

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122515\
 Data File : BF084146.D
 Acq On : 25 Dec 2015 23:31
 Operator : UM/SJ
 Sample : G4920-06
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 IM-CLAY-011

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-BF122415.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	4.978	251	253	263	rBB	240739	346345	18.57%	2.336%
2	5.424	289	292	294	rBV	1188137	1603188	85.97%	10.815%
3	6.453	380	382	385	rBV	1333077	1445155	77.49%	9.749%
4	6.579	390	393	395	rBV	1328102	1728492	92.69%	11.661%
5	6.796	410	412	415	rBV	401721	330831	17.74%	2.232%
6	6.956	424	426	428	rBV	1561751	1401587	75.16%	9.455%
7	7.367	459	462	464	rBV	1081010	1129595	60.57%	7.620%
8	8.087	522	525	527	rBV	304159	397460	21.31%	2.681%
9	9.162	616	619	621	rBV	2043459	1864888	100.00%	12.581%
10	9.550	651	653	656	rBV	265352	279461	14.99%	1.885%
11	9.836	675	678	680	rBV	460593	421882	22.62%	2.846%
12	10.270	708	716	718	rBV	41611	51623	2.77%	0.348%
13	10.488	733	735	739	rBV2	12553	23170	1.24%	0.156%
14	10.625	744	747	749	rBV	914387	871869	46.75%	5.882%
15	10.739	756	757	761	rBV	34452	49156	2.64%	0.332%
16	11.185	795	796	798	rBV2	17546	24189	1.30%	0.163%
17	11.322	806	808	810	rBV	272327	354818	19.03%	2.394%
18	12.728	929	931	933	rBV	103081	92045	4.94%	0.621%
19	12.796	933	937	939	rBV2	23504	42548	2.28%	0.287%
20	12.899	943	946	948	rBV	1376338	1209852	64.88%	8.162%
21	13.425	990	992	994	rBV	65921	69045	3.70%	0.466%
22	13.471	994	996	1001	rVV2	15083	39558	2.12%	0.267%
23	13.676	1011	1014	1016	rBV	9589	21007	1.13%	0.142%
24	13.802	1022	1025	1028	rBV2	94881	147532	7.91%	0.995%
25	13.951	1036	1038	1041	rBV	311479	354560	19.01%	2.392%
26	14.122	1050	1053	1055	rBV	41896	48007	2.57%	0.324%
27	14.419	1077	1079	1082	rBV	53404	67703	3.63%	0.457%
28	14.717	1103	1105	1108	rBV	38929	57630	3.09%	0.389%
29	15.379	1161	1163	1166	rBV	223825	294657	15.80%	1.988%
30	16.248	1237	1239	1241	rBV3	13321	24104	1.29%	0.163%
31	16.797	1285	1287	1293	rBV5	13438	31361	1.68%	0.212%

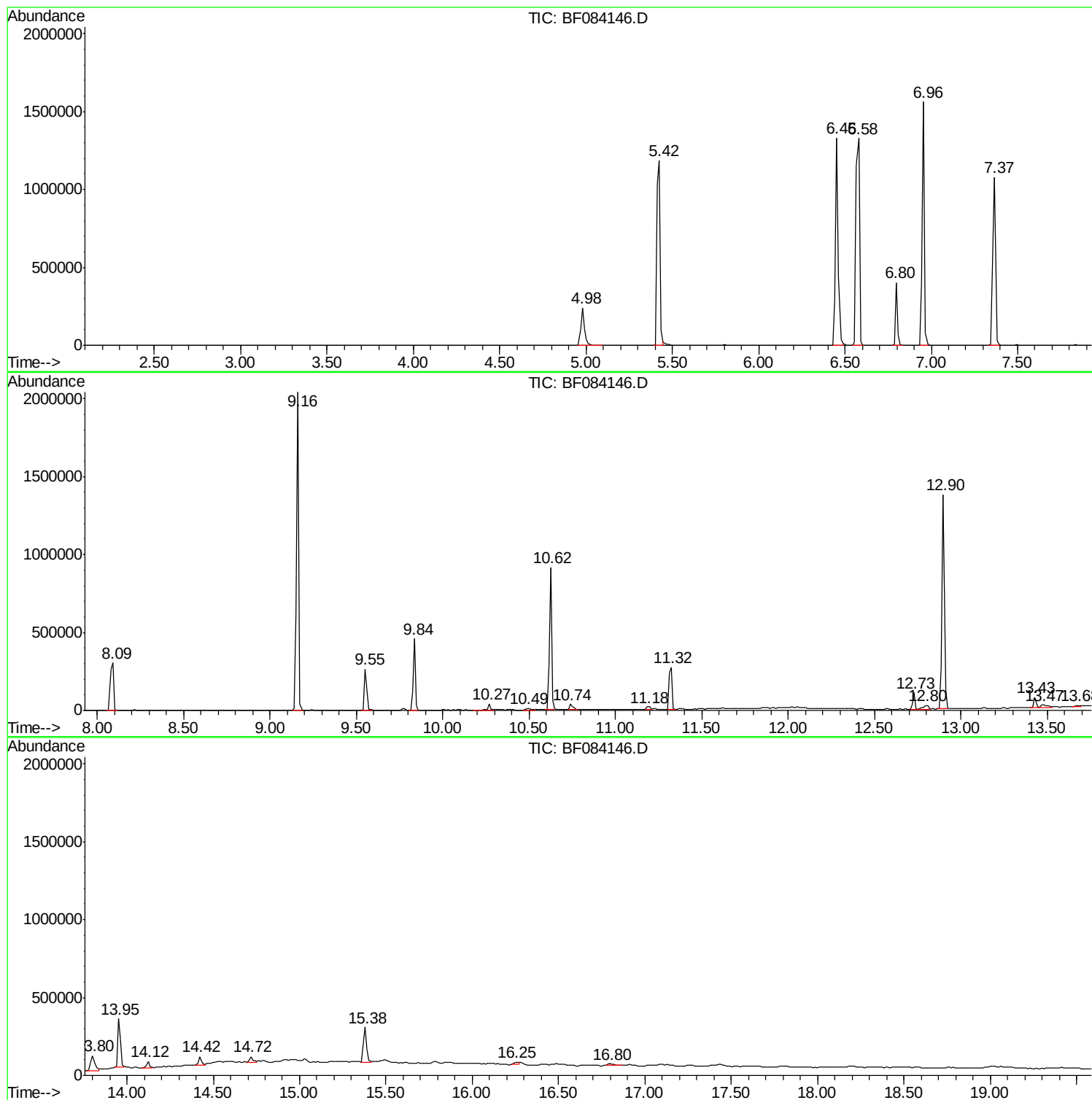
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Misc :
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Instrument :
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ClientSampleId :
IM-CLAY-011

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122415.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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Instrument :
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ClientSampled :
IM-CLAY-011

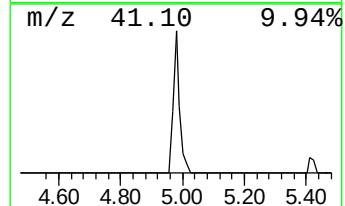
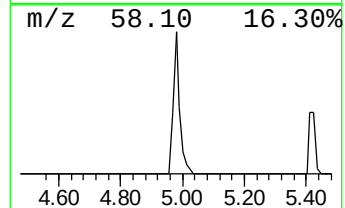
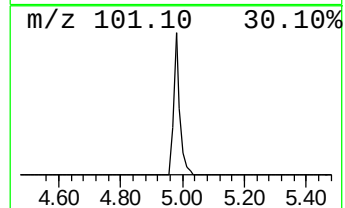
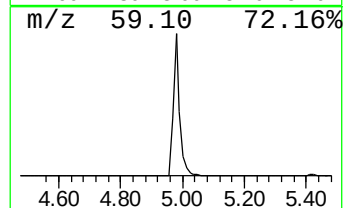
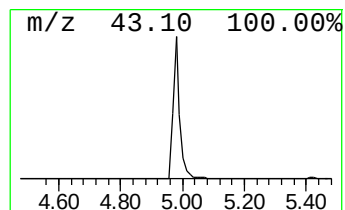
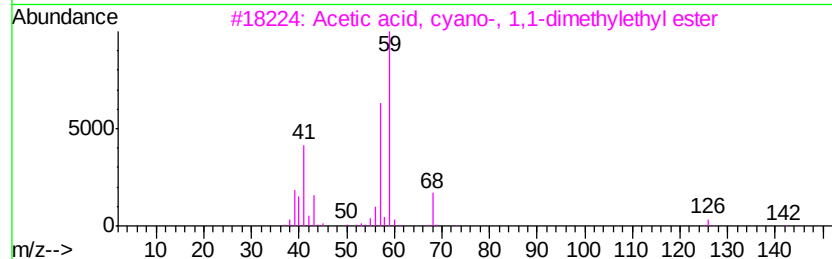
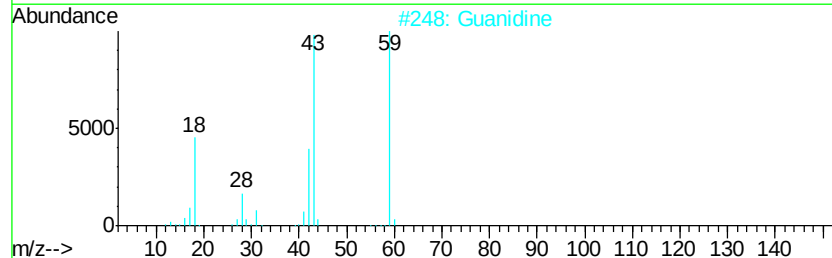
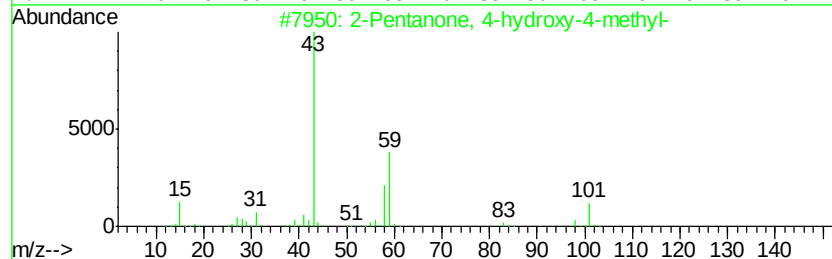
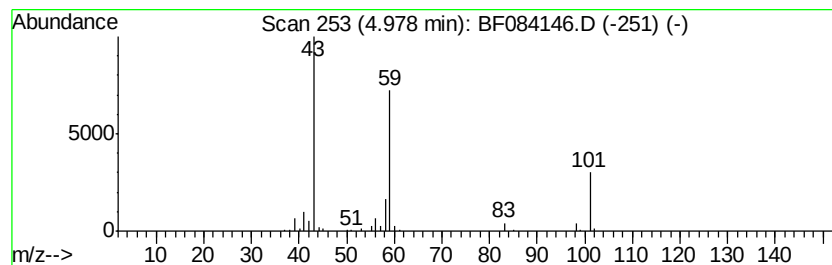
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
4.98	20.94 ng	346345	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Guanidine	59	CH5N3	000113-00-8	43
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4		Acetone	58	C3H6O	000067-64-1	10
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



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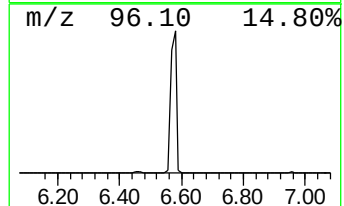
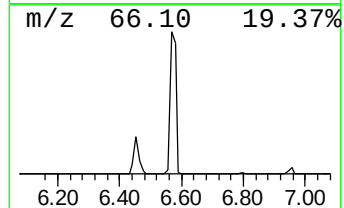
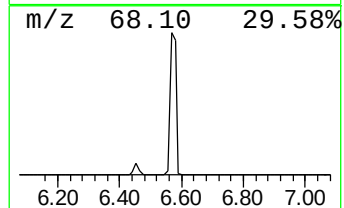
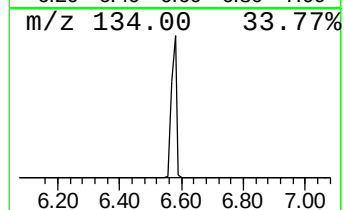
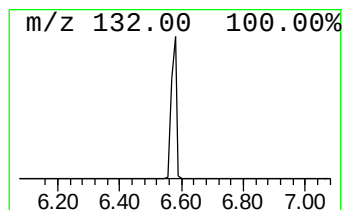
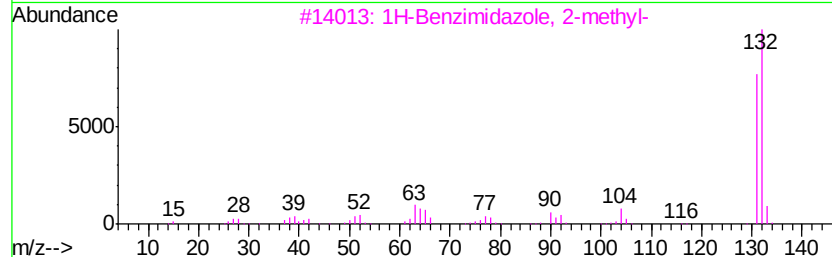
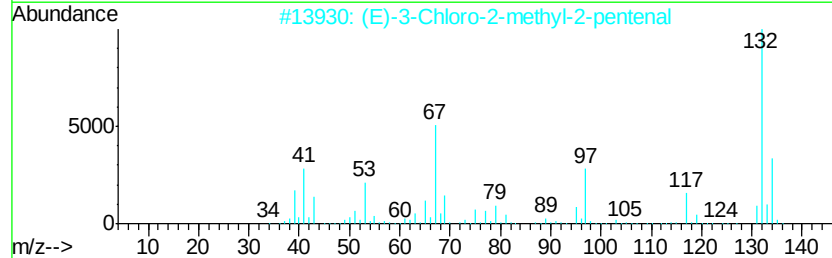
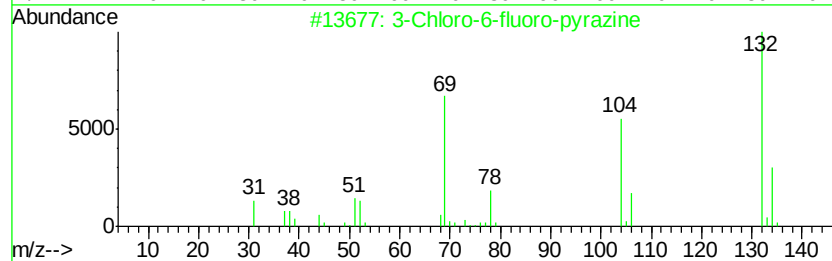
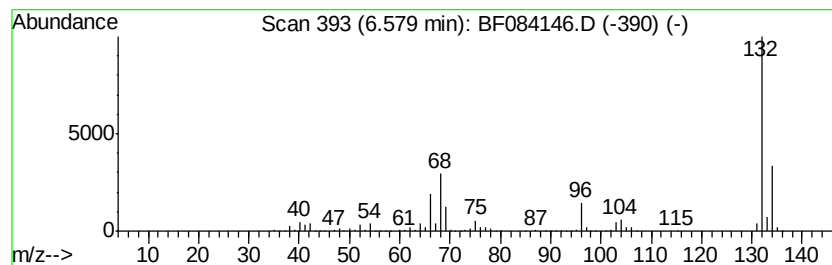
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Peak Number 2 unknown6.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	104.49 ng	1728490	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	12
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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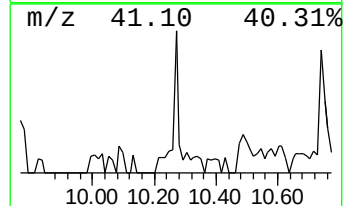
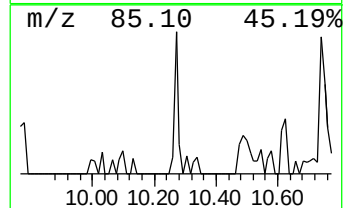
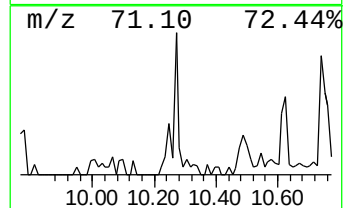
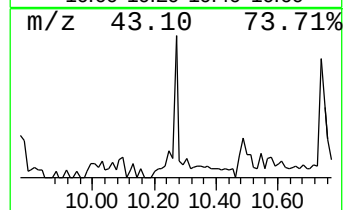
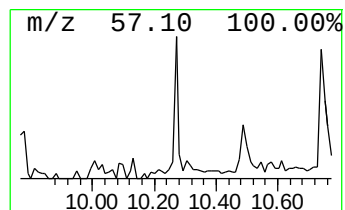
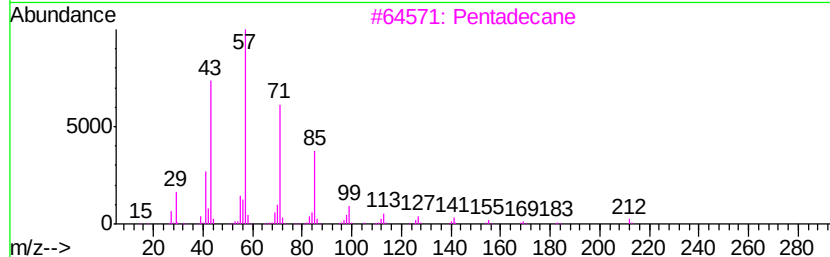
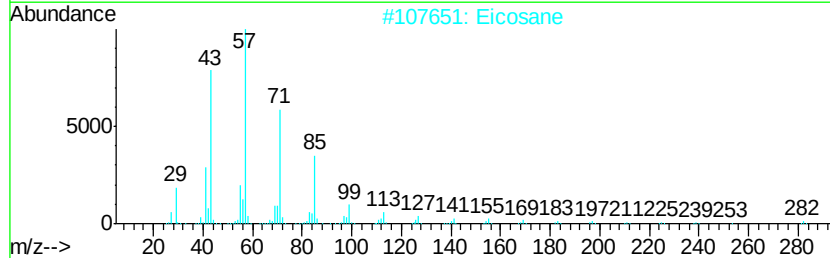
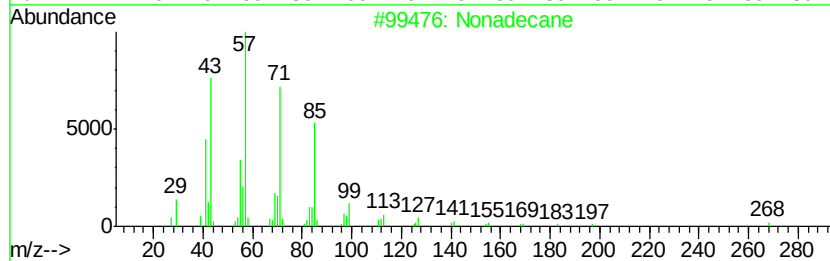
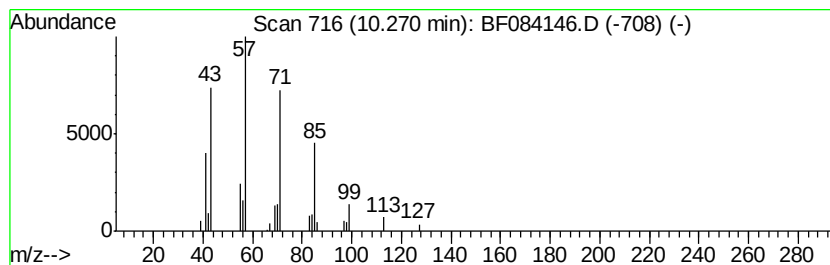
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BNA_F
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Peak Number 4 Nonadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.27	2.45 ng	51623	Acenaphthene-d10	9.84
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Nonadecane	268 C19H40	000629-92-5	91
2	Eicosane	282 C20H42	000112-95-8	86
3	Pentadecane	212 C15H32	000629-62-9	86
4	Tetradecane	198 C14H30	000629-59-4	86
5	Hexadecane	226 C16H34	000544-76-3	78



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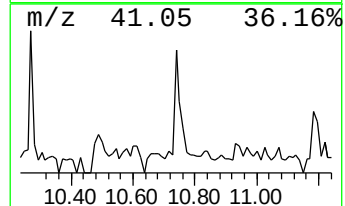
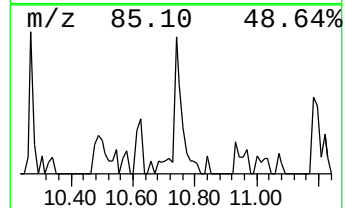
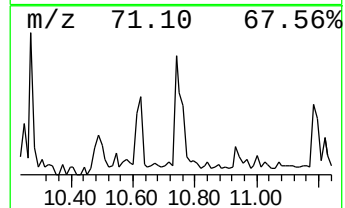
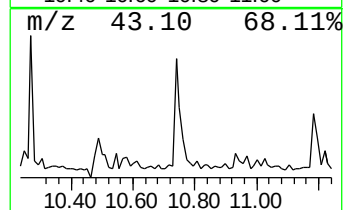
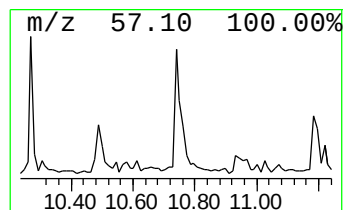
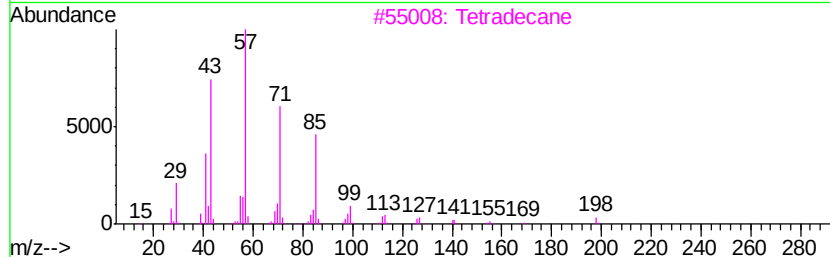
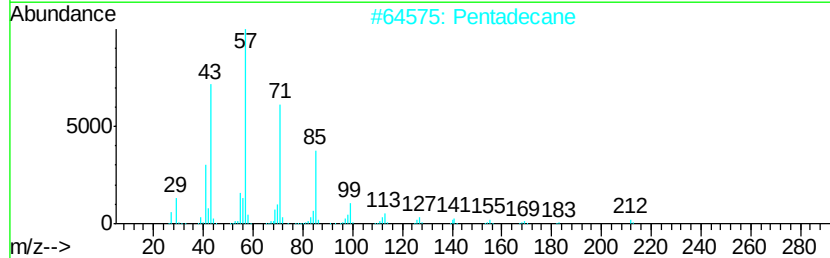
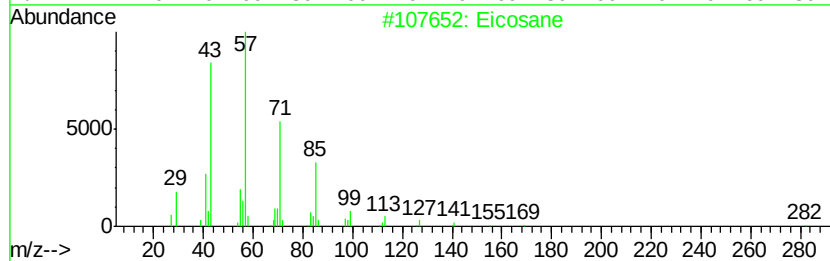
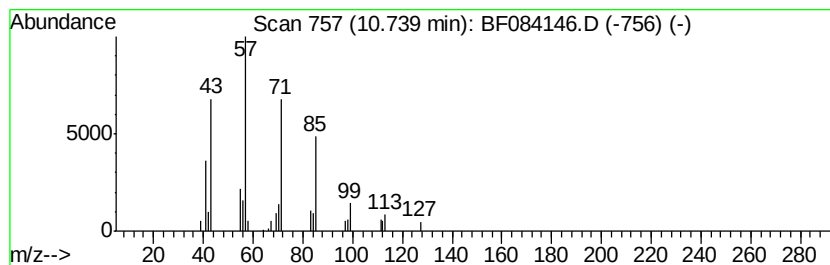
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 Eicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.74	2.77 ng	49156	Phenanthrene-d10	11.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	90
2	Pentadecane		212	C15H32	000629-62-9	90
3	Tetradecane		198	C14H30	000629-59-4	90
4	Heptadecane		240	C17H36	000629-78-7	90
5	Tridecane		184	C13H28	000629-50-5	83



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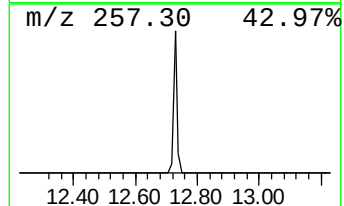
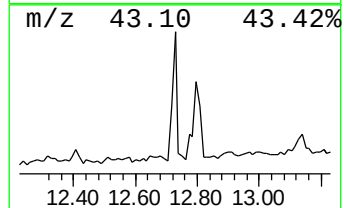
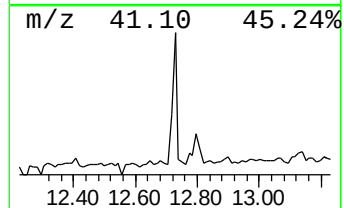
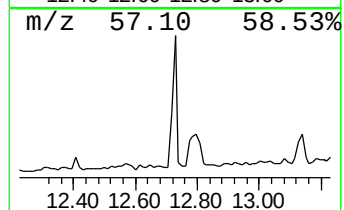
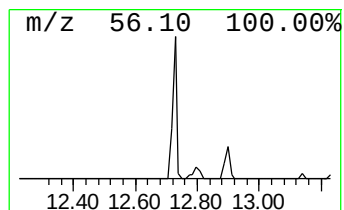
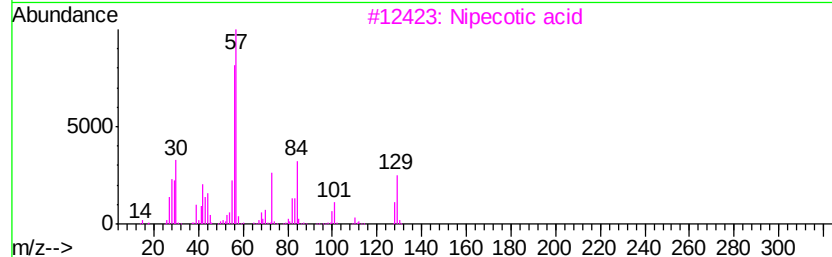
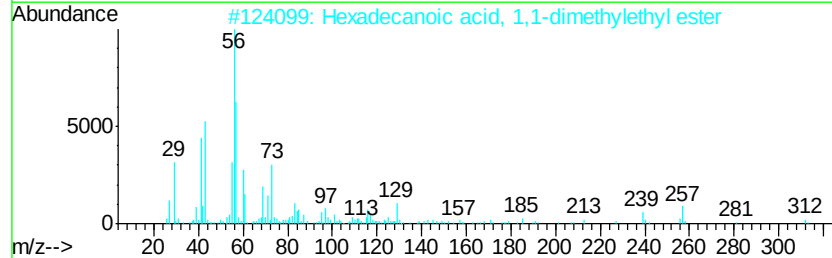
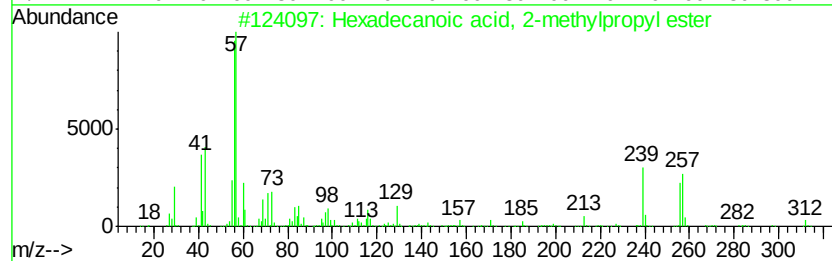
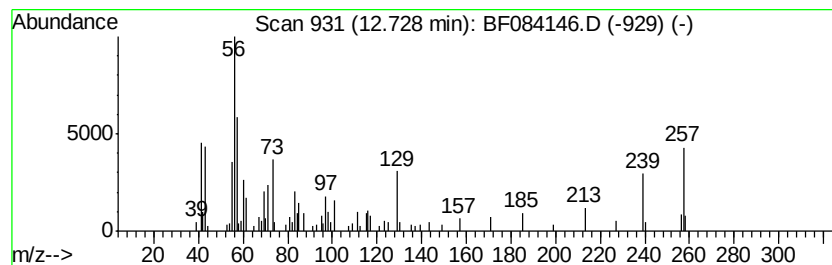
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 6 Hexadecanoic acid, 2-methyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.73	5.19 ng	92045	Chrysene-d12	13.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, 2-methylpropy...	312	C20H40O2	000110-34-9	53
2		Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	49
3		Nipecotic acid	129	C6H11NO2	000498-95-3	38
4		1-Heptene, 2-methyl-	112	C8H16	015870-10-7	25
5		Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	25



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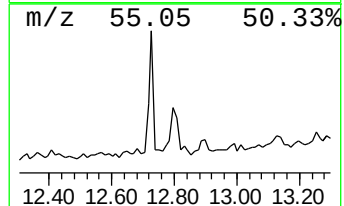
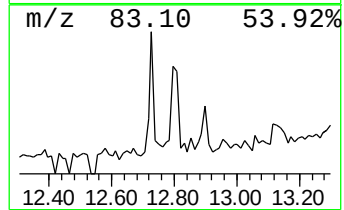
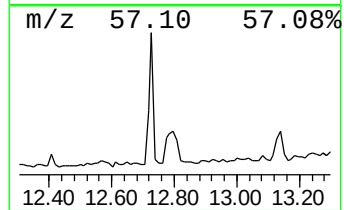
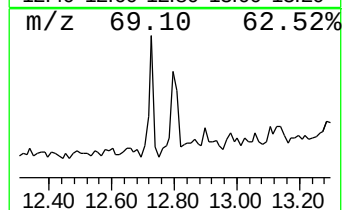
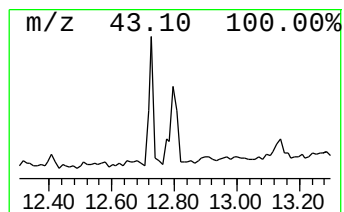
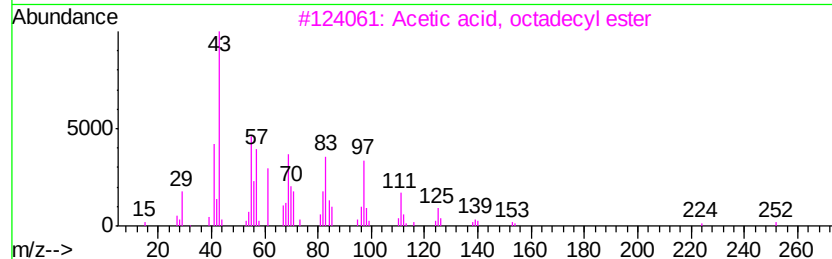
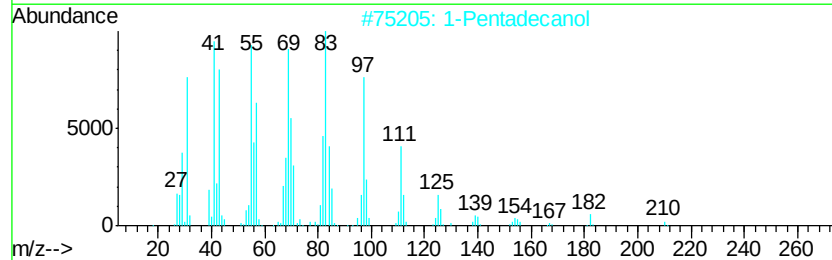
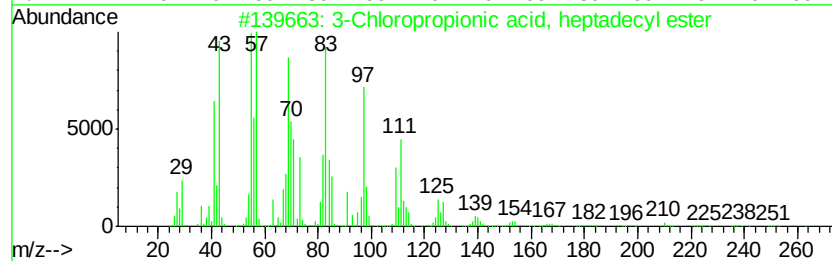
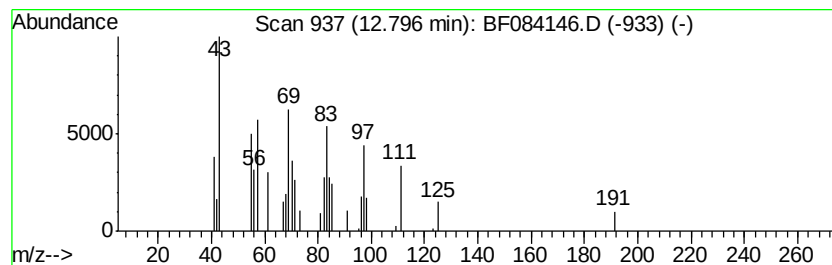
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 3-Chloropropionic acid, hep... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.80	2.40 ng	42548	Chrysene-d12	13.95		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	81
2		1-Pentadecanol	228	C15H32O	000629-76-5	74
3		Acetic acid, octadecyl ester	312	C20H40O2	000822-23-1	72
4		Chloroacetic acid, tridecyl ester	276	C15H29ClO2	018277-85-5	64
5		1-Octadecene	252	C18H36	000112-88-9	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122515\
Data File : BF084146.D
Acq On : 25 Dec 2015 23:31
Operator : UM/SJ
Sample : G4920-06
Misc :
ALS Vial : 63 Sample Multiplier: 1

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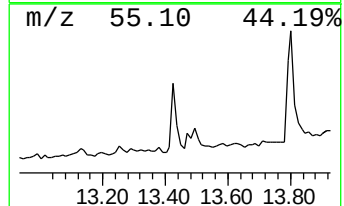
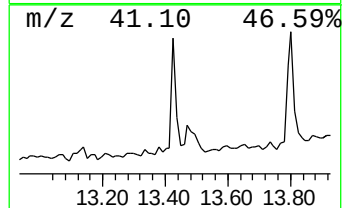
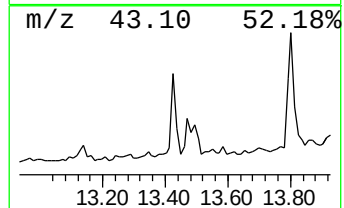
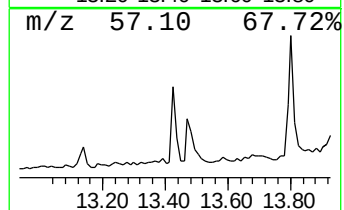
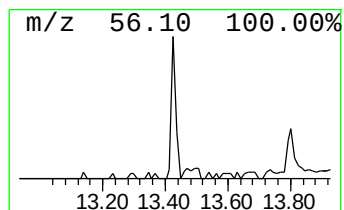
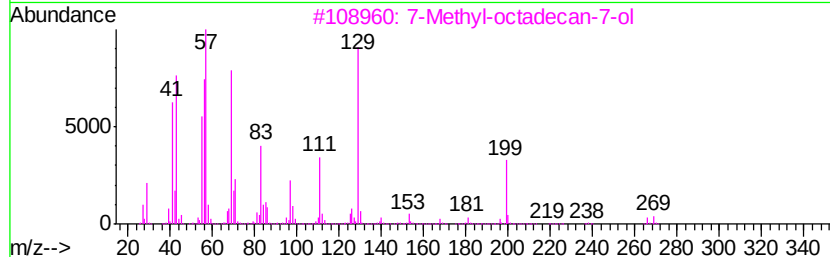
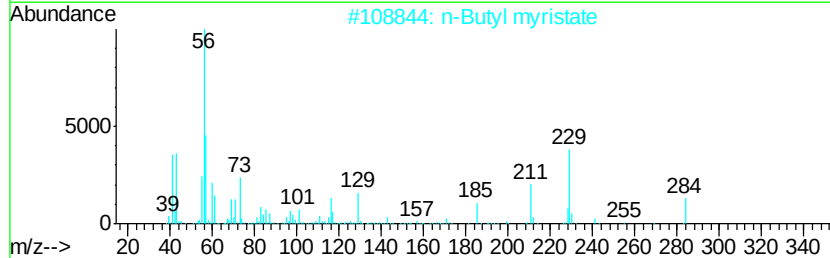
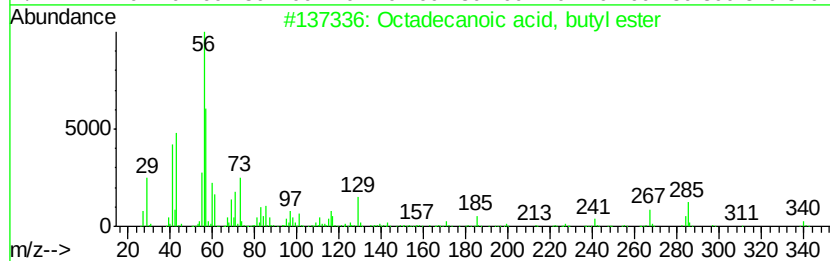
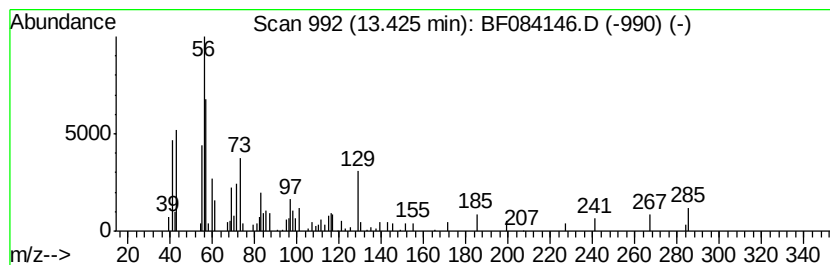
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecanoic acid, butyl ester Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.43	3.89 ng	69045	Chrysene-d12	13.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	83
2		n-Butyl myristate	284	C18H36O2	000110-36-1	56
3		7-Methyl-octadecan-7-ol	284	C19H40O	1000194-50-8	38
4		Octadecanoic acid	284	C18H36O2	000057-11-4	38
5		1-Pentene, 2,4-dimethyl-	98	C7H14	002213-32-3	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122515\
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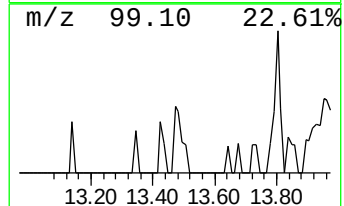
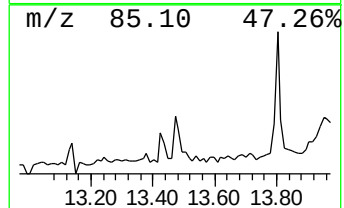
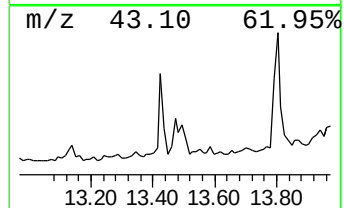
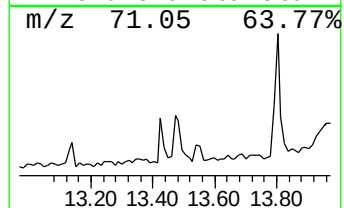
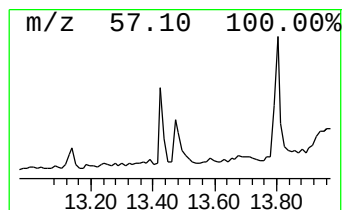
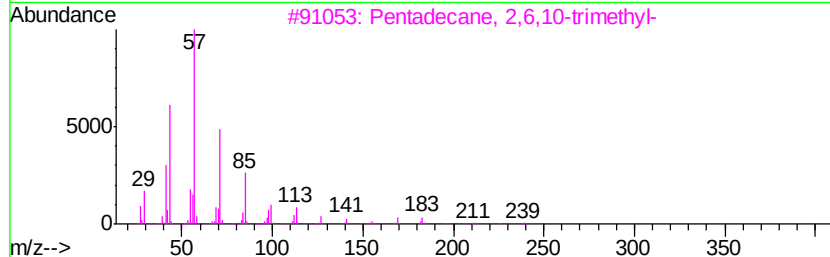
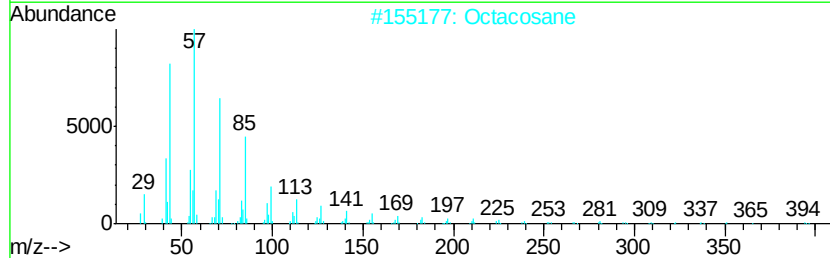
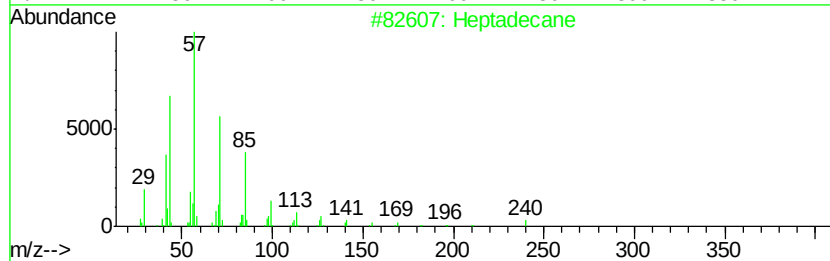
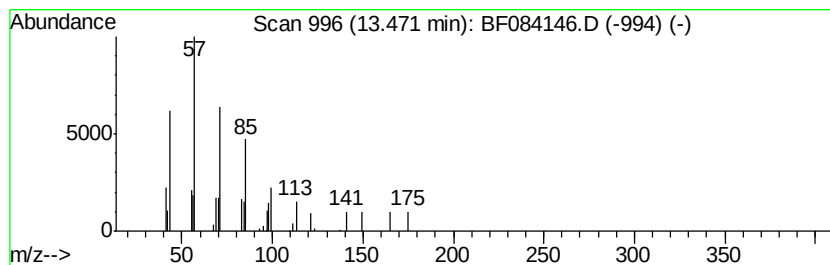
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	2.23 ng	39558	Chrysene-d12	13.95
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptadecane	240 C17H36	000629-78-7	53
2	Octacosane	394 C28H58	000630-02-4	53
3	Pentadecane, 2,6,10-trimethyl-	254 C18H38	003892-00-0	53
4	Dodecane	170 C12H26	000112-40-3	50
5	2-Bromo dodecane	248 C12H25Br	013187-99-0	50



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122515\
Data File : BF084146.D
Acq On : 25 Dec 2015 23:31
Operator : UM/SJ
Sample : G4920-06
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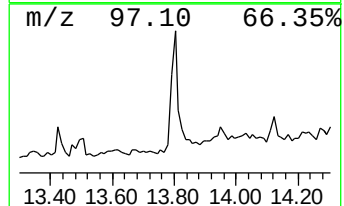
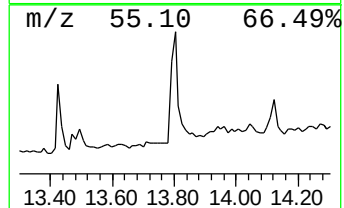
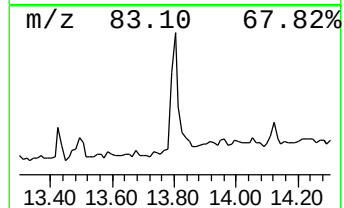
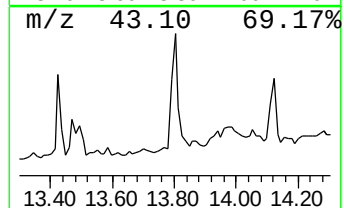
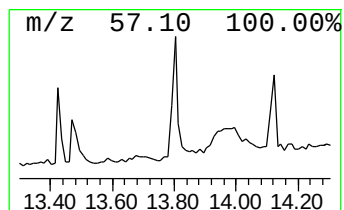
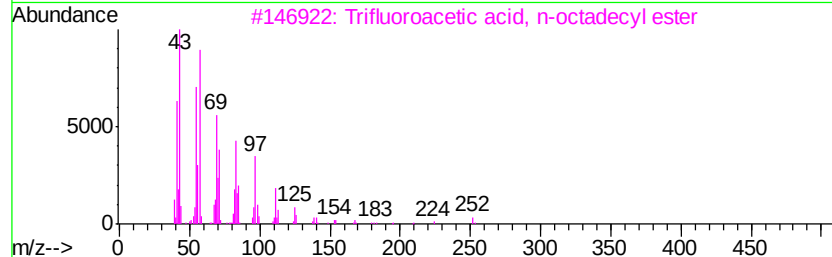
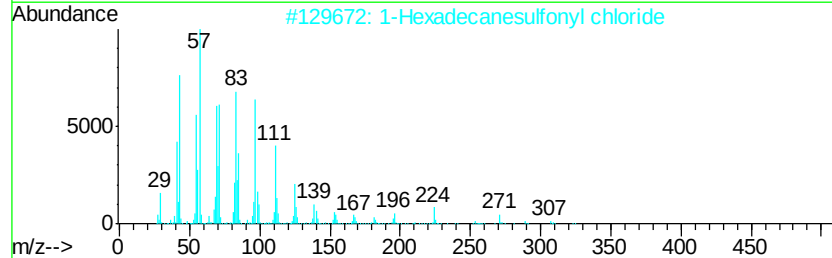
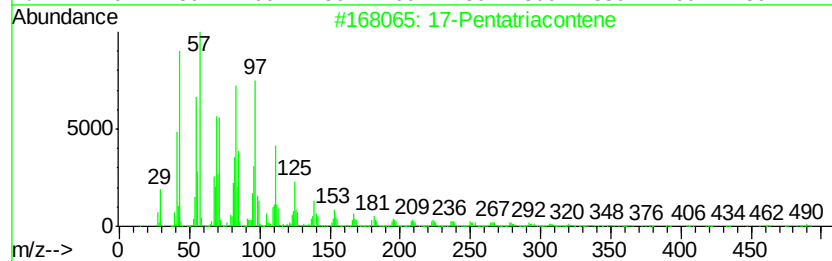
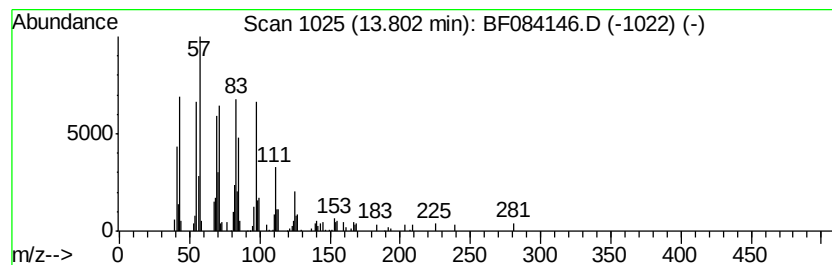
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 17-Pentatriacontene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.80	8.32 ng	147532	Chrysene-d12	13.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	17-Pentatriacontene	491	C35H70	006971-40-0	91
2		1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	87
3		Trifluoroacetic acid, n-octadecyl...	366	C20H37F3O2	079392-43-1	87
4		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	81
5		Trihexadecyl borate	735	C48H99B03	002665-11-4	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122515\
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Acq On : 25 Dec 2015 23:31
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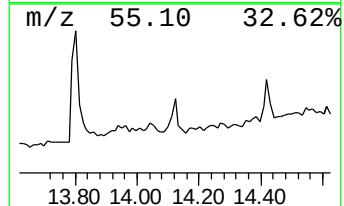
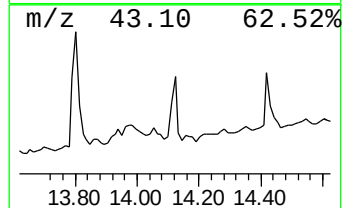
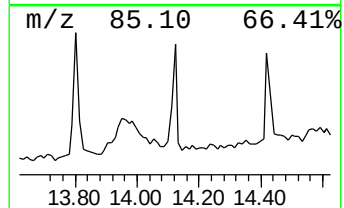
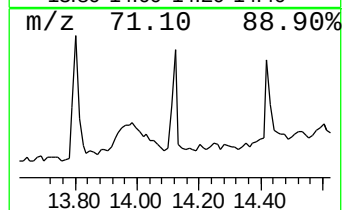
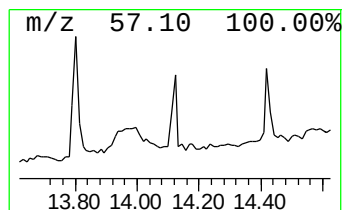
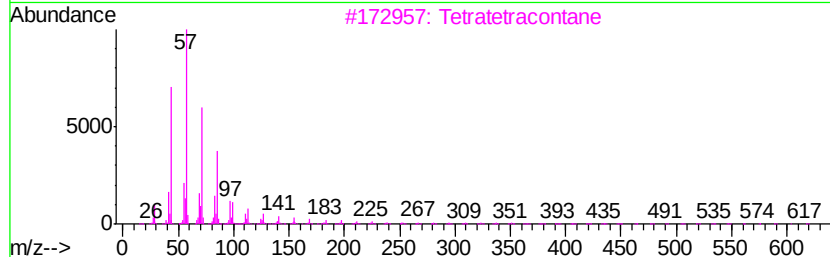
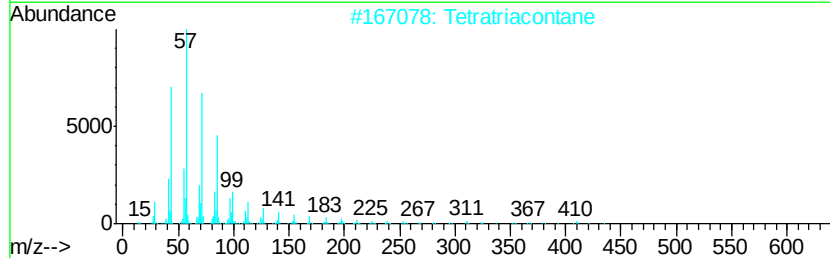
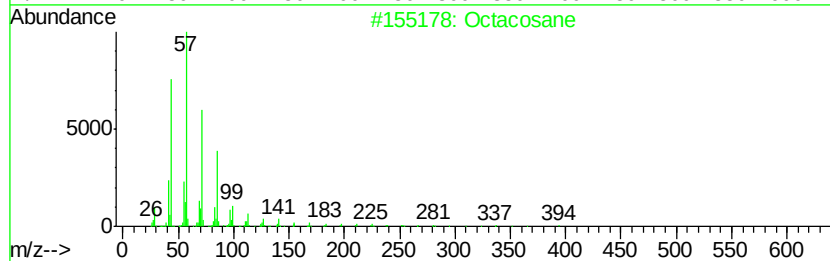
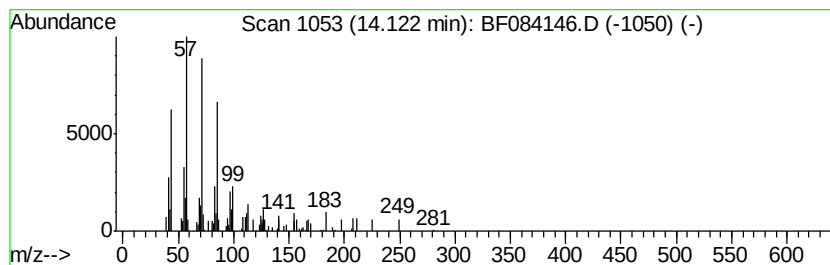
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Octacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.12	2.71 ng	48007	Chrysene-d12	13.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Tetratriacontane	479	C34H70	014167-59-0	90
3		Tetratetracontane	619	C44H90	007098-22-8	90
4		Heneicosane	296	C21H44	000629-94-7	83
5		Tetracosane	338	C24H50	000646-31-1	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122515\
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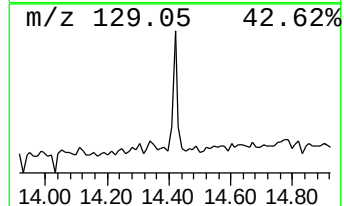
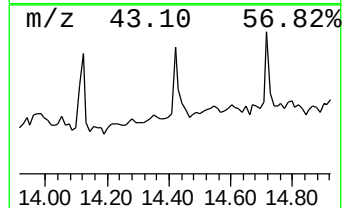
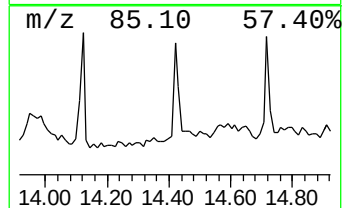
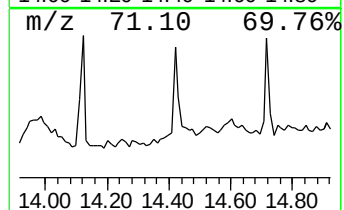
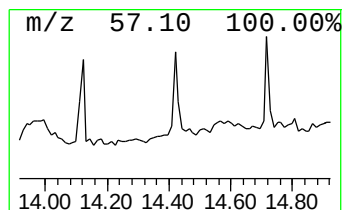
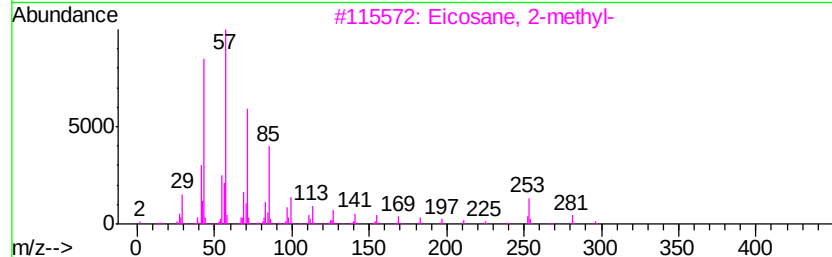
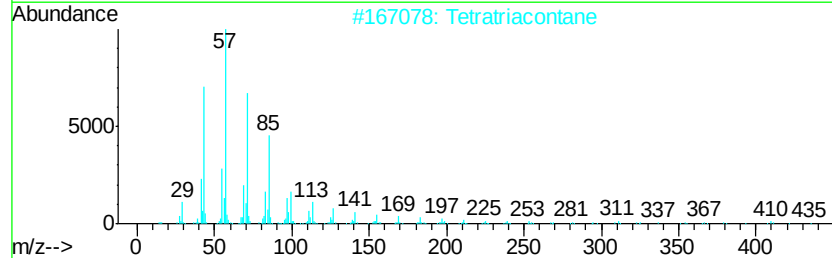
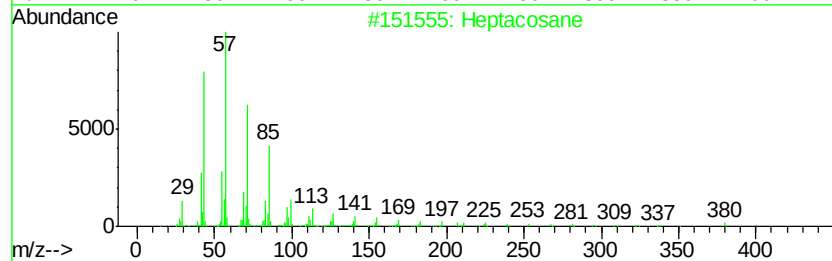
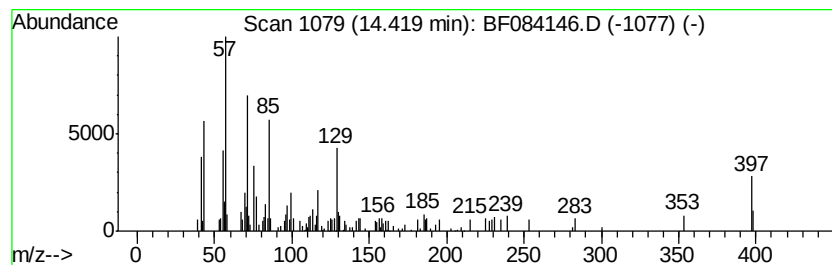
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown14.42 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.42	3.82 ng	67703	Chrysene-d12	13.95	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	46
2	Tetratriacontane	479	C34H70	014167-59-0	46
3	Eicosane, 2-methyl-	296	C21H44	001560-84-5	41
4	Eicosane	282	C20H42	000112-95-8	41
5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	41



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122515\
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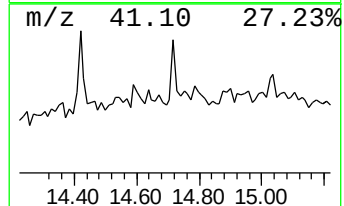
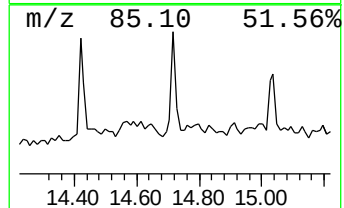
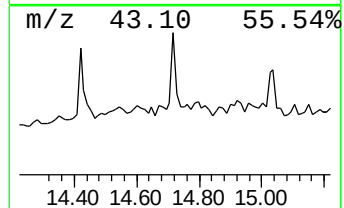
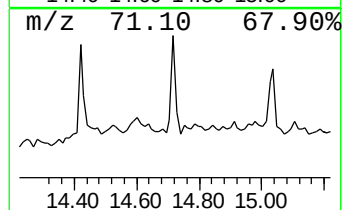
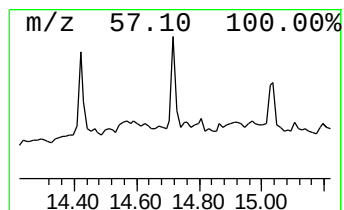
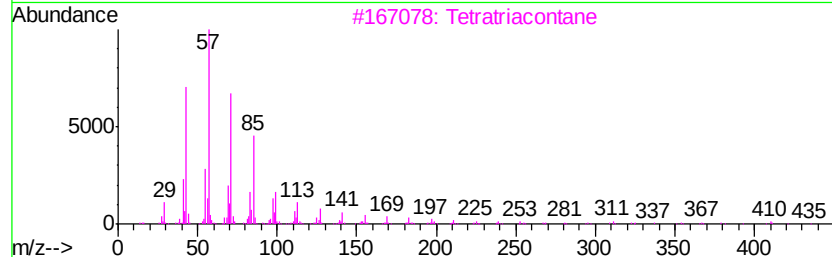
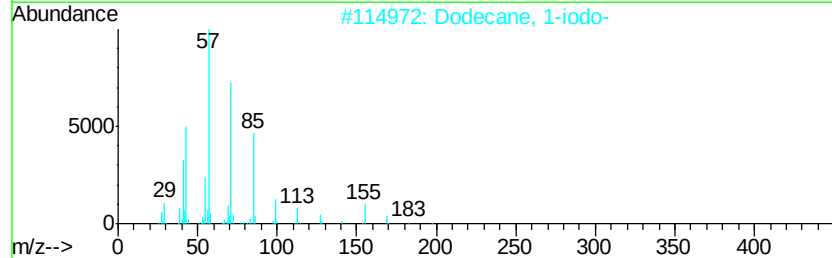
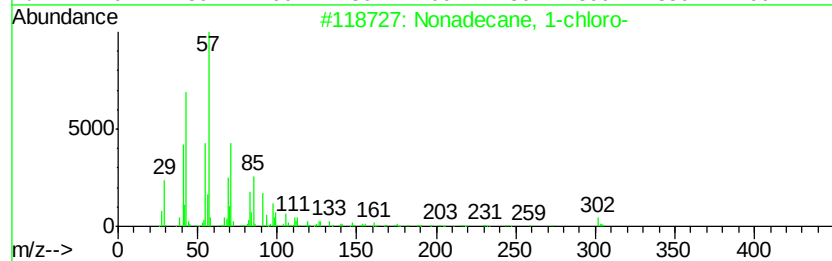
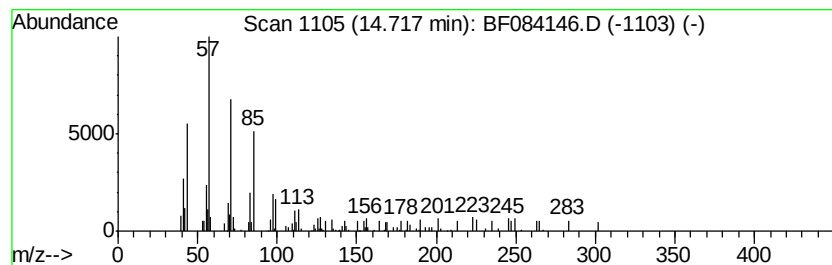
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Nonadecane, 1-chloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.72	3.91 ng	57630	Perylene-d12	15.38	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane, 1-chloro-	302	C19H39Cl	062016-76-6	68
2	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	53
3	Tetratriacontane	479	C34H70	014167-59-0	52
4	Nonadecane	268	C19H40	000629-92-5	49
5	Hexacosane	366	C26H54	000630-01-3	47



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122515\
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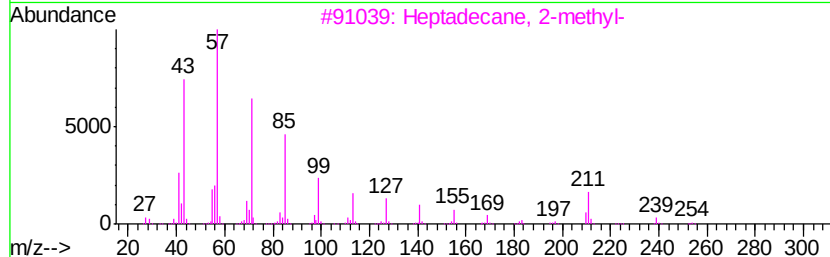
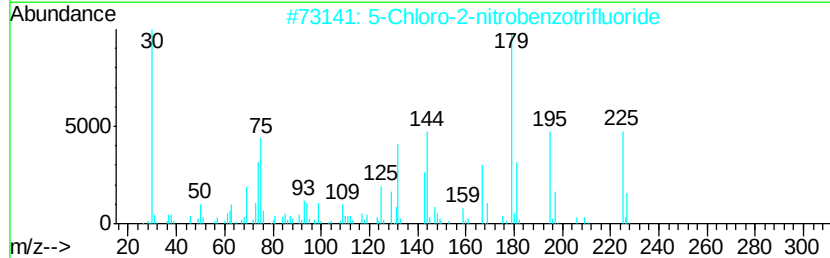
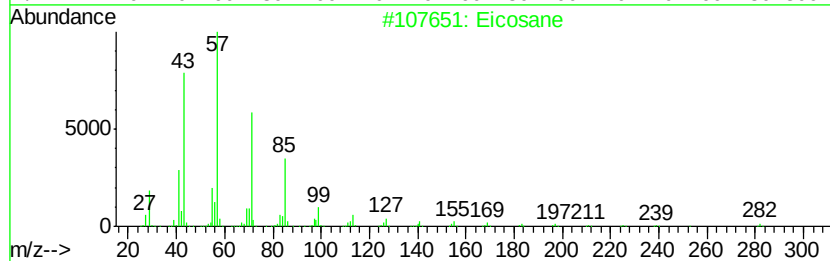
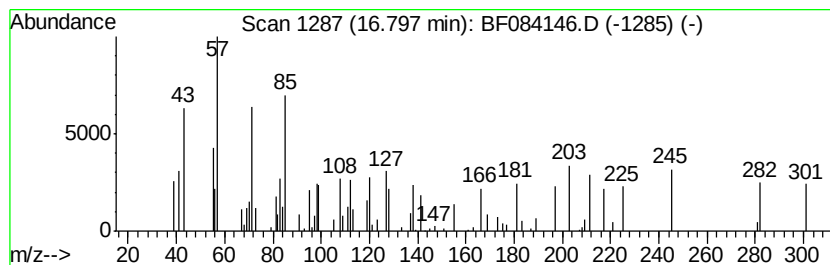
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

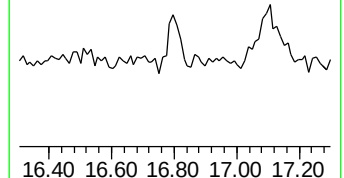
Peak Number 14 unknown16.80 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.80	2.13 ng	31361	Perylene-d12	15.38

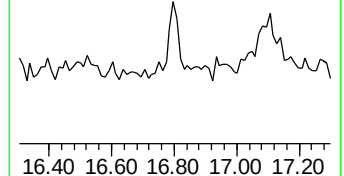
Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	35
2		5-Chloro-2-nitrobenzotrifluoride	225	C7H3ClF3NO2	000118-83-2	11
3		Heptadecane, 2-methyl-	254	C18H38	001560-89-0	10
4		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	10
5		Heneicosane	296	C21H44	000629-94-7	10



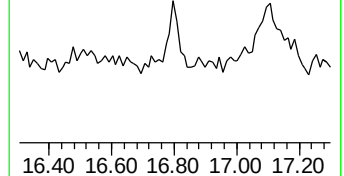
m/z 57.10 100.00%



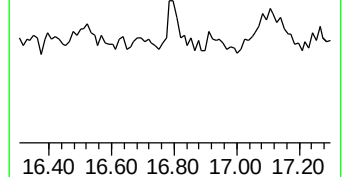
m/z 85.10 69.85%



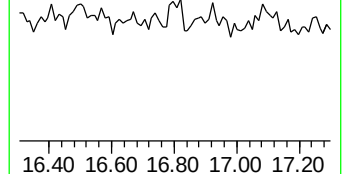
m/z 71.10 63.82%



m/z 43.10 63.17%



m/z 55.10 43.00%



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122515\
Data File : BF084146.D
Acq On : 25 Dec 2015 23:31
Operator : UM/SJ
Sample : G4920-06
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
IM-CLAY-011

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122415.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...	4.98	20.9	ng	346345	1	6.80	330831	20.0
unknown6.58	6.58	104.5	ng	1728490	1	6.80	330831	20.0
Nonadecane	10.27	2.5	ng	51623	3	9.84	421882	20.0
Eicosane	10.74	2.8	ng	49156	4	11.32	354818	20.0
Hexadecanoic acid...	12.73	5.2	ng	92045	5	13.95	354560	20.0
3-Chloropropionic...	12.80	2.4	ng	42548	5	13.95	354560	20.0
Octadecanoic acid...	13.43	3.9	ng	69045	5	13.95	354560	20.0
Heptadecane	13.47	2.2	ng	39558	5	13.95	354560	20.0
17-Pentatriacontene	13.80	8.3	ng	147532	5	13.95	354560	20.0
Octacosane	14.12	2.7	ng	48007	5	13.95	354560	20.0
unknown14.42	14.42	3.8	ng	67703	5	13.95	354560	20.0
Nonadecane, 1-chl...	14.72	3.9	ng	57630	6	15.38	294657	20.0
unknown16.80	16.80	2.1	ng	31361	6	15.38	294657	20.0