

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
 Data File : BF083031.D
 Acq On : 19 Nov 2015 2:19
 Operator : UM/IZ
 Sample : G4426-04
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-4

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-BF110715.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.898	242	246	251	rBV	1268836	1932418	42.24%	5.413%
2	5.344	282	285	287	rBV	3881524	3899652	85.24%	10.924%
3	6.396	373	377	379	rBV	3768521	4575152	100.00%	12.817%
4	6.510	384	387	389	rBV	4346719	4208314	91.98%	11.789%
5	6.727	404	406	408	rBV	564177	734739	16.06%	2.058%
6	6.887	418	420	423	rBV	2409366	2723787	59.53%	7.630%
7	7.299	453	456	459	rBV	2227067	2233465	48.82%	6.257%
8	7.424	465	467	472	rVB	75644	117849	2.58%	0.330%
9	7.767	495	497	502	rVB	51872	61202	1.34%	0.171%
10	8.019	517	519	521	rBV	1161674	955621	20.89%	2.677%
11	8.979	602	603	609	rVB	45706	47814	1.05%	0.134%
12	9.105	611	614	616	rBV	3591032	3724860	81.41%	10.435%
13	9.436	640	643	646	rBV3	33654	74674	1.63%	0.209%
14	9.493	646	648	651	rVV	921916	829914	18.14%	2.325%
15	9.722	666	668	670	rVB	138985	103810	2.27%	0.291%
16	9.767	670	672	675	rBV2	835179	941071	20.57%	2.636%
17	9.973	686	690	692	rBV4	29386	65581	1.43%	0.184%
18	10.190	704	709	710	rBV	54510	72780	1.59%	0.204%
19	10.225	710	712	713	rVB	121381	137467	3.00%	0.385%
20	10.442	728	731	732	rBV	66373	86654	1.89%	0.243%
21	10.556	739	741	744	rBV	1774796	1739451	38.02%	4.873%
22	10.693	751	753	757	rBV	186976	243459	5.32%	0.682%
23	11.139	790	792	793	rBV	122654	96507	2.11%	0.270%
24	11.253	799	802	803	rBV	760446	791346	17.30%	2.217%
25	11.322	805	808	810	rBV3	39149	51070	1.12%	0.143%
26	11.562	828	829	832	rVB	77815	89231	1.95%	0.250%
27	11.756	841	846	847	rBV3	29044	74429	1.63%	0.209%
28	11.859	853	855	860	rVB3	39445	62152	1.36%	0.174%
29	11.973	861	865	867	rBV	61653	85171	1.86%	0.239%
30	12.133	877	879	881	rBV2	33667	47009	1.03%	0.132%
31	12.248	886	889	891	rVV3	34749	77141	1.69%	0.216%
32	12.293	891	893	895	rVV2	68269	102227	2.23%	0.286%
33	12.362	895	899	900	rVB2	49869	90877	1.99%	0.255%
34	12.453	906	907	909	rBV	49131	60533	1.32%	0.170%

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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

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Peak separation: 5

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	12.636	922	923	925 rBV	40641	55663	1.22%	0.156%
36	12.682	925	927	928 rVV	62837	62548	1.37%	0.175%
37	12.728	930	931	936 rVB2	46978	57601	1.26%	0.161%
38	12.842	938	941	943 rBV	2000571	2349750	51.36%	6.582%
39	12.991	952	954	957 rBV2	54045	100077	2.19%	0.280%
40	13.082	960	962	964 rVB	60373	56214	1.23%	0.157%
41	13.208	970	973	975 rBV3	36914	67223	1.47%	0.188%
42	13.322	980	983	985 rVB	71579	93175	2.04%	0.261%
43	13.425	990	992	994 rVV	54262	84823	1.85%	0.238%
44	13.494	997	998	1001 rVV2	105806	124998	2.73%	0.350%
45	13.539	1001	1002	1005 rVB	52702	46825	1.02%	0.131%
46	13.756	1019	1021	1023 rBV	107158	137017	2.99%	0.384%
47	13.814	1023	1026	1028 rVB2	29378	56995	1.25%	0.160%
48	13.882	1030	1032	1034 rBV	770753	755011	16.50%	2.115%
49	15.265	1151	1153	1156 rVB	352135	511711	11.18%	1.433%

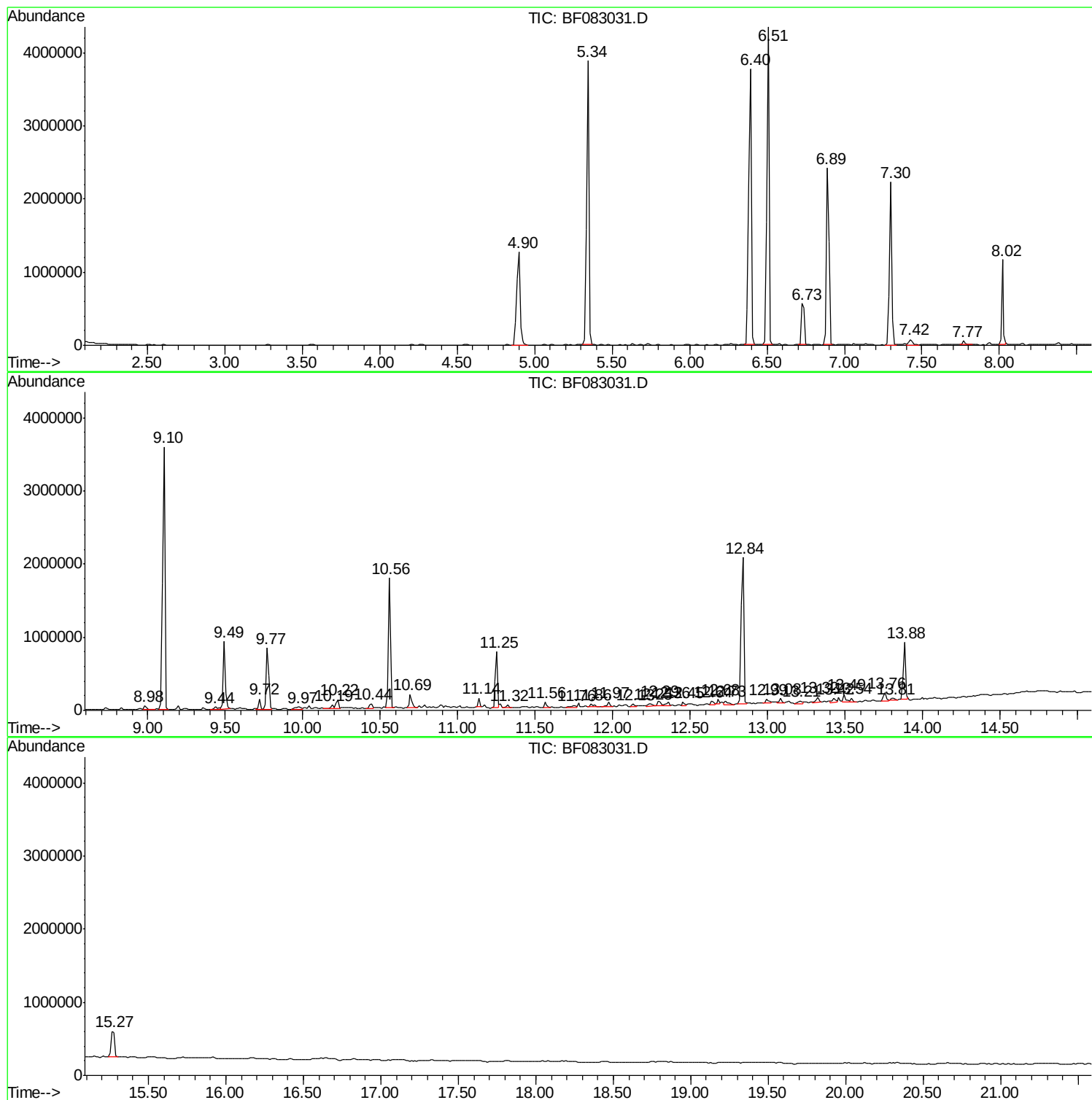
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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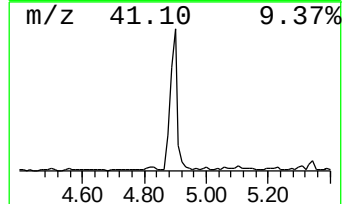
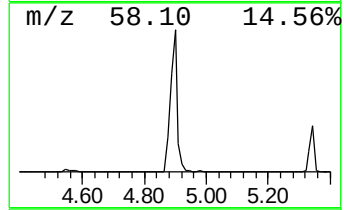
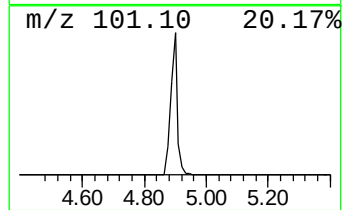
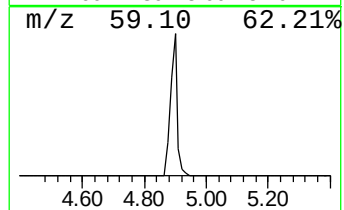
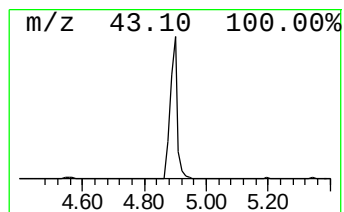
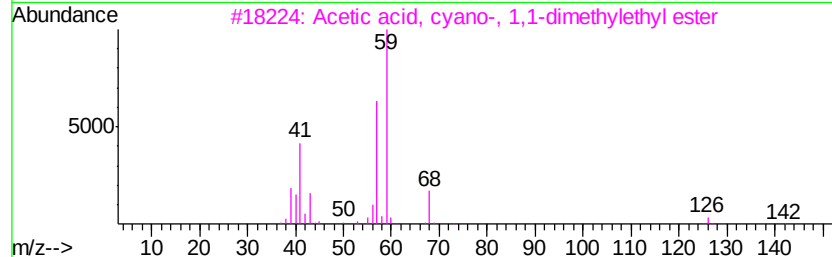
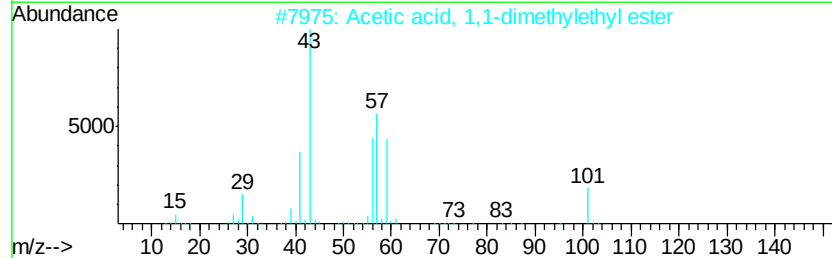
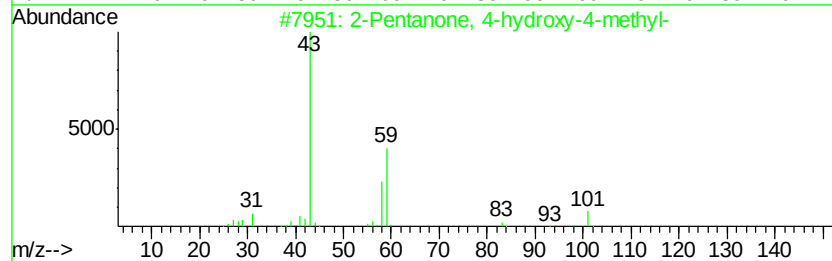
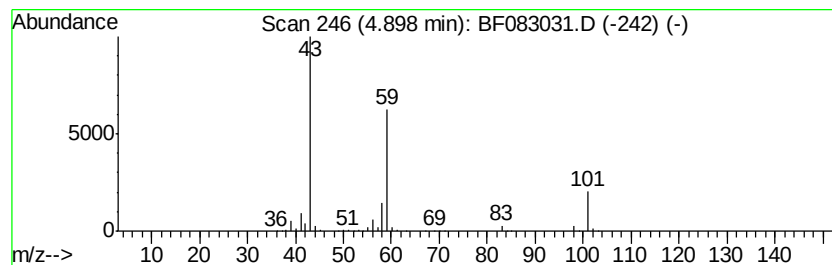
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
4.90	52.60 ng	1932420	1,4-Dichlorobenzene-d4	6.74	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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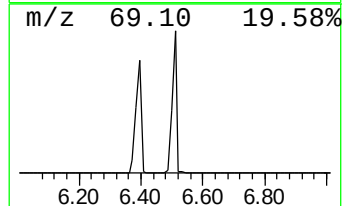
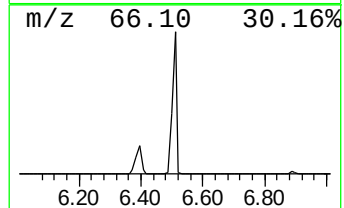
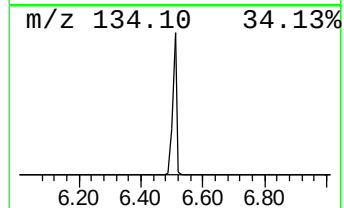
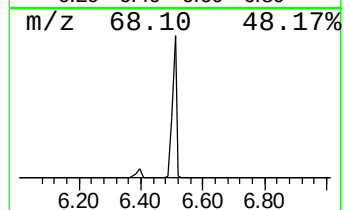
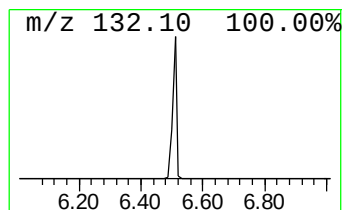
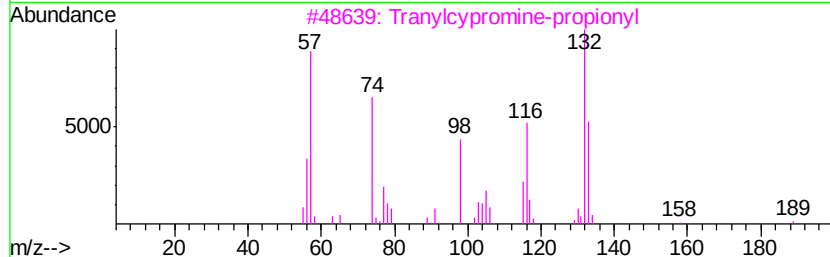
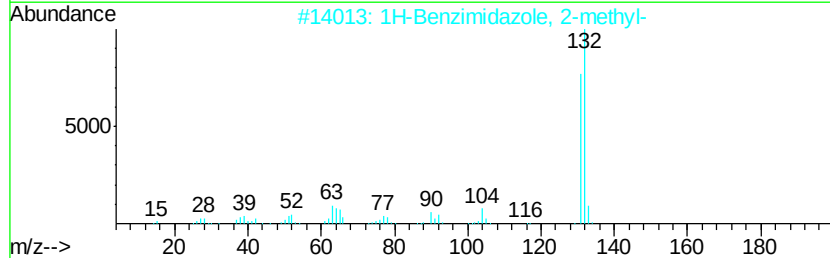
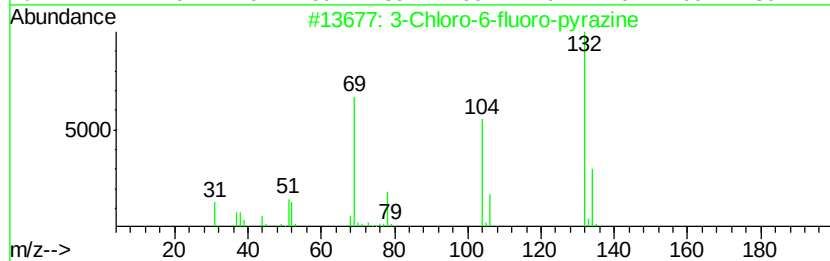
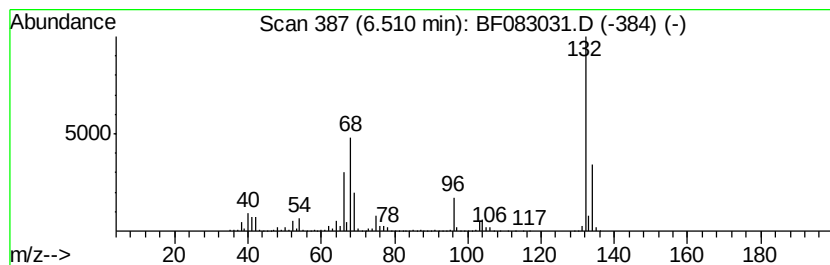
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.51 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	114.55 ng	4208310	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	12
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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Instrument :
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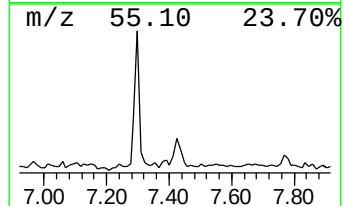
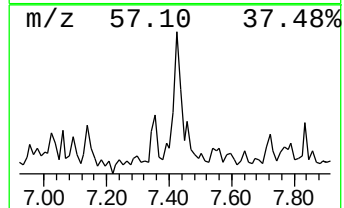
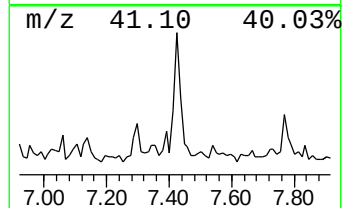
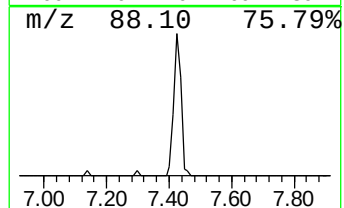
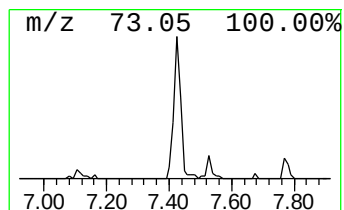
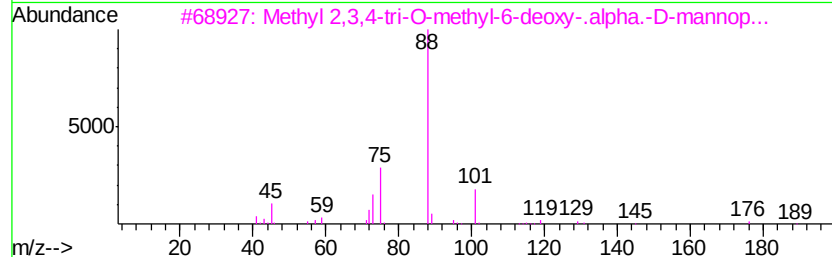
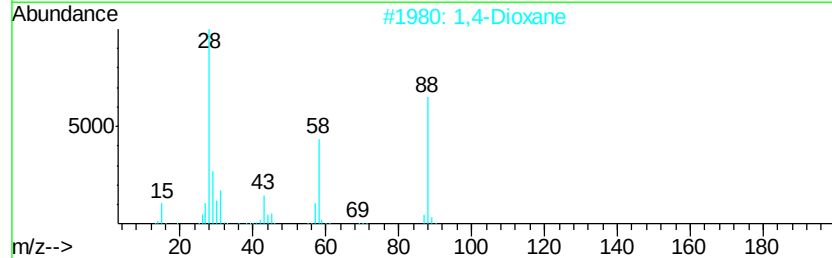
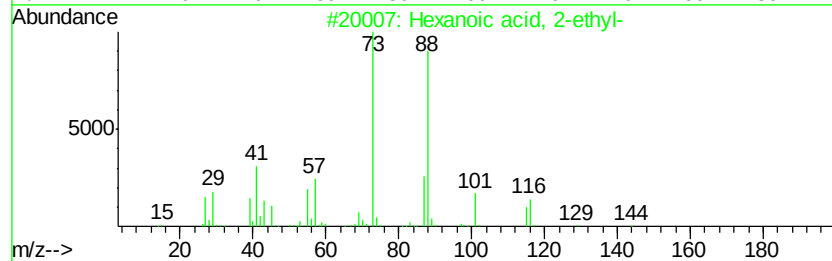
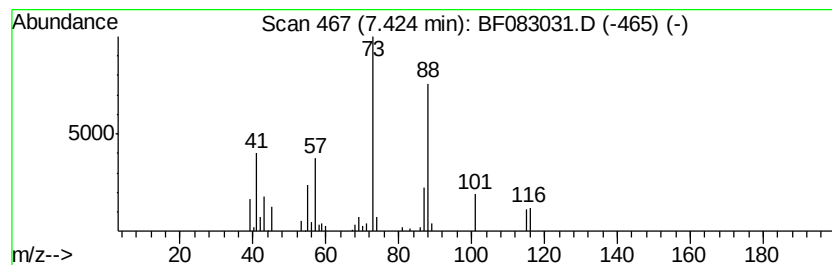
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Hexanoic acid, 2-ethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.42	2.47 ng	117849	Napthalene-d8	8.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
2		1,4-Dioxane	88	C4H8O2	000123-91-1	38
3		Methyl 2,3,4-tri-O-methyl-6-deox...	220	C10H20O5	072983-16-5	36
4		Hydrazine, 1,2-diethyl-	88	C4H12N2	001615-80-1	33
5		Urea, N,N'-dimethyl-	88	C3H8N2O	000096-31-1	32



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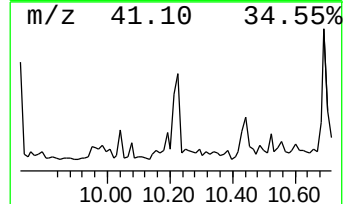
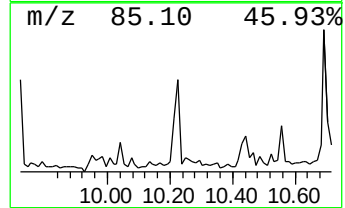
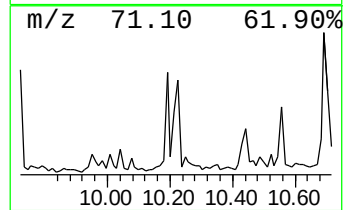
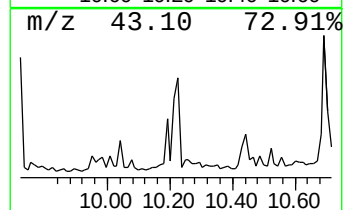
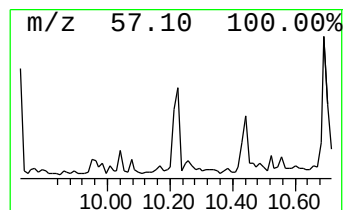
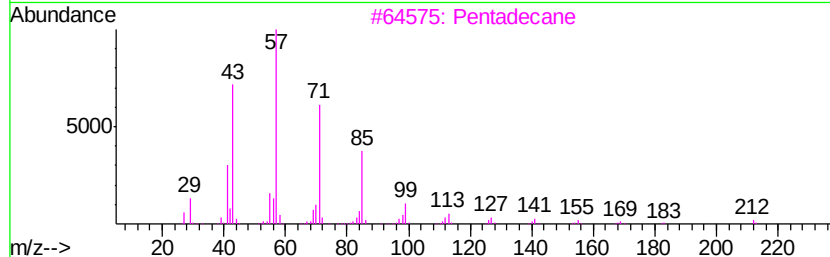
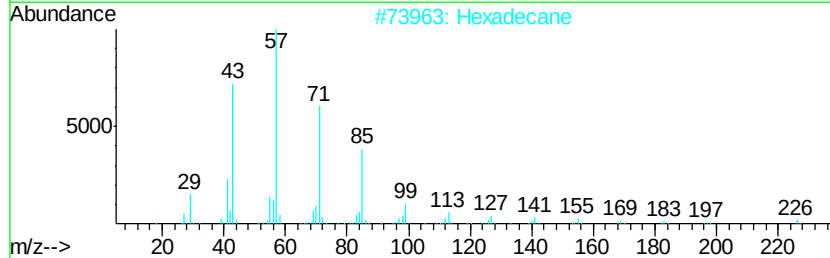
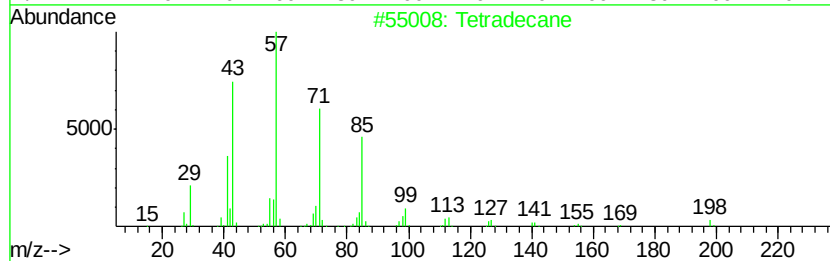
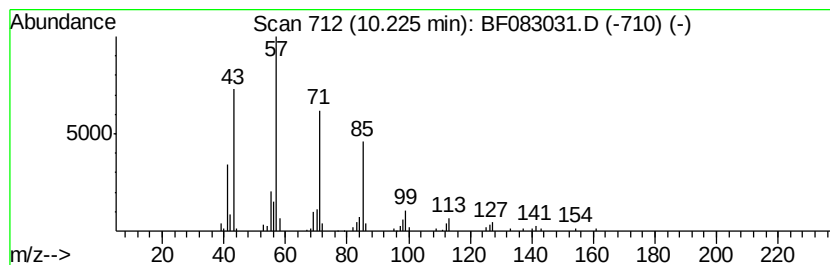
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Tetradecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.22	2.92 ng	137467	Acenaphthene-d10		9.77
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	90
2	Hexadecane	226	C16H34	000544-76-3	90
3	Pentadecane	212	C15H32	000629-62-9	90
4	Tritetracontane	605	C43H88	007098-21-7	90
5	Heptacosane	380	C27H56	000593-49-7	90



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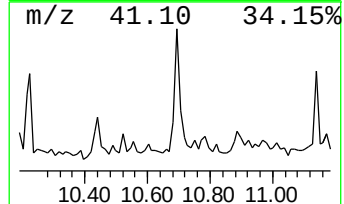
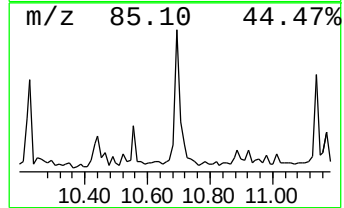
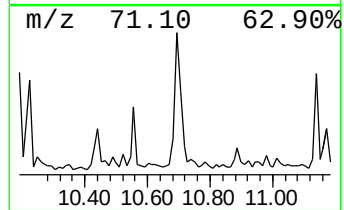
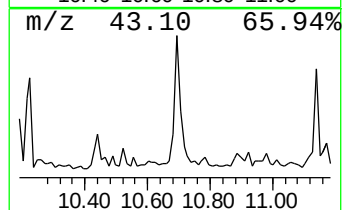
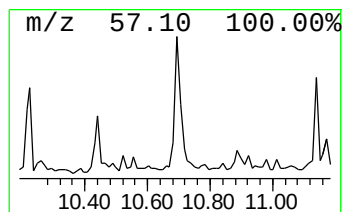
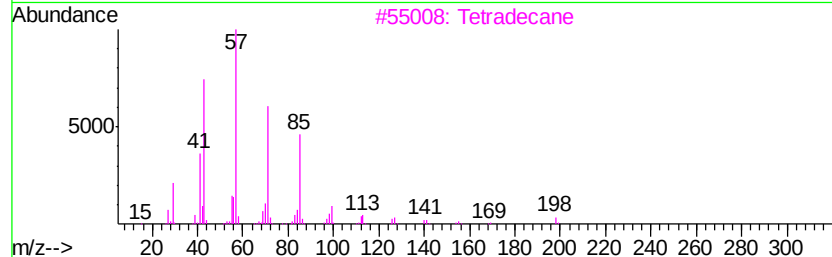
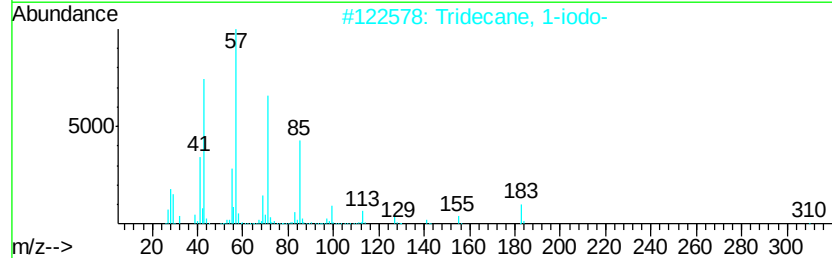
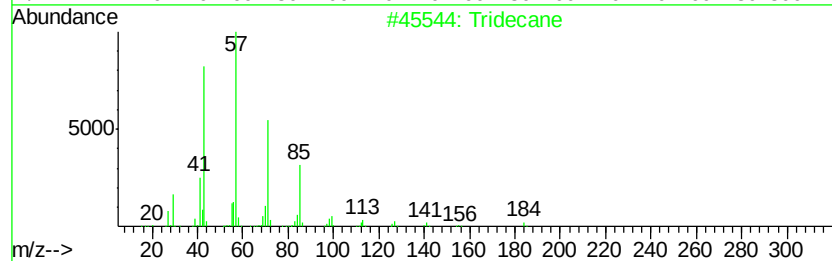
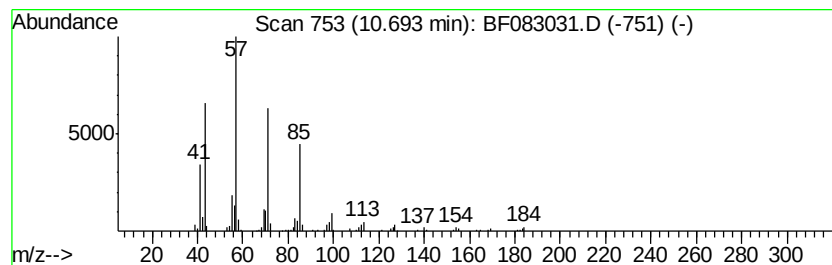
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TIC Integration Parameters: LSCINT.P

Peak Number 6 Tridecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.69	6.15 ng	243459	Phenanthrene-d10		11.25	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane	184	C13H28	000629-50-5	93
2		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	91
3		Tetradecane	198	C14H30	000629-59-4	91
4		Pentadecane	212	C15H32	000629-62-9	90
5		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
Data File : BF083031.D
Acq On : 19 Nov 2015 2:19
Operator : UM/IZ
Sample : G4426-04
Misc :
ALS Vial : 25 Sample Multiplier: 1

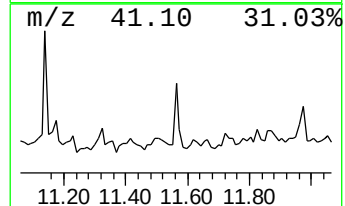
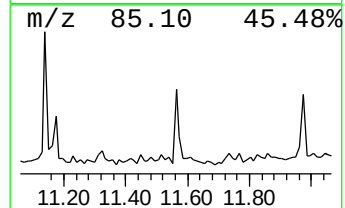
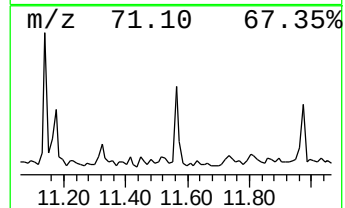
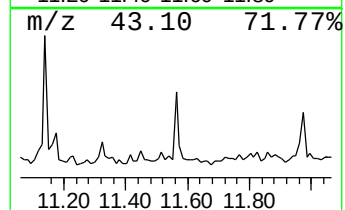
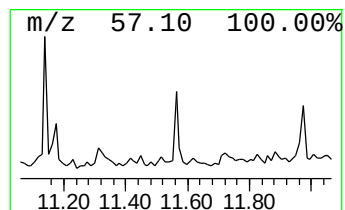
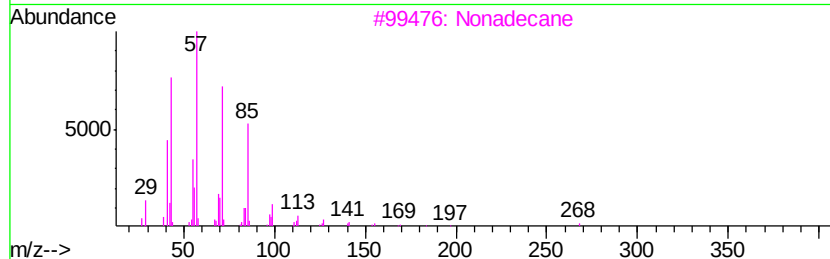
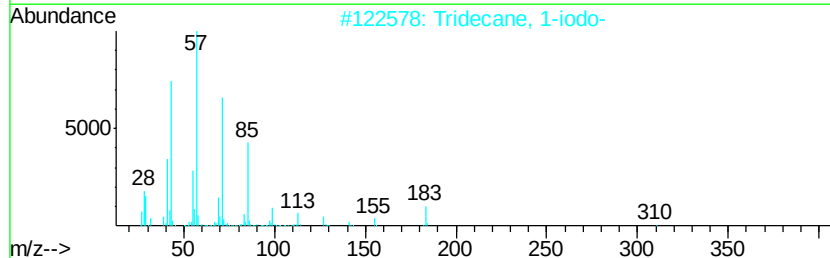
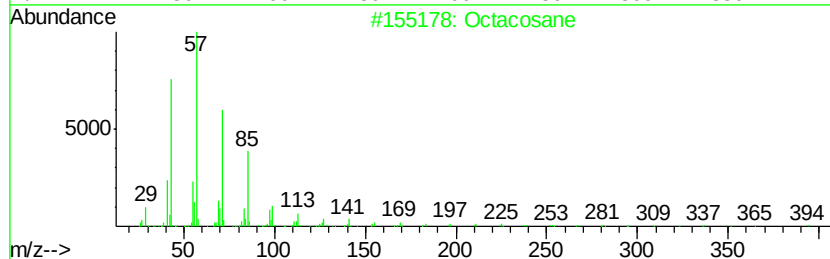
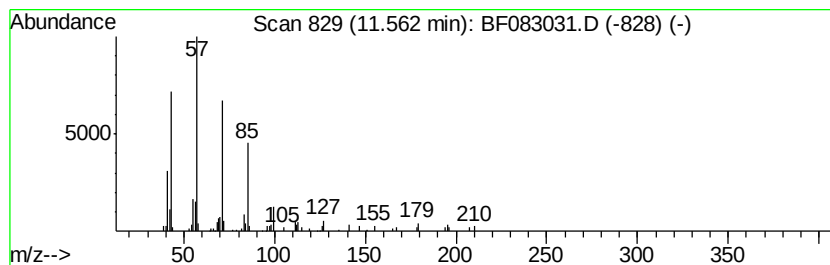
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Octacosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.56	2.26 ng	89231	Phenanthrene-d10	11.25
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octacosane	394 C28H58	000630-02-4	86
2	Tridecane, 1-iodo-	310 C13H27I	035599-77-0	80
3	Nonadecane	268 C19H40	000629-92-5	72
4	Undecane, 2,10-dimethyl-	184 C13H28	017301-27-8	72
5	Dodecane, 2-methyl-	184 C13H28	001560-97-0	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
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Sample : G4426-04
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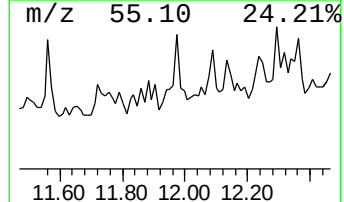
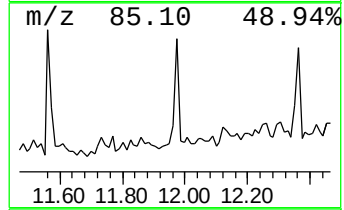
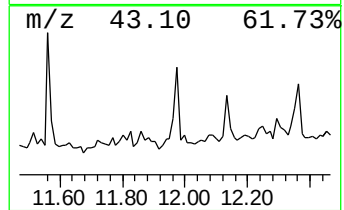
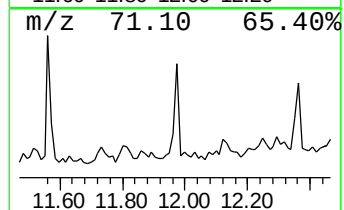
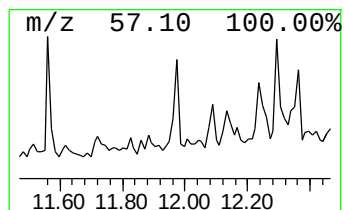
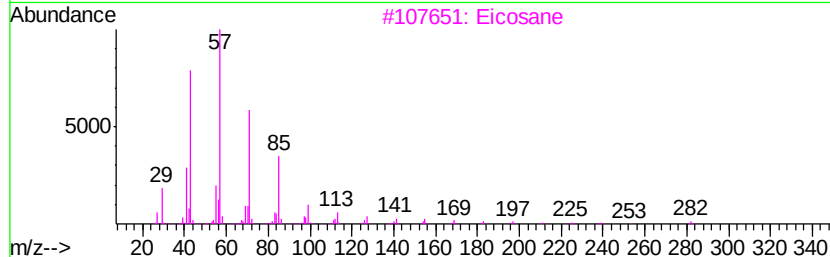
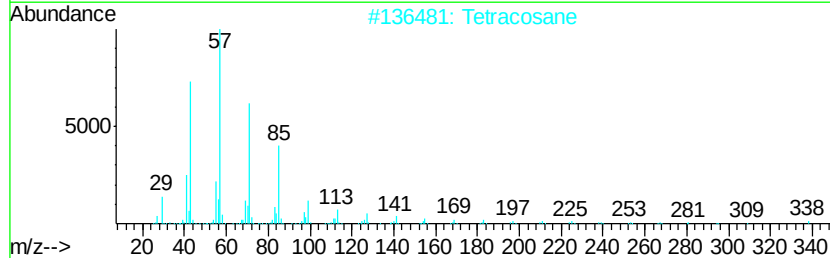
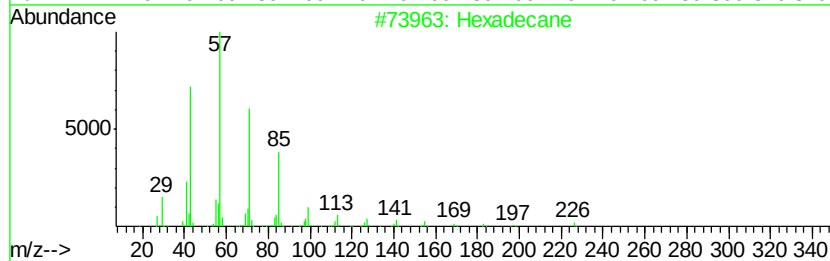
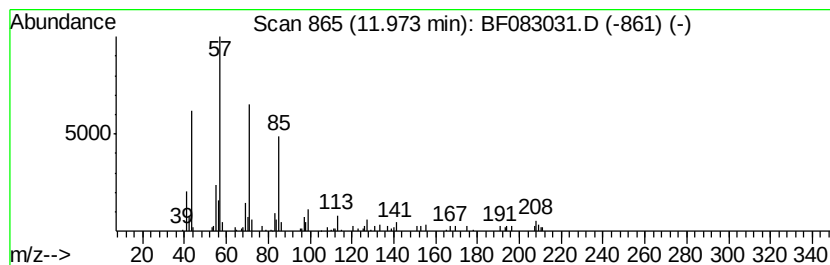
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Hexadecane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD		R.T.		
11.97	2.15 ng	85171	Phenanthrene-d10		11.25		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	83
2			Tetracosane	338	C24H50	000646-31-1	80
3			Eicosane	282	C20H42	000112-95-8	80
4			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	80
5			Pentadecane	212	C15H32	000629-62-9	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
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Acq On : 19 Nov 2015 2:19
Operator : UM/IZ
Sample : G4426-04
Misc :
ALS Vial : 25 Sample Multiplier: 1

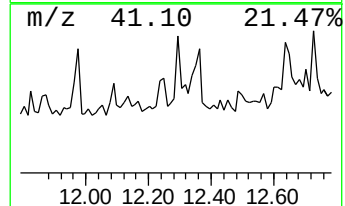
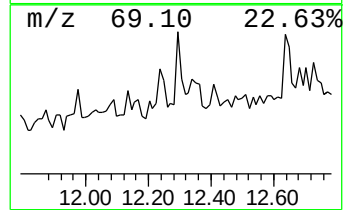
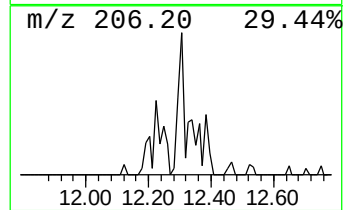
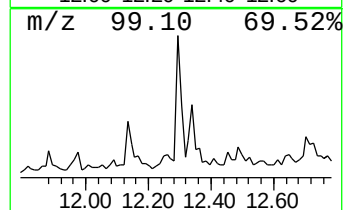
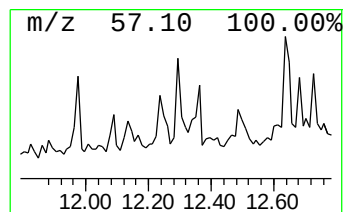
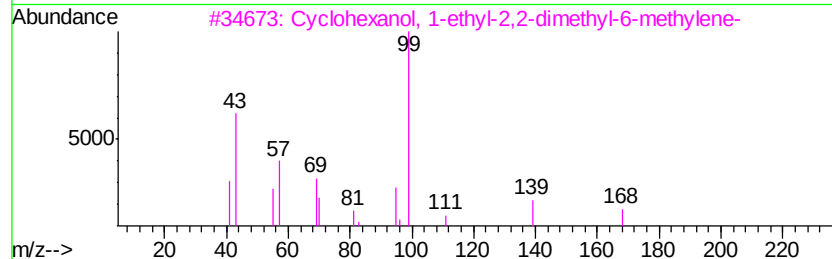
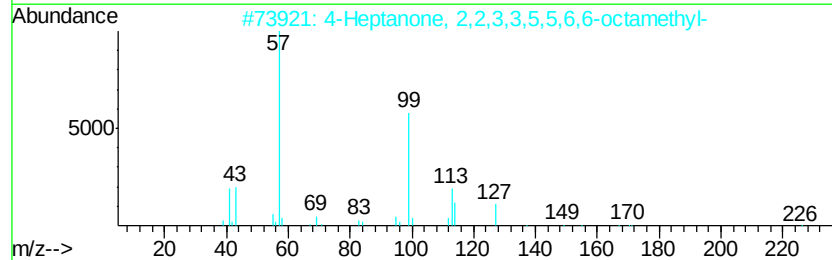
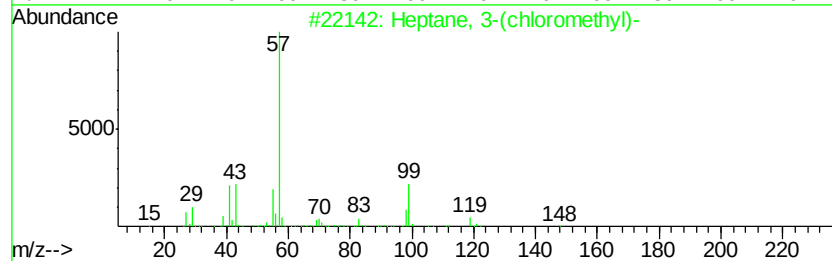
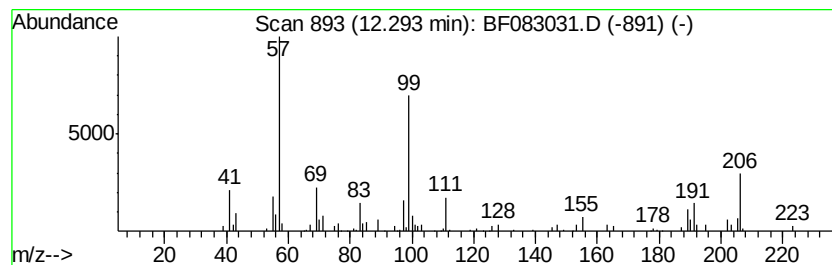
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 unknown12.29 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	2.58 ng	102227	Phenanthrene-d10	11.25
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Heptane, 3-(chloromethyl)-	148 C8H17Cl	000123-04-6 35
2		4-Heptanone, 2,2,3,3,5,5,6,6-oct...	226 C15H30O	016424-66-1 17
3		Cyclohexanol, 1-ethyl-2,2-dimeth...	168 C11H20O	054345-64-1 16
4		2-Piperidinone	99 C5H9NO	000675-20-7 14
5		2-Pyrrolidinone, 1-methyl-	99 C5H9NO	000872-50-4 14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
Data File : BF083031.D
Acq On : 19 Nov 2015 2:19
Operator : UM/IZ
Sample : G4426-04
Misc :
ALS Vial : 25 Sample Multiplier: 1

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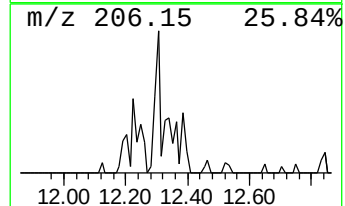
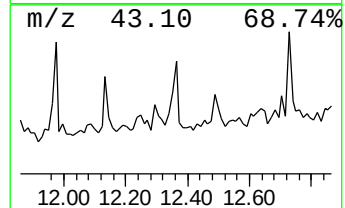
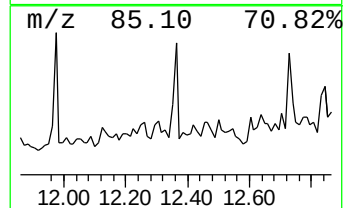
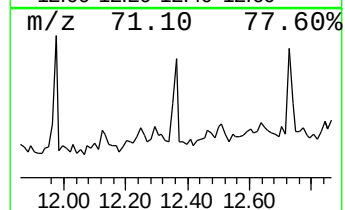
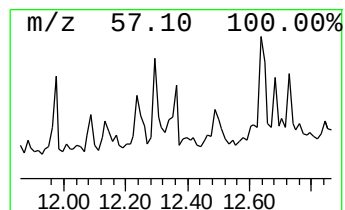
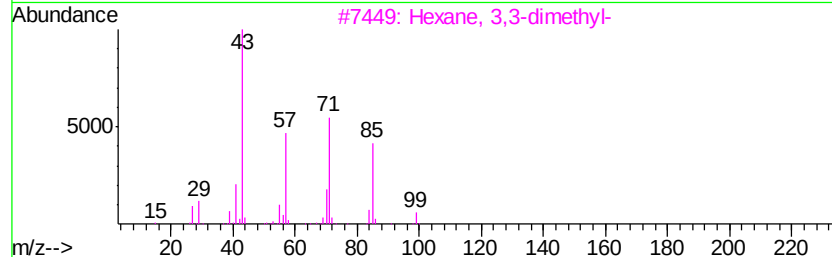
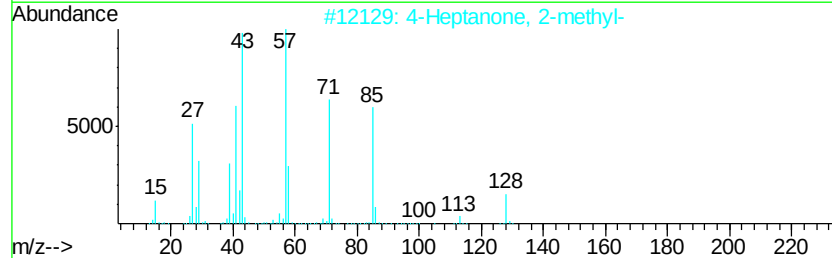
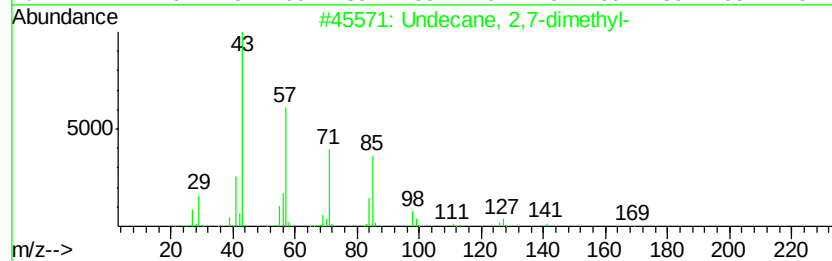
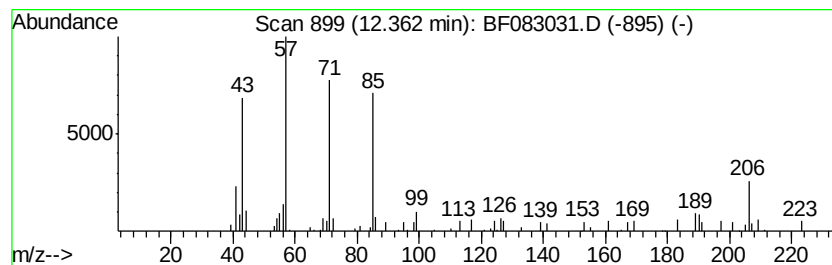
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Undecane, 2,7-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	2.30 ng	90877	Phenanthrene-d10	11.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Undecane, 2,7-dimethyl-	184	C13H28	017301-24-5	59
2		4-Heptanone, 2-methyl-	128	C8H16O	000626-33-5	53
3		Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	53
4		Tetradecane	198	C14H30	000629-59-4	43
5		4-Heptanone, 3-methyl-	128	C8H16O	015726-15-5	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
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Acq On : 19 Nov 2015 2:19
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Sample : G4426-04
Misc :
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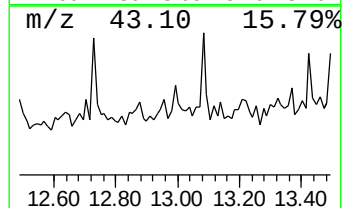
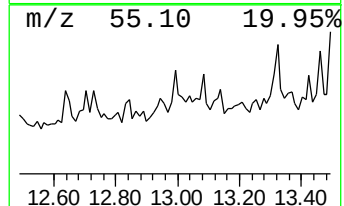
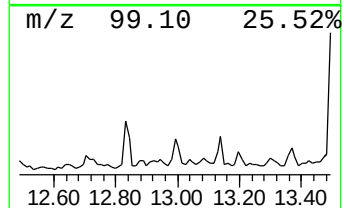
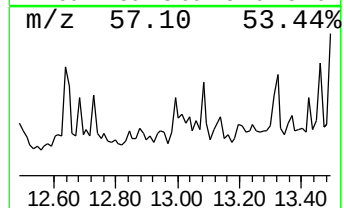
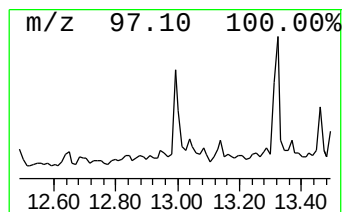
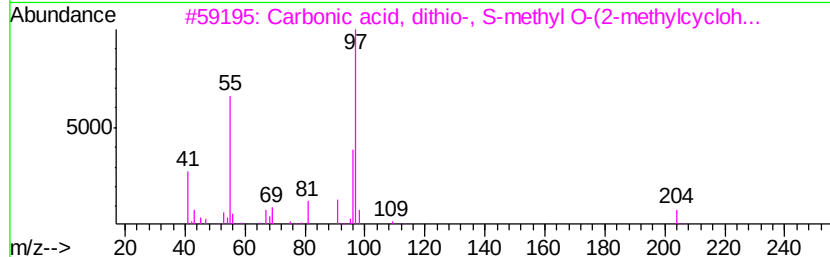
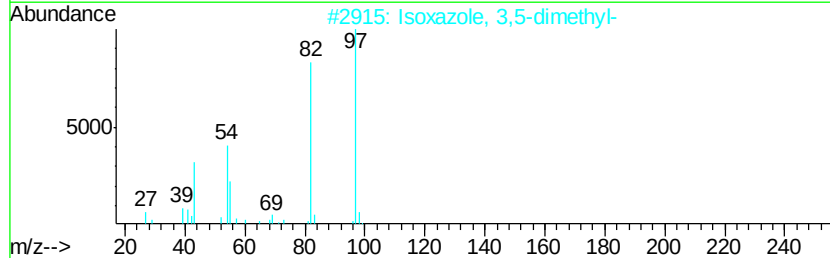
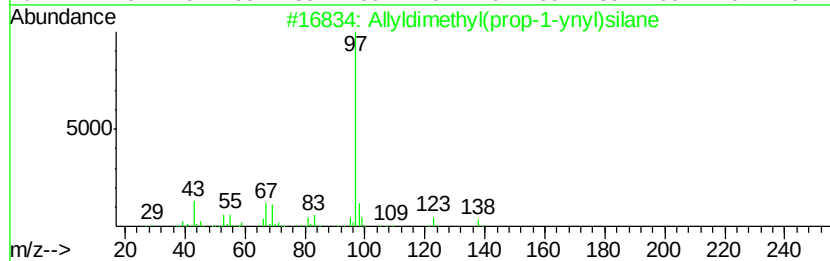
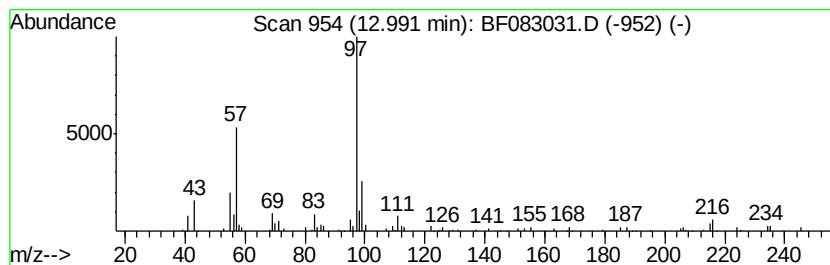
Instrument :
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Allyldimethyl(prop-1-ynyl)s... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.99	2.65 ng	100077	Chrysene-d12	13.88	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Allyldimethyl(prop-1-ynyl)silane	138	C8H14Si	1000281-75-1	53
2	Isoxazole, 3,5-dimethyl-	97	C5H7NO	000300-87-8	50
3	Carbonic acid, dithio-, S-methyl...	204	C9H16OS2	015288-12-7	42
4	2-Hexenoic acid, 2-hexenyl ester...	196	C12H20O2	054845-28-2	42
5	4H-1,2,4-Triazole, 4-ethyl-	97	C4H7N3	043183-55-7	40



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
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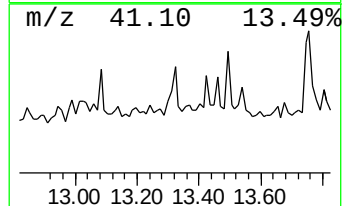
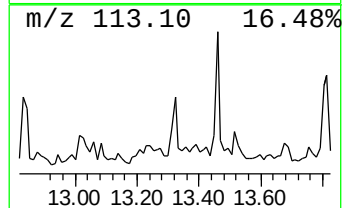
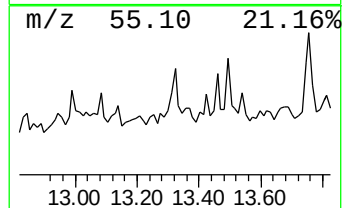
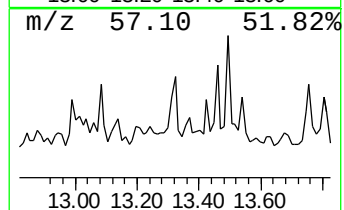
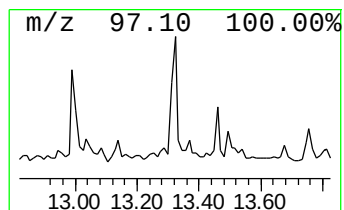
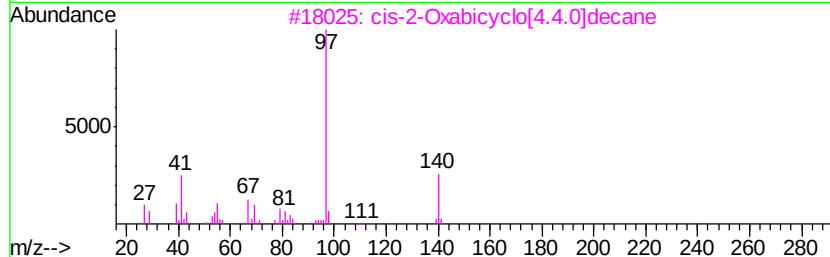
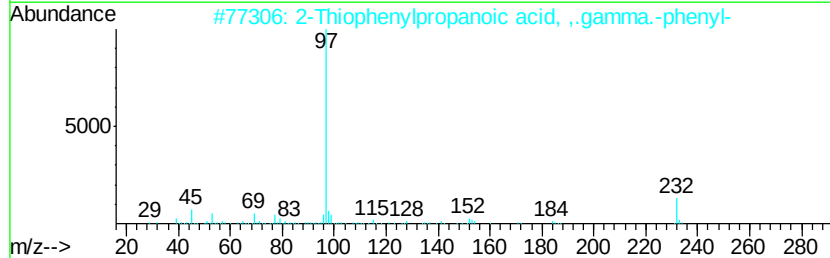
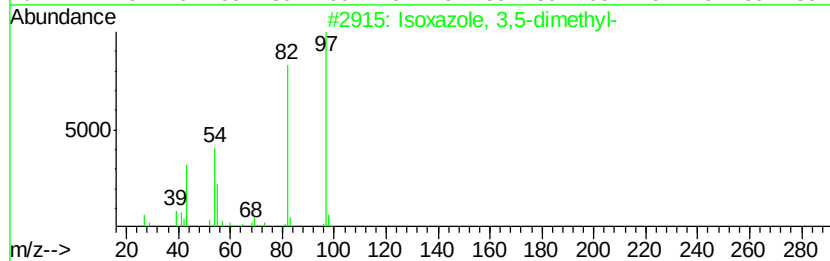
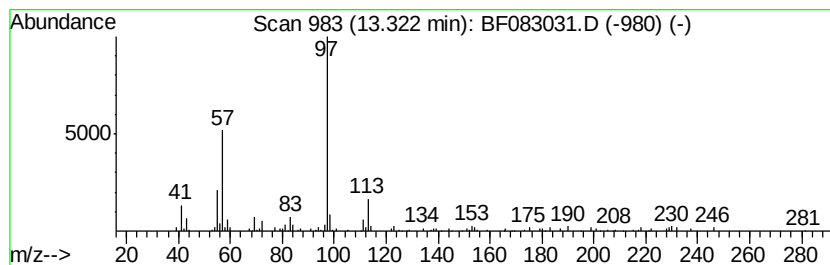
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Isoxazole, 3,5-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.32	2.47 ng	93175	Chrysene-d12	13.88
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Isoxazole, 3,5-dimethyl-	97 C5H7NO	000300-87-8 58
2		2-Thiophenylpropanoic acid, ,.ga...	232 C13H12O2S	051535-43-4 45
3		cis-2-Oxabicyclo[4.4.0]decane	140 C9H16O	060416-19-5 42
4		4-Methyl-2,3-hexadien-1-ol	112 C7H12O	1000196-09-6 42
5		Cyclohexanone, 3-butyl-	154 C10H18O	039178-69-3 42



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
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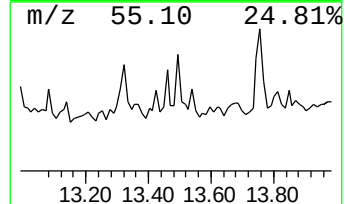
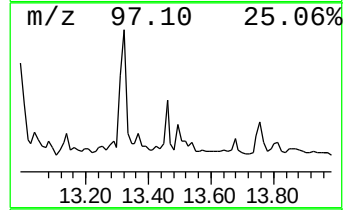
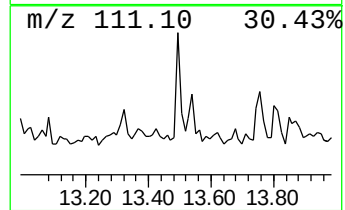
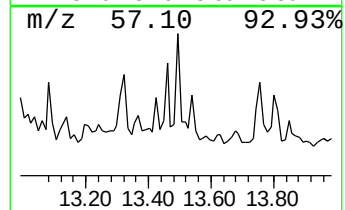
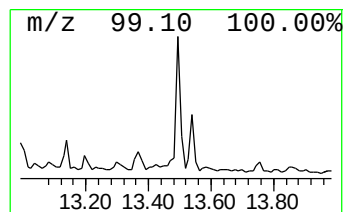
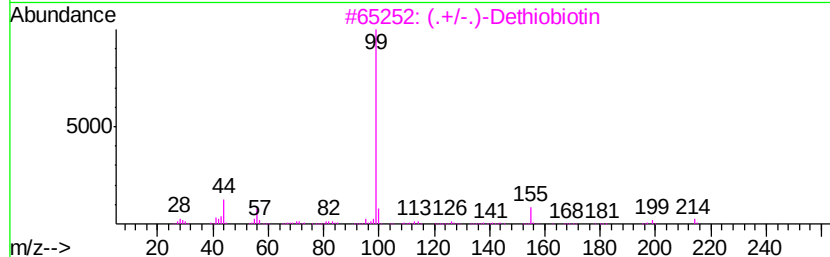
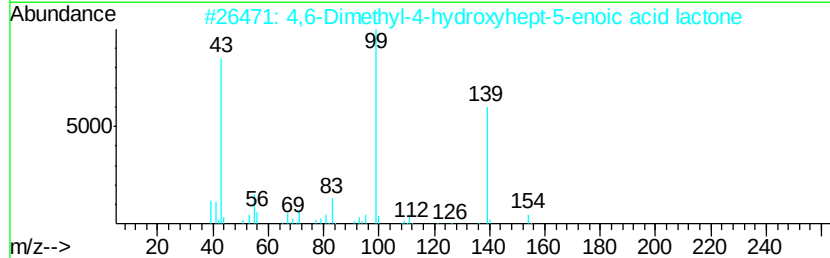
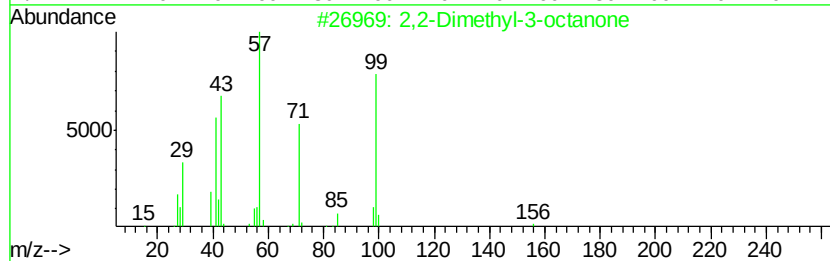
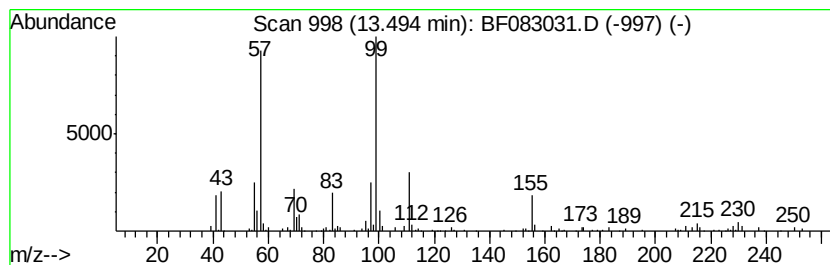
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 unknown13.49 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.49	3.31 ng	124998	Chrysene-d12	13.88	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2-Dimethyl-3-octanone	156	C10H20O	005340-64-7	47
2	4,6-Dimethyl-4-hydroxyhept-5-eno...	154	C9H14O2	1000122-47-0	37
3	(.+/-)-Dethiobiotin	214	C10H18N2O3	000636-20-4	37
4	Formaldehyde, dipropylhydrazone	128	C7H16N2	054253-56-4	37
5	2-Propylthiazole	127	C6H9NS	017626-75-4	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
Data File : BF083031.D
Acq On : 19 Nov 2015 2:19
Operator : UM/IZ
Sample : G4426-04
Misc :
ALS Vial : 25 Sample Multiplier: 1

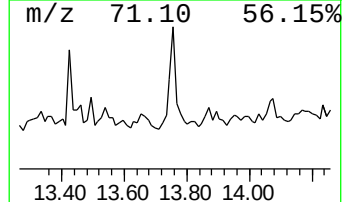
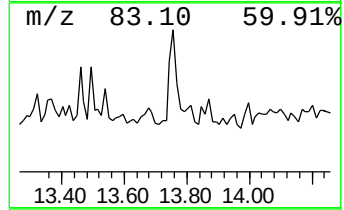
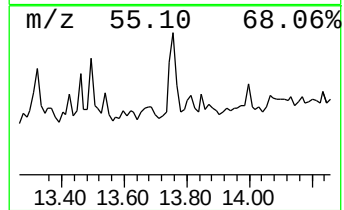
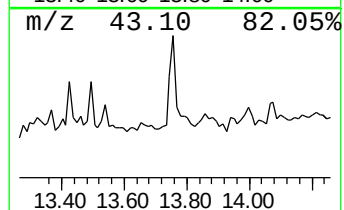
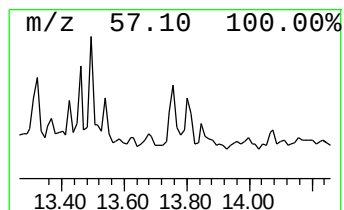
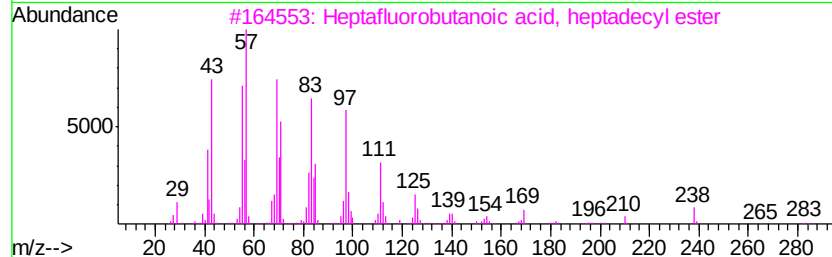
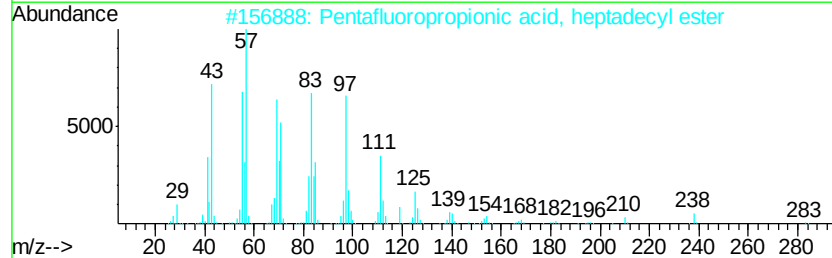
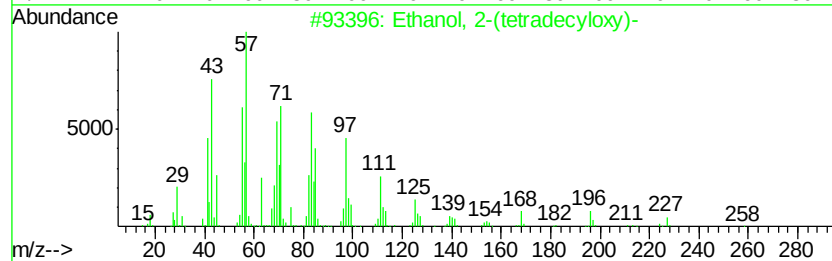
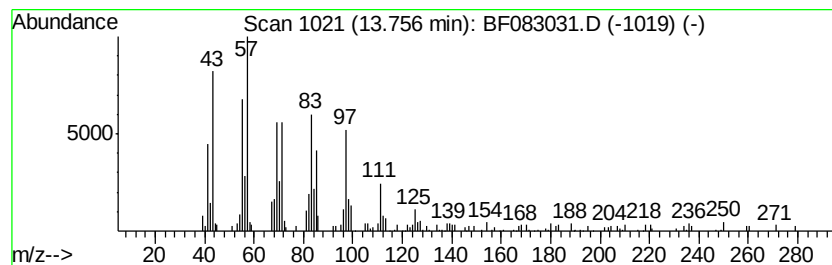
Instrument :
BNA_F
ClientSampleId :
SB-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Ethanol, 2-(tetradecyloxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.76	3.63 ng	137017	Chrysene-d12	13.88
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Ethanol, 2-(tetradecyloxy)-	258 C16H34O2	002136-70-1 87
2		Pentafluoropropionic acid, hepta...	402 C20H35F5O2	1000283-04-2 87
3		Heptafluorobutanoic acid, heptad...	452 C21H35F7O2	1000282-97-3 87
4		Dichloroacetic acid, heptadecyl ...	366 C19H36Cl2O2	1000282-98-2 76
5		17-Pentatriacontene	491 C35H70	006971-40-0 74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111815\
Data File : BF083031.D
Acq On : 19 Nov 2015 2:19
Operator : UM/IZ
Sample : G4426-04
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.90	52.6	ng	1932420	1	6.74	734739	20.0
unknown6.51	6.51	114.6	ng	4208310	1	6.74	734739	20.0
Hexanoic acid, 2-...	7.42	2.5	ng	117849	2	8.02	955621	20.0
Tetradecane	10.22	2.9	ng	137467	3	9.77	941071	20.0
Tridecane	10.69	6.2	ng	243459	4	11.25	791346	20.0
Octacosane	11.56	2.3	ng	89231	4	11.25	791346	20.0
Hexadecane	11.97	2.1	ng	85171	4	11.25	791346	20.0
unknown12.29	12.29	2.6	ng	102227	4	11.25	791346	20.0
Undecane, 2,7-dim...	12.36	2.3	ng	90877	4	11.25	791346	20.0
Allyldimethyl(pro...	12.99	2.6	ng	100077	5	13.88	755011	20.0
Isoxazole, 3,5-di...	13.32	2.5	ng	93175	5	13.88	755011	20.0
unknown13.49	13.49	3.3	ng	124998	5	13.88	755011	20.0
Ethanol, 2-(tetra...	13.76	3.6	ng	137017	5	13.88	755011	20.0