

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042817\
 Data File : BF094687.D
 Acq On : 28 Apr 2017 16:55
 Operator : SJ/MA
 Sample : I2895-10
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-08-2.0-3.0

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:\HPCHEM1\BNA_F\METHODS\8270-BF041717.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	3.942	337	341	351	rBV	473730	571185	17.90%	2.075%
2	4.331	403	407	421	rBV	3033040	3096643	97.03%	11.248%
3	4.713	469	472	477	rVB	38558	39602	1.24%	0.144%
4	5.401	584	589	592	rBV	2718050	3076109	96.39%	11.173%
5	5.525	605	610	614	rBV	3140943	3191421	100.00%	11.592%
6	5.772	648	652	659	rVB	556357	501012	15.70%	1.820%
7	5.948	677	682	685	rBV	2565877	2366411	74.15%	8.595%
8	6.425	757	763	766	rBV	1725332	1940884	60.82%	7.050%
9	7.260	901	905	912	rVB	766879	655825	20.55%	2.382%
10	8.583	1125	1130	1133	rBV	2870972	2710897	84.94%	9.847%
11	9.083	1211	1215	1224	rBV	270074	266430	8.35%	0.968%
12	9.395	1260	1268	1277	rVB2	693175	692170	21.69%	2.514%
13	9.989	1365	1369	1373	rBV	47930	47814	1.50%	0.174%
14	10.377	1430	1435	1438	rBV	1374460	1491765	46.74%	5.418%
15	10.577	1465	1469	1471	rBV	58842	61402	1.92%	0.223%
16	10.730	1491	1495	1502	rBV2	41972	41790	1.31%	0.152%
17	11.130	1559	1563	1566	rBV	45160	39280	1.23%	0.143%
18	11.218	1574	1578	1584	rVB	607600	572544	17.94%	2.080%
19	11.954	1697	1703	1710	rBV2	98973	153071	4.80%	0.556%
20	12.918	1863	1867	1873	rBV6	21406	37003	1.16%	0.134%
21	13.195	1909	1914	1918	rBV	1747043	1865802	58.46%	6.777%
22	13.512	1964	1968	1971	rBV3	28650	45501	1.43%	0.165%
23	13.542	1971	1973	1976	rVB	38393	35301	1.11%	0.128%
24	14.359	2107	2112	2120	rBV	422639	661471	20.73%	2.403%
25	14.471	2126	2131	2136	rVB	482099	523925	16.42%	1.903%
26	14.765	2177	2181	2188	rBV4	58962	99698	3.12%	0.362%
27	14.912	2202	2206	2211	rBV8	35024	44914	1.41%	0.163%
28	15.148	2241	2246	2254	rBV2	683128	1019601	31.95%	3.703%
29	15.512	2305	2308	2310	rBV3	40040	53253	1.67%	0.193%
30	15.659	2330	2333	2339	rVB5	45872	56595	1.77%	0.206%
31	15.859	2363	2367	2370	rBV	230870	346268	10.85%	1.258%
32	16.100	2404	2408	2411	rBV	332958	328014	10.28%	1.191%
33	16.471	2467	2471	2477	rBV2	91526	152022	4.76%	0.552%
34	16.589	2488	2491	2495	rVB2	76671	89414	2.80%	0.325%

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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	16.795	2524	2526	2534	rVB	83040	129172	4.05%	0.469%
36	18.342	2783	2789	2796	rVB3	261837	526938	16.51%	1.914%

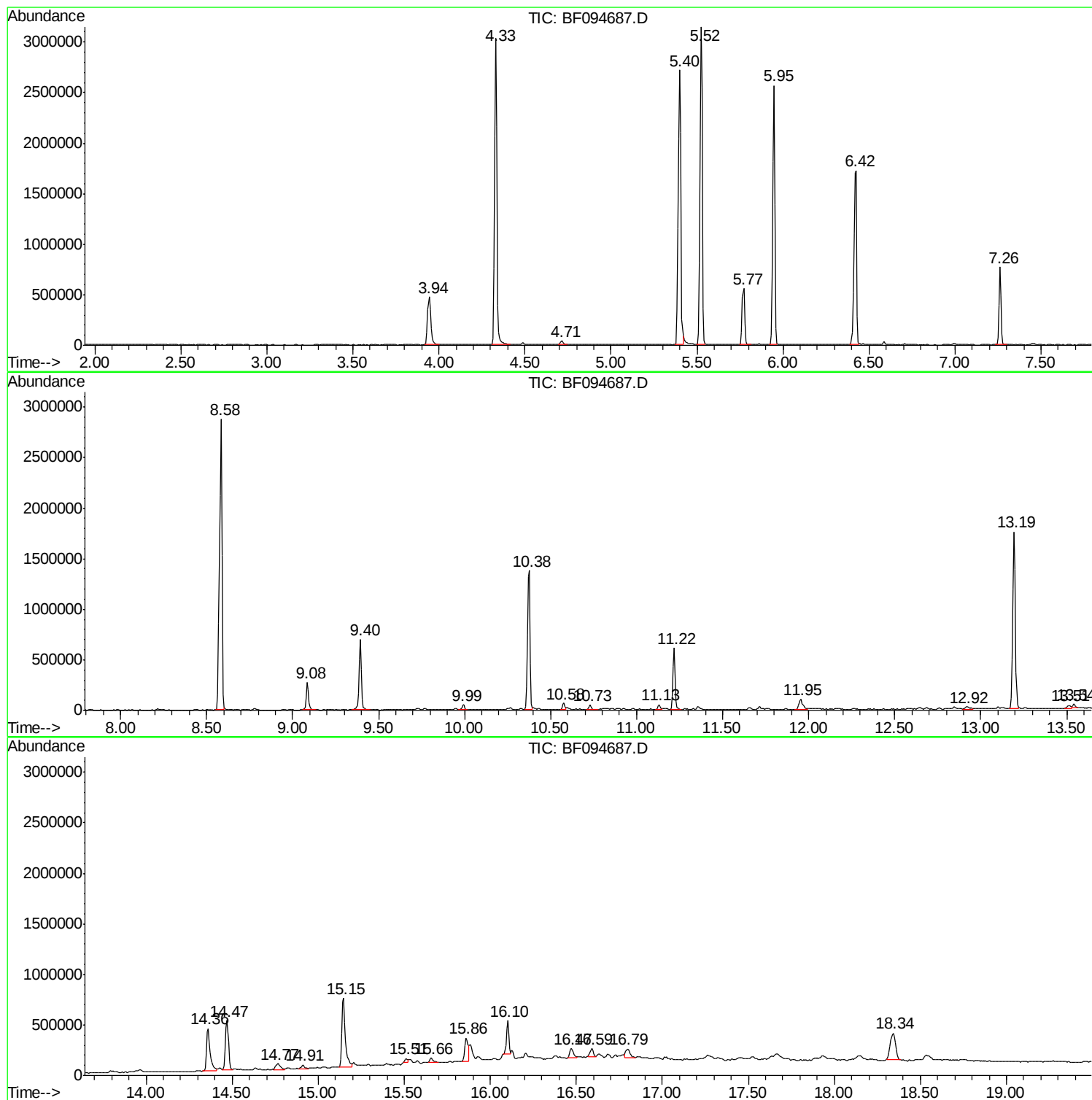
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF041717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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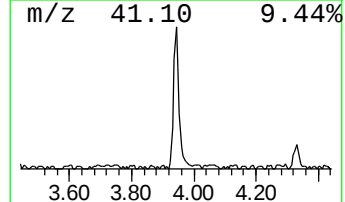
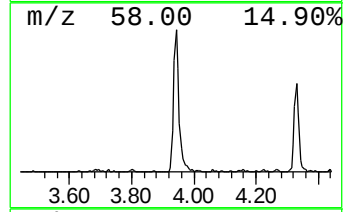
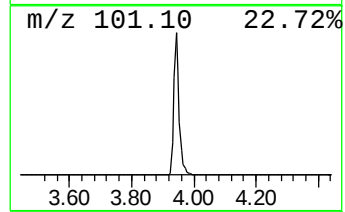
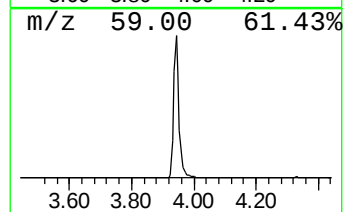
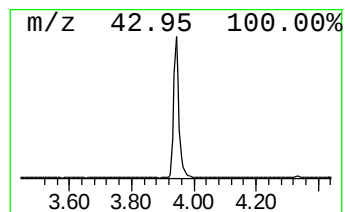
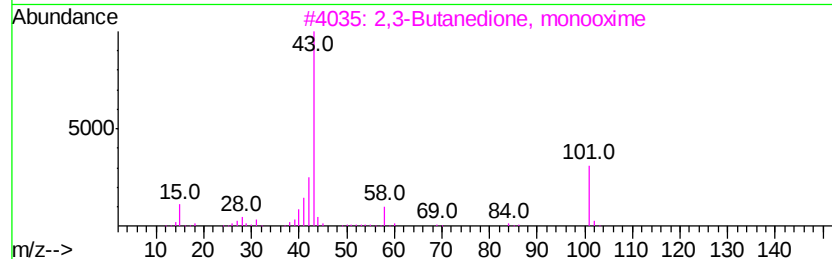
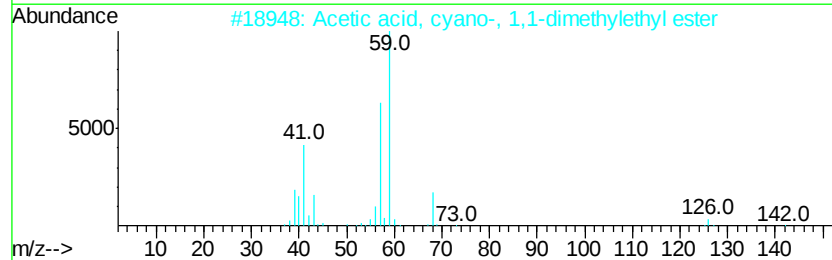
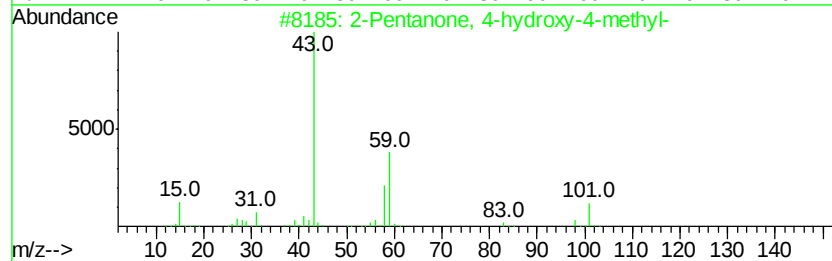
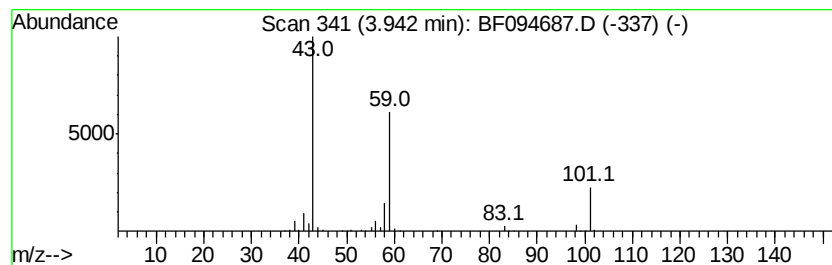
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.94	22.80 ng	571185	1,4-Dichlorobenzene-d4	5.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



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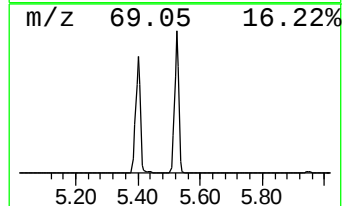
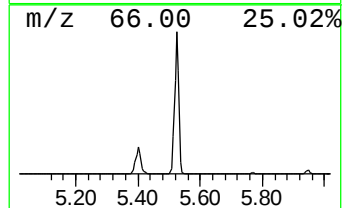
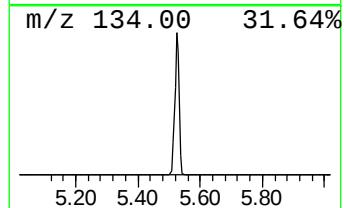
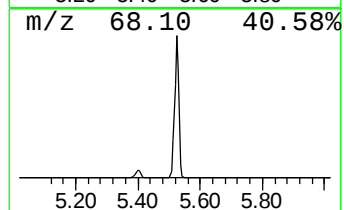
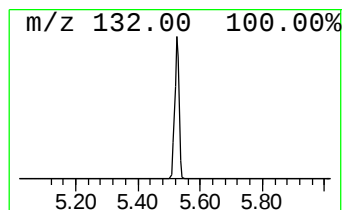
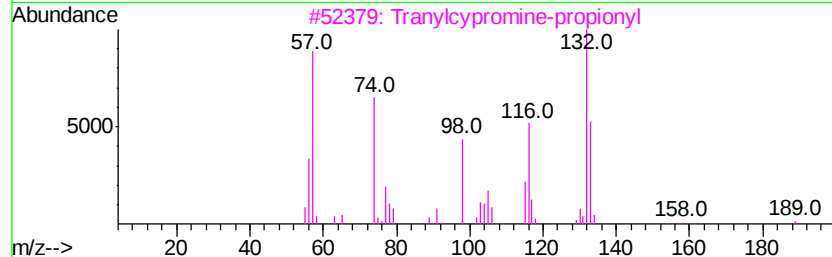
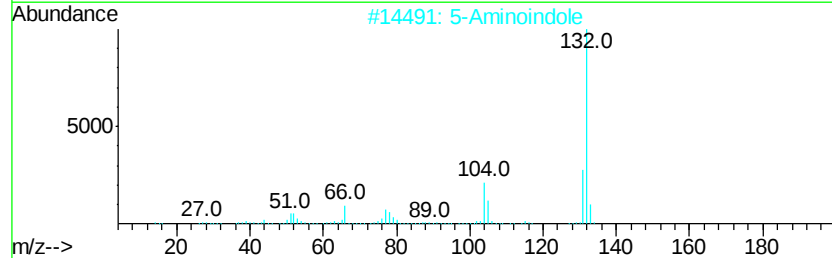
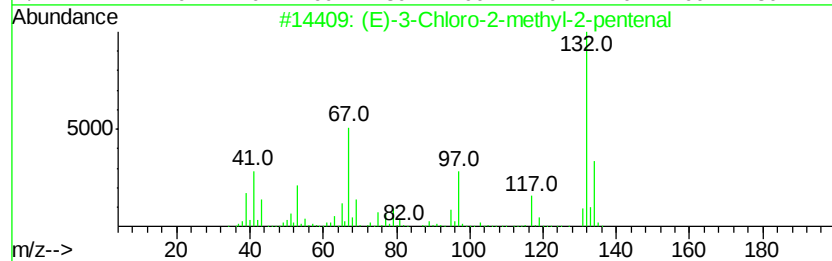
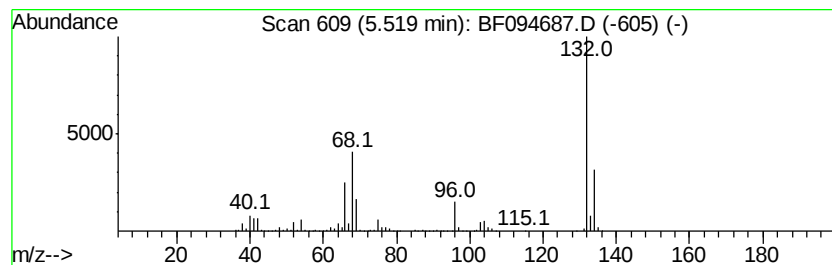
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown5.52 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.52	127.40 ng	3191420	1,4-Dichlorobenzene-d4	5.77

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16	
2	5-Aminoindole	132	C8H8N2	005192-03-0	12	
3	Tranvlcvpromine-propionyl	189	C12H15NO	1000123-86-3	10	
4	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10	
5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9	



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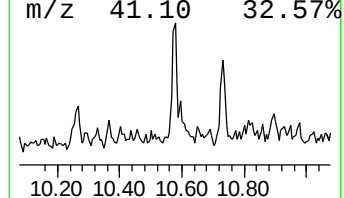
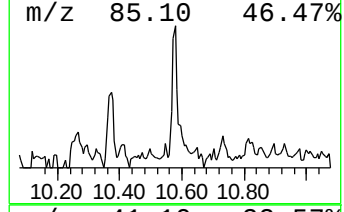
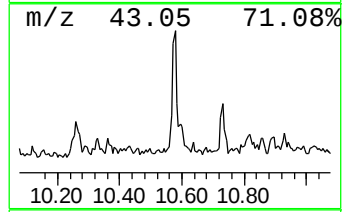
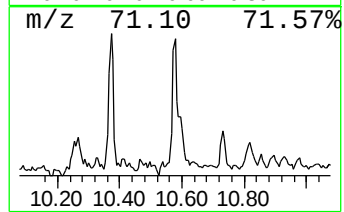
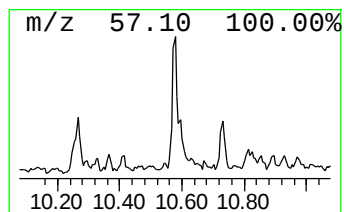
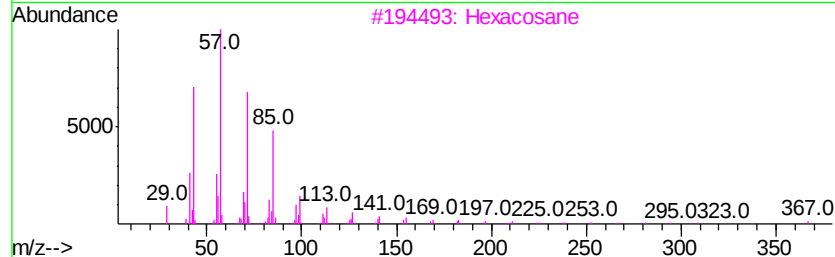
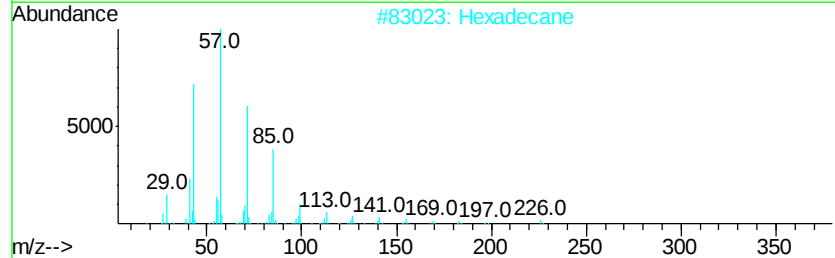
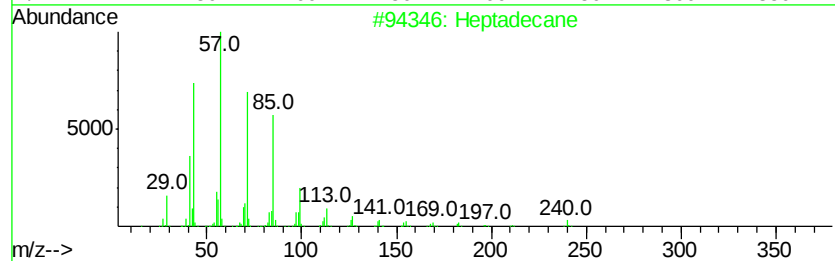
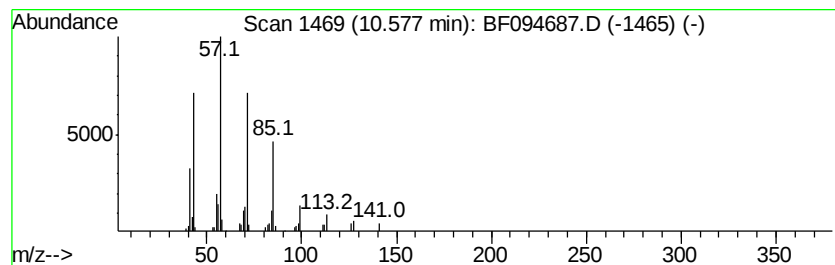
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Heptadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.58	2.14 ng	61402	Phenanthrene-d10	11.22

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	91
2	Hexadecane		226	C16H34	000544-76-3	90
3	Hexacosane		366	C26H54	000630-01-3	90
4	Heptadecane, 2,6,10,15-tetramethyl-		296	C21H44	054833-48-6	87
5	Heneicosane		296	C21H44	000629-94-7	87



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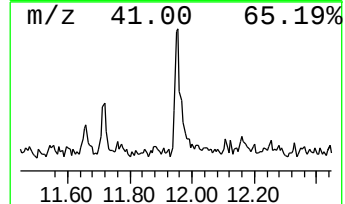
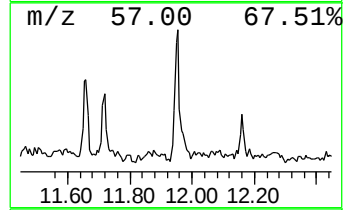
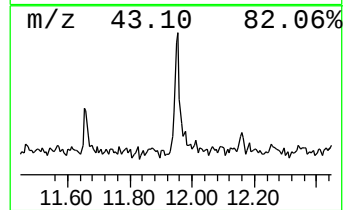
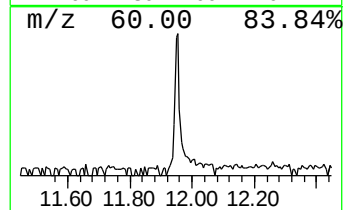
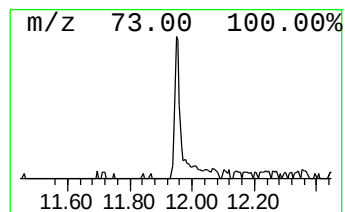
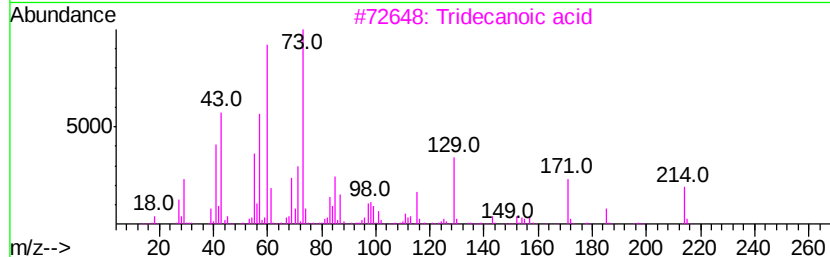
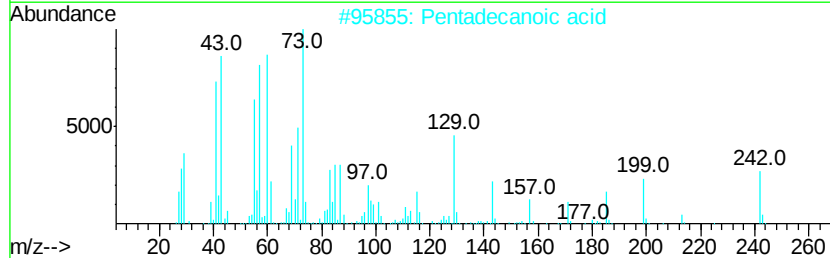
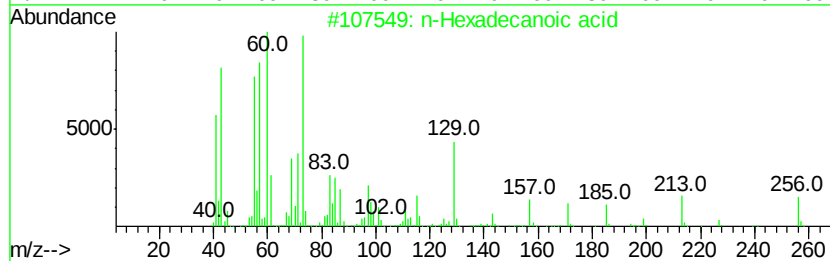
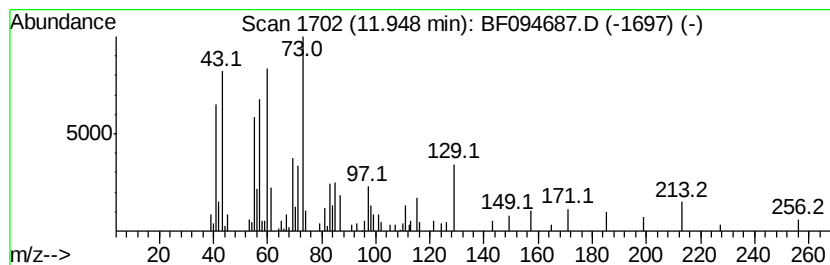
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.95	5.35 ng	153071	Phenanthrene-d10	11.22

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	86
3		Tridecanoic acid	214	C13H26O2	000638-53-9	81
4		n-Decanoic acid	172	C10H20O2	000334-48-5	62
5		Oxalic acid, dodecyl propyl ester	300	C17H32O4	1000309-26-5	25



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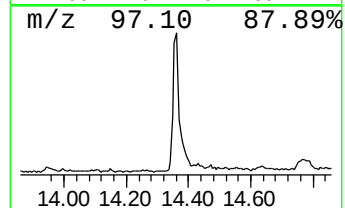
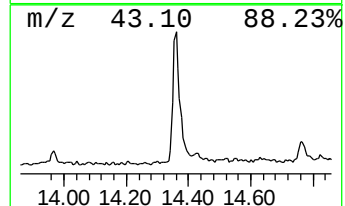
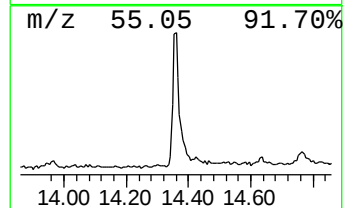
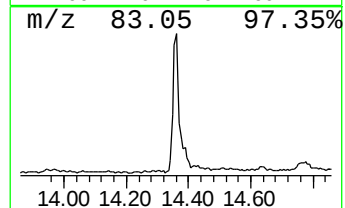
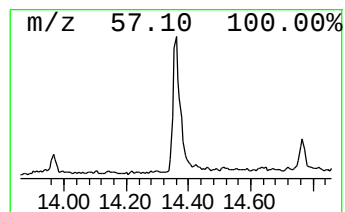
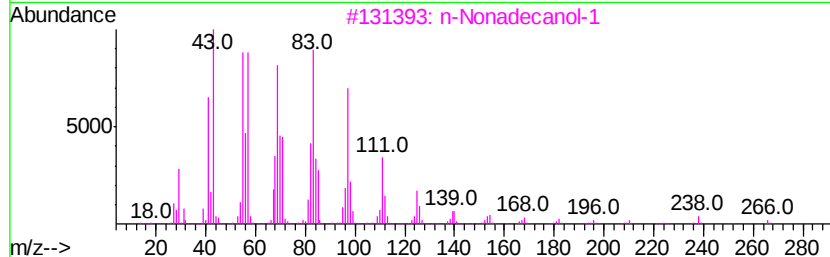
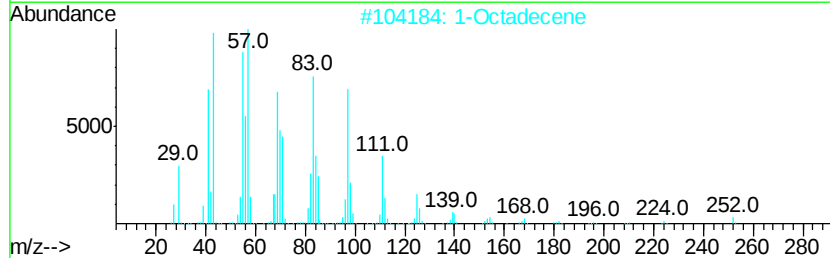
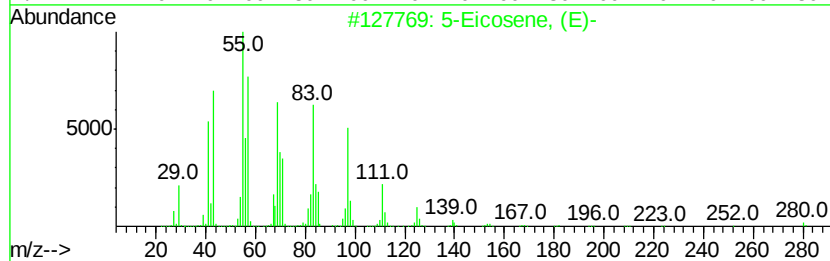
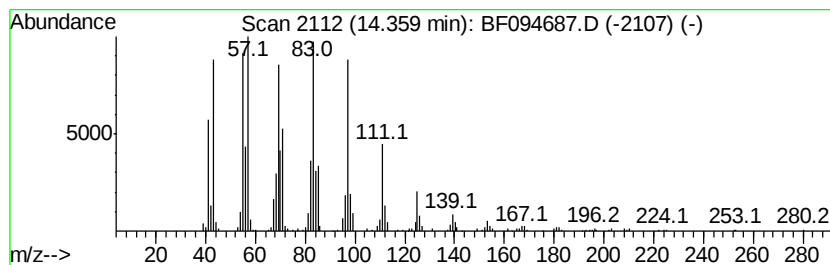
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Peak Number 6 5-Eicosene, (E)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.36	25.25 ng	661471	Chrysene-d12	14.47	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-	280	C20H40	074685-30-6	99
2	1-Octadecene	252	C18H36	000112-88-9	97
3	n-Nonadecanol-1	284	C19H40O	001454-84-8	95
4	1-Heneicosanol	312	C21H44O	015594-90-8	95
5	Cyclopentadecane	210	C15H30	000295-48-7	94



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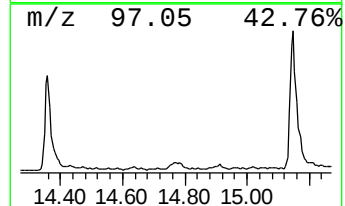
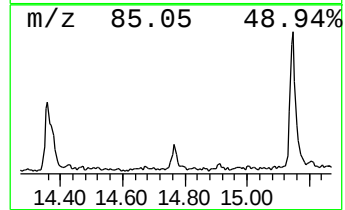
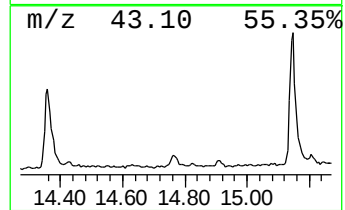
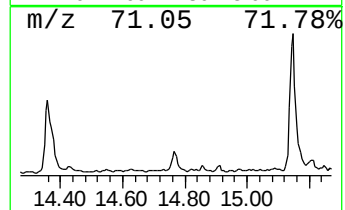
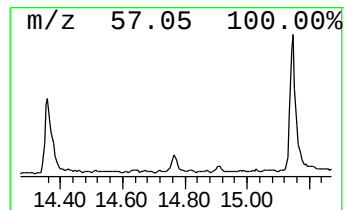
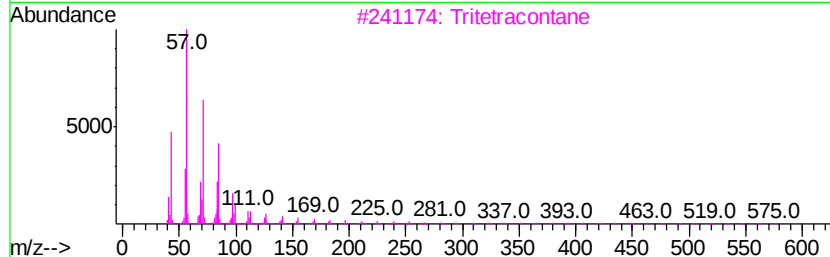
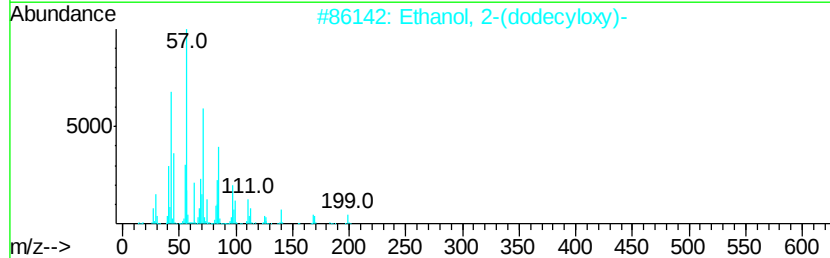
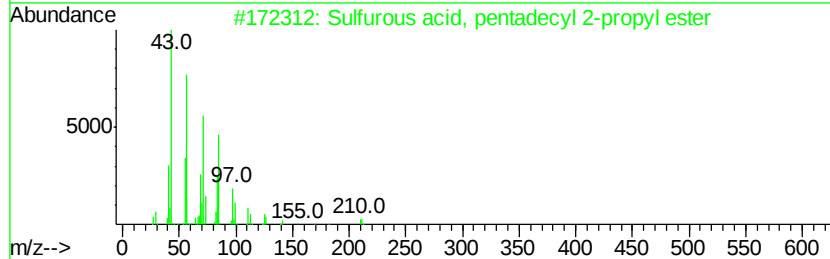
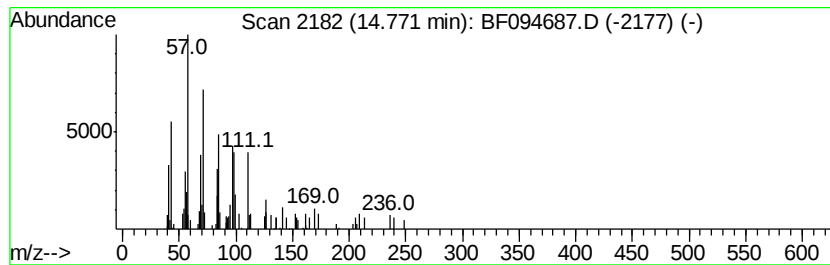
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Sulfurous acid, pentadecyl ... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.77	3.81 ng	99698	Chrysene-d12	14.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, pentadecyl 2-pro...	334	C18H38O3S	1000309-12-6	58
2			Ethanol, 2-(dodecyloxy)-	230	C14H30O2	004536-30-5	53
3			Tritetracontane	605	C43H88	007098-21-7	52
4			Tetradecane	198	C14H30	000629-59-4	49
5			Sulfurous acid, 2-propyl undecyl...	278	C14H30O3S	1000309-12-2	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042817\
Data File : BF094687.D
Acq On : 28 Apr 2017 16:55
Operator : SJ/MA
Sample : I2895-10
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-08-2.0-3.0

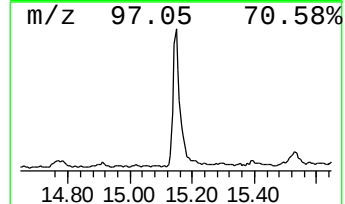
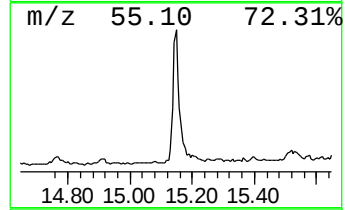
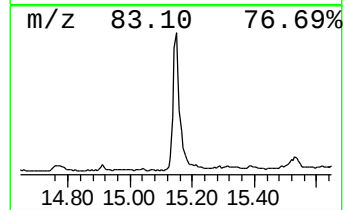
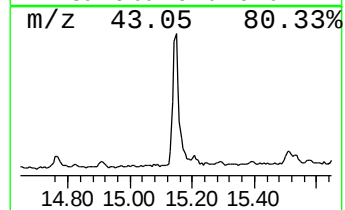
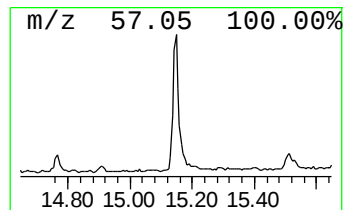
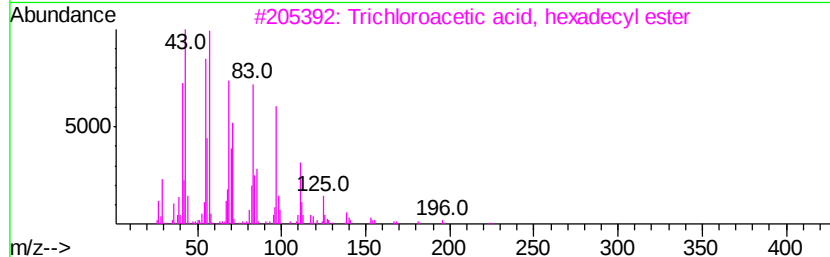
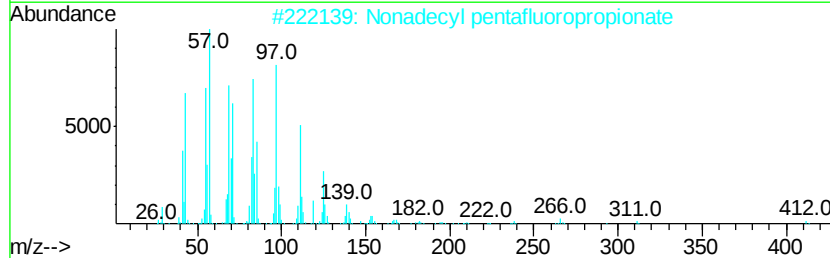
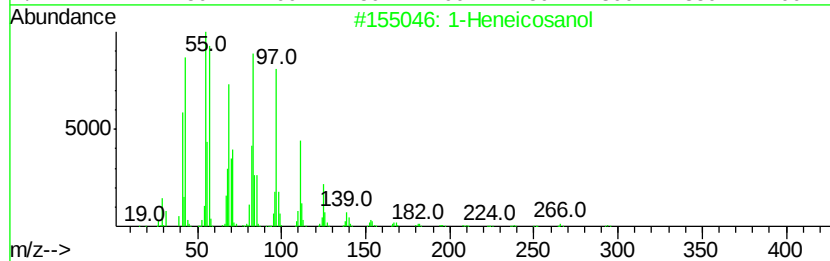
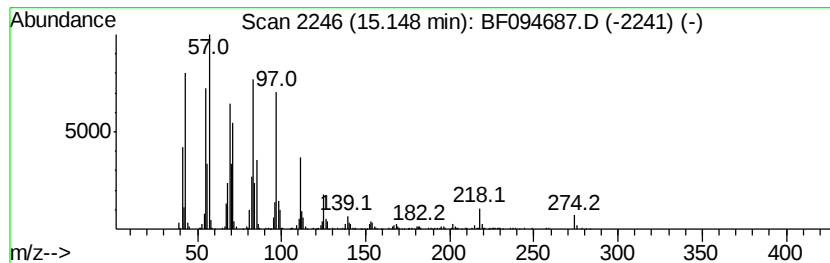
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1-Heneicosanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.15	38.92 ng	1019600	Chrysene-d12	14.47

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Heneicosanol	312	C21H44O	015594-90-8	93
2		Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
3		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
4		17-Pentatriacontene	491	C35H70	006971-40-0	91
5		Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042817\
Data File : BF094687.D
Acq On : 28 Apr 2017 16:55
Operator : SJ/MA
Sample : I2895-10
Misc :
ALS Vial : 8 Sample Multiplier: 1

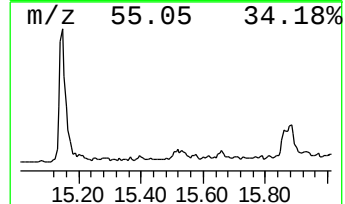
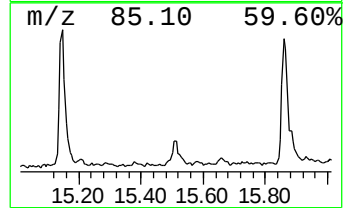
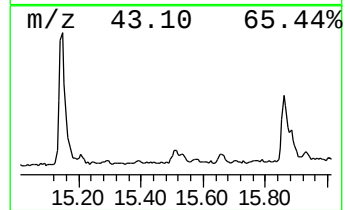
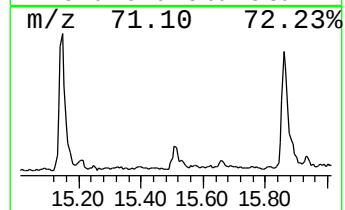
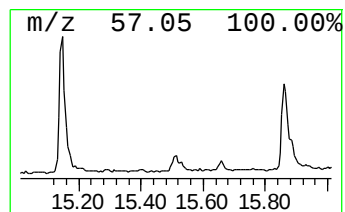
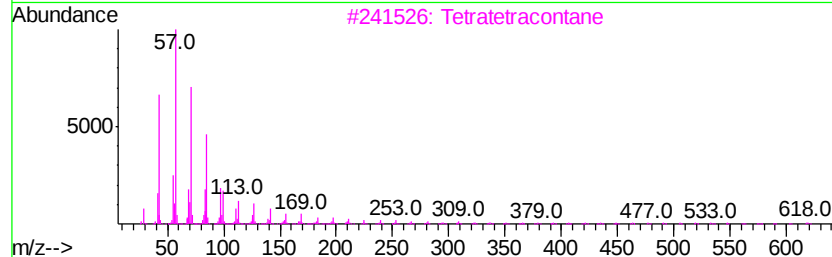
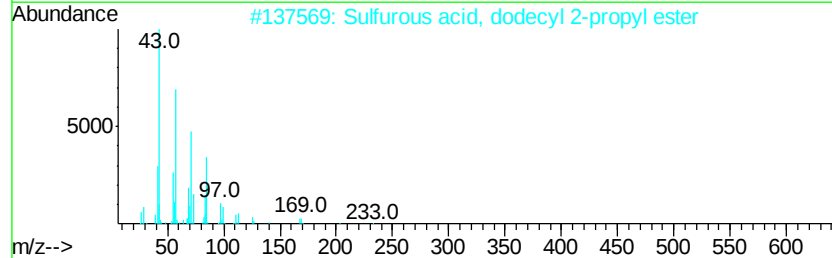
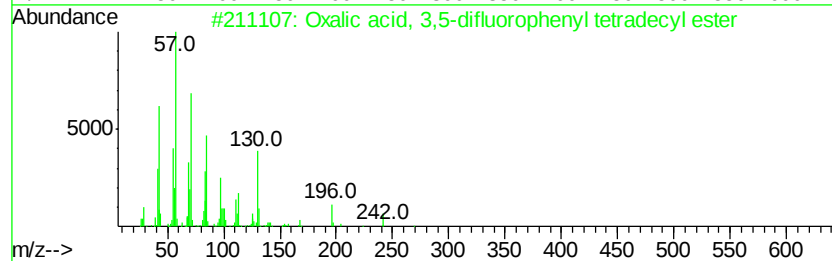
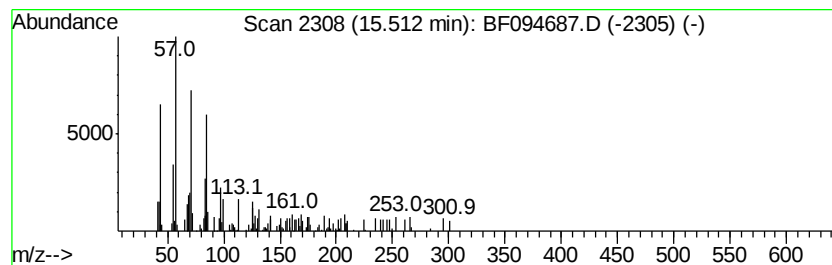
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Oxalic acid, 3,5-difluoroph... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.		
15.51	3.25 ng	53253	Perylene-d12	16.10		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxalic acid, 3,5-difluorophenyl ...	398	C22H32F2O4	1000309-67-8	64	
2	Sulfurous acid, dodecyl 2-propyl...	292	C15H32O3S	1000309-12-3	64	
3	Tetratetracontane	619	C44H90	007098-22-8	64	
4	Sulfurous acid, 2-propyl tridecyl...	306	C16H34O3S	1000309-12-4	59	
5	5,5,7,7-Tetraethylundecane	268	C19H40	1000360-42-0	50	



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042817\
Data File : BF094687.D
Acq On : 28 Apr 2017 16:55
Operator : SJ/MA
Sample : I2895-10
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-08-2.0-3.0

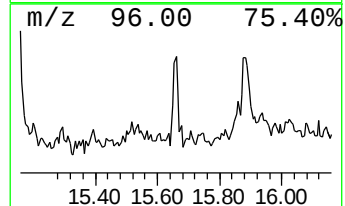
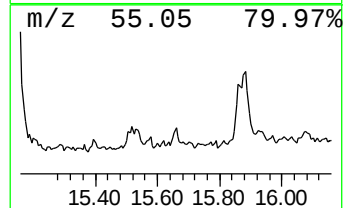
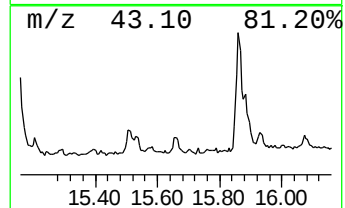
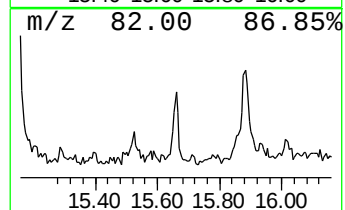
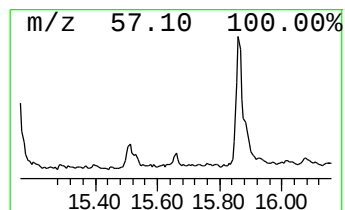
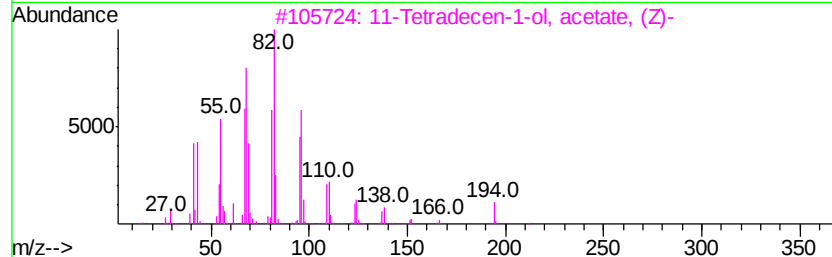
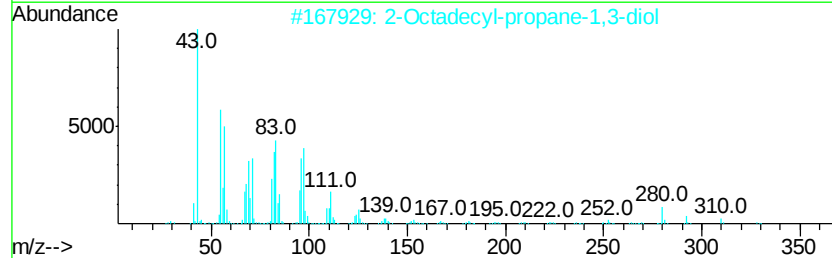
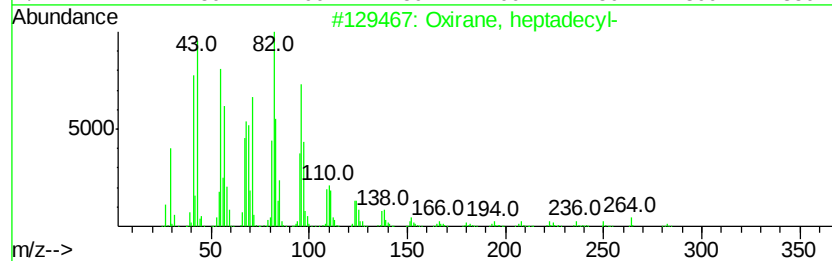
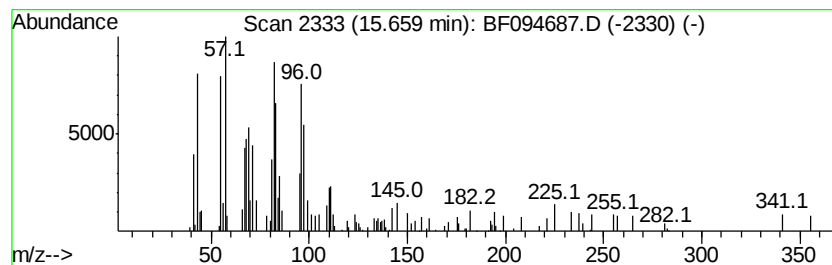
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Oxirane, heptadecyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.66	3.45 ng	56595	Perylene-d12	16.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	74
2		2-Octadecyl-propane-1,3-diol	328	C21H44O2	005337-61-1	47
3		11-Tetradecen-1-ol, acetate, (Z)-	254	C16H30O2	020711-10-8	38
4		Benzocyclodecene, tetradecahydro-	194	C14H26	061142-06-1	38
5		7-Azabicyclo[4.1.0]heptane, 1-me...	111	C7H13N	025022-25-7	30



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042817\
Data File : BF094687.D
Acq On : 28 Apr 2017 16:55
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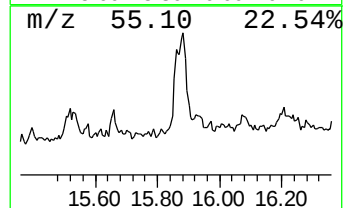
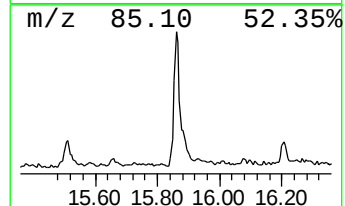
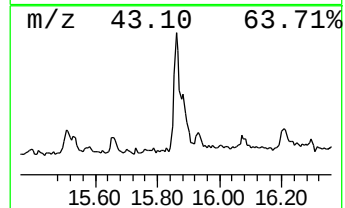
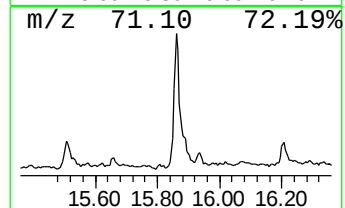
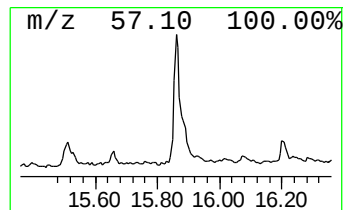
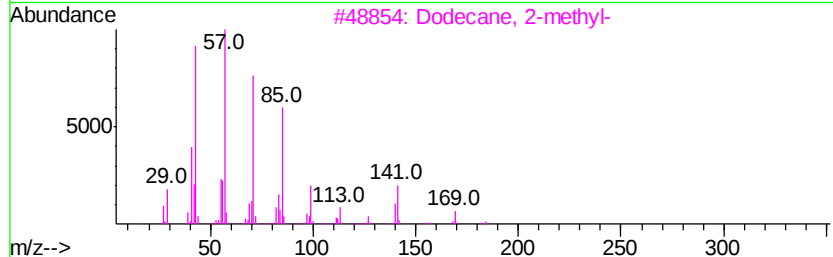
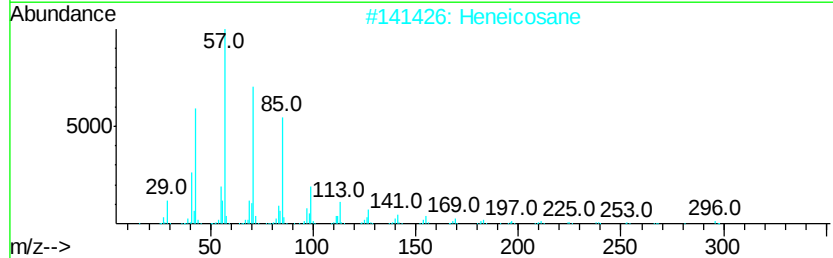
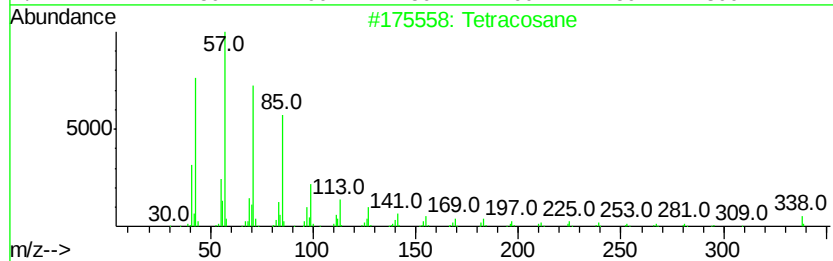
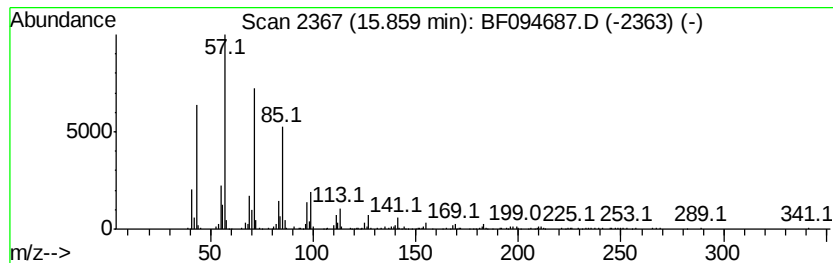
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tetracosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.86	21.11 ng	346268	Perylene-d12	16.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	91
2		Heneicosane	296	C21H44	000629-94-7	90
3		Dodecane, 2-methyl-	184	C13H28	001560-97-0	87
4		2-methyloctacosane	408	C29H60	1000376-72-8	87
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042817\
Data File : BF094687.D
Acq On : 28 Apr 2017 16:55
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ALS Vial : 8 Sample Multiplier: 1

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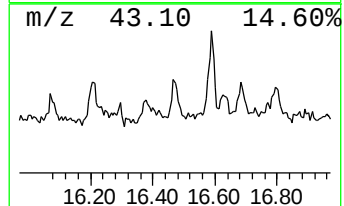
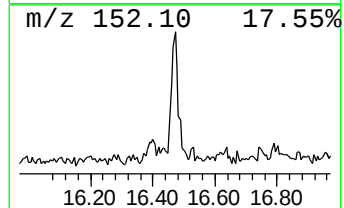
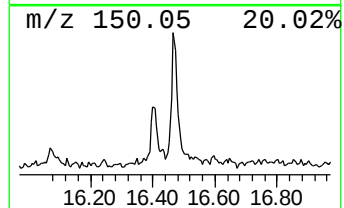
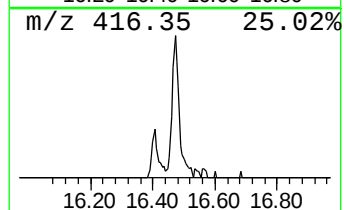
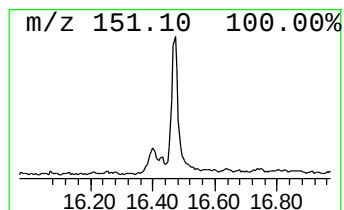
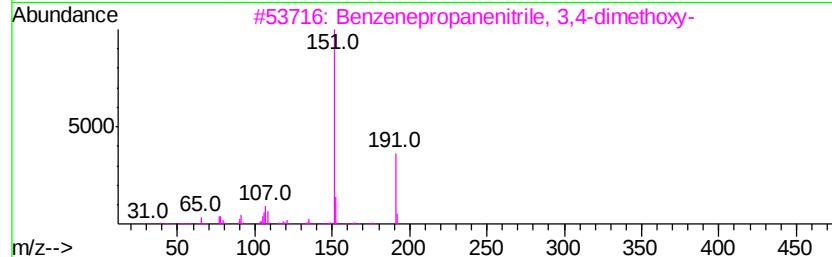
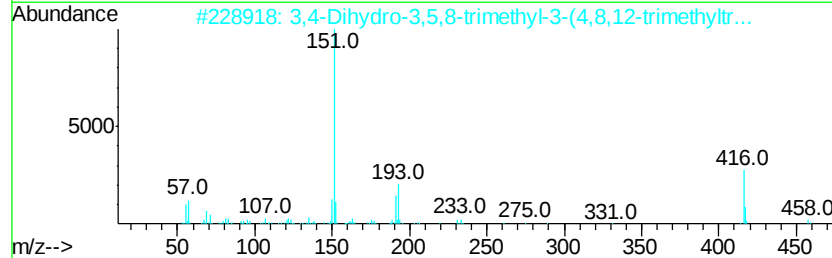
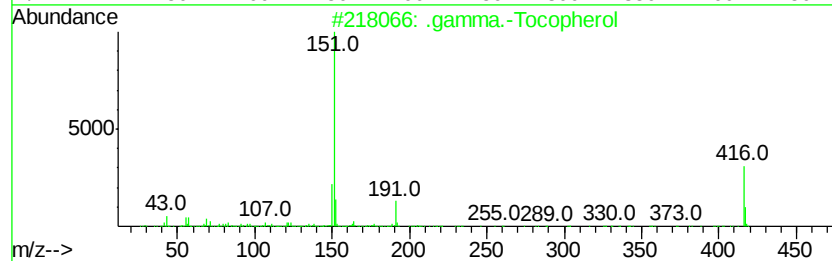
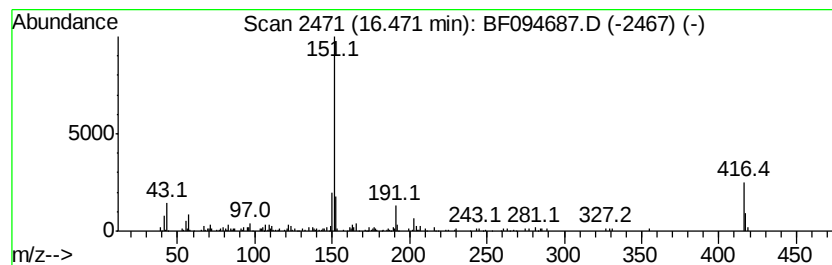
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 .gamma.-Tocopherol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.47	9.27 ng	152022	Perylene-d12	16.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.gamma.-Tocopherol	416	C28H48O2	007616-22-0	96
2		3,4-Dihydro-3,5,8-trimethyl-3-(4...	458	C30H50O3	1000105-71-9	78
3		Benzenepropanenitrile, 3,4-dimet...	191	C11H13N02	049621-56-9	52
4		Benzenethanamine, N-[(3,4-dimet...	331	C19H25N04	055124-87-3	43
5		p-tert-Octylresorcinol	222	C14H22O2	028122-52-3	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042817\
Data File : BF094687.D
Acq On : 28 Apr 2017 16:55
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ALS Vial : 8 Sample Multiplier: 1

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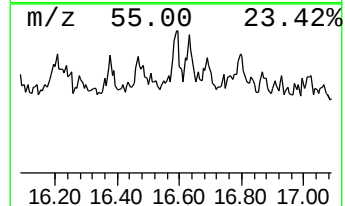
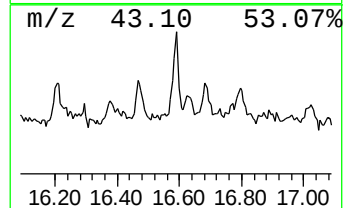
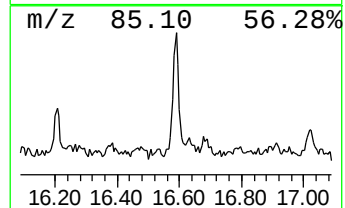
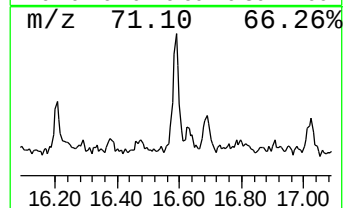
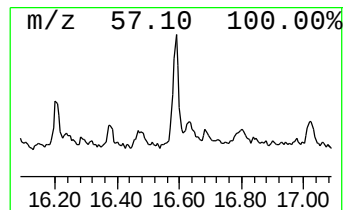
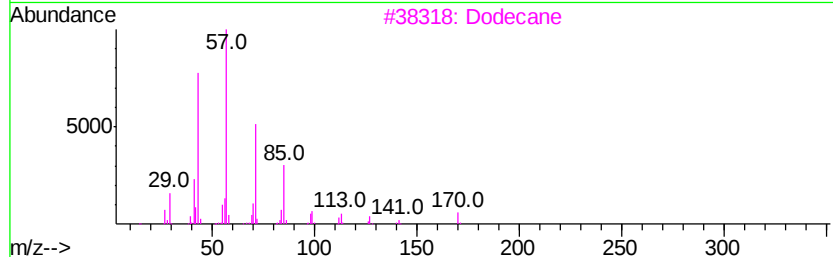
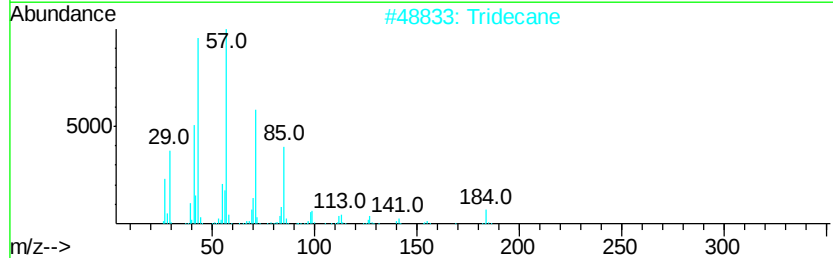
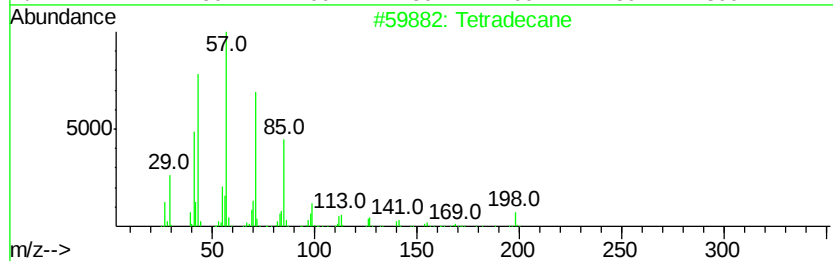
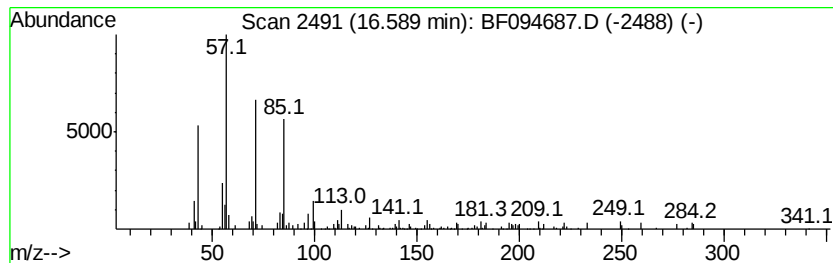
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Tetradecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.59	5.45 ng	89414	Perylene-d12	16.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	87
2		Tridecane	184	C13H28	000629-50-5	80
3		Dodecane	170	C12H26	000112-40-3	80
4		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	80
5		Heptadecane	240	C17H36	000629-78-7	80



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042817\
Data File : BF094687.D
Acq On : 28 Apr 2017 16:55
Operator : SJ/MA
Sample : I2895-10
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-08-2.0-3.0

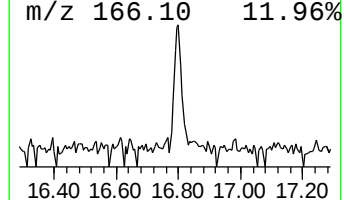
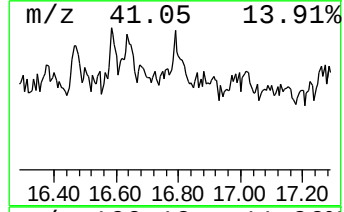
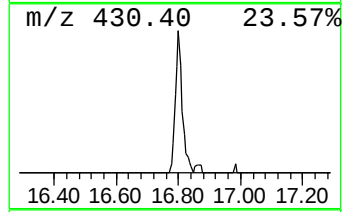
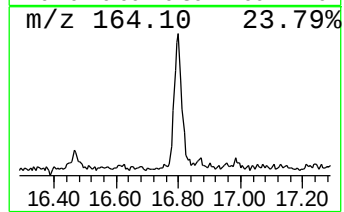
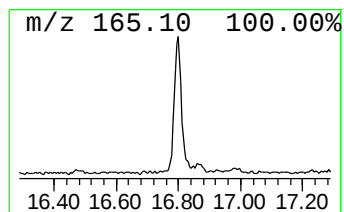
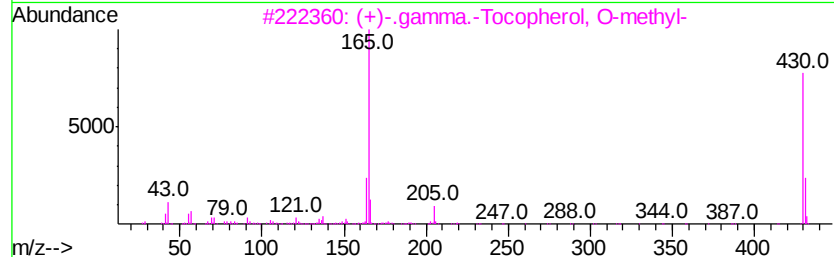
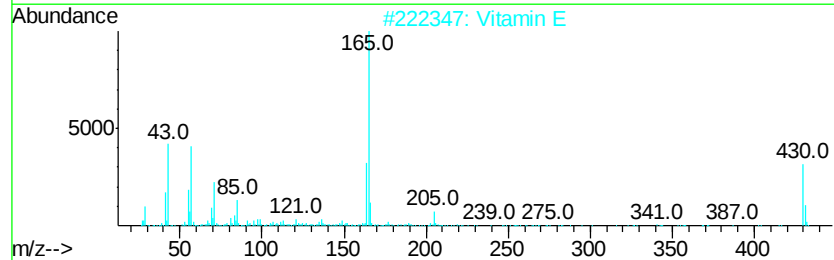
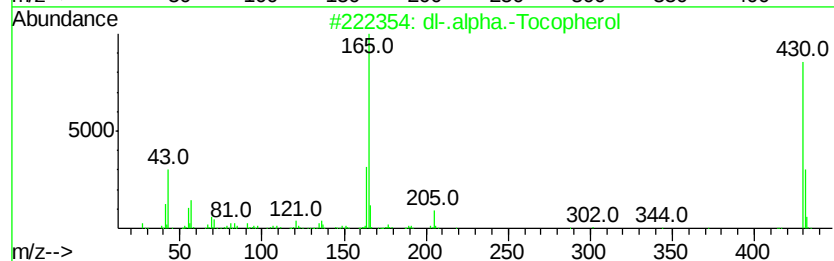
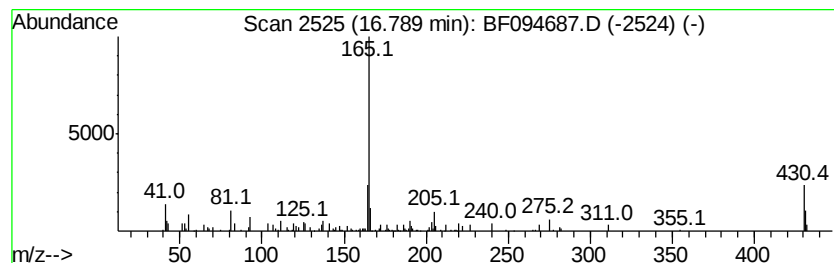
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 dl-.alpha.-Tocopherol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.79	7.88 ng	129172	Perylene-d12	16.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	dl-.alpha.-Tocopherol	430	C29H50O2	010191-41-0	94
2		Vitamin E	430	C29H50O2	000059-02-9	94
3		(+)-.gamma.-Tocopherol, 0-methyl-	430	C29H50O2	1000374-72-2	91
4		Piperazine, 1-(3,4-dimethoxybenz...	430	C26H26N2O4	333756-87-9	59
5		Furo[3,4-c]pyridine-3,4(1H,5H)-d...	165	C8H7NO3	007472-18-6	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042817\
Data File : BF094687.D
Acq On : 28 Apr 2017 16:55
Operator : SJ/MA
Sample : I2895-10
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-08-2.0-3.0

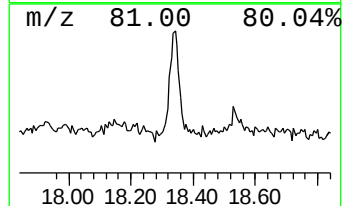
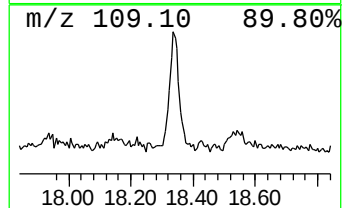
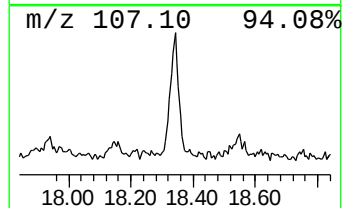
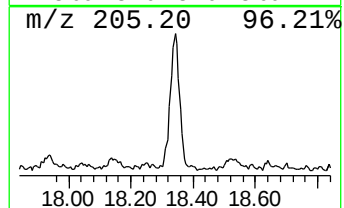
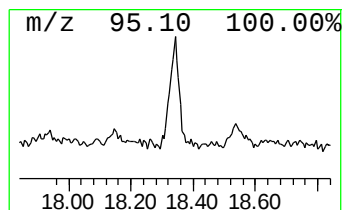
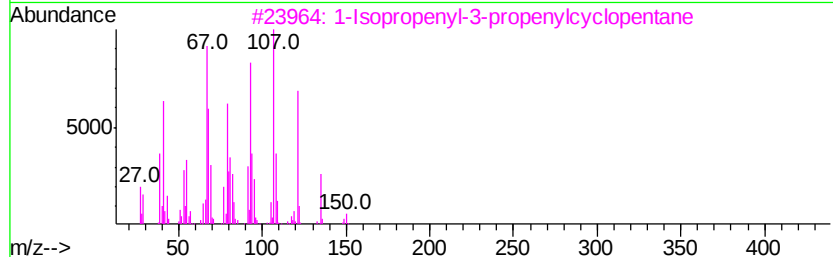
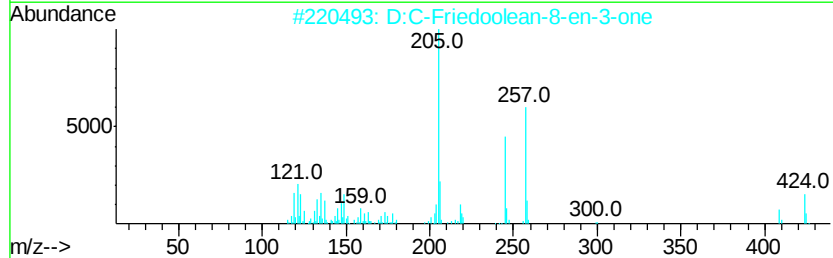
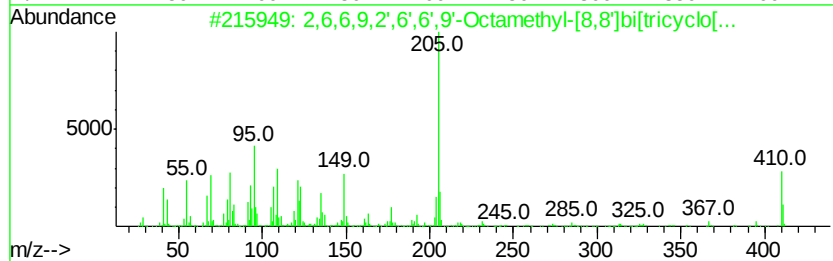
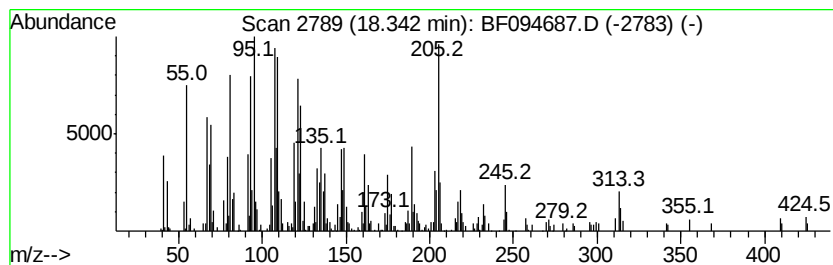
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 2,6,6,9,2',6',6',9'-Octameth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.34	32.13 ng	526938	Perylene-d12	16.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,6,9,2',6',6',9'-Octamethyl-[...]	410	C30H50	1000189-48-2	55		
2	D:C-Friedoolean-8-en-3-one	424	C30H48O	022611-26-3	49		
3	1-Isopropenyl-3-propenylcyclopentane...	150	C11H18	1000192-81-2	43		
4	But-3-enal, 2-methyl-4-(2,6,6-tr...	206	C14H22O	104900-59-6	42		
5	.gamma.-Elemene	204	C15H24	029873-99-2	41		



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042817\
 Data File : BF094687.D
 Acq On : 28 Apr 2017 16:55
 Operator : SJ/MA
 Sample : I2895-10
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-08-2.0-3.0

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF041717.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
2-Pentanone, 4-hy...	3.94	22.8	ng	571185	1	5.77	501012	20.0
unknown5.52	5.52	127.4	ng	3191420	1	5.77	501012	20.0
Heptadecane	10.58	2.1	ng	61402	4	11.22	572544	20.0
n-Hexadecanoic acid	11.95	5.3	ng	153071	4	11.22	572544	20.0
5-Eicosene, (E)-	14.36	25.3	ng	661471	5	14.47	523925	20.0
Sulfurous acid, p...	14.77	3.8	ng	99698	5	14.47	523925	20.0
1-Heneicosanol	15.15	38.9	ng	1019600	5	14.47	523925	20.0
Oxalic acid, 3,5-...	15.51	3.3	ng	53253	6	16.10	328014	20.0
Oxirane, heptadecyl-	15.66	3.5	ng	56595	6	16.10	328014	20.0
Tetracosane	15.86	21.1	ng	346268	6	16.10	328014	20.0
.gamma.-Tocopherol	16.47	9.3	ng	152022	6	16.10	328014	20.0
Tetradecane	16.59	5.5	ng	89414	6	16.10	328014	20.0
dl-.alpha.-Tocoph...	16.79	7.9	ng	129172	6	16.10	328014	20.0
2,6,6,9,2',6',6',...	18.34	32.1	ng	526938	6	16.10	328014	20.0