

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072916\
 Data File : BF089369.D
 Acq On : 28 Jul 2016 18:21
 Operator : UM/SJ
 Sample : H4242-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GW-49-10.75-15.0

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-BF072716.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	1.678	7	8	52	rVB	1154551	2491147	100.00%	17.973%
2	4.616	263	265	281	rBV	87341	139423	5.60%	1.006%
3	5.096	305	307	319	rBV	408510	526783	21.15%	3.801%
4	6.182	400	402	410	rBV	236266	414170	16.63%	2.988%
5	6.285	410	411	430	rVB	704800	993048	39.86%	7.165%
6	6.525	430	432	440	rBV	201343	259468	10.42%	1.872%
7	6.673	443	445	448	rBV	737735	858017	34.44%	6.190%
8	7.096	480	482	492	rBV	637747	785140	31.52%	5.665%
9	7.267	492	497	499	rVV	61874	152619	6.13%	1.101%
10	7.313	499	501	513	rVB2	11408	51866	2.08%	0.374%
11	7.610	525	527	537	rBV2	22586	61499	2.47%	0.444%
12	7.747	537	539	543	rBV2	47223	79961	3.21%	0.577%
13	7.805	543	544	547	rBV	240015	343123	13.77%	2.476%
14	8.239	578	582	590	rBV	45375	77772	3.12%	0.561%
15	8.776	626	629	636	rBV	67243	147876	5.94%	1.067%
16	8.890	636	639	649	rVV	1577241	1615009	64.83%	11.652%
17	9.290	672	674	676	rBV	174618	157194	6.31%	1.134%
18	9.325	676	677	692	rVB	432772	601045	24.13%	4.336%
19	9.553	695	697	700	rBV	462132	430333	17.27%	3.105%
20	10.296	759	762	764	rBV2	24894	51226	2.06%	0.370%
21	10.342	764	766	772	rVV	707606	871419	34.98%	6.287%
22	10.479	776	778	781	rBV	28926	40005	1.61%	0.289%
23	10.536	781	783	784	rBV	25580	36973	1.48%	0.267%
24	10.571	784	786	790	rVB3	17633	32335	1.30%	0.233%
25	10.822	806	808	811	rBV	20063	28729	1.15%	0.207%
26	10.914	814	816	820	rVB	185968	166569	6.69%	1.202%
27	10.971	820	821	824	rVB	45744	44495	1.79%	0.321%
28	11.028	824	826	830	rBV	418687	425294	17.07%	3.068%
29	11.588	872	875	884	rVV2	36335	62363	2.50%	0.450%
30	12.365	941	943	947	rBV	21211	28206	1.13%	0.204%
31	12.605	962	964	967	rBV	978524	1167444	46.86%	8.423%
32	13.268	1017	1022	1029	rBV	30747	86031	3.45%	0.621%
33	13.531	1043	1045	1050	rVV	39321	74120	2.98%	0.535%
34	13.645	1052	1055	1062	rVB	302241	339734	13.64%	2.451%

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Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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35	14.983	1169	1172	1178	rbV	156904	219871	8.83%	1.586%
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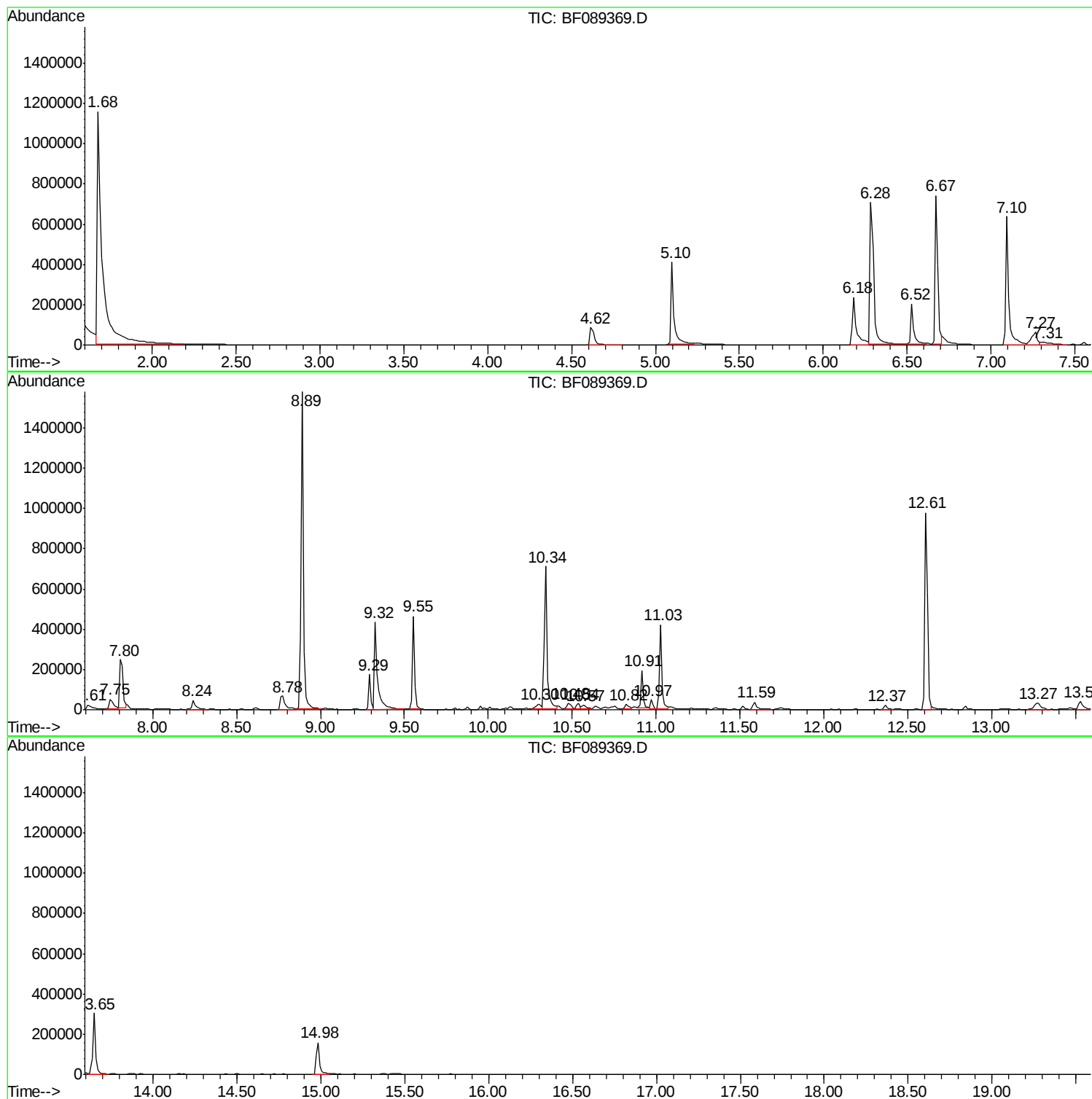
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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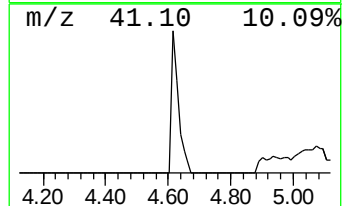
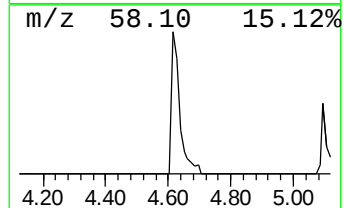
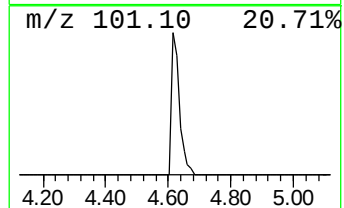
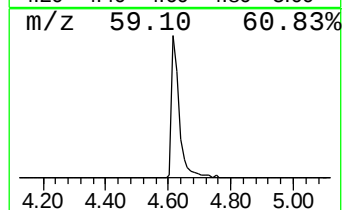
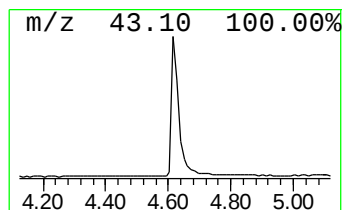
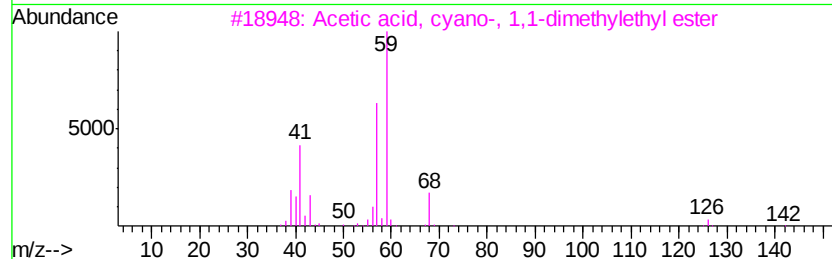
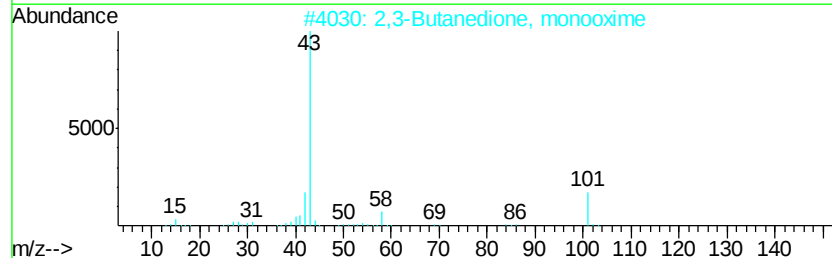
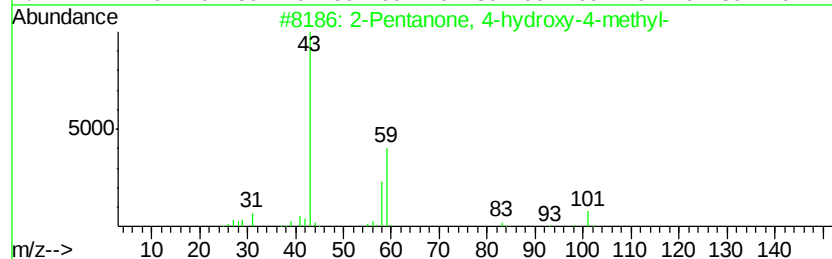
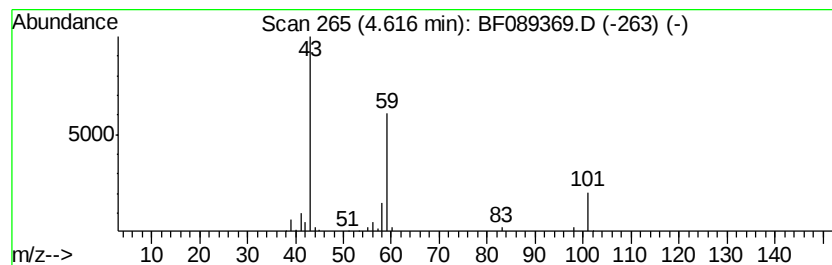
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.62	10.75 ng	139423	1,4-Dichlorobenzene-d4	6.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



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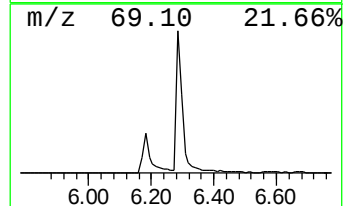
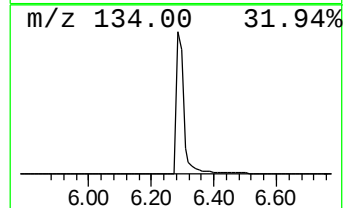
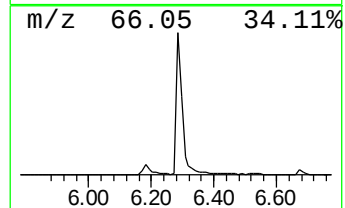
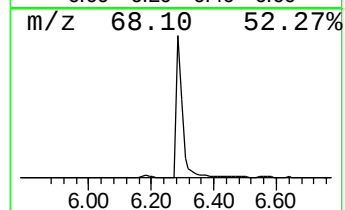
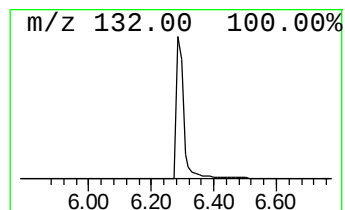
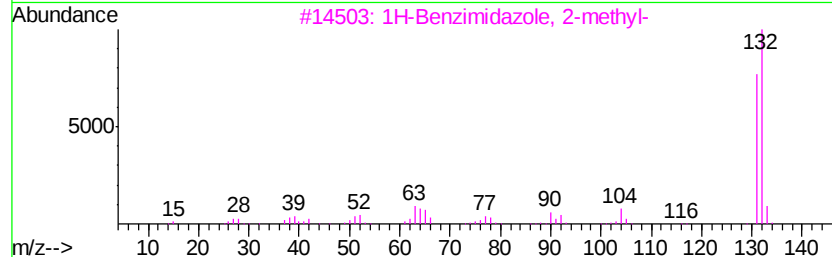
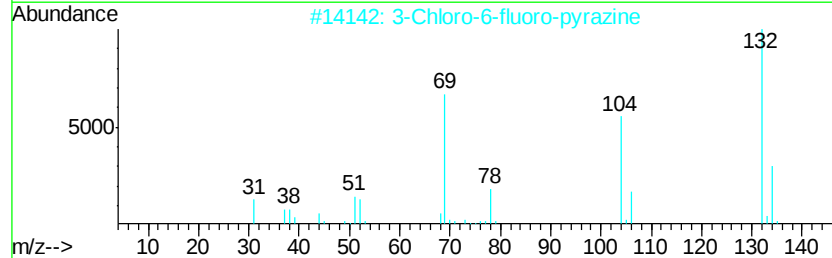
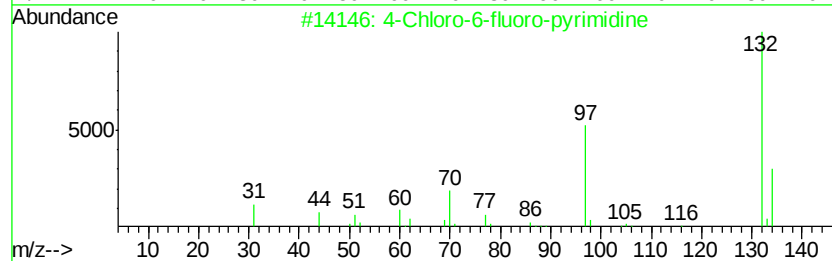
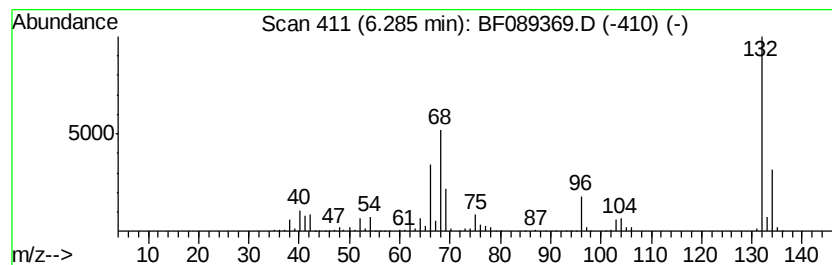
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TIC Library : C:\Database\NIST11.L
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Peak Number 3 unknown6.28 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.28	76.54 ng	993048	1,4-Dichlorobenzene-d4	6.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
5		N-(-3,4-Methylenedioxyphenylisop...	267	C17H17NO2	1000378-96-6	10



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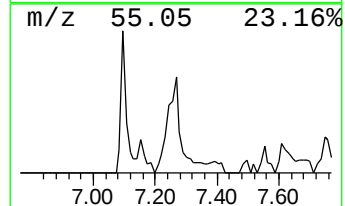
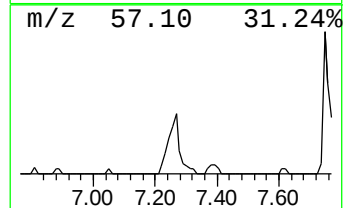
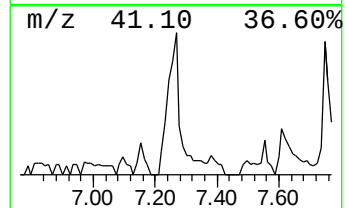
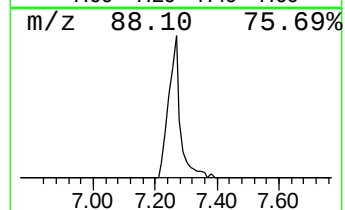
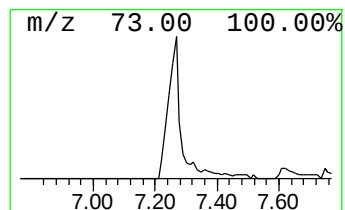
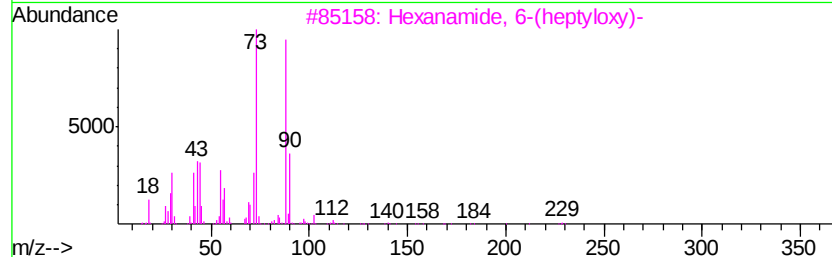
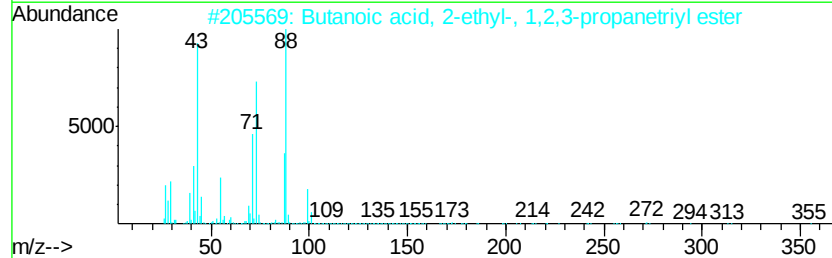
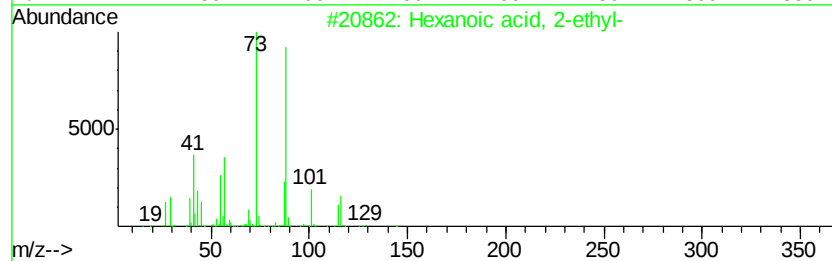
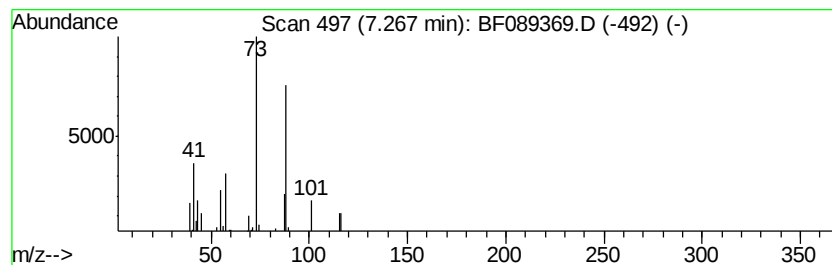
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TIC Library : C:\Database\NIST11.L
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Peak Number 4 Hexanoic acid, 2-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.27	8.90 ng	152619	Naphthalene-d8	7.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78
2		Butanoic acid, 2-ethyl-, 1,2,3-p...	386	C21H38O6	056554-54-2	42
3		Hexanamide, 6-(heptyloxy)-	229	C13H27NO2	055191-08-7	38
4		1,4-Dioxane	88	C4H8O2	000123-91-1	37
5		Heptanoic acid, ethyl ester	158	C9H18O2	000106-30-9	36



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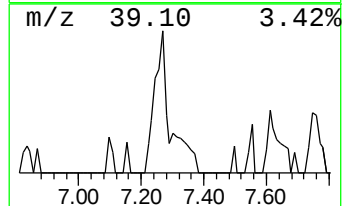
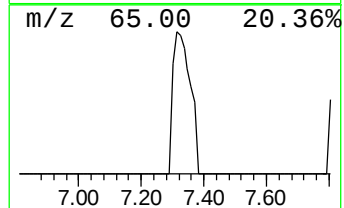
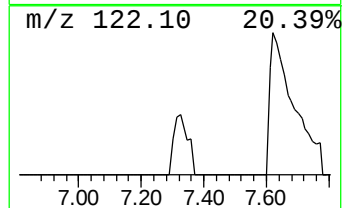
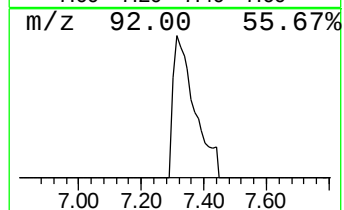
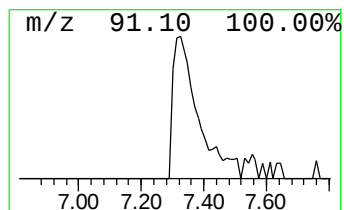
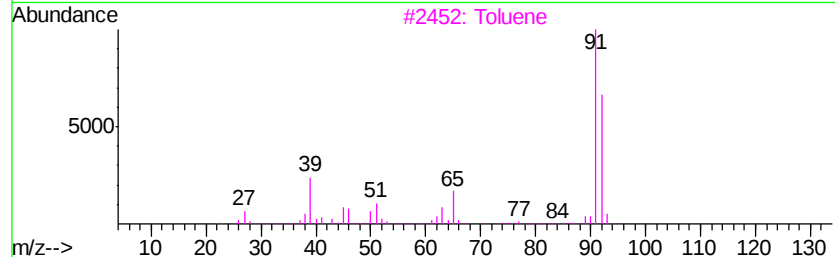
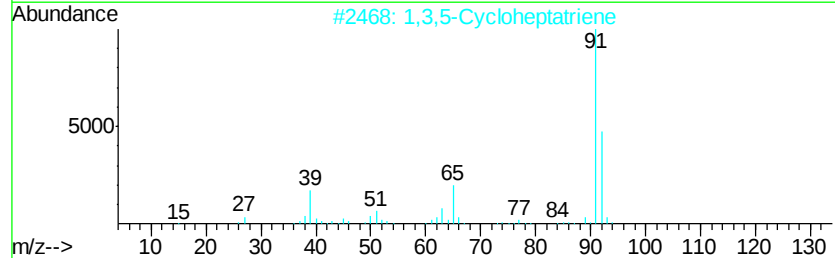
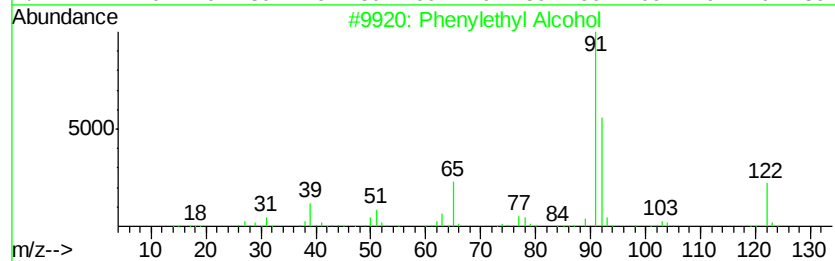
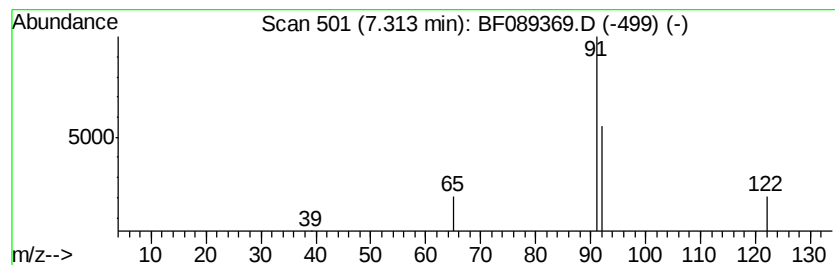
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TIC Library : C:\Database\NIST11.L
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Peak Number 5 Phenylethyl Alcohol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.31	3.02 ng	51866	Napthalene-d8	7.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenylethyl Alcohol	122	C8H10O	000060-12-8	74
2		1,3,5-Cycloheptatriene	92	C7H8	000544-25-2	9
3		Toluene	92	C7H8	000108-88-3	7
4		Spiro[2,4]hepta-4,6-diene	92	C7H8	000765-46-8	7
5		Cyclobutene, 2-propenylidene-	92	C7H8	052097-85-5	7



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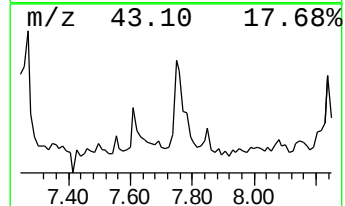
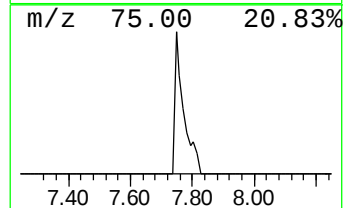
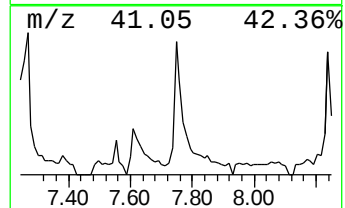
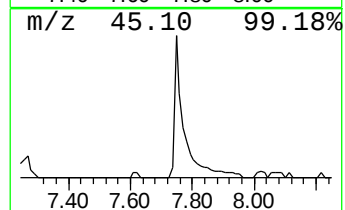
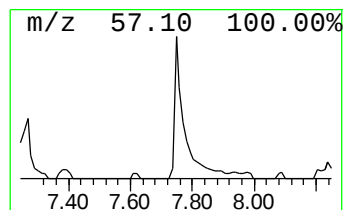
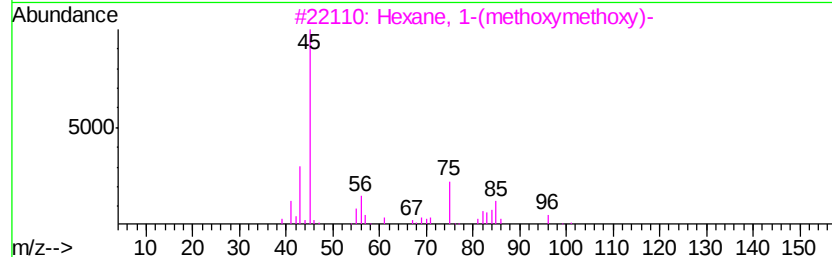
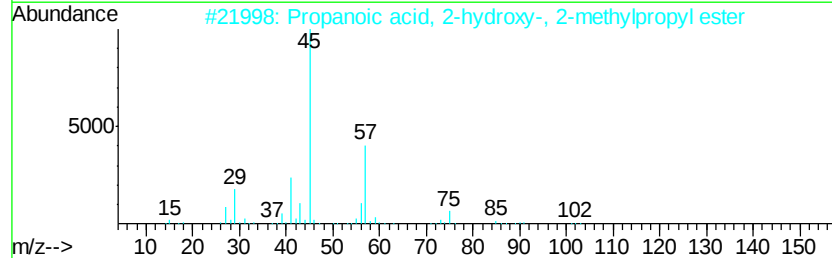
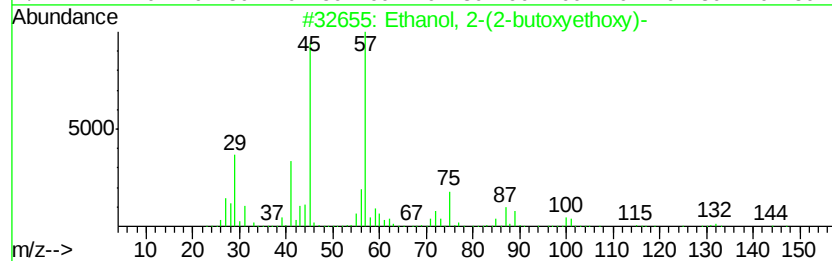
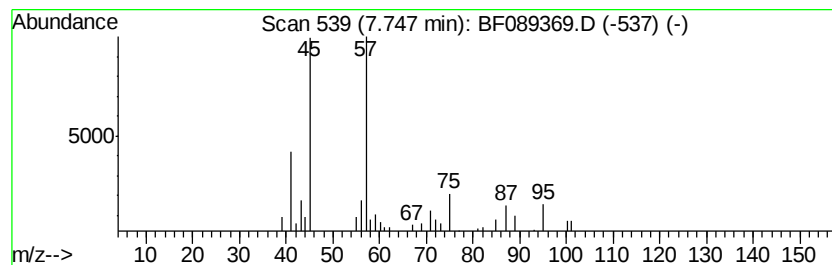
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Peak Number 6 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.75	4.66 ng	79961	Naphthalene-d8	7.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2		Propanoic acid, 2-hydroxy-, 2-me...	146	C7H14O3	000585-24-0	59
3		Hexane, 1-(methoxymethoxy)-	146	C8H18O2	066675-06-7	47
4		Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	42
5		Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	38



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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
GW-49-10.75-15.0

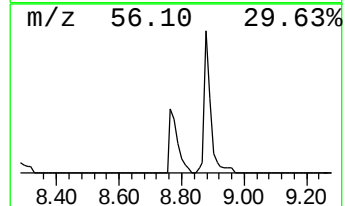
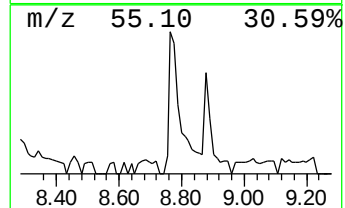
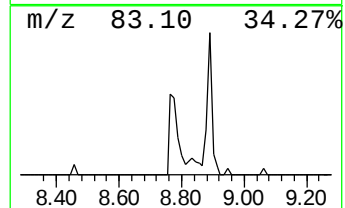
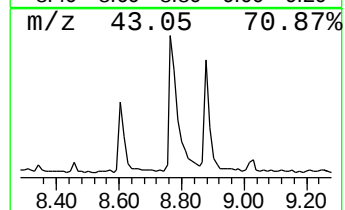
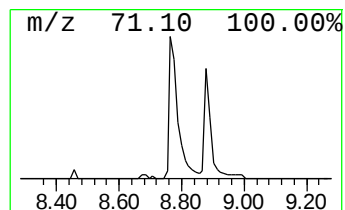
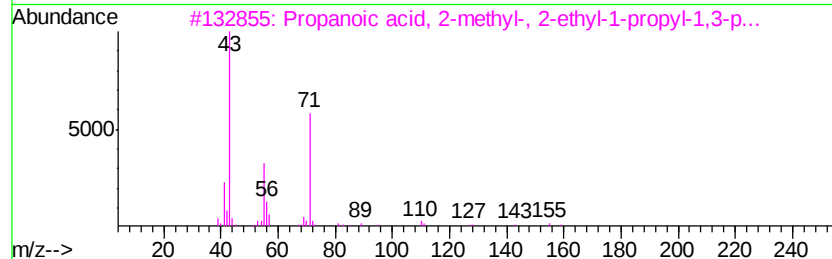
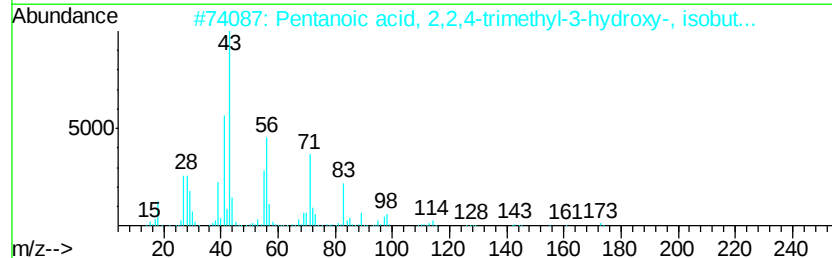
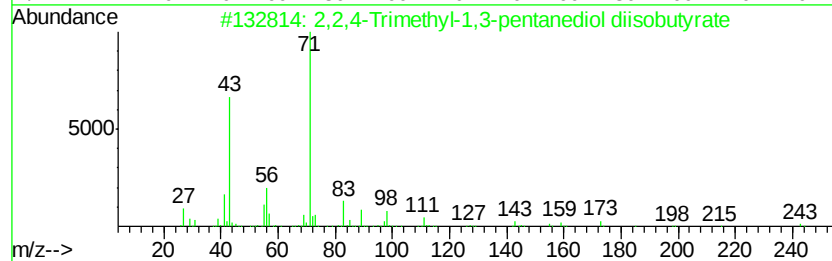
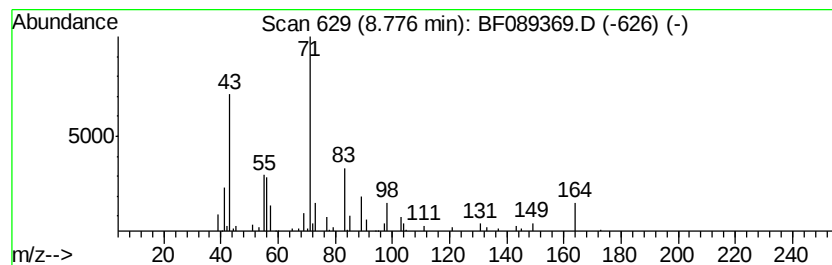
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 2,2,4-Trimethyl-1,3-pentane... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.78	6.87 ng	147876	Acenaphthene-d10	9.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2,4-Trimethyl-1,3-pentanediol ...	286	C16H30O4	006846-50-0	53
2		Pentanoic acid, 2,2,4-trimethyl-...	216	C12H24O3	244074-78-0	40
3		Propanoic acid, 2-methyl-, 2-eth...	286	C16H30O4	074367-30-9	38
4		6-Methyl-hept-2-en-4-ol	128	C8H16O	083212-30-0	35
5		Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	32



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072916\
Data File : BF089369.D
Acq On : 28 Jul 2016 18:21
Operator : UM/SJ
Sample : H4242-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

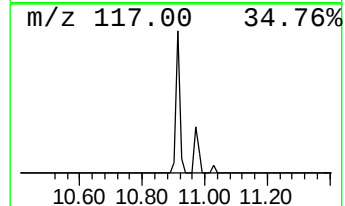
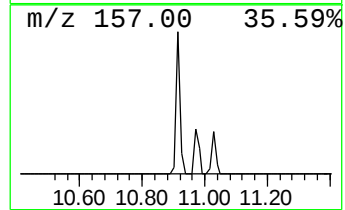
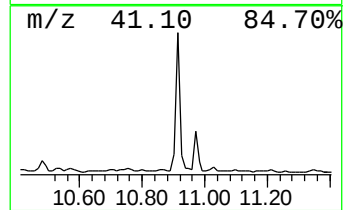
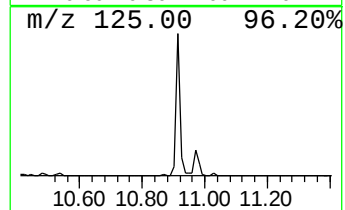
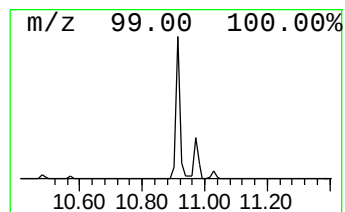
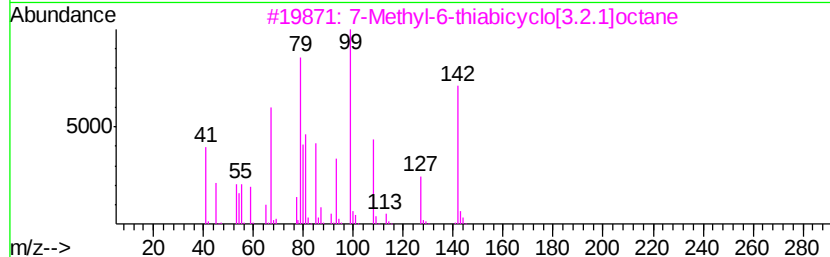
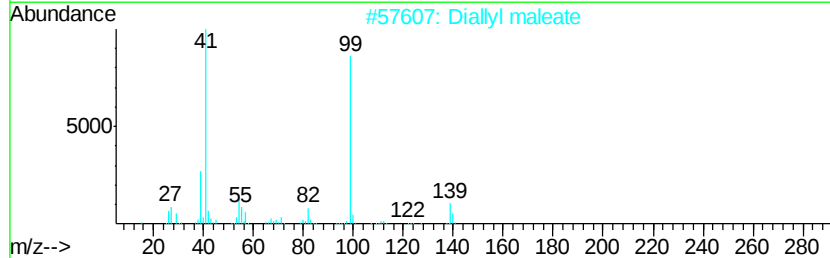
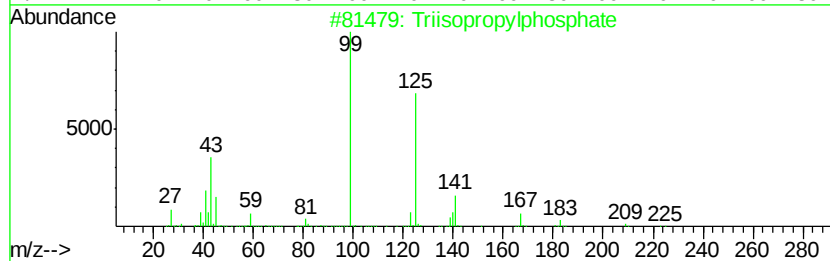
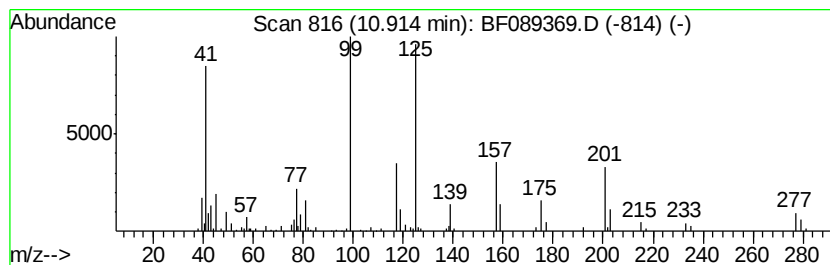
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ClientSampleId :
GW-49-10.75-15.0

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 unknown10.91 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.91	7.83 ng	166569	Phenanthrene-d10	11.03	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triisopropylphosphate	224	C9H21O4P	000513-02-0	28
2	Diallyl maleate	196	C10H12O4	000999-21-3	14
3	7-Methyl-6-thiabicyclo[3.2.1]octane	142	C8H14S	1000306-33-2	14
4	4-Chloro-2-nitrobenzoic acid	201	C7H4ClNO4	006280-88-2	10
5	Phosphonofluoridic acid, methyl-...	224	C10H22F02P	211192-74-4	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072916\
Data File : BF089369.D
Acq On : 28 Jul 2016 18:21
Operator : UM/SJ
Sample : H4242-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
GW-49-10.75-15.0

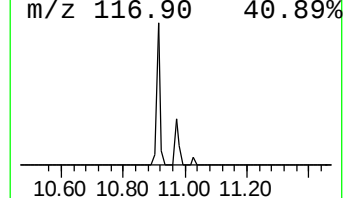
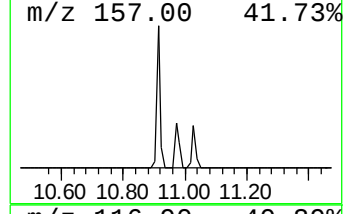
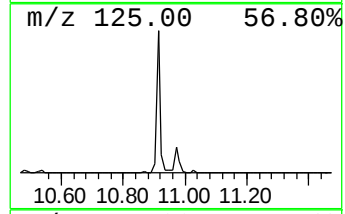
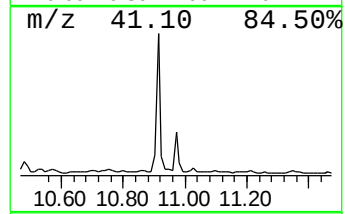
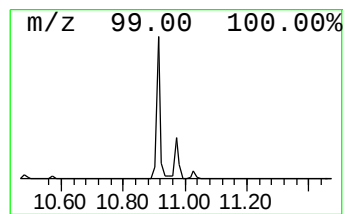
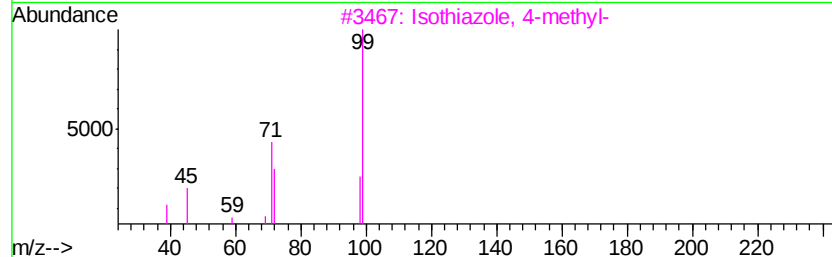
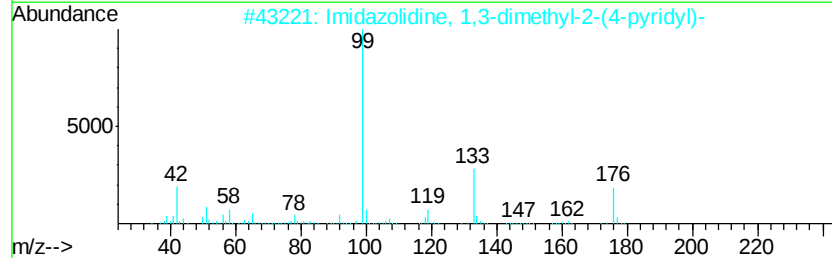
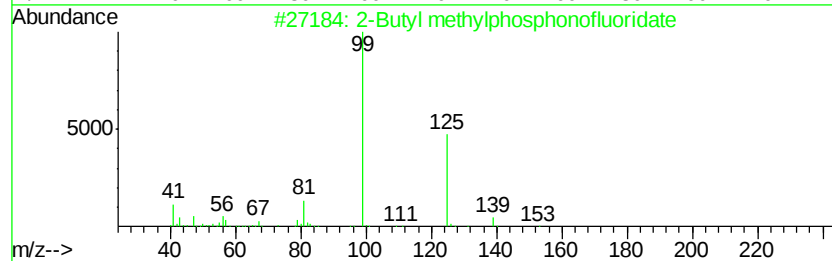
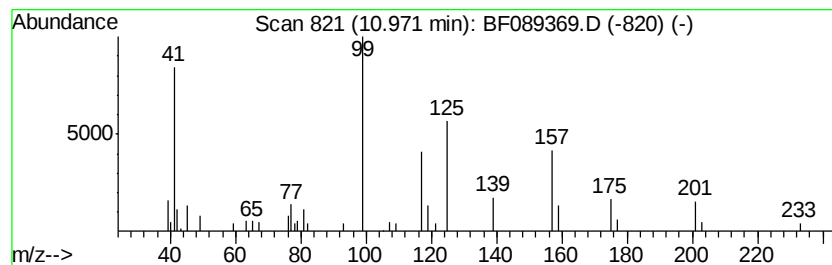
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown10.97 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.97	2.09 ng	44495	Phenanthrene-d10	11.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Butyl methylphosphonofluoridate	154	C5H12F02P	000352-52-3	23
2		Imidazolidine, 1,3-dimethyl-2-(4...	177	C10H15N3	303187-78-2	10
3		Isothiazole, 4-methyl-	99	C4H5NS	000693-90-3	9
4		Methyl 3-amino-2-thiophenecarbox...	157	C6H7N02S	022288-78-4	9
5		Cholestan-22(26)-isoepoxy	386	C27H46O	1000252-59-7	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072916\
Data File : BF089369.D
Acq On : 28 Jul 2016 18:21
Operator : UM/SJ
Sample : H4242-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
GW-49-10.75-15.0

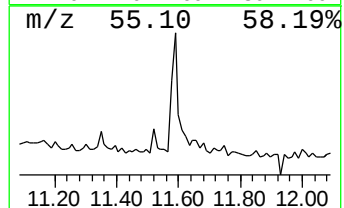
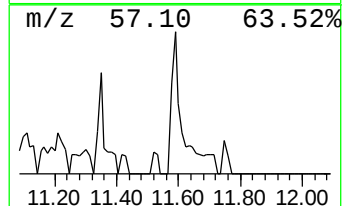
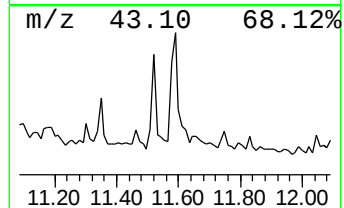
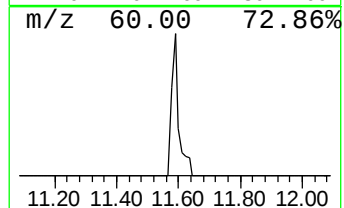
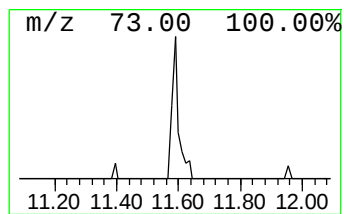
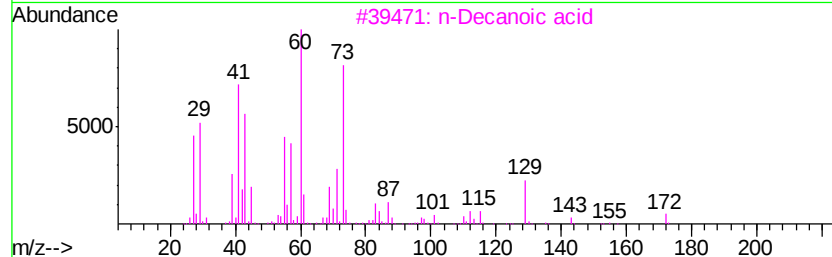
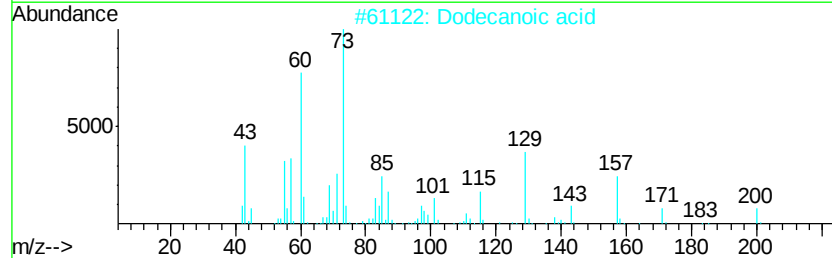
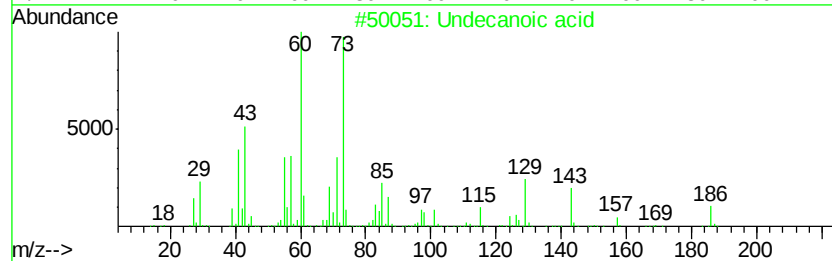
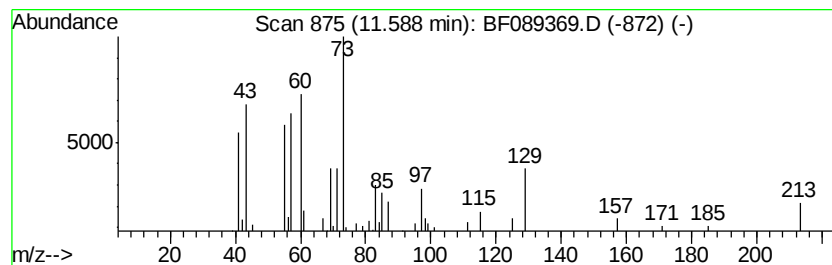
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Undecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.59	2.93 ng	62363	Phenanthrene-d10	11.03

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecanoic acid	186	C11H22O2	000112-37-8	53
2			Dodecanoic acid	200	C12H24O2	000143-07-7	50
3			n-Decanoic acid	172	C10H20O2	000334-48-5	38
4			4-Pentadecanol	228	C15H32O	109212-91-1	22
5			.beta.-D-Mannofuranoside, 1-O-(1...	332	C17H32O6	1000155-29-9	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072916\
Data File : BF089369.D
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Operator : UM/SJ
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Misc :
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Instrument :
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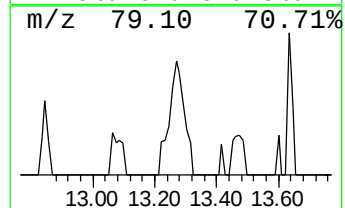
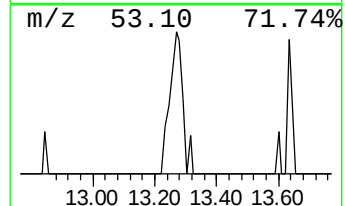
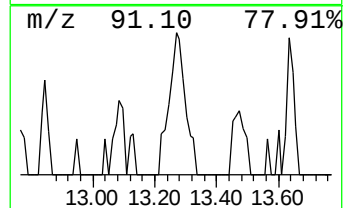
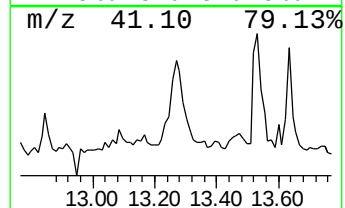
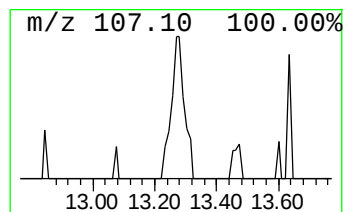
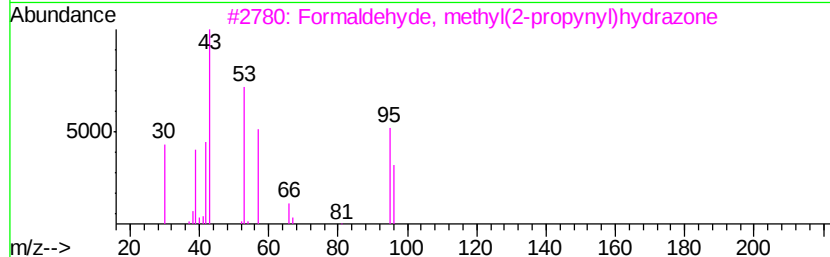
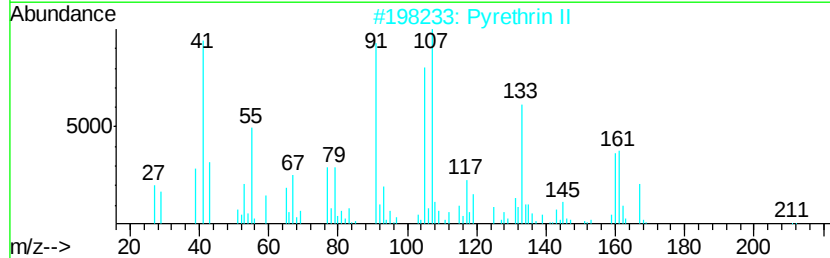
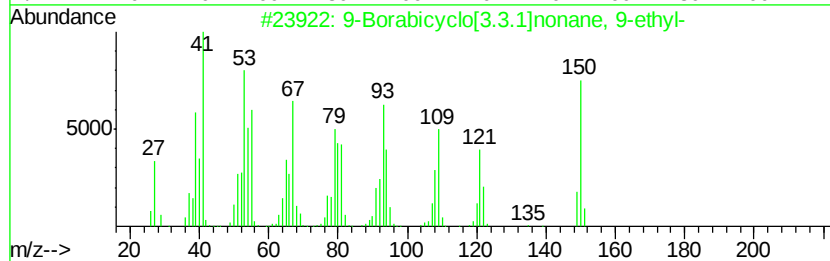
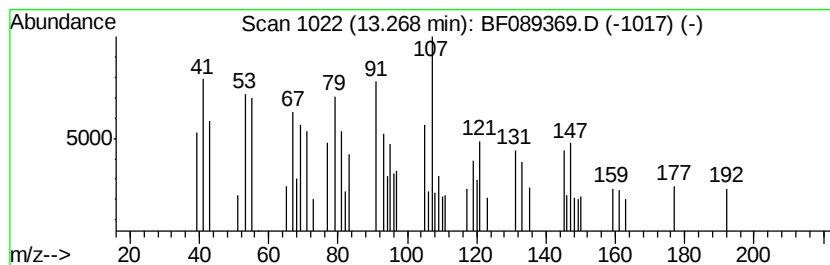
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 unknown13.27 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.27	5.06 ng	86031	Chrysene-d12	13.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Borabicyclo[3.3.1]nonane, 9-et...	150	C10H19B	052102-17-7	35
2		Pyrethrin II	372	C22H28O5	000121-29-9	25
3		Formaldehyde, methyl(2-propynyl)...	96	C5H8N2	110213-10-0	25
4		3-Octyne	110	C8H14	015232-76-5	22
5		4,5-Cyclopentene, 1-oxa-2-phosph...	146	C6H11O2P	032135-32-3	18



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072916\
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Sample : H4242-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

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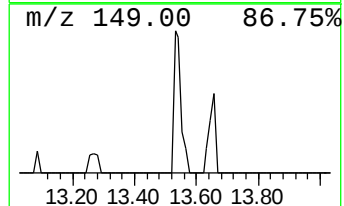
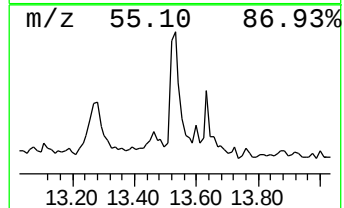
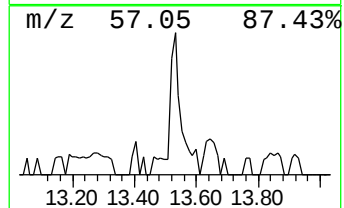
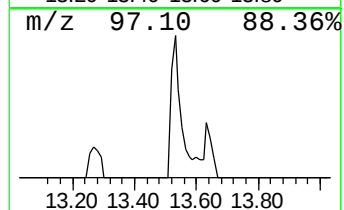
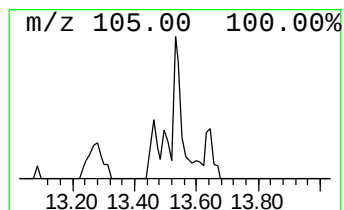
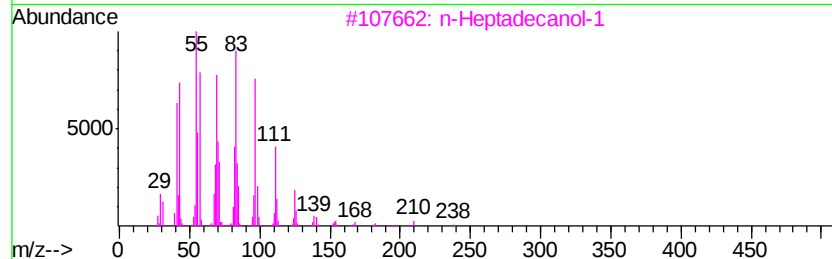
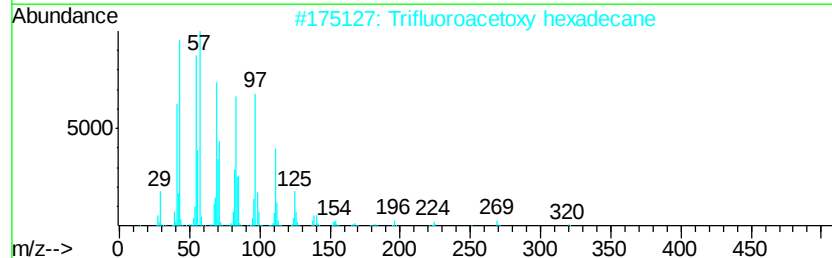
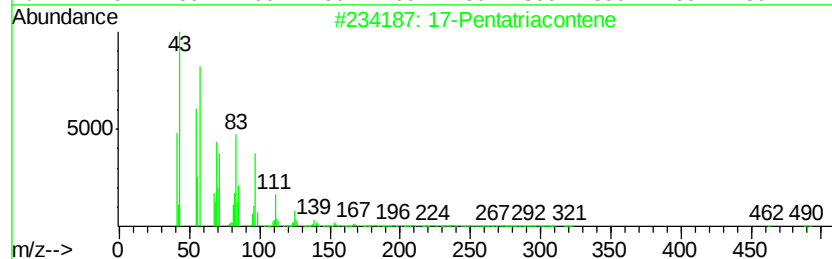
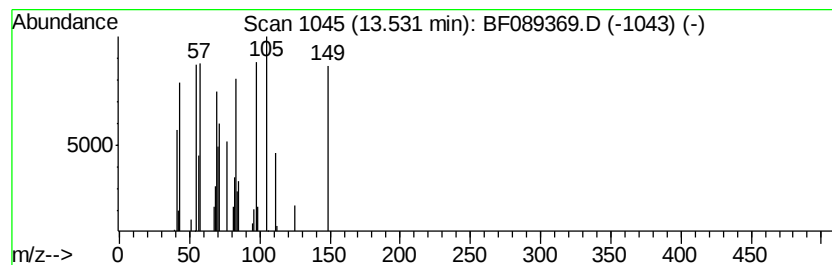
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 17-Pentatriacontene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.53	4.36 ng	74120	Chrysene-d12	13.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	17-Pentatriacontene	491	C35H70	006971-40-0	81
2		Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	81
3		n-Heptadecanol-1	256	C17H36O	001454-85-9	76
4		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	76
5		3-Eicosene, (E)-	280	C20H40	074685-33-9	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072916\
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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
GW-49-10.75-15.0

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.62	10.8	ng	139423	1	6.52	259468	20.0
unknown6.28	6.28	76.5	ng	993048	1	6.52	259468	20.0
Hexanoic acid, 2-...	7.27	8.9	ng	152619	2	7.82	343123	20.0
Phenylethyl Alcohol	7.31	3.0	ng	51866	2	7.82	343123	20.0
Ethanol, 2-(2-but...	7.75	4.7	ng	79961	2	7.82	343123	20.0
2,2,4-Trimethyl-1...	8.78	6.9	ng	147876	3	9.55	430333	20.0
unknown10.91	10.91	7.8	ng	166569	4	11.03	425294	20.0
unknown10.97	10.97	2.1	ng	44495	4	11.03	425294	20.0
Undecanoic acid	11.59	2.9	ng	62363	4	11.03	425294	20.0
unknown13.27	13.27	5.1	ng	86031	5	13.65	339734	20.0
17-Pentatriacontene	13.53	4.4	ng	74120	5	13.65	339734	20.0