

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071620\
 Data File : BP002772.D
 Acq On : 16 Jul 2020 15:06
 Operator : CG/JU
 Sample : L3315-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-16-S

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-BP071020.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	5.187	384	391	408	rVB	1996055	3091614	47.29%	7.074%
2	6.764	651	659	677	rBV	2058900	3432133	52.50%	7.853%
3	7.169	723	728	741	rBV	85105	161758	2.47%	0.370%
4	7.558	787	794	803	rBV	612714	946171	14.47%	2.165%
5	8.705	981	989	1002	rBV	1340673	2222316	33.99%	5.085%
6	10.328	1258	1265	1276	rBV	849261	1443189	22.07%	3.302%
7	12.822	1682	1689	1706	rBV	3101292	4835864	73.97%	11.065%
8	13.122	1735	1740	1748	rBV2	46237	75580	1.16%	0.173%
9	13.669	1827	1833	1844	rBV	357297	535909	8.20%	1.226%
10	14.075	1895	1902	1913	rBV2	59459	151620	2.32%	0.347%
11	14.210	1918	1925	1936	rBV2	1344917	2017961	30.87%	4.617%
12	14.669	1998	2003	2016	rBV2	62039	125582	1.92%	0.287%
13	15.716	2169	2181	2191	rBV	2363000	3728749	57.03%	8.532%
14	16.022	2221	2233	2244	rBV2	147107	281221	4.30%	0.643%
15	16.375	2288	2293	2297	rBV3	124392	227723	3.48%	0.521%
16	16.534	2316	2320	2332	rVB	53605	94017	1.44%	0.215%
17	16.969	2388	2394	2403	rBV	1671369	2345126	35.87%	5.366%
18	17.898	2547	2552	2566	rBV	439284	779866	11.93%	1.784%
19	18.151	2591	2595	2600	rBV	398902	589170	9.01%	1.348%
20	18.198	2600	2603	2618	rVB	213586	376818	5.76%	0.862%
21	18.633	2672	2677	2681	rBV	378418	469126	7.18%	1.073%
22	18.722	2688	2692	2699	rBV	61815	86052	1.32%	0.197%
23	19.557	2829	2834	2846	rBV	4958522	6538003	100.00%	14.960%
24	20.222	2942	2947	2957	rBV2	186766	413491	6.32%	0.946%
25	20.345	2965	2968	2970	rBV	62716	71588	1.09%	0.164%
26	20.445	2982	2985	2988	rBV	60073	70294	1.08%	0.161%
27	20.810	3043	3047	3057	rBV	817464	1175632	17.98%	2.690%
28	21.075	3087	3092	3098	rBV	1992759	2626157	40.17%	6.009%
29	21.245	3118	3121	3130	rBV	179972	214671	3.28%	0.491%
30	21.674	3190	3194	3204	rVB	260938	393903	6.02%	0.901%
31	22.127	3267	3271	3276	rBV	282872	364161	5.57%	0.833%
32	22.251	3288	3292	3297	rVB	211593	284015	4.34%	0.650%
33	22.627	3352	3356	3361	rVB	295792	409588	6.26%	0.937%
34	23.186	3447	3451	3458	rVB	252682	402411	6.15%	0.921%

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ClientSampleId :
TP-16-S

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-BP071020.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	23.269	3459	3465	3474	rVB2	1299025	2424940	37.09%	5.549%
36	23.821	3556	3559	3566	rVB	182272	296067	4.53%	0.677%

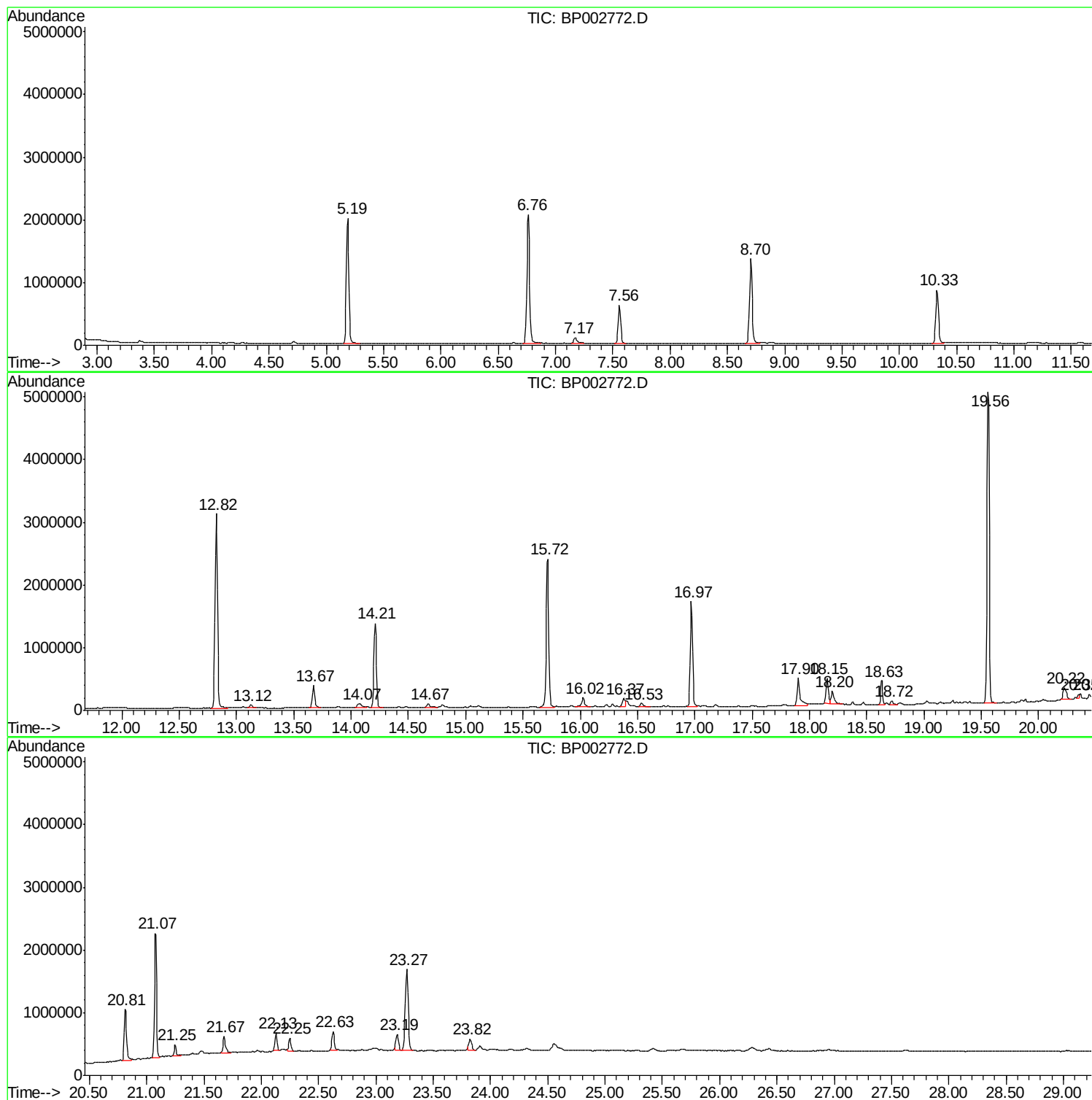
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Sample : L3315-03
Misc :
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Instrument :
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ClientSampleId :
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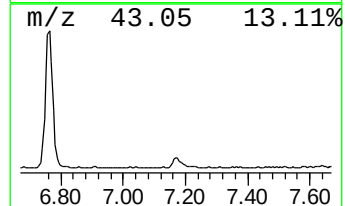
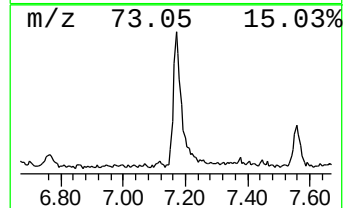
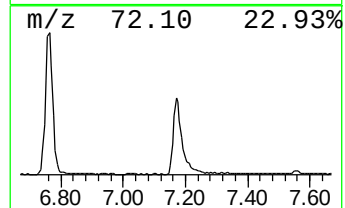
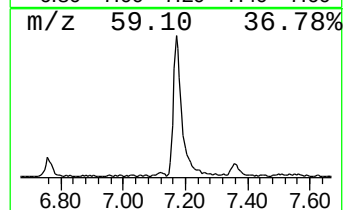
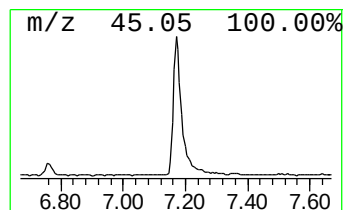
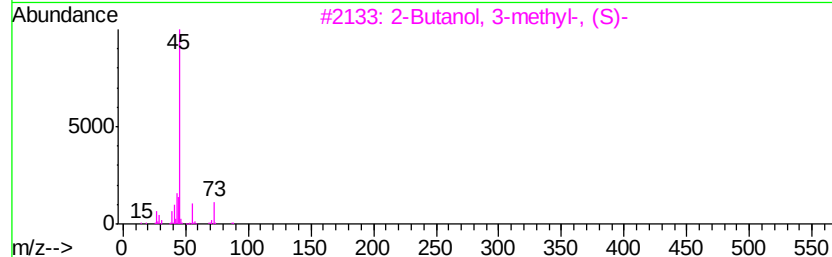
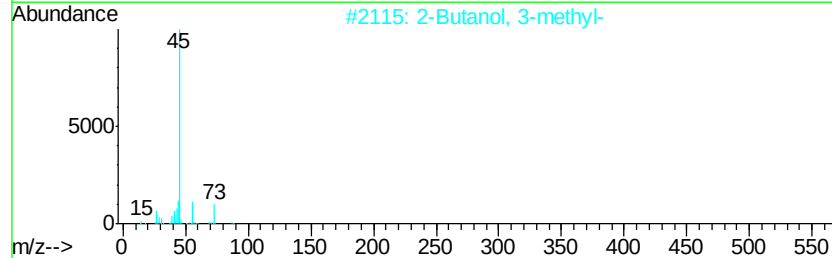
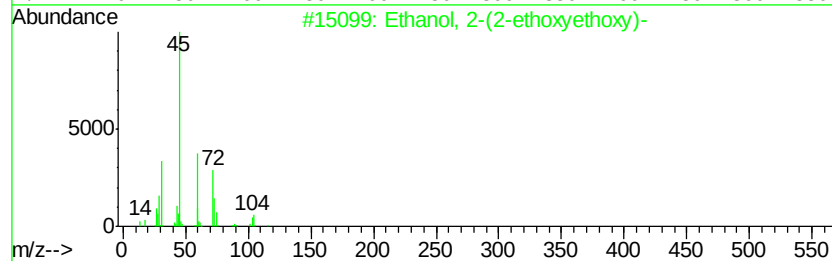
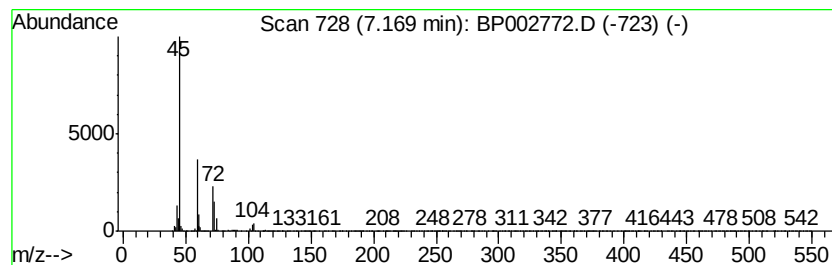
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.17	3.42 ng	161758	1,4-Dichlorobenzene-d4	7.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	90
2		2-Butanol, 3-methyl-	88	C5H12O	000598-75-4	47
3		2-Butanol, 3-methyl-, (S)-	88	C5H12O	001517-66-4	47
4		2-Propanol, 1-(1-methylethoxy)-	118	C6H14O2	003944-36-3	45
5		Ethanol, 2-[2-(2-ethoxyethoxy)et...]	178	C8H18O4	000112-50-5	43



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Misc :
ALS Vial : 6 Sample Multiplier: 1

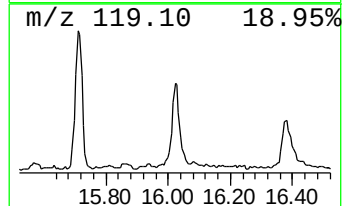
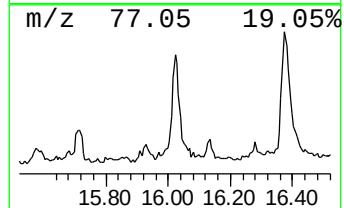
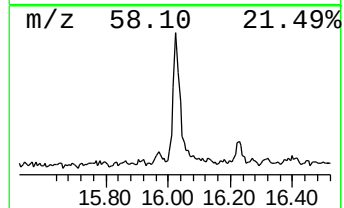
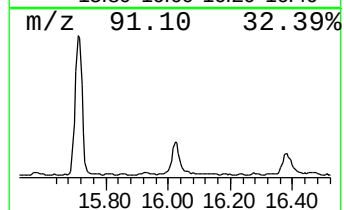
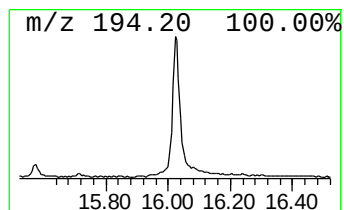
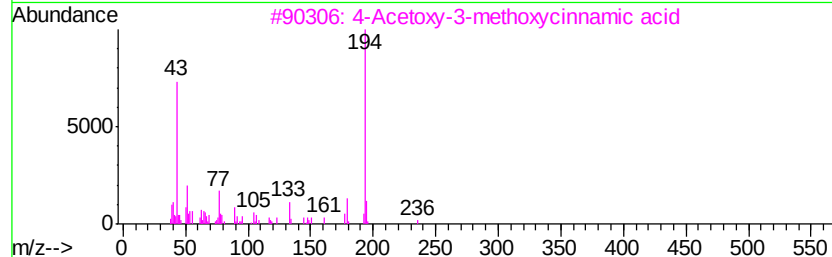
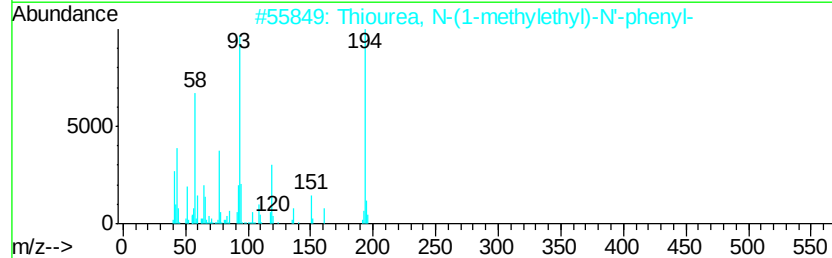
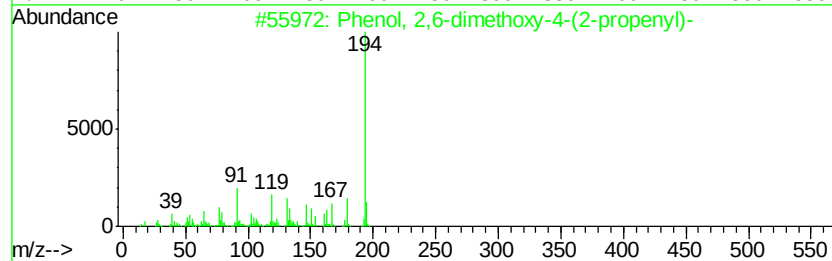
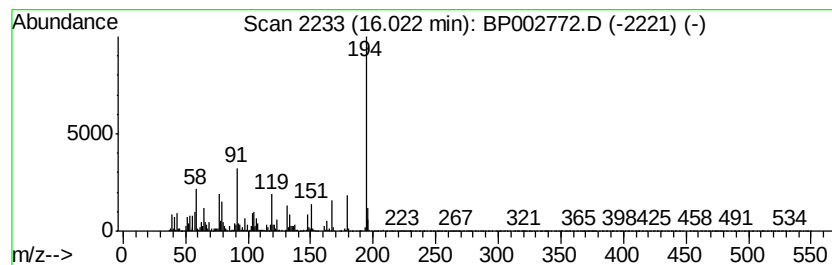
Instrument :
BNA_P
ClientSampled :
TP-16-S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Phenol, 2,6-dimethoxy-4-(2-... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.02	2.40 ng	281221	Phenanthrene-d10	16.97		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,6-dimethoxy-4-(2-prope...	194	C11H14O3	006627-88-9	89	
2	Thiourea, N-(1-methylethyl)-N'-p...	194	C10H14N2S	015093-36-4	53	
3	4-Acetoxy-3-methoxycinnamic acid	236	C12H12O5	002596-47-6	47	
4	2,5-Dimethoxyterephthalic acid	194	C10H10O4	007310-97-6	47	
5	5,7-Dimethyl-2-methylsulfanyl-[1...	194	C8H10N4S	1000300-84-3	43	



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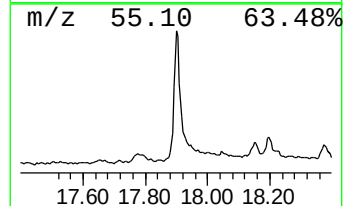
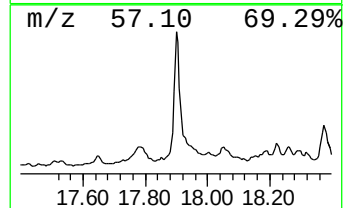
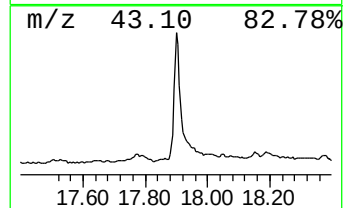
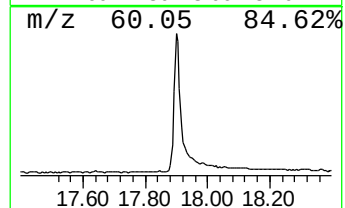
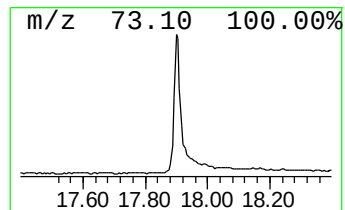
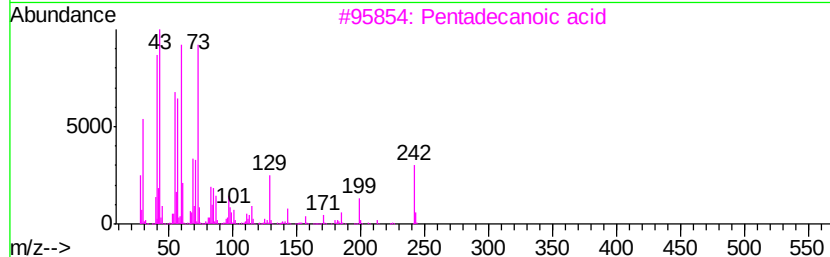
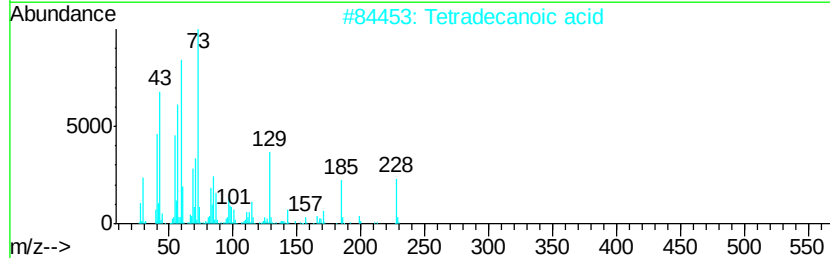
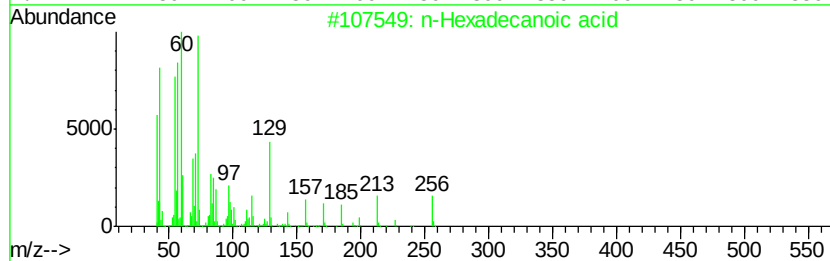
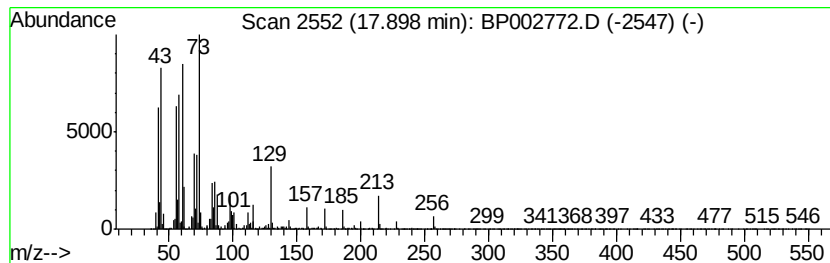
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.90	6.65 ng	779866	Phenanthrene-d10	16.97

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	93
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	86
4		Tridecanoic acid	214	C13H26O2	000638-53-9	83
5		n-Decanoic acid	172	C10H20O2	000334-48-5	59



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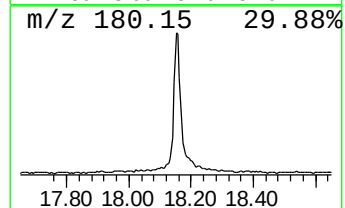
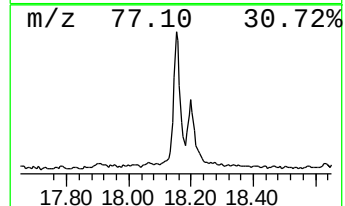
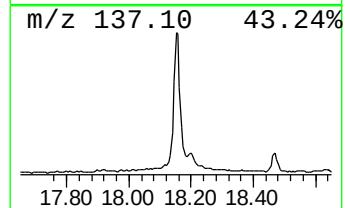
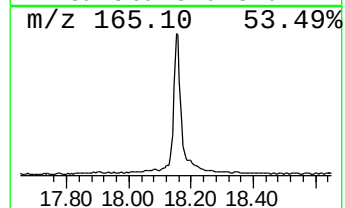
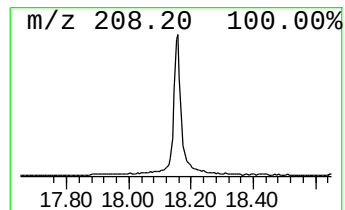
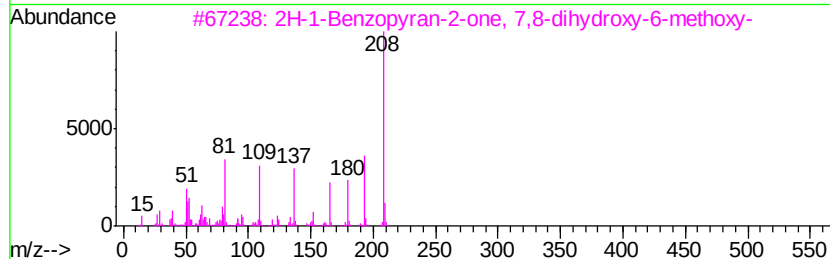
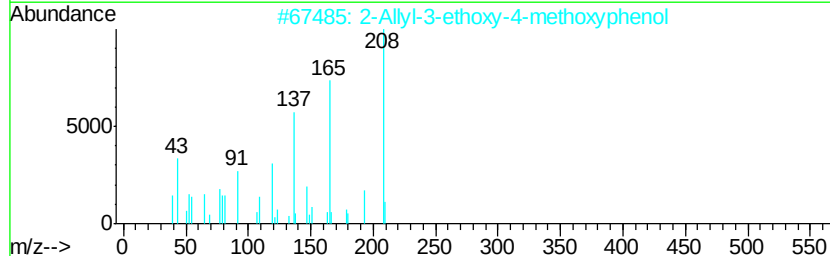
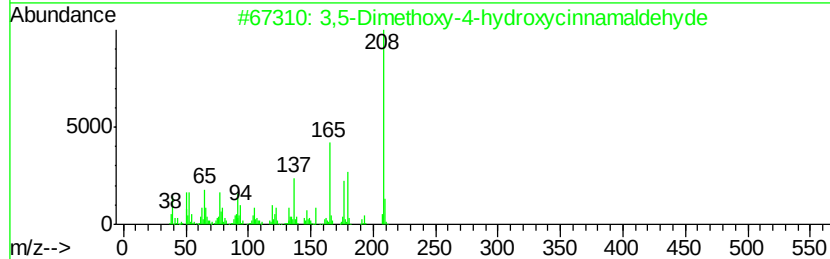
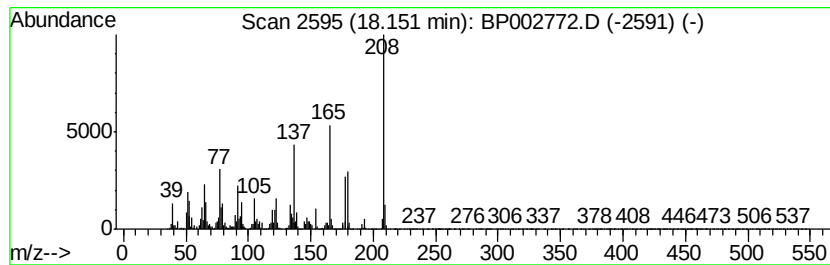
Instrument :
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 3,5-Dimethoxy-4-hydroxycinn... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.15	5.02 ng	589170	Phenanthrene-d10	16.97	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-Dimethoxy-4-hydroxycinnamaldehyde	208	C11H12O4	087345-53-7	97
2	2-Allyl-3-ethoxy-4-methoxyphenol	208	C12H16O3	1000122-70-6	58
3	2H-1-Benzopyran-2-one, 7,8-dihydroxy-	208	C10H8O5	000574-84-5	58
4	2-Propenoic acid, 3-(2,3-dimethoxy-4-hydroxyphenyl)-	208	C11H12O4	007345-82-6	45
5	1,2-Dimethoxy-4-(3-methoxy-1-propenyl)benzene	208	C12H16O3	1000313-35-0	43



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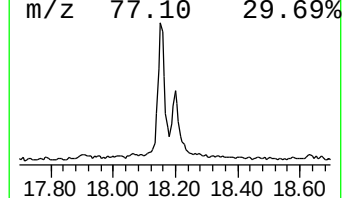
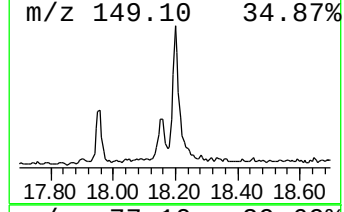
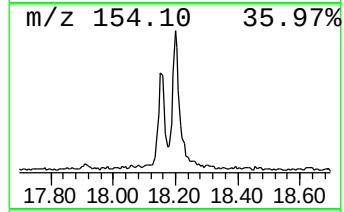
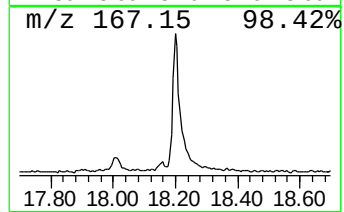
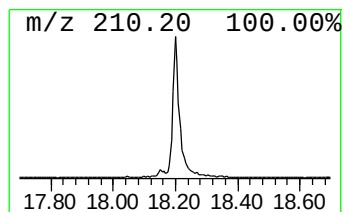
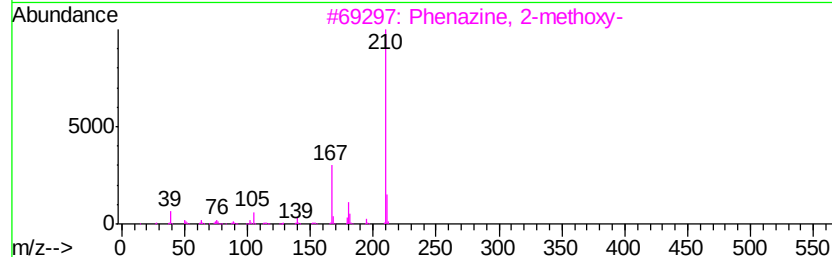
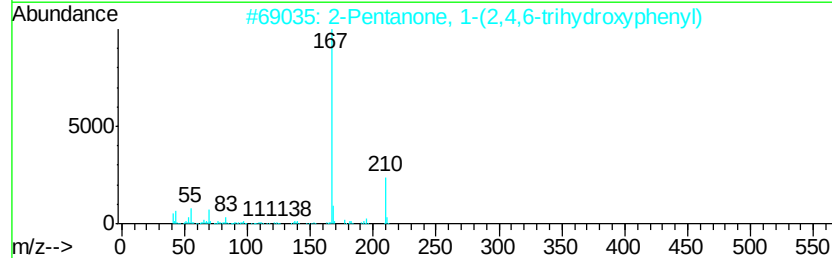
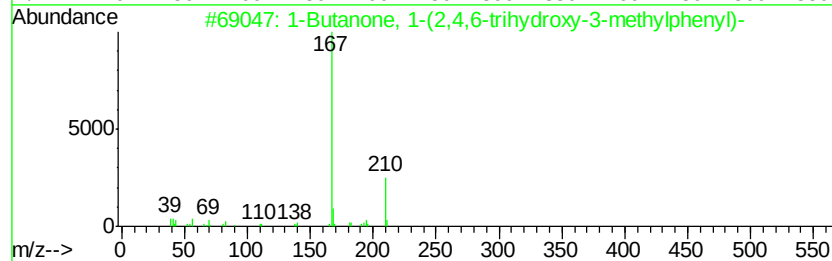
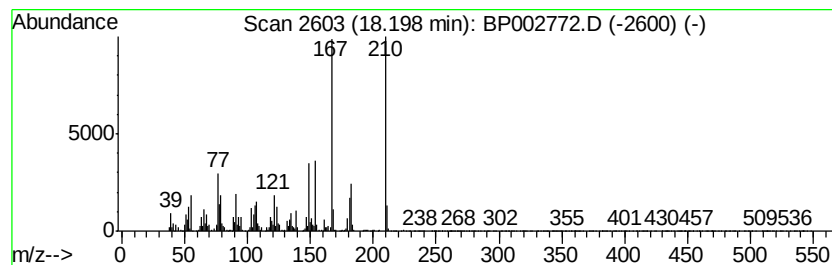
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 1-Butanone, 1-(2,4,6-trihyd... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.20	3.21 ng	376818	Phenanthrene-d10	16.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Butanone, 1-(2,4,6-trihydroxyv...	210	C11H14O4	001509-06-4	52
2			2-Pentanone, 1-(2,4,6-trihydroxy...	210	C11H14O4	1000116-22-3	50
3			Phenazine, 2-methoxy-	210	C13H10N2O	002876-18-8	38
4			5-tert-Butylpyroqallol	182	C10H14O3	020481-17-8	38
5			Imidazo[4,5-e][1,4]diazepin-8(1H...	210	C8H10N4OS	1000142-41-3	38



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Acq On : 16 Jul 2020 15:06
Operator : CG/JU
Sample : L3315-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

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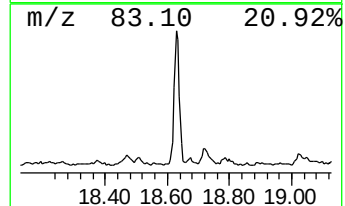
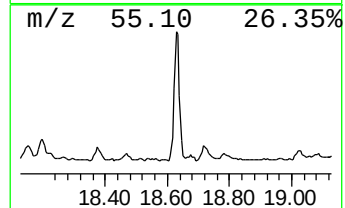
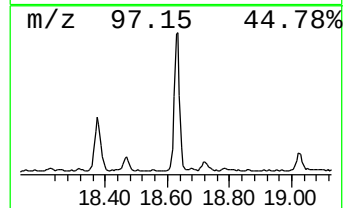
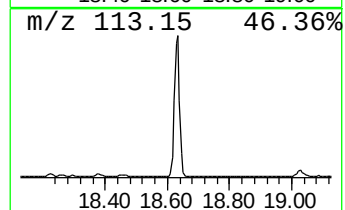
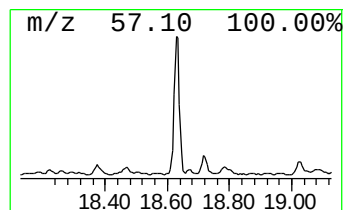
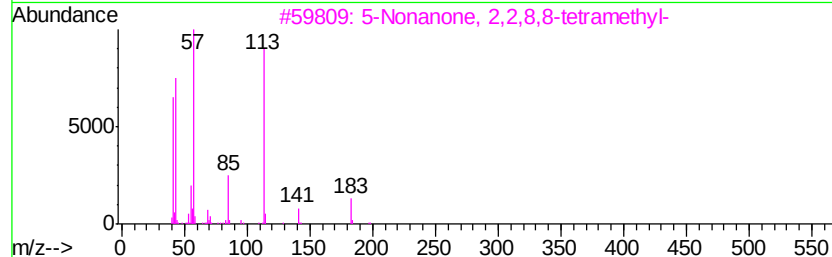
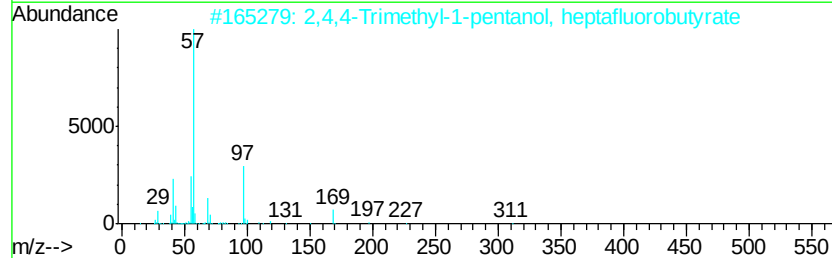
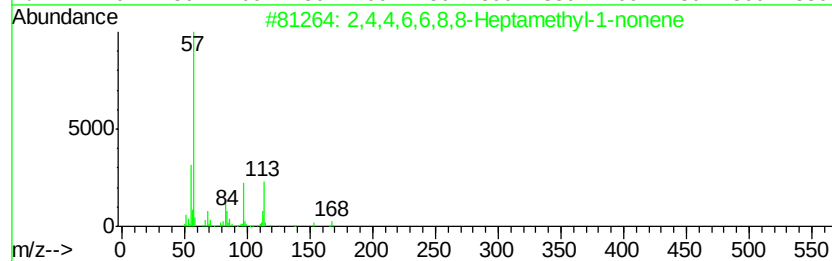
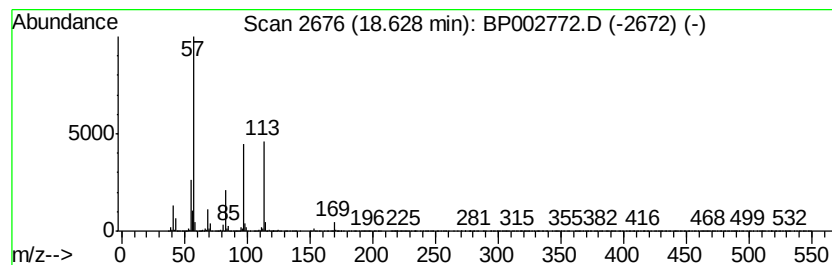
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown18.63 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.63	4.00 ng	469126	Phenanthrene-d10	16.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	38		
2	2,4,4-Trimethyl-1-pentanol, hept...	326	C12H17F7O2	1000365-01-4	37		
3	5-Nonanone, 2,2,8,8-tetramethyl-	198	C13H26O	005709-95-5	35		
4	1-Pentanol, 3-methyl-2-propyl-	144	C9H20O	054004-40-9	33		
5	2,4,4-Trimethyl-1-pentanol, pent...	276	C11H17F5O2	1000365-51-0	32		



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071620\
Data File : BP002772.D
Acq On : 16 Jul 2020 15:06
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ClientSampleId :
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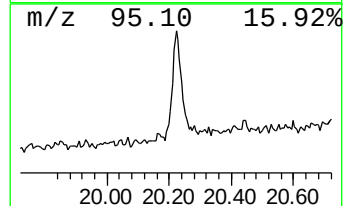
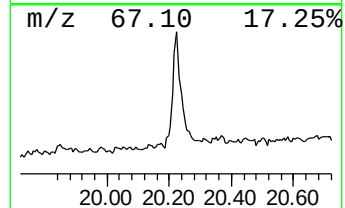
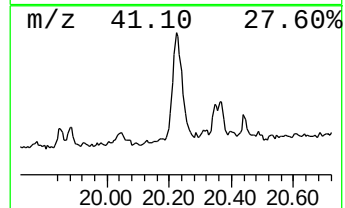
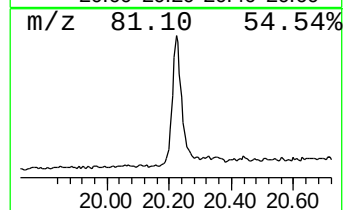
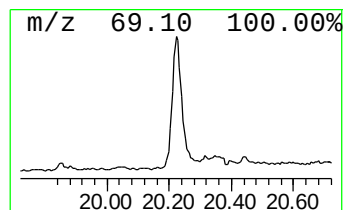
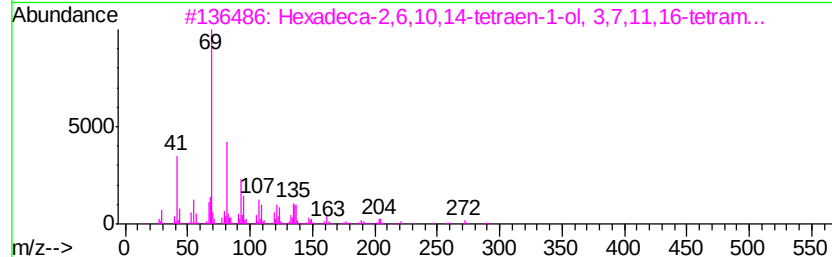
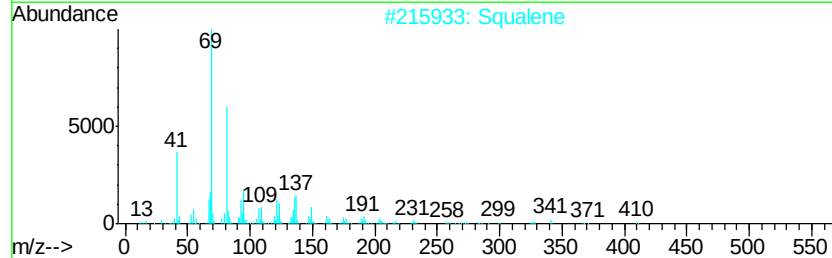
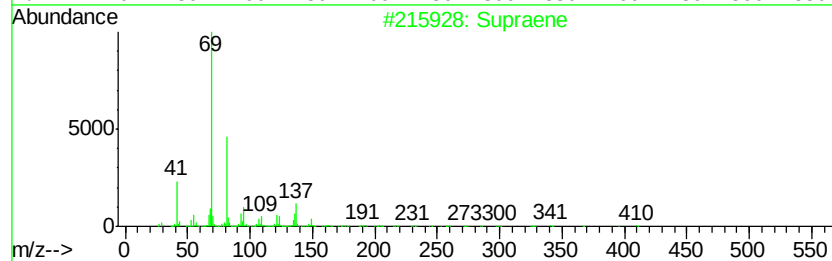
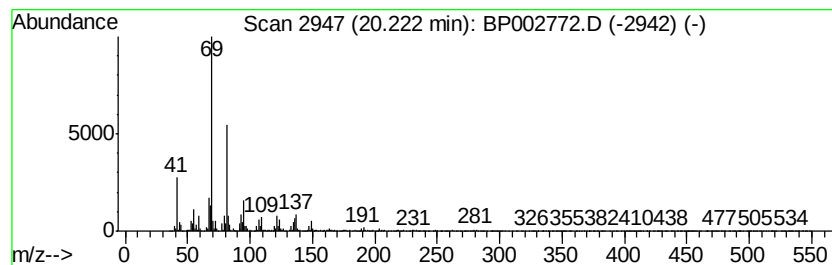
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Supraene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.22	3.15 ng	413491	Chrysene-d12	21.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Supraene	410	C30H50	007683-64-9	91
2		Squalene	410	C30H50	000111-02-4	90
3		Hexadeca-2,6,10,14-tetraen-1-ol,...	290	C20H34O	007614-21-3	72
4		2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	64
5		Bis(3-methylbut-3-enyl) phthalate	302	C18H22O4	022711-53-1	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071620\
Data File : BP002772.D
Acq On : 16 Jul 2020 15:06
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Sample : L3315-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

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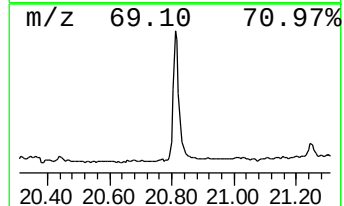
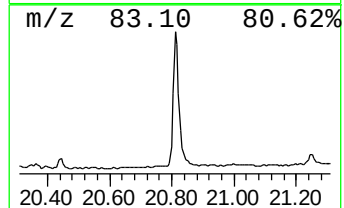
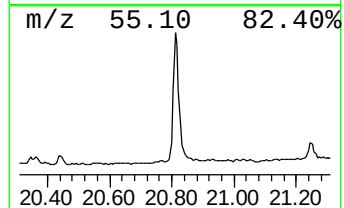
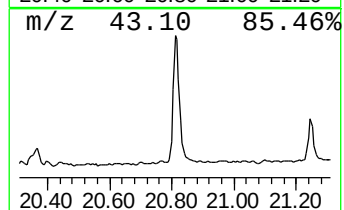
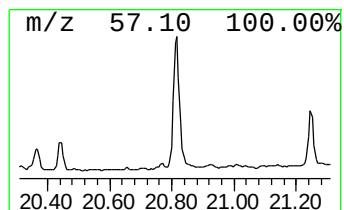
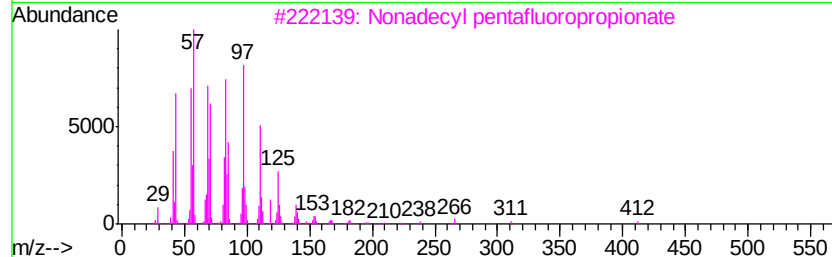
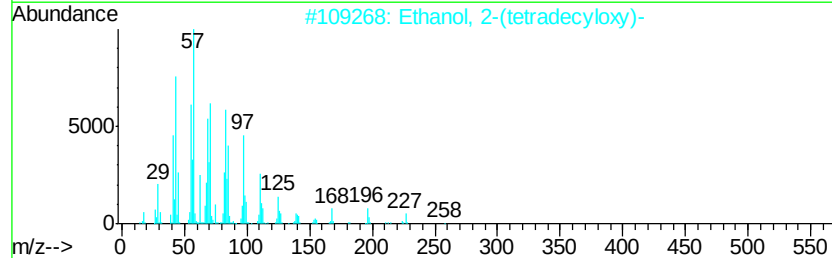
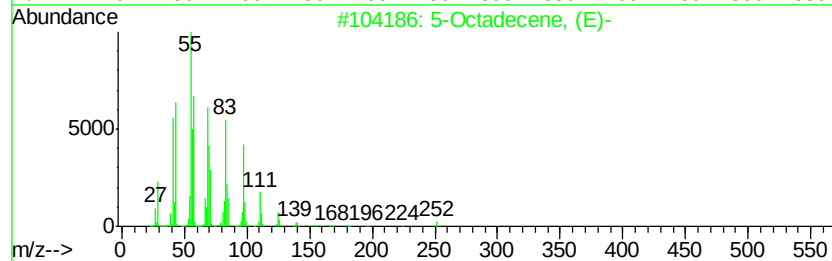
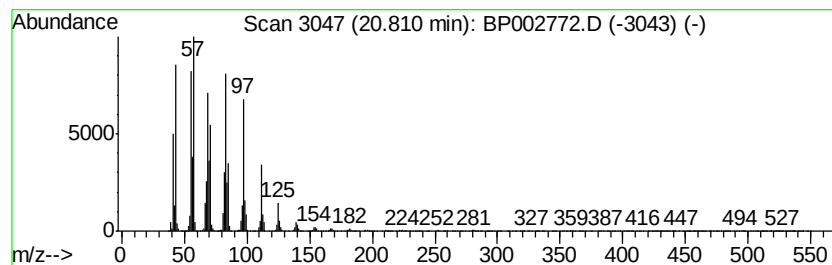
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 5-Octadecene, (E)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.81	8.95 ng	1175630	Chrysene-d12	21.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Octadecene, (E)-	252	C18H36	007206-21-5	93
2		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	93
3		Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
4		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	91
5		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071620\
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 Sample : L3315-03
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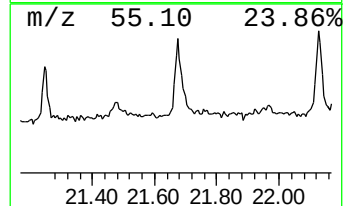
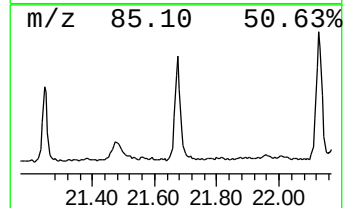
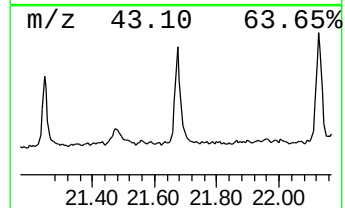
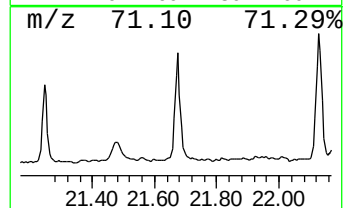
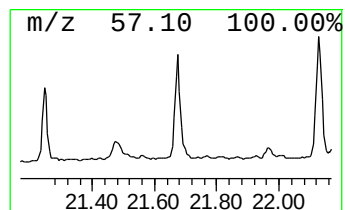
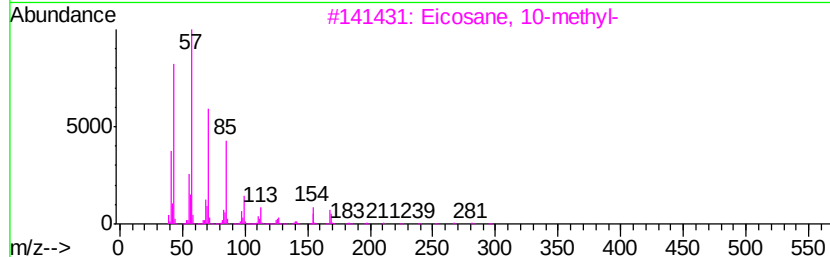
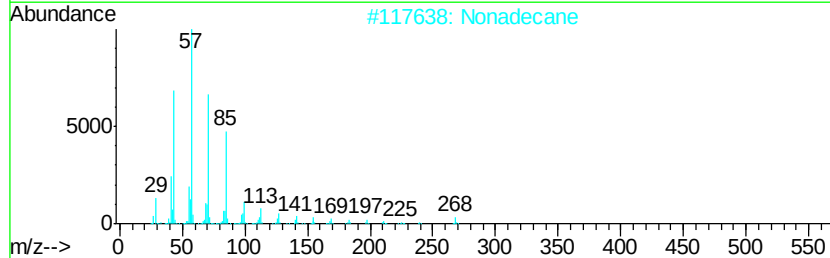
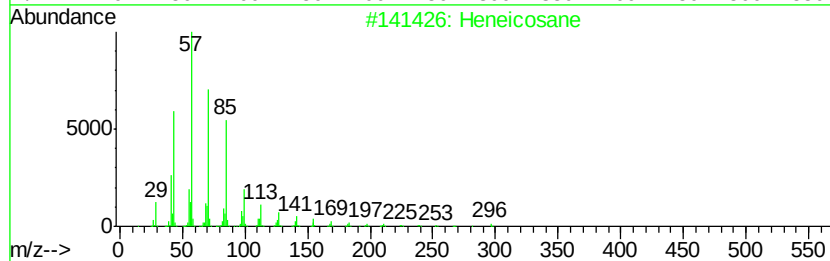
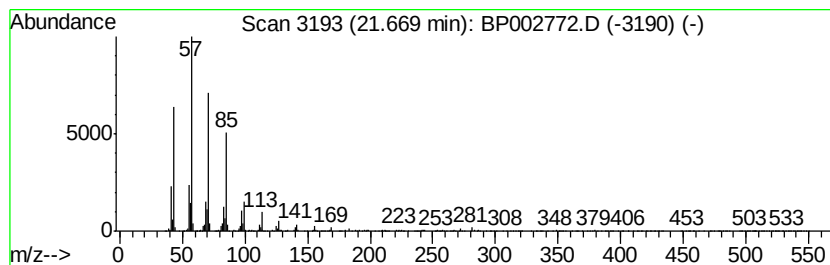
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 Heneicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.67	3.00 ng	393903	Chrysene-d12	21.07	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	97
2	Nonadecane	268	C19H40	000629-92-5	94
3	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4	Hexacosane	366	C26H54	000630-01-3	91
5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071620\
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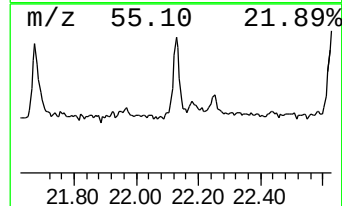
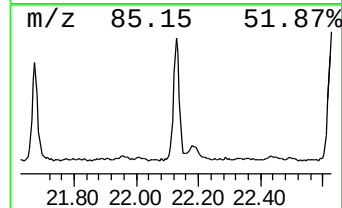
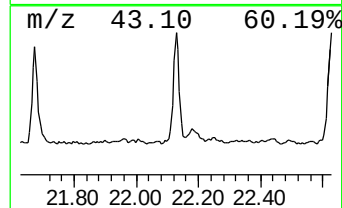
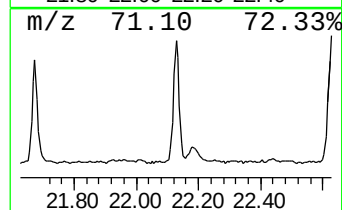
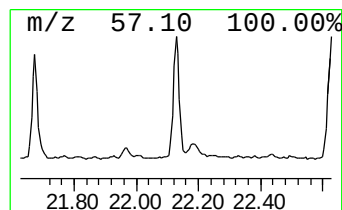
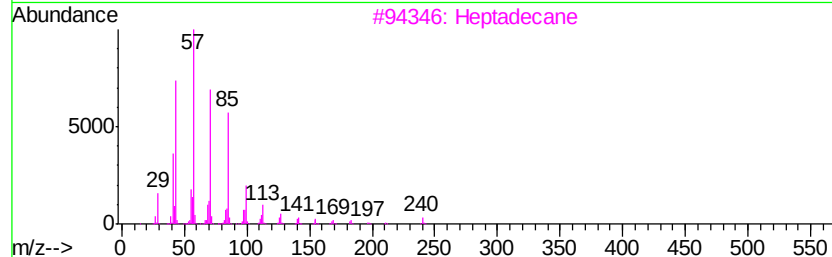
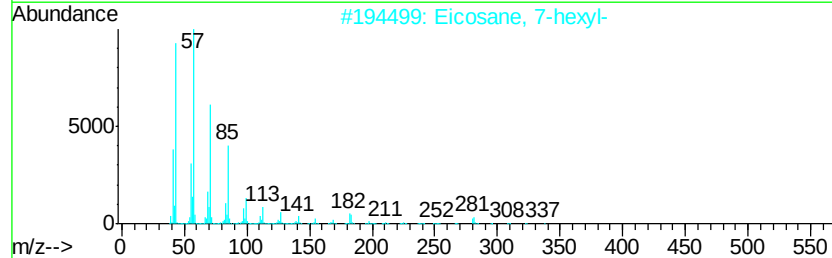
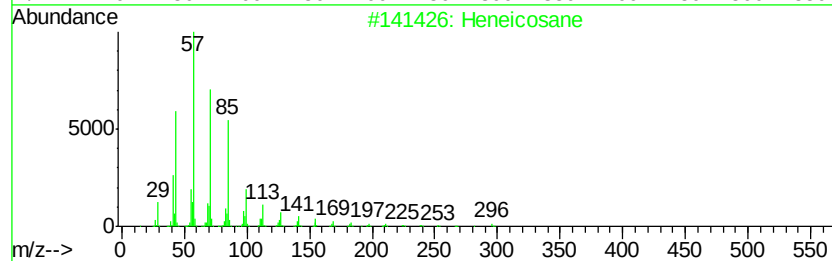
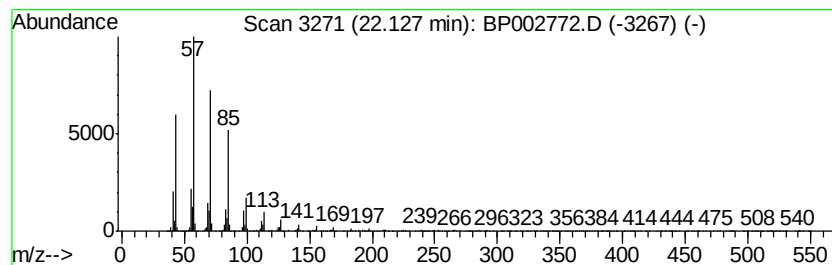
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Eicosane, 7-hexyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
22.13	2.77 ng	364161	Chrysene-d12		21.07	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heneicosane	296	C21H44	000629-94-7	91
2		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90
3		Heptadecane	240	C17H36	000629-78-7	86
4		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	83
5		Octacosane	394	C28H58	000630-02-4	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071620\
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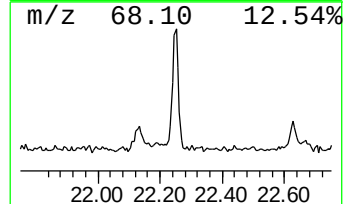
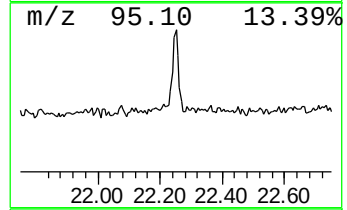
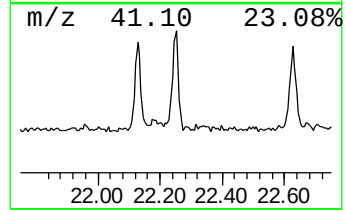
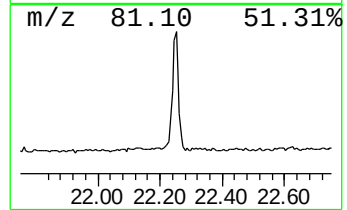
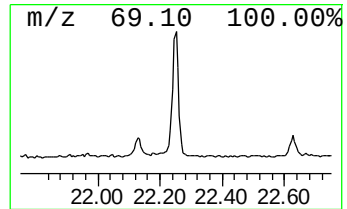
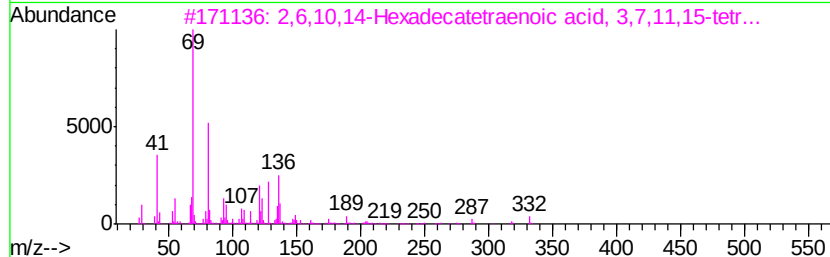
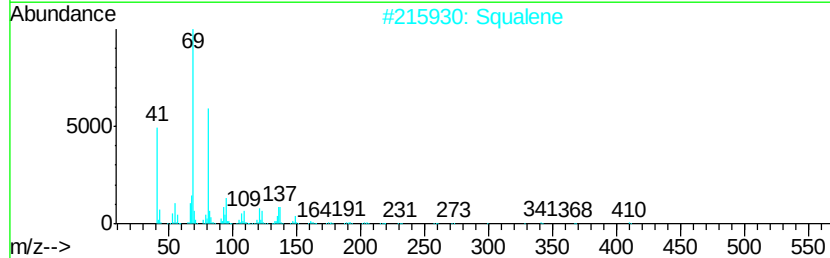
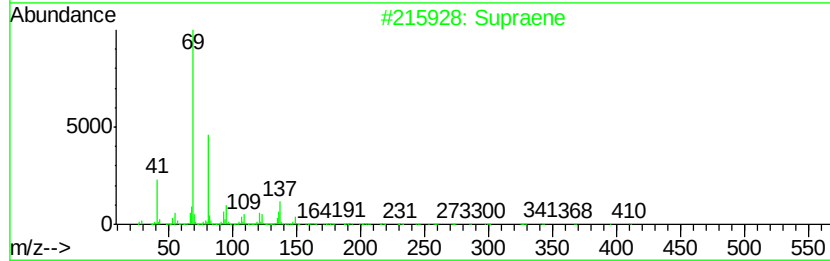
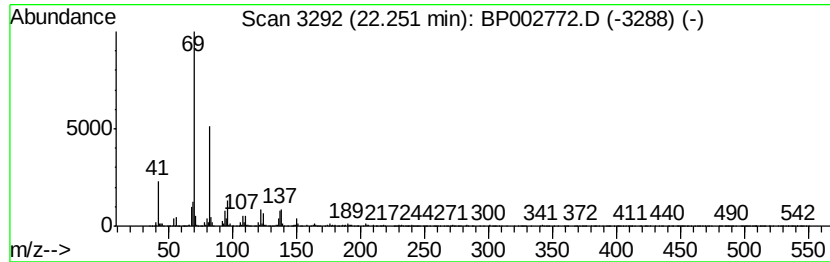
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Squalene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.25	2.34 ng	284015	Perylene-d12	23.27
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Supraene		410 C30H50	007683-64-9 91
2	Squalene		410 C30H50	000111-02-4 91
3	2,6,10,14-Hexadecatetraenoic aci...		332 C22H36O2	024035-35-6 87
4	Farnesol isomer a		222 C15H26O	1000108-92-4 72
5	2,6,10-Dodecatrien-1-ol, 3,7,11-...		264 C17H28O2	004128-17-0 72



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071620\
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ALS Vial : 6 Sample Multiplier: 1

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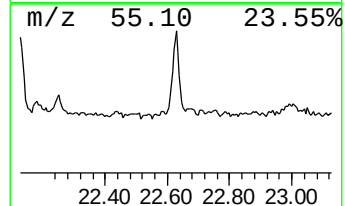
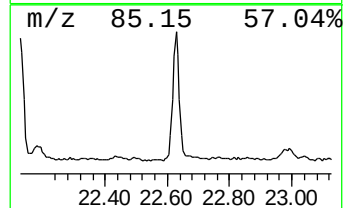
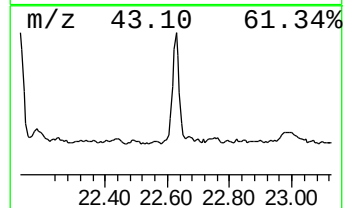
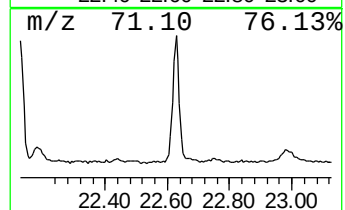
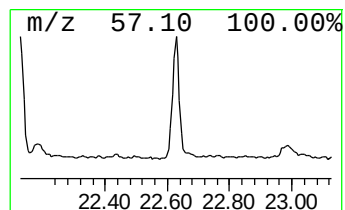
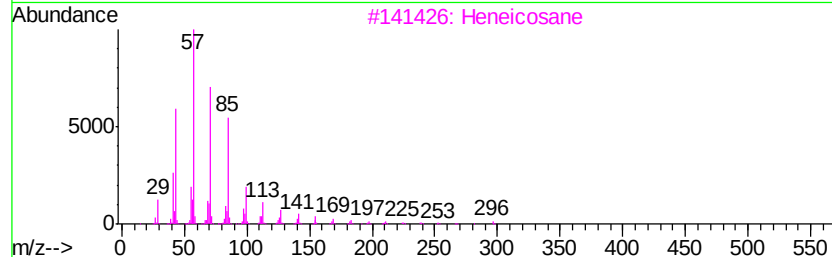
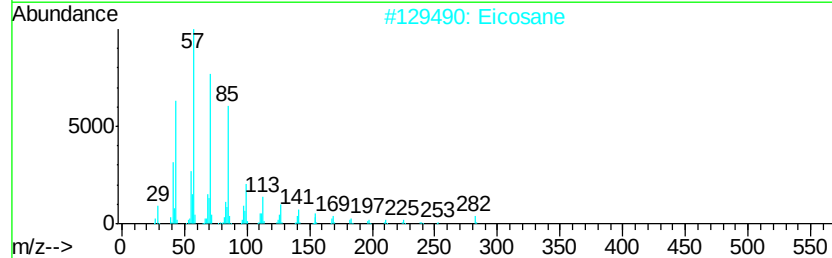
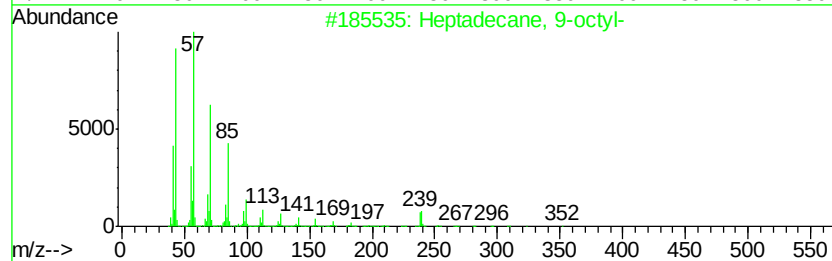
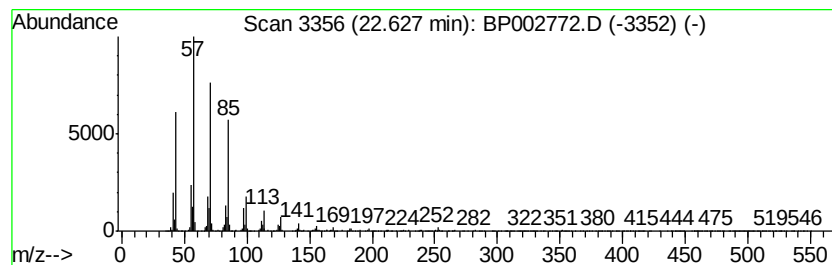
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Heptadecane, 9-octyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.63	3.38 ng	409588	Perylene-d12	23.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
2		Eicosane	282	C20H42	000112-95-8	94
3		Heneicosane	296	C21H44	000629-94-7	91
4		Docosane, 7-hexyl-	394	C28H58	055373-86-9	90
5		Tetracosane	338	C24H50	000646-31-1	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071620\
Data File : BP002772.D
Acq On : 16 Jul 2020 15:06
Operator : CG/JU
Sample : L3315-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP-16-S

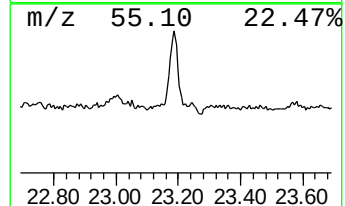
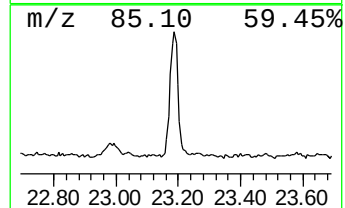
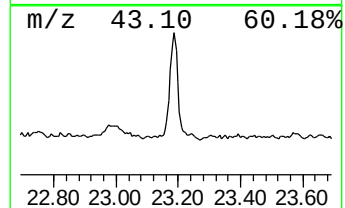
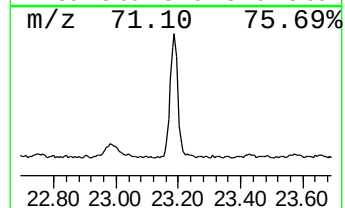
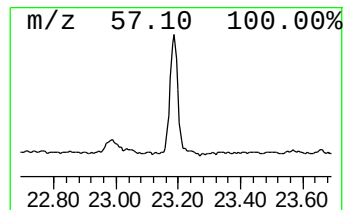
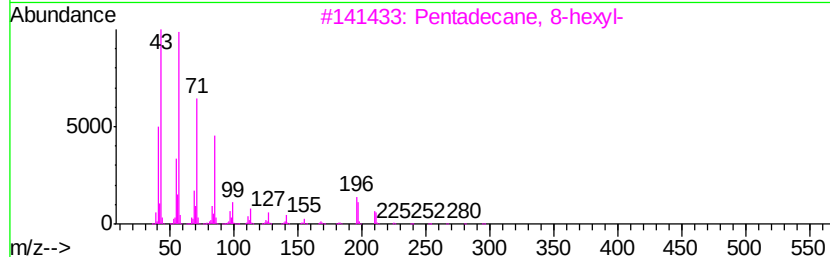
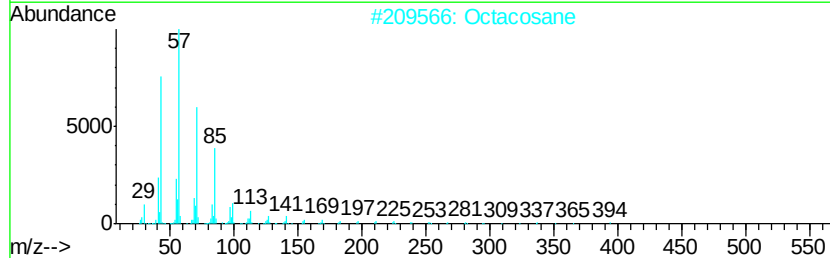
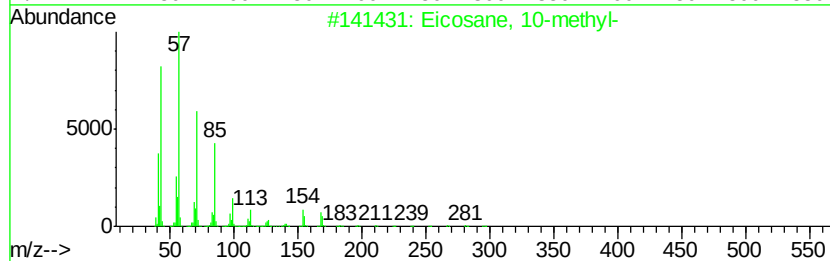
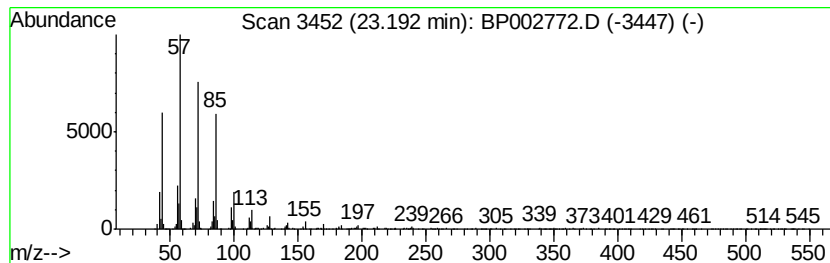
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Eicosane, 10-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.19	3.32 ng	402411	Perylene-d12	23.27

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		Tetracosane	338	C24H50	000646-31-1	91
5		Nonadecane, 2-methyl-	282	C20H42	001560-86-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071620\
Data File : BP002772.D
Acq On : 16 Jul 2020 15:06
Operator : CG/JU
Sample : L3315-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP-16-S

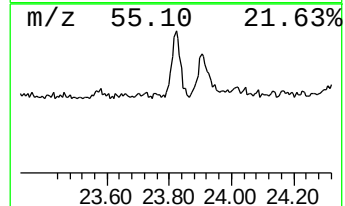
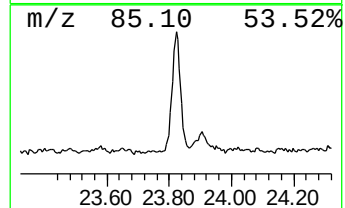
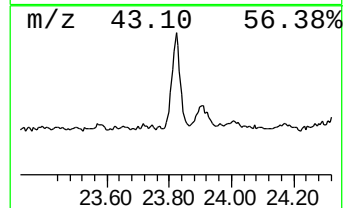
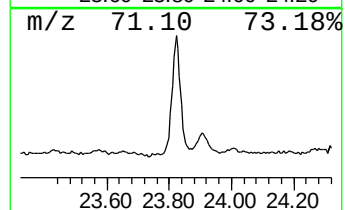
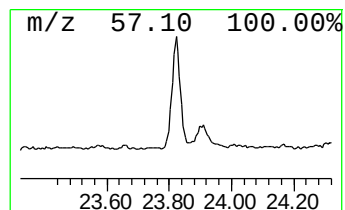
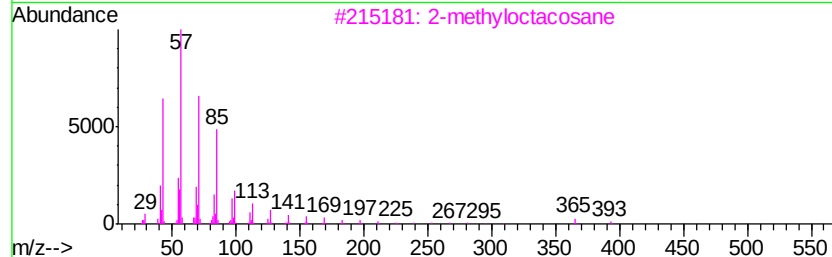
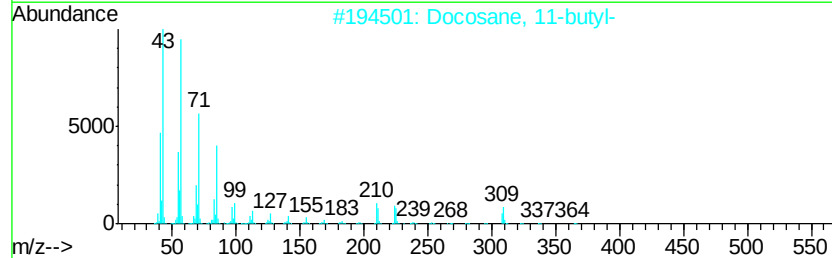
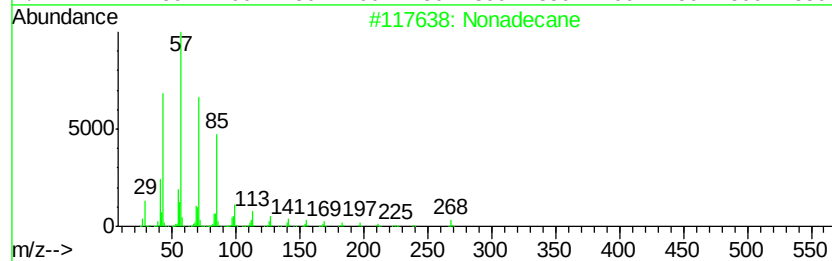
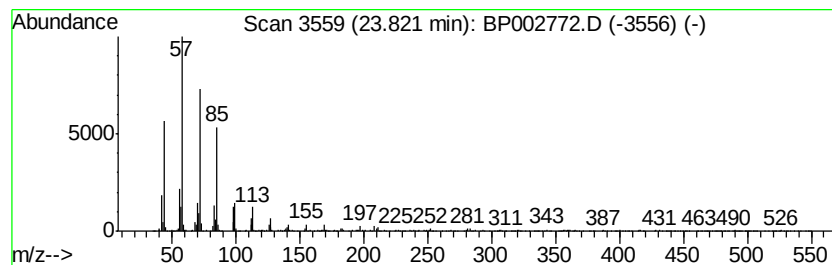
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Nonadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.82	2.44 ng	296067	Perylene-d12	23.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	94
2		Docosane, 11-butyl-	366	C26H54	013475-76-8	94
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	90
5		Pentatriacontane	493	C35H72	000630-07-9	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP071620\
 Data File : BP002772.D
 Acq On : 16 Jul 2020 15:06
 Operator : CG/JU
 Sample : L3315-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-16-S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP071020.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
Ethanol, 2-(2-eth...	7.17	3.4	ng	161758	1	7.56	946171	20.0
Phenol, 2,6-dimet...	16.02	2.4	ng	281221	4	16.97	2345130	20.0
n-Hexadecanoic acid	17.90	6.7	ng	779866	4	16.97	2345130	20.0
3,5-Dimethoxy-4-h...	18.15	5.0	ng	589170	4	16.97	2345130	20.0
1-Butanone, 1-(2,...	18.20	3.2	ng	376818	4	16.97	2345130	20.0
unknown18.63	18.63	4.0	ng	469126	4	16.97	2345130	20.0
Supraene	20.22	3.1	ng	413491	5	21.07	2626160	20.0
5-Octadecene, (E)-	20.81	8.9	ng	1175630	5	21.07	2626160	20.0
Heneicosane	21.67	3.0	ng	393903	5	21.07	2626160	20.0
Eicosane, 7-hexyl-	22.13	2.8	ng	364161	5	21.07	2626160	20.0
Squalene	22.25	2.3	ng	284015	6	23.27	2424940	20.0
Heptadecane, 9-oc...	22.63	3.4	ng	409588	6	23.27	2424940	20.0
Eicosane, 10-methyl-	23.19	3.3	ng	402411	6	23.27	2424940	20.0
Nonadecane	23.82	2.4	ng	296067	6	23.27	2424940	20.0