

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032515\
 Data File : BF077956.D
 Acq On : 25 Mar 2015 17:30
 Operator : TP/IZ
 Sample : G1612-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-1

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	1.150	21	23	26	rBV	1186473	1150590	35.51%	6.592%
2	1.356	38	41	44	rVB	179094	267841	8.27%	1.535%
3	1.516	53	55	58	rBV	210408	254308	7.85%	1.457%
4	1.744	74	75	81	rBV	56680	82882	2.56%	0.475%
5	1.973	91	95	98	rBV	1874494	3240495	100.00%	18.565%
6	2.407	130	133	142	rVB	87644	143681	4.43%	0.823%
7	3.870	258	261	266	rVB	54618	92455	2.85%	0.530%
8	4.830	344	345	350	rVB	32966	47567	1.47%	0.273%
9	5.379	390	393	396	rBV	325819	392088	12.10%	2.246%
10	5.642	413	416	418	rVB	49259	60459	1.87%	0.346%
11	5.950	442	443	446	rVB	37528	42560	1.31%	0.244%
12	6.602	498	500	502	rBV2	40184	42106	1.30%	0.241%
13	6.670	504	506	508	rVV	275154	266903	8.24%	1.529%
14	6.705	508	509	511	rVV	79621	58698	1.81%	0.336%
15	6.808	515	518	520	rVV	808409	781837	24.13%	4.479%
16	6.876	520	524	531	rVB	40514	55276	1.71%	0.317%
17	7.093	541	543	545	rBV	230625	207275	6.40%	1.188%
18	7.185	548	551	554	rVB	67132	66780	2.06%	0.383%
19	7.276	557	559	561	rVV	705539	732784	22.61%	4.198%
20	7.310	561	562	563	rVV	46278	34716	1.07%	0.199%
21	7.390	566	569	572	rVV3	52656	153983	4.75%	0.882%
22	7.459	572	575	578	rVV	53452	117000	3.61%	0.670%
23	7.791	602	604	605	rVV	743080	775362	23.93%	4.442%
24	7.836	605	608	610	rVV2	321262	688279	21.24%	3.943%
25	7.871	610	611	616	rVV2	140367	186014	5.74%	1.066%
26	7.951	616	618	622	rVB5	21106	41025	1.27%	0.235%
27	8.156	632	636	640	rVB2	70624	112655	3.48%	0.645%
28	8.351	651	653	655	rBV2	56296	69723	2.15%	0.399%
29	8.591	670	674	679	rBV4	16055	43419	1.34%	0.249%
30	8.671	679	681	685	rVB	342496	316175	9.76%	1.811%
31	9.334	736	739	742	rBV	56257	64187	1.98%	0.368%
32	9.939	791	792	796	rBV2	43110	58262	1.80%	0.334%
33	10.019	796	799	801	rVB	1447967	1253871	38.69%	7.184%
34	10.282	819	822	826	rBV	181462	217777	6.72%	1.248%

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Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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35	10.442	834	836	838	rVV	19715	32577	1.01%	0.187%
36	10.842	868	871	873	rVB	341742	355027	10.96%	2.034%
37	11.688	942	945	947	rBV	949093	891293	27.50%	5.106%
38	11.814	954	956	959	rBV2	604601	773375	23.87%	4.431%
39	12.671	1029	1031	1034	rVB	277987	360425	11.12%	2.065%
40	13.014	1059	1061	1064	rBV	478346	423857	13.08%	2.428%
41	14.671	1204	1206	1209	rBV	1170056	1553667	47.95%	8.901%
42	15.917	1312	1315	1321	rBV2	28197	57709	1.78%	0.331%
43	16.020	1321	1324	1327	rBV	400684	445814	13.76%	2.554%
44	17.940	1489	1492	1495	rBV	333845	441750	13.63%	2.531%

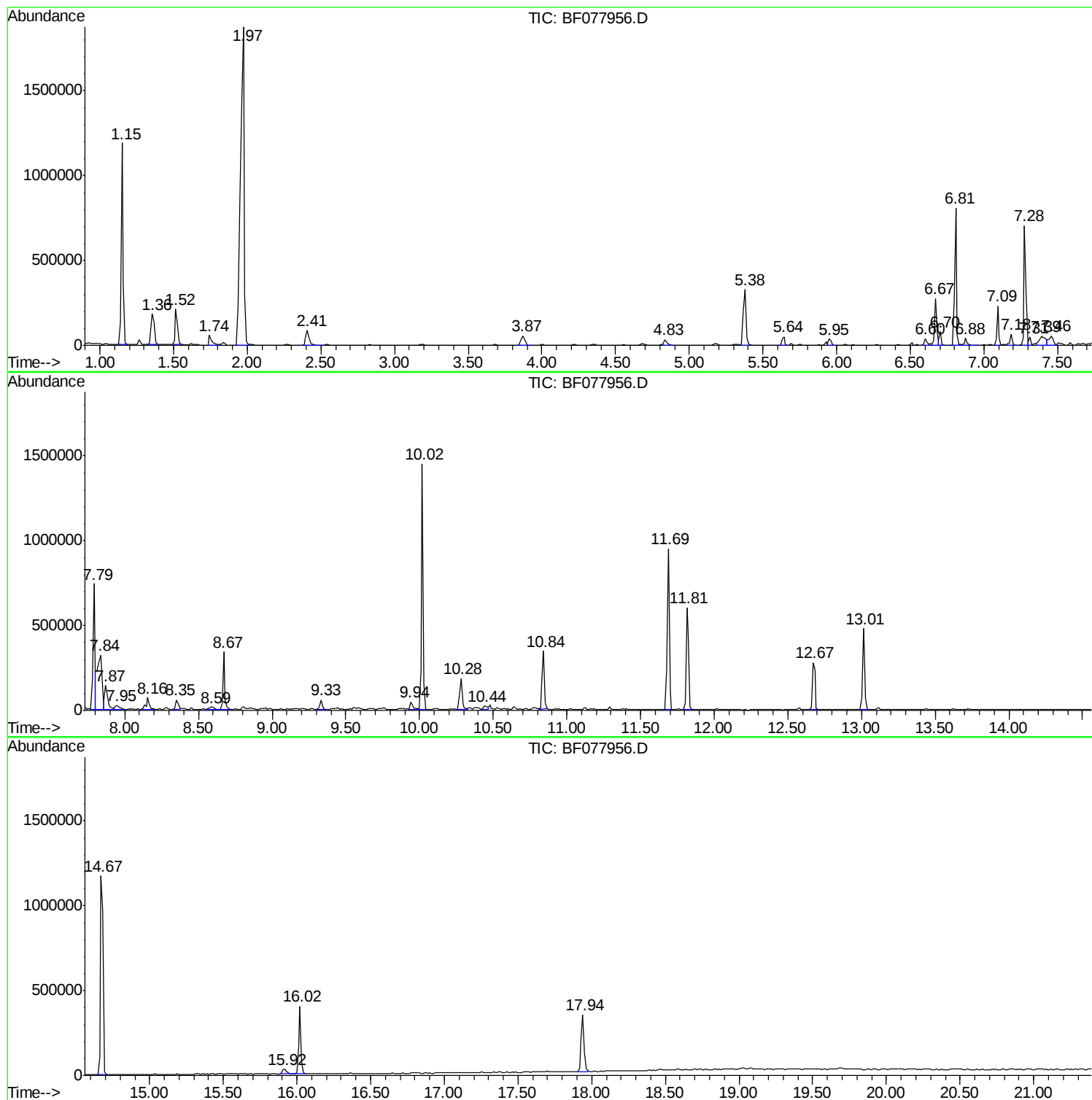
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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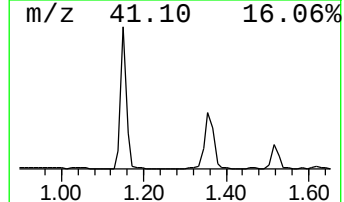
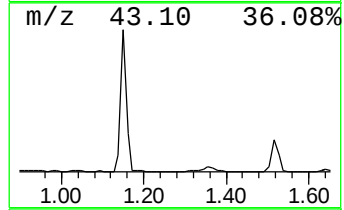
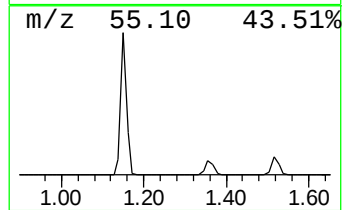
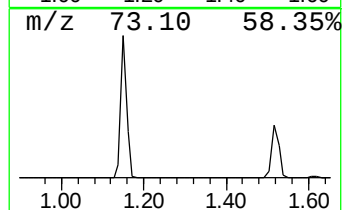
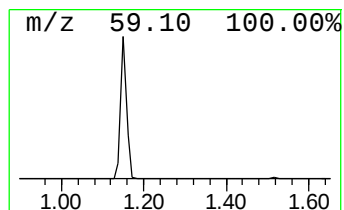
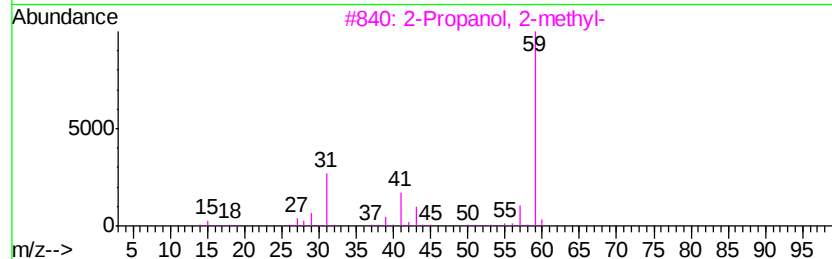
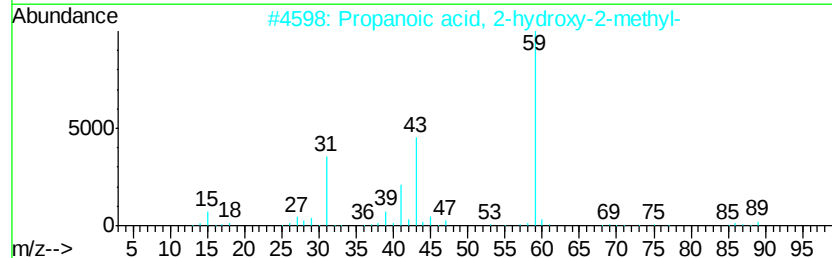
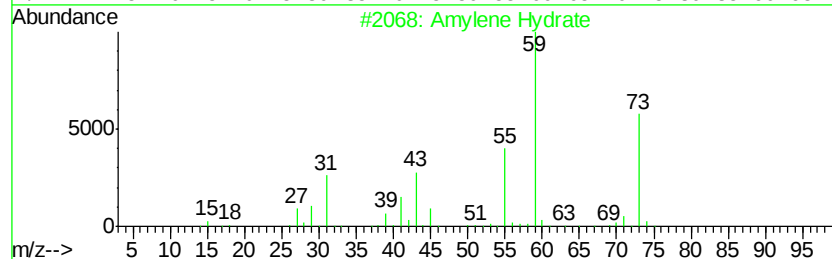
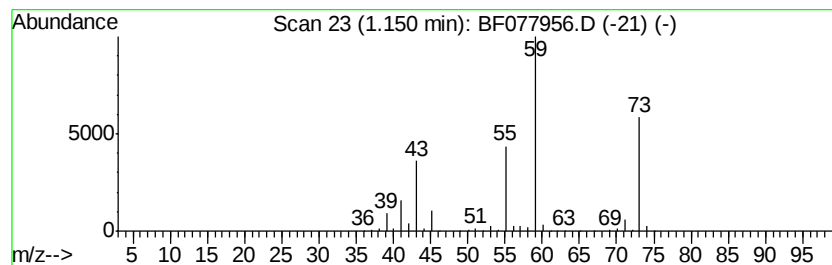
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Amylene Hydrate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.15	111.02 ng	1150590	1,4-Dichlorobenzene-d4	7.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Amylene Hydrate	88	C5H12O	000075-85-4	83
2		Propanoic acid, 2-hydroxy-2-methyl-	104	C4H8O3	000594-61-6	53
3		2-Propanol, 2-methyl-	74	C4H10O	000075-65-0	42
4		3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	40
5		Silane, trimethyl-	74	C3H10Si	000993-07-7	39



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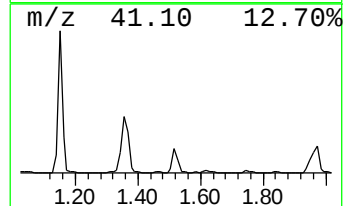
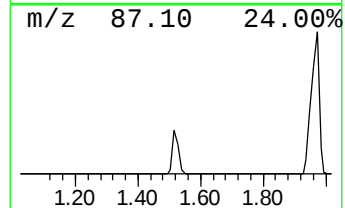
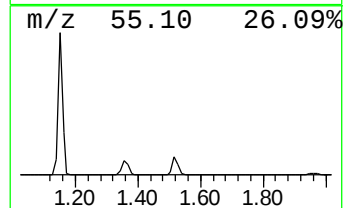
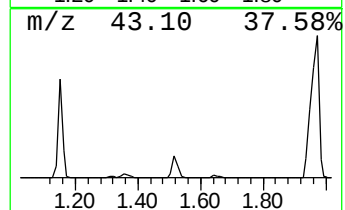
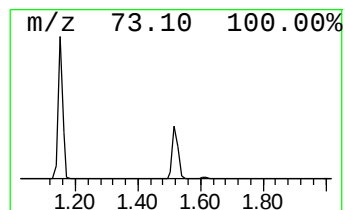
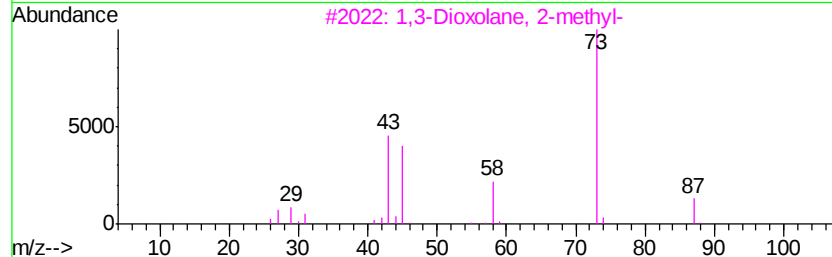
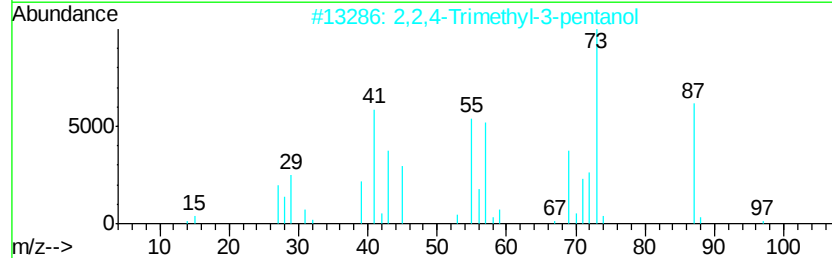
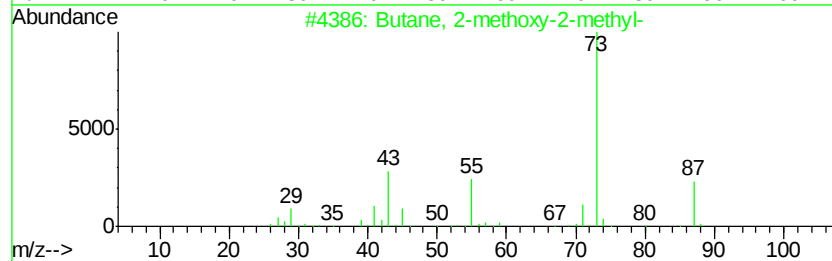
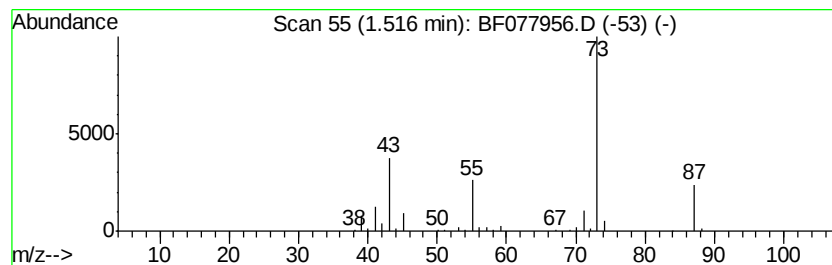
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.52	24.54 ng	254308	1,4-Dichlorobenzene-d4	7.09

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	12



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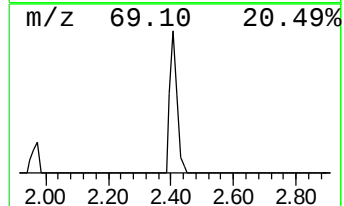
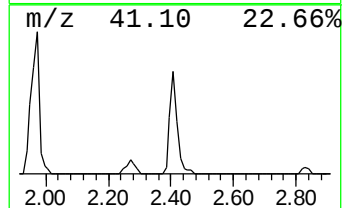
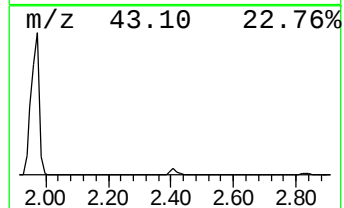
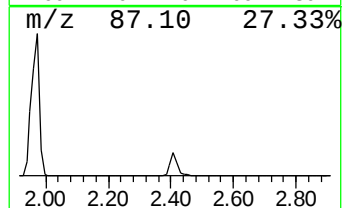
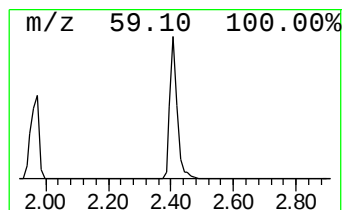
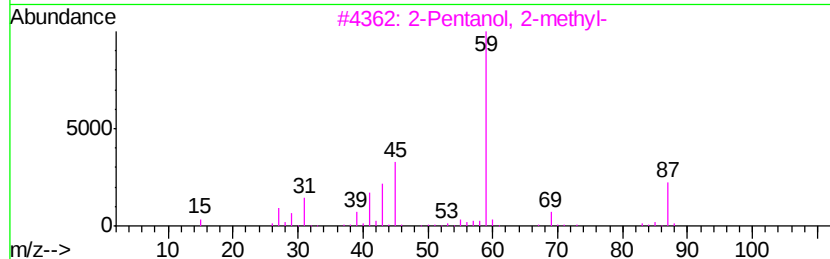
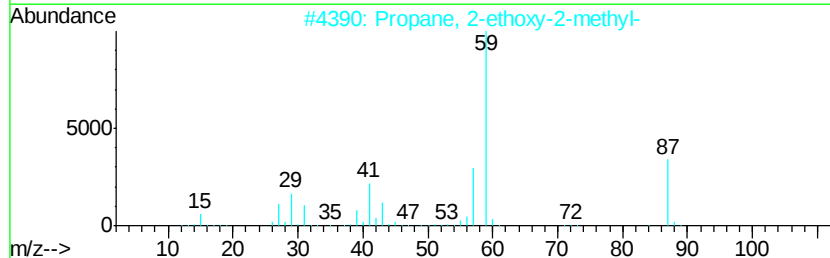
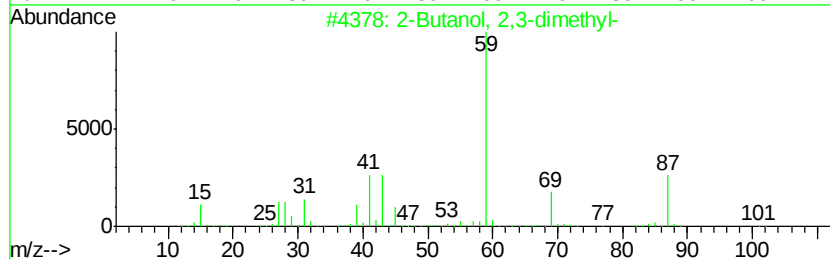
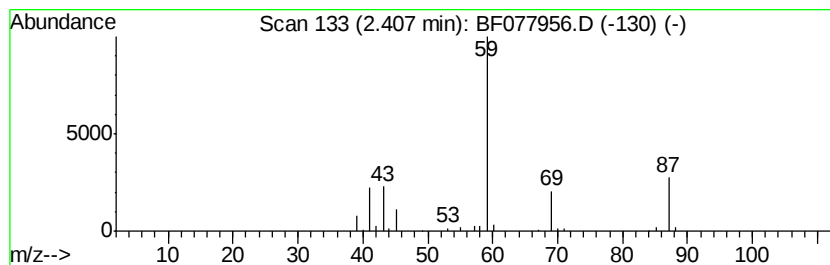
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2-Butanol, 2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.41	13.86 ng	143681	1,4-Dichlorobenzene-d4	7.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	83
2		Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	64
3		2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	56
4		Propane, 1-ethoxy-	88	C5H12O	000628-32-0	53
5		3-Pentanol, 2,2-dimethyl-	116	C7H16O	003970-62-5	40



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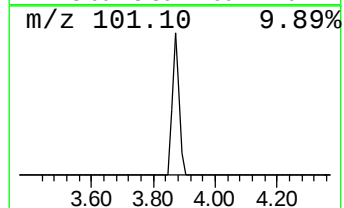
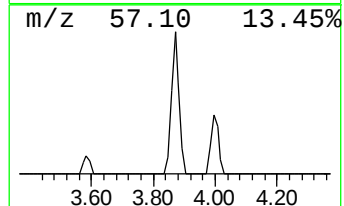
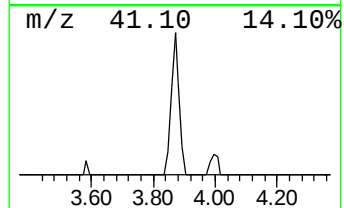
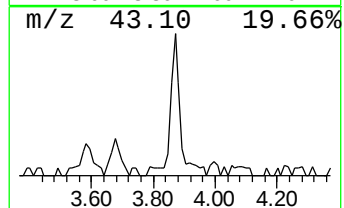
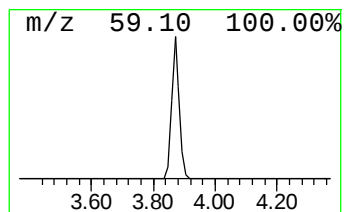
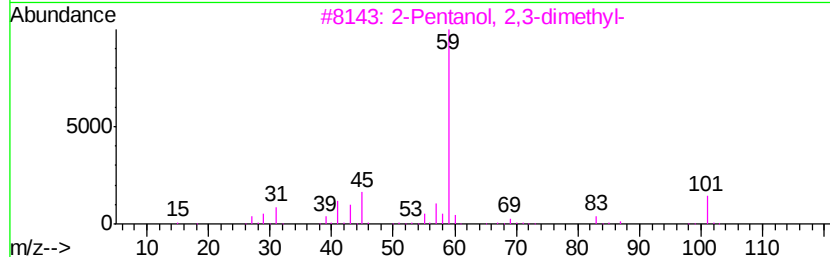
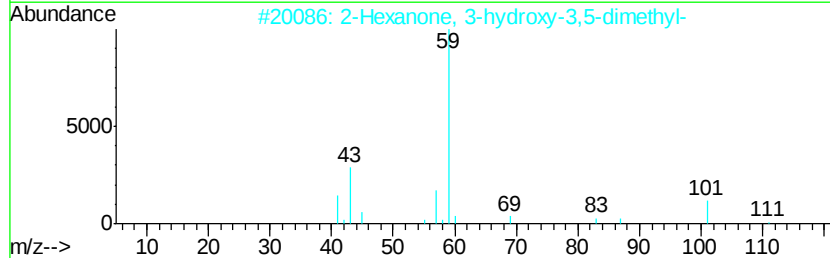
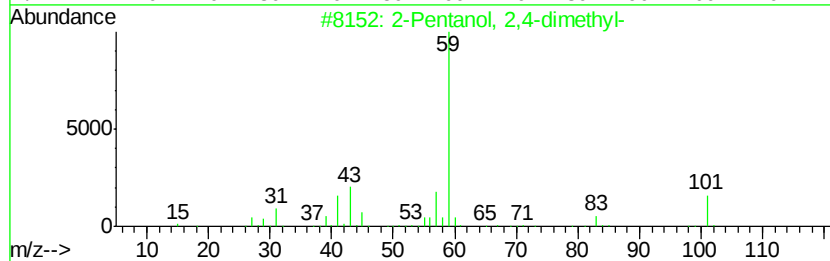
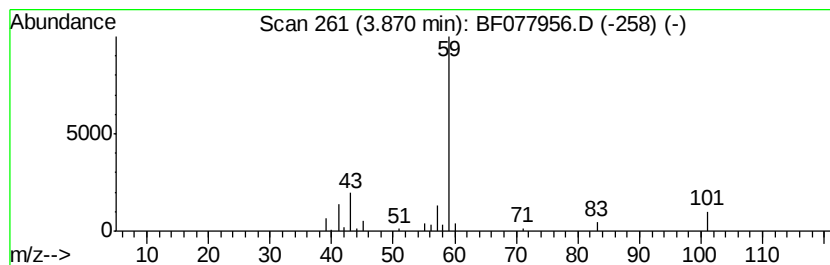
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 2-Pentanol, 2,4-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.87	8.92 ng	92455	1,4-Dichlorobenzene-d4	7.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	83
2		2-Hexanone, 3-hydroxy-3,5-dimethyl-	144	C8H16O2	006321-14-8	72
3		2-Pentanol, 2,3-dimethyl-	116	C7H16O	004911-70-0	64
4		2-Propanol, 2-methyl-	74	C4H10O	000075-65-0	50
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	40



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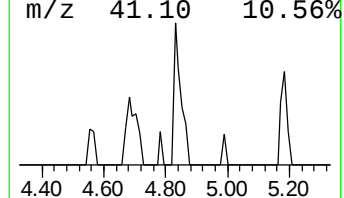
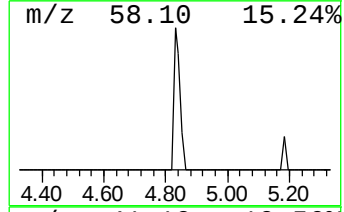
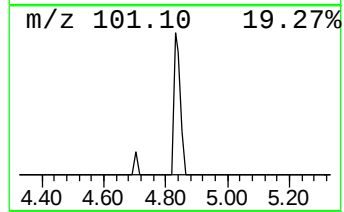
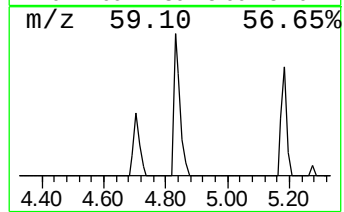
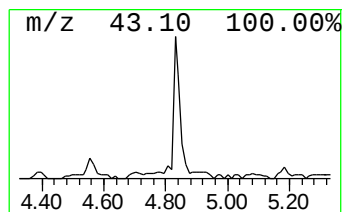
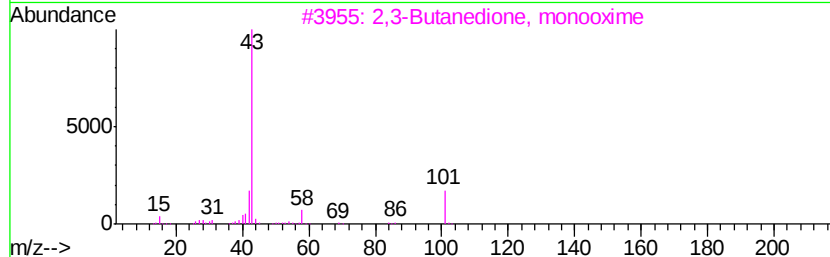
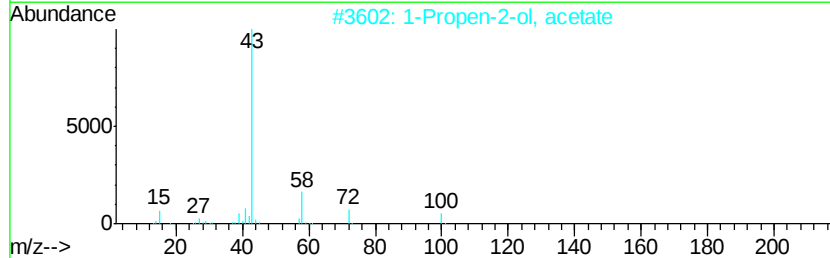
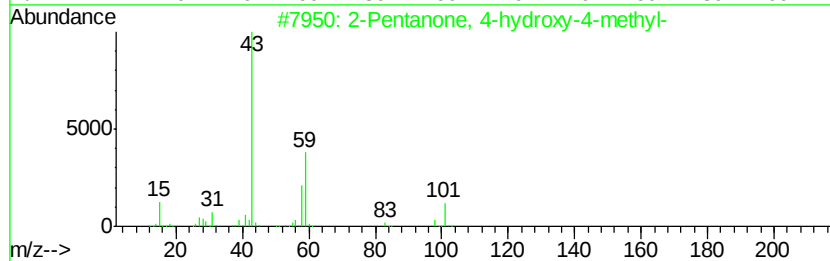
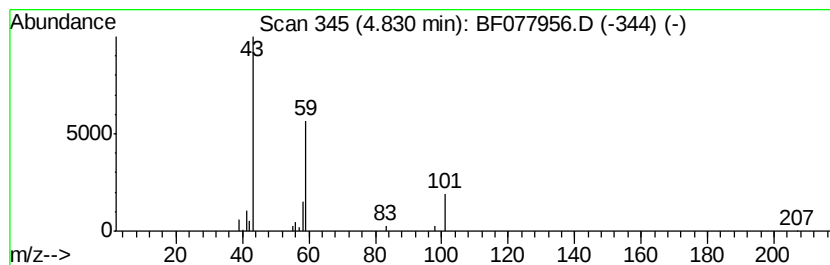
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.83	4.59 ng	47567	1,4-Dichlorobenzene-d4	7.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	7
4		Ethanone, 1-(2-methylcyclopropyl)-	98	C6H10O	000930-56-3	5
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	5



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032515\
Data File : BF077956.D
Acq On : 25 Mar 2015 17:30
Operator : TP/IZ
Sample : G1612-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

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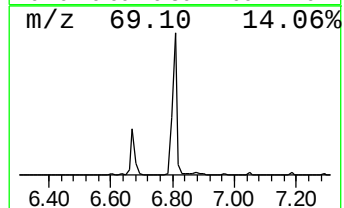
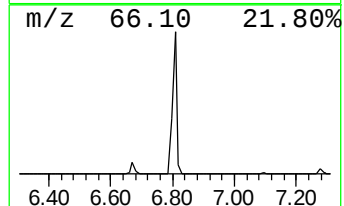
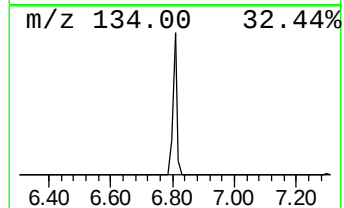
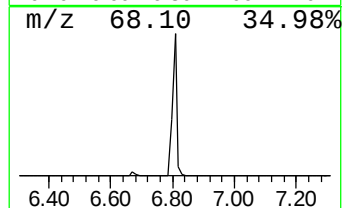
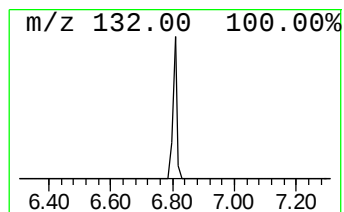
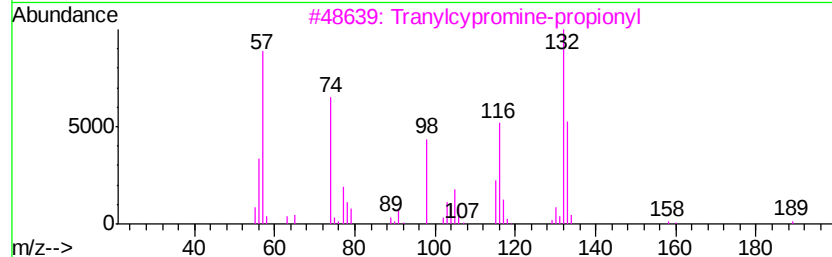
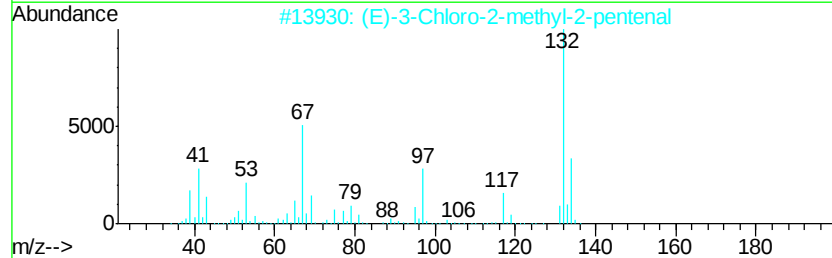
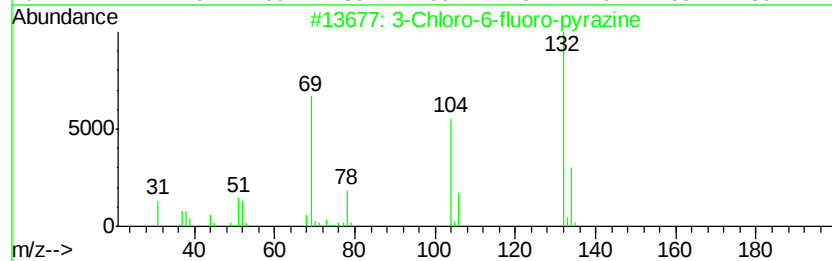
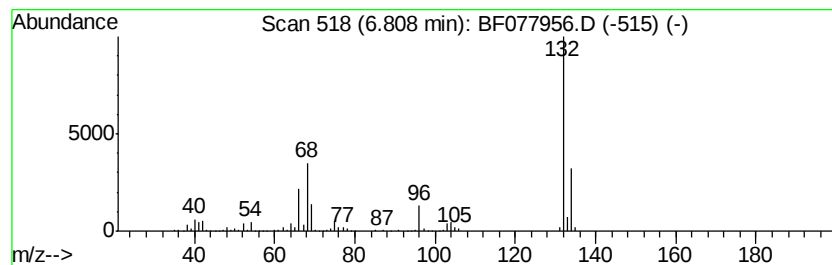
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown6.81 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.81	75.44 ng	781837	1,4-Dichlorobenzene-d4	7.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032515\
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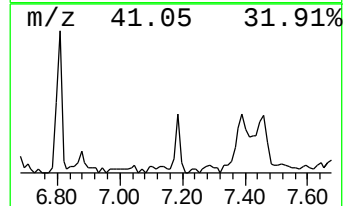
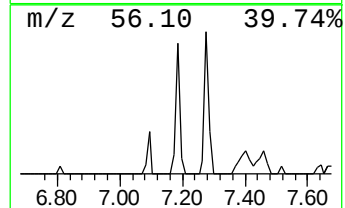
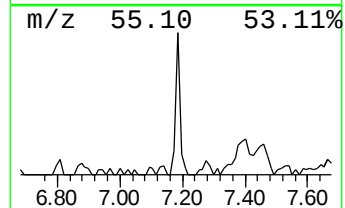
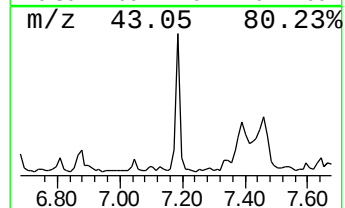
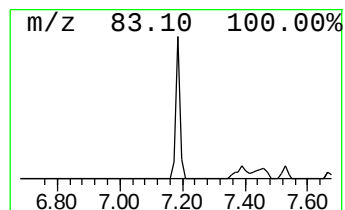
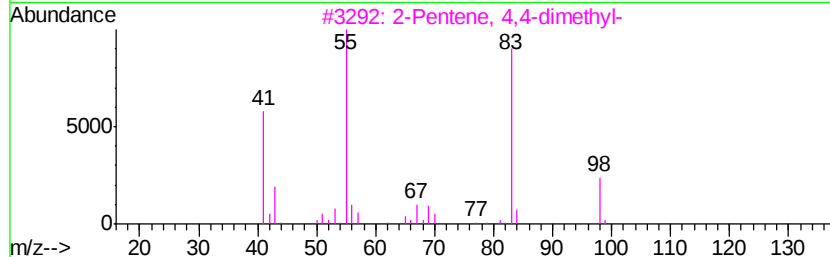
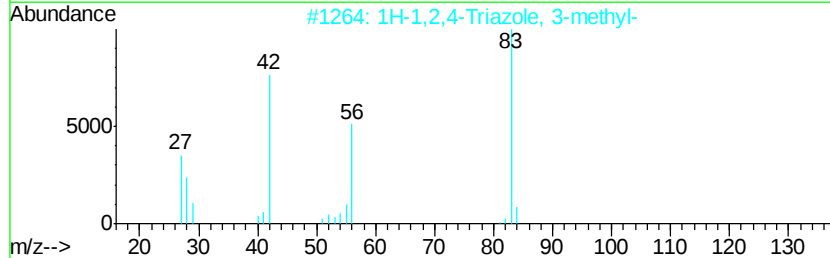
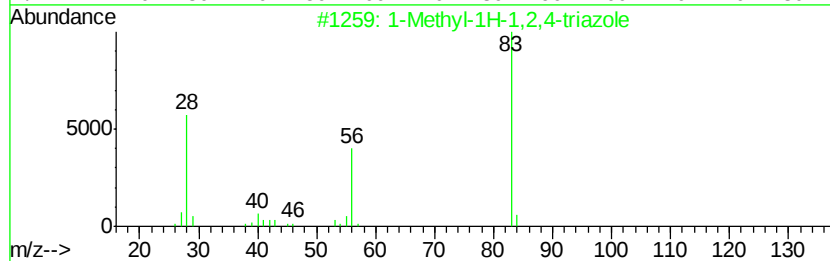
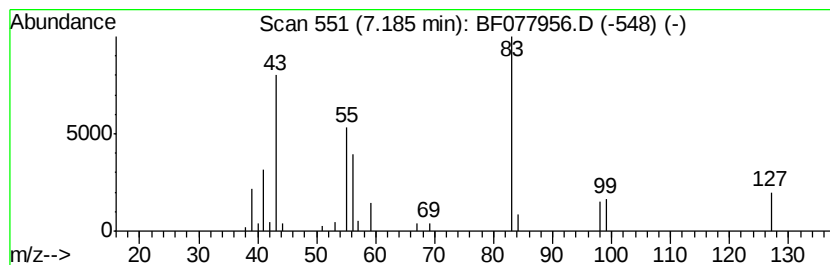
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown7.18 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.18	6.44 ng	66780	1,4-Dichlorobenzene-d4	7.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Methyl-1H-1,2,4-triazole	83	C3H5N3	006086-21-1	49
2		1H-1,2,4-Triazole, 3-methyl-	83	C3H5N3	007170-01-6	47
3		2-Pentene, 4,4-dimethyl-	98	C7H14	026232-98-4	38
4		2-Pentene, 3,4-dimethyl-	98	C7H14	024910-63-2	25
5		2-Pentene, 4,4-dimethyl-, (Z)-	98	C7H14	000762-63-0	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032515\
Data File : BF077956.D
Acq On : 25 Mar 2015 17:30
Operator : TP/IZ
Sample : G1612-01
Misc :
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ClientSampleId :
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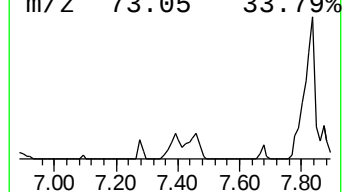
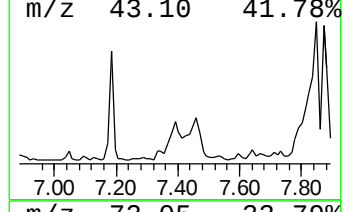
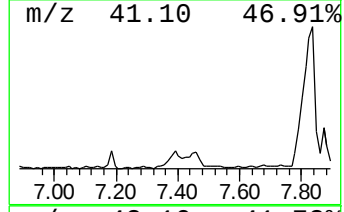
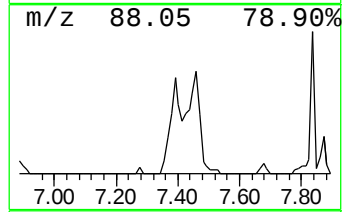
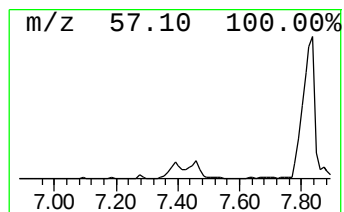
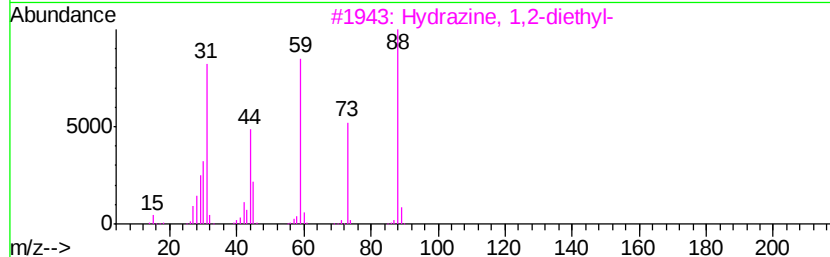
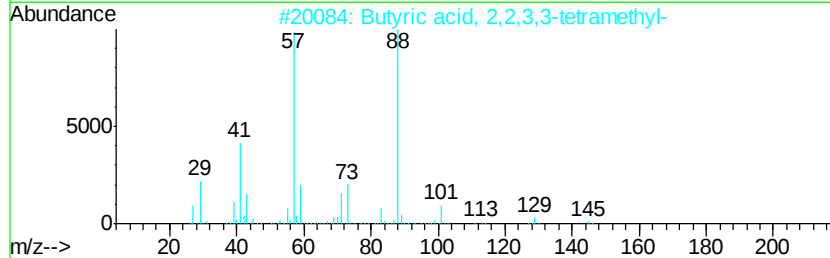
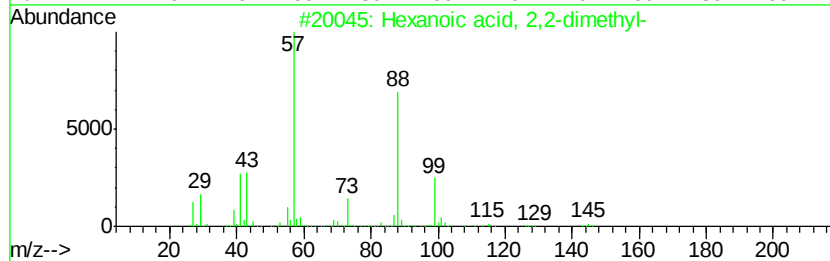
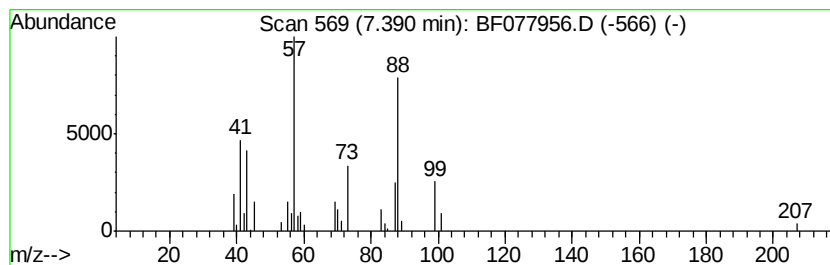
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hexanoic acid, 2,2-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.39	14.86 ng	153983	1,4-Dichlorobenzene-d4	7.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanoic acid, 2,2-dimethyl-	144	C8H16O2	000813-72-9	64
2		Butyric acid, 2,2,3,3-tetramethyl-	144	C8H16O2	030407-41-1	42
3		Hydrazine, 1,2-diethyl-	88	C4H12N2	001615-80-1	37
4		Methyl propionate	88	C4H8O2	000554-12-1	35
5		Decanoic acid, ethyl ester	200	C12H24O2	000110-38-3	32



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032515\
Data File : BF077956.D
Acq On : 25 Mar 2015 17:30
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Misc :
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Instrument :
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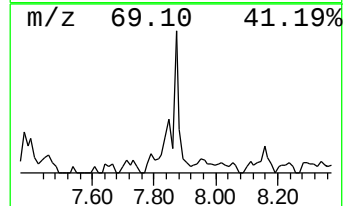
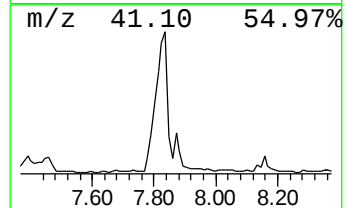
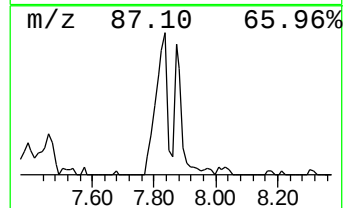
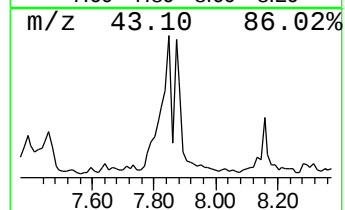
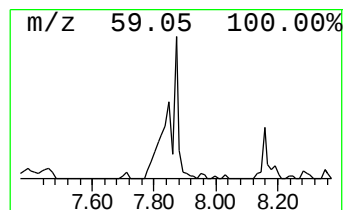
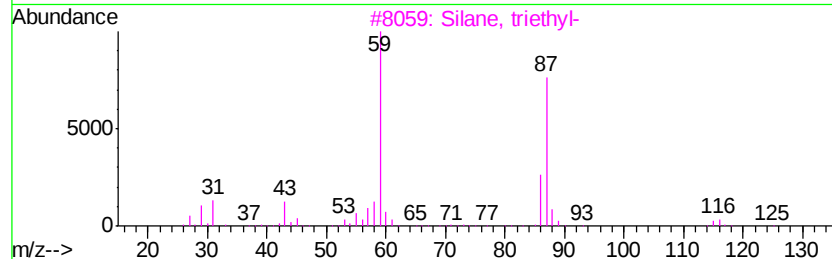
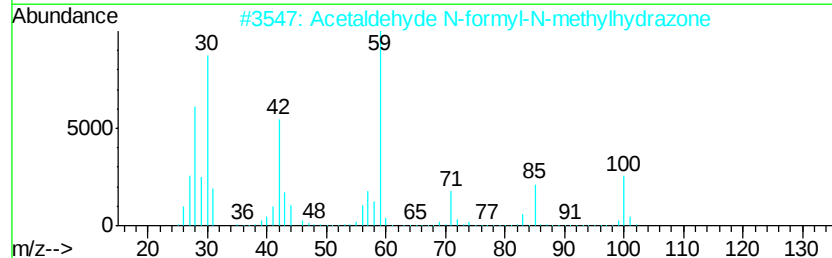
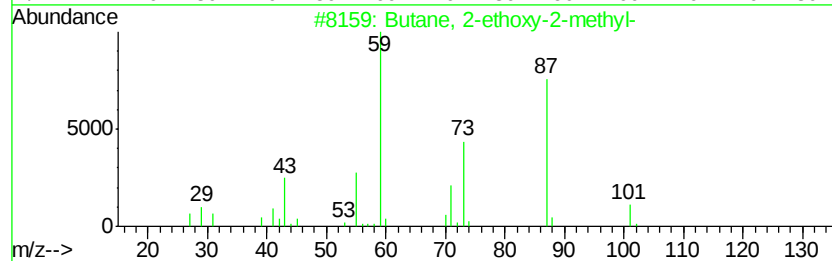
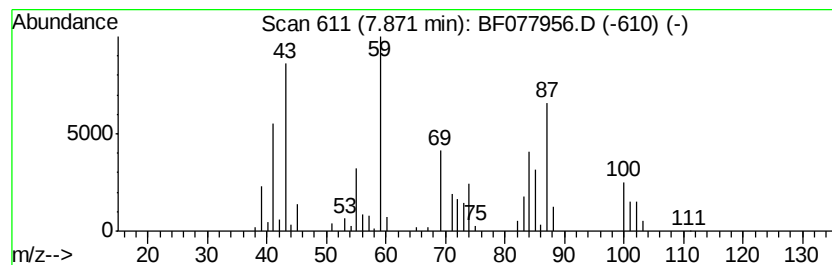
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown7.87 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.87	17.95 ng	186014	1,4-Dichlorobenzene-d4	7.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-ethoxy-2-methyl-	116	C7H16O	000919-94-8	32
2		Acetaldehyde N-formyl-N-methylhy...	100	C4H8N2O	016568-02-8	25
3		Silane, triethyl-	116	C6H16Si	000617-86-7	23
4		Propane, 1-ethoxy-	88	C5H12O	000628-32-0	22
5		Propanamide, 2-methoxy-N-methyl-	117	C5H11NO2	1000196-12-6	17



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032515\
Data File : BF077956.D
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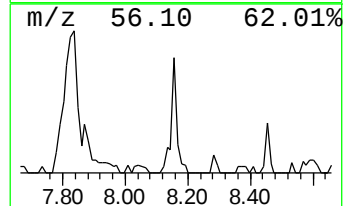
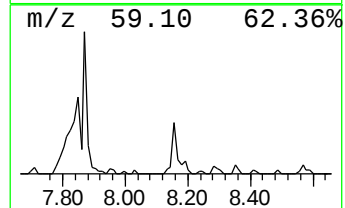
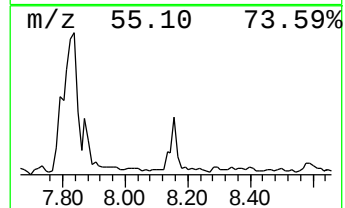
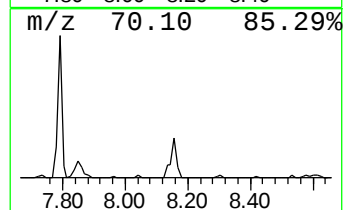
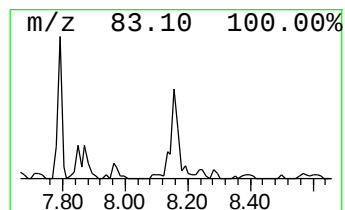
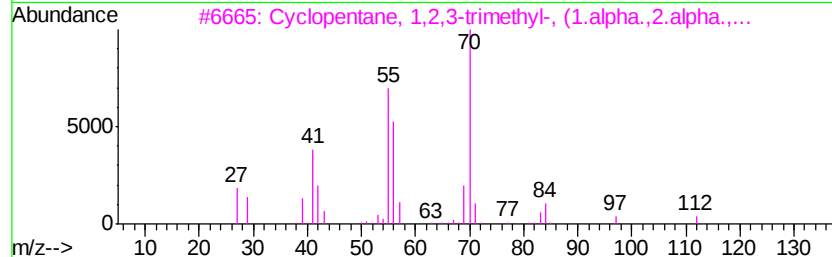
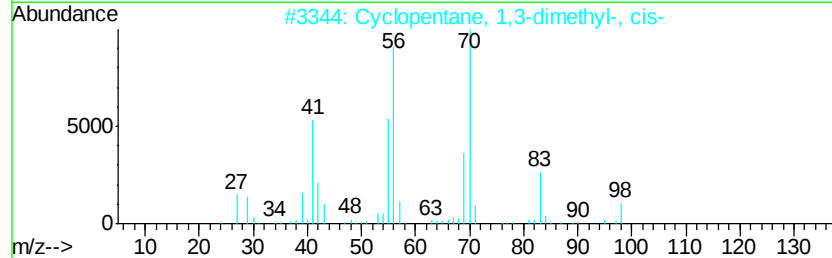
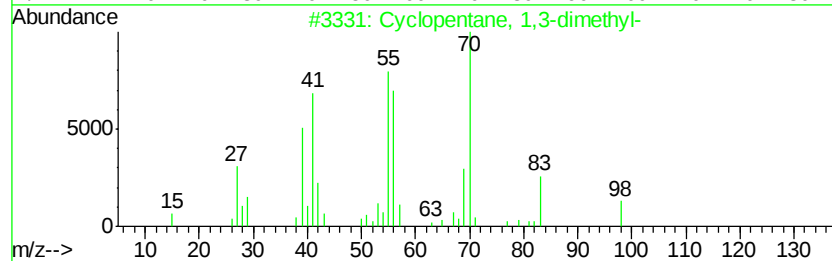
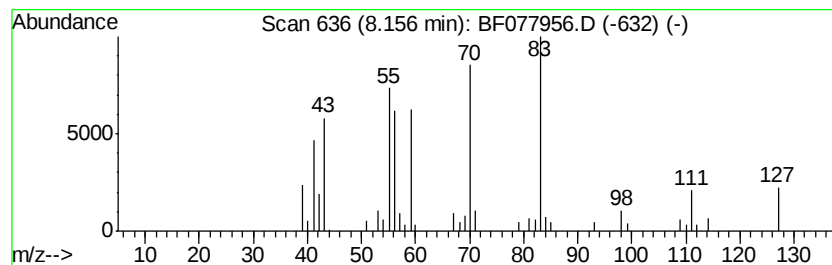
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Cyclopentane, 1,3-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.16	7.13 ng	112655	Naphthalene-d8	8.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	55
2		Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	43
3		Cyclopentane, 1,2,3-trimethyl-, ...	112	C8H16	002613-69-6	43
4		2-Pentene, 4,4-dimethyl-	98	C7H14	026232-98-4	25
5		Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	25



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032515\
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ClientSampleId :
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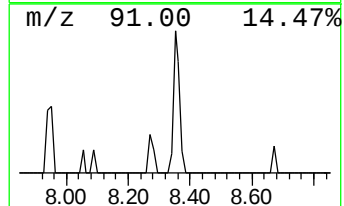
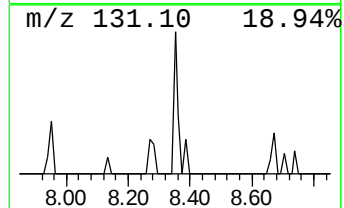
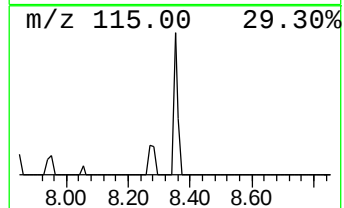
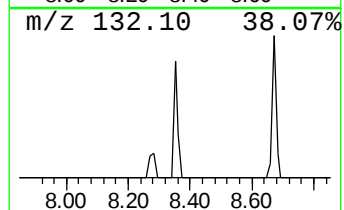
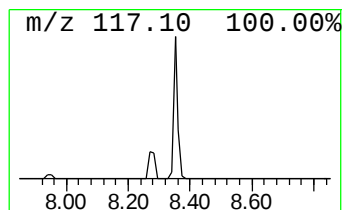
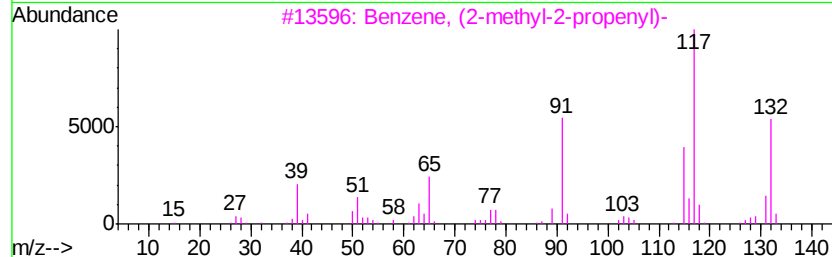
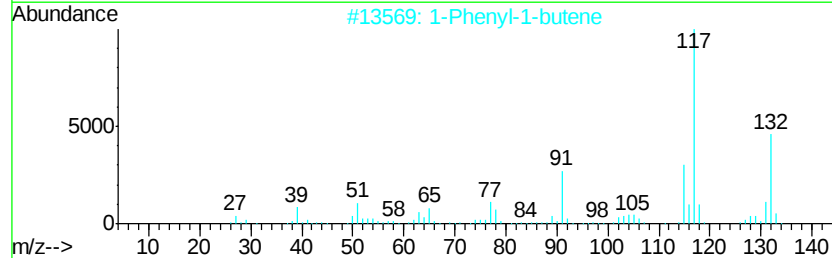
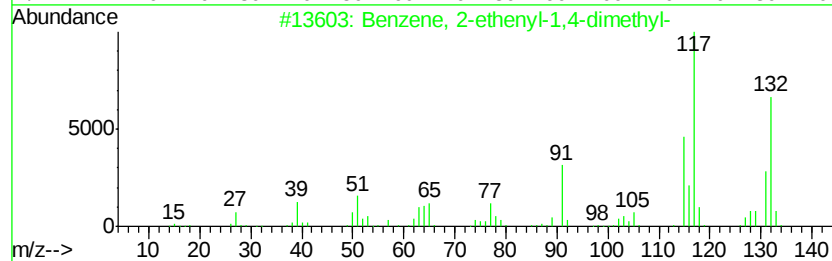
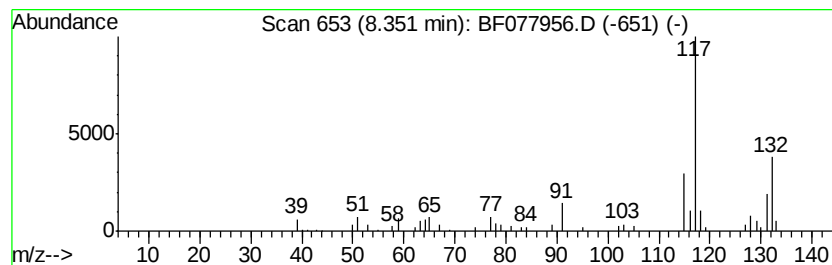
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.35	4.41 ng	69723	Naphthalene-d8	8.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2		1-Phenyl-1-butene	132	C10H12	000824-90-8	93
3		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	91
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	91
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032515\
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Sample : G1612-01
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ALS Vial : 5 Sample Multiplier: 1

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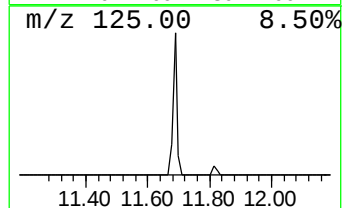
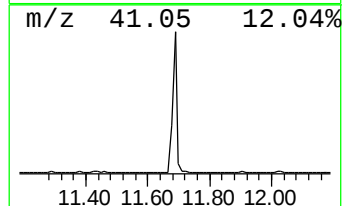
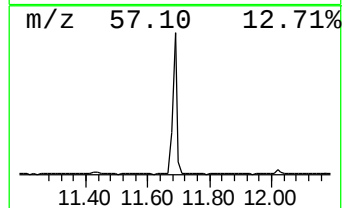
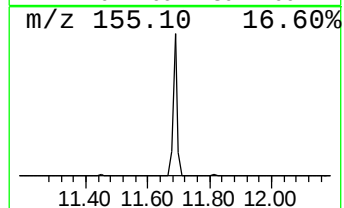
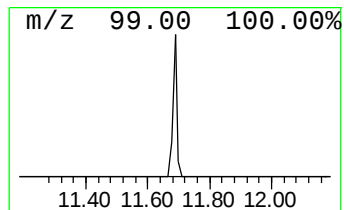
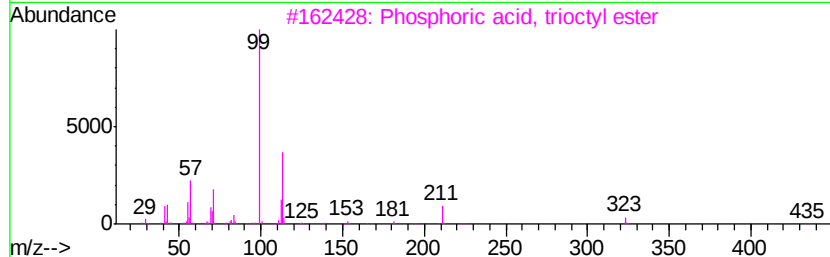
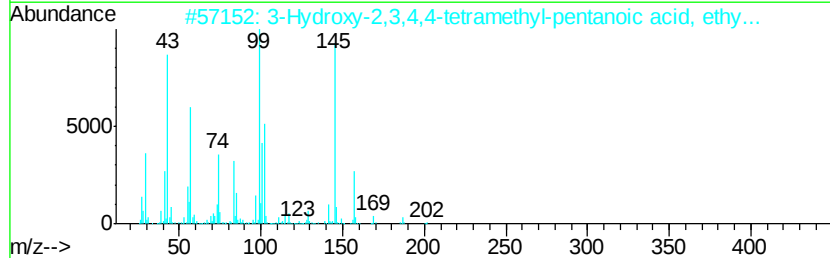
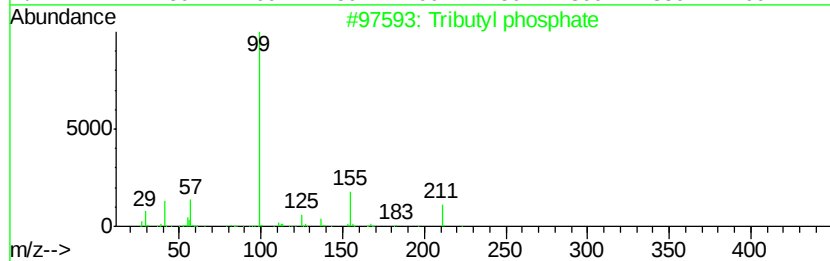
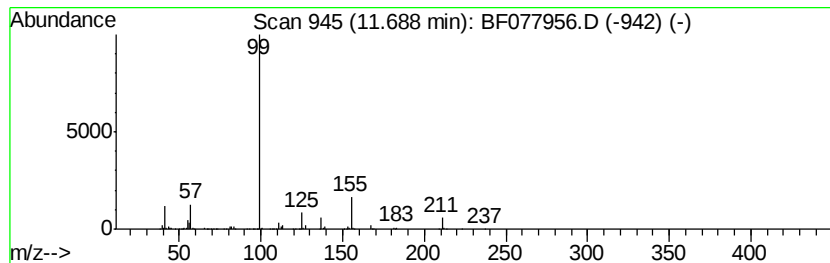
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Tributyl phosphate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.69	50.21 ng	891293	Acenaphthene-d10	10.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tributyl phosphate	266	C12H27O4P	000126-73-8	83
2		3-Hydroxy-2,3,4,4-tetramethyl-pe...	202	C11H22O3	1000194-64-5	28
3		Phosphoric acid, trioctyl ester	434	C24H51O4P	001806-54-8	9
4		Spiro[1,3-dioxolane-2,2'(1'H)-na...	196	C12H20O2	006857-86-9	9
5		n-Heptyl methylphosphonofluoridate	196	C8H18FO2P	162085-82-7	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032515\
Data File : BF077956.D
Acq On : 25 Mar 2015 17:30
Operator : TP/IZ
Sample : G1612-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-1

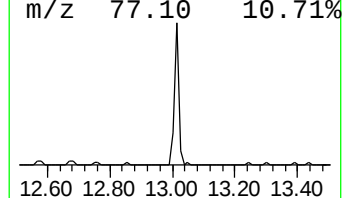
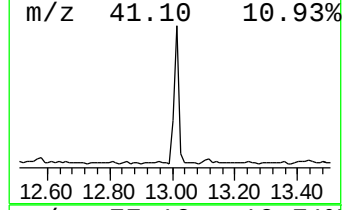
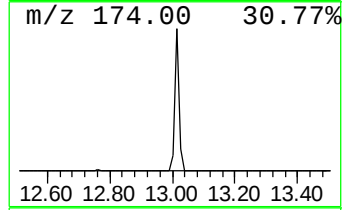
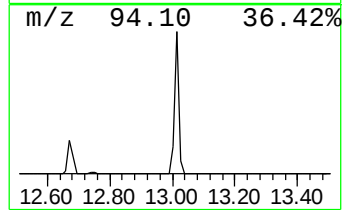
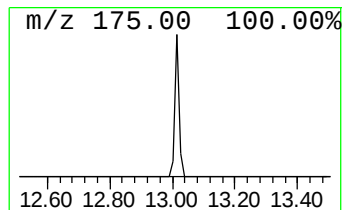
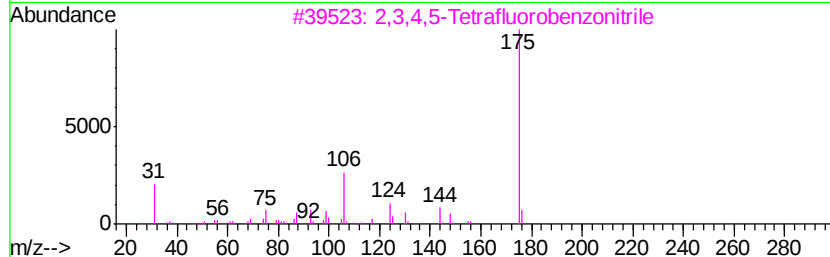
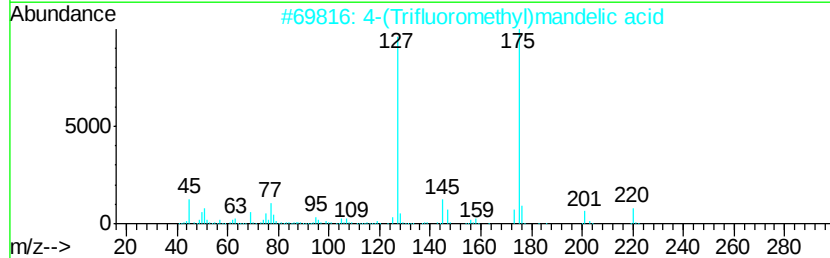
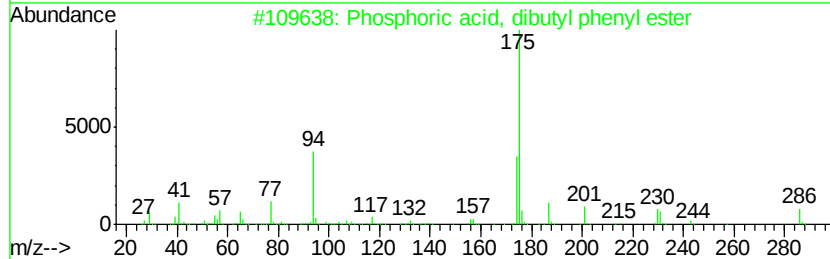
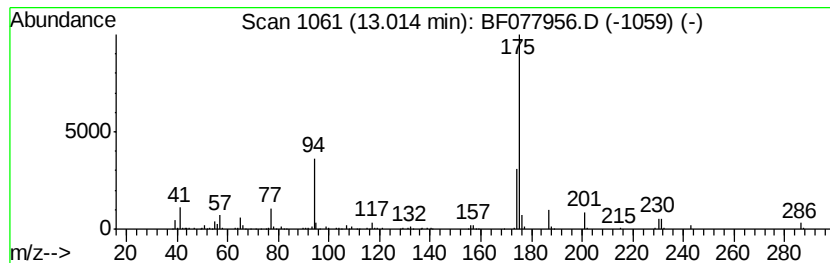
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Phosphoric acid, dibutyl ph... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.01	23.52 ng	423857	Phenanthrene-d10	12.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phosphoric acid, dibutyl phenyl ...	286	C14H23O4P	002528-36-1	95
2		4-(Trifluoromethyl)mandelic acid	220	C9H7F3O3	000395-35-7	43
3		2,3,4,5-Tetrafluorobenzonitrile	175	C7HF4N	016582-93-7	9
4		Benzene, 4-chloro-2-fluoro-1-nitro-	175	C6H3ClFN02	000700-37-8	9
5		3-Hydroxy-4-phenyl-5-butyl-1,2,4...	217	C12H15N3O	066921-14-0	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032515\
Data File : BF077956.D
Acq On : 25 Mar 2015 17:30
Operator : TP/IZ
Sample : G1612-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-1

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
Amylene Hydrate	1.15	111.0	ng	1150590	1	7.09	207275	20.0
Butane, 2-methoxy...	1.52	24.5	ng	254308	1	7.09	207275	20.0
2-Butanol, 2,3-di...	2.41	13.9	ng	143681	1	7.09	207275	20.0
2-Pentanol, 2,4-d...	3.87	8.9	ng	92455	1	7.09	207275	20.0
2-Pentanone, 4-hy...	4.83	4.6	ng	47567	1	7.09	207275	20.0
unknown6.81	6.81	75.4	ng	781837	1	7.09	207275	20.0
unknown7.18	7.18	6.4	ng	66780	1	7.09	207275	20.0
Hexanoic acid, 2,...	7.39	14.9	ng	153983	1	7.09	207275	20.0
unknown7.87	7.87	17.9	ng	186014	1	7.09	207275	20.0
Cyclopentane, 1,3...	8.16	7.1	ng	112655	2	8.67	316175	20.0
Benzene, 2-etheny...	8.35	4.4	ng	69723	2	8.67	316175	20.0
Tributyl phosphate	11.69	50.2	ng	891293	3	10.84	355027	20.0
Phosphoric acid, ...	13.01	23.5	ng	423857	4	12.67	360425	20.0