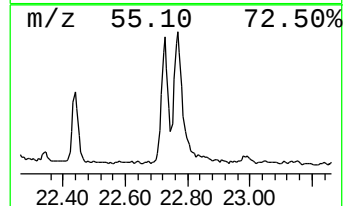
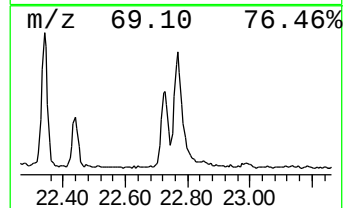
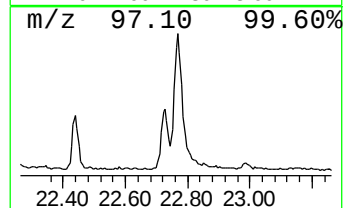
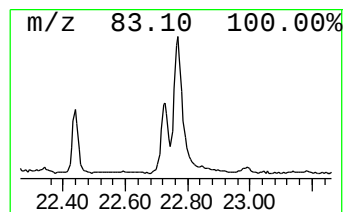
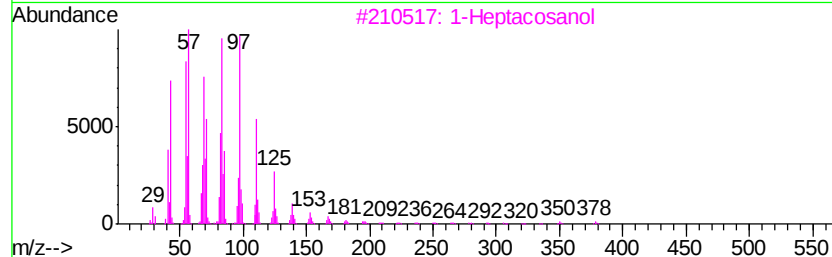
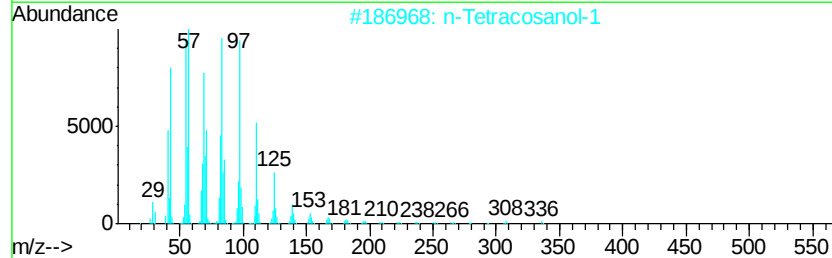
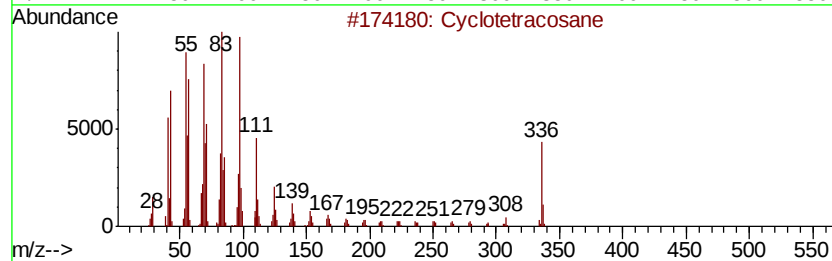
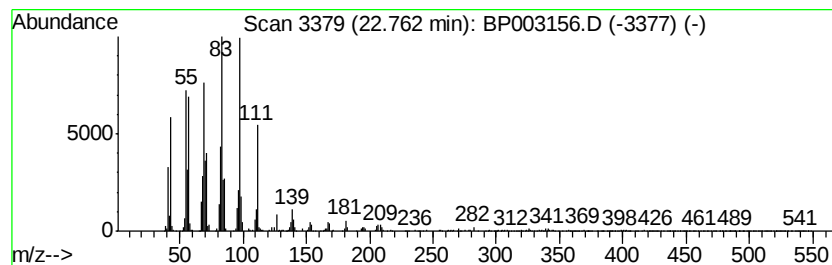
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 n-Tetracosanol-1 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.76	4.47 ng	885824	Perylene-d12	23.38	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetracosane	336	C24H48	000297-03-0	95
2	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
3	1-Heptacosanol	396	C27H56O	002004-39-9	91
4	Behenic alcohol	326	C22H46O	000661-19-8	91
5	Bacteriochlorophyll-c-stearyl	841	C52H72MgN4O4	1000164-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP082820\
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ClientSampleId :
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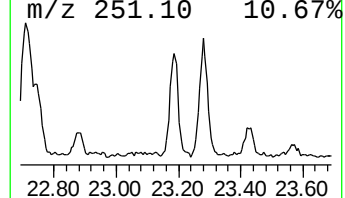
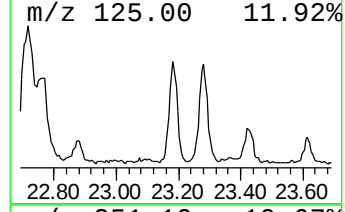
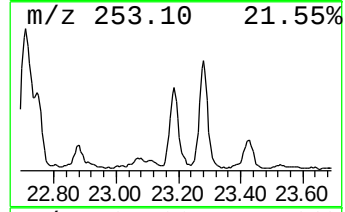
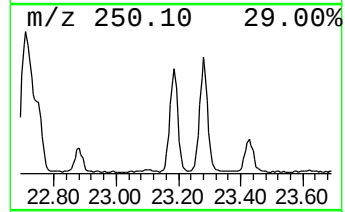
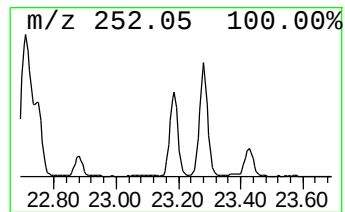
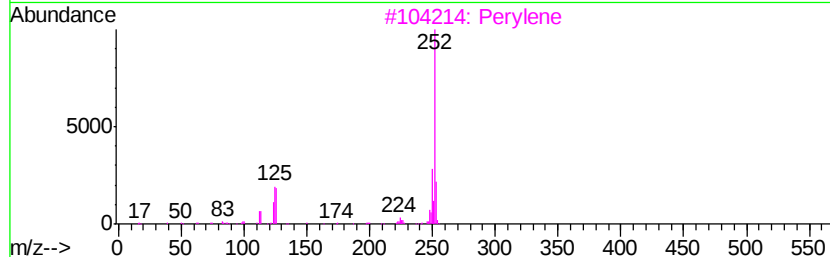
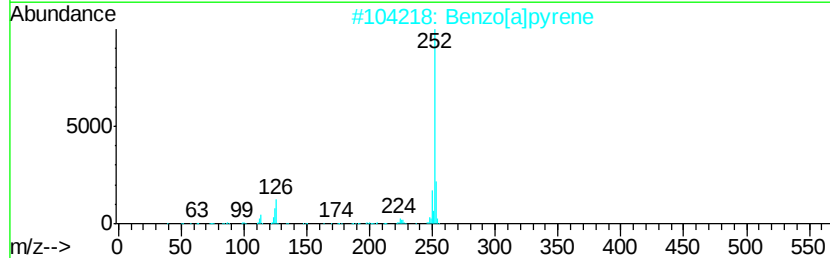
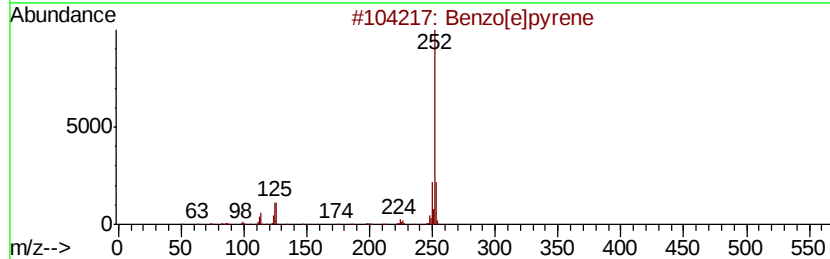
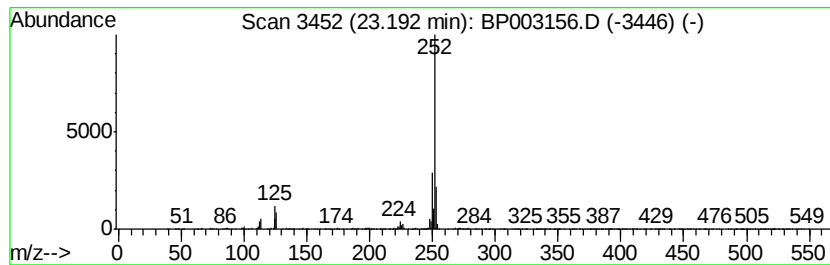
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Benzo[e]pyrene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.19	2.64 ng	523653	Perylene-d12	23.38

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[a]pyrene	252	C20H12	000050-32-8	98
3			Perylene	252	C20H12	000198-55-0	95
4			Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	94
5			Benzo[k]fluoranthene	252	C20H12	000207-08-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP082820\
Data File : BP003156.D
Acq On : 28 Aug 2020 18:15
Operator : CG/JU
Sample : L3837-13
Misc :
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ClientSampleId :
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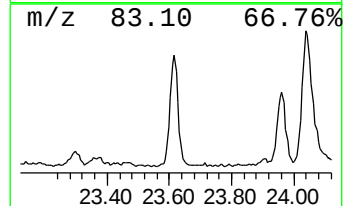
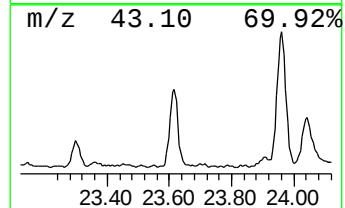
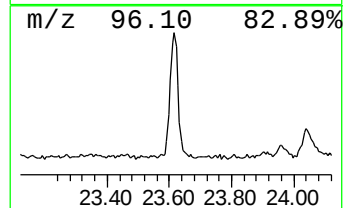
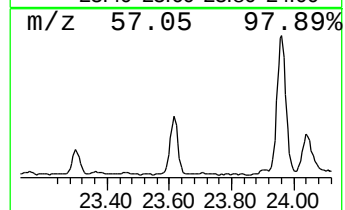
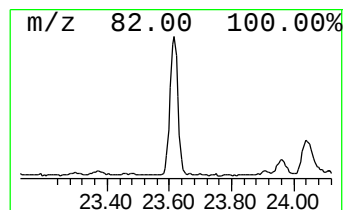
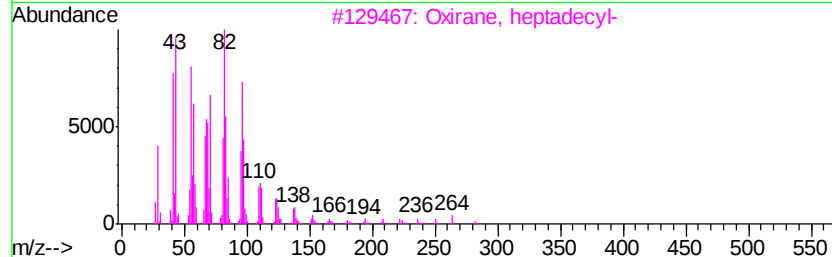
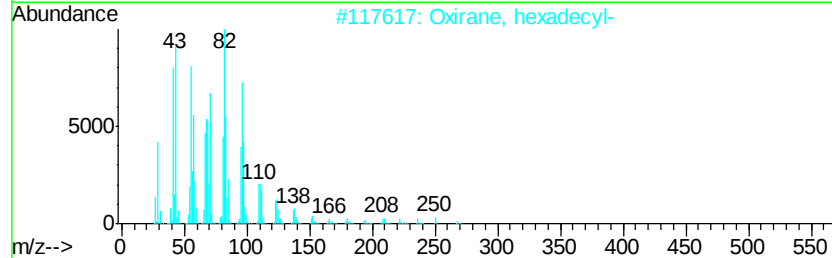
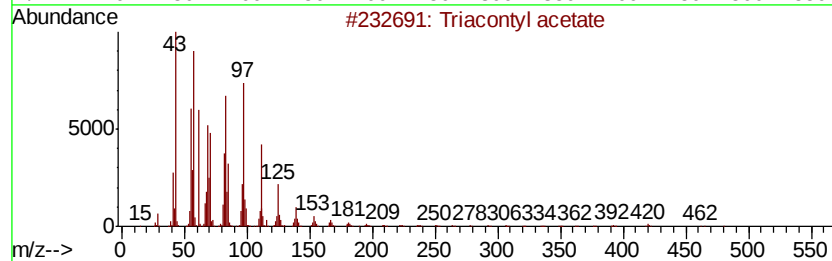
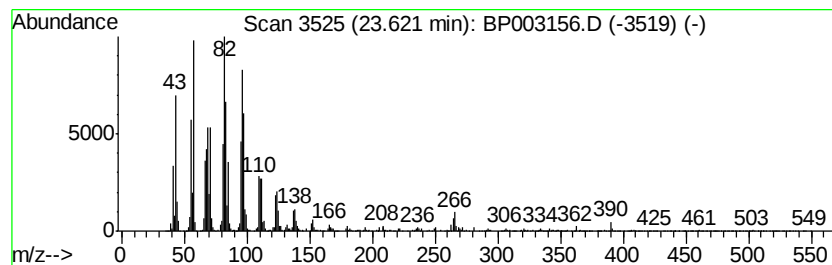
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Triacontyl acetate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.62	3.71 ng	734735	Perylene-d12	23.38

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Triacontyl acetate	480	C32H64O2	041755-58-2	93
2		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
3		Oxirane, heptadecyl-	282	C19H38O	067860-04-2	91
4		9-Octadecen-1-ol, (Z)-	268	C18H36O	000143-28-2	89
5		Octadecanal	268	C18H36O	000638-66-4	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP082820\
Data File : BP003156.D
Acq On : 28 Aug 2020 18:15
Operator : CG/JU
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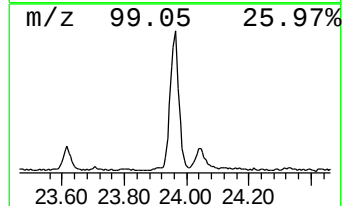
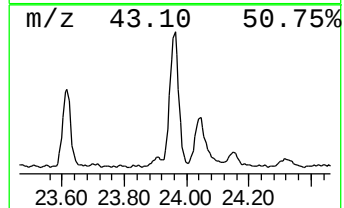
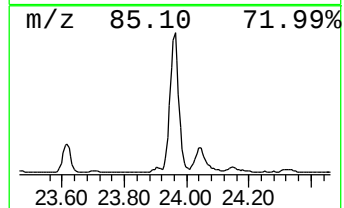
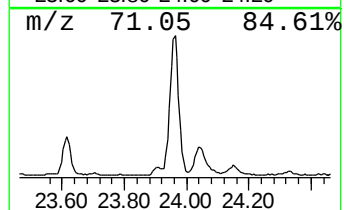
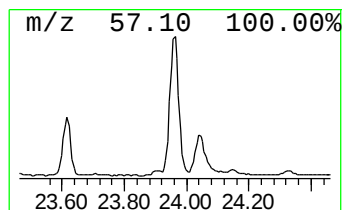
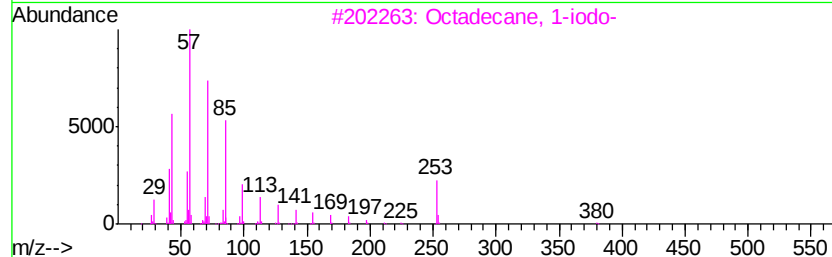
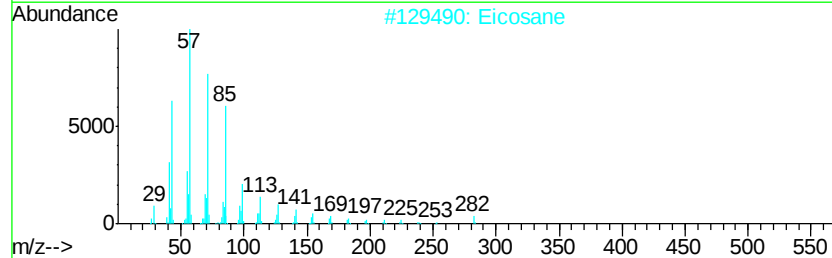
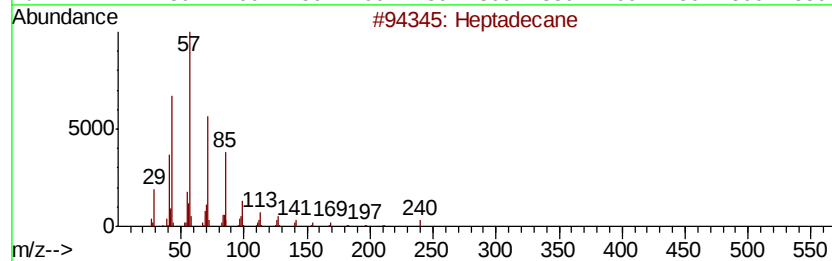
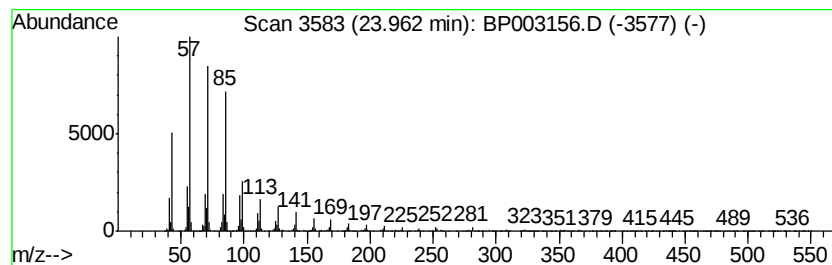
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.96	4.77 ng	943975	Perylene-d12	23.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	94
2		Eicosane	282	C20H42	000112-95-8	91
3		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	91
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
5		Pentacosane	352	C25H52	000629-99-2	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP082820\
Data File : BP003156.D
Acq On : 28 Aug 2020 18:15
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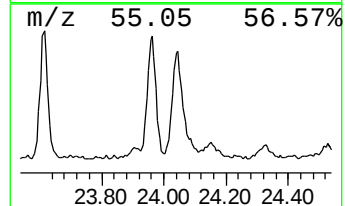
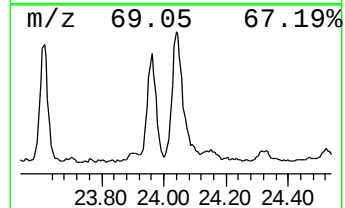
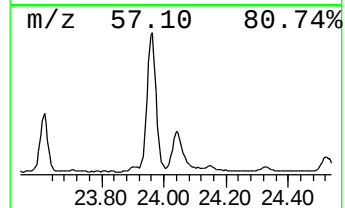
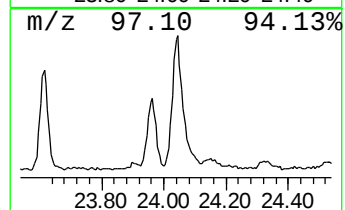
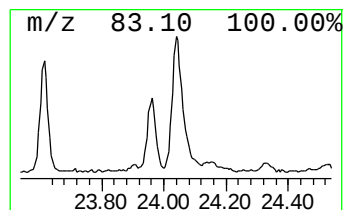
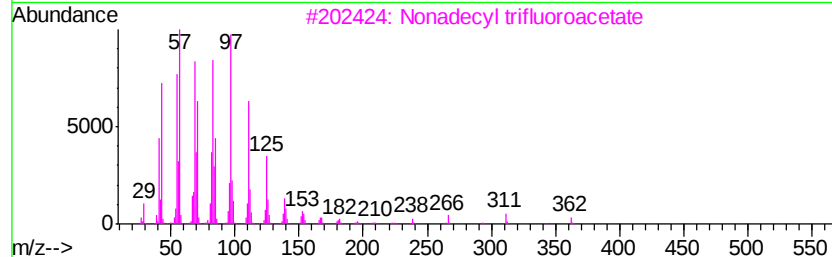
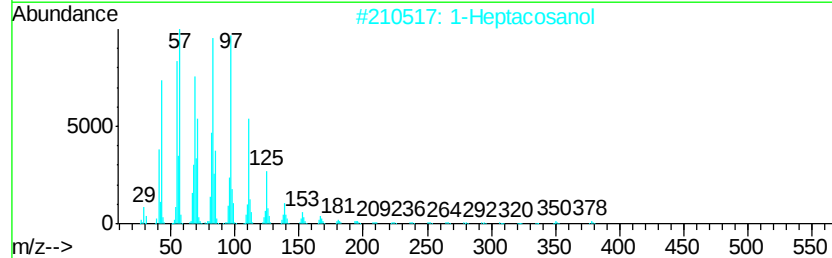
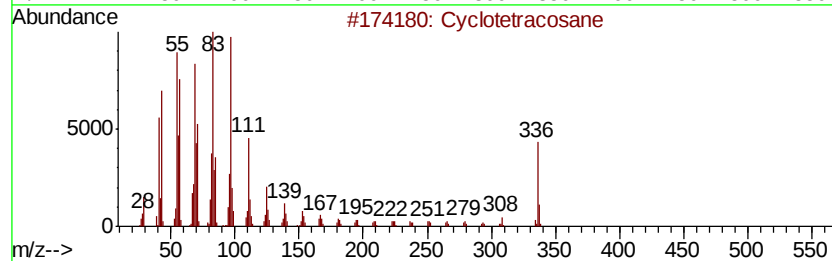
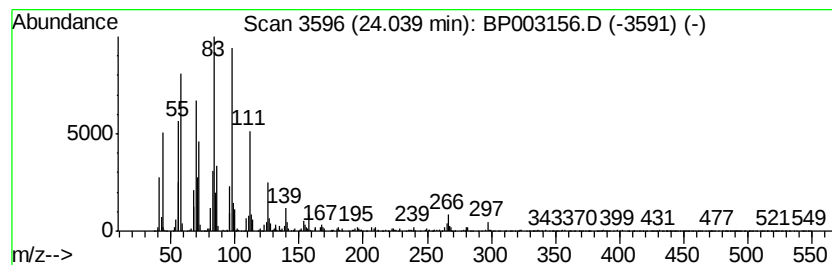
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Cyclotetracosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.04	3.11 ng	615341	Perylene-d12	23.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclotetracosane	336	C24H48	000297-03-0	96
2		1-Heptacosanol	396	C27H56O	002004-39-9	93
3		Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	92
4		9-Nonadecene	266	C19H38	031035-07-1	91
5		1-Docosanol, methyl ether	340	C23H48O	1000333-91-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP082820\
Data File : BP003156.D
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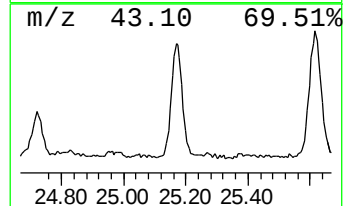
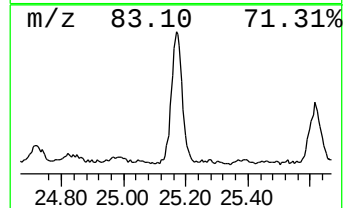
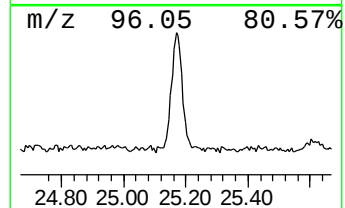
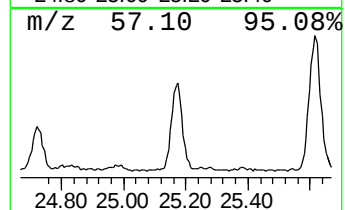
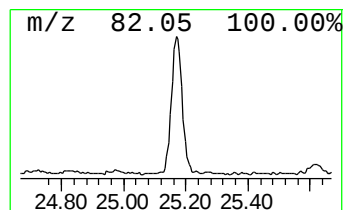
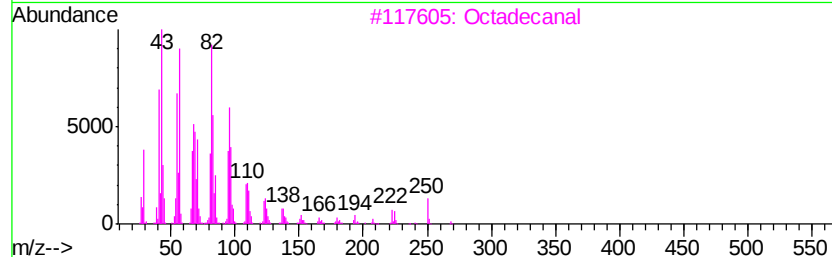
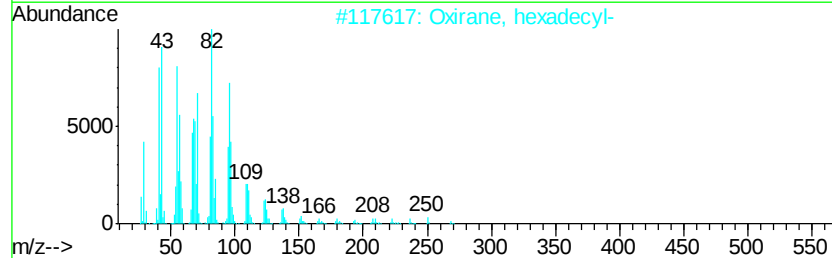
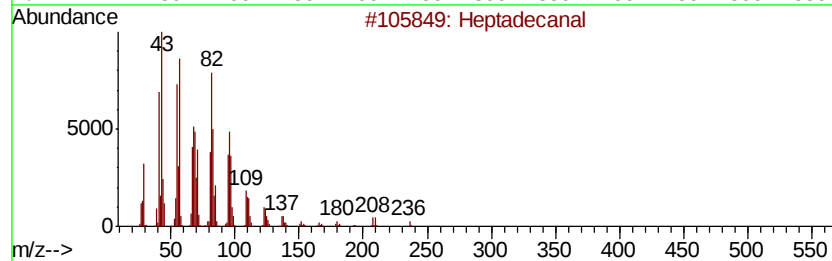
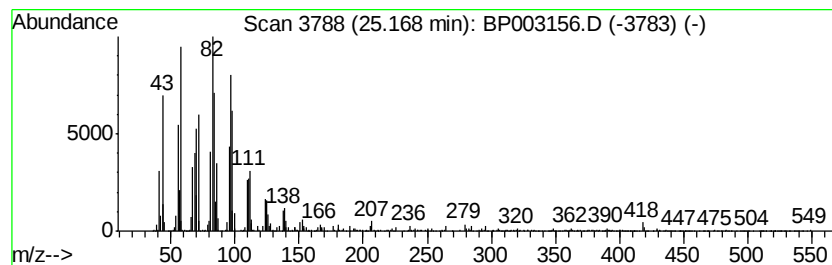
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Heptadecanal Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.17	3.38 ng	669699	Perylene-d12	23.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecanal	254	C17H34O	1000376-70-0	93
2		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
3		Octadecanal	268	C18H36O	000638-66-4	91
4		(Z)-14-Tricosenyl formate	366	C24H46O2	077899-10-6	90
5		Hexadecanal	240	C16H32O	000629-80-1	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP082820\
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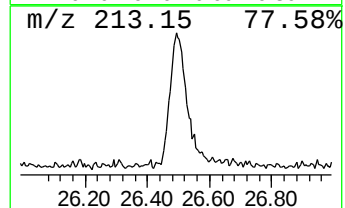
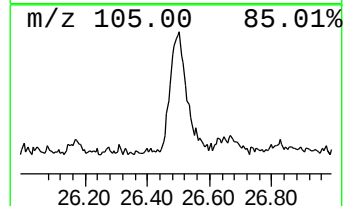
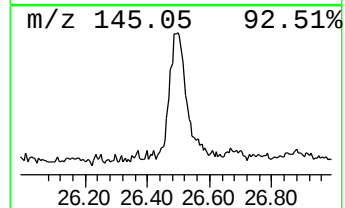
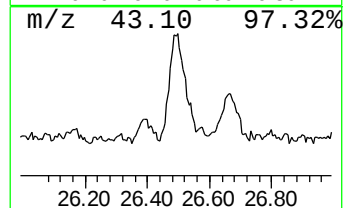
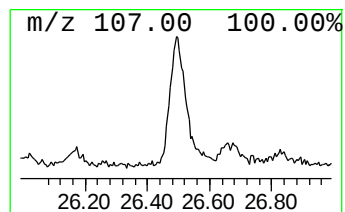
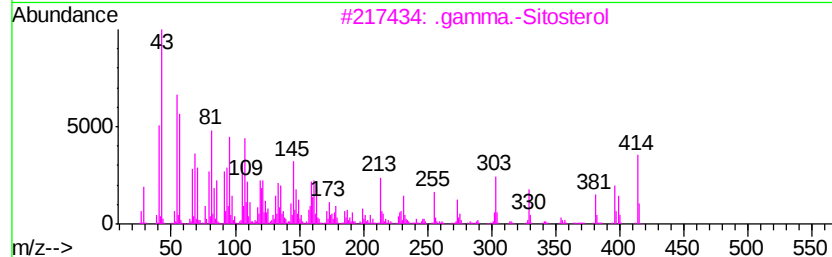
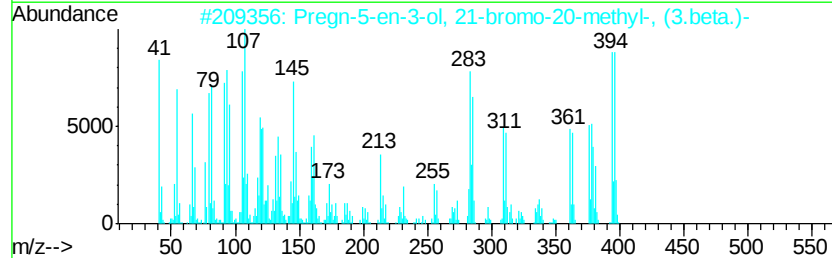
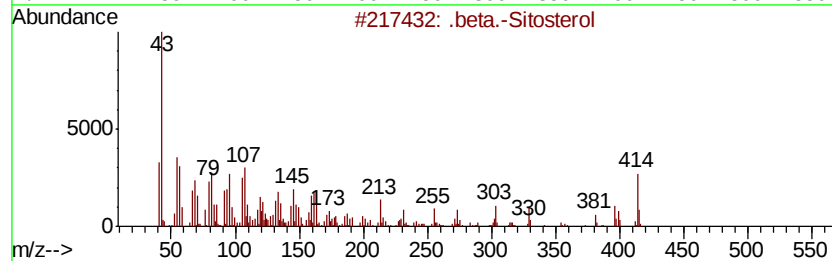
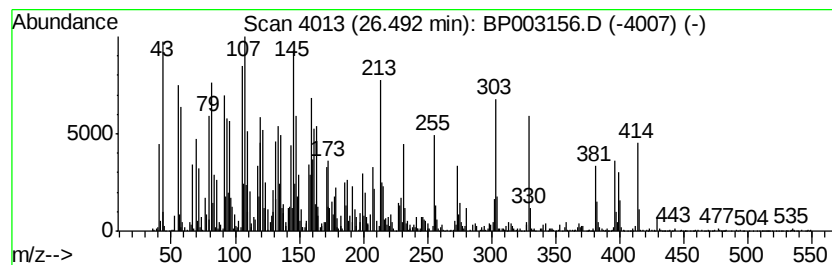
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ClientSampleId :
789UMR-TP11-0006-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 .beta.-Sitosterol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
26.49	3.93 ng	777499	Perylene-d12		23.38	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		.beta.-Sitosterol	414	C29H50O	000083-46-5	83
2		Pregn-5-en-3-ol, 21-bromo-20-met...	394	C22H35BrO	055103-80-5	81
3		.gamma.-Sitosterol	414	C29H50O	000083-47-6	64
4		Acetamide, 2-[5-(2-methyl-3-fur...	329	C16H15N3O3S	1000350-00-2	11
5		Pyridine, 2-(5-trifluoromethyl-3...	213	C9H6F3N3	003974-71-8	11



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP082820\
Data File : BP003156.D
Acq On : 28 Aug 2020 18:15
Operator : CG/JU
Sample : L3837-13
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
789UMR-TP11-0006-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP082420.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
Butane, 2-methoxy...	2.93	18.8	ng	996999	1	7.66	1060460	20.0
n-Hexadecanoic acid	17.97	3.6	ng	489068	4	17.05	2702580	20.0
Cyclohexadecane	20.88	5.1	ng	1092490	5	21.14	4312720	20.0
Eicosane	21.75	4.4	ng	958333	5	21.14	4312720	20.0
13-Docosen-1-ol, ...	22.44	2.5	ng	497414	6	23.38	3959650	20.0
n-Tetracosanol-1	22.76	4.5	ng	885824	6	23.38	3959650	20.0
Benzo[e]pyrene	23.19	2.6	ng	523653	6	23.38	3959650	20.0
Triacetyl acetate	23.62	3.7	ng	734735	6	23.38	3959650	20.0
Heptadecane	23.96	4.8	ng	943975	6	23.38	3959650	20.0
Cyclotetracosane	24.04	3.1	ng	615341	6	23.38	3959650	20.0
Heptadecanal	25.17	3.4	ng	669699	6	23.38	3959650	20.0
.beta.-Sitosterol	26.49	3.9	ng	777499	6	23.38	3959650	20.0