

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051215\
 Data File : BF079170.D
 Acq On : 12 May 2015 21:11
 Operator : TP/IZ
 Sample : G2155-08 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-43AS

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-BF043015.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	1.427	18	21	23	rVB	3319975	3795826	100.00%	19.590%
2	1.564	30	33	36	rVB	114915	168841	4.45%	0.871%
3	1.735	46	48	56	rVB2	186530	354026	9.33%	1.827%
4	1.850	56	58	94	rVB	1965275	3584806	94.44%	18.501%
5	5.084	339	341	347	rVB	25002	45060	1.19%	0.233%
6	5.587	382	385	388	rBV	631447	689383	18.16%	3.558%
7	6.845	493	495	498	rBV	799190	723496	19.06%	3.734%
8	6.993	506	508	511	rBV	838245	756123	19.92%	3.902%
9	7.279	531	533	535	rBV	312598	292699	7.71%	1.511%
10	7.462	547	549	552	rBV	421375	562978	14.83%	2.906%
11	7.976	591	594	596	rBV	342905	432193	11.39%	2.231%
12	8.856	669	671	674	rBV	474600	420085	11.07%	2.168%
13	10.205	786	789	791	rBV	959918	928976	24.47%	4.794%
14	10.708	831	833	835	rBV	54619	47193	1.24%	0.244%
15	11.028	859	861	864	rBV2	416373	463489	12.21%	2.392%
16	12.011	944	947	950	rBV	624279	598893	15.78%	3.091%
17	12.879	1020	1023	1024	rBV	444115	518077	13.65%	2.674%
18	12.902	1024	1025	1027	rVB	167555	143039	3.77%	0.738%
19	12.971	1027	1031	1032	rBV	35598	39632	1.04%	0.205%
20	13.588	1083	1085	1087	rBV2	28248	38806	1.02%	0.200%
21	13.634	1087	1089	1091	rBV2	35443	50197	1.32%	0.259%
22	14.388	1152	1155	1156	rBV	160919	219886	5.79%	1.135%
23	14.662	1176	1179	1183	rVB	247221	275084	7.25%	1.420%
24	14.868	1195	1197	1200	rBV	1022367	1086495	28.62%	5.607%
25	15.085	1210	1216	1218	rBV2	31528	71049	1.87%	0.367%
26	15.211	1225	1227	1229	rBV3	35700	48912	1.29%	0.252%
27	15.508	1250	1253	1256	rBV2	45265	95563	2.52%	0.493%
28	15.885	1284	1286	1289	rVV2	18702	40553	1.07%	0.209%
29	16.034	1297	1299	1303	rBV2	101281	184796	4.87%	0.954%
30	16.171	1307	1311	1312	rVV2	461555	769076	20.26%	3.969%
31	16.205	1312	1314	1317	rVB2	151098	222338	5.86%	1.147%
32	16.857	1369	1371	1375	rBV2	44966	81237	2.14%	0.419%
33	17.314	1409	1411	1413	rBV	35382	39772	1.05%	0.205%
34	17.451	1421	1423	1430	rBV	113495	307462	8.10%	1.587%

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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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	17.623	1436	1438	1440	rBV	55761	101342	2.67%	0.523%
36	17.771	1449	1451	1454	rVB	81252	104299	2.75%	0.538%
37	17.828	1454	1456	1460	rVB	98253	117170	3.09%	0.605%
38	17.897	1460	1462	1465	rBV	476552	600832	15.83%	3.101%
39	18.457	1509	1511	1514	rBV	53513	84117	2.22%	0.434%
40	18.514	1514	1516	1518	rBV	53389	89621	2.36%	0.463%
41	20.983	1729	1732	1739	rBV9	47337	182825	4.82%	0.944%

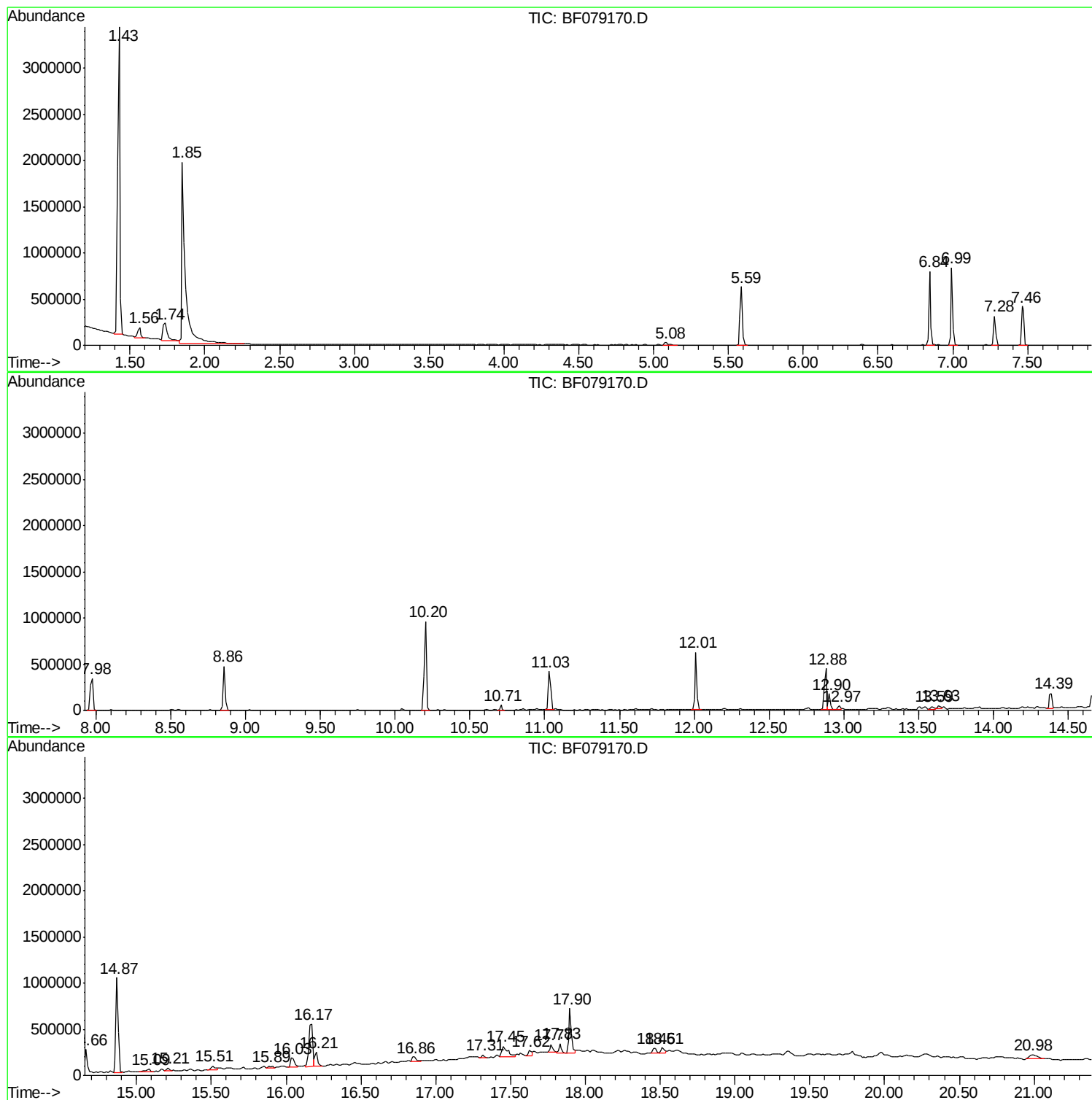
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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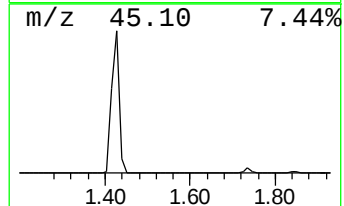
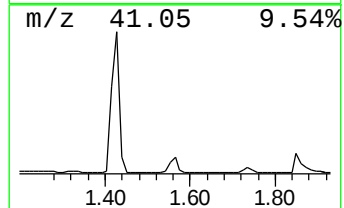
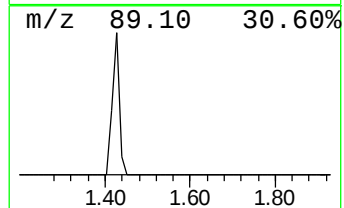
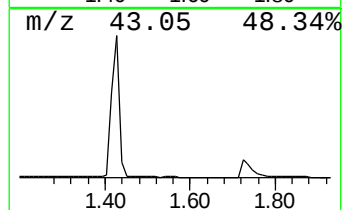
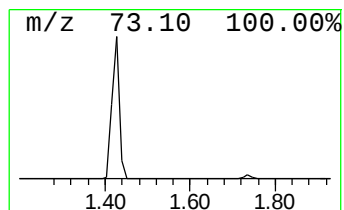
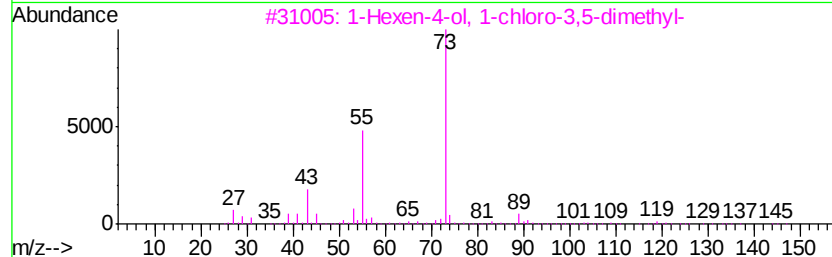
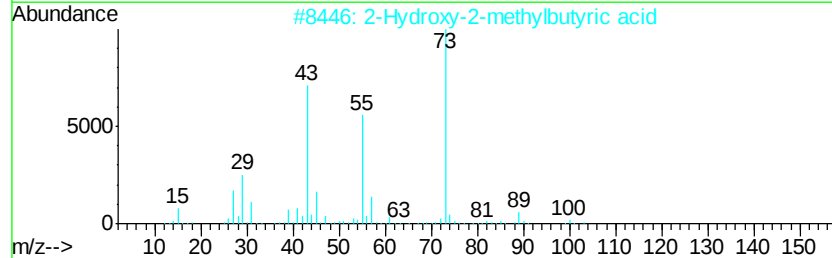
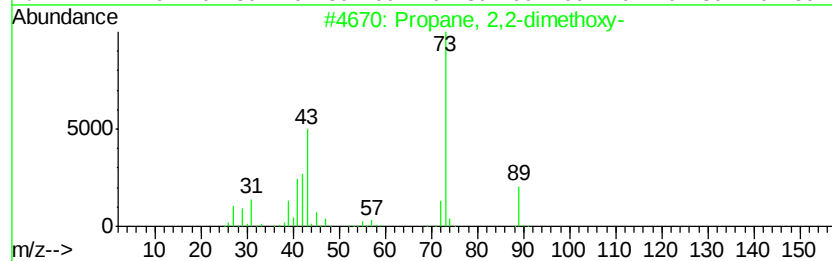
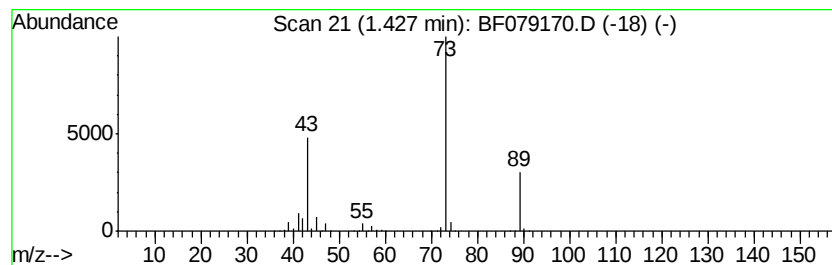
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown1.43 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.43	259.37 ng	3795830	1,4-Dichlorobenzene-d4	7.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	38
2		2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9
3		1-Hexen-4-ol, 1-chloro-3,5-dimet...	162	C8H15ClO	100244-09-5	9
4		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
5		N-Ethylformamide	73	C3H7NO	000627-45-2	5



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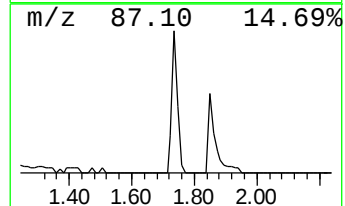
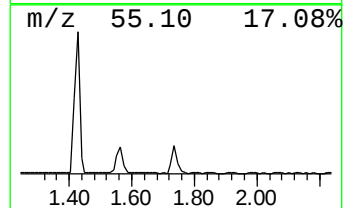
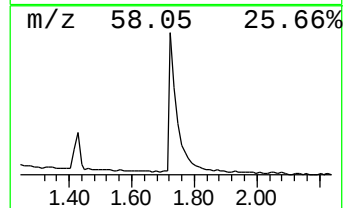
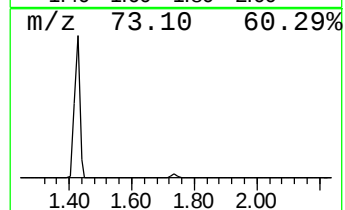
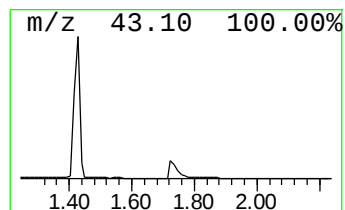
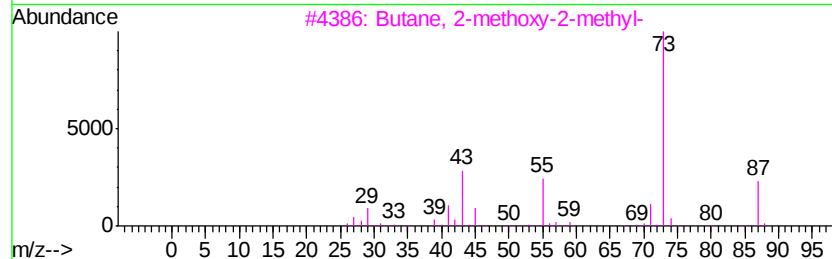
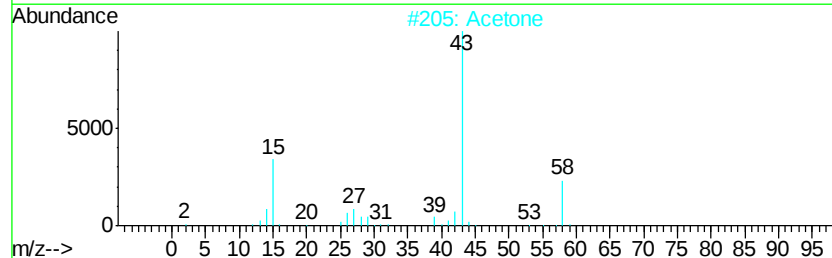
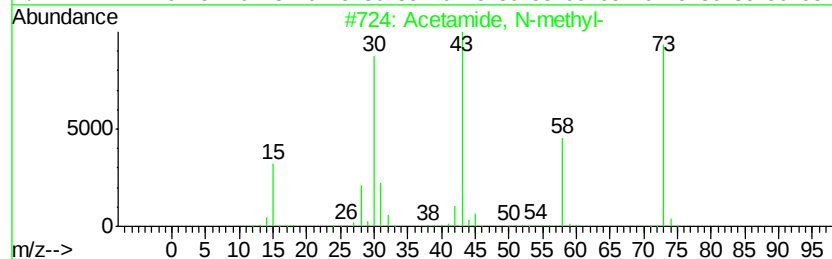
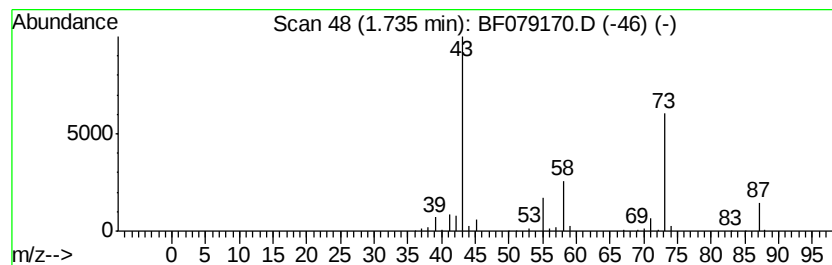
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 unknown1.74 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.74	24.19 ng	354026	1,4-Dichlorobenzene-d4	7.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Acetamide, N-methyl-	73	C3H7NO	000079-16-3	47
2		Acetone	58	C3H6O	000067-64-1	37
3		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	35
4		1-Pentanamine	87	C5H13N	000110-58-7	10
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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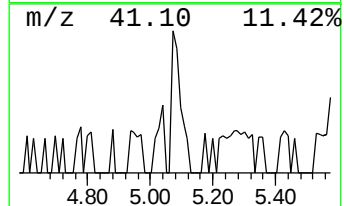
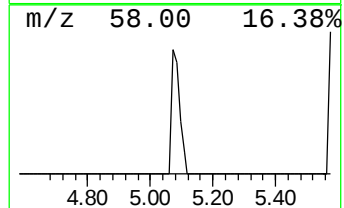
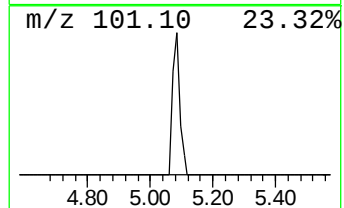
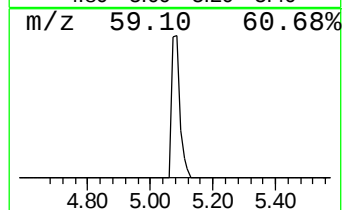
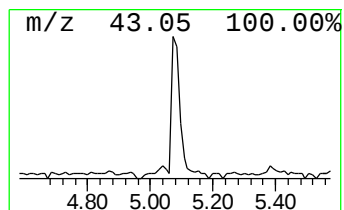
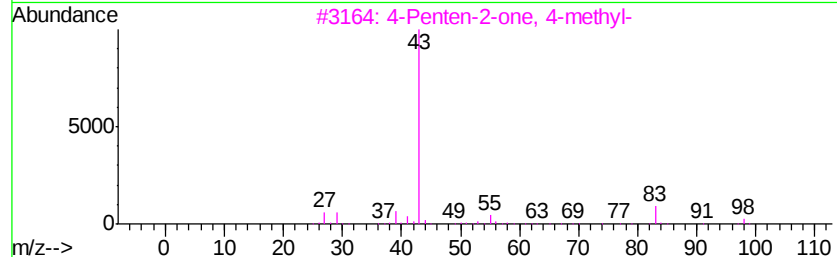
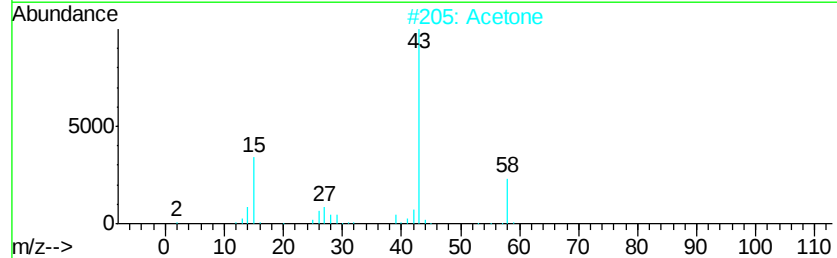
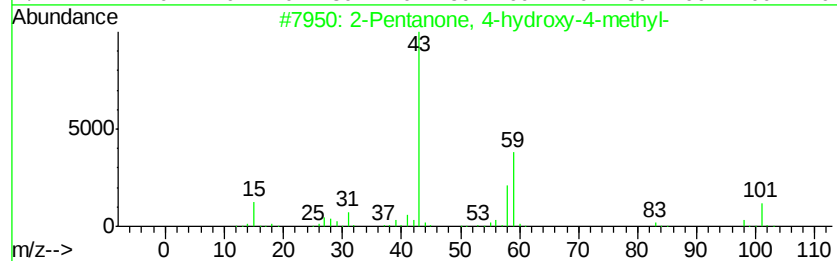
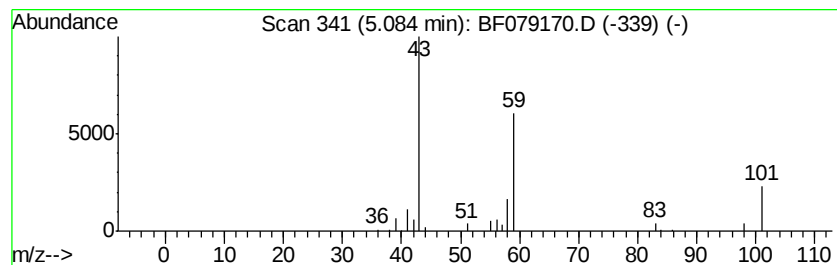
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.08	3.08 ng	45060	1,4-Dichlorobenzene-d4	7.28

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
2		Acetone	58	C3H6O	000067-64-1	9
3		4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		Ethanol, 1-(methylenecyclopropyl)-	98	C6H10O	1000152-74-6	9



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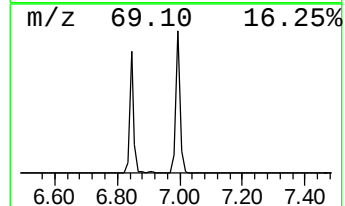
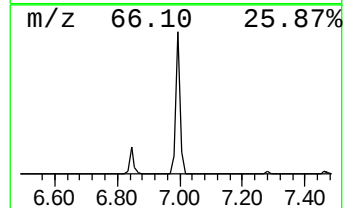
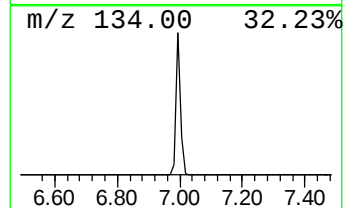
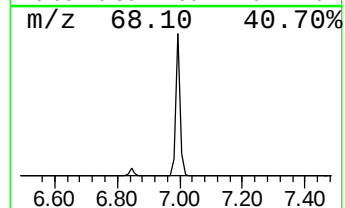
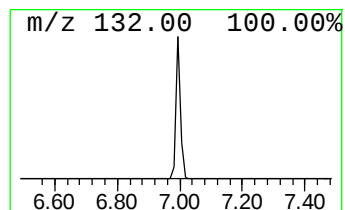
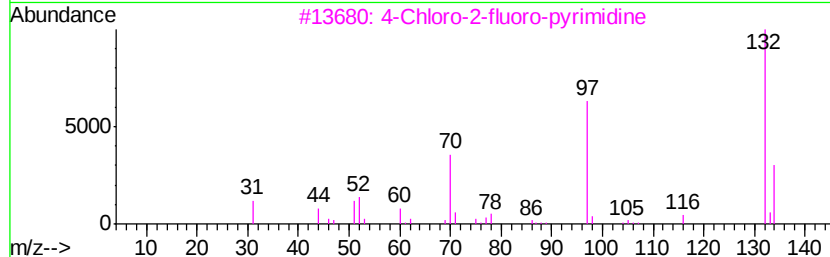
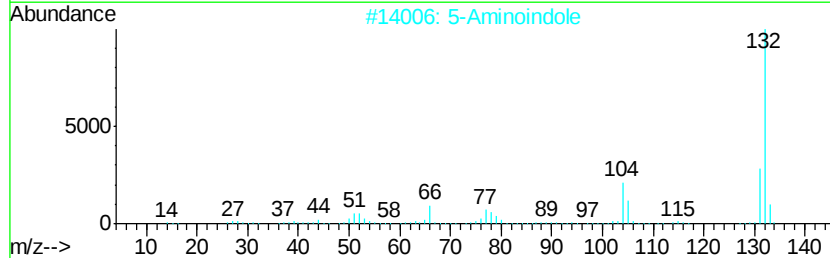
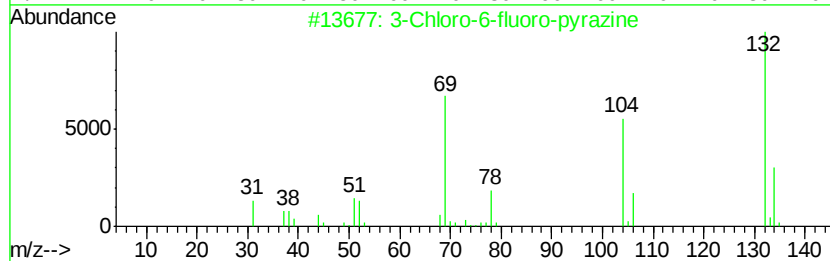
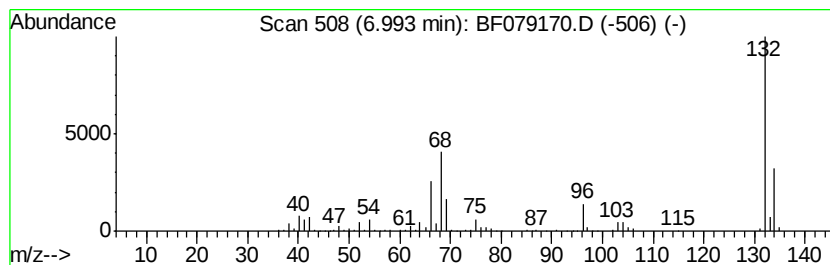
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown6.99 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.99	51.67 ng	756123	1,4-Dichlorobenzene-d4	7.28

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	5		Aminoindole	132	C8H8N2	005192-03-0	12
3	4		Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
4	(E)-3		Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5	5		Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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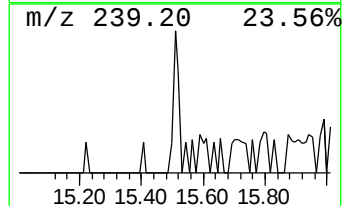
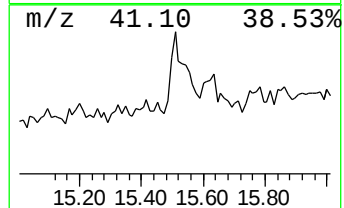
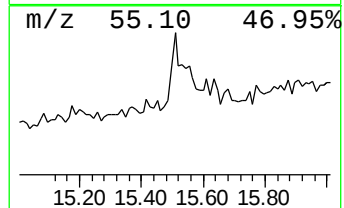
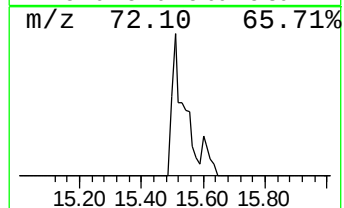
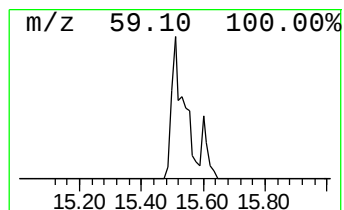
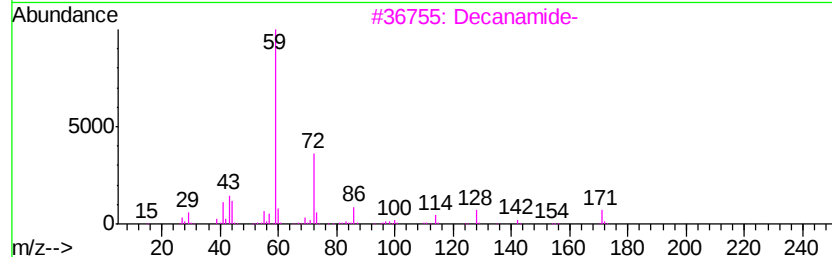
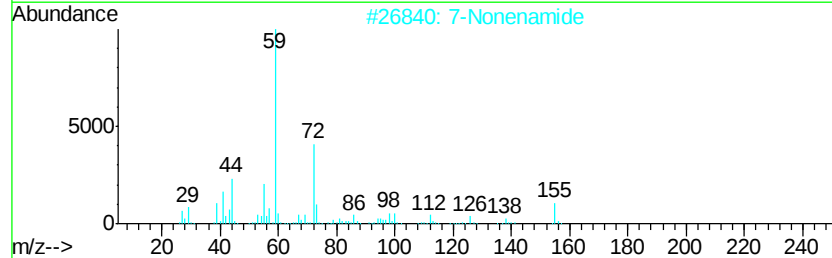
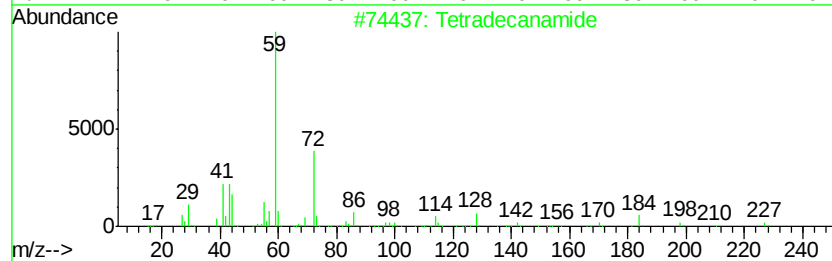
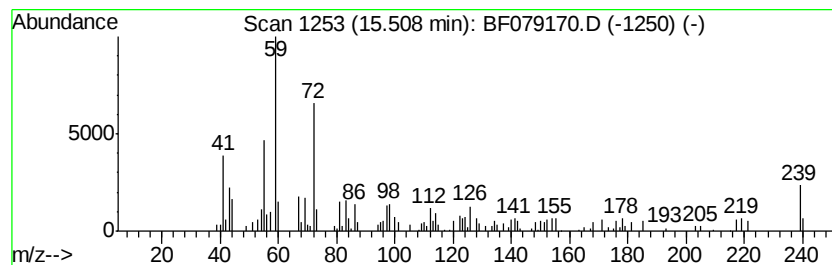
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 8 Tetradecanamide Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.51	2.49 ng	95563	Chrysene-d12	16.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecanamide	227	C14H29NO	000638-58-4	53
2		7-Nonenamide	155	C9H17NO	090949-53-4	53
3		Decanamide-	171	C10H21NO	002319-29-1	53
4		Hexadecanamide	255	C16H33NO	000629-54-9	50
5		Pentanamide	101	C5H11NO	000626-97-1	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051215\
Data File : BF079170.D
Acq On : 12 May 2015 21:11
Operator : TP/IZ
Sample : G2155-08 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-43AS

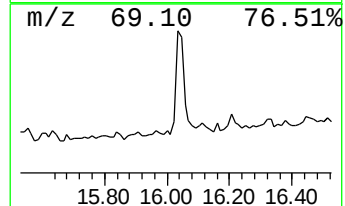
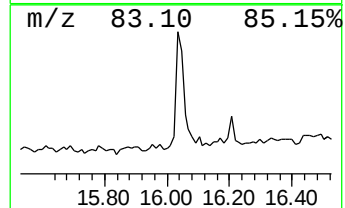
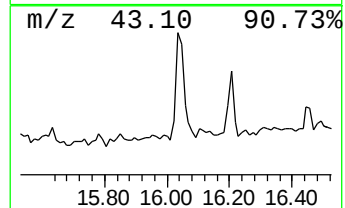
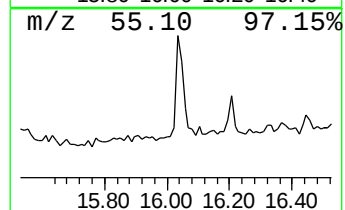
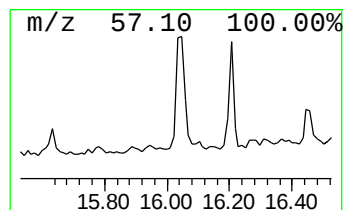
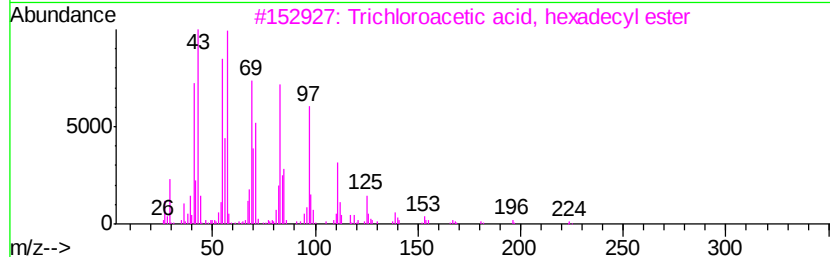
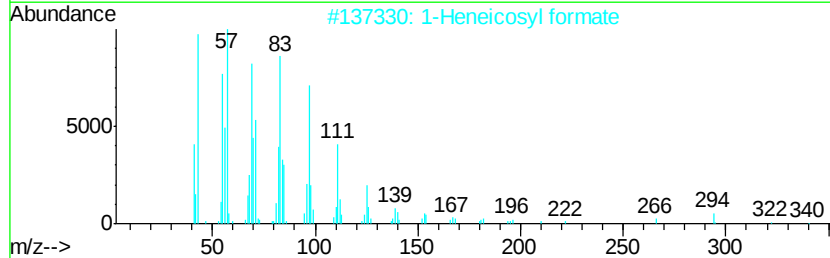
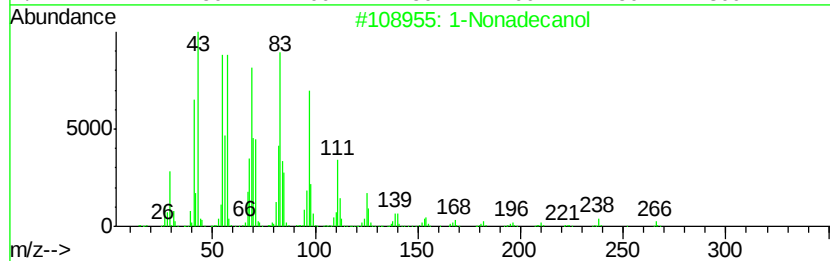
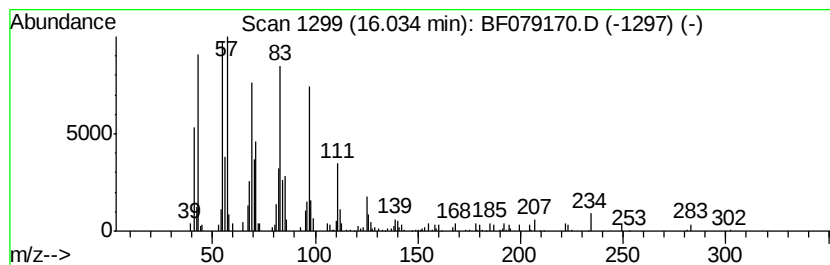
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Nonadecanol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.03	4.81 ng	184796	Chrysene-d12	16.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecanol		284	C19H40O	001454-84-8	95
2	1-Heneicosyl formate		340	C22H44O2	077899-03-7	94
3	Trichloroacetic acid, hexadecyl ...		386	C18H33Cl3O2	074339-54-1	94
4	Pentafluoropropionic acid, hepta...		402	C20H35F5O2	1000283-04-2	93
5	Heptafluorobutanoic acid, heptad...		452	C21H35F7O2	1000282-97-3	93



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
Data File : BF079170.D
Acq On : 12 May 2015 21:11
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Misc :
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Instrument :
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ClientSampleId :
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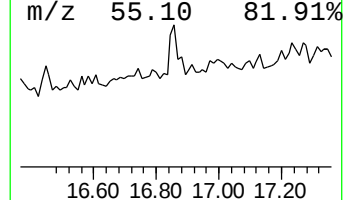
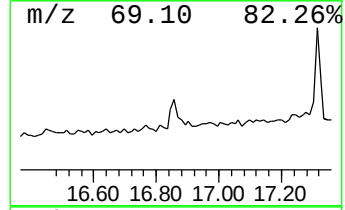
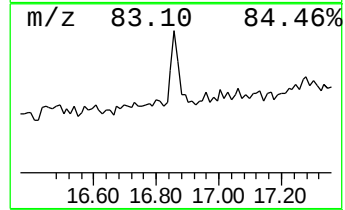
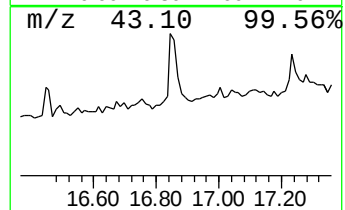
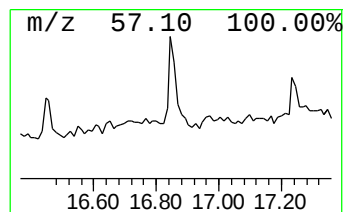
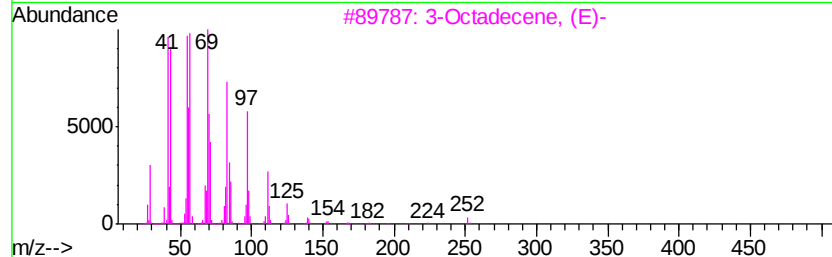
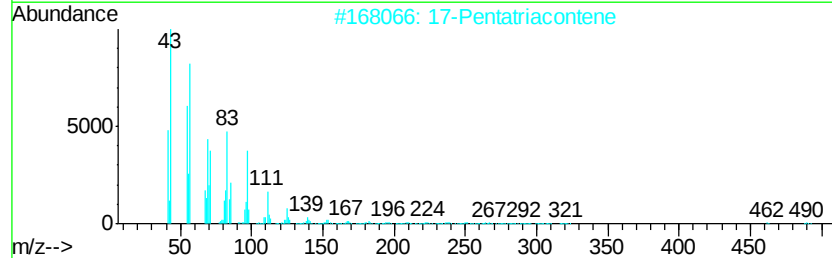
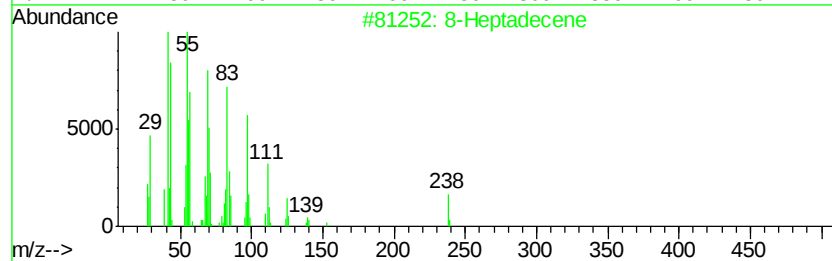
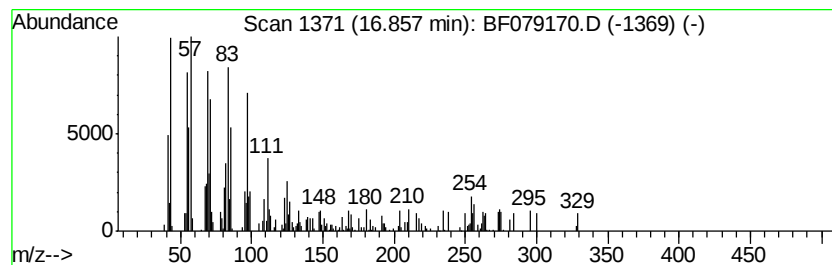
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 8-Heptadecene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.86	2.11 ng	81237	Chrysene-d12	16.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	8-Heptadecene		238	C17H34	054290-12-9	86
2	17-Pentatriacontene		491	C35H70	006971-40-0	74
3	3-Octadecene, (E)-		252	C18H36	007206-19-1	70
4	Cyclopentadecane		210	C15H30	000295-48-7	70
5	Acetic acid, chloro-, octadecyl ...		346	C20H39ClO2	005348-82-3	62



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
Data File : BF079170.D
Acq On : 12 May 2015 21:11
Operator : TP/IZ
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Misc :
ALS Vial : 15 Sample Multiplier: 1

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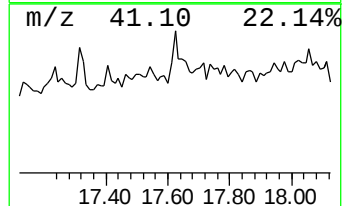
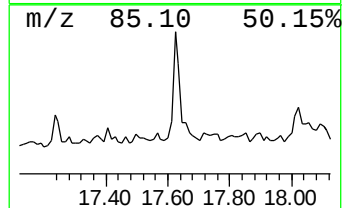
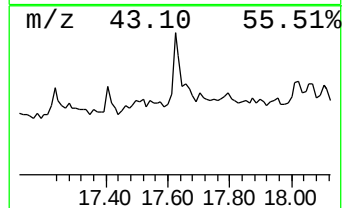
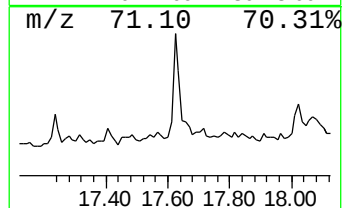
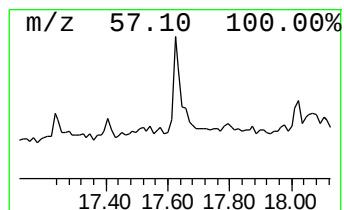
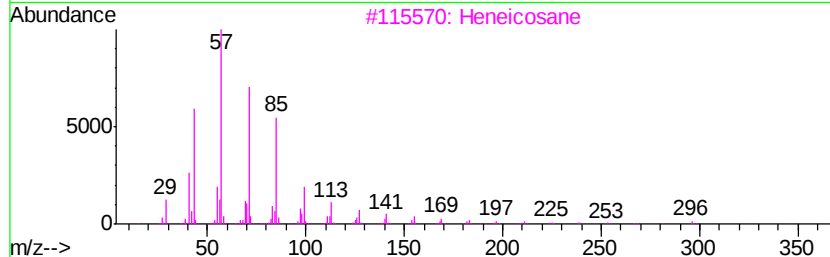
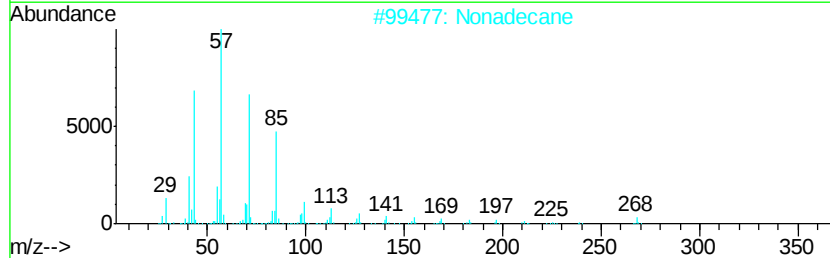
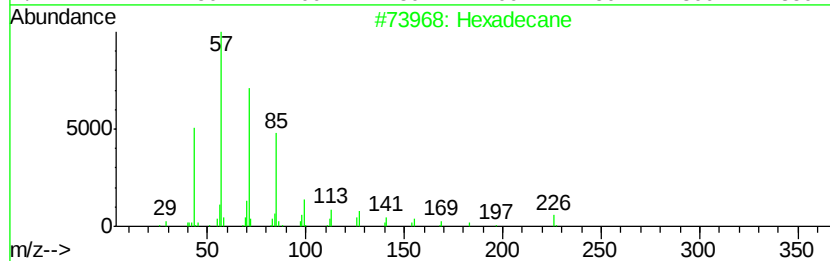
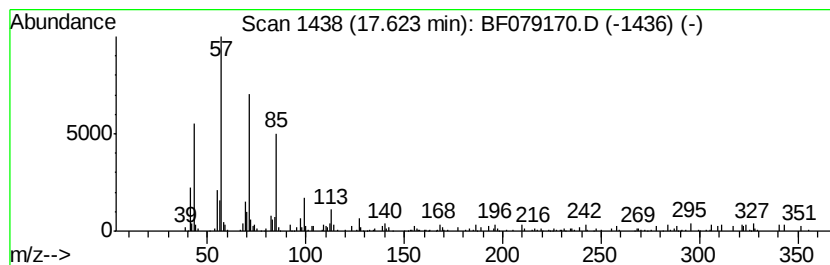
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.62	3.37 ng	101342	Perylene-d12	17.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	76
2		Nonadecane	268	C19H40	000629-92-5	76
3		Heneicosane	296	C21H44	000629-94-7	68
4		Octadecane	254	C18H38	000593-45-3	64
5		Heptadecane	240	C17H36	000629-78-7	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
Data File : BF079170.D
Acq On : 12 May 2015 21:11
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Misc :
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ClientSampleId :
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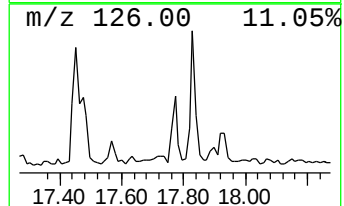
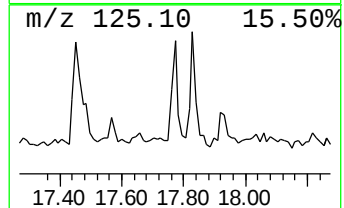
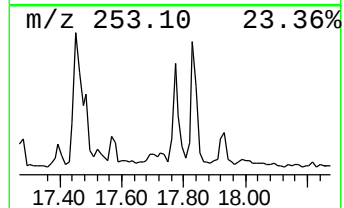
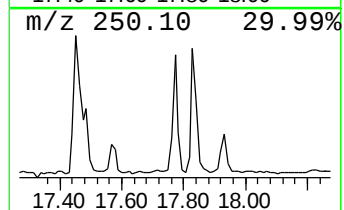
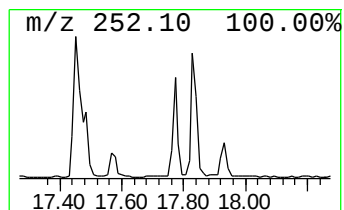
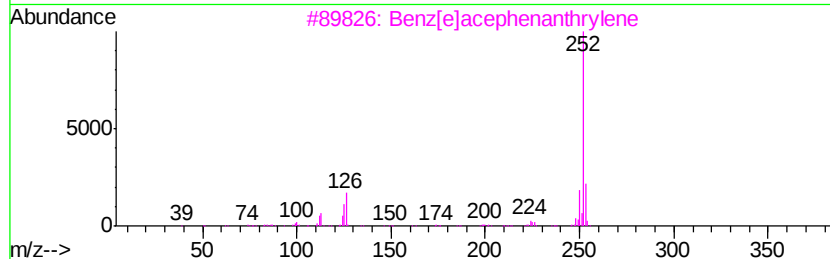
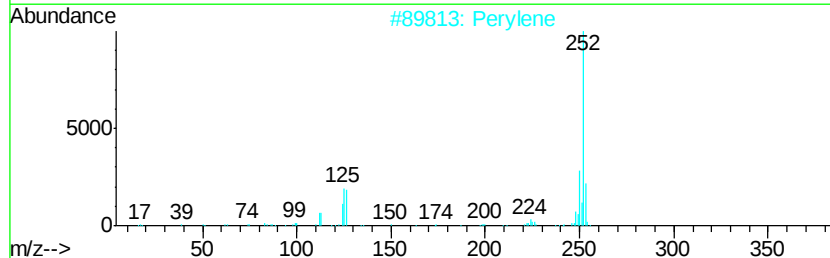
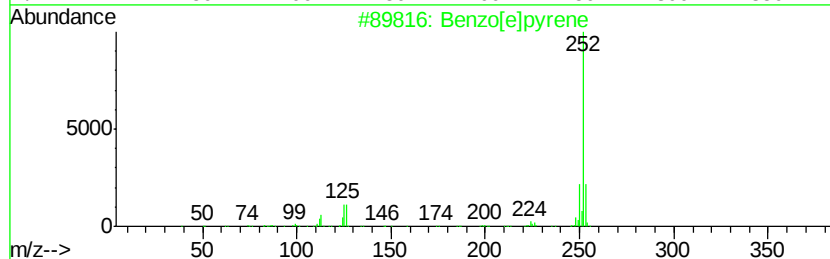
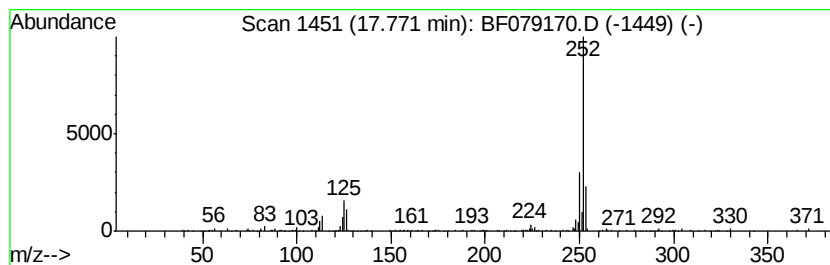
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzo[e]pyrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.77	3.47 ng	104299	Perylene-d12	17.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	96
2		Perylene	252	C20H12	000198-55-0	90
3		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	89
4		Benzo[k]fluoranthene	252	C20H12	000207-08-9	89
5		Benzo[a]pyrene	252	C20H12	000050-32-8	81



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051215\
Data File : BF079170.D
Acq On : 12 May 2015 21:11
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Sample : G2155-08 2X
Misc :
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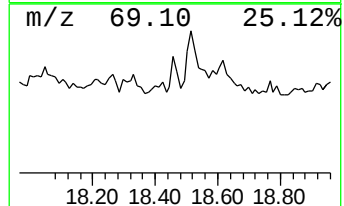
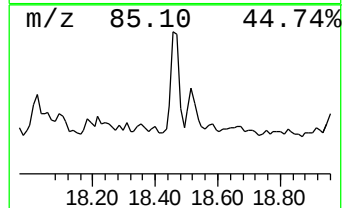
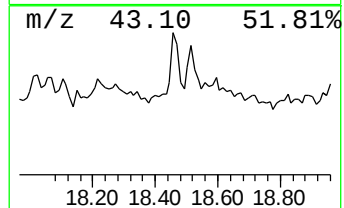
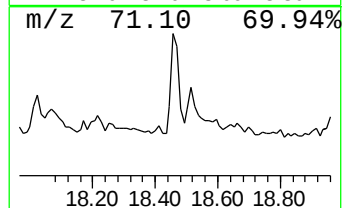
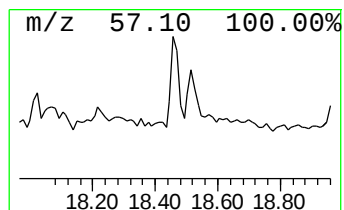
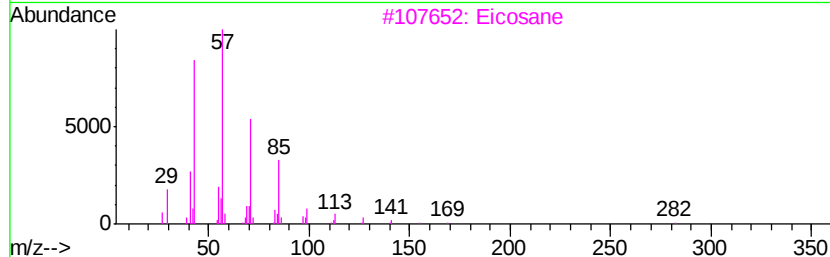
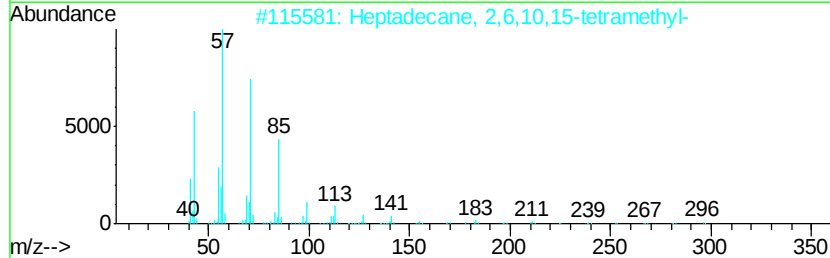
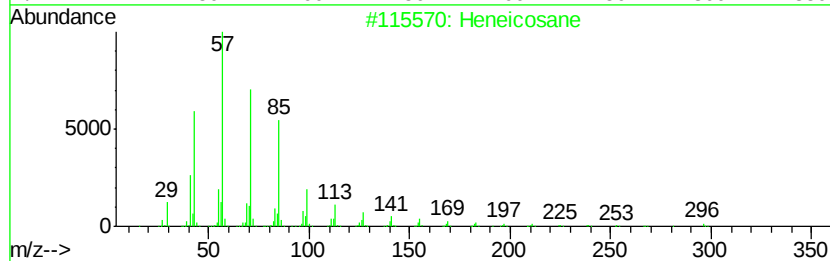
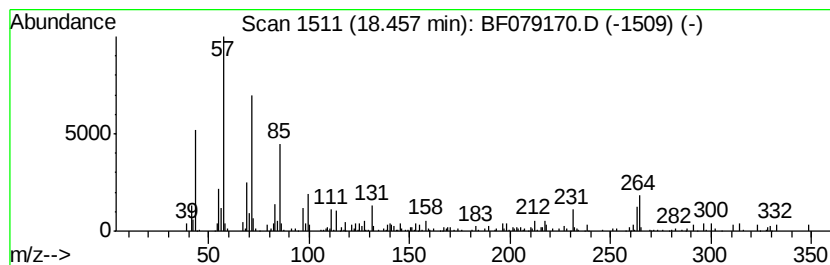
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heneicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.46	2.80 ng	84117	Perylene-d12	17.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	78
2		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	58
3		Eicosane	282	C20H42	000112-95-8	55
4		Octacosane	394	C28H58	000630-02-4	53
5		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	52



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051215\
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Acq On : 12 May 2015 21:11
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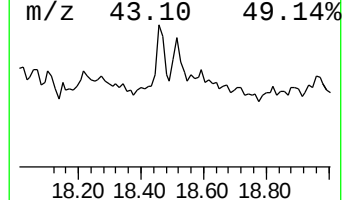
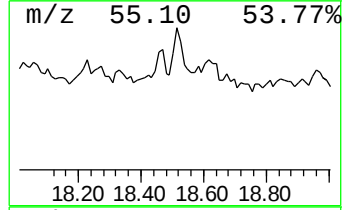
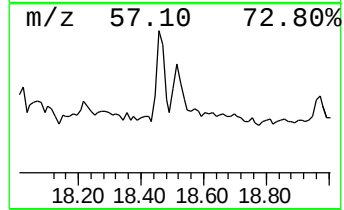
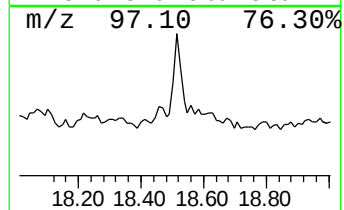
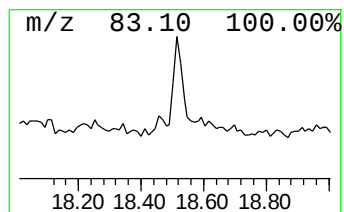
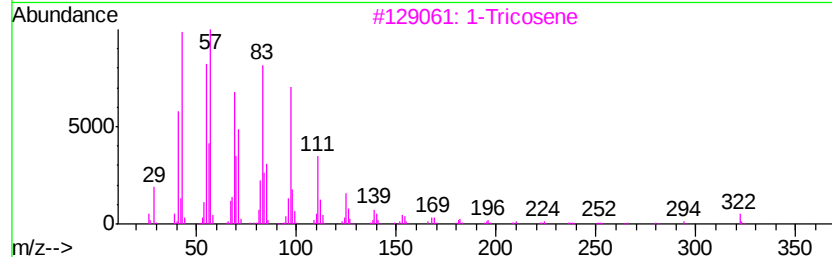
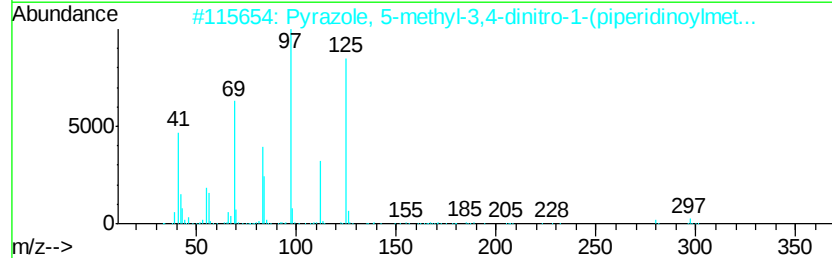
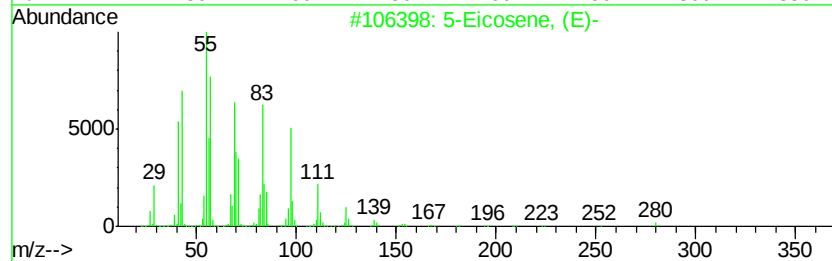
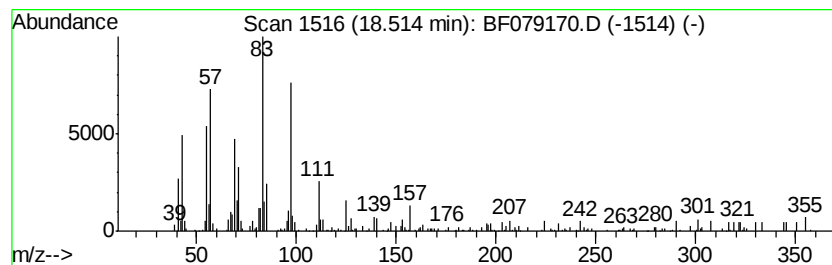
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 5-Eicosene, (E)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.51	2.98 ng	89621	Perylene-d12	17.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Eicosene, (E)-	280	C20H40	074685-30-6	80
2		Pyrazole, 5-methyl-3,4-dinitro-1...	297	C11H15N5O5	345950-83-6	38
3		1-Tricosene	322	C23H46	018835-32-0	30
4		1-Azabicyclo[2.2.2]octan-3-one	125	C7H11NO	003731-38-2	30
5		3-Decen-5-one	154	C10H18O	032064-73-6	30



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
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Acq On : 12 May 2015 21:11
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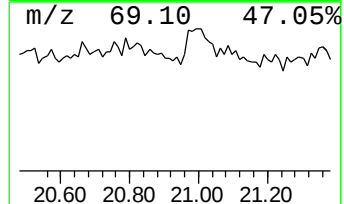
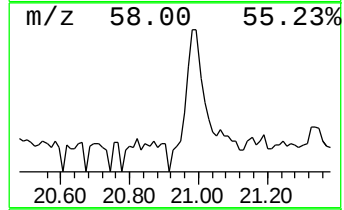
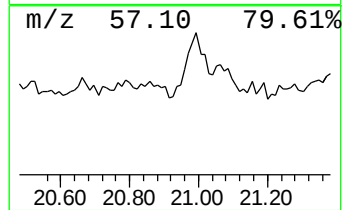
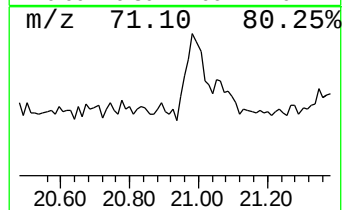
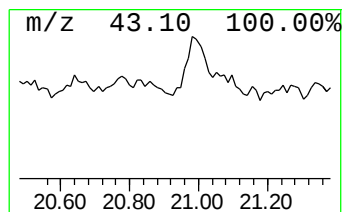
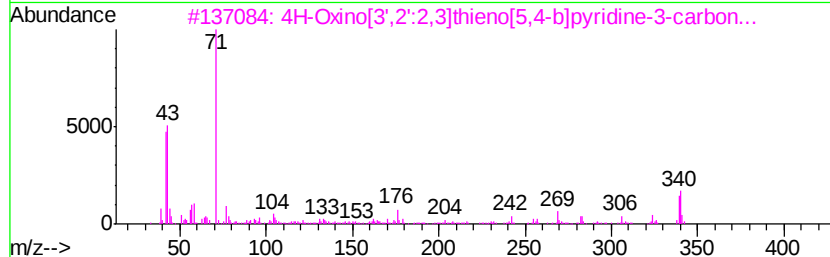
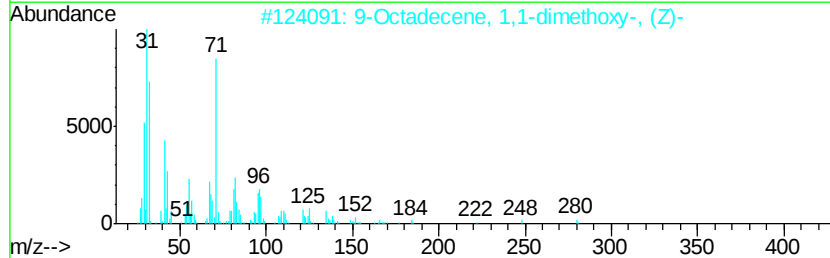
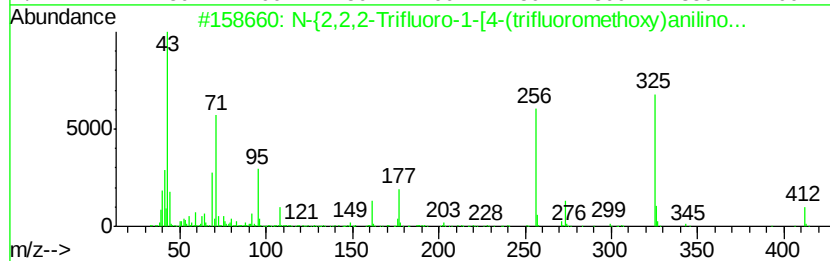
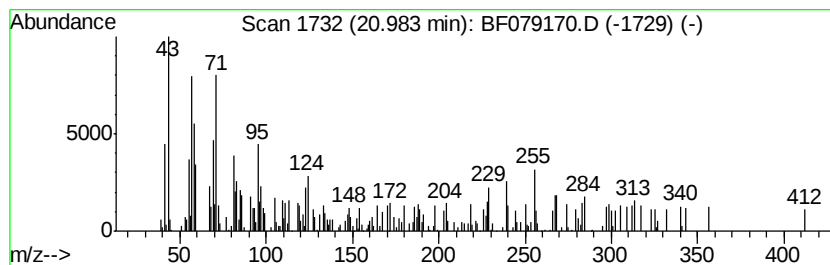
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown20.98 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.98	6.09 ng	182825	Perylene-d12	17.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	N-{2,2,2-Trifluoro-1-[4-(trifluo...	412	C14H13F9N2O2	1000225-27-5	37
2		9-Octadecene, 1,1-dimethoxy-, (Z)-	312	C20H40O2	015677-71-1	35
3		4H-Oxino[3',2':2,3]thieno[5,4-b]...	340	C18H20N4OS	1000270-48-0	35
4		Octadecane, 1,1'-[(1-methyl-1,2-...	581	C39H80O2	035545-51-8	27
5		(1H)Benzimidazole, 5-fluoro-2-(4...	340	C19H21FN4O	174468-77-0	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051215\
 Data File : BF079170.D
 Acq On : 12 May 2015 21:11
 Operator : TP/IZ
 Sample : G2155-08 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-43AS

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown1.43	1.43	259.4	ng	3795830	1	7.28	292699	20.0
unknown1.74	1.74	24.2	ng	354026	1	7.28	292699	20.0
2-Pentanone, 4-hy...	5.08	3.1	ng	45060	1	7.28	292699	20.0
unknown6.99	6.99	51.7	ng	756123	1	7.28	292699	20.0
Tetradecanamide	15.51	2.5	ng	95563	5	16.17	769076	20.0
1-Nonadecanol	16.03	4.8	ng	184796	5	16.17	769076	20.0
8-Heptadecene	16.86	2.1	ng	81237	5	16.17	769076	20.0
Hexadecane	17.62	3.4	ng	101342	6	17.90	600832	20.0
Benzo[e]pyrene	17.77	3.5	ng	104299	6	17.90	600832	20.0
Heneicosane	18.46	2.8	ng	84117	6	17.90	600832	20.0
5-Eicosene, (E)-	18.51	3.0	ng	89621	6	17.90	600832	20.0
unknown20.98	20.98	6.1	ng	182825	6	17.90	600832	20.0