

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033116\
 Data File : BF085924.D
 Acq On : 31 Mar 2016 13:39
 Operator : UM/SJ
 Sample : H1954-01
 Misc :
 ALS Vial : 3 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-BF032416.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	2.401	4	10	13	rBV	1746713	5118407	100.00%	14.580%
2	2.847	46	49	53	rBV	134985	241837	4.72%	0.689%
3	4.196	164	167	170	rBV	100530	162848	3.18%	0.464%
4	4.904	228	229	233	rVV	73961	101369	1.98%	0.289%
5	4.961	233	234	247	rVB	29168	60959	1.19%	0.174%
6	5.384	268	271	273	rBV	1232036	1477934	28.87%	4.210%
7	5.613	289	291	292	rBV	55374	61056	1.19%	0.174%
8	5.853	311	312	315	rVB	72366	79639	1.56%	0.227%
9	6.424	360	362	364	rVV	1085474	1031949	20.16%	2.940%
10	6.470	365	366	368	rVV	122308	94684	1.85%	0.270%
11	6.550	371	373	376	rBV	1982708	2497744	48.80%	7.115%
12	6.790	392	394	396	rBV	700715	700282	13.68%	1.995%
13	6.859	398	400	402	rVB	79059	110212	2.15%	0.314%
14	6.939	405	407	410	rBV	1780194	2308303	45.10%	6.575%
15	7.042	412	416	421	rVB2	161669	432416	8.45%	1.232%
16	7.362	440	444	446	rVV	2013742	2770282	54.12%	7.891%
17	7.407	446	448	452	rVV	215404	331701	6.48%	0.945%
18	7.476	452	454	458	rVB2	54260	93067	1.82%	0.265%
19	7.613	463	466	467	rBV	50046	77375	1.51%	0.220%
20	7.647	467	469	472	rVV	129692	161734	3.16%	0.461%
21	7.807	481	483	488	rBV2	101408	160157	3.13%	0.456%
22	7.990	496	499	504	rVB4	39265	79533	1.55%	0.227%
23	8.070	504	506	509	rBV	914653	949801	18.56%	2.706%
24	8.608	551	553	556	rVB	94591	96991	1.89%	0.276%
25	8.790	568	569	574	rVB	40377	63650	1.24%	0.181%
26	9.156	598	601	603	rVB	3582346	3652979	71.37%	10.406%
27	9.350	615	618	622	rBV	319910	377755	7.38%	1.076%
28	9.488	629	630	634	rBV3	52553	77501	1.51%	0.221%
29	9.545	634	635	637	rVB	144539	148415	2.90%	0.423%
30	9.648	642	644	648	rBV3	36890	58360	1.14%	0.166%
31	9.831	658	660	662	rVB	1151636	1072454	20.95%	3.055%
32	10.036	676	678	681	rBV2	41013	60075	1.17%	0.171%
33	10.173	688	690	694	rBV3	40741	52940	1.03%	0.151%
34	10.265	694	698	699	rVB2	71735	93034	1.82%	0.265%

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

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Peak separation: 5

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35	10.471	713	716	718	rBV	1023517	1261895	24.65%	3.595%
36	10.619	726	729	732	rBV	2098128	2122452	41.47%	6.046%
37	10.733	737	739	743	rVB	83447	90094	1.76%	0.257%
38	11.316	787	790	792	rBV	874612	1053093	20.57%	3.000%
39	11.534	807	809	812	rBV	600661	539842	10.55%	1.538%
40	11.842	834	836	844	rVB3	93919	135954	2.66%	0.387%
41	12.631	902	905	910	rBV	62392	99749	1.95%	0.284%
42	12.894	925	928	931	rBV	2555949	3501002	68.40%	9.973%
43	13.945	1018	1020	1023	rBV	896569	831958	16.25%	2.370%
44	15.362	1141	1144	1156	rVB	480887	612081	11.96%	1.744%

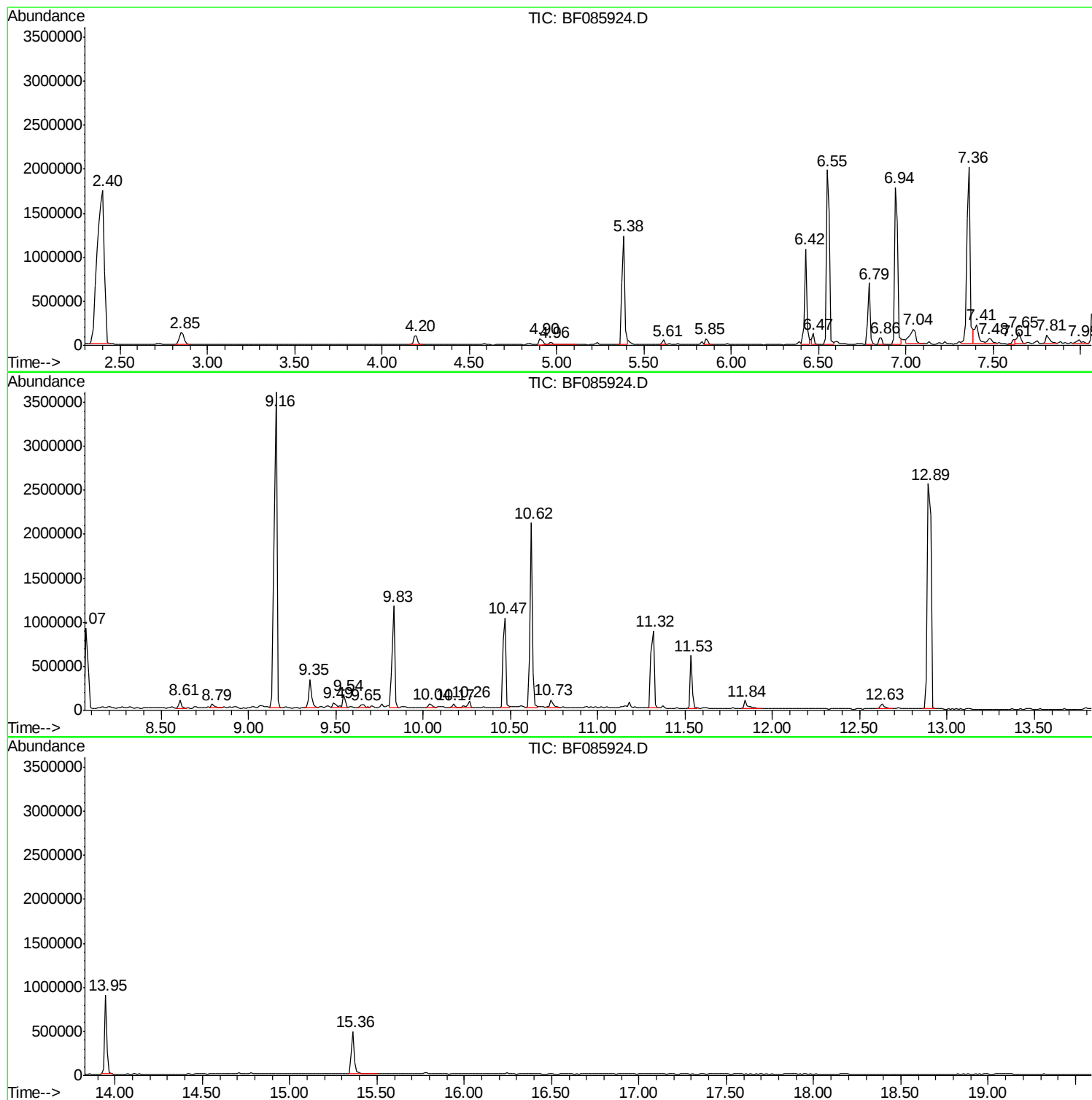
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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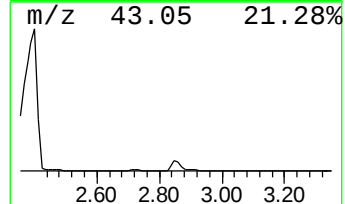
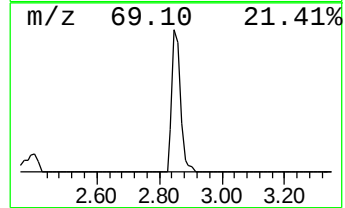
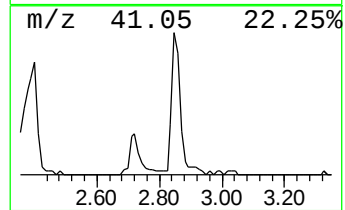
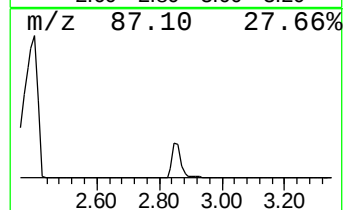
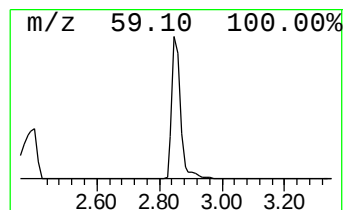
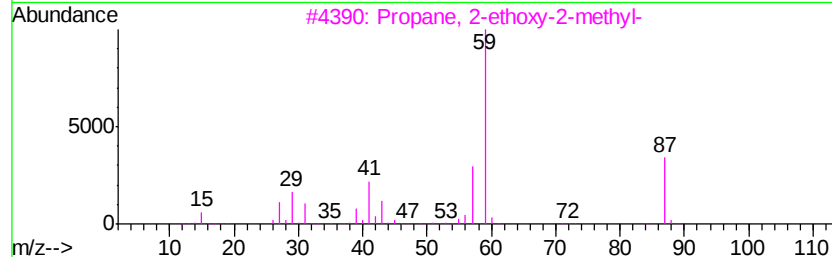
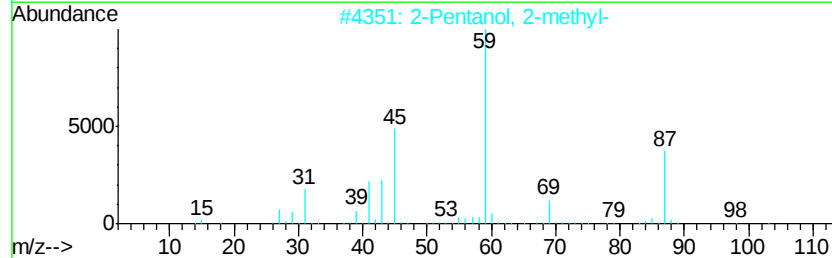
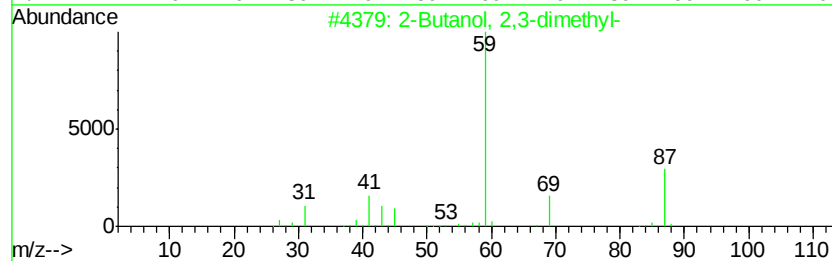
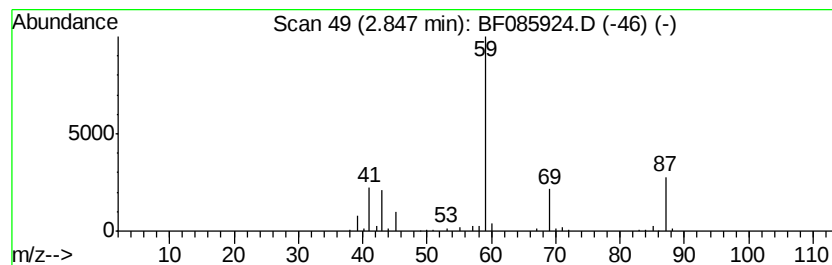
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Butanol, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.85	6.91 ng	241837	1,4-Dichlorobenzene-d4	6.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	83
2			2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	72
3			Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	64
4			Propane, 1-ethoxy-	88	C5H12O	000628-32-0	53
5			t-Butyl ethyl ether	102	C6H14O	1000118-18-4	50



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ClientSampleId :
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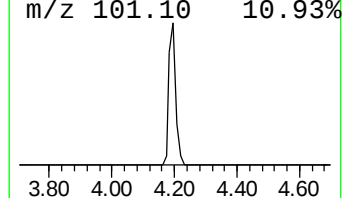
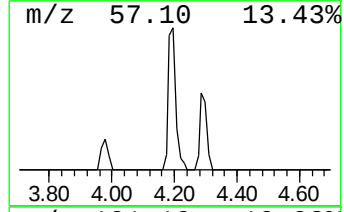
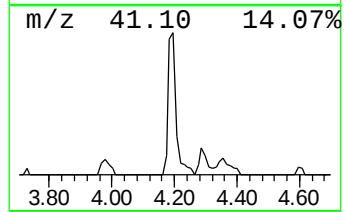
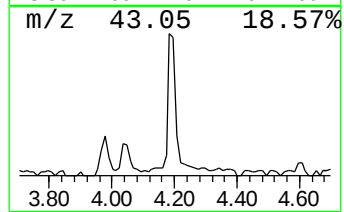
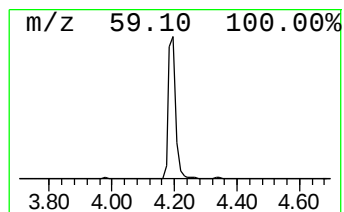
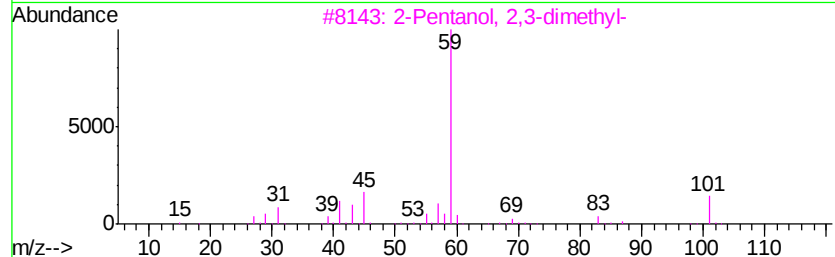
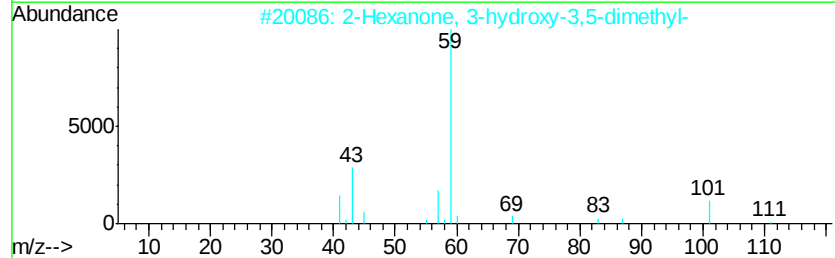
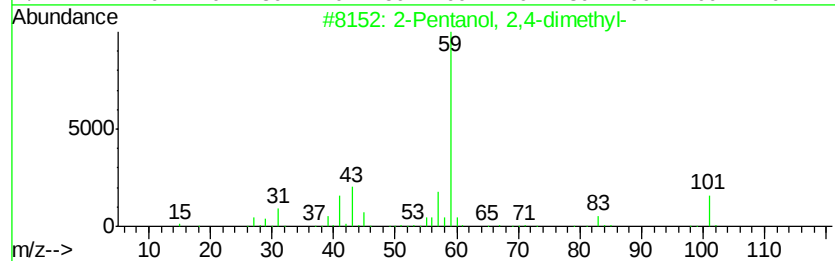
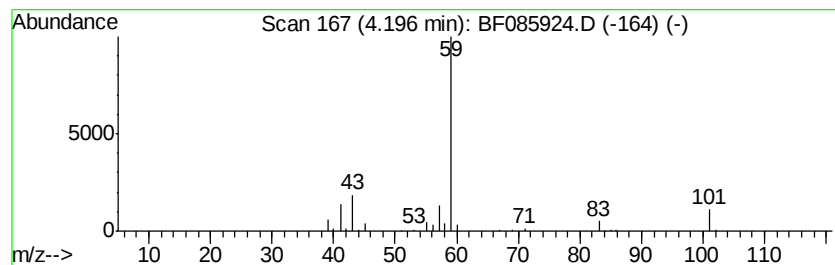
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 2-Pentanol, 2,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.20	4.65 ng	162848	1,4-Dichlorobenzene-d4	6.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	83
2			2-Hexanone, 3-hydroxy-3,5-dimethyl-	144	C8H16O2	006321-14-8	78
3			2-Pentanol, 2,3-dimethyl-	116	C7H16O	004911-70-0	64
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	56
5			3-Hexanol, 4-ethyl-	130	C8H18O	019780-44-0	40



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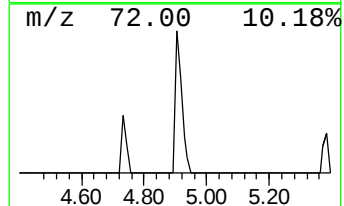
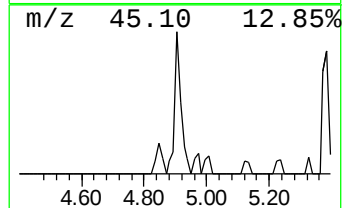
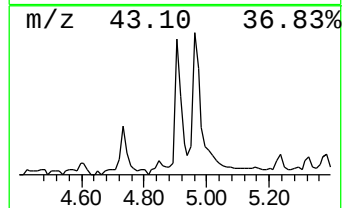
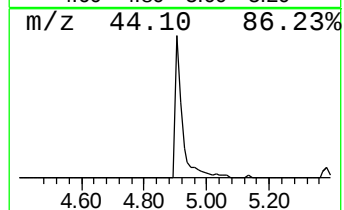
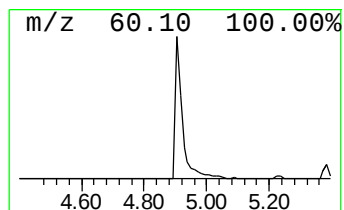
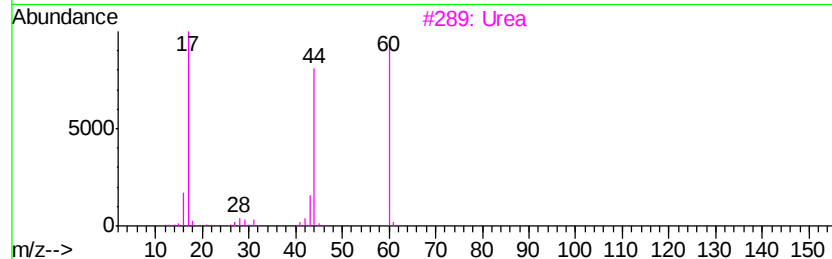
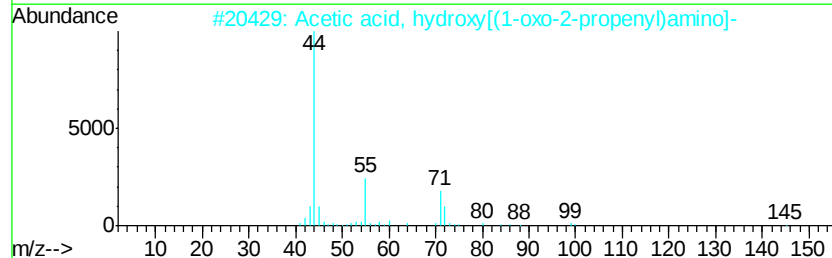
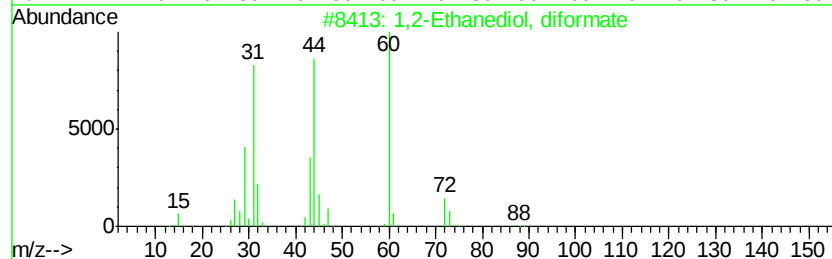
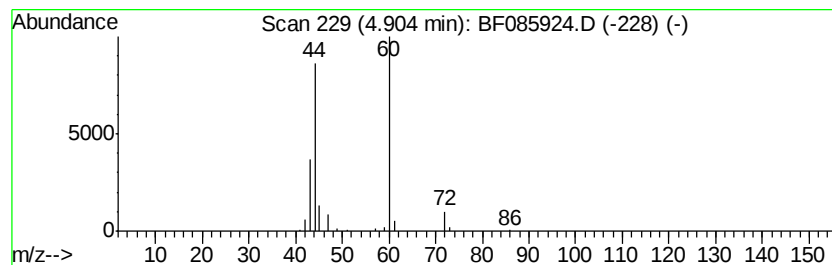
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 1,2-Ethanedio1, diformate Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.90	2.90 ng	101369	1,4-Dichlorobenzene-d4	6.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Ethanedio1, diformate	118	C4H6O4	000629-15-2	83
2			Acetic acid, hydroxy[(1-oxo-2-pr...	145	C5H7NO4	006737-24-2	25
3			Urea	60	CH4N2O	000057-13-6	9
4			Thiophene-3-ol, tetrahydro-, 1,1...	136	C4H8O3S	013031-76-0	9
5			Ethene, ethoxy-	72	C4H8O	000109-92-2	9



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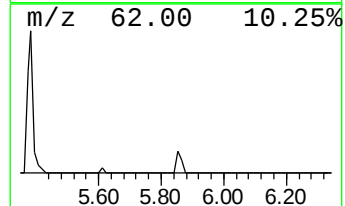
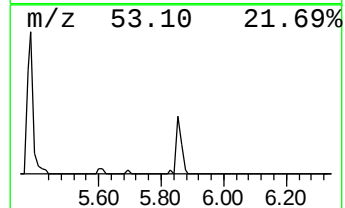
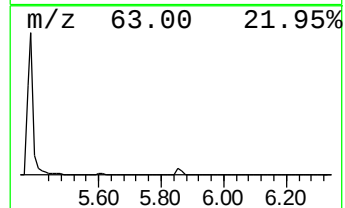
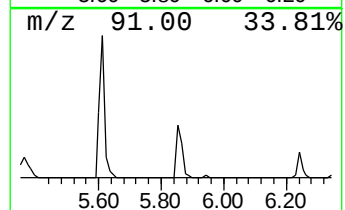
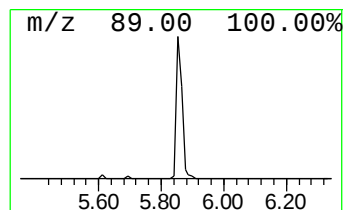
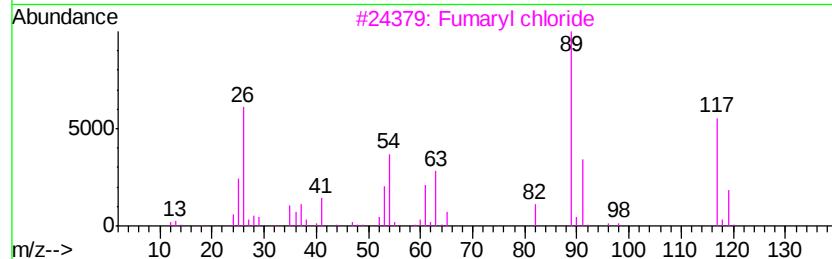
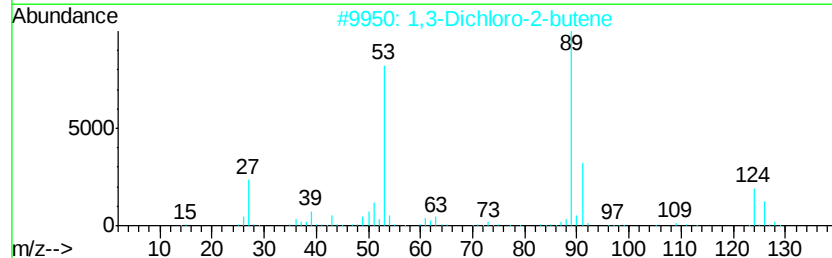
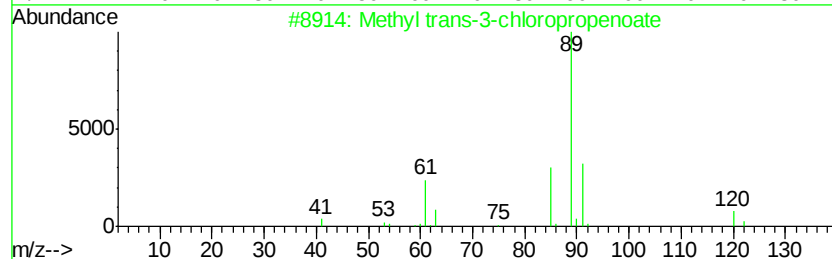
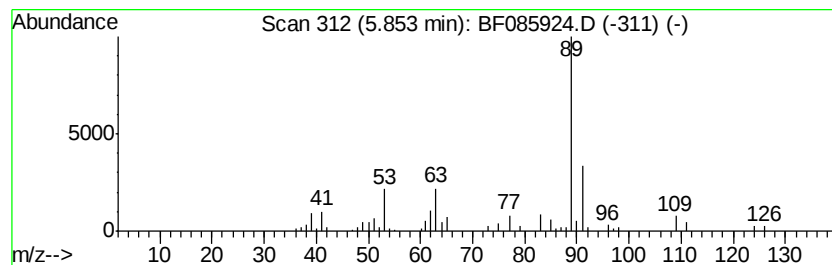
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 unknown5.85 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.85	2.27 ng	79639	1,4-Dichlorobenzene-d4	6.79

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Methyl trans-3-chloropropenoate	120	C4H5ClO2	1000144-78-8	38
2		1,3-Dichloro-2-butene	124	C4H6Cl2	000926-57-8	38
3		Fumaryl chloride	152	C4H2Cl2O2	000627-63-4	9
4		Methyl cis-3-chloropropenoate	120	C4H5ClO2	1000144-78-7	9
5		1,3,5-Trioxane	90	C3H6O3	000110-88-3	5



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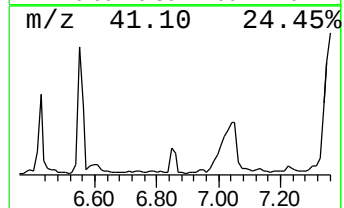
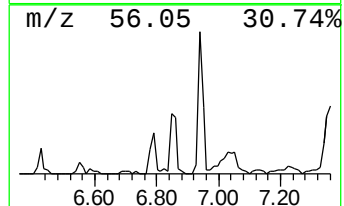
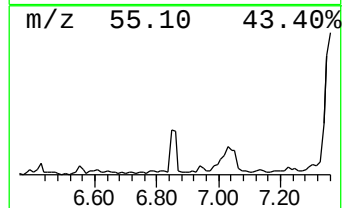
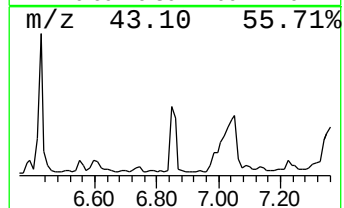
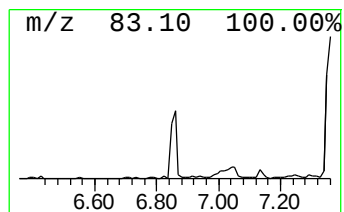
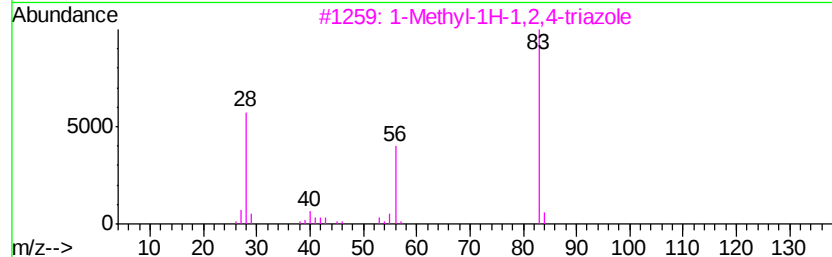
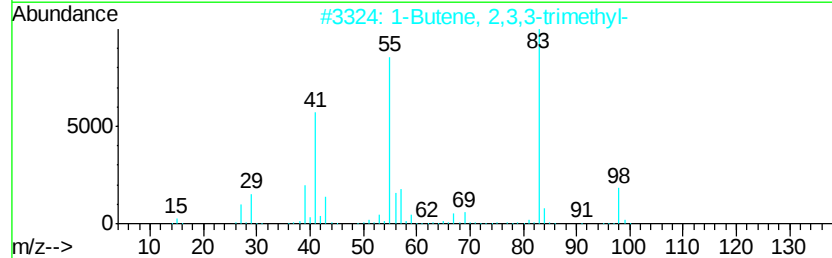
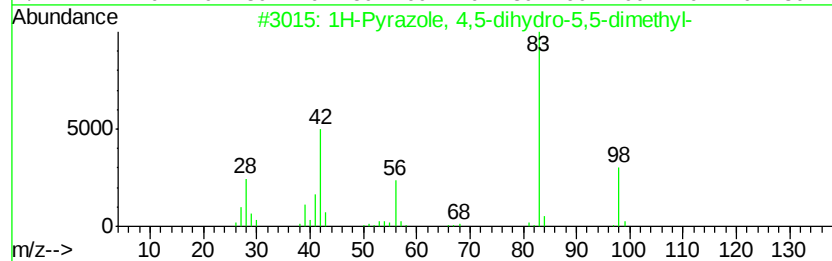
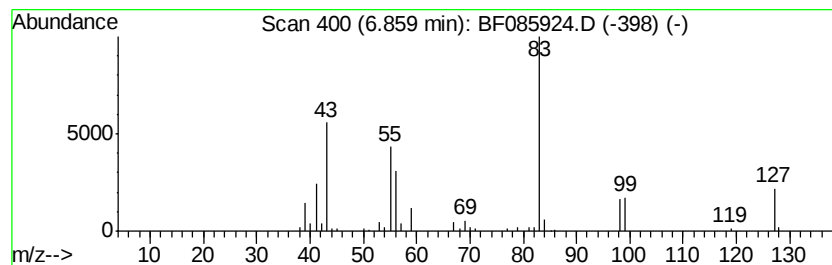
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TIC Integration Parameters: LSCINT.P

Peak Number 6 1H-Pyrazole, 4,5-dihydro-5,... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.86	3.15 ng	110212	1,4-Dichlorobenzene-d4	6.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Pyrazole, 4,5-dihydro-5,5-dim...	98	C5H10N2	004320-85-8	53
2	1-Butene, 2,3,3-trimethyl-	98	C7H14	000594-56-9	50
3	1-Methyl-1H-1,2,4-triazole	83	C3H5N3	006086-21-1	50
4	1H-Pyrazole, 4,5-dihydro-4,5-dim...	98	C5H10N2	028019-94-5	50
5	Cyclohexane, bromo-	162	C6H11Br	000108-85-0	43



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Data File : BF085924.D
Acq On : 31 Mar 2016 13:39
Operator : UM/SJ
Sample : H1954-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

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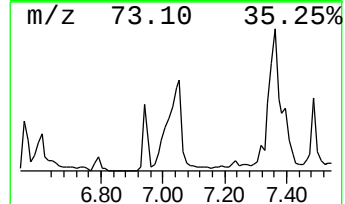
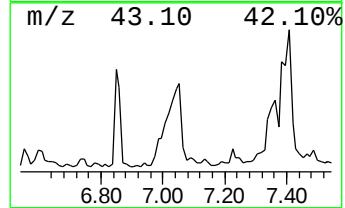
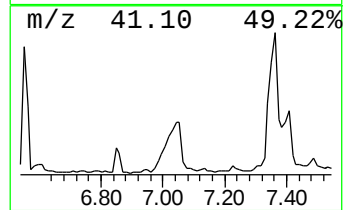
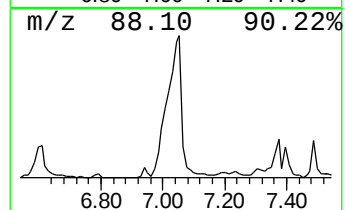
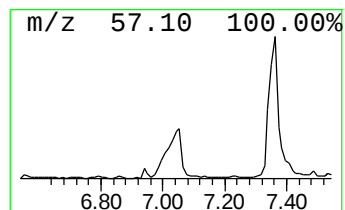
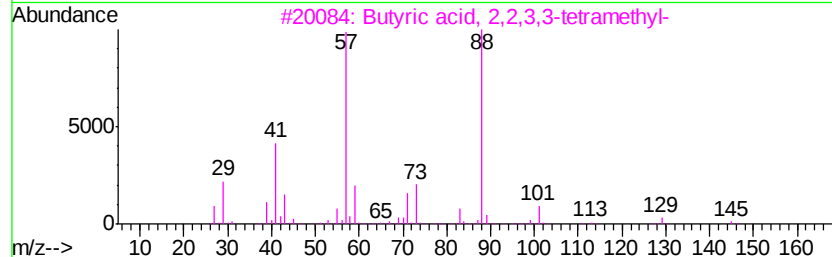
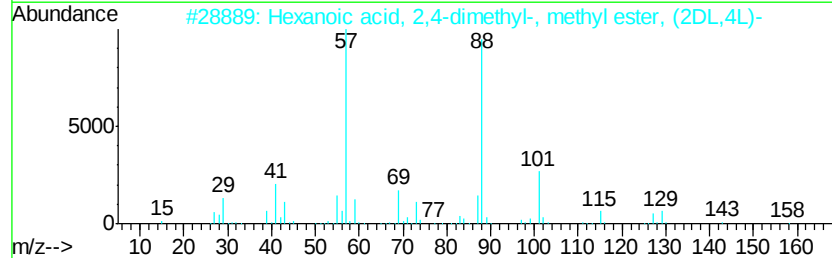
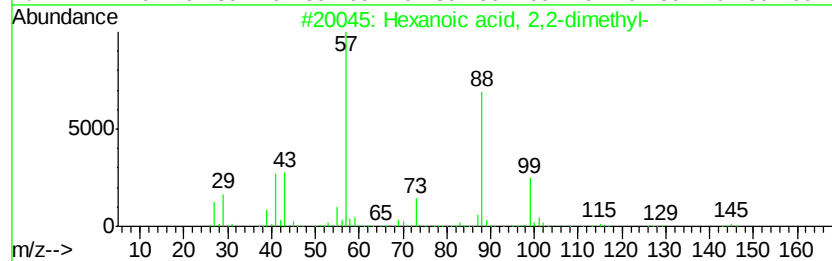
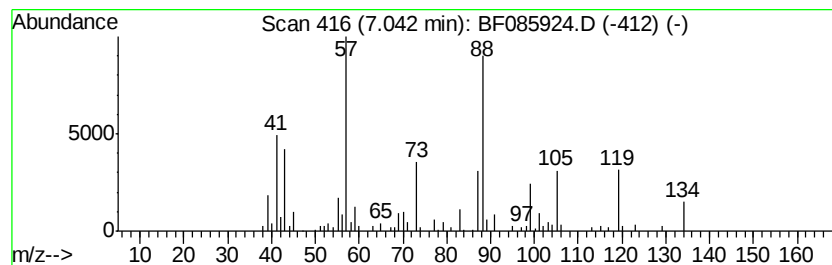
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexanoic acid, 2,2-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.04	12.35 ng	432416	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanoic acid, 2,2-dimethyl-	144	C8H16O2	000813-72-9	50
2		Hexanoic acid, 2,4-dimethyl-, me...	158	C9H18O2	014251-44-6	37
3		Butyric acid, 2,2,3,3-tetramethyl-	144	C8H16O2	030407-41-1	37
4		Butanoic acid, 2-ethyl-, 1,2,3-p...	386	C21H38O6	056554-54-2	32
5		Hydrazine, 1,2-diethyl-	88	C4H12N2	001615-80-1	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033116\
Data File : BF085924.D
Acq On : 31 Mar 2016 13:39
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Sample : H1954-01
Misc :
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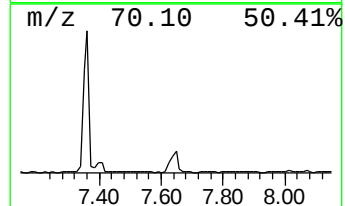
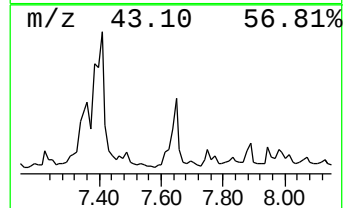
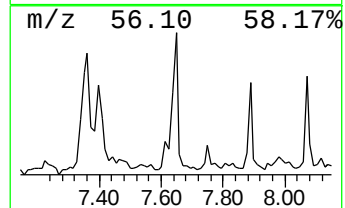
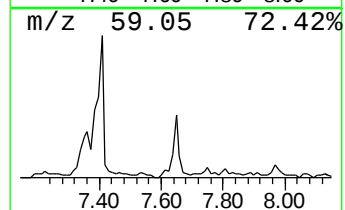
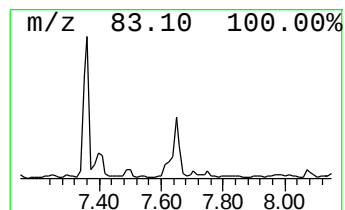
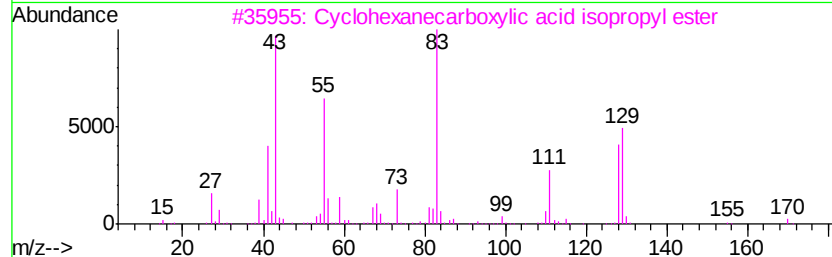
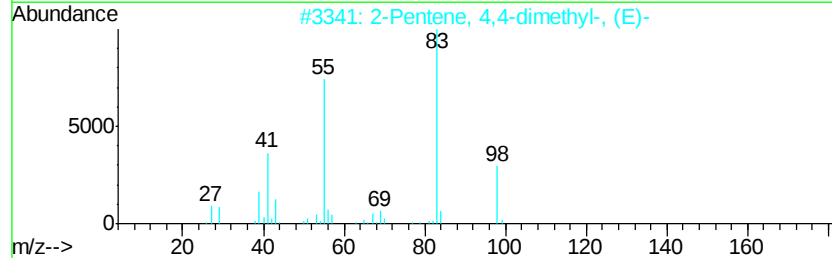
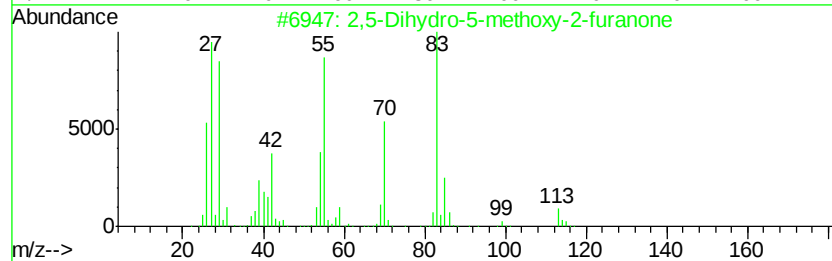
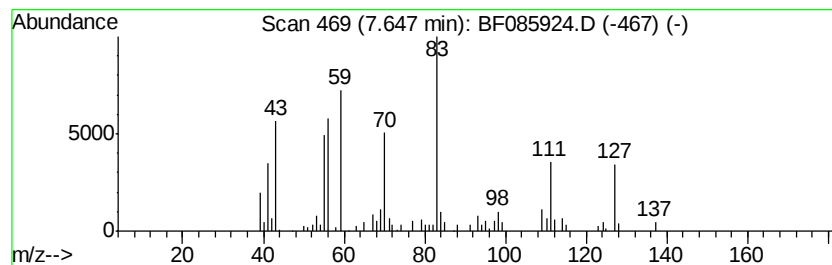
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown7.65 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.65	3.41 ng	161734	Napthalene-d8	8.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,5-Dihydro-5-methoxy-2-furanone	114	C5H6O3	010449-66-8	38
2		2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	27
3		Cyclohexanecarboxylic acid isopr...	170	C10H18O2	006553-80-6	27
4		Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	27
5		3-Amino-2-cyclohexenone	111	C6H9NO	005220-49-5	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033116\
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Misc :
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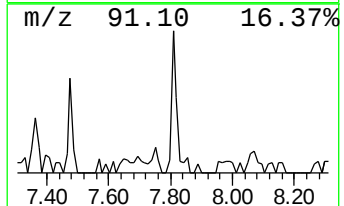
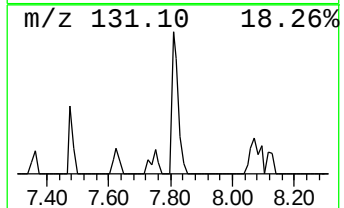
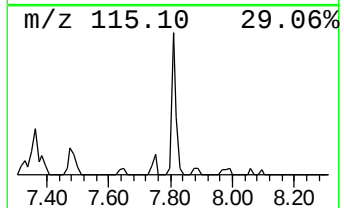
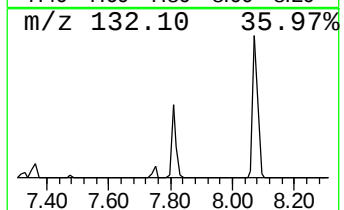
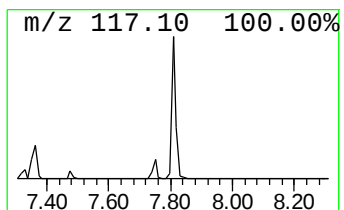
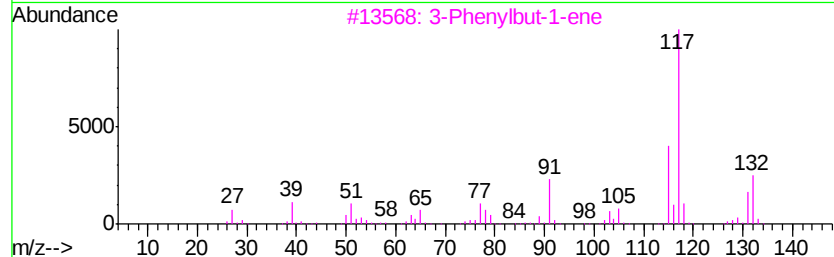
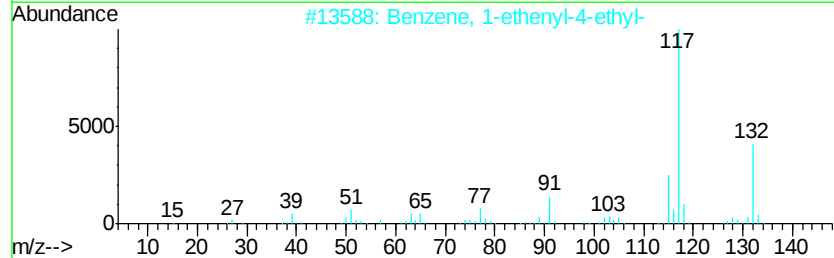
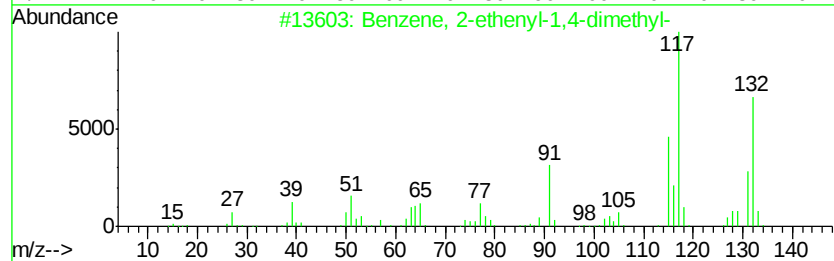
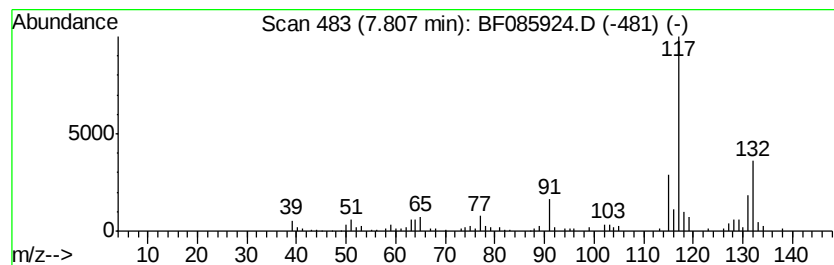
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.81	3.37 ng	160157	Naphthalene-d8	8.07

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
4			Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	91
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033116\
Data File : BF085924.D
Acq On : 31 Mar 2016 13:39
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Misc :
ALS Vial : 3 Sample Multiplier: 1

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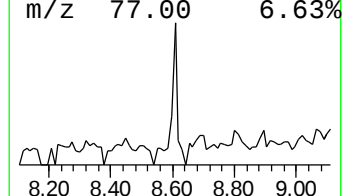
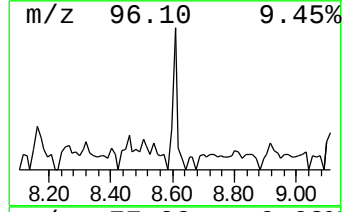
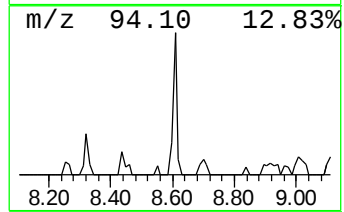
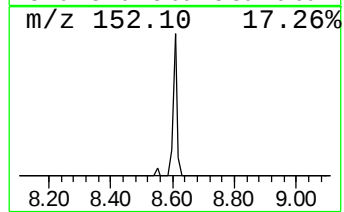
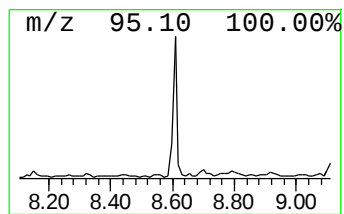
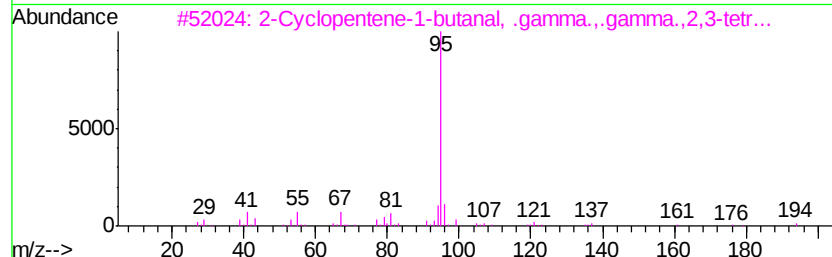
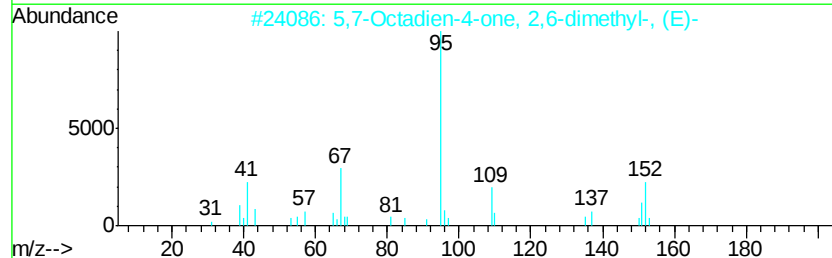
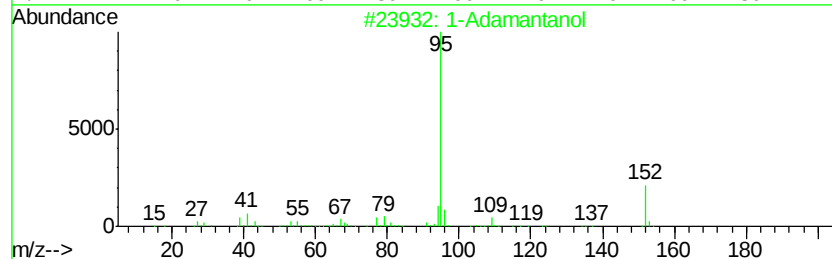
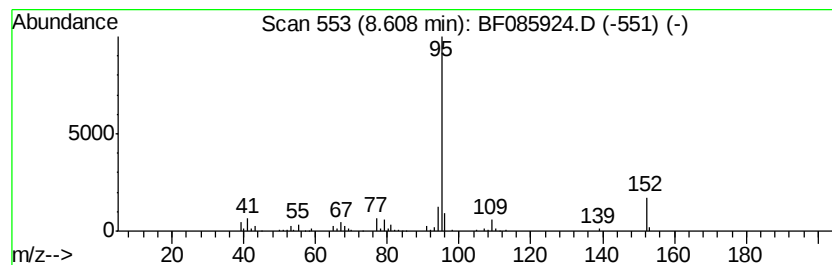
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1-Adamantanol Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.61	2.04 ng	96991	Naphthalene-d8	8.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Adamantanol	152	C10H16O	000768-95-6	90
2		5,7-Octadien-4-one, 2,6-dimethyl...	152	C10H16O	006752-80-3	59
3		2-Cyclopentene-1-butanal, .gamma...	194	C13H22O	060714-25-2	50
4		2-Furancarboxylic acid, 2-propen...	152	C8H8O3	004208-49-5	50
5		Cyclopentene, 1,4-dimethyl-5-(1-...	138	C10H18	061142-33-4	39



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033116\
Data File : BF085924.D
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Instrument :
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ClientSampleId :
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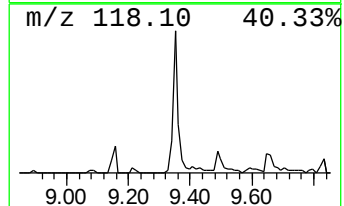
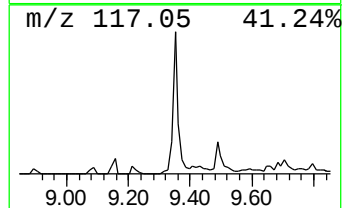
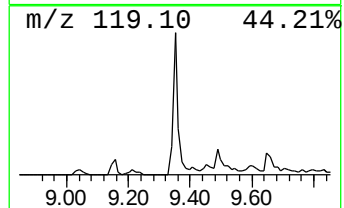
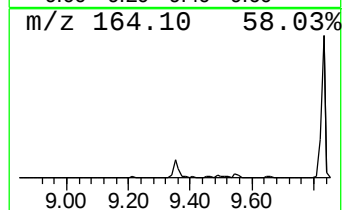
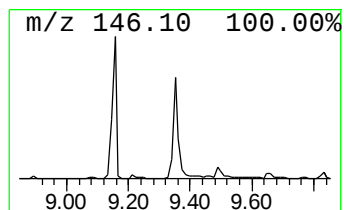
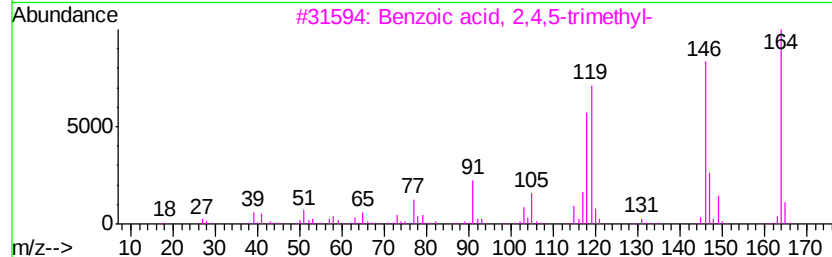
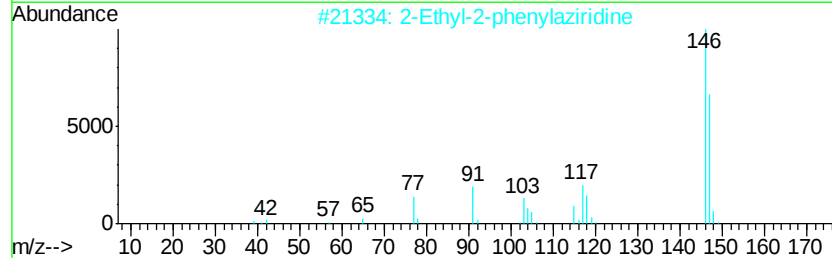
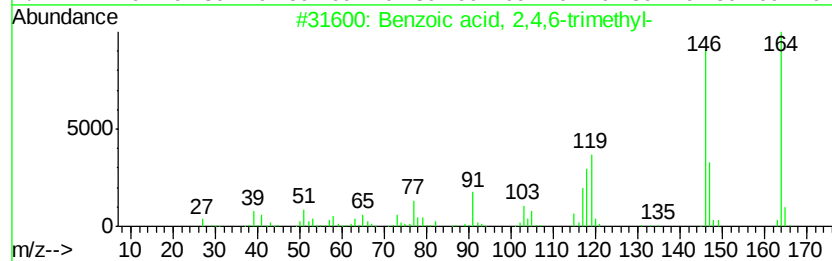
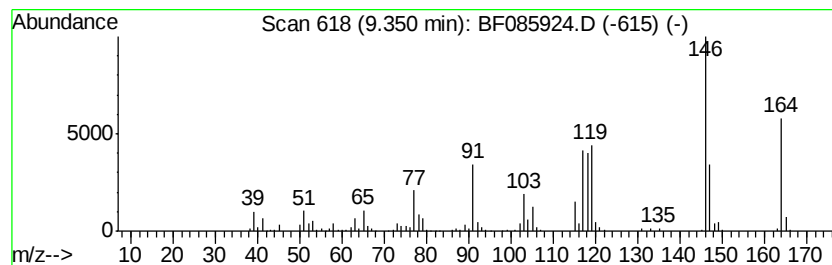
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzoic acid, 2,4,6-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.35	7.04 ng	377755	Acenaphthene-d10	9.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzoic acid, 2,4,6-trimethyl-	164	C10H12O2	000480-63-7	93
2		2-Ethyl-2-phenylaziridine	147	C10H13N	000768-82-1	50
3		Benzoic acid, 2,4,5-trimethyl-	164	C10H12O2	000528-90-5	49
4		Pteridine, 2-methyl-	146	C7H6N4	002432-20-4	27
5		4-Oxazolecarboxylic acid, 4,5-di...	233	C13H15N3O3	037592-54-4	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033116\
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Sample : H1954-01
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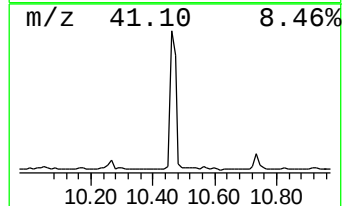
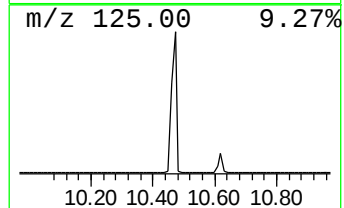
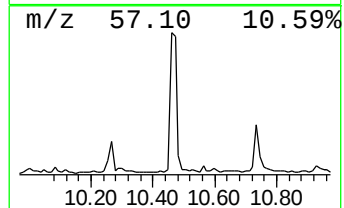
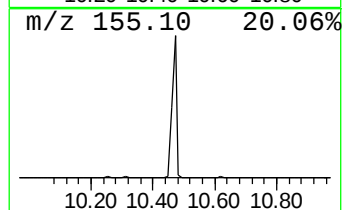
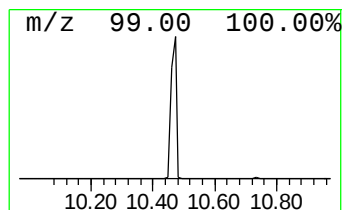
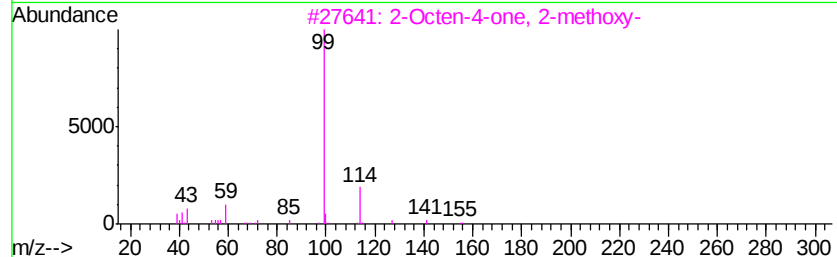
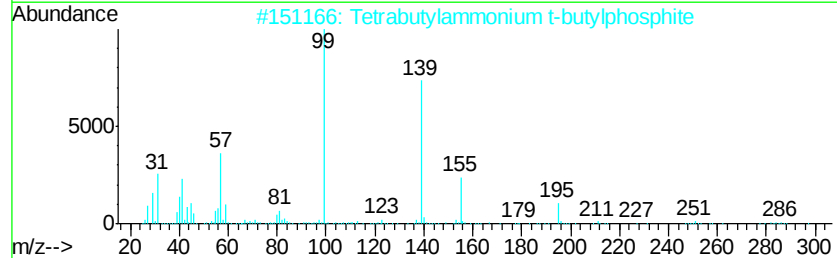
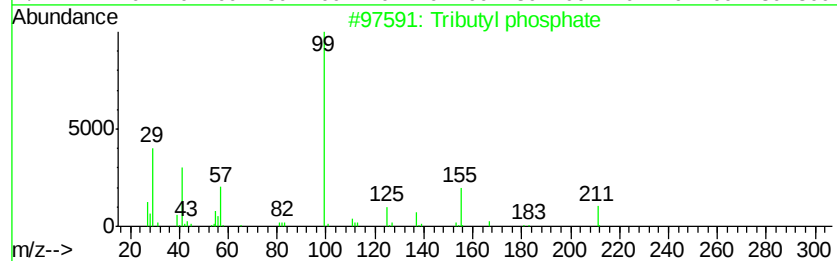
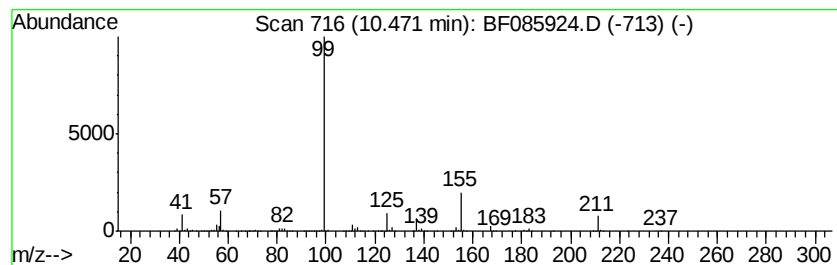
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tributyl phosphate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.47	23.53 ng	1261900	Acenaphthene-d10	9.83
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tributyl phosphate	266 C12H27O4P	000126-73-8 86
2		Tetrabutylammonium t-butylphosphite	379 C20H46N03P	1000164-81-7 38
3		2-Octen-4-one, 2-methoxy-	156 C9H16O2	024985-48-6 38
4		3-Hydroxy-2,3,4,4-tetramethyl-pe...	202 C11H22O3	1000194-64-5 25
5		Phosphoric acid, trioctyl ester	434 C24H51O4P	001806-54-8 12



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033116\
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Sample : H1954-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

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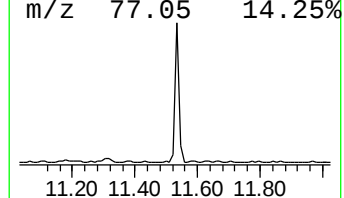
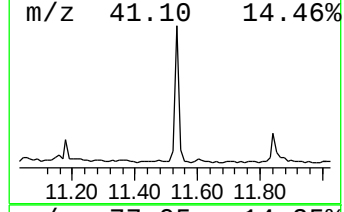
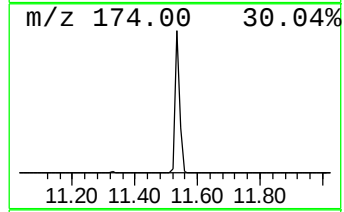
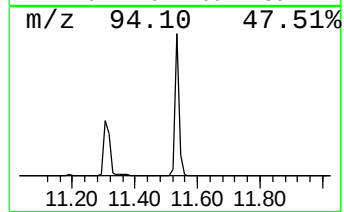
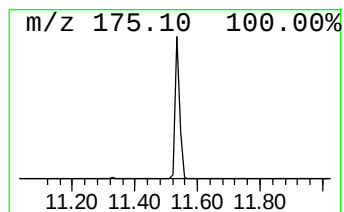
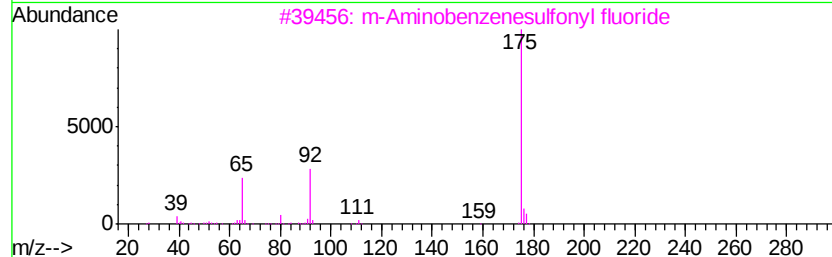
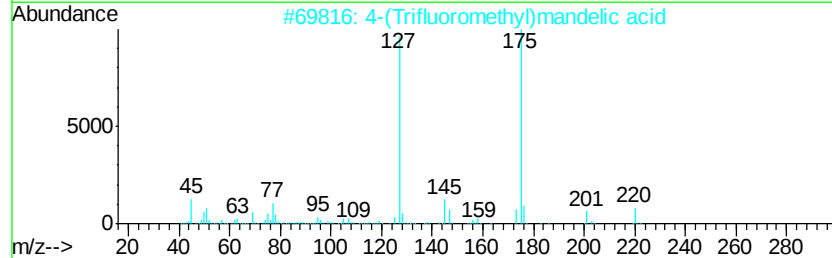
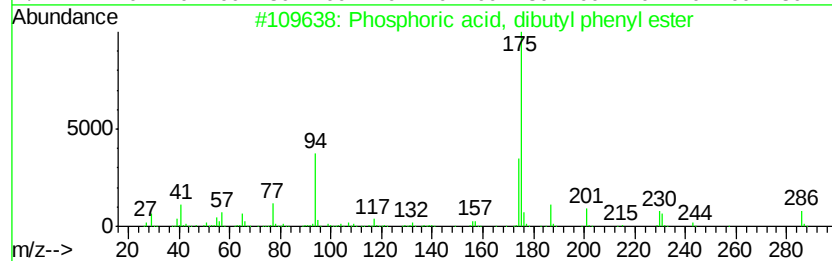
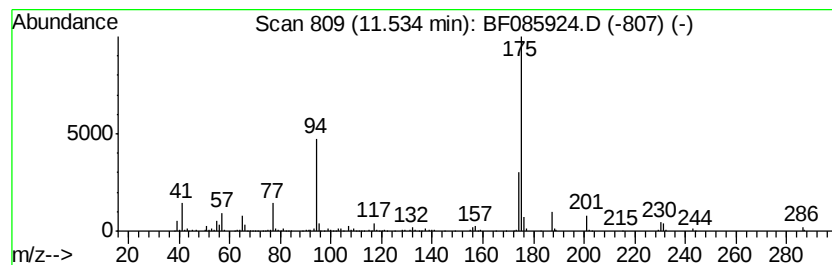
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Phosphoric acid, dibutyl ph... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.53	10.25 ng	539842	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phosphoric acid, dibutyl phenyl ...	286	C14H23O4P	002528-36-1	98
2		4-(Trifluoromethyl)mandelic acid	220	C9H7F3O3	000395-35-7	38
3		m-Aminobenzenesulfonyl fluoride	175	C6H6FN02S	000368-50-3	25
4		4-Methyl-2,7-quinolinediol	175	C10H9NO2	020513-71-7	9
5		5-Fluoro-4-chloro-2-dimethylamin...	175	C6H7ClFN3	1000070-49-0	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033116\
Data File : BF085924.D
Acq On : 31 Mar 2016 13:39
Operator : UM/SJ
Sample : H1954-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-1

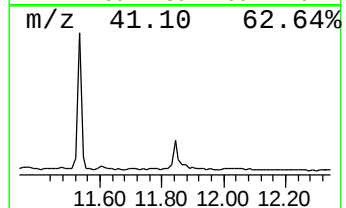
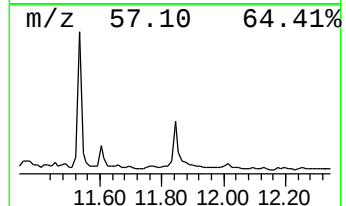
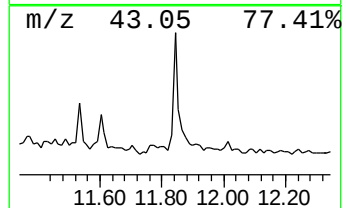
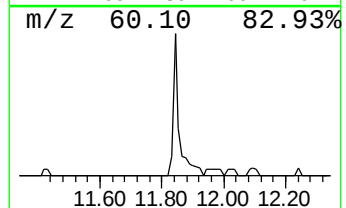
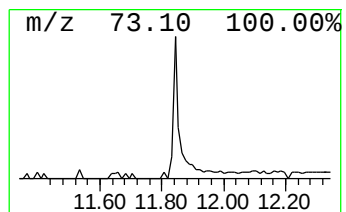
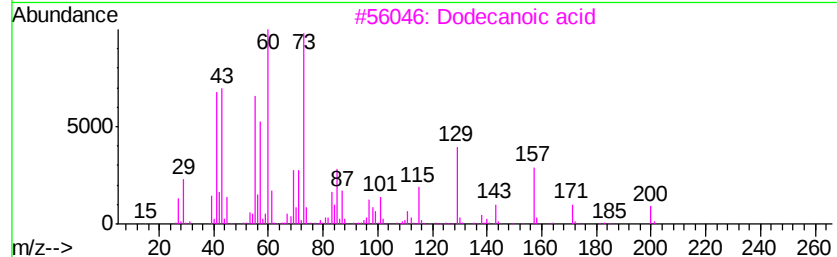
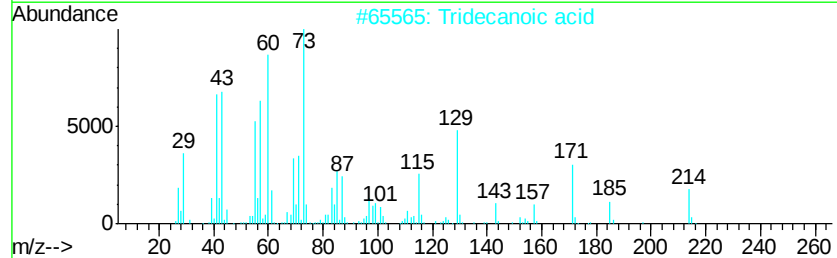
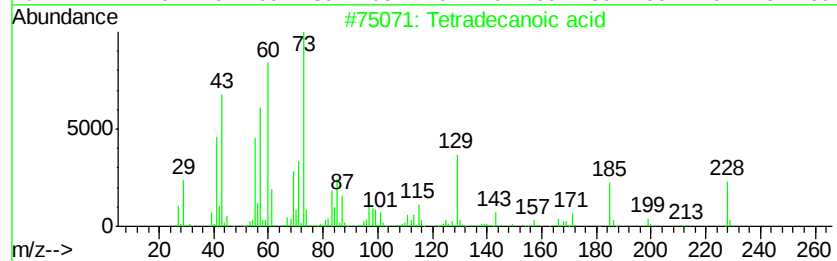
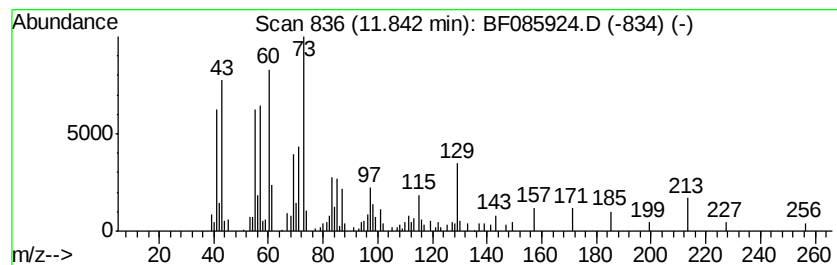
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Tetradecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.84	2.58 ng	135954	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecanoic acid	228	C14H28O2	000544-63-8	80
2		Tridecanoic acid	214	C13H26O2	000638-53-9	72
3		Dodecanoic acid	200	C12H24O2	000143-07-7	58
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	52
5		.beta.-D-Mannofuranoside, 1-O-(1...	332	C17H32O6	1000155-29-9	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033116\
 Data File : BF085924.D
 Acq On : 31 Mar 2016 13:39
 Operator : UM/SJ
 Sample : H1954-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Butanol, 2,3-di...	2.85	6.9 ng		241837	1	6.79	700282	20.0
2-Pentanol, 2,4-d...	4.20	4.7 ng		162848	1	6.79	700282	20.0
1,2-Ethanediol, d...	4.90	2.9 ng		101369	1	6.79	700282	20.0
unknown5.85	5.85	2.3 ng		79639	1	6.79	700282	20.0
1H-Pyrazole, 4,5-...	6.86	3.1 ng		110212	1	6.79	700282	20.0
Hexanoic acid, 2,...	7.04	12.4 ng		432416	1	6.79	700282	20.0
unknown7.65	7.65	3.4 ng		161734	2	8.07	949801	20.0
Benzene, 2-etheny...	7.81	3.4 ng		160157	2	8.07	949801	20.0
1-Adamantanol	8.61	2.0 ng		96991	2	8.07	949801	20.0
Benzoic acid, 2,4...	9.35	7.0 ng		377755	3	9.83	1072450	20.0
Tributyl phosphate	10.47	23.5 ng		1261900	3	9.83	1072450	20.0
Phosphoric acid, ...	11.53	10.3 ng		539842	4	11.32	1053090	20.0
Tetradecanoic acid	11.84	2.6 ng		135954	4	11.32	1053090	20.0