

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
 Data File : BF077875.D
 Acq On : 20 Mar 2015 20:54
 Operator : TP/IZ
 Sample : G1575-08
 Misc : W
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RW-9V

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-BF031115.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	1.047	13	14	17	rVB	37157	31696	1.53%	0.219%
2	1.310	35	37	39	rVV	43121	46685	2.25%	0.322%
3	1.355	39	41	42	rVV	13323	22239	1.07%	0.153%
4	1.413	42	46	48	rVB	190358	339381	16.37%	2.341%
5	1.573	58	60	63	rVB	288218	318113	15.34%	2.195%
6	1.767	75	77	85	rBV	119650	163847	7.90%	1.130%
7	1.893	85	88	92	rVB2	89905	121630	5.87%	0.839%
8	2.373	126	130	133	rVB	15349	22230	1.07%	0.153%
9	2.910	174	177	181	rBV	55129	100165	4.83%	0.691%
10	3.070	181	191	195	rVV3	13205	40625	1.96%	0.280%
11	4.201	286	290	293	rBV	336779	554852	26.76%	3.828%
12	4.407	305	308	312	rVB	53505	76827	3.71%	0.530%
13	4.876	347	349	353	rBV	88229	123169	5.94%	0.850%
14	5.013	353	361	366	rVV	26806	108159	5.22%	0.746%
15	5.150	366	373	382	rVB	17520	57996	2.80%	0.400%
16	5.413	394	396	399	rBV	608919	648213	31.26%	4.472%
17	6.567	495	497	500	rBV	58758	75044	3.62%	0.518%
18	6.705	506	509	511	rBV	421480	427749	20.63%	2.951%
19	6.842	518	521	523	rBV	1127139	1176420	56.74%	8.116%
20	7.127	543	546	548	rBV	260540	251950	12.15%	1.738%
21	7.276	557	559	560	rBV	28552	26709	1.29%	0.184%
22	7.310	560	562	565	rVB	1174535	1065420	51.38%	7.351%
23	7.607	585	588	589	rBV	19682	23429	1.13%	0.162%
24	7.825	604	607	609	rBV	783453	852510	41.11%	5.882%
25	7.893	611	613	615	rVB	82078	73311	3.54%	0.506%
26	8.430	653	660	664	rBV2	116333	316738	15.28%	2.185%
27	8.705	681	684	686	rBV	323136	347308	16.75%	2.396%
28	8.888	697	700	702	rBV	192273	209392	10.10%	1.445%
29	9.059	713	715	721	rBV	58968	76969	3.71%	0.531%
30	9.185	724	726	732	rBV2	27613	35368	1.71%	0.244%
31	9.882	785	787	790	rBV	18166	21294	1.03%	0.147%
32	10.053	799	802	804	rBV	1801621	1817995	87.68%	12.543%
33	10.865	871	873	876	rVB	299701	403529	19.46%	2.784%
34	11.185	896	901	906	rBV3	42150	51489	2.48%	0.355%

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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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35	11.848	956	959	962	rBV	1052252	1030247	49.69%	7.108%
36	12.705	1032	1034	1037	rBV	445255	421712	20.34%	2.909%
37	14.705	1206	1209	1211	rBV	1940748	2073505	100.00%	14.306%
38	15.985	1319	1321	1324	rBV	444493	472685	22.80%	3.261%
39	17.700	1468	1471	1474	rBV	390033	467843	22.56%	3.228%

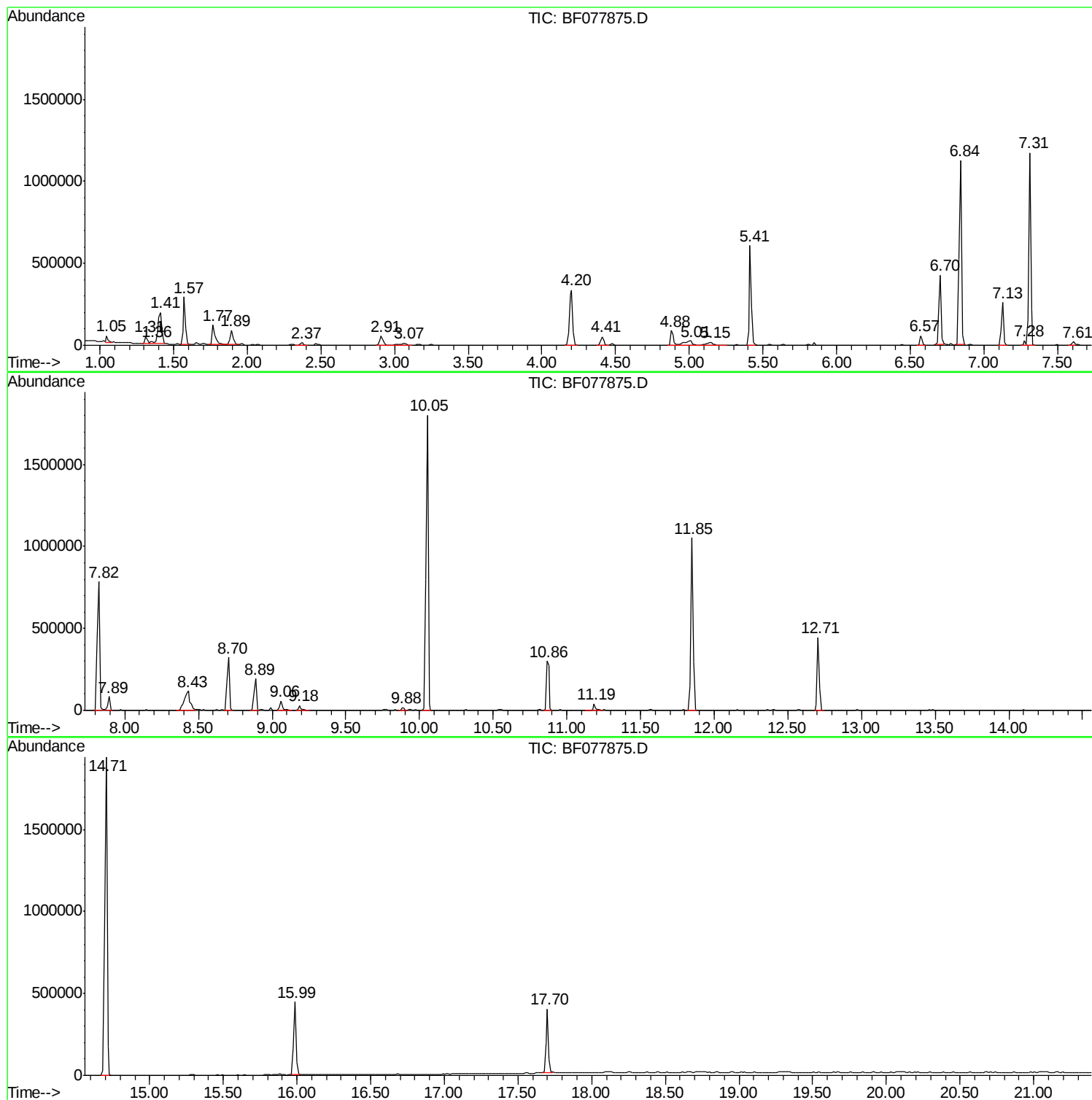
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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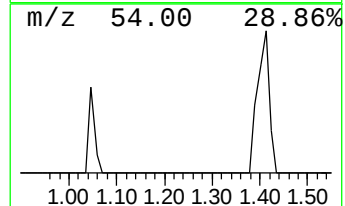
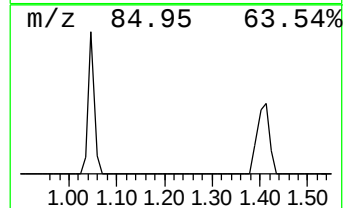
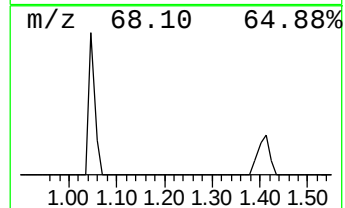
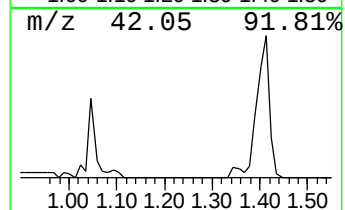
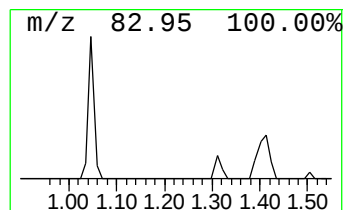
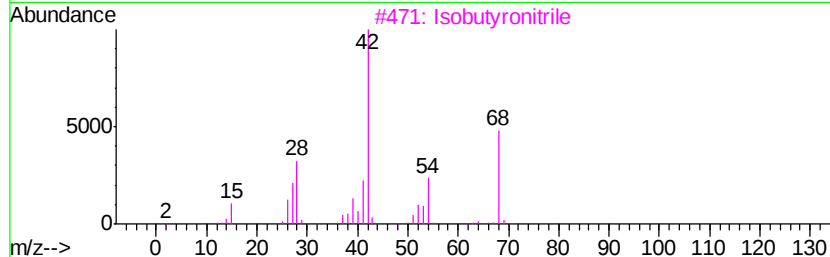
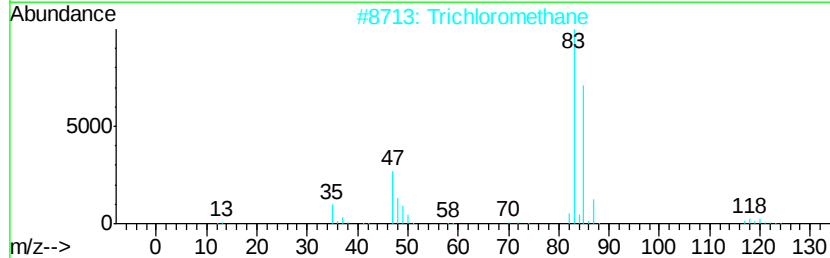
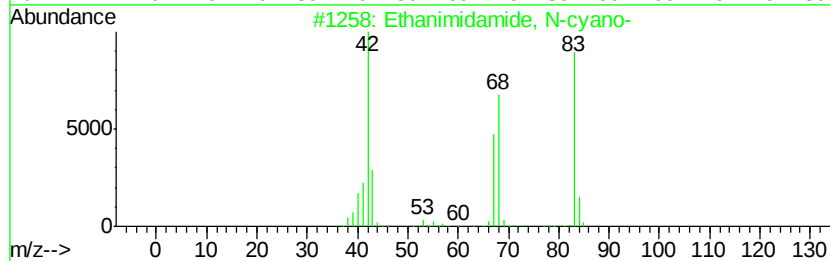
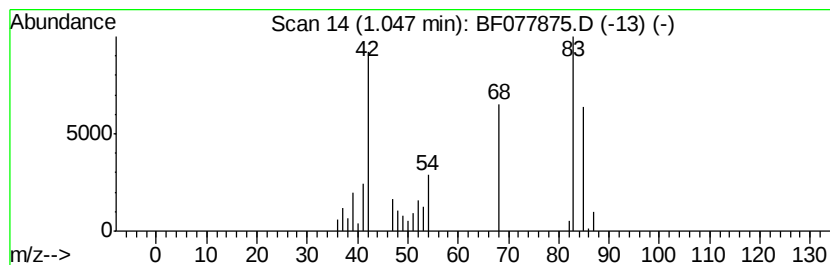
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Ethanimidamide, N-cyano- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.05	2.52 ng	31696	1,4-Dichlorobenzene-d4	7.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanimidamide, N-cyano-	83	C3H5N3	056563-07-6	50
2			Trichloromethane	118	CHCl3	000067-66-3	38
3			Isobutyronitrile	69	C4H7N	000078-82-0	27
4			Methane, oxybis[dichloro-	182	C2H2Cl4O	020524-86-1	25
5			3-Aminopyrazole	83	C3H5N3	001820-80-0	14



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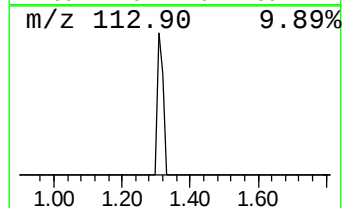
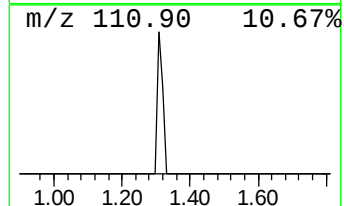
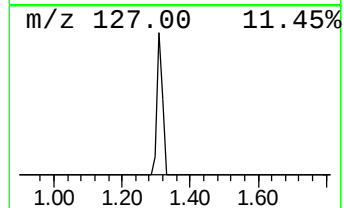
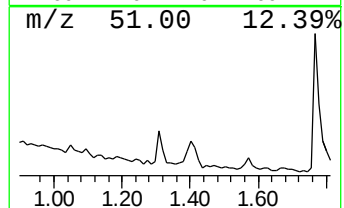
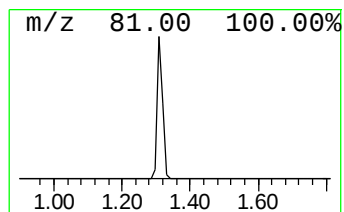
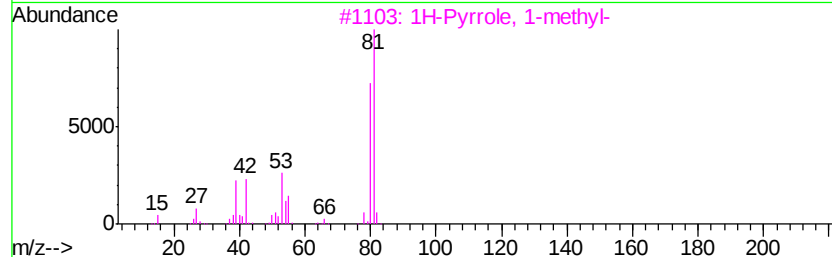
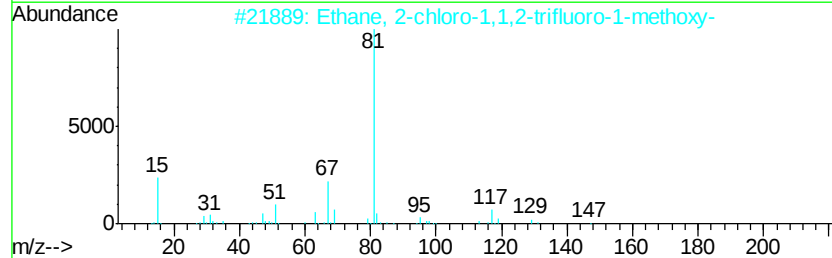
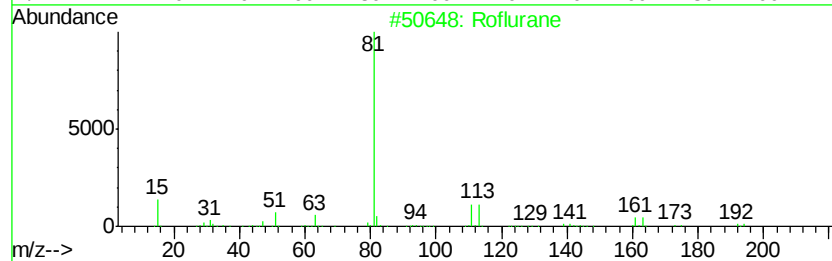
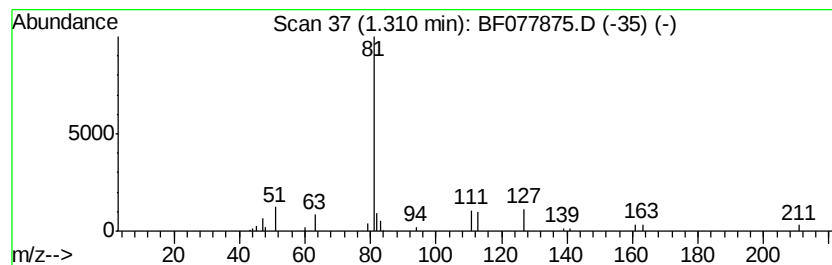
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Roflurane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.31	3.71 ng	46685	1,4-Dichlorobenzene-d4	7.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Roflurane	192	C3H4BrF3O	000679-90-3	72
2		Ethane, 2-chloro-1,1,2-trifluoro...	148	C3H4ClF3O	000425-87-6	9
3		1H-Pyrrole, 1-methyl-	81	C5H7N	000096-54-8	5
4		Cyclohexane, 1,4-bis[(tert-butyl...	344	C18H40O2Si2	1000221-82-2	4
5		Trichloroacetic acid, 4-tetradec...	358	C16H29Cl3O2	1000282-06-8	4



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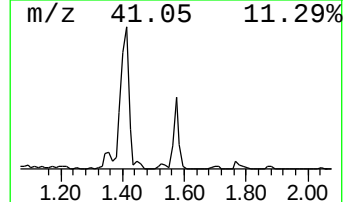
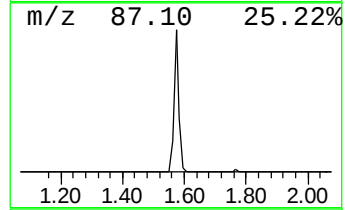
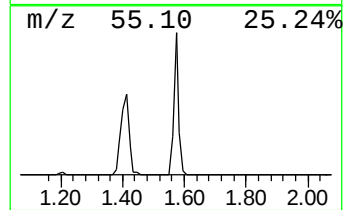
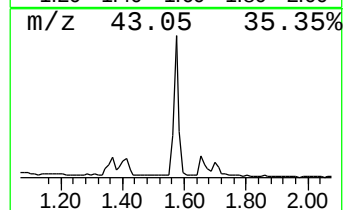
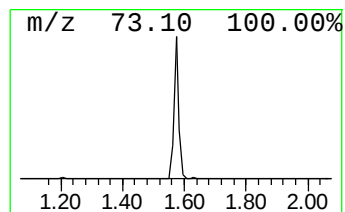
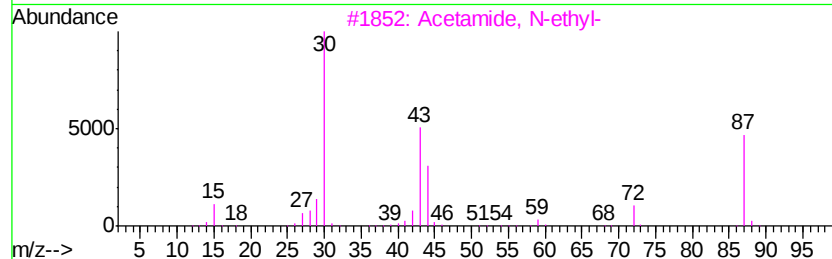
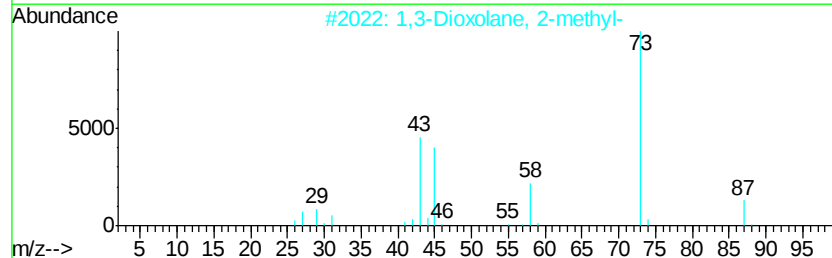
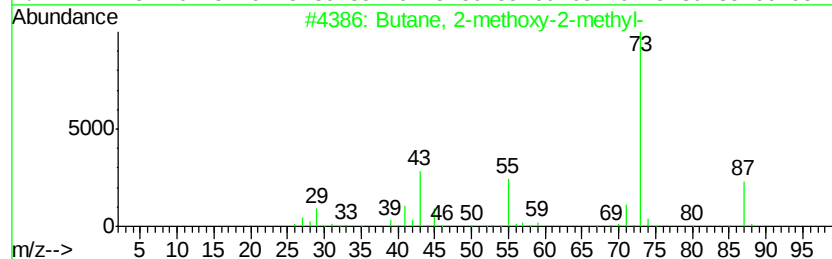
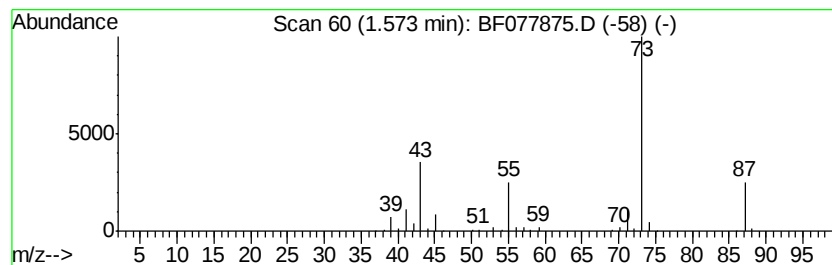
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.57	25.25 ng	318113	1,4-Dichlorobenzene-d4	7.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
3		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	12



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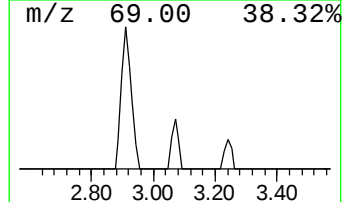
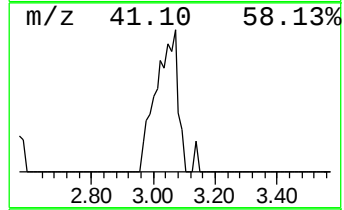
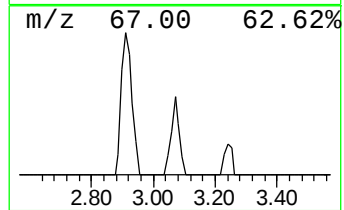
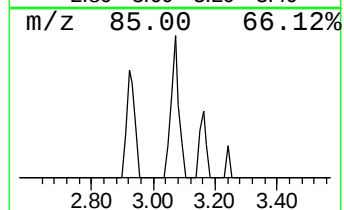
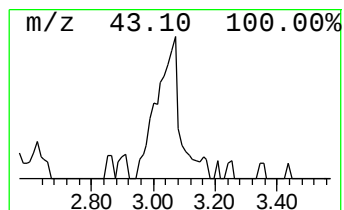
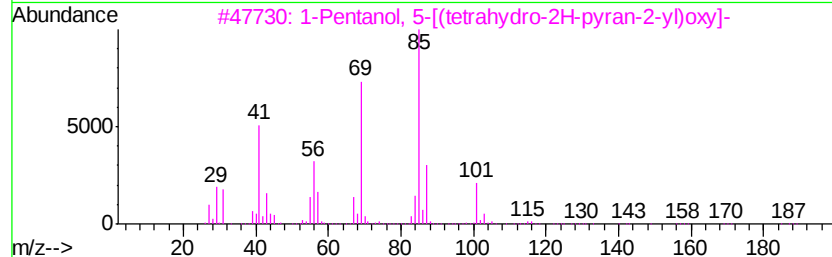
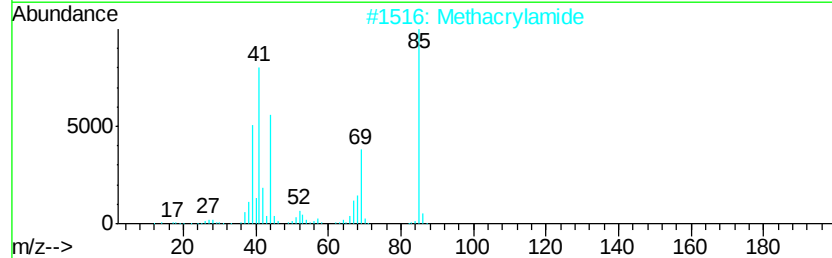
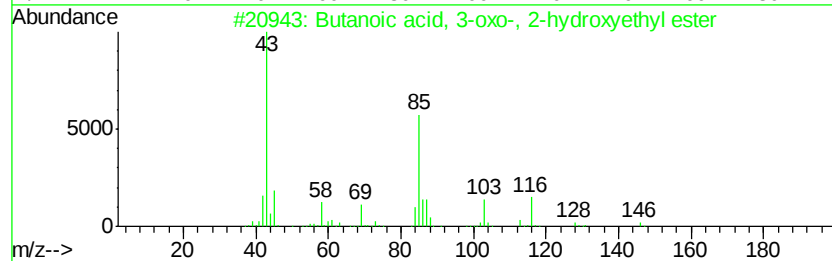
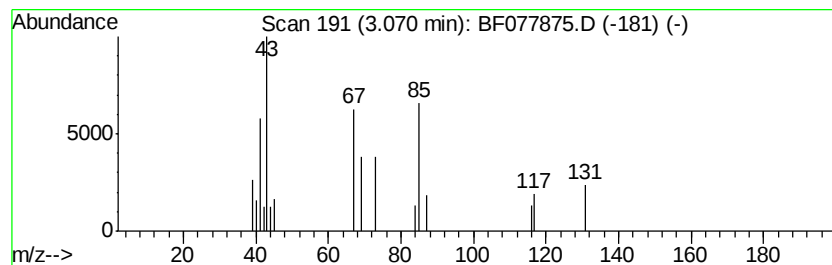
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 unknown3.07 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.07	3.22 ng	40625	1,4-Dichlorobenzene-d4	7.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, 3-oxo-, 2-hydroxy...	146	C6H10O4	033736-01-5	35
2		Methacrylamide	85	C4H7NO	000079-39-0	14
3		1-Pentanol, 5-[(tetrahydro-2H-py...	188	C10H20O3	076102-74-4	12
4		Pentane, 2-bromo-2-methyl-	164	C6H13Br	004283-80-1	10
5		2-Pyrrolidinone	85	C4H7NO	000616-45-5	9



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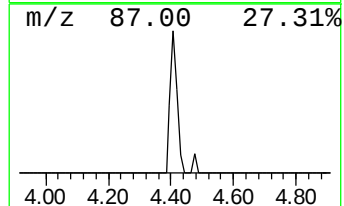
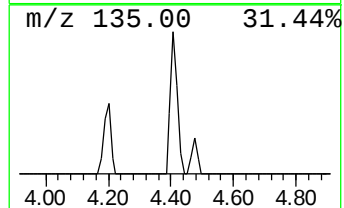
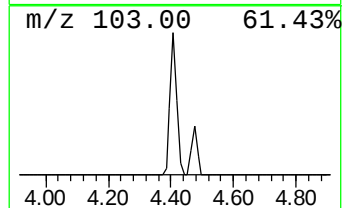
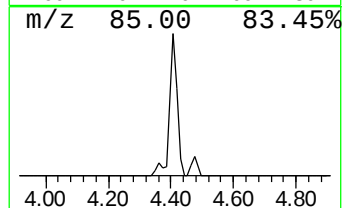
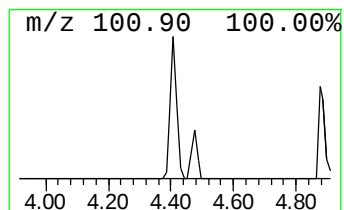
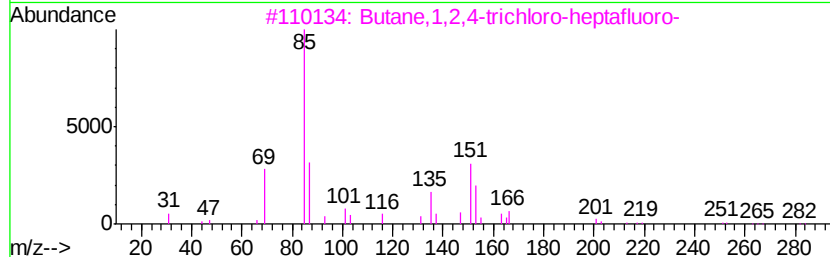
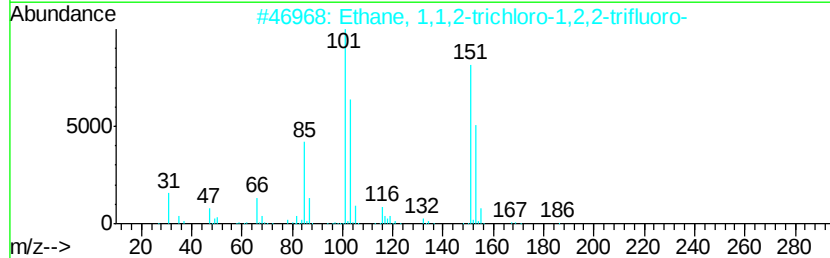
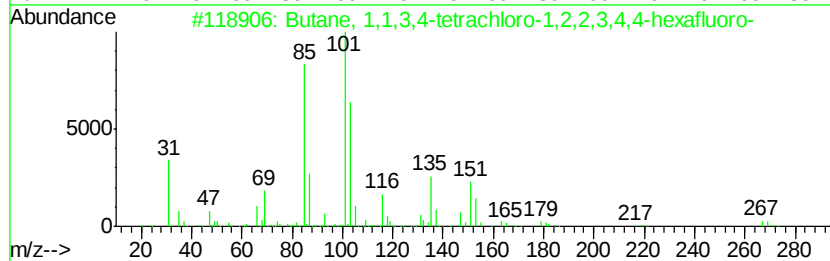
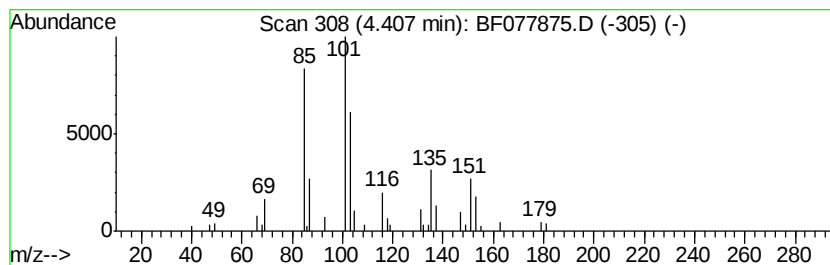
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Butane, 1,1,3,4-tetrachloro... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.41	6.10 ng	76827	1,4-Dichlorobenzene-d4	7.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 1,1,3,4-tetrachloro-1,2,...	302	C4Cl4F6	000423-38-1	91
2			Ethane, 1,1,2-trichloro-1,2,2-tr...	186	C2Cl3F3	000076-13-1	47
3			Butane,1,2,4-trichloro-heptafluoro-	286	C4Cl3F7	000335-45-5	47
4			Methane, bromodichlorofluoro-	180	CBrCl2F	000353-58-2	27
5			Ethane, 1,1,2,2-tetrachloro-1-fl...	184	C2HCl4F	000354-14-3	25



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077875.D
Acq On : 20 Mar 2015 20:54
Operator : TP/IZ
Sample : G1575-08
Misc : W
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
RW-9V

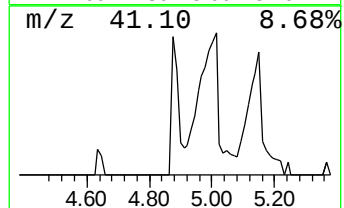
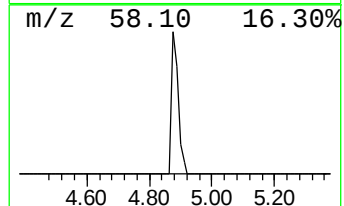
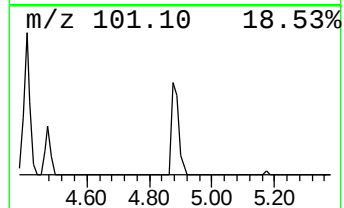
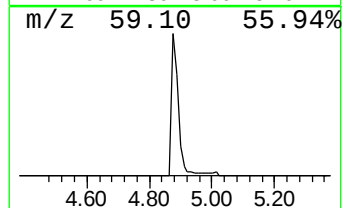
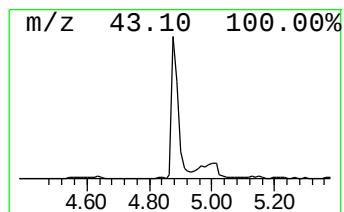
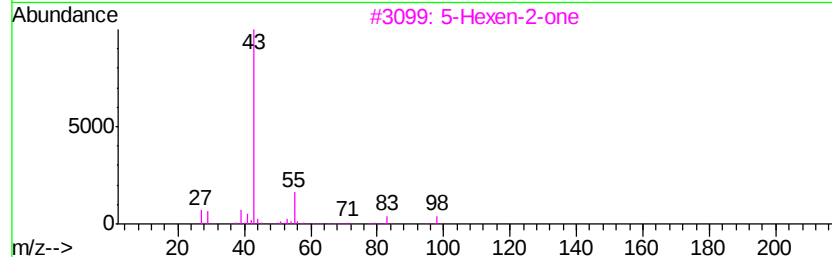
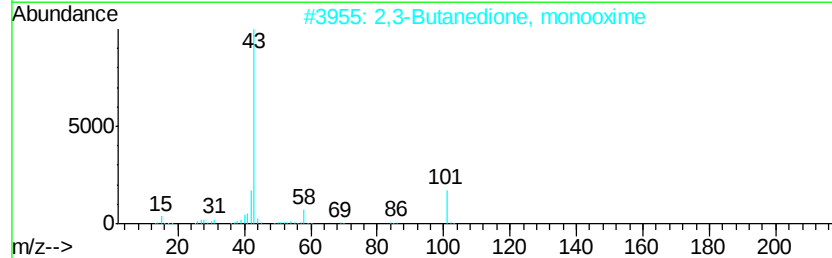
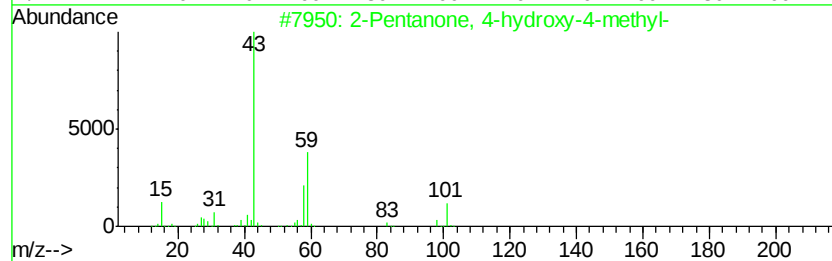
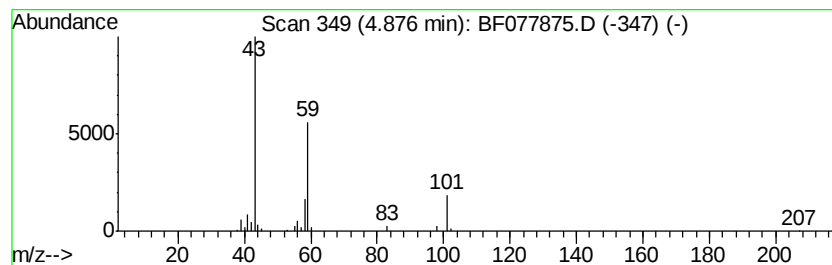
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.88	9.78 ng	123169	1,4-Dichlorobenzene-d4	7.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
3			5-Hexen-2-one	98	C6H10O	000109-49-9	9
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032015\
Data File : BF077875.D
Acq On : 20 Mar 2015 20:54
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Misc : W
ALS Vial : 14 Sample Multiplier: 1

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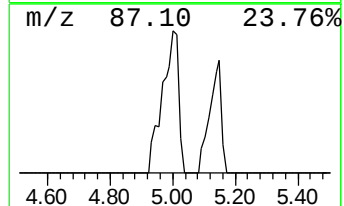
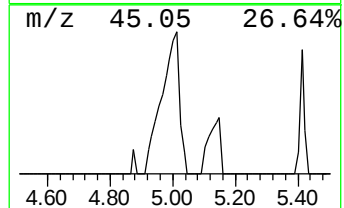
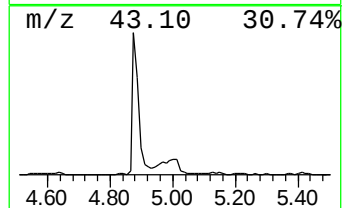
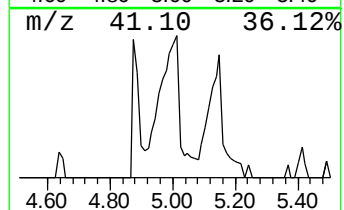
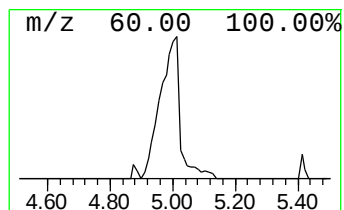
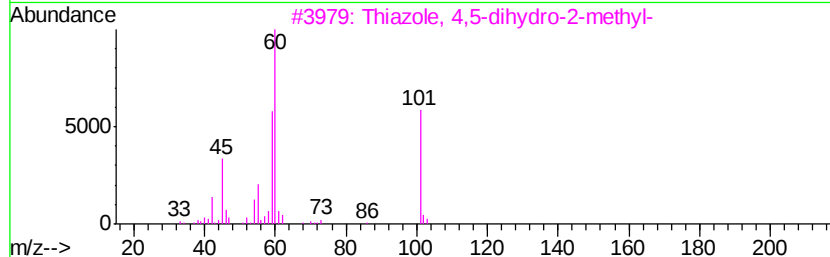
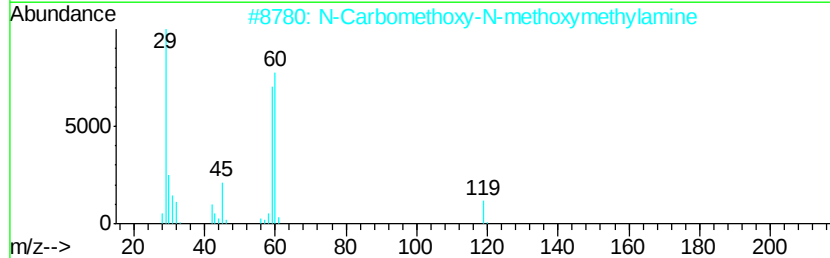
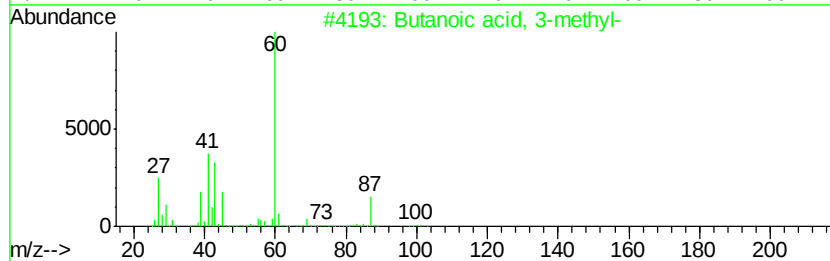
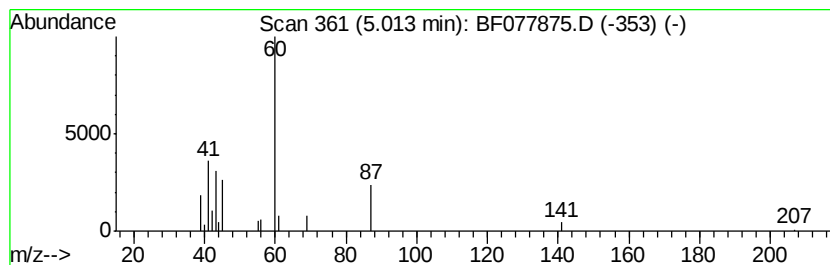
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Butanoic acid, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	8.59 ng	108159	1,4-Dichlorobenzene-d4	7.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	56
2			N-Carbomethoxy-N-methoxymethylamine	119	C4H9NO3	153654-07-0	38
3			Thiazole, 4,5-dihydro-2-methyl-	101	C4H7NS	002346-00-1	9
4			L-Mannose, 6-deoxy-	164	C6H12O5	003615-41-6	9
5			Pentanoic acid	102	C5H10O2	000109-52-4	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077875.D
Acq On : 20 Mar 2015 20:54
Operator : TP/IZ
Sample : G1575-08
Misc : W
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
RW-9V

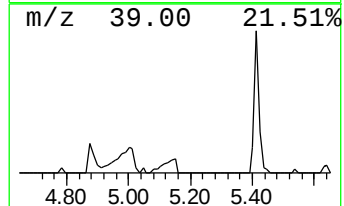
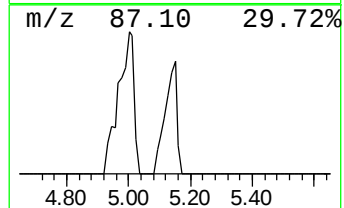
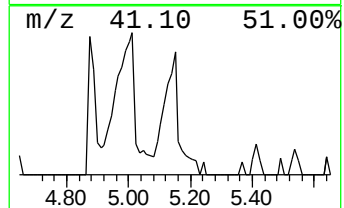
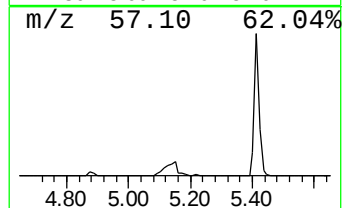
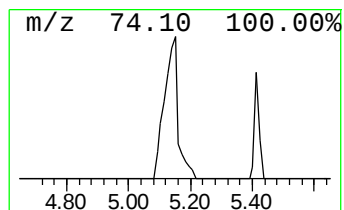
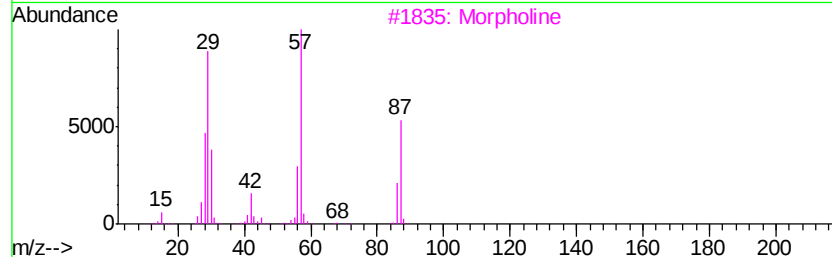
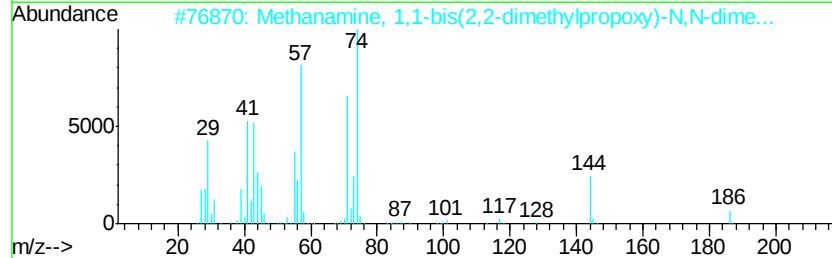
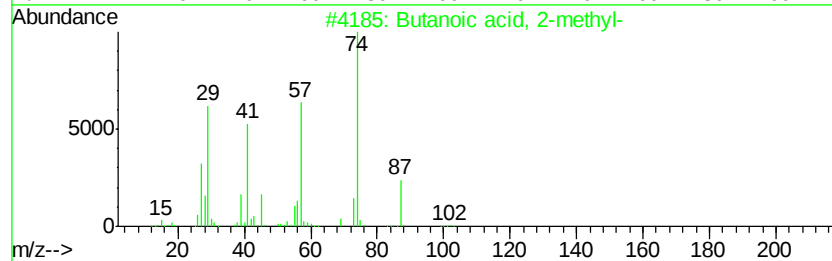
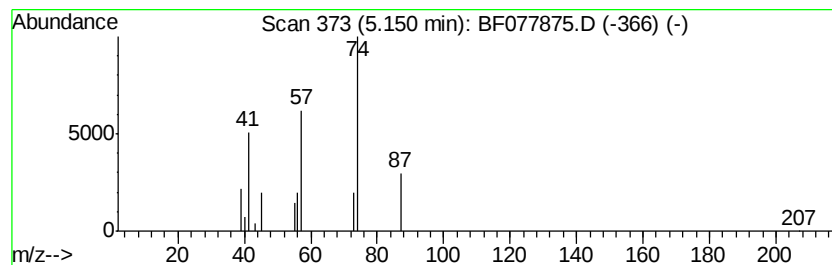
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Butanoic acid, 2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	4.60 ng	57996	1,4-Dichlorobenzene-d4	7.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, 2-methyl-	102	C5H10O2	000116-53-0	78
2			Methanamine, 1,1-bis(2,2-dimethylpropoxy)-N,N-dimethyl-	231	C13H29NO2	004909-78-8	50
3			Morpholine	87	C4H9NO	000110-91-8	9
4			2-Propenoic acid, 2-methyl-, butanoic acid, 2-methyl-	142	C8H14O2	000097-88-1	9
5			1,2-Butanediol, 3,3-dimethyl-	118	C6H14O2	059562-82-2	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077875.D
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Sample : G1575-08
Misc : W
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Instrument :
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ClientSampleId :
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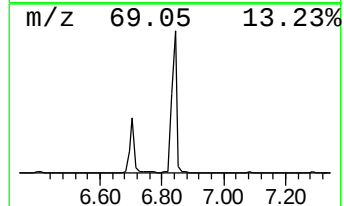
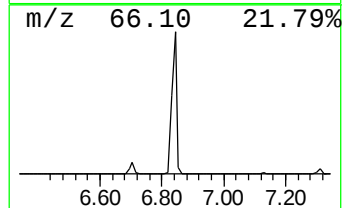
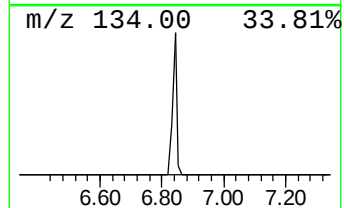
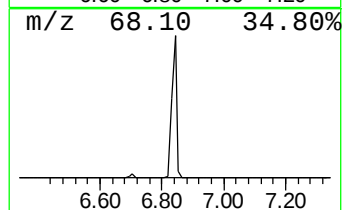
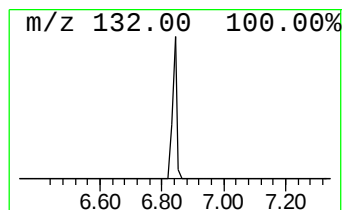
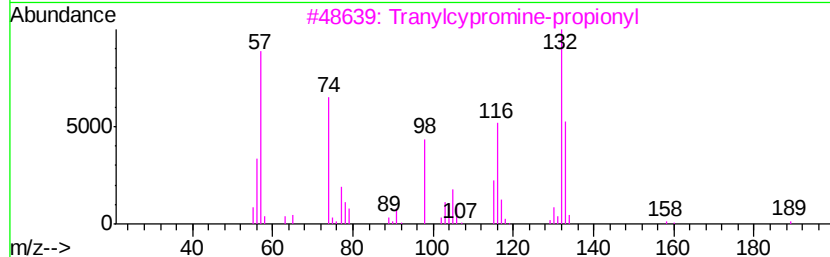
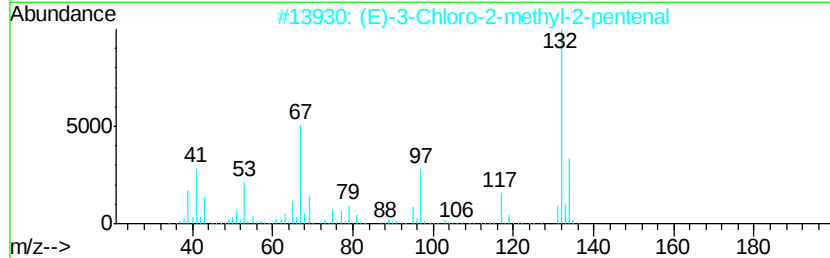
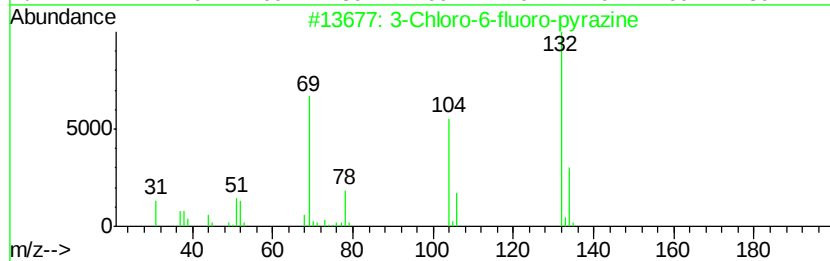
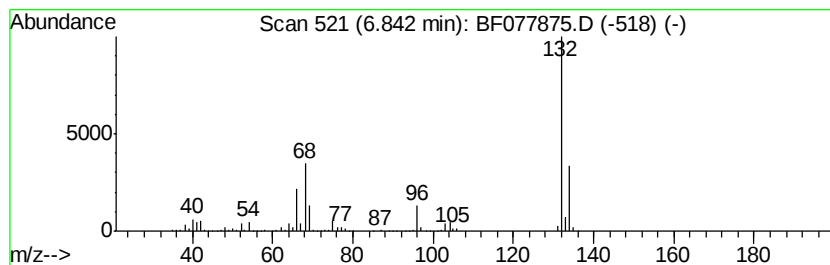
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown6.84 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.84	93.39 ng	1176420	1,4-Dichlorobenzene-d4	7.13

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	17
3		Tranvlcvpromine-propionyl	189	C12H15NO	1000123-86-3	12
4		Xenon	132	Xe	007440-63-3	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077875.D
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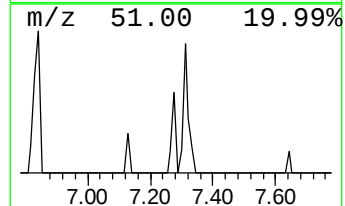
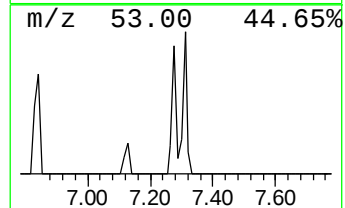
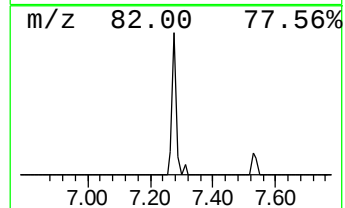
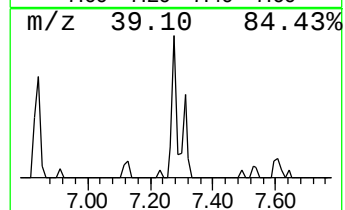
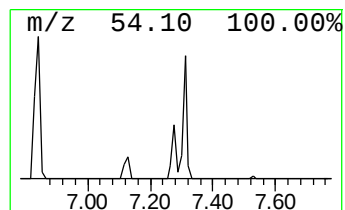
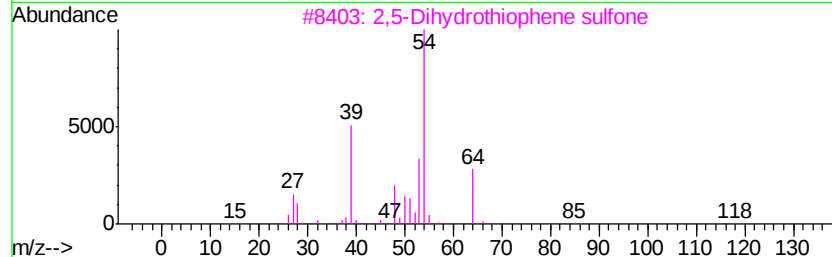
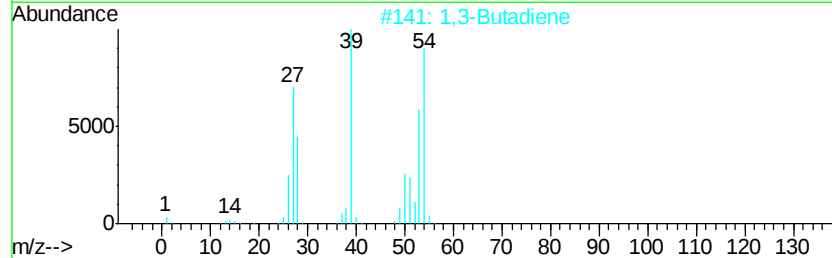
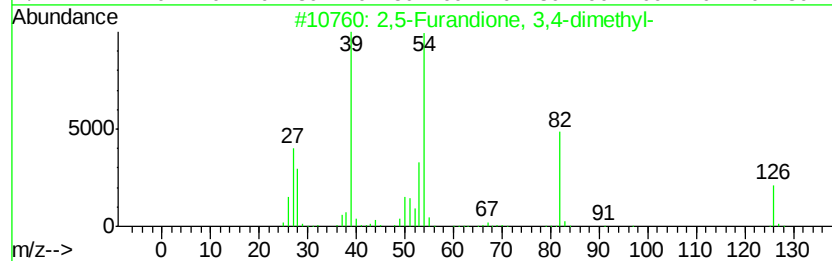
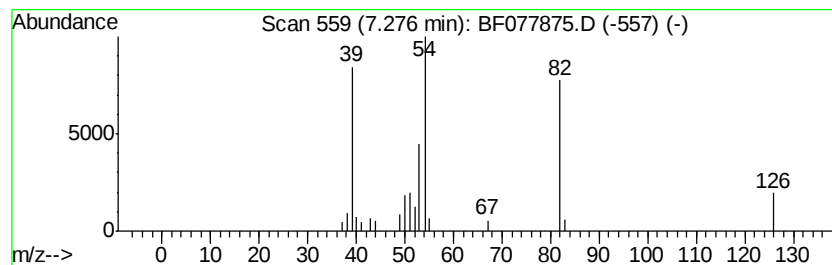
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 2,5-Furandione, 3,4-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.28	2.12 ng	26709	1,4-Dichlorobenzene-d4	7.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,5-Furandione, 3,4-dimethyl-	126	C6H6O3	000766-39-2	91
2			1,3-Butadiene	54	C4H6	000106-99-0	80
3			2,5-Dihydrothiophene sulfone	118	C4H6O2S	000077-79-2	78
4			2-Cyclopenten-1-one	82	C5H6O	000930-30-3	74
5			2-Cyclohexen-1-one, 2-methyl-	110	C7H10O	001121-18-2	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032015\
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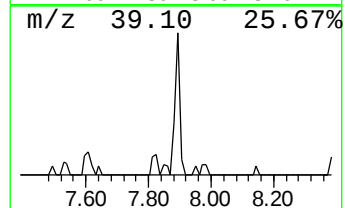
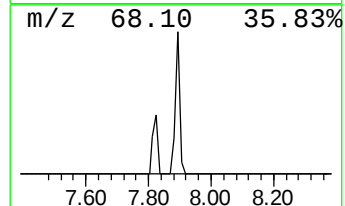
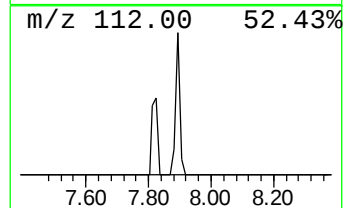
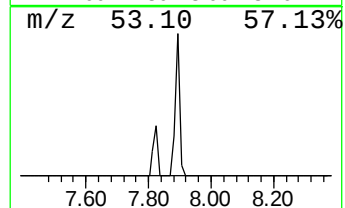
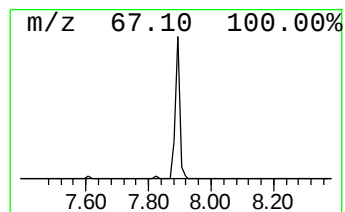
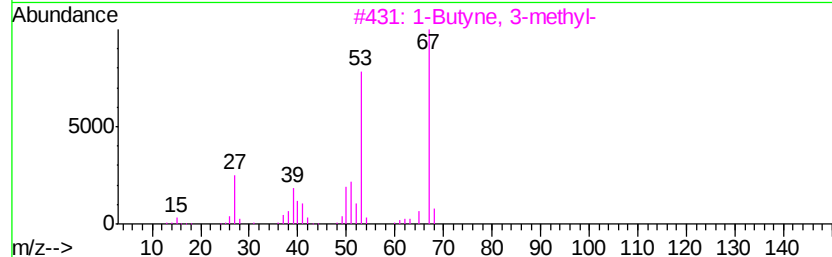
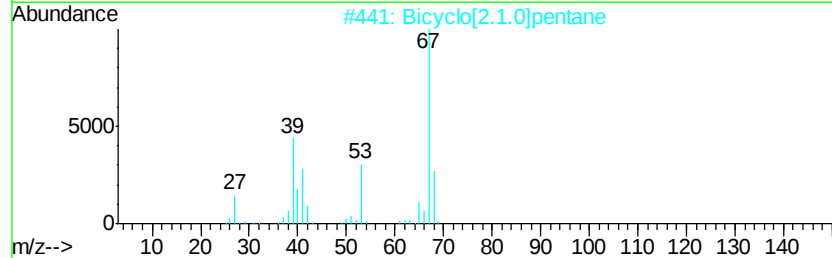
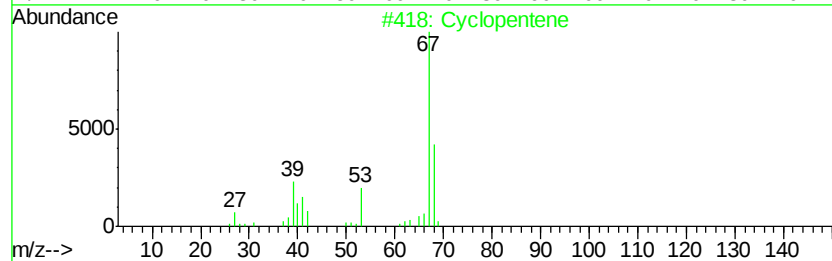
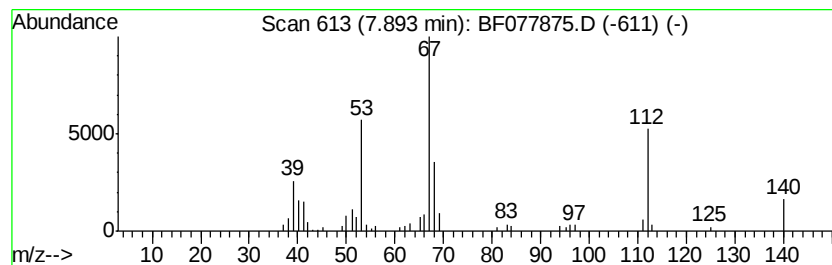
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Cyclopentene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.89	5.82 ng	73311	1,4-Dichlorobenzene-d4	7.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentene	68	C5H8	000142-29-0	72
2			Bicyclo[2.1.0]pentane	68	C5H8	000185-94-4	64
3			1-Butyne, 3-methyl-	68	C5H8	000598-23-2	59
4			1,4-Pentadiene	68	C5H8	000591-93-5	45
5			Spiropentane	68	C5H8	000157-40-4	43



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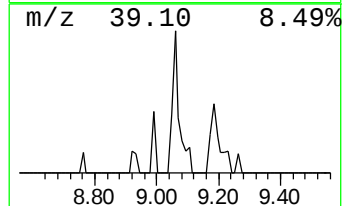
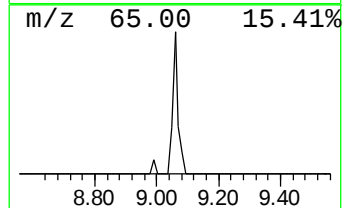
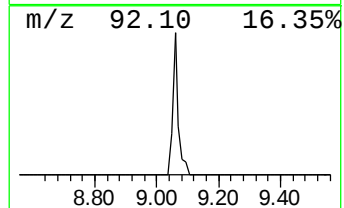
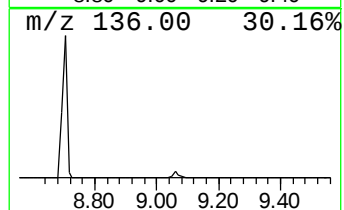
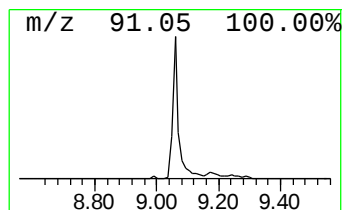
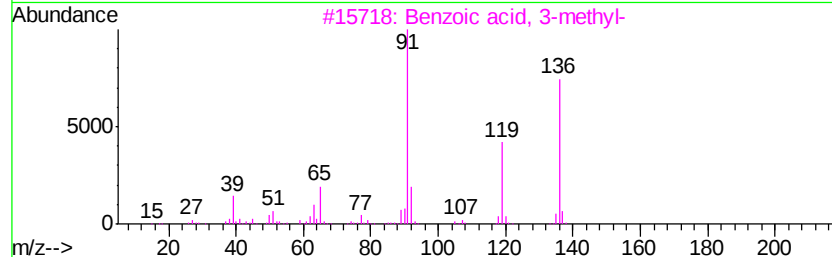
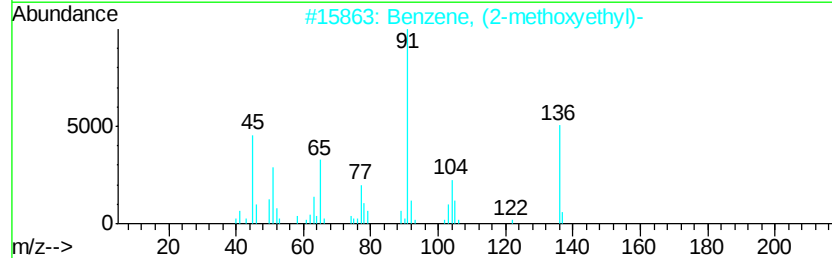
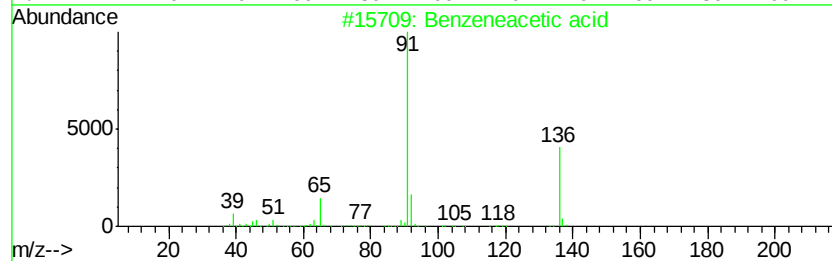
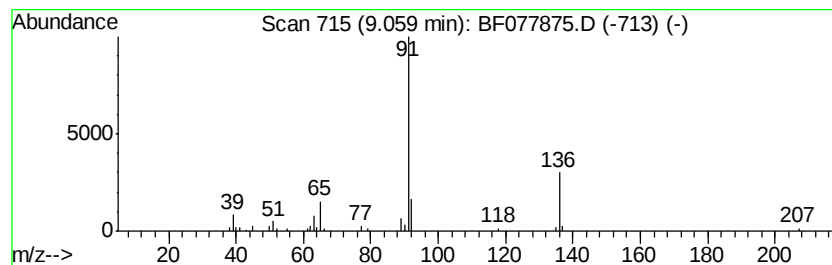
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Benzeneacetic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.06	4.43 ng	76969	Napthalene-d8	8.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzeneacetic acid	136	C8H8O2	000103-82-2	87
2		Benzene, (2-methoxyethyl)-	136	C9H12O	003558-60-9	72
3		Benzoic acid, 3-methyl-	136	C8H8O2	000099-04-7	72
4		Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	59
5		Benzene, (bromomethyl)-	170	C7H7Br	000100-39-0	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032015\
Data File : BF077875.D
Acq On : 20 Mar 2015 20:54
Operator : TP/IZ
Sample : G1575-08
Misc : W
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RW-9V

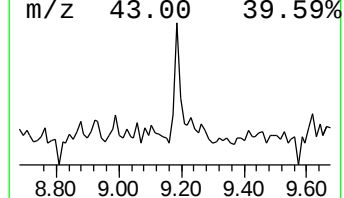
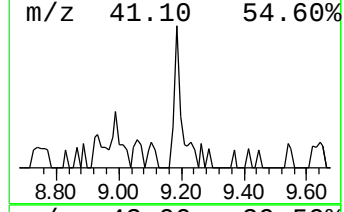
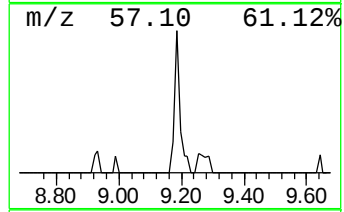
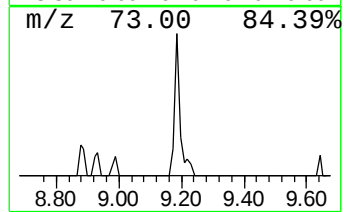
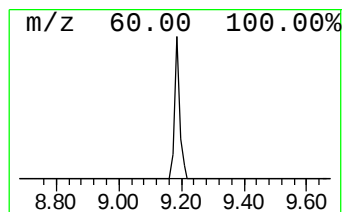
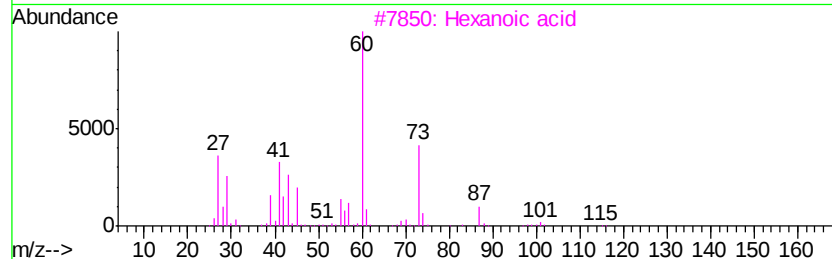
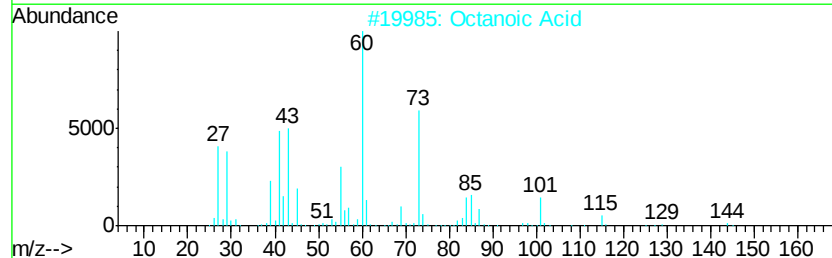
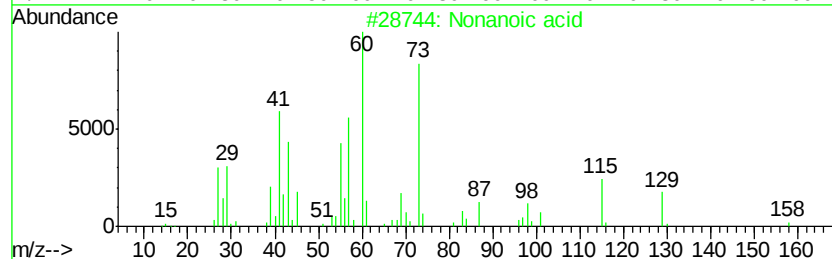
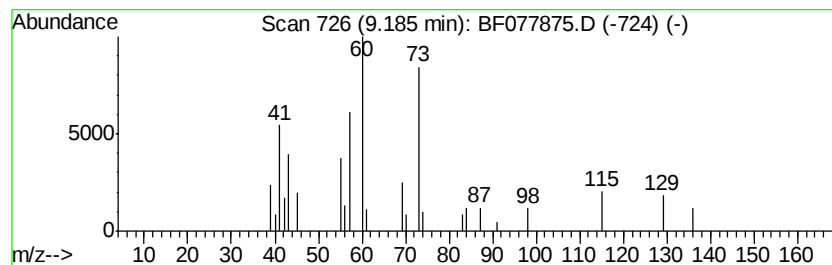
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Nonanoic acid Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.18	2.04 ng	35368	Napthalene-d8	8.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonanoic acid	158	C9H18O2	000112-05-0	64
2			Octanoic Acid	144	C8H16O2	000124-07-2	53
3			Hexanoic acid	116	C6H12O2	000142-62-1	50
4			Propanedioic acid, propyl-	146	C6H10O4	000616-62-6	38
5			D-Mannoheptulose	210	C7H14O7	003615-44-9	33



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032015\
Data File : BF077875.D
Acq On : 20 Mar 2015 20:54
Operator : TP/IZ
Sample : G1575-08
Misc : W
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RW-9V

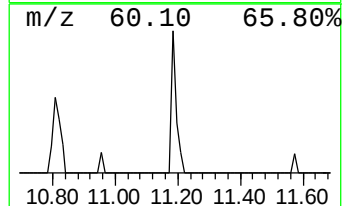
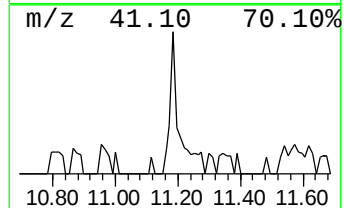
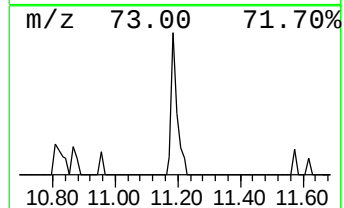
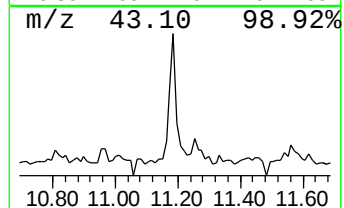
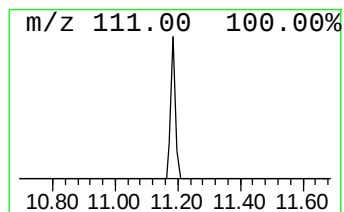
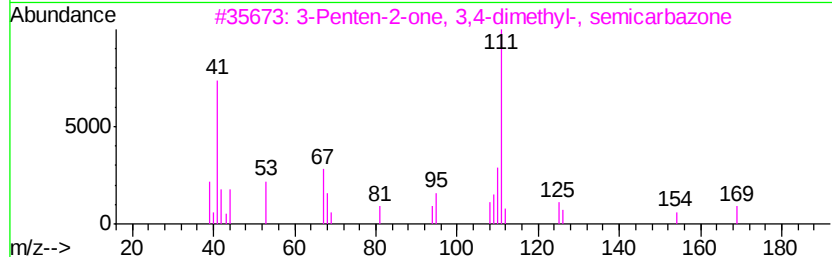
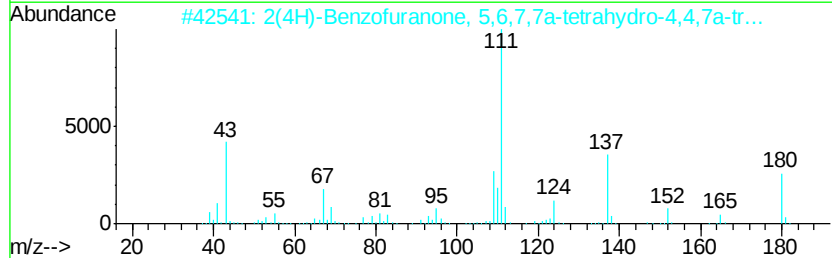
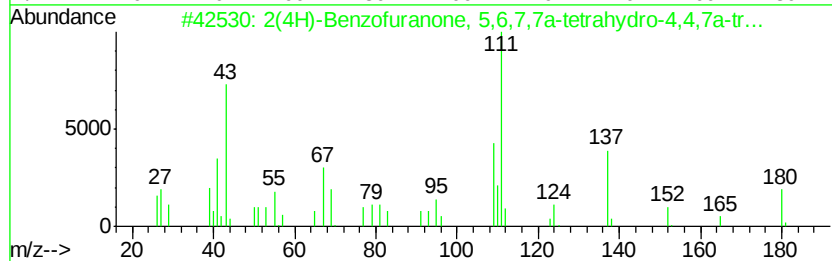
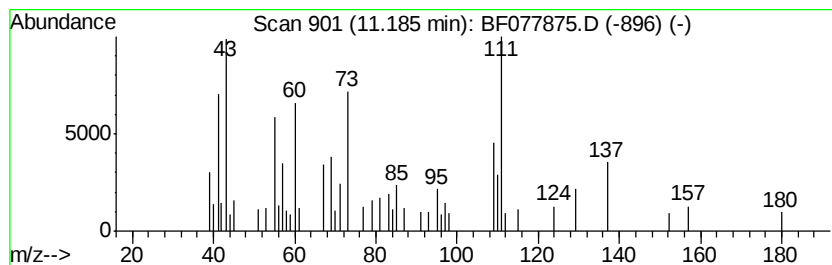
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 2(4H)-Benzofuranone, 5,6,7,... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.19	2.55 ng	51489	Acenaphthene-d10	10.87

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2(4H)-Benzofuranone, 5,6,7,7a-te...	180	C11H16O2	015356-74-8	93
2			2(4H)-Benzofuranone, 5,6,7,7a-te...	180	C11H16O2	017092-92-1	38
3			3-Penten-2-one, 3,4-dimethyl-, s...	169	C8H15N3O	016983-60-1	35
4			p-Menthan-2-one, 9-hydroxy-, (1R...	170	C10H18O2	021700-69-6	35
5			Cyclooctane, cyclohexyl-	194	C14H26	092369-78-3	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032015\
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Sample : G1575-08
Misc : W
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RW-9V

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.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
Ethanimidamide, N...	1.05	2.5	ng	31696	1	7.13	251950	20.0
Roflurane	1.31	3.7	ng	46685	1	7.13	251950	20.0
Butane, 2-methoxy...	1.57	25.3	ng	318113	1	7.13	251950	20.0
unknown3.07	3.07	3.2	ng	40625	1	7.13	251950	20.0
Butane, 1,1,3,4-t...	4.41	6.1	ng	76827	1	7.13	251950	20.0
2-Pentanone, 4-hy...	4.88	9.8	ng	123169	1	7.13	251950	20.0
Butanoic acid, 3-...	5.01	8.6	ng	108159	1	7.13	251950	20.0
Butanoic acid, 2-...	5.15	4.6	ng	57996	1	7.13	251950	20.0
unknown6.84	6.84	93.4	ng	1176420	1	7.13	251950	20.0
2,5-Furandione, 3...	7.28	2.1	ng	26709	1	7.13	251950	20.0
Cyclopentene	7.89	5.8	ng	73311	1	7.13	251950	20.0
Benzeneacetic acid	9.06	4.4	ng	76969	2	8.70	347308	20.0
Nonanoic acid	9.18	2.0	ng	35368	2	8.70	347308	20.0
2(4H)-Benzofurano...	11.19	2.5	ng	51489	3	10.87	403529	20.0